A Level Chemistry A
H432/02 Synthesis and analytical techniques
Sample Question Paper

Date – Morning/Afternoon
Time allowed: 2 hours 15 minutes

You must have:
• the Data Sheet for Chemistry A
You may use:
• a scientific or graphical calculator

INSTRUCTIONS
• Use black ink. You may use an HB pencil for graphs and diagrams.
• Complete the boxes above with your name, centre number and candidate number.
• Answer all the questions.
• Write your answer to each question in the space provided.
• Additional paper may be used if required but you must clearly show your candidate number, centre number and question number(s).
• Do not write in the bar codes.

INFORMATION
• The total mark for this paper is 100.
• The marks for each question are shown in brackets [ ].
• Quality of extended responses will be assessed in questions marked with an asterisk (*).
• This document consists of 36 pages.
1. The displayed formula of an organic compound is shown below.

What is the systematic name of this organic compound?

A  Propyl propanoate  
B  Propyl butanoate  
C  Butyl propanoate  
D  Butyl butanoate

Your answer:  

[1]
Ethanol is oxidised to ethanoic acid using acidified potassium dichromate(IV) solution. The reaction is heated under reflux using the equipment shown in the diagram below.

What is the reason for heating under reflux?

A  to ensure even heating  
B  to prevent any substances escaping  
C  to boil the mixture at a higher temperature  
D  to allow efficient mixing  

Your answer  

How many stereoisomers are there of \( CH_3CH=CHCH(OH)CH_2CH=CH_2 \)?

A  2  
B  4  
C  6  
D  8  

Your answer  

[1]
The functional group in an organic compound, \( W \), was identified by carrying out two chemical tests. The results of the tests are shown below.

<table>
<thead>
<tr>
<th>Heating with acidified sodium dichromate(VI)(aq)</th>
<th>Addition of 2,4-dinitrophenylhydrazine(aq)</th>
</tr>
</thead>
<tbody>
<tr>
<td>orange solution turns green</td>
<td>yellow/orange precipitate formed</td>
</tr>
</tbody>
</table>

Which compound could be \( W \)?

A. \( \text{CH}_3\text{CH}_2\text{CH}_2\text{OH} \)
B. \( \text{CH}_3\text{COCH}_3 \)
C. \( \text{CH}_3\text{CH(OH)}\text{CH}_3 \)
D. \( \text{CH}_3\text{CH}_2\text{CHO} \)

Your answer [ ]

Complete combustion of 40 cm\(^3\) of a gaseous hydrocarbon \( X \) requires 240 cm\(^3\) of oxygen. 160 cm\(^3\) of carbon dioxide forms. All gas volumes are at room temperature and pressure.

What is the formula of \( X \)?

A. \( \text{C}_4\text{H}_8 \)
B. \( \text{C}_4\text{H}_{10} \)
C. \( \text{C}_6\text{H}_{12} \)
D. \( \text{C}_6\text{H}_{14} \)

Your answer [ ]
6 The boiling point of butan-1-ol is 118 °C. The boiling point of 2-methylpropan-2-ol is 82 °C.

Why is the boiling point of butan-1-ol higher than that of 2-methylpropan-2-ol?

A butan-1-ol has stronger induced dipole–dipole interactions because it has more electrons

B butan-1-ol has stronger induced dipole–dipole interactions because it has a straight-chain structure

C butan-1-ol can form hydrogen bonds while 2-methylpropan-2-ol cannot

D butan-1-ol is more stable because it is a primary alcohol

Your answer [ ]

7 Hydrogen bromide reacts with 3-methylbut-1-ene.

What is the structure of the major intermediate formed in the mechanism?

A

B

C

D

Your answer [ ]
8 Two chemical tests are carried out on an aqueous solution of an aromatic organic compound Y.

The results of the tests are shown below.

<table>
<thead>
<tr>
<th>Test</th>
<th>Br₂(aq)</th>
<th>Na₂CO₃(aq)</th>
</tr>
</thead>
<tbody>
<tr>
<td>Observation</td>
<td>decolourised</td>
<td>effervescence</td>
</tr>
</tbody>
</table>

What is the minimum number of C atoms in Y?

A  6
B  7
C  8
D  9

Your answer  

9 Bromine is reacted separately with nitrobenzene and phenylamine.

Which organic products are likely to form?

<table>
<thead>
<tr>
<th>Product from nitrobenzene</th>
<th>Product from phenylamine</th>
</tr>
</thead>
<tbody>
<tr>
<td>A  2-bromonitrobenzene</td>
<td>2-bromophenylamine</td>
</tr>
<tr>
<td>B  2-bromonitrobenzene</td>
<td>3-bromophenylamine</td>
</tr>
<tr>
<td>C  3-bromonitrobenzene</td>
<td>2-bromophenylamine</td>
</tr>
<tr>
<td>D  3-bromonitrobenzene</td>
<td>3-bromophenylamine</td>
</tr>
</tbody>
</table>

Your answer  

[1]
10 Which alcohol could be used to prepare HCOOCH(CH₃)₂?

A  Propan-1-ol
B  Propan-2-ol
C  2-Methylpropan-2-ol
D  Methanol

Your answer  

11 CN⁻ ions react with haloalkanes and with carbonyl compounds.

Which row gives the correct mechanisms for the reactions?

<table>
<thead>
<tr>
<th></th>
<th>Reaction of CN⁻ with haloalkanes</th>
<th>Reaction of CN⁻ with carbonyl compounds</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>Electrophilic substitution</td>
<td>Electrophilic addition</td>
</tr>
<tr>
<td>B</td>
<td>Electrophilic substitution</td>
<td>Nucleophilic addition</td>
</tr>
<tr>
<td>C</td>
<td>Nucleophilic substitution</td>
<td>Electrophilic addition</td>
</tr>
<tr>
<td>D</td>
<td>Nucleophilic substitution</td>
<td>Nucleophilic addition</td>
</tr>
</tbody>
</table>

Your answer  

[1]
The structure of a molecule that is used as a pain reliever is shown below.

Which statement about this molecule is **not** true?

A  It has the molecular formula $\text{C}_8\text{H}_9\text{NO}_2$.

B  It reacts with bases to form salts.

C  It has a ketone functional group.

D  It can be hydrolysed with aqueous acid.

Your answer  

[1]
13 Carbonyl compounds have distinctive smells. Menthone smells of peppermint.

Menthone is reacted in a two-step synthesis shown below.

**Step 1**: A sample of menthone is added to hot acidified aqueous dichromate(VI) ions.

**Step 2**: The resulting mixture from **Step 1** is added to NaBH₄ in water.

What happens to the smell of the reaction mixture during the process?

<table>
<thead>
<tr>
<th></th>
<th>Step 1</th>
<th>Step 2</th>
</tr>
</thead>
<tbody>
<tr>
<td>A</td>
<td>Smell of peppermint remains</td>
<td>Smell of peppermint is lost</td>
</tr>
<tr>
<td>B</td>
<td>Smell of peppermint is lost</td>
<td>Smell of peppermint returns</td>
</tr>
<tr>
<td>C</td>
<td>Smell of peppermint remains</td>
<td>Smell of peppermint remains</td>
</tr>
<tr>
<td>D</td>
<td>Smell of peppermint is lost</td>
<td>Smell of peppermint does not return</td>
</tr>
</tbody>
</table>

Your answer [ ]

[1]
14 Which of the following support(s) the delocalised model for benzene rather than the Kekulé model?

1: Benzene is less reactive than cyclohexene

2: A benzene molecule has a planar, hexagonal structure

3: The enthalpy change of hydrogenation of benzene is more exothermic than predicted from the Kekulé structure

A 1, 2 and 3

B Only 1 and 2

C Only 2 and 3

D Only 1

Your answer [ ]
The structure of molecule Z is shown below.

Which of the following statements is/are true?

1: The carbon-13 NMR spectrum of Z shows four peaks
2: The proton NMR spectrum of Z shows five peaks
3: The proton NMR spectrum of Z run in D$_2$O shows three peaks

A 1, 2 and 3
B Only 1 and 2
C Only 2 and 3
D Only 1

Your answer [ ]
SECTION B

Answer all the questions.

16 Methyl allyl chloride, MAC, is a chemical used in the production of insecticides. The structure of MAC is shown below.

![MAC structure](image)

MAC

(a) (i) Give the molecular formula of MAC.

........................................................................................................................................................................ [1]

(ii) Draw the skeletal formula of MAC.

........................................................................................................................................................................ [1]

(iii) MAC has several structural isomers.

State what is meant by structural isomers.

........................................................................................................................................................................ [1]

(b) MAC is highly flammable. When MAC burns, one of the products formed is a toxic gas.

1.321 g of this gas occupies 1.053 dm³ at 100 kPa and 350 K.

Use the information provided to suggest the identity of the gas.

gas = .................. [4]
(c) The flowchart below shows some reactions of MAC.

![Flowchart Diagram]

(i) Complete the flowchart above.
- Draw the structure of the product of Reaction 1.
- Draw the structure of the minor organic product of Reaction 2 (Compound B).

(ii) Reaction 2 creates a mixture of compounds. Compound A is the major product.

Draw the mechanism for the formation of compound A.

Use curly arrows and show relevant dipoles.
(iii) Explain why compound B is the minor product of Reaction 2.

..................................................................................................................................................  [1]

..................................................................................................................................................

(iv) MAC reacts with water in the presence of AgNO₃(aq) and ethanol.

Draw the structure of the organic product of this reaction.

State what you would observe in this reaction and identify the compound responsible for the observation.

..................................................................................................................................................

..................................................................................................................................................  [2]
(d) Compound A reacts slowly in humid conditions to form compound C.

![Chemical structure of Compound A and Compound C](image)

Compound C contained the following percentage composition by mass: C, 46.1%; H, 7.7%; O, 46.2%

The infrared spectrum of compound C is shown below.
Using the information on the previous page, deduce the structure of compound C.

Give your reasoning.

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structure = [5]
Molecules with more than one functional group are useful chemical ‘building blocks’.

(a) Compound D, CH₃CH(OH)CH₂NH₂, is an intermediate in the synthesis of a variety of drugs.

(i) Compound D can be synthesised from ethanal, CH₃CHO.

Devise a **two-step** synthesis of compound D from ethanal.

- Give details of appropriate reagents and relevant conditions.
- Write an equation for each step, showing clearly all organic compounds.

(ii) Explain why compound D is very soluble in water.

Use a diagram in your answer.
(iii) Compound D reacts with propanedioic acid, HOOCCH$_2$COOH, to form a condensation polymer.

Draw a possible repeat unit of this condensation polymer.

Show clearly any functional group present in the repeat unit.

(b) Serine, shown below, is an amino acid.

(i) Use electron repulsion theory to predict the shape of the bonds around atoms A and B.

Give relevant bond angles around atoms A and B.

Give reasons for your answers.

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........................................................................................................................................................................... [4]
(ii) A student adds an excess of aqueous sodium hydroxide to a sample of solid serine. The student then purifies the resulting reaction mixture to obtain a pure sample of an ionic organic product.

- Draw the structure of the ionic organic compound obtained.
- Outline the steps that the student could carry out to obtain a pure sample of the organic product from the reaction mixture.

(c) Tabtoxin is a poisonous substance produced by bacteria found in lilac trees.

(i) Identify the chiral centres present in a molecule of tabtoxin.

On the structure above, mark each chiral centre with an asterisk, *.

(ii) Tabtoxin can be broken down by alkaline hydrolysis.

Draw the structures of all the organic products of the alkaline hydrolysis of tabtoxin.
18 A student investigates reactions of aromatic compounds.

(a) The student first carries out the reaction shown below.

\[
\text{quinol} + \text{Compound E} + \text{HCl}
\]

(i) The student obtains a very low yield of compound E.

The student obtains a much higher yield of a different organic product with molecular formula \(C_{14}H_{22}O_2\).

Suggest an identity for the organic product \(C_{14}H_{22}O_2\) and draw its structure below.

(ii) The student is told by a friend that the FeCl₃ catalyst is not needed because quinol is more reactive than benzene.

Explain why the student’s friend is correct.

You may draw a diagram to support your answer.
(b) 4-Nitrobenzoic acid is an important compound in chemical synthesis. The flowchart below shows a synthesis involving 4-nitrobenzoic acid.

![Flowchart](image)

**Compound F**

(i) State suitable reactant(s) and conditions for step 1.

(ii) In step 2, the –NO\(_2\) group in compound F is reduced by tin and concentrated hydrochloric acid.

Write an equation for the reduction of compound F.

Show the structures of any organic compounds involved.
α-Hydroxy acids (AHAs) are naturally occurring acids often used as cosmetics.

(a) The flowchart below shows some reactions of an AHA, \( \text{CH}_3(\text{CH}_2)_2\text{CHOHCOOH} \).

(i) Fill in the boxes to show the organic products of Reactions 1 and 2, clearly showing the relevant functional groups.

(ii) Give the full systematic name for compound G.

.................................................................................................................................................................................. [1]
(iii) Compound H is a stereoisomer of compound G.

- Suggest a structure for compound H.
- Draw the repeat unit of the addition polymer that can be formed from compound H.

![Compound H](image)

![addition polymer](image)

(iv) The addition polymer in (iii) is used widely in industry. Increasingly, waste polymers are being processed as a more sustainable option than disposal.

Apart from recycling, state two methods for usefully processing waste polymers.

[2]

.........................................................................................................................................................

.........................................................................................................................................................  [2]
(b) A student synthesises a sample of the AHA J using the following reaction scheme, starting from propane-1,2-diol.

\[
\text{propane-1,2-diol} \xrightarrow{[O]} \text{Compound J} \quad \text{Step 1}
\]

\[
\text{Compound J} \xrightarrow{\text{NaBH}_4} \quad \text{Step 2}
\]

(i) In the space below:
- state a suitable oxidising agent for Step 1
- write an equation for Step 1
- outline the mechanism for Step 2, showing curly arrows and relevant dipoles.
(ii) The reagent used in Step 2 of the synthesis in (i) was NaBH₄. NaBH₄ contains the ions Na⁺ and [BH₄]⁻.

Draw a ‘dot-and-cross’ diagram of NaBH₄ and give the full electron configuration of Na⁺.

Show outer shells of electrons only.

**full** electronic configuration of Na⁺: .................................................................

[2]
(c) Compound K is an AHA that is often used in ‘chemical face peels’.

A student wishes to identify compound K from the list of compounds below.

<table>
<thead>
<tr>
<th>Compound</th>
<th>Structure</th>
</tr>
</thead>
<tbody>
<tr>
<td>glycolic acid</td>
<td>HOCH₂COOH</td>
</tr>
<tr>
<td>malic acid</td>
<td>HOOCCH₂CHOHCOOH</td>
</tr>
<tr>
<td>mandelic acid</td>
<td>C₆H₅CHOHCOOH</td>
</tr>
<tr>
<td>pantoic acid</td>
<td>HOCH₂C(CH₃)₂CHOHCOOH</td>
</tr>
</tbody>
</table>

The student isolates compound K and analyses a sample of the compound by titration.

The student dissolves 1.89 g of compound K in water and makes the solution up to 250.0 cm³ in a volumetric flask. The student titrates 25.0 cm³ of this solution with 0.150 mol dm⁻³ NaOH(aq).

18.80 cm³ of NaOH(aq) were required for complete neutralisation.

Use the results of the student’s analysis to identify compound K from the list above.

Show all of your working.

\[
K = \text{..................................................} \quad [5]
\]
Cyclohexanone can be prepared in the laboratory by reacting cyclohexanol with concentrated sulfuric acid and sodium dichromate.

Ethanedioic acid is added to the reaction mixture to react with any excess dichromate.

The mixture is then distilled. The impure distillate is a mixture of cyclohexanone and water.

You will need to refer to some or all of the following data to answer these questions.

<table>
<thead>
<tr>
<th></th>
<th>Boiling point /°C</th>
<th>Density /g cm⁻³</th>
<th>Mr</th>
</tr>
</thead>
<tbody>
<tr>
<td>Cyclohexanol</td>
<td>161</td>
<td>0.962</td>
<td>100.0</td>
</tr>
<tr>
<td>Cyclohexanone</td>
<td>156</td>
<td>0.948</td>
<td>98.0</td>
</tr>
</tbody>
</table>

(a)* Draw a labelled diagram to show how you would safely set up apparatus for distillation and describe a method to obtain a pure sample of cyclohexanone from the distillate.
(b) Ethanedioic acid removes excess dichromate ions, $\text{Cr}_2\text{O}_7^{2-}$, as in the equation below.

$$3(\text{COOH})_2 + \text{Cr}_2\text{O}_7^{2-} + 8\text{H}^+ \rightarrow 6\text{CO}_2 + 2\text{Cr}^{3+} + 7\text{H}_2\text{O}$$

Suggest how you could tell when the excess dichromate has completely reacted with the ethanedioic acid.

.................................................................................................................................................................................
............................................................................................................................................................................... [1]

(c) A student monitors the course of this reaction using thin-layer chromatography (TLC).

Outline how TLC could be used to monitor the course of the reaction.

............................................................................................................................................................................. [2]

(d) Plan an experiment that would allow the student to confirm the identity of the pure organic product by means of a chemical test.

............................................................................................................................................................................ [3]
A chemist isolates compound L, with empirical formula C₃H₆O, and sends a sample for analysis. The analytical laboratory sends back the following spectra.

**Mass spectrum**
Molecular ion peak at \( m/z = 116.0 \).

**¹H NMR spectra**
The numbers next to each signal represent the number of \(^1\text{H}\) responsible for that signal. Two \(^1\text{H}\) NMR spectra were obtained: one without D₂O and one with D₂O added.

**¹H NMR spectrum with no D₂O:**

![⁰H NMR spectrum with no D₂O](image)

Expansion of multiplet centred at \( \delta 2.7 \text{ ppm} \)

**¹H NMR spectrum with D₂O added:**

![⁰H NMR spectrum with D₂O added](image)
Use the information provided to suggest a structure for compound L. Give your reasoning.
...day June 20XX – Morning/Afternoon
A Level Chemistry A
H432/02   Synthesis and analytical techniques

SAMPLE MARK SCHEME

Duration: 2 hours 15 minutes

MAXIMUM MARK  100

This document consists of 24 pages
MARKING INSTRUCTIONS

PREPARATION FOR MARKING

SCORIS

1. Make sure that you have accessed and completed the relevant training packages for on-screen marking: scoris assessor Online Training; OCR Essential Guide to Marking.

2. Make sure that you have read and understood the mark scheme and the question paper for this unit. These are posted on the RM Cambridge Assessment Support Portal http://www.rm.com/support/ca

3. Log-in to scoris and mark the required number of practice responses (“scripts”) and the required number of standardisation responses.

YOU MUST MARK 10 PRACTICE AND 10 STANDARDISATION RESPONSES BEFORE YOU CAN BE APPROVED TO MARK LIVE SCRIPTS.

MARKING

1. Mark strictly to the mark scheme.

2. Marks awarded must relate directly to the marking criteria.

3. The schedule of dates is very important. It is essential that you meet the scoris 50% and 100% (traditional 50% Batch 1 and 100% Batch 2) deadlines. If you experience problems, you must contact your Team Leader (Supervisor) without delay.

4. If you are in any doubt about applying the mark scheme, consult your Team Leader by telephone, email or via the scoris messaging system.
5. Work crossed out:
   a. where a candidate crosses out an answer and provides an alternative response, the crossed out response is not marked and gains no marks
   b. if a candidate crosses out an answer to a whole question and makes no second attempt, and if the inclusion of the answer does not cause a rubric infringement, the assessor should attempt to mark the crossed out answer and award marks appropriately.

6. Always check the pages (and additional objects if present) at the end of the response in case any answers have been continued there. If the candidate has continued an answer there then add a tick to confirm that the work has been seen.

7. There is a NR (No Response) option. Award NR (No Response)
   - if there is nothing written at all in the answer space
   - OR if there is a comment which does not in any way relate to the question (e.g. ‘can’t do’, ‘don’t know’)
   - OR if there is a mark (e.g. a dash, a question mark) which isn't an attempt at the question.

   Note: Award 0 marks – for an attempt that earns no credit (including copying out the question).

8. The scoris comments box is used by your Team Leader to explain the marking of the practice responses. Please refer to these comments when checking your practice responses. Do not use the comments box for any other reason.

   If you have any questions or comments for your Team Leader, use the phone, the scoris messaging system, or email.

9. Assistant Examiners will send a brief report on the performance of candidates to their Team Leader (Supervisor) via email by the end of the marking period. The report should contain notes on particular strengths displayed as well as common errors or weaknesses. Constructive criticism of the question paper/mark scheme is also appreciated.
10. For answers marked by levels of response:

Read through the whole answer from start to finish, concentrating on features that make it a stronger or weaker answer using the indicative scientific content as guidance. The indicative scientific content indicates the expected parameters for candidates’ answers, but be prepared to recognise and credit unexpected approaches where they show relevance.

Using a ‘best-fit’ approach based on the science content of the answer, first decide which set of level descriptors, Level 1, Level 2 or Level 3, best describes the overall quality of the answer using the guidelines described in the level descriptors in the mark scheme.

Once the level is located, award the higher or lower mark.

**The higher mark** should be awarded where the level descriptor has been evidenced and all aspects of the communication statement (in italics) have been met.

**The lower mark** should be awarded where the level descriptor has been evidenced but aspects of the communication statement (in italics) are missing.

**In summary:**

- The science content determines the level.
- The communication statement determines the mark within a level.

Level of response questions on this paper are 20(a) and 21.
## Annotations

<table>
<thead>
<tr>
<th>Annotation</th>
<th>Meaning</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>DO NOT ALLOW</strong></td>
<td>Answers which are not worthy of credit</td>
</tr>
<tr>
<td><strong>IGNORE</strong></td>
<td>Statements which are irrelevant</td>
</tr>
<tr>
<td><strong>ALLOW</strong></td>
<td>Answers that can be accepted</td>
</tr>
<tr>
<td>( )</td>
<td>Words which are not essential to gain credit</td>
</tr>
<tr>
<td>__</td>
<td>Underlined words must be present in answer to score a mark</td>
</tr>
<tr>
<td><strong>ECF</strong></td>
<td>Error carried forward</td>
</tr>
<tr>
<td><strong>AW</strong></td>
<td>Alternative wording</td>
</tr>
<tr>
<td><strong>ORA</strong></td>
<td>Or reverse argument</td>
</tr>
</tbody>
</table>
12. **Subject-specific Marking Instructions**

**INTRODUCTION**

Your first task as an Examiner is to become thoroughly familiar with the material on which the examination depends. This material includes:

- the specification, especially the assessment objectives
- the question paper
- the mark scheme.

You should ensure that you have copies of these materials.

You should ensure also that you are familiar with the administrative procedures related to the marking process. These are set out in the OCR booklet *Instructions for Examiners*. If you are examining for the first time, please read carefully Appendix 5 *Introduction to Script Marking: Notes for New Examiners*.

Please ask for help or guidance whenever you need it. Your first point of contact is your Team Leader.
SECTION A

<table>
<thead>
<tr>
<th>Question</th>
<th>Key</th>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>B</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>2</td>
<td>B</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>3</td>
<td>B</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>4</td>
<td>D</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>5</td>
<td>A</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>6</td>
<td>B</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>7</td>
<td>B</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>8</td>
<td>B</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>9</td>
<td>C</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>10</td>
<td>B</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>11</td>
<td>D</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>12</td>
<td>C</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>13</td>
<td>A</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>14</td>
<td>D</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>15</td>
<td>A</td>
<td>1</td>
<td></td>
</tr>
</tbody>
</table>

**Total** 15
## SECTION B

<table>
<thead>
<tr>
<th>Question</th>
<th>Answer</th>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>16 (a)</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>(i)</td>
<td>C₂H₂Cl</td>
<td>✓</td>
<td></td>
</tr>
<tr>
<td>(ii)</td>
<td>![Cl]</td>
<td>✓</td>
<td></td>
</tr>
<tr>
<td>(iii)</td>
<td>(compounds with) the same (molecular) formula AND different structures / structural formulae / arrangement of atoms / displayed formulae</td>
<td>✓</td>
<td>DO NOT ALLOW non-skeletal formulae</td>
</tr>
<tr>
<td>(b)</td>
<td>( n = \frac{pV}{RT} = \frac{(100 \times 10^3) \times (1.053 \times 10^{-3})}{8.314 \times 350} ) ✓</td>
<td>4</td>
<td>ALLOW same number of atoms of each element ALLOW different carbon backbone DO NOT ALLOW different spatial arrangement (of atoms)</td>
</tr>
<tr>
<td></td>
<td>n = 0.0362 mol ✓</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>M = ( \frac{m}{n} = \frac{1.321}{0.0362} = 36.5 \text{ (g mol}^{-1} )</td>
<td>✓</td>
<td></td>
</tr>
<tr>
<td></td>
<td>Identity HCl ✓</td>
<td></td>
<td></td>
</tr>
<tr>
<td>(c)</td>
<td></td>
<td></td>
<td></td>
</tr>
<tr>
<td>(i)</td>
<td>![Reaction 1] ✓</td>
<td>2</td>
<td>ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous</td>
</tr>
<tr>
<td></td>
<td>From Reaction 1 = ![compound B] ✓</td>
<td></td>
<td></td>
</tr>
<tr>
<td>Question</td>
<td>Answer</td>
<td>Marks</td>
<td>Guidance</td>
</tr>
<tr>
<td>----------</td>
<td>--------</td>
<td>-------</td>
<td>----------</td>
</tr>
<tr>
<td>(ii)</td>
<td><img src="Image" alt="Diagram" /></td>
<td>3</td>
<td>ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous. Curly arrow must start from covalent bonds and not atoms. DO NOT ALLOW any other partial charges e.g. shown on double bond. DO NOT ALLOW $C^{\delta^+}$ for charge on carbonium ion. Curly arrow from $Cl^-$ can start from the negative charge or the lone pair. DO NOT ALLOW delta negative, i.e. $Cl^{\delta^-}$.</td>
</tr>
<tr>
<td>(iii)</td>
<td>because the intermediate/carbocation in the formation of compound B is less stable (than the intermediate in the formation of compound A)</td>
<td>1</td>
<td></td>
</tr>
<tr>
<td>(iv)</td>
<td><img src="Image" alt="Diagram" /> (Formation of) white precipitate/solid/suspension AND (ppt is) silver chloride/AgCl</td>
<td>2</td>
<td>ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous.</td>
</tr>
<tr>
<td>Question</td>
<td>Answer</td>
<td>Marks</td>
<td>Guidance</td>
</tr>
<tr>
<td>----------</td>
<td>--------</td>
<td>-------</td>
<td>----------</td>
</tr>
<tr>
<td>(d)</td>
<td><strong>Use of elemental analysis data</strong></td>
<td>5</td>
<td></td>
</tr>
<tr>
<td></td>
<td><img src="chart" alt="Table of C, H, O percentages and ratios" /></td>
<td></td>
<td></td>
</tr>
</tbody>
</table>
|          | atom ratio with calculation ✓
|          | empirical formula = C₄H₈O₃ ✓ |       |          |
|          | **IR spectrum**
|          | (very) broad absorption 2500–3300 cm⁻¹ (COOH) ✓
|          | AND absorption 1640–1750 cm⁻¹ (C=O) ✓
|          | absorption 3450 cm⁻¹ (alcohol –OH) ✓ | | ALLOW any values given within ranges given on Data Sheet |
|          | **Identification**
|          | conclusion from data: compound contains –COOH and –OH
<p>|          | (empirical formula confirms no other C=O than in COOH) in place of the previous chlorine-containing groups |
|          | <img src="diagram" alt="Structural formula" /> ✓ | ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous |
|          | <strong>Total</strong> | 20    |          |</p>
<table>
<thead>
<tr>
<th>Question</th>
<th>Answer</th>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>17 (a)</td>
<td>(i)</td>
<td>Step 1: add HCN OR H₂SO₄/KCN ✓</td>
<td>4</td>
</tr>
<tr>
<td></td>
<td></td>
<td>CH₃CHO + HCN → CH₃CH(OH)CN ✓</td>
<td></td>
</tr>
<tr>
<td></td>
<td>(ii)</td>
<td>because (compound D) forms hydrogen bonds form with water ✓</td>
<td>3</td>
</tr>
<tr>
<td></td>
<td></td>
<td>demonstrated through diagram showing: - dashed line between —OH and (:)OH₂ ✓ - dashed line between —NH₂ and (:)OH₂ ✓</td>
<td></td>
</tr>
<tr>
<td>Question</td>
<td>Answer</td>
<td>Marks</td>
<td>Guidance</td>
</tr>
<tr>
<td>----------</td>
<td>--------</td>
<td>-------</td>
<td>----------</td>
</tr>
<tr>
<td>(iii)</td>
<td><img src="image" alt="Chemical Structure" /></td>
<td>2</td>
<td>ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous. ‘End bonds’ MUST be shown (solid or dotted). IGNORE brackets and/or n.</td>
</tr>
<tr>
<td>(b) (i)</td>
<td>Atom A: 3 bonding pairs AND 1 lone pair (therefore) pyramidal AND 107°</td>
<td>4</td>
<td>ALLOW 106–108°. ALLOW 4 bonding pairs but with 1 double/π-bond (therefore 3 bonding centres).</td>
</tr>
<tr>
<td></td>
<td>Atom B: 3 bonding centres (and 0 lone pairs) (therefore) trigonal planar AND 120°</td>
<td></td>
<td></td>
</tr>
<tr>
<td>(c) (i)</td>
<td><img src="image" alt="Chemical Structure" /></td>
<td>1</td>
<td>✓ for all four.</td>
</tr>
<tr>
<td>Question</td>
<td>Answer</td>
<td>Marks</td>
<td>Guidance</td>
</tr>
<tr>
<td>----------</td>
<td>--------</td>
<td>-------</td>
<td>----------</td>
</tr>
</tbody>
</table>
| (ii)     | **Left-hand fragment**<br>\[\begin{array}{c}
\text{O} \\
\text{O} \\
\text{H} \\
\text{N}_2 \\
\text{O} \\
\text{H}
\end{array}\] OR structure with COOH rather than COO\(^-\) ✓ |
|          |        | 4     | ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous |
|          | **Right-hand fragment**<br>\[\begin{array}{c}
\text{O} \\
\text{O} \\
\text{H}_2 \text{N} \\
\text{O} \\
\text{N}_2 \\
\text{O} \\
\text{H}
\end{array}\] OR structure with COOH rather than COO\(^-\) ✓ ✓ |
|          |        |       | ALLOW 1 mark for structure with right-hand ring still intact |
|          | Two OR three COO\(^-\) shown ✓ |

Total 21
<table>
<thead>
<tr>
<th>Question</th>
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<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>18 (a)</td>
<td>(i)</td>
<td>1</td>
<td>ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous ALLOW disubstituted compound with tert-buty1 groups adjacent</td>
</tr>
<tr>
<td></td>
<td>(ii)</td>
<td>3</td>
<td>ALLOW “the oxygen p-orbital overlaps with...” ALLOW diagrammatic answer for 1&lt;sup&gt;st&lt;/sup&gt; and 2&lt;sup&gt;nd&lt;/sup&gt; marks: 1&lt;sup&gt;st&lt;/sup&gt; mark: π system OR 6×p orbitals shown 2&lt;sup&gt;nd&lt;/sup&gt; mark: O lone pair OR O p-orbital AND interaction ALLOW undergoes electrophilic substitution more easily if 1&lt;sup&gt;st&lt;/sup&gt; and 2&lt;sup&gt;nd&lt;/sup&gt; marks achieved through diagram, conclusion must refer to diagram for 3&lt;sup&gt;rd&lt;/sup&gt; mark</td>
</tr>
<tr>
<td>(b)</td>
<td>(i)</td>
<td>1</td>
<td>ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous</td>
</tr>
</tbody>
</table>

Step 1 = (conc.) H<sub>2</sub>SO<sub>4</sub> AND CH₃CH₂OH ✓
<table>
<thead>
<tr>
<th>Question</th>
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<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>(ii)</td>
<td><img src="attachment" alt="Chemical Structure" /></td>
<td>2</td>
<td>ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous</td>
</tr>
</tbody>
</table>

BOTH organic structures balanced equation ✔ ✔
<table>
<thead>
<tr>
<th>Question</th>
<th>Answer</th>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td><strong>19</strong> (a) (i) Product from reaction 1:</td>
<td></td>
<td></td>
<td></td>
</tr>
</tbody>
</table>
| | \[
\begin{align*}
\text{CH}_3\text{CH}_2\text{C}&\text{COOH} \\
\text{CH}_3\text{COO}
\end{align*}
| | ✓ |
| Product from reaction 2: |
| | \[
\begin{align*}
\text{Br} \\
\text{CH}_3\text{CH}_2\text{C}&\text{COOH} \\
\text{H}
\end{align*}
| | ✓ |
| **(ii)** \((E)\)-pent-2-enoic acid | ✓ | 1 | ALLOW “\(E\)” with or without brackets |
| **(iii)** compound \(H\) = |
| | \[
\begin{align*}
\text{CH}_3\text{CH}_2\text{C}&\text{COOH} \\
\text{C}&\text{C} \\
\text{H} \\
\text{H}
\end{align*}
| | ✓ |
| addition polymer = |
| | \[
\begin{align*}
\text{CH}_3\text{CH}_2\text{H} \\
\text{C}&\text{C} \\
\text{H} \\
\text{COOH}
\end{align*}
<p>| | ✓ |
| <strong>(iv)</strong> combustion for energy production use as an organic feedstock for the production of plastics and other organic chemicals | ✓ | 2 | ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous |
| | | | ‘End bonds’ MUST be shown (solid or dotted) IGNORE brackets and/or (n) |</p>
<table>
<thead>
<tr>
<th>Question</th>
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<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>(b) (i)</td>
<td>Oxidising agent = acidified (potassium/sodium) dichromate(VI) ✓</td>
<td>5</td>
<td>ALLOW (\text{Cr}_2\text{O}_7^{2-}) OR (\text{K}_2\text{Cr}_2\text{O}_7) OR (\text{Na}_2\text{Cr}_2\text{O}_7) for dichromate. ALLOW (\text{H}^+) OR (conc.) sulfuric acid for “acidified”. ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous.</td>
</tr>
<tr>
<td>(Oxidation) equation</td>
<td></td>
<td></td>
<td></td>
</tr>
</tbody>
</table>
| | \[
\begin{align*}
\text{OH} & + 3[\text{O}] \rightarrow \text{COOH} + 2\text{H}_2\text{O} \\
\end{align*}
\] | ✓ | |
| (Reduction) mechanism | | | |
| | \[
\begin{align*}
\text{O} & \quad \text{OH} \\
\text{C} & \quad \text{H}^+ \\
\text{H}^+ & \quad \text{O} \\
\end{align*}
\] | ✓ | ALLOW correct structural OR displayed OR skeletal formulae OR a combination of above as long as unambiguous. |
| | \[
\begin{align*}
\text{O} & \quad \text{H}_2\text{O} \\
\text{C} & \quad \text{H}^+ \\
\text{H}^+ & \quad \text{O} \\
\end{align*}
\] | ✓ | ALLOW for second stage. IF \(\text{H}_2\text{O}\) is used it MUST show the curly arrow from the intermediate to \(\text{H}^{5+}\) in \(\text{H}_2\text{O}\) AND from the \(\text{O}–\text{H}\) bond to the \(\text{O}\). IGNORE product. IGNORE stereochemistry of intermediate. |
### Question (ii)

**Na**\(^{+}\) [\(\begin{array}{c}
\text{H} \\
\text{B} \\
\text{H}
\end{array}\)]**\(^{-}\)

1s\(^2\) 2s\(^2\) 2p\(^6\) ✓

- IGNORE inner electron shells for both ions
- Three different symbols required to identify electrons from different elements

### Question (c)

- \(n(\text{NaOH})\) used in titration = 0.150 \(\times\) 18.80/1000 = 0.00282 (mol) ✓
- \(n(\text{H}^+/\text{COOH})\) in 25.0 cm\(^3\) = 0.00282 (mol) ✓
- \(n(\text{H}^+/\text{COOH})\) in 250 cm\(^3\) = 0.0282 (mol) ✓
- ‘Molar’ mass of K = 1.89/0.0282 = 67.0 g mol\(^{-1}\) ✓
- K must be diprotic ✓
- K is malic acid/HOOCCH\(_2\)CHOHCOOH ✓

- **DO NOT ALLOW** [Ne] OR [He] 2s\(^2\) 2p\(^6\)

**Guidance**

Determined through realisation that none of the compounds listed have \(M = 67.0\) g mol\(^{-1}\)

<table>
<thead>
<tr>
<th>Question</th>
<th>Answer</th>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>(ii)</td>
<td><img src="image" alt="Diagram" /></td>
<td>2</td>
<td><strong>IGNORE</strong> inner electron shells for both ions&lt;br&gt;Three different symbols required to identify electrons from different elements</td>
</tr>
<tr>
<td>(c)</td>
<td>(n(\text{NaOH})) used in titration = 0.150 (\times) 18.80/1000 = 0.00282 (mol) ✓&lt;br&gt;(n(\text{H}^+/\text{COOH})) in 25.0 cm(^3) = 0.00282 (mol) ✓&lt;br&gt;(n(\text{H}^+/\text{COOH})) in 250 cm(^3) = 0.0282 (mol) ✓&lt;br&gt;‘Molar’ mass of K = 1.89/0.0282 = 67.0 g mol(^{-1}) ✓&lt;br&gt;K must be diprotic ✓&lt;br&gt;K is malic acid/HOOCCH(_2)CHOHCOOH ✓</td>
<td>5</td>
<td><strong>DO NOT ALLOW</strong> [Ne] OR [He] 2s(^2) 2p(^6)**</td>
</tr>
</tbody>
</table>

**Total** 19
<table>
<thead>
<tr>
<th>Question</th>
<th>Answer</th>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>20 (a)*</td>
<td>Please refer to the marking instructions on page 4 of this mark scheme for guidance on how to mark this question.</td>
<td>6</td>
<td>Indicative scientific points may include:</td>
</tr>
<tr>
<td></td>
<td>Level 3 (5–6 marks) Correctly labelled diagram of apparatus that works, with no safety problems AND Full appreciation of further two steps required to gain pure sample</td>
<td></td>
<td>Diagram</td>
</tr>
<tr>
<td></td>
<td>There is a well-developed diagram which is clear and structured. The information on further purification is detailed and relevant.</td>
<td></td>
<td>Includes following components:</td>
</tr>
<tr>
<td></td>
<td>Level 2 (3–4 marks) Labelled diagram of apparatus but with safety/procedural problems OR clear diagram of functional apparatus without labelling AND Some details of further purification steps</td>
<td></td>
<td>distillation flask</td>
</tr>
<tr>
<td></td>
<td>The diagram presents apparatus that is in the most part relevant with some correct labelling, and supported by some details of further purification steps.</td>
<td></td>
<td>heat source</td>
</tr>
<tr>
<td></td>
<td>Level 1 (1–2 marks) Diagram of apparatus drawn with no labelling OR labelled diagram with significant safety/procedural problems AND Few or imprecise details about further purification stages</td>
<td></td>
<td>thermometer at outlet (bulb level with outlet)</td>
</tr>
<tr>
<td></td>
<td>The diagram is basic and unstructured. Any mention of purification steps is limited to generic term, e.g. ‘drying’, without relevant detail.</td>
<td></td>
<td>still-head</td>
</tr>
<tr>
<td></td>
<td>Shake and leave to settle in a separating funnel</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>Separate layers by tapping off</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>Add (a small amount of) anhydrous magnesium sulfate/anhydrous calcium chloride to organic layer (in a dry conical flask)</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>(Re)distil the organic layer</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>Collect fraction distilling at (between 150 °C and) 156 °C.</td>
<td></td>
<td>water condenser (correct direction of water flow)</td>
</tr>
<tr>
<td></td>
<td>receiving vessel</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>open system.</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>Further purification</td>
<td></td>
<td></td>
</tr>
<tr>
<td>Question</td>
<td>Answer</td>
<td>Marks</td>
<td>Guidance</td>
</tr>
<tr>
<td>----------</td>
<td>------------------------------------------------------------------------------------------</td>
<td>-------</td>
<td>--------------------------------------------------------------------------</td>
</tr>
<tr>
<td></td>
<td>0 marks</td>
<td>0</td>
<td>No response or no response worthy of credit.</td>
</tr>
<tr>
<td>(b)</td>
<td>Lack of (further) effervescence ✔</td>
<td>1</td>
<td>ALLOW fizzing/bubbling stops</td>
</tr>
<tr>
<td>(c)</td>
<td>Take samples from reaction mixture at regular intervals ✔</td>
<td>2</td>
<td>ALLOW “frequent” for “regular”</td>
</tr>
<tr>
<td></td>
<td>Spot/run on a TLC plate, alongside cyclohexanol (and cyclohexanone) controls ✔</td>
<td></td>
<td>ALLOW measure/compare $R_f$ value to cyclohexanol</td>
</tr>
<tr>
<td></td>
<td>IGNORE reference to solvent or visualising chemicals/UV</td>
<td></td>
<td></td>
</tr>
<tr>
<td>(d)</td>
<td>React (sample of distillate) with 2,4-dinitrophenylhydrazine ✔</td>
<td>3</td>
<td>ALLOW (2,4-)DNPH/Brady’s reagent</td>
</tr>
<tr>
<td></td>
<td>recrystallise <strong>AND</strong> determine the melting point ✔</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>Compare melting point to known/library value for cyclohexanone (derivative) ✔</td>
<td></td>
<td></td>
</tr>
<tr>
<td></td>
<td>Total</td>
<td>12</td>
<td></td>
</tr>
</tbody>
</table>
### Question 21*

Please refer to the marking instructions on page 4 of this mark scheme for guidance on how to mark this question.

#### Level 3 (5–6 marks)
- Structure correct
- Analysed all $^1$H NMR signals with at least two supporting statements made.

*The analysis is clear and logically structured. The supporting statements are relevant to the correct structure drawn.*

#### Level 2 (3–4 marks)
- Structure has correct molecular formula AND C=O AND OH but in incorrect positions
- Analysed at least three $^1$H NMR signals with one or two supporting statements made

*The analysis is presented with some structure. The supporting statements are in the most-part relevant to the structure drawn.*

#### Level 1 (1–2 marks)
- Structure has correct molecular formula AND C=O OR OH but in incorrect positions
- Analysed at least two $^1$H NMR signals with no or one supporting statements made

*The analysis is basic and communicated in an unstructured way. The relationship of the supporting evidence to the structure may not be clear.*

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<table>
<thead>
<tr>
<th>Question</th>
<th>Answer</th>
</tr>
</thead>
<tbody>
<tr>
<td>21*</td>
<td>Please refer to the marking instructions on page 4 of this mark scheme for guidance on how to mark this question.</td>
</tr>
</tbody>
</table>

#### Marks

<table>
<thead>
<tr>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
</table>
| 6     | Indicative scientific points may be included:  
Structure  
![Structure Diagram]  
$^1$H NMR spectrum  
$\delta = 3.8$ ppm, triplet, 2H \( \text{CH}_2\text{CH}_2\text{O} \)  
$\delta = 3.7$ ppm, singlet, 1H \( \text{O}--\text{H} \)  
$\delta = 3.1$ ppm, triplet, 2H \( \text{CH}_2\text{CH}_2\text{C}=\text{O} \)  
$\delta = 2.7$ ppm, septet, 1H \( \text{(CH}_3)_2\text{CHC}=\text{O} \)  
$\delta = 1.0$ ppm, doublet, 6H \( \text{(CH}_3)_2\text{CH} \)  

Supporting statements  
$\delta = 3.7$ ppm lost after D$_2$O, indicating \( \text{—O} \)  
$\delta = 213$ ppm in $^{13}$C NMR but no $\delta = 9$–10 ppm in $^1$H NMR so ketone, not aldehyde  
\( M(C_3H_6O) = 58 \ 116/58 = 2 \rightarrow \text{C}_6\text{H}_{12}\text{O}_2 \) |
<table>
<thead>
<tr>
<th>Question</th>
<th>Answer</th>
<th>Marks</th>
<th>Guidance</th>
</tr>
</thead>
<tbody>
<tr>
<td>0 marks</td>
<td>No response or no response worthy of credit.</td>
<td>0</td>
<td></td>
</tr>
</tbody>
</table>

Total 6