**M1.**(a) Fractional distillation (under reduced pressure)

1

(b) BaSO₄ insoluble / remove by filtration

Do not allow answers which refer to reaction rate

1

(c) Both contain OH group

Allow OH stretch in ir spectrum of each compound Do not allow 'same bonds'

[3]

1

**M2.** (a) (i)

(ii) H<sub>3</sub>C—O or ROCH<sub>3</sub>;

(allow 1 if both (i) and (ii) give CH<sub>3</sub>- or H<sub>3</sub>C- only)

1

1

(iii) CH<sub>2</sub>CH<sub>2</sub> or two <u>adjacent</u> methylene groups;

1

(iv)

$$\begin{array}{c} \text{CH}_3\text{-CCH}_2\text{-CH}_2\text{-OCH}_3\\ \parallel\\ \bigcirc\\ \end{array}$$

OR

(b) (i) OH in acids or (carboxylic) acid present

(ii)

(c)

reagent	K₂Cr₂O₁/H⁺	KMnO₄ /H⁺
Y	no reaction	no reaction
Z	orange to green or turns green	purple to colourless or turns colourless

[9]

5

M3. (a) Allow 1 mark each for any correctly drawn primary, secondary and tertiary alcohol of molecular formula  $C_4H_8O$ 

3

Tertiary alcohol cannot be oxidised

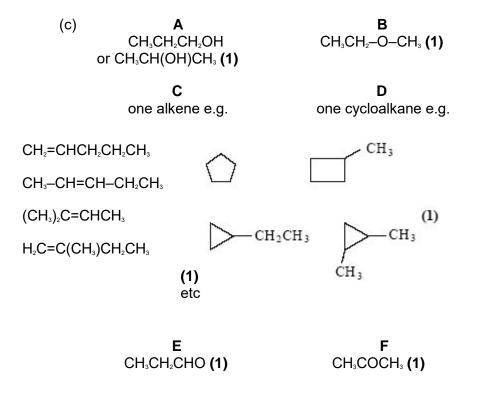
1

(b) Region 1500-400 cm<sup>-1</sup>

1

exact match to spectrum of known compound

1



M4. Functional group (isomerism) (a)

(b)

M1 Tollens' (reagent) (Credit ammoniacal silver nitrate OR Benedict's solution a description of making Tollens') (Ignore either AgNO<sub>3</sub> or [Ag(NH<sub>3</sub>)<sub>2</sub><sup>+</sup>] or "the silver mirror test" on their own, but mark M2 and M3)

M1 Fehling's (solution) or (Ignore Cu<sup>2+</sup>(aq) or CuSO4 on their own, but mark on to M2 and M3)

6

1

[12]

M2 silver mirror M2 Red solid/precipitate (Credit orange or brown solid)

OR

black solid/precipitate (NOT silver precipitate)

M3 (stays) colourless M3 (stays) blue or no change or no reaction or no change or no reaction

Mark on from an incomplete/incorrect attempt at the correct

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reagent, penalising M1
```

No reagent, CE=0

Allow the following alternatives

**M1** (acidified) potassium dichromate(VI) (solution)

M2 (turns) green

M3 (stays) orange/no change

OR

**M1** (acidified) potassium manganate(VII) (solution)

M2 (turns) colourless

M3 (stays) purple/no change

For M3

Ignore "nothing (happens)"

Ignore "no observation"

(c) (Both have) C=O **OR** a carbonyl (group)

3

(d) (i) (Free-) radical substitution ONLY

Penalise "(free) radical mechanism"

1

(ii) Initiation

 $Cl_2 \rightarrow 2Cl^{\bullet}$ 

Penalise absence of dot once only.

First propagation

CI• +  $CH_3CH_2CH_3 \rightarrow \bullet CH_2CH_2CH_3 + HCI$ OR  $C_3H_8$ 

Penalise incorrect position of dot on propyl radical once only.

Penalise C<sub>3</sub>H<sub>7</sub>• once only

Second propagation

 $Cl_2 + {\bullet}CH_2CH_2CH_3 \rightarrow CH_3CH_2CH_2CI + CI{\bullet}$ 

OR

C<sub>3</sub>H<sub>7</sub>CI

Accept CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>• with the radical dot above/below/to the side of the last carbon.

Termination (must make C<sub>6</sub>H<sub>14</sub>)

2 •  $CH_2CH_2CH_3 \rightarrow C_6H_{14}$  or  $CH_3CH_2CH_2CH_2CH_3$ 

Use of the secondary free radical might gain 3 of the four marks

4

(e) M<sub>r</sub> = 44.06352 (for propane)
M<sub>r</sub> = 43.98982 (for carbon dioxide)
Mark independently
 M1 a correct value for both of these M<sub>r</sub> values.
 M2 a statement or idea that two peaks appear (in the mass spectrum)
 OR
 two molecular ions are seen (in the mass spectrum).

[12]

2

1

**M5.** (a) Secondary **OR** 2° (alcohol);

(b)

Spectrum is for **butanone (or formula) or butan-2-one**<u>The explanation marks depend on correctly identifying butanone.</u>

If butanone is correctly identified, award any two from

- (Strong) absorption / peak at approximately 1700 (cm<sup>-1</sup>) / 1710 (cm<sup>-1</sup>) / in the range 1680 1750 (cm<sup>-1</sup>) This needs to be stated.
- (Characteristic) absorption / peak for C=O (may be shown on the spectrum in the correct place).
- No absorption / peak in range 3230 to 3550 cm<sup>-1</sup>.

1

No absorption / peak for an OH group.

Look at the spectrum to see if anything is written on it that might gain credit.

Allow the words "dip" OR "spike" OR "low transmittance" as alternatives for absorption.

2

(c) <u>Displayed structure</u> for 2-methylpropan-2-ol

Must have all bonds drawn out but ignore the bond angles

[5]

M6. (a) M1

<u>Displayed formula</u> for butan-2-ol

M1 displayed formula <u>must</u> have all bonds drawn out, including the O—H but ignore angles
Penalise "sticks"

M2 Alcohol X is

M2 structure must be clearly identifiable as 2-methylpropan-2-ol and may be drawn in a variety of ways.

**M3** Alcohol **Y** is named <u>(2)-methylpropan-1-ol</u> ONLY *M3* <u>must be correct name</u>, but ignore structures

3

(b) **M1** The infrared spectrum shows an <u>absorption/peak in the range</u> 3230 to 3550 (cm<sup>-1</sup>)(which supports the idea that an alcohol is present)

In M1, allow the words "dip", "spike", "low transmittance" and "trough" as alternatives for absorption.

**M2** Reference to the 'fingerprint region' or below 1500 (cm<sup>-1</sup>)

M3 Match with or same as known sample/database spectra

Check the spectrum to see if alcohol OH is labelled and credit.

OR

M2 Run infrared spectra (of the alcohols)

M3 Find which one matches or is the same as this spectrum.

3

(c) M1 balanced equation  $C_6H_{12}O_6 \rightarrow CH_3CH_2CH_2CH_2OH + 2CO_2 + H_2O$  or  $C_4H_6OH$ 

Or multiples for M1 and M3 In M1 and M3 penalise use of  $C_4H_{10}O$  or butan-2-ol once only

M2 Any one from

- excess/adequate/sufficient/correct amount of/enough/plenty/ a good supply of oxygen or air
- good mixing of the fuel and air/oxygen
   For M2, do <u>not</u> accept simply "oxygen" or "air" alone Ignore reference to "temperature"

M3 CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>CH<sub>2</sub>OH + 
$$\mathbf{6O}_2 \rightarrow \mathbf{4CO}_2 + \mathbf{5H}_2$$
O or C<sub>4</sub>H<sub>9</sub>OH

**M4** A biofuel is a fuel produced <u>from</u> (renewable) <u>biological (re)source(s)</u>

OR

(renewable) (re)source(s) <u>from</u> (a specified) <u>plant(s)/fruit(s)/tree(s)</u>

In M4

Ignore references to "carbon neutral"

Ignore "sugar" and "glucose"

4

- (d) **M1** butan-1-ol is a <u>primary or 1°</u> (alcohol)
  - M2 Displayed formula (ONLY) for butanal CH<sub>3</sub>CH<sub>2</sub>CH<sub>2</sub>CHO
  - M3 <u>Displayed formula</u> (ONLY) for butanoic acid CH<sub>3</sub>CH<sub>2</sub>COOH

    M2 and M3 displayed formula must have all bonds drawn out including the O—H but ignore angles.

If butanal and butanoic acid formulae are <u>both</u> correctly given but not displayed, credit one mark out of two.

## M4 Oxidation (oxidised) OR Redox

## M5 orange to green

Both colours required for M5 Ignore states

[15]

5