## P Pearson Edexcel

## Mark Scheme (Results)

January 2019

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

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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 D <br> $\boldsymbol{A}$ is not correct because not all d block metals are transition <br> elements. <br> $\boldsymbol{B}$ is not correct because the definition should refer to <br> incompletely filled d orbitals. <br> $\boldsymbol{C}$ is not correct because it must refer to ions, not just the atoms of <br> the element. | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{2}$ | The only correct answer is C |  |
| $\boldsymbol{A}$ is not correct because this sequence is typical of a Group 1 |  |  |
| element. |  |  |
| $\boldsymbol{B}$ is not correct because this sequence is typical of a Group 3 |  |  |
| element. |  |  |
| $\boldsymbol{D}$ is not correct because this sequence is typical of a Group 2 <br> element. | 1 |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{3}$ | The only correct answer is B <br> $\boldsymbol{A}$ is not correct because the oxidation numbers in columns 1 and <br> 2 are incorrect. <br> $\boldsymbol{C}$ is not correct because the oxidation number in column 2 is <br> incorrect. <br> $\boldsymbol{D}$ is not correct because the oxidation number in column 1 is <br> incorrect. | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{4}$ | The only correct answer is B | 1 |
| A is not correct because 2 nitrate ions have a total drop in <br> oxidation number of +6 so each M must increase by 2. <br> C is not correct because 2 nitrate ions have a total drop in <br> oxidation number of +6 so each M must increase by 2. | D is not correct because 2 nitrate ions have a total drop in <br> oxidation number of +6 so each M must increase by 2. |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{5 ( a )}$ | The only correct answer is C | 1 |
|  | $\boldsymbol{A}$ is not correct because $\mathrm{V}^{3+}$ is in the least positive half-cell. |  |
| $\boldsymbol{B}$ is not correct because $\mathrm{V}^{2+}$ is a reducing agent. |  |  |
|  | $\boldsymbol{D}$ is not correct because $\mathrm{Cl}^{\prime}$ is a reducing agent. |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 5(b) | The only correct answer is A |  |
| B is not correct because I- would reduce V(V) to V(IV). |  |  |
| C is not correct because $\mathrm{Cl}_{2}$ would oxidise V(IV) to V(V). |  |  |
| D is not correct because Cl' is not strong enough to reduce any <br> species in the table. | 1 |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{6}$ | The only correct answer is A |  |
| $\boldsymbol{B}$ is not correct because $E^{\theta}$ is proportional to In $K$ | 1 |  |
|  | C is not correct because $E^{\theta}$ is proportional to $\Delta S_{\text {total. }}$ |  |
|  | $\boldsymbol{D}$ is not correct because $E^{\theta}$ is proportional to $\Delta S_{\text {total. }}$ |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{7}$ | The only correct answer is D <br> $\boldsymbol{A}$ is not correct because CI has oxidation numbers above and <br> below 0 <br> $\boldsymbol{B}$ is not correct because Br has oxidation numbers above and <br> below +1 <br> $\boldsymbol{C}$ is not correct because S has oxidation numbers above and <br> below +4 | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{8}$ | The only correct answer is D <br> $\boldsymbol{A}$ is not correct because there are not different arrangements of <br> the ligands in space. <br> $\boldsymbol{B}$ is not correct because there are not different arrangements of <br> the ligands in space. <br> $\boldsymbol{C}$ is not correct because there are not different arrangements of <br> the ligands in space. | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{9}$ | The only correct answer is B | 1 |
|  | is not correct because it is not oxidised in the reaction. <br> $\boldsymbol{C}$ is not correct because products cannot be separated if they are <br> not desorbed. <br> $\boldsymbol{D}$ is not correct because metals do not form hydrogen bonds |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 0}$ | The only correct answer is B | 1 |
|  | A is not correct because none of the functional groups is ionised. <br> C is not correct because protonation of $\mathrm{NH}_{2}$ would not occur at <br> pH 12. <br> D is not correct because protonation of $\mathrm{NH}_{2}$ would not occur at <br> pH 12. |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 1}$ | The only correct answer is C <br> $\boldsymbol{A}$ is not correct because this is not a good method for separating <br> solids. <br> $\boldsymbol{B}$ is not correct because amino acids are not volatile. <br> $\boldsymbol{D}$ is not correct because a small scale method is more suitable for <br> identification purposes. | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 2}$ | The only correct answer is B | 1 |
|  | A is not correct because C2 is connected to 4 different groups. |  |
|  | C is not correct because C3 is connected to 4 different groups. |  |
|  | D is not correct because C2 is connected to 4 different groups. |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 3}$ | The only correct answer is A <br> $\boldsymbol{B}$ is not correct because the hydrogen environments on $\mathrm{CH}_{2}$ and <br> $\mathrm{CH}_{3}$ are equivalent. <br> $\boldsymbol{C}$ is not correct because the only hydrogen environments are on <br> $\mathrm{CH}_{2}$ and $\mathrm{CH}_{3}$; this answer is the number of C atoms. <br> $\boldsymbol{D}$ is not correct because the only hydrogen environments are on <br> $\mathrm{CH}_{2}$ and $\mathrm{CH}_{3}$; this answer is the number of H atoms. | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 4}$ | The only correct answer is C <br> $\boldsymbol{A}$ is not correct because this is the number of protons on each <br> atom. <br> $\boldsymbol{B}$ is not correct because it is the number of protons on the <br> carbons in the ethyl group and a singlet for the first methyl group. | 1 |
| $\boldsymbol{D}$ is not correct because there is a quartet for the $\mathrm{CH}_{2}$ but only <br> one triplet for a methyl group.The other methyl gives a singlet. |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 5}$ | The only correct answer is B <br> $\boldsymbol{A}$ is not correct because methanol not hydrogen, is the fuel in the <br> cell. <br> C is not correct because the conditions are alkaline, not acidic | 1 |
| $\boldsymbol{D}$ is not correct because this is an oxidation; it should be a <br> reduction reaction. |  |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 6}$ | The only correct answer is D | 1 |
|  | A is not correct because this compound is an acid, not an ester. |  |
| $\mathbf{B}$ is not correct because this compound is not a benzoate. |  |  |
|  | C is not correct because this compound is not a benzoate. |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 7}$ | The only correct answer is C | 1 |
|  | A is not correct because alcohols do not react with chloroalkanes. <br> $\boldsymbol{B}$ is not correct because an addition copolymer would form. <br> $\boldsymbol{D}$ is not correct because carboxylic acids do not react with <br> amides. |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 8 ( a )}$ | The only correct answer is A | 1 |
|  | B is not correct because steam distilling is needed. |  |
|  | C is not correct because steam distilling is needed. |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 8 ( b )}$ | The only correct answer is A | 1 |
| B is not correct because the C=C will decolourise acidified <br> potassium manganate(VII). <br> C is not correct because phosphorus(V) chloride does not react <br> with C=C or the aldehyde group. | $\boldsymbol{D}$ is not correct because the CHO will form silver with Tollens' <br> solution. |  |

## Section B

| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(a)(i) | $X$ : Platinum / $\mathrm{Pt}_{((\mathrm{s})}$ and <br> Y : Platinum / $\mathrm{Pt}_{((\mathrm{s}))}$ |  | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(a)(ii) | M1: |  | 2 |
|  | Iron(II) sulfate should be $2 \mathrm{~mol} \mathrm{dm}^{-3}$ |  |  |
|  | OR |  |  |
|  | Iron(III) sulfate should be replaced with |  |  |
|  | $2 \mathrm{~mol} \mathrm{dm}^{-3}$ iron(III) chloride |  |  |
|  | and |  |  |
|  | Iron(II) sulfate should be $2 \mathrm{~mol} \mathrm{dm}^{-3}$ |  |  |
|  | ALLOW |  |  |
|  | Any method that produces a equimolar |  |  |
|  | mixture of iron(II) and |  |  |
|  | iron(III) ions |  |  |
|  | eg 1 volume iron(III) sulfate +2 volumes |  |  |
|  | iron(II) sulfate of same concentration |  |  |
|  |  |  |  |
|  | M2 |  |  |
|  | Mixture should be $1 \mathrm{~mol} \mathrm{dm}^{-3}$ with respect to each iron ion (in standard electrode) |  |  |
|  | ALLOW |  |  |
|  | The mixture is equimolar with respect to each iron ion |  |  |
|  | Calculation showing concentrations are |  |  |
|  | equimolar in mixture |  |  |
|  | (1) |  |  |
|  | M2 is independent of $M 1$ |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(a)(iii) | Potassium manganate((VII))/ <br> potassium permanganate/ $\mathrm{KMnO}_{4}$ <br> (1) <br> Manganese(II) sulfate/ $\mathrm{MnSO}_{4}$ (and (dilute) sulfuric acid / $\mathrm{H}_{2} \mathrm{SO}_{4}$ ) <br> ALLOW <br> Manganese(II) nitrate/ $\mathrm{Mn}\left(\mathrm{NO}_{3}\right)_{2}$ <br> Manganese(II) chloride/ $\mathrm{MnCl}_{2}$ <br> (1) <br> IGNORE <br> $\mathrm{MnO}_{4}{ }^{-}, \mathrm{H}^{+}, \mathrm{Mn}^{2+}, \mathrm{H}_{2} \mathrm{O}$, "acidified" Dilute hydrochloric acid/ HCl | Incorrect oxidation <br> number eg <br> Potassium <br> manganate(VI)/ <br> Concentrated <br> sulfuric acid <br> Concentrated <br> hydrochloric acid <br> $\mathrm{MnO}, \mathrm{Mn}(\mathrm{OH})_{2}$ | 2 |

\(\left.$$
\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\
\text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\
\hline \text { 19(a)(iv) } & \begin{array}{l}\text { White and precipitate / ppt(e) / solid } \\
\text { (1) }\end{array} & \begin{array}{l}\text { Just "an insoluble } \\
\text { salt forms" }\end{array}
$$ \& 2 <br>
If reference made <br>

to bubbles\end{array}\right\}\)| $\mathrm{Ba}^{2+(\mathrm{aq})+\mathrm{SO}_{4}{ }^{2-}(\mathrm{aq}) \rightarrow \mathrm{BaSO}_{4}(\mathrm{~s})}$Balanced equation with (1) <br> state symbols <br> $\mathrm{M1}$ and M2 to be marked independently |
| :--- |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9 ( a ) ( v )}$ | Potassium nitrate/ $\mathrm{KNO}_{3} /$ Sodium nitrate / <br> $\mathrm{NaNO}_{3}$ | Iodides <br> Group II salts | 1 |
| ALLOW | Sodium chloride/ $\mathrm{NaCl} /$ potassium chloride / <br> $\mathrm{KCl} /$ potassium sulfate $/ \mathrm{K}_{2} \mathrm{SO}_{4} /$ sodium <br> sulfate/ $\mathrm{Na}_{2} \mathrm{SO}_{4}$ <br> If name and formulae given both must be <br> correct. |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9 ( a ) ( \mathbf { v i ) }}$ | $((+1.51)-(+0.77))=(+) \mathbf{0 . 7 4}(\mathrm{V})$ | -0.74 | 1 |
|  | ALLOW |  |  |
| .74 |  |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(b)(i) | M1 <br> $E^{\ominus}$ for item $36=+0.17(\mathrm{~V})$ <br> (and $\mathrm{Fe}^{3+} \mid \mathrm{Fe}^{2+}=+0.77(\mathrm{~V})$ ) <br> OR $\left.E_{\text {cell }}=+0.60(\mathrm{~V})\right)$ <br> M2 <br> $E_{\text {cell }}$ is positive (so the reaction is feasible/ spontaneous) <br> This depends on some data having been used to do a calculation or comparison, even if item $45(0.4 \mathrm{~V})$ or $48(+0.51 \mathrm{~V})$ has been used. <br> ALLOW <br> TE on incorrect positive value in M1 <br> The $\mathrm{SO}_{2}$ half cell is less positive than the $\mathrm{Fe}^{3+} \mid \mathrm{Fe}^{2+}$ half cell / $\mathrm{SO}_{2}$ is a more powerful reducing agent than $\mathrm{Fe}^{2+}$ (so it will work) | $+0.40(\mathrm{~V})\left(E^{\ominus}\right. \text { for }$ reduction of $\mathrm{H}_{2} \mathrm{SO}_{3}$ ) $\begin{equation*} +0.51(\mathrm{~V}) \tag{1} \end{equation*}$ | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19(b)(ii) | M1 <br> $\left.\mathrm{Mol} \mathrm{MnO}{ }_{4}^{-}=((24.50)(0.0250) / 1000)\right)$ $\begin{equation*} =6.125 \times 10^{-4} / 0.0006125 \tag{1} \end{equation*}$ <br> M2 <br> Mol Fe ${ }^{2+}$ in $25 \mathrm{~cm}^{3}=\left(6.125 \times 10^{-4} \times 5\right)$ $\begin{equation*} =3.0625 \times 10^{-3} / 0.0030625 \tag{1} \end{equation*}$ <br> M3 <br> Mol $\mathrm{Fe}_{2} \mathrm{O}_{3}$ used to make $250 \mathrm{~cm}^{3}$ solution $\begin{align*} & =\left(\left(3.0625 \times 10^{-3} \times 10\right) / 2\right) \\ & =1.53125 \times 10^{-2} / 0.0153125 \tag{1} \end{align*}$ <br> M4 <br> Mass $\mathrm{Fe}_{2} \mathrm{O}_{3}=$ $\left(159.6 \times 1.53125 \times 10^{-2}\right)$ $=2.443875 \mathrm{~g}$ <br> and $\% \mathrm{Fe}_{2} \mathrm{O}_{3}=((2.443875 / 3.00) \times 100)$ = 81.4625\% / 81.46 \% / 81.5\% <br> ALLOW <br> TE at each stage Ignore SF except 1 SF <br> 81.67 if $\mathrm{Fe}=56$ is used. |  | 4 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 19(b)(iii) | $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}+14 \mathrm{H}^{+}+6 \mathrm{e}^{-} \rightarrow 2 \mathrm{Cr}^{3+}+7 \mathrm{H}_{2} \mathrm{O}$ <br> and <br> $\mathrm{Fe}^{2+} \rightarrow \mathrm{Fe}^{3+}+\mathrm{e}^{-} \quad$ (1) <br> $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}+14 \mathrm{H}^{+}+6 \mathrm{Fe}^{2+} \rightarrow$ <br> $2 \mathrm{Cr}^{3+}+7 \mathrm{H}_{2} \mathrm{O}+6 \mathrm{Fe}^{3+}$ |  | 2 |
|  | ALLOW (1) <br> Multiples for any of the equations <br> Correct final equation scores (2) <br> lgnore state symbols even if incorrect. |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 19(b)(iv) | The colour change at the end point <br> with manganate(VII) is clearer / more <br> distinct / more obvious <br> OR <br> With dichromate(VI) the end point <br> would not be a clear change / <br> would be from greenish yellow to <br> yellowish green | Reaction occurs <br> more readily | 1 |
| ALLOW <br> MnO $4^{-}$does not need an indicator/ is <br> self indicating <br> Any reasonable colours | IGNORE <br> Potassium dichromate is toxic/ <br> Is more expensive/ <br> Is a better oxidising agent/ <br> Has a higher $E^{\ominus}$ value |  |  |

(Total for Question 19 = 18 marks)

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( a ) ( i )}$ | 2Cu <br> OR $+4 I^{-} \rightarrow 2 \mathrm{CuI}+\mathrm{I}_{2}$ <br> Multiples <br> IGNORE <br> State symbols even if incorrect | $\mathrm{Cu}_{2} \mathrm{I}_{2}$ | 1 |

$\left.\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline \text { 20(a)(ii) } & \left(1 s^{2}\right) 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{6} 3 d^{10}\left(4 s^{0}\right) & {[\text { Ar }] 3 d^{10}} & 1 \\ & \text { ALLOW } & & \\ p_{x}^{2} p_{y}{ }^{2} p_{z}^{2} \text { in } 2 p \text { and } 3 p \\ \left(1 s^{2}\right) 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{6} 3 d^{10}\left(4 s^{0}\right) \\ \left(1 s^{2}\right) 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{6} 4 s^{0} 3 d^{10}\end{array}\right]$

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 20a(iii) | $\mathrm{Zn}^{2+} / \mathrm{Ga}^{3+}$ | $\mathrm{As}^{5+}$ <br> $\mathrm{Se}^{6+}$ <br> $\mathrm{Br}^{7+}$ <br> $\mathrm{Ga}^{4+}$ | 1 |
|  | ALLOW | $\mathrm{Ge}^{4+}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 20(b)(i) | The (3)d orbitals split / (3)d sub shell <br> splits (into two groups). <br> ALLOW <br> (3)d energy level splits <br> Can be shown on a diagram | The orbital splits |  |$\quad$|  |
| :--- |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| *20(b)(ii) | M1 <br> The gap between groups of energy <br> levels is different with different <br> ligands/ <br> The 3d orbitals split to different <br> extents with different ligands <br> (1) |  | 3 |
|  | M2 <br> Electrons absorb/ gain energy of <br> specific frequencies when moving <br> from lower to higher levels |  |  |
| OR <br> Different frequencies of photons are <br> absorbed when the energy gap differs <br> (1) | Emit energy |  |  |$\quad$| M3 |
| :--- |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( b ) ( \text { iii) }}$ | Octahedral / octahedron (shape) <br> IGNORE <br> diagrams |  | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(b)(iv) | 3 ligands in an octahedral complex <br> ALLOW <br> $\mathrm{CH}_{2} \mathrm{CH}_{2}$ skeletal: $\mathrm{H}_{2} \mathrm{~N} \xrightarrow{\mathrm{NH}} \mathrm{H}_{2}$ <br> Skeletal not showing Hs on $\mathrm{NH}_{2}$ <br> (1) <br> bonds from $N$ to Cu , these can be lines, dots, wedges, arrows <br> ALLOW bond to one end of ligand only/incorrect ligand containing N <br> (1) <br> This structure scores both marks <br> IGNORE <br> Charge, brackets <br> Lone pairs on N | Two nitrogens from one ligand obviously at $180^{\circ}$ to the copper | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20(b)(v) | $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}+4 \mathrm{Cl}^{-} \rightarrow\left[\mathrm{CuCl}_{4}\right]^{2-}+6 \mathrm{H}_{2} \mathrm{O}$ <br> OR $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}+4 \mathrm{HCl} \rightarrow\left[\mathrm{CuCl}_{4}\right]^{2-}+6 \mathrm{H}_{2} \mathrm{O}+4 \mathrm{H}^{+}$ <br> OR $\begin{equation*} \left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}+4 \mathrm{HCl} \rightarrow\left[\mathrm{CuCl}_{4}\right]^{2-}+2 \mathrm{H}_{2} \mathrm{O}+4 \mathrm{H}_{3} \mathrm{O}^{+} \tag{1} \end{equation*}$ <br> IGNORE <br> state symbols even if incorrect <br> lack of [] <br> Tetrahedral <br> ALLOW <br> Square planar <br> M2 independent of M1 |  | 2 |

\(\left.$$
\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\
\text { Number }\end{array}
$$ \& Acceptable Answers \& Reject \& Mark <br>
\hline 20(b)(vi) \& \begin{array}{l}Step 1: acid-base / neutralisation <br>
Deprotonation (of complex) / protonation <br>
of ammonia <br>
ALLOW (ionic) precipitation (1) <br>
Step 2: Ligand and <br>
Exchange / substitution / replacement <br>
ALLOW <br>

'Ammonia substitutes for water'\end{array} \& electrophile\end{array}\right]\)| (1) |
| :--- |
| Final product: <br> $\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}\right]^{2+}$ <br> ALLOW <br> $\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\right]^{2+}$ <br> Round brackets, lack of [ ] brackets <br> (1) |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 20(b)(vii) | Step 1: pale blue precipitate/ solid forms <br> (1) | Step 2: (precipitate dissolves to give) deep / dark <br> blue solution (1) <br> Two correct colours with missing states can <br> score (1) <br> The blue colour in step 2 must be a darker blue <br> than the colour in step one. <br> e.g. Either pale blue step 1, blue step 2 or blue <br> step 1, dark blue in step 2 |  |



(Total for Question 20 = 20 marks)

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| *21(a) | Isomers of dichlorobenzene in which one has a <br> single bond between the C atoms bonded to Cl <br> and the other has a double bond have not been <br> found. <br> Can be shown on a Kekulé diagram. (1) | Cl can be in <br> positions other <br> than 1 or 2 | 2 |
| Cl2 would add <br> across each <br> double bond |  |  |  |
| (X-ray diffraction shows that) |  |  |  |
| all carbon-carbon bonds are the same (length) |  |  |  |
| OR intermediate between C=C and C-C (not as |  |  |  |
| in Kekulé ) |  |  |  |
| OR shows that benzene is a regular hexagon |  |  |  |
| ALLOW |  |  |  |
| All bonds are same length | The electron <br> density is even | (1) | IGNORE <br> Reference to bond angles <br> Benzene undergoes substitution reactions <br> rather than additions |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 21(b) | M1 <br> Phenol forms (2,4,6)-tribromophenol / formula ALLOW multiple substitution occurs <br> M2 <br> Phenol reacts with bromine water (at room temperature/ without heating) <br> M3 <br> Benzene forms bromobenzene/ $\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{Br}$ / one Br substitutes. <br> M4 <br> Benzene (reacts with bromine and) requires a catalyst of; iron/ iron(III) bromide/ a halogen carrier <br> ALLOW <br> Alternative M3 and M4 <br> M3 <br> Benzene reacts (with bromine) to form $\mathrm{C}_{6} \mathrm{H}_{6} \mathrm{Br}_{6}$ / 1,2,3,4,5,6 - hexabromocyclohexane /six Br add to it. <br> M4 <br> When heated in uv light <br> M2 and M4 dependent on correct or near miss for M1 and M3 respectively. | Hydroxyl benzene for phenol <br> Bromine water <br> Bromine water | 4 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 21(c) | M1 $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{Cl}+\mathrm{AlCl}_{3} \rightarrow \mathrm{AlCl}_{4}^{-}+\mathrm{C}_{2} \mathrm{H}_{5}^{+}$ <br> ALLOW <br> $\mathrm{FeBr}_{3} / \mathrm{FeCl}_{3} / \mathrm{AlBr}_{3}$ for $\mathrm{AlCl}_{3}$ <br> + on alkyl can in any position <br> M2 <br> Curly arrow from on or within the circle to $\mathrm{C}_{2} \mathrm{H}_{5}{ }^{+}$ <br> ALLOW curly arrow from anywhere within the hexagon <br> ALLOW curly arrow to any part of the $\mathrm{C}_{2} \mathrm{H}_{5}^{+}$ion, <br> including the + charge <br> TE for error in electrophile eg $\mathrm{C}_{2} \mathrm{H}_{4}{ }^{+}$ <br> M3 <br> Intermediate structure including charge with horseshoe covering at least 3 carbon atoms and facing the tetrahedral carbon atom and some part of the positive charge must be within the horseshoe <br> ALLOW dotted horseshoe <br> M4 <br> Curly arrow from $\mathrm{C}-\mathrm{H}$ bond to anywhere in the hexagon, reforming the delocalised structure <br> IGNORE <br> missing $\mathrm{H}^{+}$ <br> Reaction of $\mathrm{AlCl}_{4}^{-}$in last step <br> Correct Kekulé structures score full marks | Curly arrow on or outside the hexagon <br> Dotted bonds to H and $\mathrm{C}_{2} \mathrm{H}_{5}$ unless part of a 3D structure | 4 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 21(d)(i) | (New peak in phenylethene at) <br> $(\mathrm{C}=\mathrm{C}) 1669-1645\left(\mathrm{~cm}^{-1}\right)$ <br> OR <br> $(=\mathrm{C}-\mathrm{H}) 3095-3010\left(\mathrm{~cm}^{-1}\right)$ <br> ALLOW <br> $2962-2853\left(\mathrm{~cm}^{-1}\right)$ (alkane C-H) would not be present in <br> phenylethene. <br> If bonds are identified they must be correct. <br> IGNORE <br> Values for ethylbenzene peaks | Single <br> value <br> which <br> is not a <br> range | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |  |
| :--- | :--- | :--- | :--- | :--- |
| 21(d)(ii) |  | ALLOW <br> Bracket in polymer around side chain or round entire <br> unit |  |  |

## Section C

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 22(a)(i) | Compound A : nitrobenzene/ $\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NO}_{2}$ <br> (1) | 2 |  |
| Concentrated nitric acid + concentrated <br> sulfuric acid <br> and <br> temperature $55^{\circ} \mathrm{C}$ <br> ALLOW <br> "Concentrated nitric and sulfuric acid" <br> $50-60^{\circ} \mathrm{C}$ | Temperatures <br> above $60^{\circ} \mathrm{C}$ <br> Or less than <br> $50^{\circ} \mathrm{C}$ |  |  |
| "Heat at less than $55^{\circ} \mathrm{C}$ " | (1) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 22(a)(ii) | Tin + (concentrated) hydrochloric acid / Sn + <br> HCl <br> ALLOW <br> Iron/Fe for tin <br> IGNORE <br> Hydrogen $\mathrm{H}_{2}$ <br> followed by NaOH | Dilute HCl <br> $\mathrm{HCl}(\mathrm{Haq})$ | 1 |
| Sulfuric acid |  |  |  |$\quad$|  |
| :--- |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 22(b) | $\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NH}_{2}+2 \mathrm{CH}_{3} \mathrm{I} \rightarrow \mathrm{C}_{6} \mathrm{H}_{5} \mathrm{~N}\left(\mathrm{CH}_{3}\right)_{2}+2 \mathrm{HI}$ <br> OR $\begin{aligned} \mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NH}_{2}+2 \mathrm{CH}_{3} \mathrm{l} & \rightarrow \mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NH}\left(\mathrm{CH}_{3}\right)_{2}^{+} \\ & +\mathrm{I}^{-}+\mathrm{HI} \end{aligned}$ <br> ALLOW <br> $\mathrm{C}_{6} \mathrm{H}_{5}$ shown as delocalised ring <br> Reaction shown in 2 steps <br> Error in alkyl group if rest is correct e.g. ethyl for methyl <br> IGNORE <br> Use of molecular formulae |  | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 22(c) | $\mathrm{NaNO}_{2}+\mathrm{HCl} /$ sodium nitrite plus hydrochloric acid <br> ALLOW <br> Nitrous acid / $\mathrm{HNO}_{2}$ <br> Sulfuric acid for hydrochloric <br> A temperature in the range of $0-10\left({ }^{\circ} \mathrm{C}\right)$ <br> ALLOW $\begin{equation*} <10\left({ }^{\circ} \mathrm{C}\right) \tag{1} \end{equation*}$ <br> Mark independently | Concentrated hydrochloric acid <br> Concentrated sulfuric acid <br> Nitric acid | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 22(d)(i) |  | 1 |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 22(d)(ii) | $\mathrm{CH}_{3} \mathrm{COCl}+\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NH}_{2} \rightarrow \mathrm{CH}_{3} \mathrm{CONHC}_{6} \mathrm{H}_{5}+$ <br> OR <br> OR <br> Equation with 2 H substituted $2 \mathrm{CH}_{3} \mathrm{COCl}+\mathrm{C}_{6} \mathrm{H}_{5} \mathrm{NH}_{2} \rightarrow\left(\mathrm{CH}_{3} \mathrm{CO}\right)_{2} \mathrm{NC}_{6} \mathrm{H}_{5}+$ $2 \mathrm{HCl}$ <br> Balanced equation <br> CONH displayed, showing C=O connected to $\mathrm{N}-\mathrm{H}$ and connected to the benzene ring through N <br> ALLOW <br> NH for $\mathrm{N}-\mathrm{H}$ correct skeletal formula | $\mathrm{CH}_{3} \mathrm{NHCOC}_{6} \mathrm{H}_{5}$ | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 2 ( e ) ( i ) ~}$ | $\mathrm{C}_{9} \mathrm{H}_{11} \mathrm{NO}_{2}$ |  | 1 |
|  | ALLOW <br> Elements in any order eg $\mathrm{C}_{9} \mathrm{H}_{11} \mathrm{O}_{2} \mathrm{~N}$ <br> Answer written beside formula <br> IGNORE <br> $\mathrm{H}_{2} \mathrm{NC}_{6} \mathrm{H}_{4} \mathrm{COOCH}_{3}$ |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 22(e)(ii) | 92: $\mathrm{C}_{6} \mathrm{H}_{4} \mathrm{NH}_{2}{ }^{+}$ <br> ALLOW <br> $\mathrm{C}_{6} \mathrm{H}_{6} \mathrm{~N}^{+}$ <br> 120: : $\mathrm{C}_{6} \mathrm{H}_{4} \mathrm{NH}_{2} \mathrm{CO}^{+}$ <br> ALLOW <br> $\mathrm{C}_{7} \mathrm{H}_{6} \mathrm{NO}^{+}$ <br> OR <br> $\mathrm{C}_{6} \mathrm{H}_{4} \mathrm{CO}_{2}{ }^{+} / \mathrm{C}_{7} \mathrm{H}_{4} \mathrm{O}_{2}{ }^{+}$ <br> Penalise missing charges once only + charge can be anywhere on ion | Formulae with hexagons if number of H not clear $\mathrm{C}_{9} \mathrm{H}_{12}$ <br> Fragments with correct mass which could not form from benzocaine | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 22(e)(iii) | $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH}$ <br> OR <br> Skeletal formula, including H on OH <br> group <br> IGNORE <br> Molecular formula <br> $\mathrm{C}_{6} \mathrm{H}_{4} \mathrm{NH}_{2} \mathrm{COOH} / \mathrm{C}_{6} \mathrm{H}_{4} \mathrm{H}_{2} \mathrm{NCOOH}^{2}$ <br> OR <br> $\mathrm{C}_{6} \mathrm{H}_{4} \mathrm{NH}_{3}{ }^{+} \mathrm{COOH} / \mathrm{Cl}^{-} \mathrm{C}_{6} \mathrm{H}_{4} \mathrm{NH}_{3}{ }^{+} \mathrm{COOH}$ <br> $(1)$ | 2 |  |
| OR <br> Skeletal formula, including H on OH <br> group |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 22(f)(i) | Ignore SFs in M1,2,3 |  | 4 |
|  | M1 |  |  |
|  | 242.4 of $\mathrm{CO}_{2}$ contains |  |  |
|  | $((242.4 \times 12) / 44)=66.11 \mathrm{~g} \mathrm{C}$ |  |  |
|  | $76.30 \mathrm{~g} \mathrm{H}_{2} \mathrm{O}$ contains $((76.3 \times 2) / 18)$ |  |  |
|  | $=8.48 \mathrm{~g} \mathrm{H}$ |  |  |
|  | M2 |  |  |
|  | Mass O = |  |  |
|  | $(100-66.11-8.48-11.86)=$ | Calculation based |  |
|  | $13.55 \mathrm{~g}$ | on mass O in $\mathrm{CO}_{2}+$ |  |
|  | TE on M1 only if calculation method is correct |  |  |
|  | M3 |  |  |
|  | Moles per 100 g : |  |  |
|  | C 5.50 |  |  |
|  | H 8.48 |  |  |
|  | N 0.847 |  |  |
|  | O 0.847 |  |  |
|  | ALLOW TE from masses in M1 and M2 |  |  |
|  | (1) |  |  |
|  | M4 |  |  |
|  | $\mathrm{C}_{13} \mathrm{H}_{2} \mathrm{~N}_{2} \mathrm{O}_{2}$ |  |  |
|  | ALLOW |  |  |
|  | TE on M3 only if there are 13C |  |  |
|  | Elements in any order |  |  |
|  | (1) |  |  |


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

(Total for Question 22 = 20 marks)
Total for Section C = $\mathbf{2 0}$ marks

Total for Paper = 90 marks

