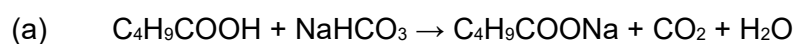


## Mark schemes

**Q1.**

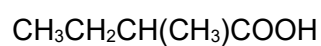
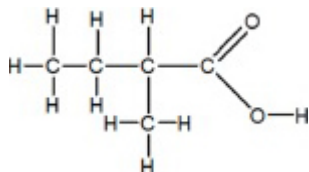
- (a) **M1** + 3 C<sub>2</sub>H<sub>6</sub>
- M2** Zeolite/Aluminosilicate/Aluminium oxide 2
- (b) Option B 1
- (c) Alkenes 1
- (d) **M1** Initial volume O<sub>2</sub> =  $0.21 \times 1350 = 283.5 \text{ (cm}^3\text{)}$   
*Alternative route:*  
*M1 Vol Air decreases by  $6.5 \times 20 = 130 \text{ cm}^3$*
- M2** Volume of O<sub>2</sub> remaining =  $M1 - (6.5 \times 20) = 153.5 \text{ cm}^3$   
*M2 =  $1220 \text{ cm}^3$*
- M3** Volume of CO<sub>2</sub> formed =  $20 \times 4 = 80 \text{ cm}^3$   
*M3 Vol CO<sub>2</sub> produced =  $4 \times 20 = 80 \text{ cm}^3$*
- M4** Total volume of gas left =  $M2 + M3 + (0.79 \times 1350) = 1300 \text{ cm}^3$   
*M4 Total Vol Air + CO<sub>2</sub> =  $1220 + 80 = 1300 \text{ cm}^3$*  4
- (e) **M1** Acid rain  
*M1 Allow damages (limestone) buildings or statues/death of aquatic organisms/air pollution*
- M2** CaO or CaCO<sub>3</sub> 2

**[10]**

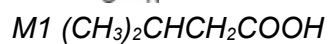
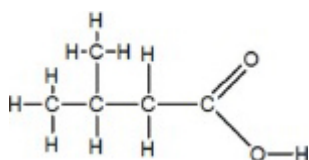
**Q2.**

1

(b)



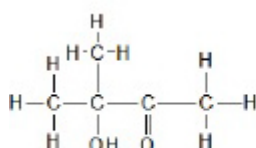
1

(c) **M1**

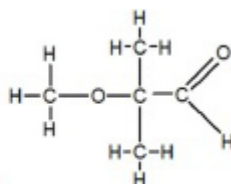
*M2* 6:1:2:1 (Any order)

*M2* Allow ECF for a 5 carbon carboxylic acid

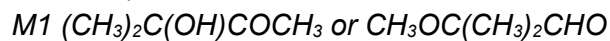
2

(d) **M1**

Or

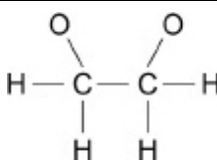
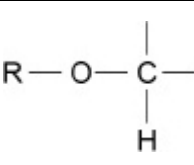
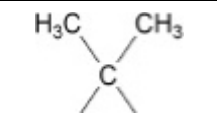
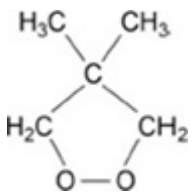
**M2**

Adjacent C has no (non-equivalent) H attached (so no splitting/spin-spin coupling takes place)



2

(e)

	Scores M1 and M2	Allow M1 for	
	Scores M3 and M4	Allow M3 for	RCH <sub>3</sub>
	Scores M5		<p>This structure also scores M5</p> 
<sup>13</sup> C peaks	= 3	M6	Allow ECF from their M5 of C <sub>5</sub> H <sub>10</sub> O <sub>2</sub>

6

[12]

**Q3.**

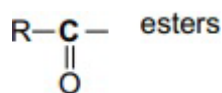
(a) C=O

1

(b) Tick in the box for 7 ONLY

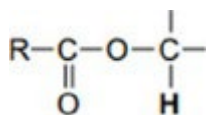
1

(c)

*Ignore acids*

1

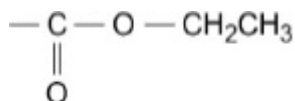
(d) M1 (Quartet) because neighbouring C has 3H

M2 (At  $\delta = 4.1$  ppm) because connected to single bonded O of ester or

M3 (Triplet) because neighbouring C has 2H

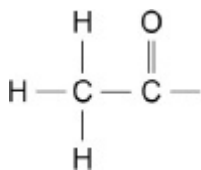
M4 (At  $\delta = 1.26$  ppm) because  $\text{R}_2\text{CH}_2$  or  $\text{RCH}_3$ 

M5

*Ignore use of integration*

5

(e)



1

(f) Cannot deduce splitting patterns of peaks (at about  $\delta = 2.60$ )

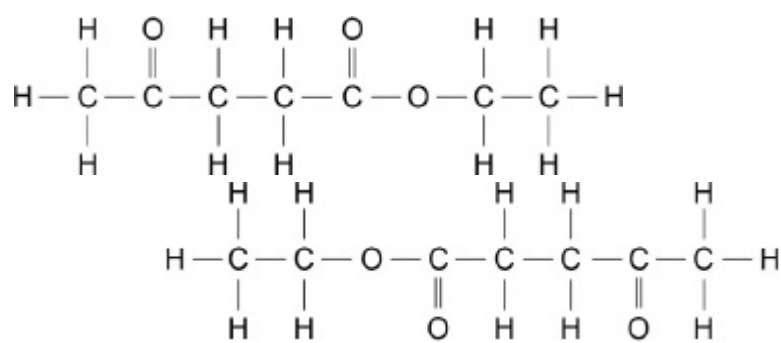
Or

No integration values

*Allow**Peaks at  $\delta = 2.60$  and  $\delta = 2.56$  ppm overlap**OR**spectrum at  $\delta = 2.60$  is second order*

1

(g)



1

[11]

**Q4.**

This question is marked using Levels of Response. Refer to the Mark Scheme Instructions for Examiners for guidance.

<b>Level 3</b> <b>5-6 marks</b>	<p>All stages are covered and each stage is generally correct and virtually complete.</p> <p>Answer is communicated coherently and shows a logical progression from Stage 1 to Stages 2 and 3</p> <p>Covers at least 1 point for stage 1, 3 for stage 2 and 3 for stage 3.</p>
<b>Level 2</b> <b>3-4 marks</b>	<p>All stages are covered but stage(s) may be incomplete or may contain inaccuracies</p> <p>Covers at least 1 point for stage 1 stage 2 and stage 3.</p> <p>OR</p> <p>two stages are covered and are generally correct and virtually complete.</p> <p>Covers at least 1 point for stage 1, and 3 for stage 2 or stage 3 OR 3 for stage 2 and 3 for stage 3</p> <p>Answer is communicated mainly coherently and shows a logical progression from Stage 1 to Stages 2 and 3.</p>
<b>Level 1</b> <b>1-2 marks</b>	<p>Two stages are covered but stage(s) may be incomplete or may contain inaccuracies OR only one stage is covered but is generally correct and virtually complete.</p> <p>Answer includes isolated statements but these are not presented in a logical order.</p>
<b>0 marks</b>	Insufficient correct chemistry to gain a mark.

**Indicative Chemistry content****Stage 1:** infrared

1a) (broad peak) at  $3400\text{ cm}^{-1}$  (any value from 3230-3550) indicates OH in alcohols

1b) peak at  $1720\text{ cm}^{-1}$  (any value from 1680-1750) indicates C=O

**Stage 2:**  $^1\text{H}$  nmr

2a) peak at 3.9 ppm integration 1 so 1 H-C-O AND quartet so adjacent to  $\text{CH}_3$  (stated or shown)

2b) peak at 3.7 ppm integration 1 so HO-C-(stated or shown)

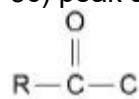
2c) peak at 2.1 ppm integration 3 so  $\text{H}_3\text{C-C=O}$  AND singlet so no adjacent H (stated or shown)

2d) peak at 1.2 ppm integration 3 so  $\text{H}_3\text{C-}$  AND doublet so adjacent to CH (stated or shown)

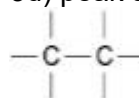
2e) sum of integration values = 8 Hence  $\text{C}_4\text{H}_8\text{O}_2$

**Stage 3:**  $^{13}\text{C}$  nmr3a) peak at 210 ppm  $\text{C}=\text{O}$  aldehydes or ketones3b) peak at 75 ppm  $\text{C}-\text{O}$  (alcohols, ethers or esters)

3c) peak at 25 ppm



3d) peak at 20 ppm



3e) structure

