## Mark Scheme (Results)

October 2020

Pearson Edexcel International Advanced Level In Chemistry (WCH05)
Paper 1: General Principles of Chemistry II Transition Metals and Organic Nitrogen Chemistry (including synoptic assessment)

## Edexcel and BTEC Qualifications

Edexcel and BTEC qualifications are awarded by Pearson, the UK's largest awarding body. We provide a wide range of qualifications including academic, vocational, occupational and specific programmes for employers. For further information visit our qualifications websites at www.edexcel.com or www.btec.co.uk. Alternatively, you can get in touch with us using the details on our contact us page at www.edexcel.com/contactus.

## Pearson: helping people progress, everywhere

Pearson aspires to be the world's leading learning company. Our aim is to help everyone progress in their lives through education. We believe in every kind of learning, for all kinds of people, wherever they are in the world. We've been involved in education for over 150 years, and by working across 70 countries, in 100 languages, we have built an international reputation for our commitment to high standards and raising achievement through innovation in education. Find out more about how we can help you and your students at: www.pearson.com/uk

October 2020
Publications Code WCH05_01_2010_MS
All the material in this publication is copyright
© Pearson Education Ltd 2020

## 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.


## Section A (multiple choice)

| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1}$ | The only correct answer is C (fuming sulfuric acid) | (1) |
|  | $\boldsymbol{A}$ is not correct because this mixture is for the nitration of benzene |  |
|  | $\boldsymbol{B}$ is not correct because this acid does not contain the electrophile |  |
|  | D is not correct because this gaseous mixture will not work |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{2}$ | The only correct answer is C (burns with a smoky flame) | (1) |
|  | $\mathbf{A}$ is not correct because bromine water is not decolourised |  |
|  | B is not correct because the mixture remains orange <br> D is not correct because this is not evidence for unsaturation |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{3}$ | The only correct answer is B | (1) |
|  | A is not correct because bromine does not substitute at every available <br> position on the ring <br> C is not correct because excess bromine substitutes at positions 2,4 and 6 |  |


| Question <br> Number | Correct Answer | Mark |
| :---: | :---: | :---: |
| 4 | The only correct answer is D <br> A is not correct because this is a cyclic amide <br> $\mathbf{B}$ is not correct because this is an amide <br> $\mathbf{C}$ is not correct because this is a substituted amide | (1) |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{5 ( a )}$ | The only correct answer is A (HNO2) | (1) |
|  | B is not correct because this is the wrong acid <br> $\boldsymbol{C}$ is not correct because these acids are used for nitration not diazotisation <br> D is not correct because this acid can be present to stabilise the diazonium <br> salt but not to produce it |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| 5(b) | The only correct answer is D (ice-cold alkaline solution) |  |
|  | A is not correct because the temperature is too high and the diazo <br> compound would break down | (1) |
| B is not correct because halogen carriers are used for substitution of the |  |  |
| benzene ring and not azo dye formation |  |  |
| C is not correct because the azo dye is not formed from reduction |  |  |$\quad$


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{6}$ | The only correct answer is B $\left(\mathrm{CH}_{2} \mathrm{CHCONH}_{2}\right)$ | (1) |
|  | $\boldsymbol{A}$ is not correct because there is no $C=C$ for addition polymerisation |  |
| C is not correct because an addition and/or condensation polymer could |  |  |
| form |  |  |
| D is not correct because this molecule would not polymerise |  |  |$\quad$.


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{7 ( a )}$ |  | The only correct answer is A <br> B is not correct because there is no - NH-in the backbone of the chain <br> a single-bonded oxygen in inside the brackets |
| D is not correct because the single-bonded oxygen should not be inside the <br> brackets |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 7(b) |  | (1) |
|  | The only correct answer is $\mathbf{C}$ <br> acyl chloride |  |
| B is not correct because in excess both amine groups will react |  |  |
| Dis not correct because the chlorine does not replace the OH to make an <br> acyl chloride |  |  |

$\left.\begin{array}{|l|l|c|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Correct Answer } & \text { Mark } \\ \hline \text { 7(c) } & \text { The only correct answer is C (ninhydrin). } & \text { (1) } \\ & \boldsymbol{A} \text { is not correct because this is used to detect aldehydes and ketones } \\ & \boldsymbol{B} \text { is not correct because this is used to detect starch } \\ \boldsymbol{D} \text { is not correct because this is used to detect aldehydes }\end{array}\right)$

| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| 7(d) | The only correct answer is C (proton transfer within the molecules results <br> in ionic bonding) |  |
|  | A is not correct because London forces would not result in these properties <br> B is not correct because hydrogen bonding would not result in these <br> properties |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{8}$ | The only correct answer is B (aluminium forms a stable oxide coating) | (1) |
|  | A is not correct because aluminium does react under suitable conditions | C is not correct because reference to non-standard conditions does not <br> explain why it is corrosion-resistant |
| D is not correct because the reaction is thermodynamically feasible |  |  |$\quad$.


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{9}$ | The only correct answer is D (made from fossil fuels) | $\mathbf{( 1 )}$ |
|  | $\mathbf{A}$ is not correct because liquid methanol is easily stored |  |
| B is not correct because liquid methanol is easily transported |  |  |
| $\mathbf{C}$ is not correct because all fuels are flammable |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 0}$ | The only correct answer is D (promotion of electrons between split d <br> orbitals) | (1) |
|  | A is not correct because electrons are not promoted to a higher quantum <br> shell <br> B is not correct because this is what happens in flame tests |  |
| C is not correct because one orbital is not split but rather the d sub-shell |  |  |$\quad$|  |
| :--- |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1 ( a )}$ | The only correct answer is B (ligand exchange) | (1) |
|  | $\boldsymbol{A}$ is not correct because there is no acidic and basic behaviour |  |
| C is not correct because no oxidation has occurred |  |  |
|  | D is not correct because no reduction has occurred |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1 ( b )}$ | The only correct answer is D (6, 4) | (1) |
|  | $\boldsymbol{A}$ is not correct because this is ratio of the two chromium ions |  |
| B is not correct because this is the numerical values of the charges |  |  |
| C is not correct because the coordination number is different in the two <br> species |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 2}$ | The only correct answer is C (the process is gentler with temperature- <br> sensitive substances). <br> $\boldsymbol{A}$ is not correct because steam distillation is more expensive due to the <br> additional equipment needed <br> $\boldsymbol{B}$ is not correct because the neutrality of water is not an issue | (1) |
|  | D is not correct because a safety valve is needed to release excess steam |  |$\quad$.


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 3}$ | The only correct answer is D (phenylamine < ammonia < butylamine) | (1) |
|  | A is not correct because ammonia is a weaker base than butylamine <br> $\boldsymbol{B}$ is not correct because phenylamine is a weaker base than ammonia <br> C is not correct because butylamine is the strongest base/phenylamine is <br> the weakest base |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 4}$ |  |  |
|  | The only correct answer is A |  |
|  | B is not correct because ethane-1,2-diamine is bidentate |  |
| C is not correct because the acetyl acetonate ion is bidentate |  |  |
| D is not correct because the ethanedioate ion is bidentate |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 5}$ | The only correct answer is A (directly proportional - directly <br> proportional | (1) <br> B is not correct because $\ln K_{c}$ is directly proportional to $E^{\theta}$ cell <br> C is not correct because $\Delta S_{\text {system }}$ is directly proportional to $E_{\text {cell }}$ <br> $\mathbf{D}$ is not correct because both $\Delta S_{\text {system }}$ and ln $K_{c}$ are directly proportional <br> to $E^{\theta}$ cell |

## Section B

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(a)(i) | Diagram to include: <br> Labelled salt bridge which must dip into solutions <br> Allow filter paper soaked with $\mathrm{KNO}_{3}$ <br> Allow just $\mathrm{KNO}_{3} / \mathrm{NaNO}_{3}$ <br> Voltmeter /V <br> Completed circuit with gold and silver electrodes Ignore absence of $1 \mathrm{~mol} \mathrm{dm}^{-3}$ <br> Allow just 'silver ions <br> Example of diagram <br> ALLOW half cells to be drawn the opposite way round | e.g Gold(I) <br> sulfate <br> $\mathrm{Au}^{+}$ <br> AgCl <br> Gold ions <br> With other <br> metal ions <br> Incorrect <br> temperatures <br> Electrodes <br> not dipping <br> into the <br> solution | (3) |

$\left.\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline \mathbf{1 6 ( a ) ( i i )} & (\text { Equation }) \\ & \mathrm{Au}^{3+}(\mathrm{aq})+3 \mathrm{Ag}(\mathrm{s}) \rightarrow \mathrm{Au}(\mathrm{s})+3 \mathrm{Ag}^{+}(\mathrm{aq}) & \text { (1) } & \\ E^{\theta} \text { cell }(=+1.41-0.80)=(+) 0.61(\mathrm{~V}) & \text { (2) } \\ \text { Standalone marks }\end{array}\right)$.

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i i i )}$ | Either <br> The electrode potential/ $E^{\theta}$ for the reduction of gold(I) is <br> more positive than that for the oxidation of gold(I) <br> Allow reduction $E^{\theta}>$ oxidation $E^{\theta}$ <br> Or <br> $E_{\text {cell }}(=+1.69-1.29)=(+) 0.4(0)(V)$ which is positive so <br> the reaction is feasible <br> Allow answers using the anticlockwise rule. | Reference <br> to Ag <br> scores $(0)$ | (2) |
| (Equation) $3 \mathrm{Au}^{+} \rightarrow \mathrm{Au}^{3+}+2 \mathrm{Au}$ | (1) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( b ) ( i ) ~}$ | Linear (shape) <br> ALLOW unambiguous incorrect spellings such as <br> lenear/linnear | (1) |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( b ) ( i i ) ~}$ | Zinc is the reducing reagent because it is oxidised/loses <br> electrons <br> Allow just <br> Zinc is the reducing agent/ is oxidised/ zinc loses electrons/ <br> Zinc's oxidation number or state increases/ goes from 0 to <br> +2 | (1) | (2) |
| (Half equation) $\mathrm{Zn}+4 \mathrm{CN}^{-} \rightarrow[\mathrm{Zn}(\mathrm{CN}) 4]^{2-}+2 \mathrm{e}(-)$ <br> Allow <br> $\mathrm{Zn} \rightarrow \mathrm{Zn}^{2+}+2 \mathrm{e}(-)$ <br> Ignore state symbols even if incorrect | (1) |  |  |

(Total for Question 16 = 10 Marks)

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( a )}$ | A metal which forms one (or more stable) ions with an <br> incompletely/partially filled d orbital/d subshell <br> Ignore reference to different oxidation states/numbers <br> Ignore reference to specific quantum levels e.g. 3 | Just <br> 'compounds' <br> shell | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( b )}$ | $4 \mathrm{~s}^{1}$ | Answer <br> given with <br> any other <br> configuration <br> such as <br> $3 \mathrm{~d}^{5} / 3 \mathrm{~d}^{10} /[\mathrm{Ar}]$ | $\mathbf{( 1 )}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( c )}$ | Any two from $\mathrm{CrO}, \mathrm{Cr}_{2} \mathrm{O}_{3}, \mathrm{CrO}_{3}$ <br> ALLOW <br> $\mathrm{CrO}_{2}$ | $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2}$ <br> $\mathrm{CrO}_{4}{ }^{-}$ | $\mathbf{( 1 )}$ |


| Question <br> Number | Acceptable Answers | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( d )}$ | (M1) (The green precipitate formed by) $\mathrm{Cr}^{3+}$ will dissolve in <br> excess sodium hydroxide solution to form a green solution <br> (1) <br> (M2) (The green precipitate formed by) $\mathrm{Fe}^{2+}$ turns brown on <br> exposure to air <br> (and doesn't dissolve in excess sodium hydroxide ) <br> (M3) (The green precipitate formed by) $\mathrm{Ni}^{2+}$ doesn't dissolve <br> in excess sodium hydroxide | (1) |
|  | ALLOW alternative method based on adding ammonia | (1) |
| (M1) (The green precipitate formed by) $\mathrm{Cr}^{3+}$ will dissolve in |  |  |
| excess ammonia (solution) to form a green solution |  |  |
| (M2) (The green precipitate formed by) $\mathrm{Fe}^{2+}$ doesn't dissolve | (1) | (1) |
| in excess ammonia (solution) |  |  |
| (M3) (The green precipitate formed by) $\mathrm{Ni}^{2+}$ will dissolve in |  |  |
| excess ammonia (solution) to form a blue solution | (1) |  |

$\left.\begin{array}{|l|lr|l|l|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline \mathbf{1 7 ( e ) ( i )} & 3 \mathrm{D} \text { image } \\ \text { Bonds clearly to the } \mathrm{N} \text { of the ammonia molecules } \\ \text { Examples of diagrams } & \text { (1) } & \begin{array}{l}\text { (2) } \\ \mathrm{NH}_{4} \\ \text { Incorrect } \\ \text { charge } \\ \text { Coordinate } \\ \text { bond in } \\ \text { the wrong } \\ \text { direction }\end{array} \\ \text { Ignore absence of square brackets and/or charge } \\ \text { Ignore absence or presence of lone pair on nitrogen }\end{array}\right]$
$\left.\begin{array}{|l|l|l|c|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline \mathbf{1 7 ( e ) ( i i ) ~} & \begin{array}{l}\text { (Bond angle) } 109.5^{\circ} \\ \text { ALLOW } 109-110^{\circ} \\ \begin{array}{l}\text { Four areas of (bonded) electron density repel } \\ \text { (to the maximum extent/minimal repulsion) }\end{array} \\ \hline\end{array} & \text { (1) }\end{array}\right)$

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( e ) ( \text { iii) }}$ | (The ligand exchange reaction results in) <br> an increase in moles/ moles change from 2 to $7 /$ <br> $\left[\mathrm{Cr}\left(\mathrm{NH}_{3}\right)_{6}\right]^{3+}+$ EDTA $^{4} \rightarrow[\mathrm{Cr}(\text { EDTA })]^{-}+6 \mathrm{NH}_{3}$ <br> Allow <br> Particles for 'moles' <br> The entropy (of the system) increases/ $\Delta S_{\text {system is positive }}$ <br> Ignore references to stability of complex |  | (2) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :---: | :---: |
| $\mathbf{1 7 ( f ) ( i )}$ | $2 \mathrm{CrO}_{4}^{2-}+2 \mathrm{H}^{+} \rightarrow \mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}+\mathrm{H}_{2} \mathrm{O}$ |  | $\mathbf{( 1 )}$ |
|  | ALLOW $\rightleftharpoons$ <br> Ignore state symbols even if incorrect |  |  |


| Question <br> Number | Acceptable Answers | Mark |
| :---: | :---: | :---: |
| 17(f)(ii) | Suitable calculation such as $\begin{equation*} \mathrm{n}\left(\mathrm{Fe}^{2+}\right)=(0.0492 \times 0.025=) 1.23 \times 10^{-3} / 0.00123(\mathrm{~mol}) \tag{1} \end{equation*}$ $\text { 6:1 ratio so } \mathrm{n}\left(\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}\right)=\left(1.23 \times 10^{-3} \div 6\right)$ $\begin{equation*} =2.05 \times 10^{-4}(\mathrm{~mol}) / 0.000205(\mathrm{~mol}) \text { in } 10.80 \text { titre } \tag{1} \end{equation*}$ $\begin{equation*} \mathrm{n}\left(\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}\right)=2.05 \times 10^{-4} \times(250 \div 10.80)=4.7454 \times 10^{-3} \tag{1} \end{equation*}$ <br> (mol)/ $0.0047454(\mathrm{~mol})$ $\begin{align*} & \mathrm{n}\left(\mathrm{CrO}_{4}^{2-}\right)=\left(4.7454 \times 10^{-3} \times 2\right)=9.4907 \times 10^{-3} / 0.0094907(\mathrm{~mol}) \\ & \mathrm{m}\left(\mathrm{Na}_{2} \mathrm{CrO}_{4}\right)=\left(9.4907 \times 10^{-3} \times 162\right)=1.5375(\mathrm{~g})  \tag{1}\\ & \%\left(\mathrm{Na}_{2} \mathrm{CrO}_{4}\right)=(1.5375 \times 100 \div 1.59)=96.698 \tag{1} \end{align*}$ <br> Final answer without working scores (6) <br> Allow TE at each stage <br> M6 not awarded for an answer greater than $100 \%$ <br> Ignore SF except 1 SF or $>5 \mathrm{SF}$ | (6) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( g )}$ | Reference to the mechanism of action involving: <br> Adsorption to the surface of the catalyst <br> Allow active site(s) for 'surface' <br> Weakens the bonds in the reactants / <br> (chemical) reaction at the surface of the catalyst (1) | Absorption | (3) |
| Desorption from the catalyst surface |  |  |  |
| Max (2) if the catalyst surface is not mentioned <br> Allow (1) rescue mark if none other scored for general <br> catalyst definition such as: lowers the activation energy <br> by providing an alternative reaction therapy pathway |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( a )}$ | $\mathrm{n}\left(\mathrm{CO}_{2}\right)=7.20 \div 24=0.30(\mathrm{~mol})$ | (1) |  |
|  | $\mathrm{n}\left(\mathrm{H}_{2} \mathrm{O}\right)=4.50 \div 18=0.25(\mathrm{~mol})$ | (1) |  |
|  | Ratio $\mathrm{C}: \mathrm{H}=6: 10=3: 5$ | (1) |  |
|  |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(b) |  |  | (2) |
|  | X |  |  |
|  | Y |  |  |
|  | Penalise use of non-displayed formulae once only |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( c ) ( i ) ~}$ | The OH and Br must be bonded to adjacent carbons |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |  |
| :--- | :--- | :---: | :--- | :---: |
| $\mathbf{1 8 ( c ) ( i i ) ~}$ | Use of a polarimeter/use of plane polarised light | (1) |  | (2) |
|  | Rotation of the plane (of plane polarised light) | (1) | Bend for <br> rotate |  |

\(\left.$$
\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\
\text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\
\hline \text { 18(d) } & \begin{array}{l}\text { Points can be made in either order } \\
\text { Absorption 1 } \\
\text { C-H 3095-3010 }\left(\mathrm{cm}^{-1}\right) \\
\text { or } \\
\text { C-H 2962-2853 }\left(\mathrm{cm}^{-1}\right) \\
\text { or } \\
\text { C-H 1485-1365 }\left(\mathrm{cm}^{-1}\right)\end{array} & \begin{array}{l}\text { Penalise } \\
\text { single values } \\
\text { once only }\end{array} & \text { (2) } \\
\begin{array}{ll}\text { Absorption 2 } \\
\text { C=C 1669-1645 }\left(\mathrm{cm}^{-1}\right)\end{array}
$$ \& (1) \& \& <br>

Two correct absorptions without bonds scores \& (1)\end{array}\right]\)|  |
| :--- |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( e ) ( i )}$ | Hexanedioic acid <br> ALLOW <br> Hexane-1,6-dioic acid/hexan-1,6-dioic acid/ hexandioic <br> acid/ <br> hexanedioic acid/1,6-dihexanoic acid <br> Ignore punctuation | Missing 'di' | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(e)(ii) | A suitable diamine such as <br> Hexane-1,6-diamine <br> Amide bond/CONH <br> A suitable equation with polymer having continuation bonds <br> Example of answer <br> ( <br> Ignore brackets and ' $n$ ' <br> Accept displayed, structural formulae or any combination thereof | C-O-N-H <br> Bond from C to H in amide | (3) |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(e)(iii) | First two curly arrows <br> Second two curly arrows <br> ALLOW (1) for any two correct curly arrows <br> Ignore dipoles and lone pairs even if incorrect <br> Exemplar answer |  | (2) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 18(e)(iv) | OH <br> ALLOW <br> Any diagram which clearly indicates 'trans' configuration <br> The hydroxy groups are on opposite side of C=C and so unlikely to <br> condense <br> ALLOW <br> The OH groups are too far away (from each other) | (1) |  |

(Total for Question $18=18$ marks)
(Total for Section $\mathbf{B}=\mathbf{5 0}$ marks)

## Section C

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 9 ( a )}$ | Carboxylic (acid) | Just ‘COOH' <br> Combination with <br> any other functional <br> group | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 9 ( b )}$ | $\mathrm{C}_{13} \mathrm{H}_{18} \mathrm{O}_{2}$ |  |  |
| Allow elements in any order |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(c)(i) | Curly arrow from or within the circle of the benzene ring to the $\mathrm{C}+$ <br> (1) <br> Intermediate structure with horseshoe covering at least 3 carbons and facing the tetrahedral carbon with some part of the positive charge being within the horseshoe <br> (1) <br> Curly arrow from the $\mathrm{C}-\mathrm{H}$ to anywhere inside the benzene ring reforming the delocalised structure of molecule $\mathbf{A}$ to form $\mathrm{H}^{+}$ <br> ALLOW HCl for $\mathrm{H}^{+}$ <br> (1) <br> Ignore regeneration of the catalyst | Curly arrow on outside of hexagon <br> Partial bonds to H or $\mathrm{CH}_{3}$ except for dot and wedge in 3D structure | (3) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9 ( c ) ( i i ) ~}$ | $\mathrm{CH}_{3}$ <br> $\mathrm{H}_{3} \mathrm{C}-\mathrm{CH}-\mathrm{C}^{\prime}$ |  | (1) |
|  | Accept any formulae or combination of, except <br> 2-methylpropanoyl chloride <br> molecular formula | 2-methylpropanyl <br> chloride |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 19(c)(iii) | It would produce an alcohol/OH/hydroxy(l) <br> (and not a hydrocarbon) |  | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | ---: | :--- | :---: |
| *19(d)(i) | The lone pair of electrons on the oxygen are <br> incorporated into the delocalised ring of electrons <br> This increases the susceptibility of the ring to <br> electrophilic attack <br> $\left(R^{2}\right.$ <br> $\mathrm{HNO}_{3}($ aq $)$ <br> Ignore references to heat | (1) | (3) |



| Question <br> Number | Acceptable Answers |  | Mark |
| :--- | :--- | :---: | :---: |
| $\mathbf{1 9 ( e ) ( i ) ~}$ | M1: Aspirin concentration $=180 \times 0.0161=2.898 \mathrm{~g} \mathrm{dm}^{-3}$ | (1) | (3) |
|  | (Need 600 mg or 0.600 g so $)$ |  |  |
|  | M2: Volume required $=(0.600 \div 2.898) \times 1000=207.039 \mathrm{~cm}^{3}$ | (1) |  |
|  | OR | M1:Number of moles $=(0.300 \times 2) \div 180=0.003333$ | (1) |
|  | M2: Volume required $=0.00333 \div 0.0161=0.0207039\left(\mathrm{dm}^{3}\right)$ | (1) |  |
|  | M3: $207 \mathrm{~cm}^{3} / 0.207 \mathrm{dm}^{3}$ | (1) |  |
|  | Answer must be to 3 SF and include units |  |  |
|  | TE throughout |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 9 ( e ) ( i i ) ~}$ | (1) <br> Allow Kekulé structure for benzene ring <br> Allow structure without charges <br> Ignore the orientation and bond length of the substituents of <br> the ring but they must be adjacent |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9 ( e ) ( \text { (iii) }}$ | Ionic salts are (more) soluble <br> (in water than the unionised aspirin) | $\mathbf{( 1 )}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| *19(f) | Any two reasons from: | (1) | (2) |
|  | Avoid/reduce side effects <br> Allow references to 'not harmful' <br> Reduce the quantity of substance needed for the required <br> dose <br> (1) | (1) | Just <br> 'saves <br> money' |
| Reduce waste (of the undesired enantiomer) <br> Ignore reference to one isomer being inactive or one being |  |  |  |

(Total for Section C=20 marks)
(Total for paper $=\mathbf{9 0}$ marks)

