



Mark Scheme (Results)

Summer 2018

Pearson Edexcel International Advanced
Level in Chemistry (WCH04) Paper 01
General Principles of Chemistry I - Rates,
Equilibria and Further Organic Chemistry

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General Marking Guidance

- All candidates must receive the same treatment. Examiners must mark the first candidate in exactly the same way as they mark the last.
- Mark schemes should be applied positively. Candidates must be rewarded for what they have shown they can do rather than penalised for omissions.
- Examiners should mark according to the mark scheme not according to their perception of where the grade boundaries may lie.
- There is no ceiling on achievement. All marks on the mark scheme should be used appropriately.
- All the marks on the mark scheme are designed to be awarded. Examiners should always award full marks if deserved, i.e. if the answer matches the mark scheme. Examiners should also be prepared to award zero marks if the candidate's response is not worthy of credit according to the mark scheme.
- Where some judgement is required, mark schemes will provide the principles by which marks will be awarded and exemplification may be limited.
- When examiners are in doubt regarding the application of the mark scheme to a candidate's response, the team leader must be consulted.
- Crossed out work should be marked UNLESS the candidate has replaced it with an alternative response.

Using the Mark Scheme

Examiners should look for qualities to reward rather than faults to penalise. This does NOT mean giving credit for incorrect or inadequate answers, but it does mean allowing candidates to be rewarded for answers showing correct application of principles and knowledge. Examiners should therefore read carefully and consider every response: even if it is not what is expected it may be worthy of credit.

The mark scheme gives examiners:

- an idea of the types of response expected
- how individual marks are to be awarded
- the total mark for each question
- examples of responses that should NOT receive credit.

/ means that the responses are alternatives and either answer should receive full credit.

() means that a phrase/word is not essential for the award of the mark, but helps the examiner to get the sense of the expected answer.

Phrases/words in **bold** indicate that the meaning of the phrase or the actual word is **essential** to the answer.

ecf/TE/cq (error carried forward) means that a wrong answer given in an earlier part of a question is used correctly in answer to a later part of the same question.

Candidates must make their meaning clear to the examiner to gain the mark. Make sure that the answer makes sense. Do not give credit for correct words/phrases which are put together in a meaningless manner. Answers must be in the correct context.

Quality of Written Communication

Questions which involve the writing of continuous prose will expect candidates to:

- write legibly, with accurate use of spelling, grammar and punctuation in order to make the meaning clear
- select and use a form and style of writing appropriate to purpose and to complex subject matter
- organise information clearly and coherently, using specialist vocabulary when appropriate.

Full marks will be awarded if the candidate has demonstrated the above abilities.

Questions where QWC is likely to be particularly important are indicated (QWC) in the mark scheme, but this does not preclude others.

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 1 | <p>The only correct answer is B</p> <p><i>A is not correct because this term does not relate to gas chromatography</i></p> <p><i>C is not correct because this term is not used</i></p> <p><i>D is not correct because this term is not used</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 2 | <p>The only correct answer is D</p> <p><i>A is not correct because this would not work</i></p> <p><i>B is not correct because this would not work</i></p> <p><i>C is not correct because this would not work</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 3(a) | <p>The only correct answer is A</p> <p><i>B is not correct because UV does not heat chemicals</i></p> <p><i>C is not correct because radio waves only affect nuclei</i></p> <p><i>D is not correct because UV does not heat chemicals and radio waves only affect nuclei</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 3(b) | <p>The only correct answer is D</p> <p><i>A is not correct because IR only makes bonds vibrate/heats up /accelerates reaction</i></p> <p><i>B is not correct because microwaves only heat up /accelerates reaction</i></p> <p><i>C is not correct because radio waves only affect nuclei</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|--|------------|
| 4(a) | <p>The only correct answer is C</p> <p><i>A is not correct because neither W nor X are methyl ketones</i></p> <p><i>B is not correct because Y is not the only molecule which is a methyl ketone</i></p> <p><i>D is not correct because Z is not the only molecule which is a methyl ketone</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 4(b) | <p>The only correct answer is B</p> <p><i>A is not correct because there are two hydrogen atoms joined to one carbon of the C=C</i></p> <p><i>C is not correct because Y is not an alkene</i></p> <p><i>D is not correct because X does have geometric isomers</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 4(c) | <p>The only correct answer is C</p> <p><i>A is not correct because X also gives a positive result</i></p> <p><i>B is not correct because W also gives a positive result</i></p> <p><i>D is not correct because neither Y nor Z gives a positive result</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|--|------------|
| 4(d) | <p>The only correct answer is B</p> <p><i>A is not correct because neither molecule W nor X would reduce to pentan-2-ol</i></p> <p><i>C is not correct because molecule Z would still have the C=C</i></p> <p><i>D is not correct because molecule Z would still have the C=C</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|--|------------|
| 4(e) | <p>The only correct answer is D</p> <p>A is not correct because a straight-chain is produced</p> <p>B is not correct because a straight-chain is produced</p> <p>C is not correct because molecule Z would also form a branched molecule</p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 4(f) | <p>The only correct answer is D</p> <p>A is not correct because Y has a CH_3CH_2 likely to produce a fragment $m/e = 29$</p> <p>B is not correct because W has a HCO likely to produce a fragment $m/e = 29$</p> <p>C is not correct because Z would not be expected to produce a fragment $m/e = 29$</p> | (1) |

| Question Number | Answer | Mark |
|-----------------|--|------------|
| 5(a) | <p>The only correct answer is A</p> <p>B is not correct because the concentrations should be squared and not doubled</p> <p>C is not correct because the squared concentrations should be multiplied and not added</p> <p>D is not correct because the concentrations should be squared and not doubled, and multiplied not added.</p> | (1) |

| Question Number | Answer | Mark |
|-----------------|--|------------|
| 5(b) | <p>The only correct answer is D</p> <p>A is not correct because dichromate(VI) is not reduced</p> <p>B is not correct because NaOH does not decolorise the solution</p> <p>C is not correct because the equilibrium position shifts to the right</p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 6 | <p>The only correct answer is A</p> <p><i>B is not correct because the units are inverted</i></p> <p><i>C is not correct because the decimetres are to the power +3</i></p> <p><i>D is not correct because the solids are absent from K_c</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|--|------------|
| 7(a) | <p>The only correct answer is A</p> <p><i>B is not correct because total pressure is multiplied by the mole fraction and not moles of chlorine</i></p> <p><i>C is not correct because total pressure is multiplied and not divided by the mole fraction</i></p> <p><i>D is not correct because total pressure is not divided by the moles of chlorine</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 7(b) | <p>The only correct answer is B</p> <p><i>A is not correct because K_p has been used as total pressure and the two partial pressures quoted are subtracted from it</i></p> <p><i>C is not correct because the two partial pressures have been multiplied by K_p instead of dividing by K_p</i></p> <p><i>D is not correct because this is the inverse value</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 8(a) | <p>The only correct answer is B</p> <p><i>A is not correct because low pressure favours the reactants</i></p> <p><i>C is not correct because both high temperature and low pressure favour the reactants</i></p> <p><i>D is not correct because high temperature favours the reactants</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|--|------------|
| 8(b) | <p>The only correct answer is C</p> <p><i>A is not correct because the value of K_c is a ratio not a quantity of a substance</i></p> <p><i>B is not correct because 1 mol of CO cannot form more than 1 mol of methanol</i></p> <p><i>D is not correct because this ignores the CO</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 9(a) | <p>The only correct answer is B</p> <p><i>A is not correct because the -148 has been omitted</i></p> <p><i>C is not correct because the -148 has been omitted and the value has not been halved</i></p> <p><i>D is not correct because the enthalpy value has to be halved for one mole of chloride ions</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 9(b) | <p>The only correct answer is D</p> <p><i>A is not correct because hydration is from gaseous ions not those in a lattice</i></p> <p><i>B is not correct because hydration is nothing to do with the forming of ions</i></p> <p><i>C is not correct because potassium ions are larger</i></p> | (1) |

| Question Number | Answer | Mark |
|-----------------|---|------------|
| 10 | <p>The only correct answer is C</p> <p><i>A is not correct because these are the base and its conjugate acid</i></p> <p><i>B is not correct because these are not conjugate pairs</i></p> <p><i>D is not correct because these are not conjugate pairs</i></p> | (1) |

TOTAL FOR SECTION A = 20 MARKS

Section B

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------------------------|------------|
| 11(a) | $[H^+] = \sqrt{(1.34 \times 10^{-5} \times 0.0500)} / = 8.185... \times 10^{-4} \text{ (1)}$ $pH = (-\log 8.185... \times 10^{-4} = 3.08696) = 3.09 \text{ (1)}$ Correct answer without working scores 2 Allow TE from incorrect $[H^+]$ only if $pH < 7$ | 3 / 3.1 ≥3dp -3.09 | (2) |
| 11(b)(i) | $pH = (14 - (-\log 0.04)) = 12.60206/12.60/12.6$ OR $pH = -\log (1.0 \times 10^{-14} \div 0.04) = 12.60206/12.60/12.6$ IGNORE SF except 1 or 2 | 13 | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--------|------------|
| 11(b)(ii) | <p>Sketch continues vertical until between 10-11 then plateau at pH~12-12.6</p> <p>ALLOW TE for plateau at answer to (b)(i)</p> <p>Exemplar graph</p> | | (1) |

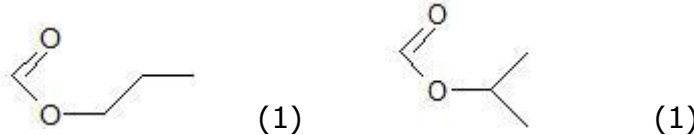
| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|-------------------------------------|------------|
| 11(b)(iii) | <p>pH at half equivalence point = $pK_a = 4.9$ to 5.0 (1)</p> <p>$K_a = (\text{inv log } -\text{pH} = \text{inv log } -4.9 = 1.26 \times 10^{-5} / 0.0000126)$ or ($\text{inv log } -\text{pH} = \text{inv log } -5.0 = 1.00 \times 10^{-5} / 0.0000100)$</p> <p>$= 1.0 \times 10^{-5} / 0.000010 /$ $1.3 \times 10^{-5} / 0.000013 (\text{mol dm}^{-3})$ (1)</p> <p>Correct final answer scores (2)</p> | <p>Not 2 SF Incorrect units</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------|------------|
| 11(b)(iv) | <p>Indicator is thymolphthalein</p> <p>ALLOW Alizarin yellow R if the pH has been given as 13 in (b)(i) (1)</p> <p>Because the pH range of the indicator must lie within the large / vertical pH change</p> <p>ALLOW the pH range overlaps with the vertical part of the sketch / the equivalence point/neutralisation point/end point lies with the pH range of the indicator / pK_{in} value is within the vertical part of the sketch (1)</p> <p>ALLOW Max (1) if a weak acid-weak base titration curve given in (b)(ii) and the comment made that no indicator would be suitable because there is no vertical section of the curve</p> <p>Marking point 2 is dependent on marking point 1</p> | | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--------|------------|
| 11 (c) | <p>Method 1</p> <p>Marking point 1 Initial moles of butanoic acid = $0.025 \times 0.06 = 1.5 \times 10^{-3}$ (mol) and moles of alkali = butanoate = $0.015 \times 0.08 = 1.2 \times 10^{-3}$ (mol) (1)</p> <p>Marking point 2 Equilibrium moles of butanoic acid in buffer = $(1.5 \times 10^{-3} - 1.2 \times 10^{-3}) = 3(.0) \times 10^{-4}$ (mol) (1)</p> <p>Marking point 3 Concentration of hydrogen ions $[H^+] = (K_a \times (3 \times 10^{-4} \div 0.04) \div (1.2 \times 10^{-3} \div 0.04))$ $= 3.75 \times 10^{-6}$ (mol dm⁻³) (1)</p> <p>Marking point 4 $pH = (-\log 3.75 \times 10^{-6}) = 5.42597/5.426/5.43/5.4$ (1) 5</p> <p>Method 2 (use of the Henderson-Hasselbalch equation) $pH = pK_a + \log ([A^-] \div [HA])$</p> <p>Marking point 1 Initial moles of butanoic acid = $0.025 \times 0.06 = 1.5 \times 10^{-3}$ (mol) and moles of alkali = butanoate = $0.015 \times 0.08 = 1.2 \times 10^{-3}$ (mol) (1)</p> <p>Marking point 2 Equilibrium moles of butanoic acid in buffer = $(1.5 \times 10^{-3} - 1.2 \times 10^{-3}) = 3(.0) \times 10^{-4}$ (mol) (1)</p> <p>Marking point 3 [butanoic acid] in buffer = $(3 \times 10^{-4} \div 0.04) = 7.5 \times 10^{-3}$ (mol dm⁻³) and [butanoate] = $(1.2 \times 10^{-3} \div 0.04) = 0.030$ (mol dm⁻³) and $pK_a = 4.8239 \dots$ (1)</p> <p>Marking point 4 $pH = (4.8239 \dots + \log (0.030 \div 7.5 \times 10^{-3}))$ $= 5.42597/5.426/5.43/5.4$ (1) 5</p> <p>Correct answer without working scores (4) Ignore SF except 1SF</p> | | (4) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|-----------------------------|------------|
| *11(d) | <p>Marking point 1 Reference to (large) reservoirs/excess of both ethanoic acid and ethanoate ions/sodium ethanoate</p> <p>ALLOW Reference to (large) reservoirs/excess of both the acid and its conjugate base/salt (1)</p> <p>Marking point 2 Reference to added OH⁻ ions reacting with ethanoic acid OR CH₃COOH + OH⁻ → CH₃COO⁻ + H₂O</p> <p>ALLOW Reference to added OH⁻ ions reacting with hydrogen ions to form water / H⁺ + OH⁻ → H₂O followed by further dissociation of the ethanoic acid (1)</p> <p>Marking point 3 The concentrations of ethanoic acid and ethanoate ions do not change (significantly and hence the pH doesn't change appreciably) / the ratio of [ethanoic acid]:[ethanoate ion] does not change (significantly and hence the pH doesn't change appreciably)</p> <p>ALLOW the ratio of ethanoic acid : ethanoate ion does not change (significantly and hence the pH doesn't change appreciably) (1)</p> <p>IGNORE Keeps pH constant</p> | Reservoir of H ⁺ | (3) |

(Total for Question 11 = 15 marks)

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------|------------|
| 12 (a)(i) |  <p>ALLOW One mark for two correct isomers if drawn with displayed/structural formulae</p> <p>IGNORE Other formulae if correct skeletal formulae given Orientation/bond angles</p> | | (2) |



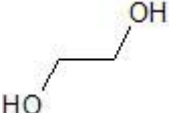
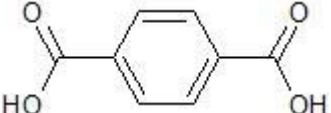

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------------|------------|
| 12(a)(ii) | $\% C = (((12 \times 4) \div 88) \times 100 =) 54.545/54.55/ 54.5/55\%$ IGNORE SF except 1 | 54/ 54.54 | (1) |

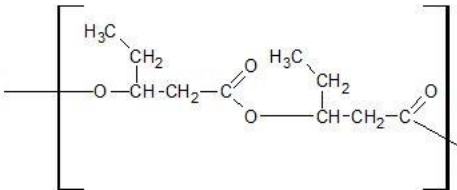
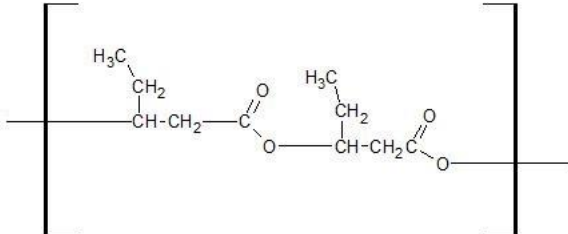
| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--|------------|
| 12(b) | (Reagent) methanol Allow methan-1-ol IGNORE Strong acid such as sulfuric acid/H ₂ SO ₄ (1) (C ₂ H ₅ COOH +) CH ₃ OH (⇌ C ₂ H ₅ COOCH ₃ +) H ₂ O (1) IGNORE state symbols even if incorrect Mark independently No TE on incorrect reagent | NaOH C ₄ H ₁₀ O | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|---|------------|
| 12 (c)(i) | <p>Accept formulae and names in any order (Equation) $\text{CH}_3\text{COOC}_2\text{H}_5 + \text{NaOH} \rightarrow \text{CH}_3\text{COO}^{(-)}\text{Na}^{(+)} + \text{C}_2\text{H}_5\text{OH}$</p> <p>Allow ionic equation $\text{CH}_3\text{COOC}_2\text{H}_5 + \text{OH}^- \rightarrow \text{CH}_3\text{COO}^- + \text{C}_2\text{H}_5\text{OH}$ (1)</p> <p>(Names) (Sodium) ethanoate and ethanol Allow ethan-1-ol (1)</p> <p>IGNORE state symbols even if incorrect</p> <p>Mark independently</p> <p>No TE on incorrect ester and its hydrolysis products</p> | <p>$\text{CH}_3\text{COO}-\text{Na}$</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|--|---|------------|
| 12 (c)(ii) | <p>(It is a catalyst because) alkali speeds up the reaction (compared to water) (1)</p> <p>IGNORE comments on activation energy</p> <p>(It is not a catalyst because) the amount of alkali present would decrease/ would be used up / is not regenerated (1)</p> <p>IGNORE references to the alkali taking part in the reaction</p> | <p>Reaction goes to completion/ Yield is increased</p> | (2) |

| Question Number | Acceptable Answers | Reject | Mark | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|------------------------------|--|----------------------------|-----------------|-------------------|------------------------------|------------------|--|--|-----------------------------------|--|--|-----------------|--|--|---------------------------------------|--|--|---|--|--|--|--|-----------------------------|--|--|--------------|-------------------------|-------------------------|--------------|-------------------------|-------------------------|------------------|-------------------------|-------------------------|----------------------------|-------------------------|-------------------------|------------------------------|----------------------------|----------------------------|------------------------------|----------------------------|----------------------------|--|-----|
| *12 (d) | <p>For marking points 1 to 4, use the table below for reference and award 1 mark for each complete row.</p> <table border="1"> <thead> <tr> <th></th> <th>Ethyl ethanoate</th> <th>Methyl propanoate</th> </tr> </thead> <tbody> <tr> <td>Similarities (any 2 from)</td> <td colspan="2" style="text-align: center;">Both have</td> </tr> <tr> <td></td> <td colspan="2">3 hydrogen or proton environments</td> </tr> <tr> <td></td> <td colspan="2">Peak area 3:2:3</td> </tr> <tr> <td></td> <td colspan="2">Singlet: quartet: triplet (any order)</td> </tr> <tr> <td></td> <td colspan="2">Triplet at 0.1 - 1.9 / peak area 3 at 0.1 - 1.9</td> </tr> <tr> <td></td> <td colspan="2">Peaks at $\delta = 0.1 - 1.9, 1.7 - 3.0$ and $3.0 - 4.2$</td> </tr> <tr> <td>Differences (any 2 from)</td> <td></td> <td></td> </tr> <tr> <td>Quartet peak</td> <td>at $\delta = 3.0 - 4.2$</td> <td>at $\delta = 1.7 - 3.0$</td> </tr> <tr> <td>Singlet peak</td> <td>at $\delta = 1.7 - 3.0$</td> <td>at $\delta = 3.0 - 4.2$</td> </tr> <tr> <td>Peak with area 2</td> <td>at $\delta = 3.0 - 4.2$</td> <td>at $\delta = 1.7 - 3.0$</td> </tr> <tr> <td>(Singlet) peak with area 3</td> <td>at $\delta = 1.7 - 3.0$</td> <td>at $\delta = 3.0 - 4.2$</td> </tr> <tr> <td>Peak at $\delta = 3.0 - 4.2$</td> <td>is a quartet has area 2</td> <td>is a singlet has area 3</td> </tr> <tr> <td>Peak at $\delta = 1.7 - 3.0$</td> <td>is a singlet has area 3</td> <td>is a quartet has area 2</td> </tr> </tbody> </table> <p>Marking point 5 Suitable explanation of one splitting pattern to include reference to (n+1) rule (1) Example of annotated diagrams:</p> <div style="display: flex; justify-content: space-around;"> <div style="text-align: center;"> <p>ETHYL ETHANOATE</p> <p>Singlet Area=3 H-C-C=O $\delta = 1.7-3.0$</p> <p>Quartet Area=2 H-C-O $\delta = 3.0-4.2$</p> <p>Triplet Area=3 H-C-C $\delta = 0.1-1.9$</p> </div> <div style="text-align: center;"> <p>METHYL PROPANOATE</p> <p>Triplet Area=3 H-C-C $\delta = 0.1-1.9$</p> <p>Quartet Area=2 H-C-C=O $\delta = 1.7-3.0$</p> <p>Singlet Area=3 H-C-O $\delta = 3.0-4.2$</p> </div> </div> <p>These diagrams provide evidence for marking points 1 to 4 inclusive and would thus score (4)</p> | | Ethyl ethanoate | Methyl propanoate | Similarities (any 2 from) | Both have | | | 3 hydrogen or proton environments | | | Peak area 3:2:3 | | | Singlet: quartet: triplet (any order) | | | Triplet at 0.1 - 1.9 / peak area 3 at 0.1 - 1.9 | | | Peaks at $\delta = 0.1 - 1.9, 1.7 - 3.0$ and $3.0 - 4.2$ | | Differences (any 2 from) | | | Quartet peak | at $\delta = 3.0 - 4.2$ | at $\delta = 1.7 - 3.0$ | Singlet peak | at $\delta = 1.7 - 3.0$ | at $\delta = 3.0 - 4.2$ | Peak with area 2 | at $\delta = 3.0 - 4.2$ | at $\delta = 1.7 - 3.0$ | (Singlet) peak with area 3 | at $\delta = 1.7 - 3.0$ | at $\delta = 3.0 - 4.2$ | Peak at $\delta = 3.0 - 4.2$ | is a quartet has area 2 | is a singlet has area 3 | Peak at $\delta = 1.7 - 3.0$ | is a singlet has area 3 | is a quartet has area 2 | | (5) |
| | Ethyl ethanoate | Methyl propanoate | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Similarities (any 2 from) | Both have | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | 3 hydrogen or proton environments | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | Peak area 3:2:3 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | Singlet: quartet: triplet (any order) | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | Triplet at 0.1 - 1.9 / peak area 3 at 0.1 - 1.9 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| | Peaks at $\delta = 0.1 - 1.9, 1.7 - 3.0$ and $3.0 - 4.2$ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Differences (any 2 from) | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Quartet peak | at $\delta = 3.0 - 4.2$ | at $\delta = 1.7 - 3.0$ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Singlet peak | at $\delta = 1.7 - 3.0$ | at $\delta = 3.0 - 4.2$ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Peak with area 2 | at $\delta = 3.0 - 4.2$ | at $\delta = 1.7 - 3.0$ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| (Singlet) peak with area 3 | at $\delta = 1.7 - 3.0$ | at $\delta = 3.0 - 4.2$ | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Peak at $\delta = 3.0 - 4.2$ | is a quartet has area 2 | is a singlet has area 3 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| Peak at $\delta = 1.7 - 3.0$ | is a singlet has area 3 | is a quartet has area 2 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------|------------|
| 12 (e)(i) | <p>(monomer structures in any order)</p>  <p>OR</p>  <p style="text-align: right;">(1)</p>  <p style="text-align: right;">(1)</p> <p>ACCEPT structural, displayed and /or Kekulé formulae, or any combination, such as</p> <p>HOOC-C₆H₄-COOH</p> <p>ClOC-C₆H₄-COCl</p>   <p>for the diacid / diacyl chloride</p> <p>OR</p> <p>HOCH₂CH₂OH / (CH₂OH)₂ for the diol</p> | | (2) |

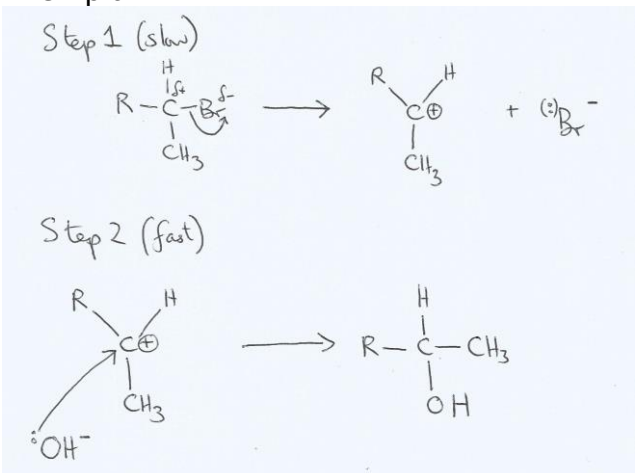
| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------------|------------|
| 12(e)(ii) | <p>Suitable diagram, e.g.</p>  <p>Ester link and continuation 'bonds' (1)</p> <p>Rest of repeat unit correct (1)</p> <p>ACCEPT use of displayed/skeletal formulae</p> <p>ALLOW C-O- at either end i.e.</p>  <p>ALLOW more than two repeat units</p> <p>ALLOW C₂H₅ for CH₃CH₂</p> <p>IGNORE missing brackets or use of 'rounded' brackets</p> | Missing 'H's | (2) |

(Total for Question 12 = 18 marks)

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--|------------|
| 13(a) | <p>Any two correct formulae scores 1 All three correct formulae scores 2</p> <p>Primary – $\text{CH}_3(\text{CH}_2)_2\text{CH}_2\text{I}$ / $\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{I}$ / $(\text{CH}_3)_2\text{CHCH}_2\text{I}$</p> <p>Secondary – $\text{CH}_3\text{CHICH}_2\text{CH}_3$ / $\text{CH}_3\text{CHIC}_2\text{H}_5$</p> <p>Tertiary – $(\text{CH}_3)_3\text{CI}$</p> <p>COMMENT Penalise skeletal / displayed formulae once only but ALLOW 'semi-structural' e.g. $\text{CH}_3-\text{CH}_2-\text{CH}_2-\text{CH}_2-\text{I}$ and</p> $\begin{array}{c} \text{CH}_3 \\ \\ \text{CH}_3-\text{C}-\text{CH}_3 \\ \\ \text{I} \end{array}$ <p>IGNORE missing brackets i.e. allow $\text{CH}_3\text{CI}(\text{CH}_3)\text{CH}_3$ or $\text{CH}_3\text{CICH}_3\text{CH}_3$</p> | $\text{CH}_3\text{CH}_3\text{CH}_3\text{CI}$ | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|--------|------------|
| 13(b)(i) | <p>First order with respect to $\text{C}_3\text{H}_7\text{Br}$ (1)</p> <p>Comparing Expts 2 and 3 the rate doubles as the concentration doubles (and $[\text{OH}^-]$ does not change) (1)</p> <p>First order with respect to OH^- (1)</p> <p>Comparing Expts 1 and 2 the rate quadruples as the concentration quadruples (and $[\text{C}_3\text{H}_7\text{Br}]$ does not change) (1)</p> <p>ALLOW comparison of Expts 1 and 3 if the order w.r.t. to one reactant has already been found</p> | | (4) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|--------|------------|
| 13(b)(ii) | They are spectator ions / they do not take part in the reaction ALLOW (Sodium and potassium ions) are not involved in the rate determining step IGNORE reference to zero order or taking part in a fast step | | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|--|------------|
| 13(c)(i) | <p>Suitable reaction mechanism to include:</p> <p>Marking point 1 Curly arrow from C-Br bond to/just beyond Br and dipole on C-Br bond (1)</p> <p>Marking point 2 Formation of carbocation and bromide ion (1) ALLOW bromide ion / NaBr shown as a final product</p> <p>Marking point 3 Curly arrow from lone pair on hydroxide ion to C+ on carbocation and formation of alcohol product (1)</p> <p>Exemplar</p>  <p>ALLOW (for a correct S_N2 reaction) M1 for correct dipole and curly arrow showing C-Br bond breaking M3 for correct curly arrow from a lone pair on the OH⁻ ion to the δ⁺ carbon, either in a transition state or not, and the alcohol product</p> | <p>Incorrect starting bromoalkane</p> <p>Curly arrow from the charge on OH⁻</p> | (3) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|---|------------------|------------|
| 13(c)(ii) | <p>Slow step only involves one species / only 2-bromopentane is involved in the rate-determining step / there are two steps, first slow and second fast / OH⁻ is not involved in the slow step</p> <p>and</p> <p>so the reaction is first order (so consistent with the rate equation) / so only 2-bromopentane is in the rate equation</p> | S _N 2 | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|---|--------------------------|------------|
| 13(c)(iii) | <p>Carbocation is (trigonal) planar about the C⁺</p> <p>ALLOW</p> <p>Reaction site is (trigonal) planar (1)</p> <p>EITHER</p> <p>Attack by nucleophile/OH⁻ can be (equally) from the top or bottom/either side (of the carbocation) (1)</p> <p>Forming a racemate/racemic mixture/equal number of moles of each optical isomer (1)</p> <p>OR</p> <p>The bromide ion obstructs attack by the hydroxide ion (1)</p> <p>Forming mixture of two optical isomers in which one predominates (1)</p> <p>TE on S_N2 in (c)(i):</p> <p>Attack by nucleophile is only from one side of the molecule (1)</p> <p>The bromide ion leaves from the other side of the molecule / molecule inverts (1)</p> <p>Forming only one optical isomer (1)</p> | Halogenoalkane is planar | (3) |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---|------------|
| 13(d) | Gradient = $(-2.5 \div 0.0004 =) -6250 \pm 50$ (K) (1) Activation Energy = $(-(-6250 \times 8.31))$ $= +51937.5/51900/52000$ J mol ⁻¹ OR $+ 51.9 / 52$ kJ mol ⁻¹ Value in the range 51-53 (1) Sign and units (1) ALLOW Max (2) for TE for incorrect slope $\times -8.31$ IGNORE SF | K ⁻¹ Incorrect units Negative value Incorrect units | (3) |

(Total for Question 13 = 17 marks)
Total for Section B = 50 marks

Section C

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---|------------|
| 14(a)(i) | Any two from: Bubble(s)/fizzing/effervescence (1) (White) Liquid/paste produced ALLOW Solution forms (1) Temperature decreases (1) Pungent smell (of ammonia)/pungent gas (1) IGNORE Gas produced Chemical test for ammonia and water | Incorrect chemical tests and incorrect observations e.g. white ppte / misty fumes | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|---|------------|
| 14(a)(ii) | (High) positive value is expected because: 3 moles → 13 moles ALLOW More moles of product (then reactant) (1) (two) solids → gas and/or liquid (+ one solid) (1) <u>Alternative answer</u> (High) positive value is expected because: Entropy of the surroundings for an endothermic reaction is negative (1) The reaction is feasible with the total entropy being positive (so entropy of the system has to be a large enough positive value for this to be true) (1) | Negative value scores (0) Incorrect numbers of moles quoted Molecules | (2) |

Allow TE throughout all parts (b)(i)-(iv) even if the calculation results in an incorrect sign

Ignore SF, except 1SF, in (b)(i), (b)(iii) and (b)(iv)

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--------|------------|
| 14 (b)(i) | $\Sigma S^{\circ}_{(products)} = ((2 \times 192.3) + (2 \times 69.9) + 213.8 =)$ $(+738.2 \text{ J K}^{-1} \text{ mol}^{-1} \quad (1)$ $\Sigma S^{\circ}_{(reactants)} = ((2 \times 151.1) + 99.7 =) (+)401.9 \text{ J K}^{-1} \text{ mol}^{-1}$ (1) $\Delta S^{\circ}_{system} = (738.2 - 401.9 =) +336.3 \text{ J K}^{-1} \text{ mol}^{-1} \quad (1)$ Correct answer without working scores (3) | | (3) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--------|------------|
| 14(b)(ii) | $\Sigma \Delta H^{\circ}_{(products)} = ((2 \times -285.8) + (2 \times -46.1) + -992.1 =)$ $-1655.9 \text{ (kJ mol}^{-1}\text{)}$ and $\Sigma \Delta H^{\circ}_{(reactants)} = ((2 \times -365.6) + -944.7 =)$ $-1675.9 \text{ (kJ mol}^{-1}\text{)} \quad (1)$ $\Delta H^{\circ}_{reaction} = (-1655.9 - -1675.9 =) +20 \text{ kJ mol}^{-1} \quad (1)$ Correct answer without working scores (2) | | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|---|--------|------------|
| 14(b)(iii) | $\Delta S^{\circ}_{surroundings} = (- \Delta H \div T) = - \frac{20\,000 \text{ J mol}^{-1}}{298 \text{ K}} \quad (1)$ $= - 67.11409... / -67.1 / -67 \text{ J K}^{-1} \text{ mol}^{-1} \quad (1)$ Correct answer without working scores (2) | | (2) |

| Question Number | Acceptable Answers | Reject | Mark |
|------------------|--|--------|------------|
| 14(b)(iv) | $\Delta S^{\circ}_{total} = \Delta S^{\circ}_{system} + \Delta S^{\circ}_{surroundings}$ $\Delta S^{\circ}_{total} = +336.3 - 67.1$ $= +269.2 \text{ J K}^{-1} \text{ mol}^{-1}$ | | (1) |

| Question Number | Acceptable Answers | Reject | Mark |
|-------------------|---|------------------------|------------|
| 14 (c) (i) | Molecules of chlorine disperse / diffuse / spread / mix into the top jar ALLOW Chlorine/it fills the (new) gas jar (1) Therefore (the system has) greater entropy / greater randomness / greater disorder (1) M2 is dependent on M1 | Reacts Enthalpy | (2) |

| Question Number | Acceptable Answers | Reject | Mark | | | | | | | | | | | | |
|------------------|---|------------|------------|-----|---|---|-----|---|---|---|---|---|---|--|------------|
| 14(c)(ii) | <table border="1" style="margin-left: auto; margin-right: auto;"> <thead> <tr> <th>Molecule A</th> <th>Molecule B</th> </tr> </thead> <tbody> <tr> <td>(2</td> <td>2</td> </tr> <tr> <td>3</td> <td>1)</td> </tr> <tr> <td>1</td> <td>3</td> </tr> <tr> <td>4</td> <td>0</td> </tr> <tr> <td>0</td> <td>4</td> </tr> </tbody> </table> The bottom three rows can be completed in any order | Molecule A | Molecule B | (2 | 2 | 3 | 1) | 1 | 3 | 4 | 0 | 0 | 4 | | (1) |
| Molecule A | Molecule B | | | | | | | | | | | | | | |
| (2 | 2 | | | | | | | | | | | | | | |
| 3 | 1) | | | | | | | | | | | | | | |
| 1 | 3 | | | | | | | | | | | | | | |
| 4 | 0 | | | | | | | | | | | | | | |
| 0 | 4 | | | | | | | | | | | | | | |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|--|---------------------|------------|
| 14(d) | General shape of increase from left to right ALLOW straight line (1) Two vertical stages for melting and boiling (1) Labelling to include the use of 1074K for melting and 1686K for boiling temperature (1) | Horizontal sections | (3) |
| | | | |

| Question Number | Acceptable Answers | Reject | Mark |
|-----------------|---|---|------------|
| *14 (e) | <p>Any two from:</p> <p>In the solid / lattice, the ions are fixed (1)</p> <p>But in the solution the ions are mobile /more dispersed / more spread out/ more random (so entropy has increased) (1)</p> <p>(Dissolving results in) an increase in the number of moles of particles and from one to two/ $\text{NaCl} \rightarrow \text{Na}^+ + \text{Cl}^-$ (1)</p> <p>The hydrogen bonds between the water molecules are disrupted (1)</p> | <p>Sodium chloride molecules</p> <p>Due to having more energy</p> | (2) |

TOTAL FOR SECTION C (QUESTION 14) = 20 MARKS

TOTAL FOR PAPER = 90 MARKS

