## Pearson Edexcel

## Mark Scheme (Results)

Summer 2022

Pearson Edexcel International Advanced Level In Chemistry (WCH14)
Paper 01: 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.
- Mark schemes will indicate within the table where, and which strands of QWC, are being assessed. The strands are as follows:
i) ensure that text is legible and that spelling, punctuation and grammar are accurate so that meaning is clear
ii) select and use a form and style of writing appropriate to purpose and to complex subject matter
iii) organise information clearly and coherently, using specialist vocabulary when appropriate


## 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.

## Section A

| Question <br> Number | Answer |
| :--- | :--- | :--- |
| $\mathbf{1 ( a )}$ | The only correct answer is B (two) |
|  | A is not correct because 2-methylpropan-2-ol has a peak for the $3 \mathrm{CH}_{3}$ groups and one for the OH group making 2 in total |
|  | C is not correct because 2-methylpropan-2-ol has a peak for the $3 \mathrm{CH}_{3}$ groups and one for the OH group making 2 in total |
|  | D is not correct because 2-methylpropan-2-ol has a peak for the $3 \mathrm{CH}_{3}$ groups and one for the OH group making 2 in total |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 ( b )}$ | The only correct answer is A (propanal) |  |
|  | B is not correct because propane has 2 peaks in the ratio 3:1 |  |
|  | C is not correct because propan-1-ol has 4 peaks in the ratio 3:2:2:1 <br>  | D is not correct because propan-2-ol has 3 peaks in the ratio is 6:1:1 |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 ( c )}$ | The only correct answer is C (butanal) |  |
|  | A is not correct because butanoic acid has a singlet due to the COOH |  |
|  | B is not correct because butanone has a singlet due to the $\mathrm{CH}_{3}$ adjacent to the $\mathrm{C}=\mathrm{O}$ |  |
|  | D is not correct because butan-1-ol has a singlet due to the OH |  |


| Question <br> Number | Answer | Mark |
| :---: | :---: | :---: |
| 2 | The only correct answer is $\mathbf{B}\left(\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CHO}\right)$ <br> $\boldsymbol{A}$ is not correct because there will not be a peak at m/z 29.0390 <br> C is not correct because there will not be a molecular ion peak a m/z 58.0417 <br> $\boldsymbol{D}$ is not correct because there will not be molecular ion peak at $\mathrm{m} / \mathrm{z} 58.0417$ nor a peak at $\mathrm{m} / \mathrm{z} 29.0390$ | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{3}$ | The only correct answer is D (octan-1-ol, octanal, octane,) |  |
|  | A is not correct because octan-1-ol is the most polar so would have the shortest retention time <br> $\boldsymbol{B}$ is not correct because octan-1-ol is more polar than octanal and so would have a shorter retention time <br> C is not correct because octane is non-polar so would have the longest retention time |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{4}$ | The only correct answer is C (0.75) |  |
|  | B is not correct because the calculation has used the distance from the solvent front to the sample not the baseline <br> B is not correct because the calculation has been inverted |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{5}$ | The only correct answer is C (alkaline hydrolysis of an ester) | (1) |
|  | A is not correct because this reaction will produce a carboxylic acid |  |
|  | B is not correct because this reaction will produce a carboxylic acid |  |
|  | D is not correct because this reaction will produce a carboxylic acid |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{6}$ | The only correct answer is C (3-methylpentan-3-ol) |  |
|  | A is not correct because this is a primary alcohol and so can be formed by the reduction of an aldehyde <br> B is not correct because this is a secondary alcohol and so can be formed by the reduction of a ketone <br> D is not correct because this is a secondary alcohol and so can be formed by the reduction of a ketone |  |


| Question <br> Number | Answer | Mark |
| :---: | :---: | :---: |
| 7 | The only correct answer is B (4.17) <br> A is not correct because this is the -log of the concentration <br> C is not correct because this is the -log of the $K_{a}$ <br> $\mathbf{D}$ is not correct because this is the - log of the $K_{a}$ multiplied by the concentration | (1) |

$\left.\begin{array}{|l|l|c|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Answer } & \text { Mark } \\ \hline \mathbf{8} & \text { The only correct answer is B (13.43) } & \text { (1) } \\ & \text { A is not correct because this is the }-\log \left[\mathrm{OH}^{-}\right] \\ \text {C is not correct because it does not produce } 2 \times \mathrm{OH}^{-} \\ \text {D is not correct because the }-\log \left[\mathrm{OH}^{-}\right] \text {has been added to } \mathrm{p} K_{w}\end{array}\right]$

| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| 9(a) | The only correct answer is D (hydrochloric acid added to ammonia) | (1) |
|  | A is not correct because it is a weak acid and strong base |  |
| $\boldsymbol{B}$ is not correct because it is a strong acid and strong base |  |  |
| C is not correct because it is a weak acid and weak base |  |  |


| Question | Answer | Mark |
| :--- | :--- | :---: |
| Number | The only correct answer is C (methyl red) | (1) |
| 9(b) | A is not correct because malachite green would change colour at about pH 1 |  |
|  | B is not correct because methyl yellow would change colour at about pH 3.5 |  |
| D is not correct because thymol blue would change colour at about pH 9 |  |  |


| Question <br> Number | Answer | Mark |
| :---: | :---: | :---: |
| 10(a) | The only correct answer is C (Graph 3) <br> $\boldsymbol{A}$ is not correct because it is a rate $v$ concentration graph for a second order reaction $\boldsymbol{B}$ is not correct because it is a concentration v time graph for a zero order reaction <br> $\mathbf{D}$ is not correct because it is a rate v concentration graph for a first order reaction | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 0 ( b )}$ | The only correct answer is D (Graph 4) |  |
|  | A is not correct because it is a graph of rate against concentration for a second order reaction |  |
| B is not correct because it is a graph of concentration against time for a zero order reaction |  |  |
|  | C is not correct because it is a graph of rate of reaction against concentration of the reactant for a zero order reaction |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1 ( a )}$ | The only correct answer is A (colorimetry) | (1) |
|  | B is not correct because the solution would not go cloudy |  |
|  | C is not correct because there is no base to titrate against |  |
|  | D is not correct because starch is an indicator and would immediately turn blue-black |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1 ( b )}$ | The only correct answer is B (1.98) |  |
|  | A is not correct because the concentration of the acid has been increased three times |  |
| C is not correct because the concentration of the acid has been decreased six times |  |  |
|  | D is not correct because the pH has been multiplied by three |  |$\quad$| (1) |
| :---: |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 2}$ | The only correct answer is $\mathbf{A}\left(+38.8 \mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ |  |
|  | $\boldsymbol{B}$ is not correct because the units are incorrect |  |
| $\boldsymbol{C}$ is not correct because the gradient has been divided by $R$ |  |  |
|  | $\boldsymbol{D}$ is not correct because the gradient has been divided by $R$ and the units are incorrect |  |$\quad$| (1) |
| :---: |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 3}$ | A is not correct because vanadium oxide is a heterogeneous catalyst <br> B is not correct because decreasing pressure would decrease the equilibrium yield of sulfur trioxide <br> C is not correct because increasing the surface area of the catalyst will affect the rate not the equilibrium yield of <br> sulfur trioxide |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 4}$ | The only correct answer is $\mathbf{A}\left(\mathrm{CaO}(\mathrm{s})<\mathrm{H}_{2} \mathrm{O}(\mathrm{l})<\mathrm{CO}_{2}(\mathrm{~g})<\mathrm{SO}_{2}(\mathrm{~g})\right)$ |  |
|  | B is not correct because $\mathrm{SO}_{2}(g)$ has a greater standard molar entropy than $\mathrm{CO}_{2}(\mathrm{~g})$ |  |
| C is not correct because $\mathrm{SO}_{2}(g)$ has the greatest standard molar entropy |  |  |
| D is not correct because $\mathrm{SO}_{2}(g)$ has the greatest standard molar entropy |  |  |$\quad(1)$


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 5}$ | The only correct answer is B (PS) |  |
|  | A is not correct because $R$ is smaller than $S$ |  |
| C is not correct because $Q$ is larger than $P$ and $R$ is smaller than $S$ |  |  |
|  | D is not correct because $Q$ is larger than $P$ | $(1)$ |

## Section B

| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 16(a)(i) | $K_{\mathrm{p}}=\frac{p^{2} \mathrm{NH}_{3}}{p \mathrm{~N}_{2} p^{3} \mathrm{H}_{2}}$ | Allow round or no brackets <br> Allow upper case <br> Allow pp/PP <br> Allow <br> $p\left(\mathrm{NH}_{3}\right)^{2} \quad p \mathrm{NH}_{3}{ }^{2} \quad$ etc <br> Ignore units even if incorrect <br> Do not award square brackets | (1) |


| Question | Answer | Additional Guidance |  |  | Mark <br> (3) |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 16(a)(ii) | - mole fraction of $\mathrm{N}_{2}$ <br> - mole fraction of $\mathrm{H}_{2}$ <br> - both partial pressures | Example of completed table |  |  |  |
|  |  | Substance | Mole fraction | Partial pressure/atm |  |
|  |  | $\mathrm{N}_{2}$ | 0.18 | 36 |  |
|  |  | $\mathrm{H}_{2}$ | 0.54 | 108 |  |
|  |  | $\mathrm{NH}_{3}$ | 0.28 | 56 |  |
|  |  | TE for M3 on calculated mole fractions multiplied by 200 |  |  |  |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 16(a)(iii) | An answer that makes reference to the following points: <br> - correct use of $K_{p}$ expression <br> (1) <br> - correct answer and 1 or 2 SF <br> - correct units | Example of calculation $56^{2} \div\left(108^{3} \times 36\right)$ <br> 7 or $6.9 \times 10^{-5} / 0.00007$ or 0.000069 <br> Allow 3SF <br> $6.92 \times 10^{-5} / 0.0000692$ <br> Do not award $7.0 \times 10^{-5} / 0.000070$ <br> $\mathrm{atm}^{-2}$ <br> Allow TE from (a)(i) and (a)(ii) <br> If mole fractions are used for the calculation max score 1 for the correct answer and 1-3 SF <br> Correct answer with or without working scores 3 | (3) |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 16(a)(iv) | An answer that makes reference to the following points: <br> - the (forward) reaction is/ must be exothermic <br> - (more ammonia shows that) the equilibrium has moved/shifted to the right <br> OR (more ammonia shows that) a new $K_{\mathrm{p}}$ is established which is larger | Allow reverse argument <br> Allow favours forward reaction/shifts to the product side <br> Allow $K_{\mathrm{p}}$ increases/eqm constant increases <br> Ignore just 'more ammonia produced' or 'yield increases’ | (2) |


| Question <br> Number | Answer | Additional Guidance |
| :--- | :---: | :---: | :---: |
| $\mathbf{1 6 ( b ) ( i )}$ | $\bullet \mathrm{NH}_{4}{ }^{+}(\mathrm{aq})+\mathrm{Cl}^{-}(\mathrm{aq})$ | (1) <br> Do not award $\mathrm{NH}_{4}{ }^{+} \mathrm{Cl}^{-}(\mathrm{aq})$ <br> Do not award $\mathrm{NH}_{4} \mathrm{Cl}(\mathrm{aq})$ |
|  |  | Do not award any other state symbols |


| Question <br> Number | Answer | Additional Guidance |
| :--- | :---: | :--- |
| 16(b)(ii) | $\bullet \Delta_{\text {sol }} H=\Delta_{\text {hyd }} H-$ Lattice Energy | Allow LE for Lattice Energy |
|  | or | Allow $\Delta_{\text {sol }} H=-$ Lattice Energy + hydration enthalpies |
|  | $\Delta_{\text {sol }} H=-$ Lattice Energy $+\Delta_{\text {hyd }} H$ | Allow $\Delta_{\text {sol }} H=-$ Lattice Energy + hydration enthalpies |
|  | Allow $\Delta H_{\text {hyd }}$ etc |  |
| Ignore standard signs |  |  |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 16(b)(iii) | - enthalpy change of hydration of ammonium chloride <br> - enthalpy change of solution | (1) <br> (1) | Example of calculation $-307+(-378)=-685\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> Allow (kJ mol${ }^{-}$) $705+(-685)=(+) 20\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> Allow TE on arithmetical errors <br> Do not award use of incorrect expression <br> Correct answer with or without working scores 2 <br> Units are not required but if wrong penalise only once. | (2) |



| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 16(c) | - $\mathrm{NH}_{4}{ }^{+}+\mathrm{H}_{2} \mathrm{O} \longrightarrow \mathrm{NH}_{3}+\mathrm{H}_{3} \mathrm{O}^{+}$ | $\begin{aligned} & \text { Accept } \mathrm{NH}_{4}^{+} \longrightarrow \mathrm{NH}_{3}+\mathrm{H}^{+} \\ & \text {Allow eqm sign } \\ & \text { Ignore } \mathrm{NH}_{4} \mathrm{Cl} \longrightarrow \mathrm{NH}_{4}^{+}+\mathrm{Cl}^{-} \\ & \text {Ignore } \mathrm{NH}_{4} \mathrm{Cl}+\mathrm{aq} \longrightarrow \mathrm{NH}_{4}^{+}+\mathrm{Cl}^{-} \\ & \text {Ignore state symbols even if incorrect } \\ & \text { Do not award } \mathrm{NH}_{4} \mathrm{Cl}+\mathrm{aq} \longrightarrow \mathrm{NH}_{3}+\mathrm{HCl} \\ & \text { Do not award } \mathrm{NH}_{4} \mathrm{Cl} \longrightarrow \mathrm{NH}_{3}+\mathrm{HCl} \\ & \text { Do not award } \mathrm{NH}_{4} \mathrm{Cl} \longrightarrow \mathrm{NH}_{3}+\mathrm{H}^{+}+\mathrm{Cl}^{-} \\ & \text {Do not award } \mathrm{NH}_{4}^{+}+\mathrm{OH}^{-} \longrightarrow \mathrm{NH}_{3}+\mathrm{H}_{2} \mathrm{O} \end{aligned}$ | (1) |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 17(a) | An answer that makes reference to the following points: <br> - $\mathbf{A}$ <br> - B <br> - C | Allow structural or skeletal formulae for max 2 marks <br> Ignore any names even if incorrect <br> Ignore bond angles/lengths <br> Penalise missing Hs only once | (3) |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :--- | :--- | :--- | :---: |
| 17(b)(i) | - 2-hydroxybutanenitrile | Allow 2-hydroxy(l)buta(n)nitrile <br> Allow 2-hydroxy(l)butane-1-nitrile | (1) |
|  |  | Do not award 2-hydroxobutanenitrile <br> Do not award 2-oxobutanenitrile <br> Do not award cyanides or other non IUPAC names |  |


| Question <br> Number | Answer | Additional Guidance |
| :--- | :--- | :--- | :---: |
| 17(b)(ii) | (1) <br> one isomer rotates (the plane of monochromatic) <br> plane-polarised light in one direction and the other in <br> the opposite direction/ the isomers rotate (the plane <br> of) plane-polarised light in opposite <br> directions/clockwise and anticlockwise | Do not award bends <br> Allow different directions <br> Allow for plane polarised light |
| Allow the direction of rotation of plane polarised light |  |  |
| Allow see which way the sample rotates PPL |  |  |


| Question <br> Number | Answer | Additional Guidance | Mark |  |
| :--- | :--- | :--- | :--- | :---: |
| 17(b)(iii) | An answer that makes reference to the following points: | (1) | Do not award just propanal is planar <br> Do not award planar intermediate/carbocation <br> Do not award any reference to nucleophilic <br> substitution (S $\left.\mathrm{S}_{\mathrm{N}} 1 / \mathrm{S}_{\mathrm{N}} 2\right)$ | (2) |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 17(c)(i) |  | Ignore displayed or structural formulae Ignore bond lengths and bond angles <br> Allow <br> Ignore connectivity if vertical bond | (1) |


| Question <br> Number | Answer | Additional Guidance |
| :--- | :--- | :--- | :--- |
| 17(c)(ii) | Mark <br> - no carbon atom has 4 different groups <br> (central) carbon atom is bonded to two CH3/same <br> groups <br> or <br> no asymmetric/chiral carbon atom <br> or <br> the compound is superimposable on its mirror <br> image <br> or <br> it does not have a chiral centre | (1) <br> Ignore symmetrical <br> attached to the carbon atom |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 17(d)(i) | - correct chemical shift and carbon environment |  | (1) |
|  |  | Chemical shift range Carbon environment |  |
|  |  | 190-225 (ppm) C=O |  |
|  |  | OR |  |
|  |  | 0-60 (ppm) $\quad$ C-C |  |
|  |  | Both range and carbon environment required. Allow the full range or a number/ smaller range within the range. |  |
|  |  | Ignore any splitting patterns |  |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 17(d)(ii) | • Propanal: 3/three |  | (1) |
|  | and |  |  |
|  | Propanone: 2/two |  |  |

(Total for Question 17 = 11 Marks)


The following table shows how the marks should be awarded for structure and lines of reasoning.

|  | Number of marks awarded for <br> structure of Answer and <br> sustained lines of <br> reasoning |
| :--- | :--- |
| Answer shows a coherent <br> logical structure with linkages <br> and fully sustained lines of <br> reasoning demonstrated <br> throughout | 2 |
| Answer is partially structured <br> with some linkages and lines of <br> reasoning | 1 |
| Answer has no linkages <br> between points and is <br> unstructured | 0 |



| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 19(a)(i) | An answer that makes reference to the following point: <br> - $\mathrm{HCOOH}+\mathrm{KOH} \longrightarrow \mathrm{HCOOK}+\mathrm{H}_{2} \mathrm{O}$ | Allow $\mathrm{HCOO}^{-} \mathrm{K}^{+} / \mathrm{HCOO}^{-}+\mathrm{K}^{+}$ <br> Allow $\mathrm{HCOOH}+\mathrm{OH}^{-} \longrightarrow \mathrm{HCOO}^{-}+\mathrm{H}_{2} \mathrm{O}$ <br> Allow Na in place of K <br> Ignore state symbols even if incorrect <br> Do not award HCOO - K | (1) |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 19(a)(ii) | - correct volume read off the graph <br> - correct concentration | (1) <br> (1) | Example of calculation $22\left(\mathrm{~cm}^{3}\right)$ <br> This may be noted on the graph $25.0 \times 0.15 / 22.0=0.17045\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> Ignore SF except 1SF <br> Allow TE on wrong volume <br> Correct answer scores 2 | (2) |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 19(a)(iii) | An answer that makes reference to the following points: <br> - volume at half-neutralisation <br> - pH value at half-neutralisation <br> - calculation of $K_{\mathrm{a}}$ | $11 \mathrm{~cm}^{3}$ (Allow TE from volume in (a)(ii)) $\mathrm{pH}=3.8( \pm 0.1)$ <br> (Hydrogen ion concentration $=10^{-3.8}$ ) $K_{\mathrm{d}}=1.5849 \times 10^{-4} / 0.00015849\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ Correct answer with no working scores 3 <br> Allow TE throughout Ignore SF <br> If 3.9 used <br> (Hydrogen ion concentration $=10^{-3.9}$ ) <br> $K_{\mathrm{a}}=1.2589 \times 10^{-4} / 0.00012589\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> If 3.7 used <br> (Hydrogen ion concentration $=10^{-3.7}$ ) <br> $K_{\mathrm{a}}=1.9953 \times 10^{-4} / 0.00019953\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> Allow TE from wrong pH | (3) |



|  | Alternative method 2 (using pH of the methanoic acid at the start) <br> - pH at the start <br> - convert pH into $\mathrm{H}^{+}$concentration <br> - calculation of $K_{\mathrm{a}}$ <br> 2.0 gives a value of $6.667 \times 10^{-4}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> 2.1 gives a value of $4.206 \times 10^{-4}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> 2.2 gives a value of $2.654 \times 10^{-4}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> 2.3 gives a value of $1.674 \times 10^{-4}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> 2.4 gives a value of $1.057 \times 10^{-4}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ <br> 2.5 gives a value of $3.162 \times 10^{-3}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ | 2.3 <br> Allow 2.0-2.5 $\begin{aligned} & \text { Hydrogen ion concentration }=10^{-2.3} \\ & =5.0119 \times 10^{-3} / 0.0050119\left(\mathrm{~mol} \mathrm{dm}^{-3}\right) \\ & K_{\mathrm{a}}=\frac{\left(5.0119 \times 10^{-3}\right)^{2}}{0.15}=1.6746 \times 10^{-4} / 0.00016746\left(\mathrm{~mol} \mathrm{dm}^{-3}\right) \end{aligned}$ <br> Correct answer with no working scores 3 <br> Allow TE from wrong pH (i.e. not in the range of 2.0-2.5) <br> Ignore SF |
| :---: | :---: | :---: |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 19(b) | An answer that makes reference to the following points: <br> - calculation of $\left[\mathrm{H}^{+}\right]$ <br> (1) <br> - correct ratio <br> (1) | Example of calculation $\left[\mathrm{H}^{+}\right]=K_{\mathrm{a}} \times \frac{[\mathrm{HA}]}{\left[\mathrm{A}^{-}\right]} \text {OR } \frac{\left[\mathrm{H}^{+}\right]}{K_{\mathrm{a}}}=\frac{[\mathrm{HA}]}{\left[\mathrm{A}^{-}\right]}$ $2.5119 \times 10^{-5} / 0.000025119$ $\frac{[\mathrm{HA}]}{\left[\mathrm{A}^{-}\right]}=\frac{2.5118864 \times 10^{-5}}{1.3 \times 10^{-5}}=1.9322: 1$ <br> Correct answer with no working scores 2 <br> Allow just 1.9322 <br> Allow rounding to 2:1 <br> Ignore SF <br> Reciprocal ratio correctly identified $0.5175: 1$ scores 2 Correct answer with no working scores 2 <br> Allow Henderson-Hasselbach equation $\begin{align*} & \mathrm{pH}=\mathrm{p} K \mathrm{a}-\log \frac{[\mathrm{HA}]}{\left[\mathrm{A}^{-}\right]} \\ & 4.6=4.8861-\log \frac{[\mathrm{HA}]}{\left[\mathrm{A}^{-}\right]}  \tag{1}\\ & {[\mathrm{HA}]=1.9322: 1} \tag{1} \end{align*}$ <br> [ $\mathrm{A}^{-}$] <br> Allow just 1.9322 <br> Ignore SF <br> Reciprocal ratio correctly identified $0.5175: 1$ scores 2 | (2) |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 20(a)(i) | An answer that makes reference to two of the following points: <br> - 2-bromobutane: first order as doubling the concentration (in experiments 1 and 2 where $\mathrm{OH}^{-}$is constant) the rate doubles <br> - hydroxide ions: zero order as doubling the concentration (in experiments 1 and 3 where 2-bromobutane is constant) the rate does not change <br> OR <br> hydroxide ions: zero order as doubling the concentration (in experiments 2 and 3) where the concentration of 2-bromobutane is halved the rate halves. | Two correct orders with no or incorrect reasoning scores 1 <br> Note the reasoning can be shown on the table | (2) |


| Question <br> Number | Answer | Additional Guidance |
| :--- | :--- | :--- | :--- |
| 20(a)(ii) | $\bullet$ rate $/ \mathrm{r}=k\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]$ | (1) |
|  |  | TE on (i) <br> Allow displayed or structural formulae <br> Allow rate $=k\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]^{1}\left[\mathrm{OH}^{-}\right]^{0}$ <br> Allow upper case K <br> Allow reactants in any order <br> Do not award round brackets |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 20(a)(iii) | An answer that makes reference to the following points: <br> - correct calculation <br> - correct units | Allow the calculation from any experiment Example of calculation from experiment 1 $1.01 \times 10^{-3} / 0.100=0.0101 / 1.01 \times 10^{-2}$ <br> TE on (ii) Ignore SF $\mathrm{s}^{-1}$ <br> Allow s- <br> TE on (ii) | (2) |




## Section C

| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 21(a)(i) | - correct use of enthalpy data <br> - correct enthalpy change | (1) <br> (1) | Example of calculation $\begin{aligned} & -(-824.2)+(3 \times-110.5) \\ & =(+) 492.7\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right) \end{aligned}$ <br> Correct answer with or without working scores 2 The following score 1 for a single error: <br> $(+) 713.7\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ not x 3 <br> $-492.7\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ signs reversed <br> Allow 3SF <br> Penalise wrong units once only in (a)(i) and (ii) | (2) |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 21(a)(ii) | - $\quad \sum S$ products <br> - $\quad \sum S$ reactants <br> - $\Delta S_{\text {system }}=\sum S$ products $-\sum S$ reactants | (1) <br> (1) <br> (1) | Example of calculation $\begin{aligned} & S=\text { products }(2 \times 27.3)+(3 \times 197.6)= \\ & 647.4\left(\mathrm{~J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1}\right) \end{aligned}$ $S=\text { reactants } 87.4+(3 \times 5.7)=104.5\left(\mathrm{~J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1}\right)$ $\Delta S_{\text {system }}=647.4-104.5=(+) 542.9\left(\mathrm{~J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1}\right)$ <br> Correct answer with no working scores 3 <br> Allow TE for M3 | (3) |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 21(a)(iii) | - use of $\Delta S_{\text {surroundings }}=-\frac{\Delta H}{T}$ <br> - at equilibrium $\Delta S_{\text {total }}=0=\Delta S_{\text {surroundings }}+\Delta S_{\text {system }}$ <br> - calculation of temperature | (1) <br> (1) <br> (1) | Example of calculation $\begin{aligned} & \Delta S_{\text {surroundings }}=-(+492.7) \times 1000 / T \\ & 0=-492.7 \times 1000 / T+542.9 \\ & \geq 907.53(\mathrm{~K}) \\ & 0.90753 \text { scores } 2(\text { not } \times 1000) \\ & \text { Ignore SF } \\ & \text { TE on (a)(i) and (a)(ii) } \end{aligned}$ <br> Correct answer based on ai and aii without working scores 3 <br> Allow use of $\Delta G=\Delta H-T \Delta S_{\text {system }}$ | (3) |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 21(b)(i) | An answer that makes reference to the following points: <br> - $\Delta S_{\text {surroundings }}$ and $\Delta S_{\text {system }}$ are positive <br> - so $\Delta S_{\text {total }}$ will always be positive (so reaction will be feasible) <br> OR <br> Using $\Delta G=\Delta H-T \Delta S_{\text {system }}$ <br> - Allow $\Delta H$ is negative and $\Delta S_{\text {system }}$ is positive <br> - so $\Delta G$ will always be negative (so reaction will be feasible) | (1) <br> (1) <br> (1) <br> (1) | Allow $\Delta H$ is negative/reaction exothermic and $\Delta S_{\text {system }}$ is positive <br> M2 dependent on M1 <br> M2 dependent on M1 | (2) |


| Question <br> Number | Answer | Additional Guidance |  |
| :--- | :---: | :---: | :---: |
| 21(b)(ii) | An answer that makes reference to the following points: |  |  |
|  | • at a higher temperature $\Delta S_{\text {surroundings will decrease }}$ | (1) | Ignore reference to $\Delta S_{\text {total }}=$ Rlnk |
|  | • $\Delta S_{\text {system }}$ does not change (significantly) | (1) |  |
|  | • so $\Delta S_{\text {total }}$ will decrease/become less positive | (1) |  |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 21(c)(i) | All 7 correct 3 marks <br> $4-6$ correct 2 marks <br> $2-3$ correct 1 mark | G | Allow values instead of letters | (3) |



| Question <br> Number | Answer | Additional Guidance | Mark |  |
| :--- | :---: | :---: | :--- | :---: |
| 21(c)(iii) | - the electron is being added to a negative ion | (1) | This can be shown by an equation | (2) |
|  | • and so there is repulsion (so energy is required) | (1) | Allow repulsion between the electrons |  |

## (Total for Question $21=20$ Marks) <br> (Total for Section C = 20 Marks) <br> TOTAL FOR PAPER $=\mathbf{9 0}$ MARKS

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