Mark Scheme (Results)

October 2018

Pearson Edexcel International
Advanced Level
In Chemistry (WCH05)
Paper 01 Transition of Metals and Organic Nitrogen

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October 2018
Publications Code WCH05_01_1810_MS*
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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 |
| :--- | :--- | :---: |
| $\mathbf{1}$ (a) | The only correct answer is A <br> B is not correct because Fe(III) acts as an oxidising <br> agent | (1) |
| C is not correct because $\mathrm{Cr}(\mathrm{II})$ loses electrons so is <br> oxidised <br> $\mathbf{D}$ is not correct because $\mathrm{Cr}(\mathrm{II})$ loses electrons so is <br> oxidised |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1}$ (b) | The only correct answer is B | (1) |
|  | A is not correct because they should both be positive |  |
|  | C is not correct because they should both be positive |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 ~ ( c )}$ | The only correct answer is D | (1) |
|  | A is not correct because carbonate ions might react <br> B is not correct because hydroxide ions might react <br> C is not correct because iodide ions might react |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 ~ ( d )}$ | The only correct answer is A | (1) |
|  | B is not correct because dividing by 10 for $100 \mathrm{~cm}^{3}$ but <br> not dividing by 2 for $\mathrm{Fe}_{2}\left(\mathrm{SO}_{4}\right)_{3}$ <br> C is not correct because dividing by 10 for $100 \mathrm{~cm}^{3}$ but <br> multiplying by 2 and not dividing by 2 <br> D is not correct because dividing by 2 for $\left.\mathrm{Fe}_{2( } \mathrm{SO}_{4}\right)_{3}$ but <br> not dividing by 10 |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{2}$ | The only correct answer is C | $(1)$ |
|  | A is not correct because Fe has an oxidation state of +6 |  |
|  | B is not correct because $C r$ has an oxidation state of +6 |  |
|  | D is not correct because $W$ has an oxidation state of +6 |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{3}$ | The only correct answer is A <br> B is not correct because it is oxidised at the negative <br> electrode <br> C is not correct because it is oxidised not reduced | (1) |
| D is not correct because it is oxidised not reduced at <br> the negative electrode |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{4 ( a )}$ | The only correct answer is C | (1) |
|  | A is not correct because this is the second step <br> B is not correct because this is the second step with an <br> incorrect product |  |
| $\mathbf{D}$ is not correct because this is the first step |  |  |$\quad$|  |
| :--- |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| 4 (b) | The only correct answer is C | (1) |
|  | A is not correct because the 3d subshell does split | B is not correct because the 3d subshell is full <br> D is not correct because there is no movement of <br> electrons in the 3d subshell |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{5}$ | The only correct answer is B <br> A is not correct because orbitals are occupied singly <br> before pairing | (1) |
| C is not correct because the 4s electrons are lost first to <br> form an ion | D is not correct because the 4s electrons are lost first <br> to form an ion |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{6}$ | The only correct answer is A |  |
| B is not correct because X ray diffraction provides no <br> evidence for this | 1 |  |
| C is not correct because X ray diffraction provides no <br> evidence for this <br> D is not correct because it is not a true statement |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{7}$ | The only correct answer is B | (1) |
| $\mathbf{A}$ is not correct because this is the reverse order <br> C is not correct because phenylamine has a lower pH <br> than ammonia <br> $\mathbf{D}$ is not correct because diethylamine has a higher pH <br> than ethylamine |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{8 ( a )}$ | The only correct answer is D | $(1)$ |
|  | A is not correct because this is not a reducing agent <br> B is not correct because this does not produce the <br> amine <br> C is not correct because this is an oxidising agent |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{8 ( b )}$ | The only correct answer is D <br> A is not correct because this is not used to separate <br> phenylamine | (1) |
| B is not correct because this is not used to separate <br> phenylamine | C is not correct because this is not used to separate <br> phenylamine |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{8 ( c )}$ | The only correct answer is C | $(1)$ |
|  | A is not correct because this has an extra amine group | B is not correct because use of phenol would leave an - <br> OH in the molecule |
| D is not correct because this is a product of the <br> reaction of 1,4-diaminobenzene with nitrous acid in <br> hydrochloric acid |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{9}$ | The only correct answer is D <br> A is not correct because this is the mass of the <br> intermediate | (1) |
| $\mathbf{B}$ is not correct because this is the overall percentage |  |  |
| $\mathbf{C}$ is not correct because this is the overall percentage |  |  |
| by mass |  |  |$\quad$


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 0}$ | The only correct answer is $\mathbf{A}$ <br> B is not correct because this is not the correct <br> momomer <br> $\mathbf{C}$ is not correct because this is not the correct <br> momomer <br> D is not correct because this is not the correct <br> momomer | (1) |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1}$ (a) | The only correct answer is A | (1) |
|  | B is not correct because 2,4-dinitrophenylhydrazine <br> does react with X | C is not correct because 2,4-dinitrophenylhydrazine <br> does react with Y |
| D is not correct because 2,4-dinitrophenylhydrazine <br> does react with Z |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1 ( b )}$ | The only correct answer is D | $(1)$ |
|  | A is not correct because W does not react with either <br> B is not correct because X reacts with acidified <br> potassium dichromate(VI) but not Tollens' reagent | C is not correct because Y does not react with either |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 1}$ (c) | The only correct answer is D | (1) |
|  | A is not correct only W does not react |  |
|  | B is not correct only W does not react |  |
| $\mathbf{C}$ is not correct only W does not react |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 2}$ | The only correct answer is B | (1) |
|  | A is not correct has a chiral carbon |  |
|  | $\mathbf{C}$ is not correct has a chiral carbon |  |
| D is not correct has a chiral carbon |  |  |

(Total for Section A = 20 marks)

## Section B

| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 13(a) |  |  | 1 |
|  | E / V |  |  |
|  |  |  |  |
|  |  |  |  |
|  |  |  |  |
|  |  |  |  |
|  |  |  |  |
|  | Fe ${ }^{3+}(\mathrm{aq})+\mathrm{e}^{(-)} \rightleftharpoons \mathrm{Fe}^{2+}(\mathrm{aq}) \mathrm{c}^{(+0.77} \mathrm{)}$ | Incorrect state |  |
|  |  | symbols |  |
|  |  |  |  |
|  | $\left(\mathrm{Cl}_{2}(\mathrm{aq})+2 \mathrm{e}^{-} \rightleftharpoons 2 \mathrm{Cl}(\mathrm{aq})\right){ }^{\text {a }}+1.36$ | 1.36 |  |
|  | ALLOW <br> Single arrow instead of reversible arrows | $\begin{aligned} & \text { without + } \\ & (+) 0.68 / \end{aligned}$ |  |
|  | Single arrow instead of reversible arrows | $(+) 2.72$ |  |
|  | IGNORE <br> Missing state symbols |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 3 ( b ) ( i )}$ | Zn / Zinc / Zn(s)/Zinc(s) | Zn <br> Zinc(II) | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 3 ( b ) ( i i )}$ | $\mathrm{SO}_{3}^{2-}(\mathrm{aq}) /$ sulfate(IV) (ions) / sulfite <br> (ions) | $\mathrm{SO}_{4}^{2-}(\mathrm{aq}) /$ <br> Sulfate(VI) <br> /sulfate | 1 |
|  | ALLOW <br> $\mathrm{SO}_{3}^{2-}(\mathrm{aq})+\mathrm{H}_{2} \mathrm{O}(\mathrm{I})$ |  |  |
|  | IGNORE <br> $\mathrm{H}^{+}$ <br> missing state symbols |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 13(c)(i) | $+2 / 2+$ <br> This can be scored if there is no answer in the space and $\mathrm{V}^{2+}$ is shown as the product in the equation, even if the equation is incorrect <br> ALLOW $\begin{align*} & \mathrm{V}^{2+} /+\mathrm{II} / \mathrm{II}+  \tag{1}\\ & \mathrm{VO}_{2}^{+}+4 \mathrm{H}^{+}+3 \mathrm{e}^{-} \rightarrow \mathrm{V}^{2+}+2 \mathrm{H}_{2} \mathrm{O} \end{align*}$ <br> Must be a half-equation not a full equation with zinc <br> ALLOW <br> Multiples <br> IGNORE <br> State symbols even if incorrect <br> No TE for equations on incorrect values of | $\begin{aligned} & 2 \text { / II / } \\ & \text { V(II) } \end{aligned}$ | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 13(c)(ii) | M1 (Recognition of oxidation by air) <br> (Vanadium(II) / V ${ }^{2+}$ / Vanadium(III) / $\mathrm{V}^{3+} /$ solution) is oxidised by / reacts with oxygen (in the air) <br> This can be scored if an equation is given showing reaction of $\mathrm{V}^{2+}$ or $\mathrm{V}^{3+}$ with $\mathrm{O}_{2}$ <br> M2 (Formation of V(III) from V(II)) $4 \mathrm{~V}^{2+}+\mathrm{O}_{2}+4 \mathrm{H}^{+} \rightarrow 4 \mathrm{~V}^{3+}+2 \mathrm{H}_{2} \mathrm{O}$ <br> ALLOW $\begin{equation*} \mathrm{V}^{2+} \rightarrow \mathrm{V}^{3+}+\mathrm{e}^{-} \quad / \mathrm{V}^{3+}+\mathrm{e}^{-} \rightleftharpoons \mathrm{V}^{2+} \tag{1} \end{equation*}$ <br> and <br> $\mathrm{V}^{2+}$ becomes $\mathrm{V}^{3+}$ which is green <br> M3 (Formation of V(IV)) $\begin{equation*} 4 \mathrm{~V}^{3+}+\mathrm{O}_{2}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow 4 \mathrm{VO}^{2+}+4 \mathrm{H}^{+} \tag{1} \end{equation*}$ <br> IGNORE <br> State symbols even if incorrect <br> ALLOW <br> $2 \mathrm{~V}^{2+}+\mathrm{O}_{2} \rightarrow 2 \mathrm{VO}^{2+}$ Scores 1 (of M2 and M3) <br> M4 (Calculation of $\mathrm{E}_{\text {cell }}$ values) <br> $\mathrm{E}_{\text {cell }}$ for M 2 equation $=(+) 1.49(\mathrm{~V})$ <br> and <br> $\mathrm{E}_{\text {cell }}$ for M 3 equation $=(+) 0.89(\mathrm{~V})$ | V(III) from $\mathrm{VO}^{2+}$ | 4 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 13(c)(iii) | $\mathrm{E}_{\text {cell }}$ is $(+) 0.23(\mathrm{~V})$ so the oxidation of $\mathrm{VO}^{2+}$ to $\mathrm{VO}_{2}^{+}$is feasible <br> ALLOW <br> $\mathrm{E}_{\text {cell }}$ is (slightly) positive <br> EITHER <br> the activation energy is too large / kinetically inert <br> OR <br> concentration of oxygen is too low <br> IGNORE <br> Non-standard conditions <br> Mark each as stand alone | Rate is slow | 2 |



| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 4 ( b ) ( i )}$ | $\mathrm{C}_{10} \mathrm{H}_{12} \mathrm{O}$ | $\mathrm{C}_{10} \mathrm{H}_{11} \mathrm{OH}$ | 1 |
| Ignore names e.g. Anethole |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 4 ( b ) ( i i )}$ | Restricted rotation around a carbon- <br> carbon double bond |  | 2 |
|  | ALLOW <br> No rotation around a carbon-carbon (1) <br> double bond | Two different groups attached to each <br> carbon (1) | Mark independently |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 4 ( c ) ( i )}$ | Propanoyl chloride <br> Ignore formulae as working <br> ALLOW <br> 1-propanoyl chloride | Propanyl <br> chloride | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 14(c)(ii) | First mark |  | 4 |
|  | $\mathrm{RCOCl}+\mathrm{AlCl}_{3} \longrightarrow \mathrm{RCO}^{+}+\mathrm{AlCl}_{4}^{-}$ |  |  |
|  | OR |  |  |
|  | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COCl}+\mathrm{AlCl}_{3} \longrightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CO}^{+}+\mathrm{AlCl}_{4}^{-}$ |  |  |
|  | ALLOW any acyl chloride or halogenoalkane from (c)(i) |  |  |
|  |  |  |  |
|  | Second mark <br> Curly arrow from on or within the circle towards the C of $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CO}^{+} / \mathrm{RCO}^{+}$ <br> ALLOW curly arrow from anywhere within the hexagon ALLOW curly arrow to any part of the $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CO}^{+}$ including to the + charge | Curly arrow on or outside the hexagon |  |
|  | ALLOW TE for any R group attached to $\mathrm{CO}^{+}$ <br> Third mark | hexagon |  |
|  | Intermediate structure including charge with horseshoe covering at least 3 carbon atoms <br> and <br> facing the tetrahedral carbon <br> and <br> some part of the positive charge must be within the horseshoe <br> ALLOW dotted horseshoe | Dotted bonds to H and RCO unless part of a 3-D shape |  |
|  |  |  |  |
|  | Incorrect orientation of product at this marking point Fourth mark |  |  |
|  | Curly arrow from $\mathrm{C}-\mathrm{H}$ bond to anywhere in the hexagon, reforming the correct delocalised structure (and $\mathrm{H}^{+}$) | Curly arrow from H |  |
|  | IGNORE any involvement of $\mathrm{AlCl}_{4}^{-}$in the final step <br> Correct Kekulé / skeletal structures score full marks |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 4 ( c ) ( \text { iii) }}$ Lone pair (of electrons) on the oxygen <br> (of the methoxy- group) <br> AND <br> Overlaps with the n / delocalised <br> electrons in the benzene ring / <br> delocalised system <br> OR <br> Feeds into / donates to / interacts with <br> the delocalised electrons / delocalised <br> system / n system of the benzene ring <br> ALLOW <br> Increases the electron density of the (1) <br> benzene ring <br> Making it more susceptible to <br> electrophilic attack / attack by propanoyl <br> cation / RCO 2 <br> ALLOW <br> Making it a better nucleophile <br> Mark each point independently (1) |  |  |  |



| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| $\mathbf{1 4 ( e ) ( i )}$ |  | Any other <br> hydrogens <br> labelled | 1 |
|  | ALLOW <br> Any other unambiguous identification of the <br> three hydrogens |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |  |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\mathbf{1 4 ( e ) ( i i )}$ |  |  |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 14(f) | (Reduction using) Lithium tetrahydridoaluminate((III)) / Lithium aluminium hydride / Lithal / $\mathrm{LiAlH}_{4}$ in (dry) ether <br> ALLOW <br> Sodium tetrahydridoborate((III)) / Sodium borohydride / $\mathrm{NaBH}_{4}$ <br> IGNORE <br> Heat / reflux / distillation <br> ALLOW <br> Skeletal formula <br> (Substitution using) $\mathrm{PCl}_{5}$ <br> OR <br> $\mathrm{NaCl} / \mathrm{KCl}$ and concentrated / conc. $\mathrm{H}_{2} \mathrm{SO}_{4}$ <br> ALLOW <br> $\mathrm{PCl}_{3} / \mathrm{SOCl}_{2} /$ concentrated hydrochloric acid <br> (Substitution using) $\mathrm{PBr}_{3} / \mathrm{P}$ and $\mathrm{Br}_{2}$ (giving bromoalkane) <br> (Substitution using) $\mathrm{PI}_{3} /$ (red) P and $\mathrm{I}_{2}$ (giving iodoalkane) | Hydrogen and nickel | 5 |


|  <br> Or bromo- or iodo- compounds as appropriate <br> (Elimination using) ethanolic / alcoholic <br> / EtOH / alc. sodium/potassium hydroxide <br> AND <br> Heat / boil / heat under reflux <br> Marking consequential on correct intermediates but <br> ALLOW for max 3 a two step synthesis using step 1 as above and then <br> Conc. $\mathrm{H}_{2} \mathrm{SO}_{4} / \mathrm{H}_{3} \mathrm{PO}_{4} / \mathrm{Al}_{2} \mathrm{O}_{3}$ <br> AND <br> Heat / boil / heat under reflux / $170^{\circ} \mathrm{C}$ giving anethole |  |  |
| :---: | :---: | :---: |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 15(a)(i) | Moles of thiosulfate |  | 2 |
|  | $=21.60 \times 10^{-3} \times 3 \times 10^{-3}$ |  |  |
| $=6.48 \times 10^{-5} / 0.0000648$ (mol) (1) |  |  |  |
| Moles of $\mathrm{Cu}^{2+}$ in $100 \mathrm{~cm}^{3}=$ moles of |  |  |  |
| thiosulfate $\times 10$ |  |  |  |
| $=6.48 \times 10^{-4} / 0.000648$ (mol) (1) |  |  |  |
| If M1 is scored, then there is no |  |  |  |
| further attempt, the second mark |  |  |  |
| can be scored in (a)(ii) |  |  |  |
| Ignore SF except 1 SF |  |  |  |
| Correct answer with no working |  |  |  |
| scores 2 |  |  |  |$\quad$| ( |
| :--- |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 5 ( a ) ( i i )}$ | Mass of $\mathrm{Cu}=6.48 \times 10^{-4} \times 63.5$ |  | 1 |
|  | $=0.041148 / 4.1148 \times 10^{-2}(\mathrm{~g})$ |  |  |
|  | $=0.041 / 4.1 \times 10^{-2}(\mathrm{~g})$ |  |  |
|  | Answer must be to 2 SF |  |  |


| Question | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 15(b) | First two marks are stand alone |  | 5 |
|  | In 0.500 g |  |  |
|  | Mass of water $=0.07(0)$ |  |  |
|  | AND |  |  |
|  | $\begin{align*} & \text { Moles of water }=\frac{0.07(0)}{18} \\ & =0.0038889(\mathrm{~mol}) / 3.8889 \times 10^{-3}(\mathrm{~mol}) \tag{1} \end{align*}$ |  |  |
|  | $\begin{align*} & n=\frac{\text { moles of water }}{\text { moles of } \mathrm{Cu}^{2+}} \\ & =\frac{0.0038889}{6.48 \times 10^{-4}}=6(.0014) \tag{1} \end{align*}$ |  |  |
|  | Method 1 <br> moles of sulfate $=2 \times$ moles of $\mathrm{Cu}^{2+}$ $=0.001296 / 1.296 \times 10^{-3}(\mathrm{~mol})$ |  |  |
|  | Mass of sulfate $=$ moles of sulfate $\times 96.1$ $\begin{equation*} =0.12455(\mathrm{~g}) \tag{1} \end{equation*}$ |  |  |
|  | $\begin{align*} & \text { Mass of } M=0.500-\text { mass of copper - mass } \\ & \text { of sulfate - mass of water } \\ & \qquad=0.500-0.12455-0.070- \\ & =0.26430(\mathrm{~g}) \end{align*}$ |  |  |
|  | $\begin{aligned} & \text { Atomic mass of } M=\frac{\text { mass of } M}{\text { moles of } M} \\ & =2 \times \frac{0.264}{6.48 \times 10^{-4}}=203.94 \end{aligned}$ |  |  |
|  | So compound is $\mathrm{Tl}_{2} \mathrm{Cu}\left(\mathrm{SO}_{4}\right)_{2} \cdot 6 \mathrm{H}_{2} \mathrm{O}$ |  |  |
|  | Method 2 |  |  |


(Total for Question 15 = 8 marks)

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i )}$ | (2-)aminobutan(e)dioic acid | Answers with <br> dibutan(e) in <br> the name e.g. <br> aminodibutanoic <br> acid | 1 |
|  | (2-)aminobutan(e)-1,4-dioic acid <br> (2-)aminebutan(e)dioic acid <br> (2-)aminebutan(e)-1,4-dioic acid | IGNORE <br> Punctuation marks (e.g. hyphens, <br> commas, full stops etc) in either <br> version of the answer so for example <br> 2 aminobutandioic acid would score. |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i i )}$ |  |  | Neutral structure <br> Single negative <br> ion |
|  | ALLOW <br> Structural formulae / displayed <br> formulae e.g. <br> $-\mathrm{OOCCH}_{2} \mathrm{CHNH}_{2} \mathrm{COO}^{-}$ | 1 |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(a)(iii) | Diagram <br> ALLOW <br> $\mathrm{NH}_{3}{ }^{+}$ <br> Phenylalanine molecules are held to each other by ionic bonds / strong electrostatic attractions between oppositely charged ions (so high melting temperature) / held in (giant) ionic lattice <br> IGNORE <br> Strong electrostatic attractions between molecules without mention of ionic bonds or between oppositely charged ions <br> Reference to hydrogen bonds | Internal ionic bonds specifically mentioned but assume intermolecular if not specific. | 2 |
| Question Number | Acceptable Answers | Reject | Mark |
| 16(b)(i) | Methanol/ $\mathrm{CH}_{3} \mathrm{OH}$ |  | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(b)(ii) | Put spots of the amino acid mixture / hydrolysis products (and known amino acids) <br> AND <br> on a tlc plate / filter paper / <br> chromatography paper <br> AND <br> in a (suitable) solvent / run with a (suitable) solvent <br> ALLOW <br> Labelled diagram <br> Use ninhydrin (to make amino acids visible) <br> ALLOW <br> Iodine vapour in place of ninhydrin <br> Compare distance travelled of mixture components with known amino acids <br> OR <br> Compare $R_{f}$ / formula of $R_{f}$ / description of $R_{f}$ with data book values (1) | Just already separated amino acids <br> Amino acids dissolved in mobile phase solvent <br> just 'paper' <br> Ni as an abbreviation <br> Just 'compare with data book values' <br> Just 'Calculate $\mathrm{R}_{\mathrm{f}}$ values' | 3 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( b ) ( i i i )}$ | Heat causes hydrolysis <br> OR <br> Amino acids are not sweet <br> ALLOW <br> Decomposition / breakdown / unstable on <br> heating | Just 'cooking' <br> without 'heat' | 1 |
| IGNORE <br> Methanol is toxic <br> Changes to flavour without mention of <br> sweetness | Decomposition / <br> breakdown / <br> unstable without <br> heat |  |  |

(Total for Question 16 = 10 marks)
(Total for Section $B=50$ marks)

## Section C

| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(a) | Two calculations which must be accompanied by a correct statement about toxicity. <br> Several approaches are possible. <br> e.g. <br> Mass of CO released $=0.35 \times 28=9.8(\mathrm{~g})$ <br> Mass per $\mathrm{m}^{3}=\frac{9.8}{200}=0.049(\mathrm{~g}) / 49(\mathrm{mg})$ <br> Which is greater than the toxicity limit <br> OR $\begin{equation*} \text { Max. mass }=43.2 \times 200=8640(\mathrm{mg}) \tag{1} \end{equation*}$ <br> Maximum moles allowed $=\frac{8640}{1000 \times 28}$ $=0.30857143$ <br> Which is less than was released (so not within the limits) <br> OR <br> Max. moles per $\mathrm{m}^{3}=\frac{43.2 \times 10^{-3}}{28}$ $\begin{equation*} =0.0015429 / 1.5429 \times 10^{-3}(\mathrm{~mol}) / \tag{1} \end{equation*}$ <br> 1.5429 (mmol) <br> Moles per $\mathrm{m}^{3}$ released $=\frac{0.35}{200}=0.00175(\mathrm{~mol})$ <br> Which is more that the toxicity limit <br> OR <br> Moles per $\mathrm{m}^{3}$ released $=\frac{0.35}{200}=0.00175(\mathrm{~mol})$ <br> Mass per $\mathrm{m}^{3}$ released $=0.00175 \times 28$ $\begin{equation*} =0.049(\mathrm{~g}) / 4.9 \times 10^{-2}(\mathrm{~g}) / 49(\mathrm{mg}) \tag{1} \end{equation*}$ <br> Which is more than the toxicity limit <br> ALLOW TE only on a suitable attempt at a calculation of mass or moles in M1 <br> Other approaches may be possible |  | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(b)(i) | If the name of a shape is given award M1 for the name if correct. Do not negate with and incorrect diagram. <br> Trigonal bipyramid(al) <br> ALLOW <br> Pyramidal / bipyramidal if a correct diagram is given <br> ALLOW <br> If no name is given, a three dimensional diagram showing three bonds in plane (straight lines) and two bonds out of plane, either a wedge and dots (which may also be wedged, but ignore the direction of this wedge) or two oppositely directed wedges (one fat at Fe and another fat at CO) <br> One angle labelled $120^{\circ}$ and one angle labelled $90^{\circ}$, which may be shown as the symbol ' $r$ '. <br> ALLOW <br> If no other mark has been scored, a diagram with no dots and wedges which has at least one correct $90^{\circ}$ and one correct $120^{\circ}$ angle scores <br> IGNORE <br> Point of attachment of CO to Fe | Square based pyramids <br> Just <br> 'pyramidal' or <br> 'bipyramidal' <br> Any additional angles which are labelled incorrectly but not the correct $180^{\circ}$ angle | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(b)(ii) | Dative covalent bond from C to Fe <br> AND <br> Ione pair on O <br> Triple bond between C and O with one dative covalent bond. <br> ALLOW <br> Crosses for carbon and dots for oxygen <br> Dative covalent bond to Fe , double bond between C and O and two lone pairs on O scores (1) <br> IGNORE <br> Circles for electron shells / lines as well as dots and crosses to show bonds / lone pairs on the Fe |  | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(c)(i) | Moles of CO $=\underline{4.8}=0.2$ moles |  | 3 |
|  | Mass of CO $=0.2 \times 28=5.6 \mathrm{~g}$ |  |  |
|  | AND |  |  |
|  | Mass of $\mathrm{Mn}=7.8-5.6 \mathrm{~g}=2.2$ |  |  |
|  | Moles of $\mathrm{Mn}=\frac{2.2}{54.9}=0.04007286$ |  |  |
|  | Ratio is $0.04: 0.2$ <br> 1 : $5 \quad\left(\mathrm{SoMn}(\mathrm{CO})_{5}\right)$ |  |  |
|  | If Ar Mg used instead (24 / 24.3) final answer of $0.091667: 0.2 / 0.090535: 0.2$ |  |  |
|  | or 1:2 can score M1 and M2. |  |  |
|  | ALLOW |  |  |
|  | TE for incorrect mass of manganese in M2 for ratio M3. |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(c)(ii) | Empirical formula mass $=194.9$ $390=2 \times 194.9$ <br> so molecular formula $=\mathrm{Mn}_{2}(\mathrm{CO})_{10}$ <br> ALLOW <br> Just $\mathrm{Mn}_{2}(\mathrm{CO})_{10}$ without working or a structure with 2 Mn and 10 CO <br> ALLOW <br> Any sensible structure of two Mn and ten CO covalently bonded at any angle to each other <br> IGNORE <br> Connectivity of the CO group |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( d ) ( i )}$ | Sulfuric acid is a catalyst <br> IGNORE <br> Sulfuric acid / H |  |  |
|  | Hydrogen is an electrophile (reacts with 2- <br> methylpropene and) is regenerated at the <br> end / in the last step of the reaction / <br> takes part in the reaction but is still <br> present at the end | Just 'sulfuric acid is <br> chemically <br> unchanged after <br> the reaction' <br> 'Not participating in <br> the overall reaction' | 2 |
| M2 dependent on M1. | (1) |  |  |



| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( d ) ( i i i )}$ | Sulfuric acid is corrosive | Just cost | 1 |
|  | OR |  |  |
|  | Difficult to recover the sulfuric acid <br> IGNORE <br> Irritant <br> Burns skin / toxic / discussion of yield |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( e )}$ | The (two) reactants are / carbon monoxide is <br> adsorbed onto the surface / active sites of the (1) <br> catalyst <br> The activation energy for the reaction is lowered <br> / bonds are weakened in the reactant molecules <br> $(1)$ | absorbed | 3 |
|  | The products are desorbed from / diffuse from / <br> leave the catalyst |  |  |

(Total for Question 17 = 20 marks) (Total for Section $C=20$ marks) Total for Paper $=90$ marks

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