# Pearson Edexcel 

## Mark Scheme (Results)

## Summer 2023

Pearson Edexcel International Advanced
Subsidiary Level In Chemistry (WCH15)
Paper 01
Unit 5: Transition Metals and Organic Nitrogen
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.

## Section A

| Question Number | Answer | Mark |
| :---: | :---: | :---: |
| 1 | The only correct answer is $\mathbf{D}\left(\mathrm{Pt}(\mathrm{s})\left\|\mathrm{V}^{2+}(\mathrm{aq}), \mathrm{V}^{3+}(\mathrm{aq})\right\|\left\|\mathrm{Cu}^{2+}(\mathrm{aq})\right\| \mathrm{Cu}(\mathrm{s})\right)$ <br> $\boldsymbol{A}$ is not correct because the $V^{3+}(a q) / V^{2+}(a q)$ half-cell should have a platinum electrode and should show oxidation and the $\mathrm{Cu}^{2+}(a q) / \mathrm{Cu}(\mathrm{s})$ half-cell should show reduction <br> B is not correct because the $V^{3+}(a q) / V^{2+}($ aq $)$ half-cell should have a platinum electrode <br> $\boldsymbol{C}$ is not correct because the $\mathrm{V}^{3+}(a q) / V^{2+}(a q)$ half-cell should show oxidation and the $\mathrm{Cu}^{2+}(a q) / \mathrm{Cu}(\mathrm{s})$ half-cell should show reduction | (1) <br> Computer |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{2}$ | The only correct answer is $\mathbf{D ~}\left(\mathrm{Mg}+2 \mathrm{Ce}^{4+} \rightarrow \mathrm{Mg}^{2+}+2 \mathrm{Ce}^{3+}\right)$ | $\mathbf{( 1 )}$ |
|  | $\boldsymbol{A}$ is not correct because $C$ e is a weaker reducing agent than Mg |  |
| $\boldsymbol{B}$ is not correct because $C e^{3+}$ is a weaker reducing agent than Ce |  |  |
| $\boldsymbol{C}$ is not correct because $\mathrm{Mn}^{2+}$ is a weaker reducing agent than Mn | Computer |  |


| Question <br> Number | Answer | Mark |
| :---: | :---: | :---: |
| 3 | The only correct answer is $\mathbf{D}\left(\Delta S^{\ominus}{ }_{\text {total }}\right)$ <br> $\boldsymbol{A}$ is not correct because $E^{\theta}$ cell is directly proportional to $\ln K_{c}$ <br> $\boldsymbol{B}$ is not correct because $E^{\theta}{ }_{\text {cell }}$ is directly proportional to $\Delta S^{\theta}{ }_{\text {total }}$ and not $\Delta H^{\theta}$ <br> $\boldsymbol{C}$ is not correct because $E^{\theta}$ cell is directly proportional to $\Delta S^{\theta}{ }_{\text {total }}$ and not $\Delta S^{\theta}{ }_{\text {system }}$ | (1) <br> Computer |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{4}$ | The only correct answer is C (the reactants are thermodynamically unstable with respect to the products) | (1) |
|  | $\boldsymbol{A}$ is not correct because the reaction is thermodynamically feasible so will occur under certain conditions |  |
| $\boldsymbol{B}$ is not correct because the $E^{\theta}$ cell value is a thermodynamic and not a kinetic property | Computer |  |
| $\boldsymbol{D}$ is not correct because the reaction may be kinetically inert and the conditions may be non-standard |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{5}$ | The only correct answer is A $\left(\mathrm{H}_{2}+2 \mathrm{OH}^{-} \rightarrow 2 \mathrm{H}_{2} \mathrm{O}+2 \mathrm{e}^{-}\right)$ | $\mathbf{( 1 )}$ |
|  | $\boldsymbol{B}$ is not correct because hydrogen is consumed and not produced in a hydrogen-oxygen fuel cell |  |
| $\boldsymbol{C}$ is not correct because oxygen is reduced at the positive electrode in a hydrogen-oxygen fuel cell |  |  |
| $\boldsymbol{D}$ is not correct because oxygen is consumed and not produced in a hydrogen-oxygen fuel cell |  |  |


| Question Number | Answer | Mark |
| :---: | :---: | :---: |
| 6 | The only correct answer is $\mathbf{D}\left(\mathrm{Cu} \quad[\mathrm{Ar}] 3 \mathrm{~d}^{10} 4 \mathrm{~s}^{1}\right)$ | (1) |
|  | $\boldsymbol{A}$ is not correct because the 4 s electrons are removed before the 3 d electrons | Computer |
|  | $\boldsymbol{B}$ is not correct because the electronic configuration of chromium is [Ar] $3 d^{5} 4 s^{1}$ |  |
|  | $\boldsymbol{C}$ is not correct because the 4s electrons are removed before the 3d electrons |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{7}$ | The only correct answer is C (six) | $\mathbf{( 1 )}$ |
|  | $\boldsymbol{A}$ is not correct because it only takes into account water ligands |  |
|  | $\boldsymbol{B}$ is not correct because it only takes into account ethanoate ions |  |
| $\boldsymbol{D}$ is not correct because the coordination numbers of the two chromiums have been added together |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{8}$ | The only correct answer is A $\left(\left[\mathrm{Fe}(\mathrm{CN})_{6}\right]^{4-}\right)$ | (1) |
|  | $\boldsymbol{B}$ is not correct because $\mathrm{C}_{2} \mathrm{O}_{4}{ }^{2-}$ is a bidentate ligand |  |
| $\boldsymbol{C}$ is not correct because $\mathrm{EDTA}^{4-}$ is a hexadentate ligand |  |  |
| $\boldsymbol{D}$ is not correct because $\mathrm{NH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{NH}_{2}$ is a bidentate ligand |  |  | Computer.


| $\begin{array}{l}\text { Question } \\ \text { Number }\end{array}$ | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{9}$ | The only correct answer is $\mathbf{A}\left(\left[\mathrm{CuCl}_{4}\right]^{2-}\right)$ | $\mathbf{( 1 )}$ |
|  | $\boldsymbol{B}$ is not correct because this complex is octahedral with a bond angle of $90^{\circ}$ |  |
| $\boldsymbol{C}$ is not correct because this complex is linear with a bond angle of $180^{\circ}$ |  |  |
| $\boldsymbol{D}$ is not correct because this complex is square planar with a bond angle of $90^{\circ}$ |  |  |$)$ Computer |  |
| :--- |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: | :---: |
| $\mathbf{1 0}$ | The only correct answer is $\mathbf{B}\left(\left[\mathrm{Co}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}\right)$ | (1) |
|  | $\boldsymbol{A}$ is not correct because $\mathrm{VO}^{2+}$ is blue | Computer |
|  | $\boldsymbol{C}$ is not correct because $\left[\mathrm{Cr}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}$ is blue |  |
| D is not correct because $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}$ is blue |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1}$ | The only correct answer is $\mathbf{D}\left(2 \mathrm{Cr}(\mathrm{OH})_{3}+3 \mathrm{H}_{2} \mathrm{O}_{2}+4 \mathrm{KOH} \rightarrow 2 \mathrm{~K}_{2} \mathrm{CrO}_{4}+8 \mathrm{H}_{2} \mathrm{O}\right)$ | $\mathbf{( 1 )}$ |
|  | $\boldsymbol{A}$ is not correct because $\mathrm{FeCl}_{2}$ forms a green solution | Computer |
|  | $\boldsymbol{B}$ is not correct because this is not a redox reaction |  |
| $\boldsymbol{C}$ is not correct because this is not a redox reaction |  |  |


| Question Number | Answer | Mark |
| :---: | :---: | :---: |
| 12 | The only correct answer is $\mathbf{C}\left(\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}+4 \mathrm{NH}_{3} \rightarrow\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}\right]^{2+}+4 \mathrm{H}_{2} \mathrm{O}\right)$ <br> $\boldsymbol{A}$ is not correct because this is the ionic equation describing the deprotonation when ammonia is not in excess <br> $\boldsymbol{B}$ is not correct because four water ligands are exchanged by ammine ligands when ammonia is in excess <br> $\boldsymbol{D}$ is not correct because four water ligands are exchanged by ammine ligands when ammonia is in excess | (1) <br> Computer |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 3}$ | The only correct answer is A $\left(\left[\mathrm{Zn}(\mathrm{OH})_{4}\right]^{2-}+2 \mathrm{H}_{3} \mathrm{O}^{+} \rightarrow\left[\mathrm{Zn}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}(\mathrm{OH})_{2}\right]\right)$ | (1) |
|  | $\boldsymbol{B}$ is not correct because $\left[\mathrm{Zn}(\mathrm{OH})_{4}\right]^{2-}$ is a soluble complex ion |  |
| $\boldsymbol{C}$ is not correct because $\left[\mathrm{Cr}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{3+}$ is a soluble complex ion |  |  |
| D is not correct because $\left[\mathrm{Cr}(\mathrm{OH})_{6}\right]^{3-}$ is a soluble complex ion | Computer |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 4}$ | The only correct answer is $\mathbf{C}(+5 \rightarrow+4 \rightarrow+5)$ | (1) |
|  | $\boldsymbol{A}$ is not correct because the oxidation state in $V_{2} O_{5}$ is +5 not +2 | Computer |
|  | $\boldsymbol{B}$ is not correct because the oxidation state in $V_{2} O_{5}$ is +5 not +2 |  |
| $\boldsymbol{D}$ is not correct because the vanadium cannot be oxidised from +5 to +6 |  |  |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 5}$ | The only correct answer is C (14) | $\mathbf{( 1 )}$ |
|  | $\boldsymbol{A}$ is not correct because each carbon contributes one electron from a p-orbital |  |
|  | $\boldsymbol{B}$ is not correct because each carbon contributes one electron from a p-orbital |  |
| $\boldsymbol{D}$ is not correct because each carbon contributes one electron from a p-orbital |  |  |


| Question Number | Answer | Mark |
| :---: | :---: | :---: |
| 16 | The only correct answer is $\mathbf{C}$ (a lone pair of electrons on oxygen in phenol is delocalised into the ring) <br> $\boldsymbol{A}$ is not correct because the polarity of the $O-H$ bond does not increase the electron density of the benzene ring <br> $\boldsymbol{B}$ is not correct because the electronegativity of the oxygen atom does not increase the electron density of the benzene ring <br> $\boldsymbol{D}$ is not correct because there is a greater electron density in the ring in phenol than in benzene | (1) <br> Computer |


| Question <br> Number | Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 7}$ | The only correct answer is $\mathbf{B}\left(\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{NH}_{2} \xrightarrow[\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{Br}]{\longrightarrow}\right)$ | $\mathbf{( 1 )}$ |
|  | $\boldsymbol{A}$ is not correct because the reduction of a nitrile forms a primary amine |  |
| $\boldsymbol{C}$ is not correct because the products would be a tertiary amine and a quaternary ammonium salt |  |  |
| $\boldsymbol{D}$ is not correct because the alkaline hydrolysis of this amide forms a primary amine |  |  |


| Question |  | Mark |
| :--- | :--- | :---: | :---: |
| $\mathbf{N u m b e r ~}$ | The only correct answer is $\mathbf{D}$ ( |  |
| $\mathbf{A}$ is not correct because this is not an azo dye | (1) | Computer |
|  | $\boldsymbol{B}$ is not correct because this is not an azo dye |  |
| C is not correct because this azo dye could only form if the reagents were not in excess |  |  |


| Question Number | Answer | Mark |
| :---: | :---: | :---: |
| 19 | The only correct answer is $B$ ( <br> $\boldsymbol{A}$ is not correct because this amino acid contains one acidic group and one basic group <br> $\boldsymbol{C}$ is not correct because this amino acid contains one acidic group and two basic groups <br> D is not correct because this amino acid contains one acidic group and two basic groups | (1) <br> Computer |


| Question <br> Number | The only correct answer is $\mathbf{A}$ |
| :--- | :--- | :--- |
| $\mathbf{2 0}$ | $\boldsymbol{B}$ is not correct because these reagents would lead to the formation of compound $\boldsymbol{B}$ |
| $\boldsymbol{C}$ is not correct because these reagents would lead to the formation of compound $\boldsymbol{B}$ |  |
|  | $\boldsymbol{D}$ is not correct because these reagents would lead to the formation of compound $\boldsymbol{B}$ |

## Section B

| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 21(a) | An explanation that makes reference to the following points: <br> - $\left(\mathrm{Hg}^{+}\right.$is $\left.[\mathrm{Xe}] 4 \mathrm{f}^{14}\right) 5 \mathrm{~d}^{10} 6 \mathrm{~s}^{1}$ <br> and <br> $\left(\mathrm{Hg}^{2+}\right.$ is $\left.[\mathrm{Xe}] 4 \mathrm{f}^{14}\right) 5 \mathrm{~d}^{10}\left(6 \mathrm{~s}^{0}\right)$ <br> - (d-block element as last) electron goes into a (5)d-orbital(s) (when the electronic configuration is written according to the Aufbau principle) <br> - (not transition element as) $\mathrm{Hg}^{+}$and $\mathrm{Hg}^{2+} /$ (stable) ions do not have incompletely filled (5)d-orbital(s) | Accept use of d-subshell for d-orbital(s) <br> Allow use of d-shell for d-subshell <br> Penalise use of just d-block for d-shell once only <br> Penalise use of $3 \mathrm{~d} / 4 \mathrm{~d}$ for 5 d once only <br> Allow Hg loses (only) its $\mathbf{6 s}$ electrons (when forming ions/compounds) <br> Do not award answer in terms of the electronic configuration of an ion of mercury <br> Allow $\mathrm{Hg}^{+}$and $\mathrm{Hg}^{2+} /($ stable) ions have completely full (5)d-orbital(s) <br> Ignore any reference to d-d transitions / other transition element properties <br> Do not award answer in terms of the electronic configuration of the element / an Hg atom | (3) <br> Expert |



| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 21(b)(ii) | An answer that makes reference to the following points: <br> - ionic half-equation for oxidation of mercury <br> - ionic half-equation for reduction of nitrate | Allow multiples and $\rightleftharpoons$ for $\rightarrow$ <br> Ignore state symbols, even if incorrect <br> Examples of ionic half-equations: $\begin{equation*} \mathrm{Hg} \rightarrow \mathrm{Hg}^{2+}+2 \mathrm{e}^{(-)} \tag{1} \end{equation*}$ <br> Allow $\mathrm{Hg}-2 \mathrm{e}^{(-)} \rightarrow \mathrm{Hg}^{2+}$ <br> Do not award half-equation including $\mathrm{HNO}_{3} / \mathrm{NO}_{3}{ }^{-}$ $4 \mathrm{H}^{+}+\mathrm{NO}_{3}^{-}+3 \mathrm{e}^{(-)} \rightarrow \mathrm{NO}+2 \mathrm{H}_{2} \mathrm{O}$ <br> Allow $3 \mathrm{H}^{+}+\mathrm{HNO}_{3}+3 \mathrm{e}^{(-)} \rightarrow \mathrm{NO}+2 \mathrm{H}_{2} \mathrm{O}$ <br> Allow $4 \mathrm{HNO}_{3}+3 \mathrm{e}^{(-)} \rightarrow \mathrm{NO}+2 \mathrm{H}_{2} \mathrm{O}+3 \mathrm{NO}_{3}^{-}$ | (2) <br> Expert |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :--- | :--- | :---: |
| 21(b)(iii) | An answer that makes reference to the following point: | Example of completed equation: | (1) |
|  | $\bullet$ balanced equation | $\underline{\mathbf{3} H g(1)+\underline{\mathbf{8}} \mathrm{HNO}_{3}(\mathrm{aq}) \rightarrow \underline{\mathbf{3} H g}\left(\mathrm{NO}_{3}\right)_{2}(\mathrm{aq})+\underline{\mathbf{2}} \mathrm{NO}(\mathrm{g})+\underline{\mathbf{4}} \mathrm{H}_{2} \mathrm{O}(1)}$ | Clerical |
|  |  | Allow multiples |  |


| Question <br> Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 21(c)(i) | An answer that makes reference to the following point: <br> - correct species <br> - balanced equation | (1) (1) | Example of completed equation: $\mathrm{Hg}\left(\mathrm{NO}_{3}\right)_{2}+3 \mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH} \rightarrow \mathrm{Hg}(\mathrm{CNO})_{2}+2 \mathrm{CH}_{3} \mathrm{CHO}+5 \mathrm{H}_{2} \mathrm{O}$ <br> Ignore state symbols even if incorrect <br> Do not award molecular formulae eg $\mathrm{C}_{2} \mathrm{H}_{4} \mathrm{O}$ for $\mathrm{CH}_{3} \mathrm{CHO}$ <br> Do not award $\mathrm{CH}_{3} \mathrm{COH}$ for $\mathrm{CH}_{3} \mathrm{CHO}$ <br> Allow multiples <br> No TE on M1 except on correct molecular formulae and on $\mathrm{CH}_{3} \mathrm{COH}$ | (2) <br> Graduate |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 21(c)(ii) |  |  | Correct answer with some working scores (3) | (3) |
|  |  |  | Ignore SF except 1SF throughout | Expert |
|  |  |  | Example of calculation: |  |
|  | - moles of $\mathrm{Hg}(\mathrm{CNO})_{2}$ | (1) | $\mathrm{n}=\frac{1.00}{284.6}=0.0035137 / 3.5137 \times 10^{-3}$ |  |
|  | - moles of gas produced | (1) | $\begin{aligned} & \mathrm{n}=0.0035137 \times 2=0.0070274 / 7.0274 \times 10^{-3} \\ & \text { TE on M1 } \end{aligned}$ |  |
|  | - volume of gas produced | (1) | $\begin{aligned} & \mathrm{v}=0.0070274 \times 24000=168.66\left(\mathrm{~cm}^{3}\right) \\ & \text { Accept } 0.16866 \mathbf{d m}^{3} \\ & \mathrm{TE} \text { on M2 } \end{aligned}$ |  |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{2 1 ( d ) ( i ) ~}$ | An answer that makes reference to the following point: | Allow KCl for $\mathrm{Cl}^{-}$throughout | (1) |
|  | $\bullet$ (to provide a) constant concentration (of $\mathrm{Cl}^{-}$) | Allow to keep the solution $/ \mathrm{Cl}^{-}$saturated | Expert |
|  |  | Ignore just to provide $\mathrm{Cl}^{-}$ <br> Ignore stated concentrations <br> Do not award salt bridge $/$to complete the circuit <br> Do not award catalyst |  |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :--- | :---: |
| 21(d)(ii) | An answer that makes reference to the following point: |  | (1) |
|  | - $(0.24-0.37=)-0.13(\mathrm{~V})$ | Ignore working, even if incorrect | Expert |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 21(d)(iii) | An answer that makes reference to the following point: <br> - $\mathrm{Hg}_{2} \mathrm{Cl}_{2}+\mathrm{Sn} \rightarrow 2 \mathrm{Hg}+\mathrm{Sn}^{2+}+2 \mathrm{Cl}^{-}$ | Allow $\mathrm{Hg}_{2} \mathrm{Cl}_{2}+\mathrm{Sn} \rightarrow 2 \mathrm{Hg}+\mathrm{SnCl}_{2}$ <br> Allow multiples <br> Allow $\rightleftharpoons$ for $\rightarrow$ <br> Ignore state symbols even if incorrect <br> Ignore half-equations even if incorrect <br> Ignore use of cell diagrams <br> Do not award uncancelled electrons <br> Do not award $2 \mathrm{Hg}^{+}\left(+2 \mathrm{Cl}^{-}\right)$for $\mathrm{Hg}_{2} \mathrm{Cl}_{2}$ <br> If answer to (d)(ii) is $\mathbf{+ 0 . 6 1}(\mathrm{V}) / \boldsymbol{+ 0 . 3 7}(\mathrm{V}) /$ greater than $+0.24(\mathrm{~V})$, equation must be reversed: <br> $2 \mathrm{Hg}+\mathrm{Sn}^{2+}+2 \mathrm{Cl}^{-} \rightarrow \mathrm{Hg}_{2} \mathrm{Cl}_{2}+\mathrm{Sn}$ <br> OR <br> $2 \mathrm{Hg}+\mathrm{SnCl}_{2} \rightarrow \mathrm{Hg}_{2} \mathrm{Cl}_{2}+\mathrm{Sn}$ | (1) <br> Expert |



| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 21(d)(v) | An answer that makes reference to any one of the following points: <br> - (calomel electrode) does not require a (separate) salt bridge <br> OR <br> (calomel electrode) does not require a continuous supply of hydrogen / gas <br> OR <br> platinum/Pt (of hydrogen electrode is) easily poisoned <br> OR <br> difficult to ensure hydrogen electrode is at equilibrium | Ignore calomel electrode is quicker to use / easier to set up / done in the same container / more portable <br> Accept does not require a hydrogen / gas generator <br> Ignore just does not require hydrogen / gas <br> Ignore any reference to pressure <br> Ignore hydrogen is flammable / explosive / <br> difficult to store <br> Ignore (calomel electrode is) safer <br> Ignore platinum is expensive <br> Ignore (calomel electrode) is cheaper <br> Allow (calomel electrode) reaches equilibrium sooner <br> Allow (calomel electrode gives) more stable (reading) <br> Ignore (calomel electrode is) more accurate <br> Ignore calomel electrode potential is more positive | (1) <br> Expert |


| Question <br> Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 22 | An answer that makes reference to the following points: |  | Correct answer to 2SF or 3SF with some working scores (6) | (6) |
|  |  |  | Ignore SF except 1SF | Expert |
|  |  |  | Example of calculation: |  |
|  | - moles of $\mathrm{FeSO}_{4}$ | (1) | $\mathrm{n}=0.0500 \times \frac{25.95}{1000}=0.0012975 / 1.2975 \times 10^{-3}$ |  |
|  | - moles of excess $\mathrm{MnO}_{4}{ }^{-}$ | (1) | $\mathrm{n}=0.0012975 \div 5=0.0002595 / 2.595 \times 10^{-4}$ <br> TE on moles of $\mathrm{FeSO}_{4}$ |  |
|  | - initial moles of $\mathrm{MnO}_{4}^{-}$ |  | $\mathrm{n}=0.0100 \times \frac{50.0}{1000}=0.0005 / 5 \times 10^{-4}(\text { Allow } 1 \mathrm{SF})$ |  |
|  | and moles of $\mathrm{MnO}_{4}^{-}$reacted | (1) | and $\mathrm{n}=0.0005-0.0002595=0.0002405 / 2.405 \times 10^{-4}$ |  |
|  |  |  | TE on moles of excess $\mathrm{MnO}_{4}{ }^{-}$provided answer is positive |  |
|  | - moles of $\mathrm{C}_{2} \mathrm{O}_{4}{ }^{2-}$ | (1) | $\begin{aligned} & \mathrm{n}=0.0002405 \times 2.5=0.00060125 / 6.0125 \times 10^{-4} \\ & \mathrm{TE} \text { on moles of } \mathrm{MnO}_{4}^{-} \text {reacted } \end{aligned}$ |  |
|  | - mass of $\mathrm{CaC}_{2} \mathrm{O}_{4}$ | (1) | $\text { mass }=0.00060125 \times 128.1=0.077020(\mathrm{~g})$ <br> TE on moles of $\mathrm{C}_{2} \mathrm{O}_{4}{ }^{2-}$ |  |
|  | - percentage by mass of $\mathrm{CaC}_{2} \mathrm{O}_{4}$ and answer to 2 SF or 3 SF | (1) | $\begin{aligned} & \% \text { mass }=\frac{0.077020}{11.4} \times 100=0.67562(\%) \\ & =0.68 / 0.676(\%) \end{aligned}$ <br> TE on mass of $\mathrm{CaC}_{2} \mathrm{O}_{4}$ provided positive value to $2 \mathrm{SF} / 3 \mathrm{SF}$ and $<100 \%$ Allow use of 128 for $M_{\mathrm{r}}$ of $\mathrm{CaC}_{2} \mathrm{O}_{4}$ giving 0.675 (\%) |  |

(Total for Question 22 = 6 marks)


## Indicative points:

- IP1: thermochemical data calculation
(enthalpy of hydrogenation of 1,3,5-cyclohexatriene / benzene is) expected to be $-360\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$
- IP2: thermochemical data comparison
(enthalpy of hydrogenation is) less exothermic / less negative than expected (for 1,3,5-cyclohexatriene)


## less exothermic / more stable by $152\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ scores IP1 and IP2

- IP3: X-ray diffraction data
(carbon-carbon) bond lengths in benzene are equal
Ignore any reference to bond strength / bond angle in IP3 and IP4
- IP4: X-ray diffraction data (carbon-carbon) bond length in benzene is longer than (localised) $\mathrm{C}=\mathrm{C}$ (in cyclohexene)
- IP5: Bromination data
(product for benzene is formed by electrophilic) substitution
- IP6: Bromination data
(benzene $\pi$-bonds less reactive than localised $\pi$-bonds and) requires $\left(\mathrm{FeBr}_{3}\right)$ catalyst (and heat)

Allow (enthalpy of hydrogenation is) expected to be three times the value for cyclohexene
Allow (enthalpy of hydrogenation is) different by $152\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$
Accept reverse argument
Ignore higher/lower for less exothermic
Ignore benzene more exothermic than cyclohexene
Ignore just benzene more stable than expected
Do not award enthalpy required/needed

Allow (carbon-carbon) bond lengths are not different Allow cyclohexene (carbon-carbon) bond lengths are different

Accept (carbon-carbon) bond length is in between $\mathrm{C}=\mathrm{C}$ and $\mathrm{C}-\mathrm{C}$ (in cyclohexene)
Allow (carbon-carbon) bond length is shorter than $\mathrm{C}-\mathrm{C}$ (in cyclohexene)

Allow (benzene) does not react by addition
Allow cyclohexene/localised $\pi$-bonds react by addition Ignore any equations / mechanisms
Do not award nucleophilic (substitution / addition)
Accept cyclohexene does not require a catalyst
Allow halogen carrier for catalyst
Ignore just benzene does not decolourise bromine water Do not award Fe catalyst

| Question <br> Number | Answer | Additional Guidance | Mark |
| :--- | :--- | :--- | :---: |
| 24(a)(i) | An answer that makes reference to the following point: | Allow capital letters and spaces | (1) |
|  | - prop-2-enamide / 2-propenamide | Ignore omission of hyphen <br> Allow propenamide <br> Allow 'ene' for 'en' <br> Allow propyl for prop |  |
|  |  | Graduate <br> Do not award propan for prop <br> Do not award N- prefix <br> Do not award cis/trans/E/Z- prefix |  |


| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 24(a)(ii) | An explanation that makes reference to the following points: <br> - (PAM can form many) hydrogen bonds with water <br> - H-bonds (with water) can form at $\mathrm{NH}_{2}$ and ( $\mathrm{C}=$ ) O <br> - diagram of (at least one) hydrogen bond between a water molecule and any amide group | (1) <br> (1) <br> (1) | Allow M1 from a labelled diagram <br> Ignore PAM reacts with water / acts as a base / accepts a proton from water / forms $\mathrm{RNH}_{3}{ }^{+}$ <br> M2 can be awarded from a diagram <br> diagram must include: <br> H -bond to lone pair on O or N <br> and <br> $\delta+\mathrm{H}$ atom <br> Ignore bond angle <br> Do not award H-bond shown as coordinate bond / solid line (ie covalent bond) <br> Example of diagram scoring (3): <br> Allow H -bond between lone pair on N of $\mathrm{NH}_{2}$ and $\delta+\mathrm{H}$ of water | (3) <br> Expert |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 24(a)(iii) | An explanation that makes reference to the following points: <br> - carboxylate / $\mathrm{COO}^{-}$(above pH 8 ) <br> - repulsion between negative charges (above pH 8 ) | Allow carboxylic acid/COOH/OH groups are deprotonated / donate $\mathrm{H}^{+}$/ become anions <br> Allow $\mathrm{OH}^{-}$removes H atoms involved in hydrogen bonds <br> Ignore just PAA is deprotonated / donates $\mathrm{H}^{+}$/ becomes anion <br> Ignore just salt is formed <br> Do not award zwitterion is formed <br> Allow ( $\mathrm{COO}^{-}$) cannot form (intramolecular) hydrogen bonds Allow (all) hydrogen bonds break <br> Ignore hydrogen bonds weaken <br> Ignore fewer hydrogen bonds <br> Ignore any reference to denaturation <br> Ignore any reference to intermolecular hydrogen bonds | (2) <br> Expert |


| Question <br> Number | Answer | Additional Guidance |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 4 ( b ) ( i )}$ | An answer that makes reference to the following <br> point: | Allow any combination of skeletal, structural or displayed <br> formulae |
| (1) |  |  |
|  | estructure of vinylpyrrolidone monomer |  |


| Question <br> Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 24(b)(ii) | An answer that makes reference to the following points: <br> - molar mass of vinylpyrrolidone monomer / PVP repeat unit <br> - number of monomers per polymer and answer to nearest whole number | (1) (1) | Correct answer with some working scores (2) <br> Example of calculation: $M\left(\mathrm{C}_{6} \mathrm{H}_{9} \mathrm{NO}\right)=111.0 / 111\left(\mathrm{~g} \mathrm{~mol}^{-1}\right)$ <br> TE on (b)(i) if molar mass is not $111.0 / 111$ $\begin{aligned} & \text { number }=90000 \div 111.0=810.81=811 \\ & \text { TE on M1 } \end{aligned}$ | (2) <br> Expert |



| Question Number | Answer |  | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 24(c)(i) | An answer that makes reference to the following points: <br> - (polymer is a very) large molecule <br> OR <br> (polymer is formed from) large number of / many monomers <br> - (condensation as) splitting off of a (small) molecule | (1) <br> (1) | Allow long-chain (molecule) <br> Allow macromolecule <br> Allow repeating for many <br> Ignore 2 or more / several / different for many <br> Ignore (formed by) addition <br> Allow with loss/elimination of $\mathrm{H}_{2} \mathrm{O} / \mathrm{HCl}$ Ignore forms byproduct | (2) <br> Expert |



| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 24(c)(iii) | An answer that makes reference to the following points: <br> - amide linkage (within polymer chain) <br> - two repeat units | Allow any combination of skeletal, structural or displayed formulae <br> Allow -NHCO- / -CONH- / -HNCO- / -OCNH- <br> Ignore omission of square brackets <br> Ignore $n$ <br> Examples of two repeat units: $-\mathrm{CO}\left(\mathrm{CH}_{2}\right)_{5} \mathrm{NHCO}\left(\mathrm{CH}_{2}\right)_{5} \mathrm{NH}-$  $-\mathrm{NH}\left(\mathrm{CH}_{2}\right)_{5} \mathrm{CONH}\left(\mathrm{CH}_{2}\right)_{5} \mathrm{CO}-$ | (2) <br> Expert |

(Total for Question 24 = 18 marks)
Total for Section B=50 marks

## Section C

| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 25(a) | An answer that makes reference to the following points: <br> - ester <br> - (primary) amine <br> and arene/benzene/phenyl <br> OR <br> phenylamine | Ignore any structures / formulae <br> Ignore carbonyl <br> Do not award ketone / aldehyde / carboxylic acid <br> Do not award ether <br> Allow amino <br> Allow aryl <br> Ignore alkyl/alkane <br> Do not award alkene <br> Do not award phenol <br> Allow aniline <br> Allow aromatic amine | (2) <br> Graduate |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 25(b)(i) | An answer that makes reference to the following points: <br> - molecular formulae of procaine and HCl <br> - molecular formula of procaine monohydrochloride | Ignore non-molecular formulae $\begin{equation*} \mathrm{C}_{13} \mathrm{H}_{20} \mathrm{~N}_{2} \mathrm{O}_{2}+\mathrm{HCl} \tag{1} \end{equation*}$ <br> Allow elements in any order $\begin{equation*} \mathrm{C}_{13} \mathrm{H}_{21} \mathrm{Cl}^{(-)} \mathrm{N}_{2}{ }^{(+)} \mathrm{O}_{2} \tag{1} \end{equation*}$ <br> Allow elements in any order TE on molecular formula of procaine <br> Ignore position of charges <br> Do not award separate $\mathrm{C}_{13} \mathrm{H}_{21} \mathrm{~N}_{2}{ }^{+} \mathrm{O}_{2}$ and $\mathrm{Cl}^{-}$ions Do not award any additional product(s) <br> Example of equation: $\mathrm{C}_{13} \mathrm{H}_{20} \mathrm{~N}_{2} \mathrm{O}_{2}+\mathrm{HCl} \rightarrow \mathrm{C}_{13} \mathrm{H}_{21} \mathrm{ClN}_{2} \mathrm{O}_{2} \text { scores (2) }$ | (2) <br> Graduate |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 25(b)(ii) | An explanation that makes reference to the following points: <br> - identification of tertiary amine nitrogen and effect of ethyl groups OR benzene ring <br> EITHER <br> ethyl / alkyl (groups) are electron donating <br> OR <br> lone pair (on N of $\mathrm{NH}_{2}$ partially) delocalised into (aromatic) $\pi$-bond(s) <br> - second effect | Ignore just comparison of electron density on N atoms <br> Ignore just comparison of ability of $(\mathrm{N})$ lone pairs to accept $\mathrm{H}^{+}$ <br> Accept any unambiguous identification <br> Accept ethyl / alkyl has positive inductive effect <br> Allow ethyl / alkyl are electron pushing / electron releasing <br> Allow methyl / R / attached groups for ethyl / alkyl <br> Accept non-bonding pair for lone pair <br> Allow electron pair for lone pair <br> Allow overlaps with / interacts with / released into / drawn into for delocalised into <br> Allow p-orbitals / ring for (aromatic) $\pi$-bond(s) <br> Ignore just benzene for (aromatic) $\pi$-bond(s) <br> Ignore just ring is electron withdrawing (with no mention of electron pair) <br> If no other mark awarded, tertiary / aliphatic amine is more basic <br> OR <br> aromatic / primary amine is less basic <br> scores (1) | (2) <br> Expert |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{2 5 ( c )}$ | An answer that makes reference to the following point: | Allow any combination of skeletal, structural or <br> displayed formulae <br> Example of structure: | (1) |
|  |  |  |  |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 25(d)(i) | An answer that makes reference to the following points: <br> - equation for formation of nitronium ion <br> - curly arrow from within hexagon to anywhere on $\mathrm{NO}_{2}{ }^{+}$ <br> - structure of intermediate ion <br> - curly arrow from $\mathrm{C}-\mathrm{H}$ bond to within ring and correct product and $\mathrm{H}^{+}$ | Ignore omission or incorrect placement of methyl groups in M2 and M3 $\begin{equation*} \mathrm{HNO}_{3}+\mathrm{H}_{2} \mathrm{SO}_{4} \rightarrow \mathrm{NO}_{2}^{+}+\mathrm{HSO}_{4}^{-}+\mathrm{H}_{2} \mathrm{O} \tag{1} \end{equation*}$ <br> OR $\mathrm{HNO}_{3}+2 \mathrm{H}_{2} \mathrm{SO}_{4} \rightarrow \mathrm{NO}_{2}^{+}+2 \mathrm{HSO}_{4}^{-}+\mathrm{H}_{3} \mathrm{O}^{+}$ <br> OR $\mathrm{HNO}_{3}+\mathrm{H}_{2} \mathrm{SO}_{4} \rightarrow \mathrm{H}_{2} \mathrm{NO}_{3}{ }^{+}+\mathrm{HSO}_{4}^{-} \text {and } \mathrm{H}_{2} \mathrm{NO}_{3}{ }^{+} \rightarrow \mathrm{NO}_{2}{ }^{+}+\mathrm{H}_{2} \mathrm{O}$  <br> TE on electrophile from M1 provided positively charged Do not award lone pair on N of $\mathrm{NO}_{2}{ }^{+}$ <br> Allow any part of gap in 'horseshoe' facing tetrahedral carbon and covering at least three carbons with some part of positive sign within 'horseshoe'. 'Horseshoe' may be dashed <br> TE on electrophile from M2 <br> Do not award $\mathrm{NO}_{2}-\mathrm{C}$ connectivity <br> Do not award dashed $\mathrm{C}-\mathrm{H}$ and $\mathrm{C}-\mathrm{N}$ bonds unless 3D structure | (4) <br> Expert |



| Question <br> Number | Answer | Additional Guidance | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{2 5 ( d ) ( i i i )}$ | An answer that makes reference to the following point: |  | (1) |
|  | $\bullet$ tin and (concentrated) hydrochloric acid | Accept Sn and $\mathrm{HCl}((\mathrm{aq}))$ <br> Ignore heat $/$ reflux <br> Ignore NaOH in second step <br> Do not award NaOH with Sn and HCl in the same step <br> Do not award any reference to catalyst | Graduate |


| Question | Answer | Additional Guidance | Mark |  |
| :--- | :---: | :---: | :---: | :---: |
| Number | An answer that makes reference to the following point: | Ignore non-skeletal formulae <br> Ignore bond angles and bond lengths | Graduate |  |
|  |  | skeletal formula of 2-chloroethanoyl chloride |  |  |


| Question <br> Number | Answer | Additional Guidance | Mark |
| :--- | :---: | :--- | :---: |
| $\mathbf{2 5 ( d ) ( v )}$ | An answer that makes reference to the following point: |  | (1) |
|  | $\bullet$ nucleophilic substitution | Allow $\mathrm{S}_{\mathrm{N}} 2 / \mathrm{S}_{\mathrm{N}} 1$ | Clerical |
|  |  | Do not award any other answer |  |


| Question <br> Number | Answer | Additional Guidance | Mark |  |
| :--- | :--- | :--- | :---: | :---: |
| 25(e) | An answer that makes reference to the following point: | (1) | Graduate | articaine |


| Question Number | Answer | Additional Guidance | Mark |
| :---: | :---: | :---: | :---: |
| 25(f) | An answer that makes reference to the following points: | Example of calculation: | (3) |
|  | - number of half-lives in 4 hours | $\text { half-lives }=\frac{(4 \times 60)}{20}=12$ | Expert |
|  | - mass of articaine remaining in mg | $\text { mass }=100 \times 0.5^{12}=0.024414(\mathrm{mg})$ <br> TE on M1 <br> Ignore SF except 1 SF |  |
|  | - conversion of mg to $\mu \mathrm{g}$ | $\text { mass }=0.024414 \times 1000=24.414(\mu \mathrm{~g})$ <br> TE on M1 and M2 |  |

(Total for Question $25=20$ marks)
Total for Section C=20 marks
Total for Paper $=90$ marks

