### **M1.** (a)

NB The bonds shown in the structure must be correct

1

1

1

1

1

Isomerism: E-Z isomerism

If written answer is correct, ignore incorrect labelling of structures.

If no written answer, allow correctly labelled structures.

Both COOH groups must be on the same side/ close together/ cis

No rotation about C=C axis

#### Structure

Allow

(b)  $Br_2 / HBr / H_2SO_4 / H^+ / Br^+ / NO_2^+ (Mark M1)$ 

HOOC COOH H COOH HOOC COOH

$$C = C \xrightarrow{(l) M2} M4(l) + C - C - H \longrightarrow Br - C - C - H$$

$$H \xrightarrow{(l) M3} H \xrightarrow{(l) M3} :Br$$

NB If electrophile H<sup>+</sup> / Br<sup>+</sup> / NO<sub>2</sub><sup>+</sup> allow M1, M2 and M4 If the acid is incorrect, M2 and M3 can still be scored Allow M4 consequentially if repeat error from part (a)

4

(c) e.g. 2NaOH + HO₂CCHCHCO₂H → NaO₂CCHCHCO₂Na + 2H₂O
 Both H replaced

1

Balanced for atoms and charges

1

NB Allow ionic equations and  $2NaOH + C_4H_4O_4 \rightarrow C_4H_2O_4Na_2 + 2H_2O$ 

Allow one if structure incorrect but molecular formula correct

Allow one for a correct equation showing one H replaced

(d) M1 Two peaks

1

M2 No splitting or singlets

1

M3 (Two) non-equivalent protons or two proton environments

1

M4 No adjacent protons

1

M5 Same area under the two peaks or same relative intensity

1

NB Doublet could score M1 and M3 or M5 (Max 2)

More than two peaks CE = 0

Apply the "list principle" to incorrect answers if more than 3 given

Max 3

[15]

**M2.**C

[1]

**M3.**B

[1]

**M4.**D

[1]

**M5.**D

M6.

[1]

(a)  $H \subset CH_3 \subset CH_3 \subset CH_3$ 

May circle 4 C's separately

1

(b)  $\begin{array}{c}
H^{+}\text{can score M1 + M2} \\
H_{2}\text{SO}_{4} \text{ only M1 - see diagram not M2}
\end{array}$   $\begin{array}{c}
H \\
C = C \\
H
\end{array}$   $\begin{array}{c}
C + CH_{3} \\
C + CH_{2}CH_{3}
\end{array}$   $\begin{array}{c}
H \\
C - C - CH_{2}CH_{3}
\end{array}$   $\begin{array}{c}
H \\
C - C - CH_{2}CH_{3}
\end{array}$   $\begin{array}{c}
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H \\
C - CH_{2}CH_{3}$   $\begin{array}{c}
H \\
C -$ 

Ignore  $\delta$ + and  $\delta$ - unless wrong

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(c) Reagent: H<sub>2</sub>O or water **OR steam, Or dilute sulphuric acid (1)**Condition: heat, or warm, or boil or reflux [50-100°C] **(1)**Name of compound **C**: 2-methylbutan-2-ol **(1)** 

Allow 2-methylbutan<u>e</u>-2-ol Penalise hydroxy-2-methylbutane and 2-methylbut-2-ol once only in the paper

3

3

(d) Identity of alcohol D: 2-methylbutan-1-ol (1), OR its structure, could describe structure

Explanation: C formed via t-carbocation; D via p-carbocation, (1) tertiary more stable than primary (1)

If have wrong carbocation can still score stability mark

[11]

**M7.**B

[1]

M8. (a) Identity of X; 2-methylpropene (1)
Absorption at 1650 cm<sup>-1</sup> indicates an alkene present (1)

OR a chemical answer e.g. Br<sub>2</sub> (aq) brown to colourless

2

(b) Reagents

Step 1 KOH (allow NaOH) (1) alcoholic (1) warm (1)

Only allow solvent and warm if reagent correct

Step 2 HBr (1)

Mechanism:  $A \rightarrow X$ 

$$CH_{3} - CH_{2} CH_{2} CH_{2} CH_{3} CH_{3} C = C CH_{3}$$

$$CH_{3} CH_{3} C = C CH_{3}$$

$$CH_{3} CH_{3} C = C CH_{3}$$

$$CH_{3} CH_{3} C = C CH_{3}$$

Or a carbocation mechanism

(c) A gives three peaks (1) B gives one peak (1)

Allow one for "A has more peaks than B" when no number of peaks is given

[15]

11

2

M9. (a) (i)

(1) M4
arrow
$$Br\Theta$$

H<sub>3</sub>C = C H

H<sub>3</sub>C - C - CH<sub>3</sub>
H

M3
(1) carbocation

(1) M4
(1) C - C - CH<sub>3</sub>
(1) C -

If wrong carbocation, lose structure mark
If wrong alkene, lose structure mark
Can still score ¾ i.e. penalise M3
Penalise M2 if polarity included incorrectly
no bond between H and Br
bond is shown as ••• or ••••

Page 6

credit secondary carbocation here if primary carbocation has been used in (i)

Ignore attack on this carbocation by Bro

5

#### (b) (i) Structure:

$$\begin{array}{c} \text{OH} \\ \text{I} \\ \text{H}_3\text{C} - \text{CH} - \text{CH}_3 \end{array} \text{ (1)} \qquad \begin{bmatrix} \text{in sist on} \\ \text{C} - \text{OH bond} \end{bmatrix}$$

No credit for propan-1-ol even when named correctly Credit propane-2-ol

Name: propan-2-ol (1)
Not 2-hydroxypropane

# (ii) Name of mechanism: nucleophilic substitution (1) (both words) (NOT S₁1 orS₂2)

Mechanism:

penalise incorrect polarity on C - Br (M1) Credit the arrows even if incorrect haloalkane If  $S_{N}1$ , both marks possible

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## (c) (i) elimination (1) Ignore nucleophylic elimination

### Penalise electrophilic elimination

(ii) base **(1)** 

OR proton acceptor NOT nucleophile (base)

[12]

2