

# **A-Level Chemistry**

## **Elimination of Alcohols**

## **Mark Scheme**

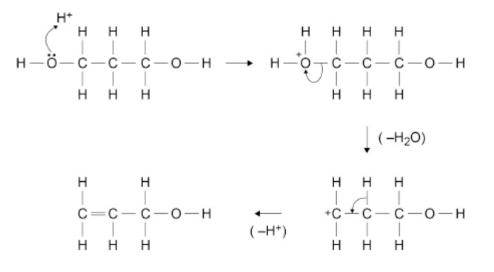
## Time available: 54 minutes Marks available: 50 marks

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#### Mark schemes

### **1.** (a) nucleophilic substitution

(b) M1 elimination



M2 arrow from lone pair on O to H<sup>+</sup>

1.1+

M3 1<sup>st</sup> intermediate and arrow from C–O<sup>+</sup>H<sub>2</sub> bond to O

**M4** 2nd intermediate (carbocation) **and** arrow from a correct C–H bond to correct C–C to form C=C

*Max 2 of 3 marks (M2-4)* for wrong organic reactant (ignore structure of product)

M3 and M4 can be scored in one concurrent step:

M3 for correct intermediate and arrow from  $C-O^+H_2$  bond to O

M4 for arrow from a correct C–H bond to correct C–C to form C=C 1

$$\begin{array}{c} \begin{pmatrix} \mathbf{r}^{\mathsf{T}} \mathsf{H} & \mathsf{H} \\ \mathsf{H} - \ddot{\mathsf{O}} - \ddot{\mathsf{C}} - \ddot{\mathsf{C}} - \ddot{\mathsf{C}} - \ddot{\mathsf{O}} - \mathsf{O} & \mathsf{H} & \longrightarrow & \mathsf{H} - \overset{\mathsf{I}}{\overset{\mathsf{O}}} - \overset{\mathsf{I}}{\overset{\mathsf{C}}} - \overset{\mathsf{I}}{\overset{\mathsf{C}}} - \overset{\mathsf{I}}{\overset{\mathsf{O}}} - \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{I}}{\overset{\mathsf{O}}} + \overset{\mathsf{I}}{\overset{\mathsf{O}}} + \overset{\mathsf{I}}{\overset{\mathsf{O}}} + \overset{\mathsf{I}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{I}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{O}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{O}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}} - \overset{\mathsf{O}}{\overset{\mathsf{O}}} + \overset{\mathsf{O}}{\overset{\mathsf{O}}}$$

1

1

1

1

H H - C - C -H CH<sub>2</sub> OH Any correct structural representation Ignore any brackets and/or n

Structure in any form

(C)

(d)

2.

 $\begin{array}{c} H & H & O \\ | & | & | \\ H - C - C - C - H \end{array}$ 

н

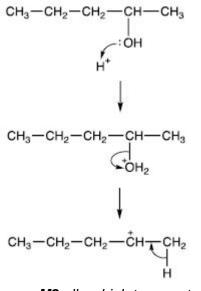
H

1

[7]

			-
(a)	M1	idea that pentan-2-ol has stronger intermolecular forces <i>M1</i> idea that hydrogen bonds are stronger than van der Waals' forces	
		Penalise <b>M1</b> for any reference to idea of breaking covalent bonds	1
	M2	pent-1-ene has van der Waals' forces (only)	
		<b>M2</b> allow London forces or temporary/induced dipole forces or vdW forces for van der Waals' forces	1
	M3	pentan-2-ol (also) has hydrogen bonds	
		M3 Ignore reference to dipole-dipole forces in pentan-2-ol	
			1
(b)	M1	reagent = <u>conc</u> sulfuric acid or <u>conc</u> phosphoric acid	
		<i>M1</i> penalise incorrect name or formula (even if both name and formula are given)	
			1

M2 condition = hot / temperature in range  $150-200^{\circ}(C)$ 



M2 allow high temperature

**M2** reagent must indicate an acid in some way in order for **M2** to be awarded

**M1/2** allow 1 mark if  $H_2SO_4/H_3PO_4$  given as reagent and conc(entrated) given as condition

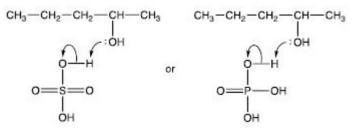
M3 curly arrow from lone pair on alcohol O to H<sup>+</sup>

#### M3-5

penalise **M3/4/5** for any additional arrow(s) in addition to the correct one at each stage

If incorrect reactant (or product if shown), maximum 2 marks of **M3-5** 

#### Alternatives for M3



1

1

1

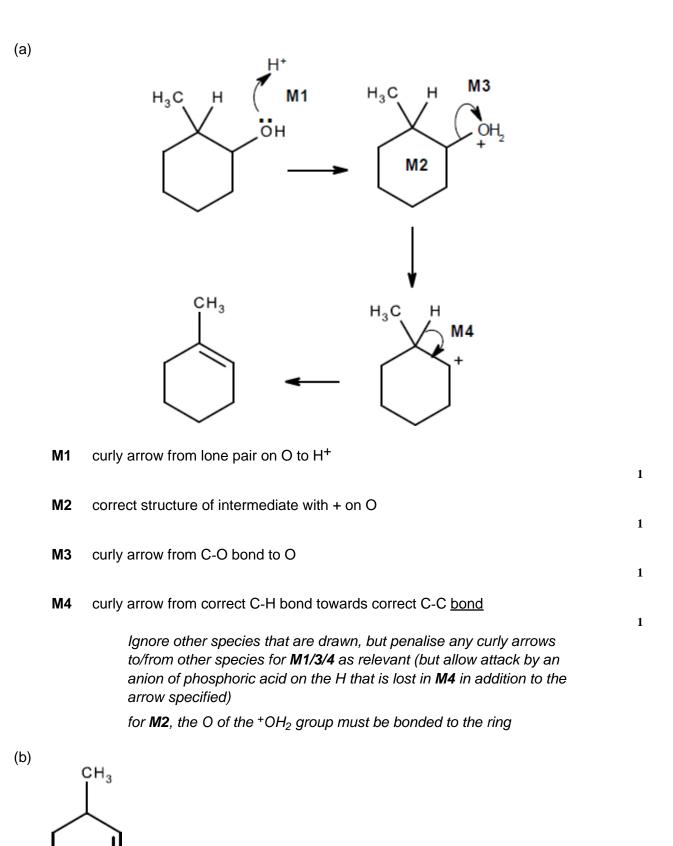
- M4 curly arrow from C-O bond to O on correct intermediate
- M5 arrow from C-H bond on C1 to C-C bond between C1 and C2 on correct carbocation allow M4 and M5 concurrent:

 $CH_3 - CH_2 -$ 

[8]

1

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3.

Any correct structural representation

(c) M1 more stable (carbocation formed) For M1 penalise more stable product

#### M2 changes from secondary to tertiary (carbocation)

For **M2** allow explanation via inductive effect with more alkyl / C groups attached or inductive effect from methyl group as alternatives

Allow 2° or 2<sup>y</sup> for secondary and 3° or 3<sup>y</sup> for tertiary

(d) CH<sub>2</sub>

#### Any correct structural representation

(e) M1 cyclohexene : van der Waals' forces (between molecules) 1 M2 cyclohexanol : hydrogen bonds (between molecules) 1 M3 phosphoric acid: hydrogen bonds (between molecules) 1 M4 idea that cyclohexene has weakest forces 1 M5 separated by (simple / fractional) distillation 1 M6 cyclohexene has lowest boiling point / boils off first Extended response Maximum of 5 marks for answers which do not refer to the van der Waals forces or hydrogen bonds being between molecules in some way M1 penalise reference to presence of other intermolecular forces M1 allow vdW forces (on this occasion) M1/2/3 penalise reference to breaking covalent bonds M2 & M3 ignore reference to van der Waals and/or (permanent) dipole-dipole forces **M2** allow use of term H bonds (on this occasion) M4 allow converse argument M4 & M6 – allow correct comparison of cyclohexene forces and boiling point to one of the other two compounds if only one of cyclohexanol or phosphoric acid discussed

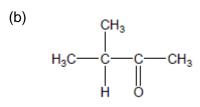
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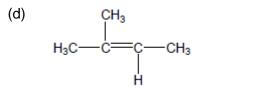
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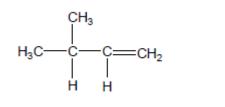
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4.





Allow 
$$(CH_3)_2C=CHCH_3$$



(g) CH<sub>3</sub> | H<sub>3</sub>C-C-CH<sub>2</sub>CH<sub>2</sub>CH<sub>3</sub> | OH

(h) CH<sub>3</sub> | H<sub>3</sub>C---C---CH<sub>2</sub>OH | CH<sub>3</sub>

Allow (CH<sub>3</sub>)<sub>3</sub>CCH<sub>2</sub>OH

1

1

1

1

1

1

1

#### (a) (i) 2-methylpropan-2-ol (1) OR the second one

(ii) Dehydrating agent:  $\operatorname{conc} H_2 SO_4$  OR  $\operatorname{conc} H_3 PO_4$  OR  $\operatorname{Al}_2O_3$  (1)

Equation: 
$$CH_3 \longrightarrow CH_3 \longrightarrow CH_3 \xrightarrow{CH_3} (1)$$

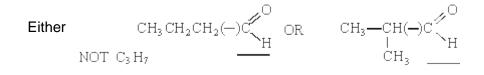
Allow  $C_4H_9OH$  in equation provided RHS is correct if b(i) is blank, b(ii) equation must be full for credit i.e. NOT  $C_4H_9OH$ Mark consequential on b(i)

(b) (i) *Isomer*: butan-2-ol OR <u>the fourth one</u> [look at name in table] wrong isomer = CE

Structure of the ketone:

(ii) *Isomer*: butan-1-ol OR <u>the first one</u> OR 2-methylpropan-1-ol OR <u>the third one</u> [look at name in table]

> Wrong isomer = CE Structure of the aldehyde:



3

5.

Reagent	M1	Tollen's (AgNO <sub>3</sub> /NH <sub>3</sub> )	Fehling's
Observation with ketone	M2	Stays colourless no change	stays blue no change
Observation with aldehyde	М3	Silver mirror black ppt	<u>red solid</u> orange/ <u>red</u> brown/ <u>red</u> <u>ppt/solid</u>

Other include(\*)  $K_2Cr_2O_7 / H_2SO_4$   $KMnO_4/H_2SO_4$ Schiff's Benedict's Wrong reagent R No reagent = CE Penalise AgNO<sub>3</sub> [Ag(NH<sub>3</sub>)<sub>2</sub>] but allow M2 and M3 sequentially.

	(*)	$K_2Cr_2O_7$ / $H_2SO_4$ acidified	<u>ketone</u>	<u>aldehyde</u>				
				green				
		KMnO <sub>4</sub> /H <sub>2</sub> SO <sub>4</sub> acidified	purple no change	colourless (v. Pale pink)				
	В	enedict's ≡ Fehling's ; S v	Schiff's colouless – violet		7			
Equation: $CH_3CH_2CH_2CH_2OH$ (or $C_4H_9OH$ ) + 2[O] $\rightarrow CH_3CH_2CH_2COOH$ (or $C_3H_7COOH$ ) + $H_2O$ (1)								
	Name of product. butanoic acid (1)							

Accept butaneoic acid

(iii)

(c)

54.5