## Pearson Edexcel

## Mark Scheme (Results)

## Summer 2019

Pearson International Advanced Level In Chemistry (WCH05) Paper 01General Principles of Chemistry II - Transition Metals and Organic Nitrogen Chemistry

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## General Marking Guidance

- $\quad$ 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.

## 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)

| $\begin{array}{l}\text { Question } \\ \text { Number }\end{array}$ | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1}$ | The only correct answer is C | (1) |
|  | $\boldsymbol{A}$ is not correct because Al has an oxidation number of +3 |  |
| $\boldsymbol{B}$ is not correct because Cr has an oxidation number of +3 |  |  |
| $\boldsymbol{D}$ is not correct because $V$ has an oxidation number of +4 |  |  |$]$


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{2}$ | The only correct answer is C <br> A is not correct because this is the number of moles of iodate(V) <br> ions needed to react with 1 mol of hydrogensulfite ions | (1) |
| $\boldsymbol{B}$ is not correct because the reaction is not 1:1 |  |  |
| $\boldsymbol{D}$ is not correct because this is the number of moles of |  |  |
| hydrogensulfite ions needed to react with 2 mol of iodate(V) ions |  |  |
| (giving an equation with integer coefficients). |  |  |$\quad$.


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{3 ( a )}$ | The only correct answer is A <br> $\mathbf{B}$ is not correct because manganese would react to form a cell <br> system with $\mathrm{Mn}^{2+}$ <br> C is not correct because $\mathrm{Mn}^{2+}$ is the reducing agent in the system <br> $\mathbf{D}$ is not correct because $\mathrm{Mn}^{2+}$ is the reducing agent in the system <br> and manganese would react to form a cell system with $\mathrm{Mn}^{2+}$ | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| 3(b) | The only correct answer is D <br> A is not correct because all the substances in an electrode system <br> must be present in a half-cell <br> B is not correct because all the substances in an electrode system <br> must be present in a half-cell <br> C is not correct because all the substances in an electrode system <br> must be present in a half-cell | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{4}$ | The only correct answer is D <br> $\boldsymbol{A}$ is not correct because $E_{\text {cell }}$ ® <br> nor $K$ | (1) proportional to neither $\Delta S_{\text {system }}$ |
| $\boldsymbol{B}$ is not correct because $E_{\text {cell }}{ }^{\circledR}$ is proportional to InK but not to |  |  |
| $\Delta S_{\text {system }}$ |  |  |
| $\boldsymbol{C}$ is not correct because $E_{\text {cell }}$ is proportional to $\Delta S_{\text {total }}$ but not to $K$ |  |  |$\quad$.


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{5}$ | The only correct answer is A | (1) |
|  | B is not correct because oxidation always occurs at the anode <br> C is not correct because ethanol is oxidised in this cell <br> $\boldsymbol{D}$ is not correct because ethanol is oxidised in this cell |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{6}$ | The only correct answer is A <br> B is not correct because manganese in the +4 oxidation state will <br> have partially filled d orbitals <br> C is not correct because iron in the +4 oxidation state will have <br> partially filled d orbitals <br> $\boldsymbol{D}$ is not correct because copper in the +4 oxidation state will <br> have partially filled d orbitals | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{7}$ | The only correct answer is B <br> $\boldsymbol{A}$ is not correct because oxygen is not a reducing agent and <br> sulfur dioxide is not an oxidising agent <br> C is not correct because the highest stable oxidation number of <br> vanadium is +5 | (1) |
| D is not correct because oxygen is not a reducing agent and <br> sulfur dioxide is not an oxidising agent and because the highest <br> stable oxidation number of vanadium is +5 |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{8}$ | The only correct answer is B <br> A is not correct because this shows benzene at a higher energy <br> level than cyclohexa-1,3,5-triene and at a lower energy level than <br> cyclohexane. <br> C is not correct because this shows cyclohexane at the highest <br> energy level instead of the lowest. | (1) |
| D is not correct because this shows benzene at a higher energy <br> level than cyclohexa-1,3,5-triene. |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{9}$ | The only correct answer is C <br> A is not correct because this compound does not form an ionic <br> compound with sodium hydroxide or react with ethanol | (1) |
| $\boldsymbol{B}$ is not correct because this compound does not decolorise |  |  |
| bromine water |  |  |
| $\boldsymbol{D}$ is not correct because this compound does not react with |  |  |
| ethanol |  |  |$\quad$|  |
| :--- |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 0}$ | The only correct answer is C <br> A is not correct because ethanal cannot form the required <br> electrophile <br> B is not correct because ethanoic acid cannot form the required <br> electrophile <br> D is not correct because propanone cannot form the required <br> electrophile | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1}$ | The only correct answer is B <br> $\boldsymbol{A}$ is not correct because alkenes form poly(alkenes) from single <br> monomers <br> C is not correct because $\boldsymbol{F}$ reacts with $\boldsymbol{G}$ to form a polyamide <br> $\boldsymbol{D}$ is not correct because $\boldsymbol{F}$ reacts with $\boldsymbol{H}$ to form a polyester | (1) |


| Question Number | Answer | Mark |
| :---: | :---: | :---: |
| 12(a) | The only correct answer is C <br> $\boldsymbol{A}$ is not correct because the nitrogen atom furthest left in the structure is in an amine group <br> B is not correct because the other two nitrogen atoms are part of amide groups <br> D is not correct because the hydroxyl group attached directly to the benzene ring is phenolic | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| 12(b) | The only correct answer is A <br> B is not correct because the phenolic OH group is very much less <br> basic than the amine group | (1) |
| C is not correct because the phenolic OH group will not lose a <br> proton in acidic conditions | D is not correct because the carboxylic acid group is very much <br> less basic than the amine group |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 3 ( a )}$ | The only correct answer is C <br> A is not correct because covalent bonds are not broken when <br> amino acids melt <br> B is not correct because hydrogen bonds cannot form <br> between zwitterions <br> D is not correct because London forces are much weaker <br> than ionic bonds | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 3 ( b )}$ | The only correct answer is D <br> A is not correct because only alanine has a chiral carbon and <br> exists as optical isomers | (1) |
| B is not correct because alanine has a chiral carbon and exists as <br> optical isomers <br> C is not correct because glycine does not have a chiral carbon <br> and does not exist as optical isomers |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| 13(c) | The only correct answer is D <br> $\boldsymbol{A}$ is not correct because neither structure has a peptide (CONH) <br> link | (1) |
| B is not correct because the right-hand structure does not have a <br> peptide (CONH) link <br> C is not correct because the left-hand structure does not have a <br> peptide (CONH) link |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 4}$ | The only correct answer is B <br> A is not correct because sulfuric acid is not involved in the <br> hydrolysis so does not act as a catalyst <br> C is not correct because the sodium hydroxide effects the <br> hydrolysis <br> $\boldsymbol{D}$ is not correct because this is not the function of the sulfuric <br> acid | (1) |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 5}$ | The only correct answer is A <br> $\boldsymbol{B}$ is not correct because this is the number of protons in each <br> environment, not the splitting pattern | (1) |
|  | C is not correct because this considers the splitting of the C2 <br> protons to be due only to the C1 protons |  |
| D is not correct because this considers the splitting of the C2 <br> protons to be due only to the C1 protons and additionally shows <br> the hydroxy proton peak being split |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 6}$ | The only correct answer is D <br> A is not correct because it will not form a white solid with dilute <br> sulfuric acid. <br> $\boldsymbol{B}$ is not correct because it will not form a pale yellow precipitate <br> when warmed with iodine and sodium hydroxide <br> C is not correct because it will not form a pale yellow precipitate <br> when warmed with iodine and sodium hydroxide or a white solid <br> with sulfuric acid. | (1) |

Section B

| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(a)(i) | $E_{\text {cell }}^{\ominus}=+0.40-(-0.74)=(+) 1.14(\mathrm{~V})$ <br> No TE on incorrect half equations <br> $4 \mathrm{Cr}+3 \mathrm{O}_{2}+6 \mathrm{H}_{2} \mathrm{O} \rightleftharpoons 4 \mathrm{Cr}^{3+}+12 \mathrm{OH}^{-}$ <br> OR <br> $4 \mathrm{Cr}+3 \mathrm{O}_{2}+6 \mathrm{H}_{2} \mathrm{O} \rightleftharpoons 4 \mathrm{Cr}(\mathrm{OH})_{3}$ <br> ALLOW <br> Multiples <br> $\rightarrow$ in place of $\rightleftharpoons$ <br> Species and equation in correct direction <br> Balancing <br> ALLOW <br> TE on incorrect half equations from the table <br> eg $4 \mathrm{Cr}^{2+}+\mathrm{O}_{2}+2 \mathrm{H}_{2} \mathrm{O} \rightleftharpoons 4 \mathrm{Cr}^{3+}+4 \mathrm{OH}^{-}$ <br> IGNORE state symbols even if incorrect <br> COMMENT <br> Correct $E^{\ominus}$ cell value may be credited in (a)(ii) | $\begin{equation*} -1.14(\mathrm{~V}) \tag{1} \end{equation*}$ <br> uncancelled electrons | (3) |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(a)(ii) | TE on any positive value in 17(a)(i) <br> The positive $E^{\theta}$ cell value indicates that the corrosion of chromium is (thermodynamically) feasible <br> ALLOW <br> spontaneous for feasible <br> $E^{\circ}$ cell value indicates chromium and oxygen <br> should react / chromium corrodes <br> TE for M1 only on any negative value in 17(a)(i) <br> So the corrosion is kinetically unfavourable / has a high activation energy / slow OR chromium forms a stable / unreactive oxide coating (that protects the metal from corrosion) <br> ALLOW hydroxide for oxide | Just 'needs high energy' <br> Any reference to sacrificial protection | (2) |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(b)(i) | Correct equation with $E^{\ominus}{ }_{\text {cell }}$ value scores (2) <br> Route 1 <br> Zinc / Zn <br> IGNORE <br> Acids / H ${ }^{+}$ <br> Ionic half-equation $\begin{align*} & \left(\mathrm{Zn}(\mathrm{~s})+2 \mathrm{Cr}^{3+}(\mathrm{aq}) \rightleftharpoons 2 \mathrm{Cr}^{2+}(\mathrm{aq})+\mathrm{Zn}^{2+}(\mathrm{aq})\right) \\ & E_{\text {cell }}^{9}[=-0.41-(-0.76)]=(+) 0.35(\mathrm{~V}) \tag{1} \end{align*}$ <br> Route 2 <br> Chromium / Cr $\begin{equation*} E_{\text {cell }}^{\theta}=-0.41-(-0.74)=(+) 0.33(\mathrm{~V}) \tag{1} \end{equation*}$ | Other additional reagents | (2) |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :---: | :--- | :--- | :---: |
| 17(b)(ii) | Standalone marks <br> Chromium(II) <br> (is readily oxidised back to chromium(III)) <br> (1) | Oxidation to <br> any other <br> oxidation state | (2) |
|  | oxidised by oxygen (in the air) <br> ALLOW <br> by air <br> IGNORE just 'reacts with oxygen (in the air)' | (1) |  |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :---: | :--- | :---: | :---: |
| $\mathbf{1 7 ( b ) ( i i i ) ~}$ | From green to blue |  | (1) |


|  | ALLOW <br> violet to blue <br> IGNORE <br> modifiers (eg pale) |  |  |
| :--- | :--- | :--- | :--- |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(b)(iv) | M1 (energy gap) <br> There is a different energy gap between the <br> (3)d orbitals / in the (3)d subshell <br> ALLOW <br> different d-d splitting <br> (3)d orbitals split differently <br> M2 (explanation of energy gap) <br> Because of the (different) charge / oxidation state / radius / size / charge density / electronic structure / numbers of (d)electrons of the ions <br> M3 (effect of energy gap) <br> So different frequencies / wavelengths of (visible) light / radiation / energy are absorbed / reflected / transmitted OR <br> Photons of different energy are absorbed / reflected / transmitted <br> IGNORE <br> Colour for (visible) light / radiation / energy General explanations of the colour of transition metal complexes even if incorrect | (3)d orbital <br> Just different ions/ligands <br> emitted <br> emitted | (3) |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 17(c)(i) | In 17(c)(i) and (ii) <br> two correct formulae scores (1) in (c)(i) |  | (1) |


|  | Ignore omission of square brackets <br> $\left[\mathrm{CrCl}_{4}\right]^{-}$and tetrahedral |  |  |
| :--- | :--- | :--- | :--- |


| $\begin{array}{l}\text { Question } \\ \text { Number }\end{array}$ | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( c ) ( \text { (ii) }}$ | $\begin{array}{l}{\left[\mathrm{Cr}\left(\mathrm{NH}_{3}\right)_{6}\right]^{3+} /\left[\mathrm{Cr}\left(\mathrm{NH}_{3}\right)_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}\right]^{3+}} \\ \text { and } \\ \text { octahedral }\end{array}$ | $\left[\mathrm{Cr}\left(\mathrm{NH}_{3}\right)_{4}\right]^{3+}$ | (1) |
| square planar |  |  |  |$]$


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( d ) ( i )}$ | $2 \mathrm{CrO}_{4}{ }^{2-}+2 \mathrm{H}^{+} \rightarrow \mathrm{Cr}_{2} \mathrm{O}_{7}^{2-}+\mathrm{H}_{2} \mathrm{O}$ |  | (2) |
|  | OR <br> $\rightleftharpoons$ in place of $\rightarrow$ <br> ALLOW <br> Multiples <br> Formulae of both chromium species (in <br> any equation) <br> Correct balanced equation | (1) | Additional <br> chromium species |
| IGNORE <br> State symbols even if incorrect |  |  |  |
| No TE on incorrect chromium species |  |  |  |$\quad$|  |
| :--- |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |



| Question <br> Number | Acceptable Answer | Reject | Mark |
| :---: | :--- | :--- | :---: |
| 17(d)(iii) | Colourless <br> and <br> $\mathrm{Cr}^{6+}$ has no (3)d electrons <br> ALLOW <br> Colourless <br> and <br> (3)d subshell /orbitals empty /(3)d ${ }^{0}$ | White <br> (3)d orbital is <br> empty <br> no (3)d orbitals / <br> no (3)d subshell | (1) |
| IGNORE <br> No d-d transitions <br> No d-d splitting |  |  |  |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( e ) ( i ) ~}$ | Marks are standalone | (1) |  |
| Starch (solution) <br> Blue-black / blue / black <br> and <br> to colourless / green <br> IGNORE <br> clear | (1) | yellow |  |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(e)(ii) | M1 $\begin{align*} \mathrm{mol} \mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-} & =10 \times 0.0495 \times 10^{-3} \\ & =4.95 \times 10^{-4} / 0.000495 \tag{1} \end{align*}$ <br> M2 $\begin{align*} & \mathrm{mol}_{2}=3 \times \mathrm{mol} \mathrm{Cr}_{2} \mathrm{O}_{7}^{2-} \\ &= 3 \times 4.95 \times 10^{-4}=1.485 \times 10^{-3} \\ & \quad / 0.001485 \tag{1} \end{align*}$ <br> M3 $\begin{align*} \mathrm{mol} \mathrm{~S}_{2} \mathrm{O}_{3}{ }^{2-} & =2 \times \mathrm{mol} \mathrm{I}_{2} \\ & =2 \times 1.485 \times 10^{-3} \\ & =2.97 \times 10^{-3} / 0.00297 \tag{1} \end{align*}$ <br> M4 <br> concentration of $\mathrm{S}_{2} \mathrm{O}_{3}{ }^{2-}$ $\begin{align*} & =1000 \times 2.97 \times 10^{-3} / 19.50 \\ & =0.1523 \mathrm{~mol} \mathrm{dm}^{-3} \tag{1} \end{align*}$ <br> Ignore SF except 1 SF <br> ALLOW <br> TE at each stage <br> Do not penalise correct intermediate rounding to 2 SF <br> Correct answer with some / no working scores (4) | incorrect or missing units | (4) |

(Total for Question 17 = 27 marks)
Some Common Incorrect Answers

| Error | Final Answer | Mark $/ 4$ |
| :---: | :---: | :---: |
| $\mathrm{~mol} \mathrm{~S}_{2} \mathrm{O}_{3}{ }^{2-}=\mathrm{mol} \mathrm{I}_{2} / 2$ | $0.03808 / 3.808 \times 10^{-2}$ | 3 |
| $\mathrm{~mol} \mathrm{~S}_{2} \mathrm{O}_{3}^{2-}=\mathrm{mol} \mathrm{I}_{2}$ | $0.07615 / 7.615 \times 10^{-2}$ | 3 |
| $\mathrm{~mol} \mathrm{~S}_{2} \mathrm{O}_{3}{ }^{2-}=\mathrm{mol} \mathrm{I}_{2} \times 2 / 3$ | $0.05077 / 5.077 \times 10^{-2}$ | 3 |
| $\mathrm{~mol} \mathrm{~S}_{2} \mathrm{O}_{3}{ }^{2-}=\mathrm{mol} \mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-} \times 2$ | $0.05077 / 5.077 \times 10^{-2}$ | 3 |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(a) | Step 1 <br> Add bromine <br> ALLOW <br> Chlorine <br> Intermediate is 1,2-dibromoethane <br> OR <br> $\mathrm{CH}_{2} \mathrm{BrCH}_{2} \mathrm{Br} /$ displayed / skeletal formula <br> OR <br> 1,2-dichloroethane <br> OR <br> $\mathrm{CH}_{2} \mathrm{ClCH}_{2} \mathrm{Cl} /$ displayed / skeletal formula <br> IGNORE $\mathrm{C}_{2} \mathrm{H}_{4} \mathrm{Br}_{2} / \mathrm{C}_{2} \mathrm{H}_{4} \mathrm{Cl}_{2}$ <br> ALLOW feasible alternatives for Step1 eg acidic/alkaline $\mathrm{KMnO}_{4}$ <br> and <br> forms $\mathrm{CH}_{2} \mathrm{OHCH}_{2} \mathrm{OH}$ and using $\mathrm{P} / \mathrm{I}_{2}$ <br> 1,2-diiodoethane <br> Step 2 <br> TE on any halogenoalkane from Step 1 <br> Ammonia ((gas)) <br> ALLOW <br> Conc $\mathrm{NH}_{3}(\mathrm{aq})$ <br> Dissolved in ethanol / alcohol (and heat) <br> OR <br> heat in a sealed tube / heat under pressure <br> (add alkali / NaOH / KOH) <br> If $\mathrm{NH}_{3}(\mathrm{aq})$ used M4 may be scored | $\mathrm{Br}_{2}(\mathrm{aq}) / \mathrm{Br}$ UV light Additional reagents | (4) |



| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(c)(i) |  <br> OR displayed / structural formulae <br> ALLOW <br> Two or three 'en' molecules attached without showing octahedral shape provided nitrogens correctly bonded to copper <br> IGNORE <br> Omission of water <br> Omission of square brackets <br> Omission of charge / incorrect charge |  | (1) |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 18(c)(ii) | The number of particles / molecules / <br> moles increases | Ammonia is <br> released as a <br> gas | (2) |
|  | Do not penalise the use of specific <br> numbers showing an increase, even if <br> incorrect | IGNORE <br> Just 'more product' | system is positive / increases (with no / <br> negligible change in $\Delta S_{\text {surroundings so the }}$ <br> reaction is favoured) <br> ALLOW <br> $\Delta S_{\text {total }} /$ entropy increases <br> IGNORE <br> Disorder increases <br> References to stability constants |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(d)(i) | Ethanoyl chloride / $\mathrm{CH}_{3} \mathrm{COCl}$ <br> OR <br> Ethanoic anhydride / ( $\left.\mathrm{CH}_{3} \mathrm{CO}\right)_{2} \mathrm{O}$ <br> OR <br> Displayed / skeletal formula <br> ALLOW <br> Acetic anhydride <br> Acetic acid anhydride | $\mathrm{CH}_{3} \mathrm{CICO}$ <br> Ethanoic acid / $\mathrm{CH}_{3} \mathrm{COOH}$ <br> Additional reagents | (1) |


| Question <br> Number | Acceptable Answer | Reject | Mark |  |
| :--- | :--- | :--- | :--- | :--- |
| 18(d)(ii) | The (four) protons on the central carbon <br> chain are equivalent <br> The (twelve) protons on the four methyl <br> groups are equivalent <br> Two peaks (because there are two proton <br> environments) <br> and <br> The relative peak areas are <br> $4: 12 / 1: 3$ <br> Clearly labelled diagram scores M1 and M2 <br> eg |  | (3) |  |



| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(b)(i) | $m / e=77\left(\right.$ is $\mathrm{C}_{6} \mathrm{H}_{5}{ }^{+}$) so phenyl group/ $\mathrm{C}_{6} \mathrm{H}_{5}$ present ALLOW <br> $\mathbf{M}$ is an arene / aromatic / contains benzene (ring) <br> Effervescence / reaction with $\mathrm{NaHCO}_{3}$ is typical of an acid so carboxylic acid/COOH present <br> Decolourisation of reaction with $\mathrm{KMnO}_{4} / \mathrm{H}^{+}$indicates a carbon-carbon double bond / C=C / alkene <br> (1) <br> Three groups correctly identified with no explanation scores (1) | phenol $\begin{equation*} \mathrm{C}_{6} \mathrm{H}_{6} \tag{1} \end{equation*}$ | (3) |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :---: | :---: | :---: | :---: |
| 19(b)(ii) | (3) |  |  |

(Total for Question 19 = 10 marks)
Total for Section B=51 marks

## Section C

| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(a) | In (a) and (b) award marks for correct intermediates even if prepared incorrectly and correct reagents used with correct functional groups even if the molecule is wrong. <br> Penalise omission of conditions once only in (a) <br> Step 1 <br> KCN / NaCN in ethanol / alcohol <br> (reflux/heat) <br> To form butanenitrile / $\mathrm{C}_{3} \mathrm{H}_{7} \mathrm{CN}$ <br> Step 2 <br> Reflux with sulfuric acid / $\mathrm{H}_{2} \mathrm{SO}_{4}$ <br> OR <br> Reflux with NaOH and followed by addition of sulfuric acid / $\mathrm{H}_{2} \mathrm{SO}_{4}$ <br> ALLOW <br> Any strong acid / HCl for $\mathrm{H}_{2} \mathrm{SO}_{4}$ <br> Heat / boil for reflux <br> IGNORE <br> Concentration of acid <br> To form butanoic acid / $\mathrm{C}_{3} \mathrm{H}_{7} \mathrm{COOH}$ <br> Step 3 <br> Lithium tetrahydridoaluminate(III) / lithium aluminium hydride / $\mathrm{LiAlH}_{4}$ in (dry) ether / ethoxyethane (under reflux) (followed by dilute strong acid) <br> ALLOW <br> Lithal in (dry) ether <br> IGNORE <br> Name or formula of butan-1-ol, even if incorrect | HCN | (5) |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(b) | Step 1 <br> $\mathrm{KOH} / \mathrm{NaOH}$ in ethanol / alcohol <br> / alcoholic (reflux) <br> To form but-1-ene $/ \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CHCH}_{2}$ / <br> skeletal / displayed formula <br> IGNORE <br> butene <br> If name and formula are given, both <br> must be correct <br> Step 2 (depends on an alkene as the organic reactant) <br> Add hydrogen bromide / HBr <br> / hydrobromic acid / $\mathrm{NaBr} \& \mathrm{H}_{2} \mathrm{SO}_{4}$ | Aqueous ethanol <br> but-2-ene | (3) |


| Question Number | Acceptable Answer | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(c)(i) | These marks are standalone <br> Sodium nitrite /sodium nitrate(III) / $\mathrm{NaNO}_{2}$ and hydrochloric acid / $\mathrm{HCl}((\mathrm{aq}))$ <br> OR <br> Potassium nitrite / $\mathrm{KNO}_{2}$ and $\mathrm{HCl}((\mathrm{aq}))$ <br> OR <br> Sulfuric acid (for hydrochloric acid) <br> ALLOW <br> Nitrous acid / $\mathrm{HNO}_{2} / \mathrm{HONO}$ <br> IGNORE <br> Concentration of acid hydrochloric acid / sulfuric acid with nitrous acid <br> $0-10^{\circ} \mathrm{C} /$ ice(-water) bath <br> ALLOW <br> $\angle 10^{\circ} \mathrm{C} / \angle 5^{\circ} \mathrm{C}$ | sodium nitrate | (2) |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 20(c)(ii) | Left-hand curly arrow <br> ALLOW <br> Left-hand curly arrow going to the positive charge <br> Right-hand curly arrow and lone pair <br> ALLOW any type of connecting arrow or none <br> Penalise half-arrows once only <br> IGNORE any additional lone pairs <br> COMMENT <br> Penalise additional curly arrows once only | (1) |  |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( c ) ( \text { iii) }}$ |  | (2) |  |
|  | OR <br> $1,2 / 1,3$ structures <br> Electrophilic substitution <br> ALLOW <br> Electrophilic coupling | (1) |  |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 20(d)(i) | (nucleophilic substitution of groups attached directly to <br> a benzene ring is normally very difficult because) <br> High electron density of the ring repels nucleophiles <br> OR <br> benzene ring sterically hinders the approach of <br> nucleophiles | (2) |  |
|  | ALLOW <br> pi / delocalised electrons repel nucleophiles <br> OR <br> Nucleophilic attack is difficult because of the high <br> electron density of the ring <br> (benzenediazonium ions readily undergo nucleophilic <br> substitution because) <br> Nitrogen is a very good leaving group <br> ALLOW <br> Nitrogen is a gas so the entropy change (of the system) <br> is (very) positive <br> OR <br> nitrogen gas is very stable <br> OR <br> N2 <br> electron density on the benzene ring <br> IGNORE <br> References just to the positive charge on the nitrogen | (1) |  |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :---: | :---: |
| 20(d)(ii) | The (main) oxidation states of copper are readily <br> interchanged <br> ALLOW <br> (Copper(I) ions) are (easily) changed into copper(II) ions <br> and copper <br> OR <br> (Copper(I) ions) are (easily) oxidised and reduced <br> OR <br> Copper has variable oxidation states / variable valency | (1) |  |
| IGNORE <br> Copper is a transition element / metal <br> References to partially filled 3(d) orbitals / subshell <br> References to surface catalysis <br> COMMENT <br> ALLOW <br> Copper(I) / Cu/ It / they have variable oxidation states |  |  |  |


| Question <br> Number | Acceptable Answer | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 20(d)(iii) | Mark independently | (2) |  |

(Total for Question 20 = 19 marks)
(Total for Section C = 19 marks)
Total for PAPER = 90 Marks

