## AQA

Please write clearly in block capitals.

Centre number

|  |  |  |  |  |
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Candidate number

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Surname
Forename(s)
Candidate signature
I declare this is my own work.

## A-level

CHEMISTRY

## Paper 2 Organic and Physical Chemistry

Time allowed: 2 hours

## Materials

For this paper you must have:

- the Periodic Table/Data Booklet, provided as an insert (enclosed)
- a ruler with millimetre measurements
- a scientific calculator, which you are expected to use where appropriate.


## Instructions

- Use black ink or black ball-point pen.
- Fill in the boxes at the top of this page.
- Answer all questions.
- You must answer the questions in the spaces provided. Do not write outside the box around each page or on blank pages.
- If you need extra space for your answer(s), use the lined pages at the end of this book. Write the question number against your answer(s).
- All working must be shown.
- Do all rough work in this book. Cross through any work you do not want to be marked.

| For Examiner's Use |  |
| :---: | :---: |
| Question | Mark |
| 1 |  |
| 2 |  |
| 3 |  |
| 4 |  |
| 5 |  |
| 6 |  |
| 7 |  |
| 8 |  |
| 9 |  |
| 10 |  |
| TOTAL |  |

## Information

- The marks for questions are shown in brackets.
- The maximum mark for this paper is 105.
$\qquad$


Coconut oil contains a triester with three identical R groups.
This triester reacts with potassium hydroxide.


Type of compound
Use
$\qquad$

| $\mathbf{0}$ | $\mathbf{1}$ | B | A 1.450 g sample of coconut oil is heated with 0.421 g of KOH in aqueous ethanol |
| :--- | :--- | :--- | :--- | until all of the triester is hydrolysed.

The mixture is cooled.
The remaining KOH is neutralised by exactly $15.65 \mathrm{~cm}^{3}$ of $0.100 \mathrm{~mol} \mathrm{dm}^{-3} \mathrm{HCl}$
Calculate the percentage by mass of the triester $\left(M_{r}=638.0\right)$ in the coconut oil.
$\qquad$

| $\mathbf{0}$ | $\mathbf{1}$. | $\mathbf{4}$ | Suggest why aqueous ethanol is a suitable solvent when heating the coconut oil |
| :--- | :--- | :--- | :--- | with KOH .

Give a safety precaution used when heating the mixture.
Justify your choice.

Reason
$\qquad$
Safety precaution $\qquad$
$\qquad$
Justification $\qquad$

| $\mathbf{0}$ | $\mathbf{2} \quad$ This question is about fuels. |
| :--- | :--- |


| 0 | 2 | $\mathbf{1}$ The petrol fraction obtained from crude oil can be used as fuel in cars. |
| :--- | :--- | :--- |

State the meaning of fraction, as used in the term petrol fraction.
$\qquad$
$\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{2} .2$ | $\mathbf{2}$ Hexadecane $\left(\mathrm{C}_{16} \mathrm{H}_{34}\right)$ can be cracked at high temperature to form petrol... .0 |
| :--- | :--- | :--- |

Complete the equation to show the cracking of one molecule of hexadecane to form hexane and cyclopentane only.

Give the name of a catalyst used in this cracking reaction.

$$
\mathrm{C}_{16} \mathrm{H}_{34} \rightarrow+
$$

Catalyst $\qquad$
$\begin{array}{lllll}0 & 2 & 3 & \text { Carbon dioxide is formed when petrol is burned. }\end{array}$
Carbon dioxide acts as a greenhouse gas when it absorbs infrared radiation.
Give a reason why carbon dioxide absorbs infrared radiation.
$\qquad$
$\qquad$
$\qquad$

Question 2 continues on the next page

| $\mathbf{0}$ | $\mathbf{2} .4 \mathrm{C}$ Compound $\mathbf{Z}\left(\mathrm{HOCH}_{2} \mathrm{CH}_{2} \mathrm{NH}_{2}\right)$ can be used to remove carbon dioxide from the |
| :--- | :--- | :--- | mixture of waste gases produced in some power stations.

Figure 1 shows part of a suggested mechanism for the reaction of $\mathbf{Z}$ with carbon dioxide.

Figure 1

$\mathrm{HOCH}_{2} \mathrm{CH}_{2} \ddot{\mathrm{~N}}_{2}$


Draw two curly arrows to complete the mechanism in Figure 1.
Name compound $\mathbf{Z}\left(\mathrm{HOCH}_{2} \mathrm{CH}_{2} \mathrm{NH}_{2}\right)$
Deduce the role of $\mathbf{Z}$ in step $\mathbf{2}$ of the mechanism.

Name $\qquad$
Role $\qquad$

| $\mathbf{0}$ | $\mathbf{2}$ | $\mathbf{5}$ |
| :--- | :--- | :--- |
| HOCH | $\mathrm{CH}_{2} \mathrm{NH}_{2}$ can be represented as $\mathrm{XNH}_{2}$ |  |

$\left[\mathrm{HOCH}_{2} \mathrm{CH}_{2} \mathrm{NH}_{3}\right]^{+}$can be represented as $\left[\mathrm{XNH}_{3}\right]^{+}$
Draw the shape of $\mathrm{XNH}_{2}$ and of $\left[\mathrm{XNH}_{3}\right]^{+}$
State whether the $\mathrm{H}-\mathrm{N}-\mathrm{H}$ bond angle in $\mathrm{XNH}_{2}$ is greater than, the same as, or smaller than that in $\left[\mathrm{XNH}_{3}\right]^{+}$

Explain your answer.

Shape of $\mathrm{XNH}_{2}$
Shape of $\left[\mathrm{XNH}_{3}\right]^{+}$

Bond angle $\qquad$
Explanation $\qquad$

Question 2 continues on the next page

| $\mathbf{0}$ | $\mathbf{2}$ | $\mathbf{6}$ | Bioethanol is used as an alternative to fossil fuels. |
| :--- | :--- | :--- | :--- |

This statement appeared on a website.
"The fact that bioethanol is a carbon-neutral fuel outweighs the environmental disadvantages of producing bioethanol."

Evaluate this statement.
In your answer you should include:

- an outline of how bioethanol is produced
- relevant equations
- analysis of the environmental impacts.
$\qquad$
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| $\mathbf{0}$ | $\mathbf{3}$ A student does an experiment to determine a value for the enthalpy of combustion of |
| :--- | :--- | :--- | heptane.

Figure 2 shows some of the apparatus used.
Figure 2


| $\mathbf{0}$ | $\mathbf{3}$. | $\mathbf{1}$ Design a table to record all the readings necessary to determine an experimental |
| :--- | :--- | :--- | value for the enthalpy of combustion for heptane in this experiment.


| 0 | $\mathbf{3} .2$ | 2 |
| :--- | :--- | :--- | clamped copper calorimeter.

Suggest two disadvantages of using a glass beaker on a tripod and gauze.

Disadvantage 1 $\qquad$
$\qquad$
Disadvantage 2 $\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{3}$ | $\mathbf{3}$ Suggest two reasons why the value of enthalpy of combustion from this experiment is |
| :--- | :--- | :--- | :--- | less exothermic than a data book value.

[2 marks]
Reason 1 $\qquad$
$\qquad$
Reason 2
$\qquad$

| 0 | 3 | 4 | Suggest one addition to this apparatus that would improve the accuracy of the |
| :--- | :--- | :--- | :--- | enthalpy value obtained.

Kekulé suggested this structure for benzene.

Benzene is now represented by this structure.


Figure 3 shows the relative stability of
 compared to

Figure 3


| $\mathbf{0}$ | $\mathbf{4} .1$ |
| :--- | :--- | :--- |

Table 1

|  | $\Delta H / \mathrm{kJ} \mathrm{mol}^{-1}$ |
| :--- | :---: |
| Enthalpy of atomisation for carbon | +715 |
| Enthalpy of atomisation for hydrogen | +218 |
| Bond enthalpy (C-C) | +348 |
| Bond enthalpy (C=C) | +612 |
| Bond enthalpy (C-H) | +412 |

$\Delta H_{2}$ $\qquad$ $\mathrm{kJ} \mathrm{mol}{ }^{-1}$

| $\mathbf{0}$ | $\mathbf{4}$ | $\mathbf{2}$ Explain, in terms of structure and bonding, why |
| :--- | :--- | :--- |

 is more thermodynamically stable than

$\qquad$
$\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{4}$ | $\mathbf{3}$ | A mixture of concentrated nitric acid and concentrated sulfuric acid reacts with |
| :--- | :--- | :--- | :--- | benzene.

Figure 4 shows the incomplete mechanism for this reaction.
Name the mechanism.
Complete the mechanism in Figure 4 by adding

- any lone pairs of electrons involved in each step
- two curly arrows in step 1
- a curly arrow in step 2
- a curly arrow in step 3
- a curly arrow in step 4.

Name of mechanism
Figure 4


$\qquad$






| $\mathbf{0}$ | 5 |
| :--- | :--- | This question is about equilibrium.


| $\mathbf{0}$ | $\mathbf{5} .1$ |
| :--- | :--- | :--- | presence of a small amount of catalyst.

The mixture is left to reach equilibrium at a constant temperature.

$$
\mathrm{C}_{7} \mathrm{H}_{12} \mathrm{O}_{4}(\mathrm{I})+2 \mathrm{H}_{2} \mathrm{O}(\mathrm{I}) \rightleftharpoons 2 \mathrm{CH}_{3} \mathrm{COOH}(\mathrm{I})+\mathrm{HO}\left(\mathrm{CH}_{2}\right)_{3} \mathrm{OH}(\mathrm{I})
$$

At equilibrium, $x$ mol of ethanoic acid are present in the mixture.
Complete Table 2 by deducing the amounts, in terms of $x$, of the diester, water and diol present in the equilibrium mixture.

Table 2

| Amount in the mixture / mol |  |  |  |  |  |
| :--- | :---: | :---: | :---: | :---: | :---: |
|  | Diester | Water | Acid | Diol |  |
| At the start | 1 | 1 | 0 | 0 |  |
| At equilibrium |  |  | $x$ |  |  |

$\begin{array}{lllll}0 & 5 & 2 & \text { Deduce the structure of the diester in Question } 05.1\end{array}$

| 0 | $\mathbf{5}$ | $\mathbf{3}$ | A new equilibrium mixture of the substances from Question 05.1 is prepared at a |
| :--- | :--- | :--- | :--- | different temperature.

$$
\mathrm{C}_{7} \mathrm{H}_{12} \mathrm{O}_{4}(\mathrm{I})+2 \mathrm{H}_{2} \mathrm{O}(\mathrm{I}) \rightleftharpoons 2 \mathrm{CH}_{3} \mathrm{COOH}(\mathrm{I})+\mathrm{HO}\left(\mathrm{CH}_{2}\right)_{3} \mathrm{OH}(\mathrm{I})
$$

Table 3 shows the amount of each substance in this new equilibrium mixture.
Table 3

| Amount in the mixture / mol |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: |
|  | Diester | Water | Acid | Diol |
| At equilibrium | 0.971 | To be <br> calculated | 0.452 | 0.273 |

The value of the equilibrium constant, $K_{\mathrm{c}}$ is 0.161 at this temperature.
Calculate the amount of water, in mol, in this new equilibrium mixture. Show your working.

| $\mathbf{0}$ | 6 | This question is about isomers with the molecular formula $\mathrm{C}_{5} \mathrm{H}_{10} \mathrm{O}$ |
| :--- | :--- | :--- |


| 0 | 6 | 1 |
| :--- | :--- | :--- | $\mathrm{C}_{5} \mathrm{H}_{10} \mathrm{O}$ that is optically active.


| 0 | 6 | 2 |
| :--- | :--- | :--- | the branched chain aldehyde $\mathrm{C}_{5} \mathrm{H}_{10} \mathrm{O}$

$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$

| 0 | 6 | 3 | Draw the $E$ and $Z$ forms of a structural isomer of $\mathrm{C}_{5} \mathrm{H}_{10} \mathrm{O}$ that shows both |
| :--- | :--- | :--- | :--- | optical and geometric isomerism.


| E isomer | Z isomer |
| :---: | :---: |
|  |  |
|  |  |
|  |  |


| 0 | 6 | 4 | Isomer J is cyclic and has an ether functional group (C-O-C) |
| :--- | :--- | :--- | :--- |

Isomer J has only three peaks in its ${ }^{13} \mathrm{C}$ NMR spectrum.


Draw two other cyclic isomers of $\mathrm{C}_{5} \mathrm{H}_{10} \mathrm{O}$ that have an ether functional group and only three peaks in their ${ }^{13} \mathrm{C}$ NMR spectra.

| $\mathbf{0}$ | $\mathbf{7}$ | This question is about spectroscopy. |
| :--- | :--- | :--- |


| $\mathbf{0}$ | $\mathbf{7}$. |
| :--- | :--- | :--- |

Figure 5 shows the infrared spectrum of $\mathbf{K}$.
Figure 5


Which functional group does $\mathbf{K}$ contain?
Tick $(\checkmark)$ one box.

| Functional Group |  |  |  |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: |
| alcohol | alkene | amine | carbonyl | nitrile |  |
|  |  |  |  |  |  |

Question 7 continues on the next page

| $\mathbf{0}$ | $\mathbf{7}$ | $\mathbf{2}$ | Compound $\mathbf{L}$ has molecular formula $\mathrm{C}_{4} \mathrm{H}_{7} \mathrm{NO}$ |
| :--- | :--- | :--- | :--- |

Figure 6 shows the infrared spectrum of $\mathbf{L}$.
Figure 6

$L$ reacts with $\mathrm{H}_{2}$ in the presence of a nickel catalyst to give compound $\mathbf{M}$.
Suggest three ways in which the infrared spectrum of $\mathbf{M}$ is different from the infrared spectrum of $\mathbf{L}$.

1 $\qquad$
$\qquad$

2 $\qquad$
$\qquad$
3 $\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{7}$ | $\mathbf{3}$ | Figure $\mathbf{7}$ shows the ${ }^{1} \mathrm{H}$ NMR spectrum of $\mathbf{Q}, \mathrm{C}_{3} \mathrm{H}_{7} \mathrm{ClO}$ |
| :--- | :--- | :--- | :--- |

Figure 7


Table 4 shows the chemical shifts ( $\delta$ values) and integration values for each peak.
Table 4

| $\delta$ value / ppm | 3.95 | 3.65 | 3.35 |
| :--- | :---: | :---: | :---: |
| Integration value | 0.6 | 0.6 | 0.9 |

Deduce the structure of $\mathbf{Q}$.
Explain your answer.
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{8} \quad$ This question is about making a diester from cyclohexanol. |
| :--- | :--- | :--- |



| 0 | $\mathbf{8}$. | $\mathbf{1}$ |
| :--- | :--- | :--- |

Give the name of the reagent needed for step 1.

Type of reaction $\qquad$
Reagent $\qquad$

| $\mathbf{0}$ | $\mathbf{8}$ | $\mathbf{2}$ State the reagents needed and give equations for step $\mathbf{2}$ and step $\mathbf{3}$. |
| :--- | :--- | :--- | :--- |

Show the structure of Compound $\mathbf{G}$ in your equations.

Step 2 reagent $\qquad$
Step 2 equation
$\qquad$

Step 3 reagent $\qquad$
Step 3 equation

| 0 | $\mathbf{8}$ | $\mathbf{3}$ Cyclohexane-1,2-diol reacts with ethanedioyl dichloride. |
| :--- | :--- | :--- | :--- |

Give the name of the mechanism for this reaction.
Complete the mechanism to show the formation of one ester link in the first step of this reaction.

Mechanism name $\qquad$
Mechanism



Do not write

Mechanism



| $\mathbf{0}$ | $\mathbf{8}$ | $\mathbf{4}$ Suggest why chemists usually aim to design production methods |
| :--- | :--- | :--- | :--- |

- with fewer steps
- with a high percentage atom economy.
[2 marks]
Fewer steps $\qquad$
$\qquad$
$\qquad$
High percentage atom economy $\qquad$
$\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{9} \quad$ This question is about the ozone layer in the upper atmosphere. |
| :--- | :--- | :--- |


| $\mathbf{0}$ | $\mathbf{9}$. | $\mathbf{1}$ State why the ozone layer is beneficial for living organisms. |
| :--- | :--- | :--- |

$\qquad$
$\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{9}$. | $\mathbf{2}$ State how chlorofluorocarbons (CFCs) form chlorine atoms in the |
| :--- | :--- | :--- |

$\qquad$
$\qquad$
$\qquad$

| $\mathbf{0}$ | $\mathbf{9}$ | $\mathbf{3}$ Give equations to show how chlorine atoms catalyse the decomposition of ozone. |
| :--- | :--- | :--- |


| 0 | $\mathbf{9} .4$ Hydrochlorofluorocarbons (HCFCs) have been used in place of CFCs. |
| :--- | :--- | :--- | In the mechanism to make an HCFC from a fluoroalkane, two incomplete steps are shown.

Complete each step in the mechanism.
Give the name of the type of step shown by both these equations.
$\qquad$
$\bullet \mathrm{CHF}_{2}+\mathrm{Cl}_{2} \rightarrow$ $\qquad$
Type of step $\qquad$
-
$\qquad$
$\qquad$ show

| 1 | $\mathbf{0}$ | This question is about rates of reaction. |
| :--- | :--- | :--- | lodine and propanone react together in an acid-catalysed reaction

$$
\mathrm{CH}_{3} \mathrm{COCH}_{3}(\mathrm{aq})+\mathrm{I}_{2}(\mathrm{aq}) \rightarrow \mathrm{CH}_{3} \mathrm{COCH}_{2} \mathrm{I}(\mathrm{aq})+\mathrm{HI}(\mathrm{aq})
$$

A student completed a series of experiments to determine the order of reaction with respect to iodine.

Method

- Transfer $25 \mathrm{~cm}^{3}$ of $1.0 \mathrm{~mol} \mathrm{dm}^{-3}$ propanone solution into a conical flask.
- Add $10 \mathrm{~cm}^{3}$ of $1.0 \mathrm{~mol} \mathrm{dm}^{-3} \mathrm{HCl}(\mathrm{aq})$
- Add $25 \mathrm{~cm}^{3}$ of $5.0 \times 10^{-3} \mathrm{~mol} \mathrm{dm}^{-3} \mathrm{I}_{2}(\mathrm{aq})$ and start a timer.
- At intervals of 1 minute, remove a $1.0 \mathrm{~cm}^{3}$ sample of the mixture and add each sample to a separate beaker containing an excess of $\mathrm{NaHCO}_{3}(\mathrm{aq})$
- Titrate the contents of each beaker with a standard solution of sodium thiosulfate and record the volume of sodium thiosulfate used.

| $\mathbf{1}$ | $\mathbf{0}$. | $\mathbf{1}$ Suggest why the $1.0 \mathrm{~cm}^{3}$ portions of the reaction mixture are added to an excess of |
| :--- | :--- | :--- | $\mathrm{NaHCO}_{3}$ solution.

$\qquad$
$\qquad$
$\qquad$
$\qquad$
10.2 Suggest why the order of this reaction with respect to propanone can be ignored in this experiment.
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$

Question 10 continues on the next page

The volume of sodium thiosulfate solution used in each titration is proportional to the concentration of iodine in each beaker.

Table 5 shows the results of the experiment.
Table 5

| Time / minutes | Volume of sodium thiosulfate solution <br> $/ \mathbf{c m}^{\mathbf{3}}$ |
| :---: | :---: |
| 1 | 41 |
| 2 | 35 |
| 3 | 24 |
| 4 | 22 |
| 5 | 16 |
| 6 | 10 |


| 1 | 0 | 3 |
| :--- | :--- | :--- | Use the results in Table 5 to draw a graph of volume of sodium thiosulfate solution against time.

Draw a line of best fit.

Volume of sodium thiosulfate solution / $\mathrm{cm}^{3}$


Time / minutes

| $\mathbf{1}$ | $\mathbf{0} .4$ | 4 |
| :--- | :--- | :--- |
| Explain how the graph shows that the reaction is zero-order with respect to iodine in |  |  | the reaction between propanone and iodine.

$\qquad$
$\qquad$
$\qquad$
$\qquad$
$\qquad$

Question 10 continues on the next page

| 1 | 0 | 5 |
| :--- | :--- | :--- |

$$
\ln k=\frac{-E_{\mathrm{a}}}{R T}+\ln \mathrm{A}
$$

Figure 8 shows a graph of $\ln k$ against $\frac{1}{T}$ for the reaction

$$
2 \mathrm{HI}(\mathrm{~g}) \rightarrow \mathrm{H}_{2}(\mathrm{~g})+\mathrm{I}_{2}(\mathrm{~g})
$$

Figure 8


Use Figure 8 to calculate a value for the activation energy $\left(E_{\mathrm{a}}\right)$, in $\mathrm{kJ} \mathrm{mol}^{-1}$, for this reaction.

The gas constant $R=8.31 \mathrm{~J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1}$
$E_{a}$ $\qquad$ $\mathrm{kJ} \mathrm{mol}^{-1}$




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