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## Mark Scheme (Results)

## Summer 2018

Pearson Edexcel International
Advanced Level In Chemistry (WCH05)
Paper 01
General Principles of Chemistry II - 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 (multiple choice)

| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1}$ | The only correct answer is A <br> B is not correct because this is the other common oxidation number <br> associated with iron <br> C is not correct because this is the numerical value of the charge on <br> the complex | (1) |
| D is not correct because this is the number of cyano ligands in the <br> complex |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{2}$ | The only correct answer is B <br> A is not correct because this is calculated on the basis of a 1:2 <br> reaction and that 2 mol hydrogen atoms forms 1 mol hydrogen gas | (1) |
|  | C is not correct because this is calculated on the basis of a 1:2 <br> reaction and does not take into account that 2 mol hydrogen atoms <br> forms 1 mol hydrogen gas | D is not correct because this is calculated on the basis of a 1:3 <br> reaction but does not take into account that 2 mol hydrogen atoms <br> forms 1 mol hydrogen gas |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{3 ( a )}$ | The only correct answer is D <br> A is not correct because both oxidation states in the system must be <br> present | (1) |
| B is not correct because both oxidation states in the system must be <br> present <br> C is not correct because the electrode must be inert |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{3 ( b )}$ | The only correct answer is B | (1) |
| A is not correct because the value for the Fe(III)/Fe(II) electrode <br> reaction has been incorrectly doubled <br> C is not correct because the subtraction has been incorrectly <br> reversed <br> D is not correct because the value for the Fe(III)/Fe(II) electrode <br> reaction has been incorrectly doubled and the subtraction has been <br> incorrectly reversed |  |  |


| Question | Correct Answer | Mark |
| :--- | :--- | :---: |
| Number | The only correct answer is C | (1) |
| $\mathbf{4}$ | A is not correct because if $\mathrm{E}_{\text {cell }}$ is positive $\Delta \mathrm{S}_{\text {total }}$ must be positive but <br> $\Delta \mathrm{S}_{\text {system }}$ could be negative <br> B is not correct because if $\mathrm{E}_{\text {cell }}$ is positive $\Delta \mathrm{S}_{\text {total }}$ must be positive but <br> $\Delta \mathrm{S}_{\text {surroundings }}$ could be negative <br> $\mathbf{D}$ is not correct because if $\mathrm{E}_{\text {cell }}$ is positive $\Delta \mathrm{S}_{\text {total }}$ must be positive but <br> $\Delta \mathrm{S}_{\text {system }}$ or $\Delta S_{\text {surroundings }}$ could be negative |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{5}$ | The only correct answer is D <br> A is not correct because hydrogen is the fuel so must be oxidised, <br> hence oxygen is reduced <br> B is not correct because hydrogen is the fuel so must be oxidised, <br> hence oxygen is reduced <br> $\mathbf{C}$ is not correct because reduction occurs at the positive electrode <br> (the cathode) | (1) |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{6}$ | The only correct answer is A | (1) |
|  | B is not correct because a large number of valence electrons is a <br> factor in heterogeneous catalysis but not in homogeneous catalysis <br> C is not correct because active sites are involved in heterogeneous <br> catalysis but not in homogeneous catalysi |  |
| $\mathbf{D}$ is not correct because $\mathrm{Fe}^{3+}$ cannot oxidise peroxodisulfate ions |  |  |$\quad$|  |
| :--- |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{7}$ | The only correct answer is B <br> A is not correct because an atom of vanadium has three unpaired <br> electrons <br> $\mathbf{C}$ is not correct because an atom of manganese has five unpaired <br> electrons <br> $\mathbf{D}$ is not correct because an atom of iron has four unpaired electrons | (1) |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{8}$ | The only correct answer is B | (1) |
| A is not correct because this formula would give a peak in the graph <br> when equimolar quantities of M \& L were present (at $5 \mathrm{~cm}^{3}$ ) <br> C is not correct because this answer is based on the ratio of 3.33: 10 <br> (rather than 3.33:(10-3.33) <br> D is not correct because this is a common complex formula but <br> incorrect in this case |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{9}$ | The only correct answer is B <br> A is not correct because the addition of water has no effect apart <br> from hydrating the ions. The addition of sulfuric acid changes the <br> vanadium species but not the oxidation state. Addition of zinc <br> reduces V(V) to V(II). So there are two oxidation states | (1) |
|  | C is not correct because the addition of water has no effect apart <br> from hydrating the ions. The addition of sulfuric acid changes the <br> vanadium species but not the oxidation state. Addition of zinc <br> reduces V(V) to V(II). So there are two oxidation states | D is not correct because the addition of water has no effect apart <br> from hydrating the ions. The addition of sulfuric acid changes the <br> vanadium species but not the oxidation state. Addition of zinc <br> reduces V(V) to V(II). So there are two oxidation states |


| Question | Correct Answer | Mark |
| :--- | :--- | :---: |
| Number | The only correct answer is C | (1) |
| $\mathbf{1 0}$ | A is not correct because this value is obtained by addition of the <br> enthalpy changes given in the stem <br> B is not correct because this value is obtained by subtraction of the <br> enthalpy changes given in the stem <br> D is not correct because this value is half the correct value |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1}$ | The only correct answer is C | (1) |
|  | A is not correct because a true statement but irrelevant |  |
| B is not correct because a true statement but irrelevant |  |  |
| D is not correct because a true statement but irrelevant |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 2}$ | The only correct answer is D | (1) |
|  | A is not correct because arenes will not undergo electrophilic <br> addition reactions | B is not correct because electrophilic substitution is a characteristic <br> reaction of arenes but will not occur with dilute sulfuric acid |
| $\mathbf{C}$ is not correct because phenylamine is not hydrolysed |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 3}$ | The only correct answer is D <br> A is not correct because dipole-dipole forces are possible but the <br> high melting temperature of alanine is due to the ionic forces <br> between the zwitterions | (1) |
|  | B is not correct because London forces are present but the high <br> melting temperature of alanine is due to the ionic forces between <br> the zwitterions | C is not correct because hydrogen bonds are possible but the high <br> melting temperature of alanine is due to the ionic forces between <br> the zwitterions |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 4 ( a )}$ | The only correct answer is C <br> A is not correct because a nitrogen with two H atoms attached is an <br> amine | (1) |
| B is not correct because the group lower far right is an ester but |  |  |
| candidates might overlook the methyl group |  |  |
| D is not correct because Benzene ring is a phenyl group. Candidates |  |  |
| might mistake phenyl for phenol |  |  |$\quad$|  |
| :--- |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 4 ( b )}$ | The only correct answer is A <br> B is not correct because candidates might identify one or both of the <br> $\mathrm{CH}_{2}$ groups as asymmetric <br> C is not correct because candidates might identify one or both of the <br> $\mathrm{CH}_{2}$ groups as asymmetric | (1) |
| D is not correct because candidates might identify all of the C atoms <br> in the aliphatic chain as asymmetric |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 5}$ | The only correct answer is A | (1) |
|  | B is not correct because the amide group has been reversed <br> implying a single monomer with an acid (chloride) and an amine | C is not correct because Kevlar has benzene rings |
| Dis not correct because the amide group has been reversed <br> implying a single monomer with an acid (chloride) and an amine <br> group, and because Kevlar has benzene rings |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 6}$ | The only correct answer is A | (1) |
|  | B is not correct because the molecular ion peak will be at 73 but <br> $\mathrm{m} / \mathrm{e}=30$ requires $\mathrm{CH}_{2} \mathrm{NH}_{2}$ fragment <br> $\mathbf{C}$ is not correct because the molecular ion peak will be at 73 but <br> $\mathrm{m} / \mathrm{e}=30$ requires $\mathrm{CH}_{2} \mathrm{NH}_{2}$ fragment <br> D is not correct because because The molecular ion peak will be at <br> 73 but $\mathrm{m} / \mathrm{e}=30$ requires $\mathrm{CH}_{2} \mathrm{NH}_{2}$ fragment |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 7}$ | The only correct answer is D <br> A is not correct because equilibrium position is not affected by use <br> of a support <br> B is not correct because polymeric supports are often expensive and <br> this is a disadvantage of their use <br> $\mathbf{C}$ is not correct because side reactions can occur but this is a <br> disadvantage (unlike combinatorial chemistry) | (1) |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 8}$ | The only correct answer is D | (1) |
|  | $\mathbf{A}$ is not correct because not used at all in solvent extraction | B is not correct because not used at all in solvent extraction |
| $\mathbf{C}$ is not correct because not used at all in solvent extraction |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(a)(i) | M1 <br> Structure of $\mathrm{CO}_{2}$ and curly arrow from $\mathrm{C}=\mathrm{O}$ bond to the oxygen atom or just beyond it. <br> ALLOW dipolar electrophile <br> M2 <br> Curly arrow from on or within the circle to the carbon of the electrophile (molecule or dipolar ion) <br> ALLOW <br> Curly arrow from anywhere within the hexagon <br> Arrow to any part of the electrophile including to the $\delta+$ charge (if given). TE on any electrophile <br> M3 <br> Intermediate structure including charge with horseshoe covering at least 3 carbon atoms, and facing the tetrahedral carbon and with some part of the positive charge within the horseshoe <br> and negative charges $\mathrm{COO}^{-}$and $\mathrm{O}^{-}\left(\mathrm{OR} \mathrm{COO}^{-}\right.$and OH ) <br> ALLOW <br> dotted horseshoe <br> Do NOT penalise incorrect position of the $\mathrm{O}^{-}$/ phenol group at this marking point <br> M4 <br> Curly arrow from $\mathrm{C}-\mathrm{H}$ bond to anywhere in the benzene ring reforming delocalized structure with $\mathrm{O}^{-}$ / phenol group in the 2 position and carboxylate ion <br> Correct Kekulé structures score full marks <br> ALLOW phenol group as OH throughout | electrophile with a net charge <br> Curly arrow starting on or outside the hexagon <br> Partial bonds to H and $\mathrm{CO}_{2}{ }^{-}$ unless clearly part of a 3D structure <br> carboxylic acid group | (4) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9 ( a ) ( i i )}$ | Any strong acid by name or formula <br> e.g. sulfuric acid $/ \mathrm{H}_{2} \mathrm{SO}_{4}$ <br> hydrochloric acid $/ \mathrm{HCl}(\mathrm{aq})$ <br> ALLOW <br> nitric acid $/ \mathrm{HNO}_{3} / \mathrm{HCl}$ <br> If name and formula are given, both must <br> be correct <br> IGNORE <br> conc / dilute / H <br> just 'acid' / 'strong acid' | (1) |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(b) | Higher temperature / pressure <br> IGNORE <br> References to catalysts / use of milder / harsher (conditions) <br> Phenol reacts with electrophiles much faster / under milder conditions than benzene <br> ALLOW <br> Phenol reacts with electrophiles more easily / readily <br> OR <br> Phenol more reactive than benzene / phenol reacts faster than benzene / phenol more susceptible to attack <br> Because the electron density of the benzene ring in phenol is higher <br> Due to interaction between the (oxygen) lone pair and the $\pi$ electrons of the benzene ring <br> ALLOW <br> (Oxygen / OH) lone pair donated to the benzene ring <br> Two electrons from oxygen donated to the benzene ring <br> Or <br> Reverse argument for benzene | charge density (of ring) | (4) |
| Question Number | Acceptable Answers | Reject | Mark |
| 19(c)(i) | Esterification <br> ALLOW <br> Ester formation <br> Ethanoylation <br> Acetylation <br> Acylation <br> IGNORE <br> Addition-elimination / Condensation | Friedel-Crafts | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(c)(ii) | Any three of the following <br> Feasibility <br> Ethanoic acid does not react with phenol <br> OH groups to form esters <br> Reference to equilibrium or completion <br> Reaction with ethanoic acid would be an equilibrium / reversible / does not go to completion (so low yield) <br> OR <br> Reaction with ethanoyl chloride goes to completion (so high yield) <br> Reference to rate of reaction <br> Reaction with ethanoic acid is slow or needs a catalyst or requires heat <br> OR <br> Reaction with ethanoyl chloride is fast or does not require a catalyst or occurs at room temperature <br> Reference to products <br> Toxic / poisonous HCl is a by-product <br> ALLOW <br> Corrosive <br> IGNORE <br> Reference to thermicity <br> Reference to reactivity <br> Reference to 'vigorous' reactions <br> Explanations of reactivity, even if incorrect <br> Safety / ease of handling ethanoic acid <br> Reference to cost <br> ALLOW <br> Reverse arguments | Just ' HCl is formed' Just 'harmful' | (3) |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(a)(i) | Left-hand label: <br> $\mathrm{H}^{+}((\mathrm{aq})) / \mathrm{H}_{3} \mathrm{O}^{+}((\mathrm{aq})) /$ hydrogen ions OR <br> $\mathrm{HCl}((\mathrm{aq})) / \mathrm{H}_{2} \mathrm{SO}_{4}((\mathrm{aq})) / \mathrm{HNO}_{3}((\mathrm{aq}))$ <br> and <br> $1 \mathrm{~mol} \mathrm{dm}^{-3} /$ activity $=1$ <br> Right- hand label: <br> $\mathrm{H}_{2}((\mathrm{~g}))$ / hydrogen <br> and <br> 1 atm <br> / 101 / $100 \mathrm{kPa}\left(\mathrm{kN} \mathrm{m}^{-2}\right)$ <br> / 101000 / $100000 \mathrm{~Pa}\left(\mathrm{~N} \mathrm{~m}^{-2}\right)$ <br> / 1.01 /1.0 /1 (Bar) <br> IGNORE <br> Temperature <br> Both substances correct but any number of conditions omitted scores | $\mathrm{H}_{2} \mathrm{SO}_{4}((\mathrm{aq}))$ with concentration of 0.5 $\mathrm{mol} \mathrm{dm}{ }^{-3}$ | (2) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( a ) ( i i )}$ | Platinum is used because it is (chemically) <br> inert (and conducts electricity) <br> ALLOW <br> unreactive <br> OR <br> Platinum catalyses $2 \mathrm{H}^{+}+2 \mathrm{e} \rightleftharpoons \mathrm{H}_{2}$ <br> OR <br> Platinum catalyses $\quad \mathrm{H}_{2} \rightleftharpoons 2 \mathrm{H}^{+}+2 \mathrm{e}$ | (2) |  |
| ALLOW <br> Platinum catalyses the electrode reaction <br> Platinum is a catalyst <br> Using platinum black increases the <br> surface area (making the catalysis more <br> efficient) <br> ALLOW <br> Increase the number of active sites $\quad$ (1) <br> IGNORE <br> Reference to cost or rate of reaction |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( a ) ( i i i )}$ | $0.0($ V) /zero <br> ALLOW <br> $0(V)$ <br> IGNORE <br> Charges |  | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( a ) ( i v )}$ | It is not possible to measure the <br> potential difference (of a half cell) <br> between the metal electrode and the ion <br> solution | ALLOW <br> A potential difference requires a <br> complete circuit/two electrodes <br> OR <br> Current will not flow unless the circuit is <br> complete <br> OR <br> The potential of a single electrode can <br> only be measured by reference to / <br> comparison with / against another <br> electrode. |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( a ) ( v )}$ | M1 <br> Conditions of the reaction are not <br> standard <br> OR <br> Concentration is not $1 \mathrm{~mol} \mathrm{dm}^{-3}$ <br> OR <br> Temperature not 298 K |  | (2) |
|  | IGNORE <br> Temperature too low <br> M2 |  |  |
|  | Reaction is (very) slow  <br> OR  <br> Activation energy is high  <br> OR  <br> Reaction is kinetically unfavourable  <br> OR  <br> Reaction mixture is kinetically stable (1) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( b ) ( i )}$ | $\mathrm{Cr}_{2} \mathrm{O}_{7}^{2-}+8 \mathrm{H}^{+}+3 \mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH} \rightleftharpoons 2 \mathrm{Cr}^{3+}+$ <br> $3 \mathrm{CH}_{3} \mathrm{CHO}+7 \mathrm{H}_{2} \mathrm{O}$ <br> OR <br> Multiples <br> Correct species throughout and norraran (1) <br> electrons <br> Fully balanced and surplus $\mathrm{H}^{+}$eliminated <br> (1) |  | (2) |
|  | IGNORE <br> State symbols even if incorrect <br> Use of $\rightarrow$ instead of $\rightleftharpoons$ <br> Fully correct reverse equation scores 1 <br> mark |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( b ) ( i i )}$ | oxidation of ethanol to ethanal <br> $\mathrm{E}_{\text {cell }}=1.33-(-0.61)=(+) 1.94$ (V) (1) |  | (2) |
|  | oxidation of ethanal to ethanoic acid <br> $\mathrm{E}_{\text {cell }}=1.33-(-0.94)=(+) 2.27$ (V) (1) |  |  |
|  | Correct values with no working scores <br> both marks | If no other mark is scored <br> $\mathrm{E}_{\text {cell }}($ to ethanal $=0.72$ (V) and <br> $\mathrm{E}_{\text {cell }}($ to ethanoic acid) $=0.39$ (V) scores <br> 1 mark |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(b) (iii) | Route 1 (Depends on $\mathrm{E}_{\text {cell }}$ (to ethanoic acid) more positive than $\mathrm{E}_{\text {cell }}$ (to ethanal)) <br> The oxidation of ethanal (to ethanoic acid) is more favourable / feasible / spontaneous than the oxidation of ethanol to ethanal <br> (because the $\mathrm{E}_{\text {cell }}$ value is more positive) <br> ALLOW <br> Ethanal is oxidised more easily than ethanol <br> Route 2 (Depends on both $\mathrm{E}_{\text {cell }}$ values being positive) <br> Both oxidations are thermodynamically favourable / feasible <br> IGNORE <br> To prevent further oxidation |  | (1) |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(c)(i) | M1 <br> Relevant reaction is $\mathrm{MnO}_{4}^{-}+8 \mathrm{H}^{+}+5 \mathrm{e}^{(-)} \rightleftharpoons \mathrm{Mn}^{2+}+4 \mathrm{H}_{2} \mathrm{O}$ <br> OR <br> Multiples <br> and $\begin{equation*} \mathrm{E}^{\ominus}=+1.51(\mathrm{~V}) \tag{1} \end{equation*}$ <br> IGNORE <br> State symbols even if incorrect <br> M2 <br> More positive $\mathrm{E}^{\ominus}$ value shows that $\mathrm{MnO}_{4}^{-}$is a stronger oxidising agent than $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}$ so oxidation might proceed further / to $\mathrm{CO}_{2}$ and $\mathrm{H}_{2} \mathrm{O}$ <br> ALLOW <br> More positive $\mathrm{E}^{\ominus}$ value shows that $\mathrm{MnO}_{4}^{-}$is a stronger oxidising agent than $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}$ so only ethanoic acid formed / ethanal not formed <br> Allow higher for more positive in both No TE on incorrect manganate(VII) equation |  | (2) |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(c)(ii) | M1 and M2 $3 \mathrm{MnO}_{4}^{2-}+2 \mathrm{H}_{2} \mathrm{O} \rightleftharpoons 2 \mathrm{MnO}_{4}^{-}+\mathrm{MnO}_{2}+4 \mathrm{OH}^{-}$ <br> ALLOW $\begin{equation*} 5 \mathrm{MnO}_{4}^{2-}+8 \mathrm{H}^{+} \rightleftharpoons 4 \mathrm{MnO}_{4}^{-}+\mathrm{Mn}^{2+}+4 \mathrm{H}_{2} \mathrm{O} \tag{1} \end{equation*}$ <br> Species and no electrons <br> Balanced <br> ALLOW <br> Multiples <br> IGNORE <br> State symbols even if incorrect <br> M3 $\mathrm{E}_{\text {cell }}=(0.59-0.56)=(+) 0.03(\mathrm{~V})$ <br> OR (for ALLOW equation) $\mathrm{E}_{\text {cell }}=(1.51-0.56)=(+) 0.95(\mathrm{~V})$ <br> and in both cases <br> (Positive so disproportionation is) feasible (1) <br> Correctly balanced equation except for missing charge(s) can score M2 and M3 <br> M3 dependent on use of correct manganese species for appropriate equations in M1 Fully correct use of reverse disproportionation to give $-0.03(\mathrm{~V}) /-0.95(\mathrm{~V})$ scores $2 / 3$ |  | (3) |



| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 1 ( b ) ( i )}$ | Penalise the omission of the positive <br> charge in (b)(i) and (b)(ii) once only <br> Penalise use of structures once only <br> ALLOW <br> Atoms in any order |  | (2) |
|  | $(\mathrm{m} / \mathrm{e}=135)$  <br> $\mathrm{C}_{8} \mathrm{H}_{9} \mathrm{NO}^{+}$  <br> $(\mathrm{m} / \mathrm{e}=77)$ (1) <br> $\mathrm{C}_{6} \mathrm{H}_{5}+$  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 21(b) (ii) | $(\mathrm{m} / \mathrm{e}=43)$  <br> ALLOW $\mathrm{CH}_{3} \mathrm{CO}^{+}$ <br> OR <br> OR <br> Bracketed structures with charges outside brackets <br> ALLOW <br> Correct charge on any part of the ion <br> COMMENT <br> Allow open bonds on C or N <br> eg |  | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 21(c) | Both bond and range required <br> (Alkane) C-H stretch at 2962-2853 ( $\mathrm{cm}^{-1}$ ) <br> IGNORE <br> Methyl group / $\mathrm{CH}_{3}$ <br> (Amide) $\mathrm{C}=\mathrm{O}$ stretch 1700 - 1630 ( $\mathrm{cm}^{-1}$ ) <br> IGNORE <br> Amide N-H stretch at 3500-3140 $\left(\mathrm{cm}^{-1}\right)$ | 1485-1365 $\begin{equation*} 1700-1680 \tag{1} \end{equation*}$ <br> Additional ranges e.g. from benzene ring | (2) |


| Question Number | Acceptable Answers |  |  |  |  | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 22(a) |  C H N <br> $\%$ <br> mass 61.0 15.3 23.7 <br> Mol $61.0 / 1$ <br> 2 $15.3 / 1$ $23.7 / 1$ <br> 4 <br> Mol $=$ 5.083 15.3 1.693 <br> Ratio 3.00 9.04 1.00 <br> (1) <br> Empirical formula $=\mathrm{C}_{3} \mathrm{H}_{9} \mathrm{~N}$ <br> (1) <br> Correct formula without working scores M3 only <br> Do not penalise the use of structural formulae or $\mathrm{C}_{3} \mathrm{H}_{7} \mathrm{NH}_{2}$ here or in part (b) |  |  |  |  |  | (3) |
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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 2 ( b )}$ | $42.7 / 24000\left(=1.779 \times 10^{-3}\right)$ mol weighs <br> $0.105(\mathrm{~g})$ <br> 1 mol weighs $0.105 \times 24000 / 42.7$ <br> $=59.016 / 59 \mathrm{~g}$ <br> Correct answer with some working scores the <br> mark <br> Formula mass of $\mathrm{C}_{3} \mathrm{H}_{9} \mathrm{~N}=59$ <br> So molecular formula (is the same as the (1) <br> empirical formula) is $\mathrm{C}_{3} \mathrm{H}_{9} \mathrm{~N}$ <br> M2 depends on M1 | (2) |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 22(c) | 1 mark each for any two structures |  | (2) |
|  | $\mathrm{CH}_{3}-\mathrm{CH}_{2}-\mathrm{CH}_{2}-\mathrm{NH}_{2}$  |  |  |
|  | ALLOW |  |  |
|  | $\mathrm{CH}_{3}-\mathrm{NH}-\mathrm{CH}_{2}-\mathrm{CH}_{3}$  |  |  |
|  | (secondary and tertiary amines even though not on spec) <br> OR <br> Fully displayed <br> OR <br> Skeletal |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 22(d) |  | (2) |  |

## Section C



| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 3 ( a ) ( i i )}$ | (When the electronic structure is derived <br> according to the aufbau rules), the <br> final/last electron added is placed in a (3)d <br> orbital / the (3)d orbitals / the (3)d <br> subshell | The outermost/ <br> highest energy <br> electron is in a (3)d <br> orbital <br> Shell for subshell | (1) |
| IGNORE <br> Cu has valence electrons in the d subshell | ${ }^{\text {Cu }}$(1) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 3 ( a ) ( i i i )}$ | $\left(\mathrm{Cu}^{2+}\right)$ has a partially filled d orbital <br> OR <br> $\left(\mathrm{Cu}^{2+}\right)$ d orbitals are/ subshell is partially <br> filled <br> OR <br> $\left(\mathrm{Cu}^{2+}\right)$ has a half- filled d orbital <br> ALLOW <br> $\left(\mathrm{Cu}^{2+}\right)$ d orbitals are/ subshell is <br> incomplete <br> $\left(\mathrm{Cu}^{2+}\right)$ d orbitals are/ subshell is not filled <br> General definition of transition metal <br> subshell | empty dal(s)/ <br> COMMENT <br> Do not penalise shell for subshell | (1) |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 23(b)(i) | M1 <br> The $\mathrm{Cu}^{2+}$ ions are coordinated / surrounded by / complexed/bonded to (water) ligands <br> OR <br> In water $\mathrm{Cu}^{2+}$ exists as $\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}{ }^{2+}$ <br> M2 <br> (3)d orbitals / (3)d subshell split (by the attached ligands into two different energy levels) <br> M3 <br> Electrons absorb energy / photons of a certain frequency (in the visible region) <br> ALLOW <br> Energy / photons / light is absorbed <br> M4 <br> Electrons are promoted <br> (from lower to higher energy d orbital(s) <br> / levels) <br> OR <br> Electrons move from lower to higher energy d orbital(s) / levels) <br> ALLOW <br> Electrons excited <br> d-d transitions occur <br> M5 <br> Reflected / transmitted / remaining light is coloured / in the visible region ALLOW <br> Complementary colour seen <br> Reflected / transmitted / remaining light / frequency is seen <br> Penalise omission of (3)d once only. Ignore reference to electrons relaxing / dropping to the ground state <br> COMMENT <br> Do not penalise shell for subshell | Just 'mention of ligand' <br> orbital <br> emitted | (5) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{2 3 ( b ) ( i i )}$ | There are no ligands coordinated around <br> / bonded to / surrounding the Cu <br>  <br> (so the d subshell is not split) |  | (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{2 3 ( c ) ( i )}$ | Buffer (solution) <br> OR <br> The components of a suitable buffer e.g. <br> a weak acid and its conjugate base <br> ALLOW <br> Addition of (measured amounts of) any <br> carbonate / hydrogencarbonate / <br> hydroxide by name or formula <br> Ammonia / $\mathrm{NH}_{3}((\mathrm{aq)}$ ) <br> compounds <br> IGNORE <br> IGNORE <br> Descriptions of the buffer <br> Addition of alkali/ $\mathrm{OH}^{-}$ | (1) |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 23(c)(ii) | Amount of edta $=0.205 \times 27.50 / 1000$ * $\begin{equation*} \left(=5.6375 \times 10^{-3} / 0.0056375\right) \tag{1} \end{equation*}$ <br> Mass of copper in sample $\begin{align*} & =\text { ans* } \times 10 \times 63.5  \tag{1}\\ & =3.5798 \mathrm{~g}^{* *} \end{align*}$ <br> Proportion of copper $\begin{align*} & =100 \times \text { ans } * * / 3.63 \\ & =98.6174 / 98.6 / 99 \% \tag{1} \end{align*}$ <br> Correct answer with no working scores 3 marks <br> ALLOW use of $\mathrm{A}_{\mathrm{r}}(\mathrm{Cu})=64(\Rightarrow 99.39$ / 99.4\%) <br> Penalise arithmetical errors once only at M3 <br> TE at each stage but do not award M3 if $\% \mathrm{Cu}$ is $\geq 100$ <br> IGNORE SF except 1 SF <br> Do not penalise correct intermediate rounding |  | (3) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{2 3 ( c ) ( i i i )}$ | Nickel ions also form a complex with <br> edta |  | (1) |
|  | ALLOW <br> Nickel also forms a complex with edta <br> Nickel (ions) react(s) with edta |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 3 ( d ) ( i )}$ | Penalise failure to mention the central <br> ion once only |  | (2) |
|  | A bidentate ligand <br> occupies two coordination positions of <br> the central cation <br> OR <br> forms two dative covalent bonds with the <br> central cation (1) | The sulfur (atom) Ione pair and the <br> nitrogen (atom) lone pair of <br> cysteamine bond to the central cation / <br> Cu² ion <br> ALLOW <br> This mark from the diagram in (d)(ii) <br> (1) |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 3 d ( i i )}$ |  | Dative bond(s) from <br> the H atoms <br> covalent bonds <br> (instead of dative <br> bonds) | (1) |
|  | ALLOW <br> Two or three cysteamine ligands <br> IGNORE <br> Charge on Cu <br> Omission of lone pairs <br> length of the carbon chain | Cu $^{2+}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| *23d(iii) | Copper(I) complexes are (usually) linear <br> (1) |  | (2) |
|  | The carbon chain in cysteamine is too <br> short to give a bond angle of $180^{\circ} /$ to <br> give linear geometry <br> OR <br> The resulting 5-membered ring would <br> have bond angles of (about) $108^{\circ}$ (rather <br> than 180) <br> ALLOW <br> Two separate cysteamine molecules (1) <br> would have to bond with copper(I) (1) <br> IGNORE <br> Just 'cysteamine acts as a monodentate <br> ligand' |  |  |

(Total for Question 23 = 20 marks)
TOTAL FOR SECTION C = 20 MARKS

