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Mark Scheme (Results)
January 2013

GCE Chemistry (6CH05) Paper 01
General Principles of Chemistry II Transition Metals and Organic 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.
- Mark schemes will indicate within the table where, and which strands of QWC, are being assessed. The strands are as follows:
i) ensure that text is legible and that spelling, punctuation and grammar are accurate so that meaning is clear
ii) select and use a form and style of writing appropriate to purpose and to complex subject matter
iii) organise information clearly and coherently, using specialist vocabulary when appropriate


## Using the Mark Scheme

Examiners should look for qualities to reward rather than faults to penalise. This does NOT mean giving credit for incorrect or inadequate answers, but it does mean allowing candidates to be rewarded for answers showing correct application of principles and knowledge. Examiners should therefore read carefully and consider every response: even if it is not what is expected it may be worthy of credit.

The mark scheme gives examiners:

- an idea of the types of response expected
- how individual marks are to be awarded
- the total mark for each question
- examples of responses that should NOT receive credit.
/ means that the responses are alternatives and either answer should receive full credit.
( ) means that a phrase/word is not essential for the award of the mark, but helps the examiner to get the sense of the expected answer.
Phrases/words in bold indicate that the meaning of the phrase or the actual word is essential to the answer.
ecf/TE/cq (error carried forward) means that a wrong answer given in an earlier part of a question is used correctly in answer to a later part of the same question.

Candidates must make their meaning clear to the examiner to gain the mark. Make sure that the answer makes sense. Do not give credit for correct words/phrases which are put together in a meaningless manner. Answers must be in the correct context.

## Quality of Written Communication

Questions which involve the writing of continuous prose will expect candidates to:

- write legibly, with accurate use of spelling, grammar and punctuation in order to make the meaning clear
- select and use a form and style of writing appropriate to purpose and to complex subject matter
- organise information clearly and coherently, using specialist vocabulary when appropriate.
Full marks will be awarded if the candidate has demonstrated the above abilities. Questions where QWC is likely to be particularly important are indicated (QWC) in the mark scheme, but this does not preclude others.


## Section A (multiple choice)

| Question <br> Number | Correct Answer | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{1 ( a )}$ | C |  | $\mathbf{1}$ |
| (b) | A |  | $\mathbf{1}$ |
| Question <br> Number Correct Answer Reject  <br> $\mathbf{2 ( a )}$ C  Mark <br> (b) A  $\mathbf{1}$ <br> (c) B  $\mathbf{1}$ |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{3}$ | C |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{4}$ | C |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{5}$ | B |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{6 ( a )}$ | D |  | $\mathbf{1}$ |
| (b) | C |  | $\mathbf{1}$ |
| (c) | A |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{7}$ | B |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{8}$ | C |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{9}$ | B |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{1 0}$ | C |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| ---: | :--- | :--- | :--- |
| $\mathbf{1 1}$ | D |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 2}$ | A |  | $\mathbf{1}$ |
| Question <br> Number Acceptable Answers Reject Mark <br> $\mathbf{1 3 ( a )}$ D  $\mathbf{1}$ <br> (b) B  $\mathbf{1}$ <br> (c) A  $\mathbf{1}$ |  |  |  |

## Section B

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 14 (a) | $\mathbf{A}=\operatorname{copper}(\mathrm{II})$ hydroxide $/ \mathrm{Cu}(\mathrm{OH})_{2} /$ <br> $\mathrm{Cu}(\mathrm{OH})_{2}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}$ <br> B = copper(II) oxide / CuO <br> $\mathbf{C}=$ tetraamminecopper (II) / $\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}{ }^{2+} /$ $\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}{ }^{2+}$ <br> ALLOW Cu( $\left.\mathrm{NH}_{3}\right)_{6}{ }^{2+}$ / hexaamminecopper(II) <br> $\mathbf{D}=$ copper $/ \mathrm{Cu} / \operatorname{copper}(0) / \mathrm{Cu}(0)$ <br> $\mathbf{E}=\operatorname{copper}(\mathrm{II})$ sulfate $/ \mathrm{CuSO}_{4} / \mathrm{Cu}^{2+} /$ $\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}{ }^{2+}$ <br> $\mathbf{F}=$ diamminecopper(I) / $\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{2}{ }^{+}$ <br> ALLOW coordination numbers 1-6 in $\mathbf{F}$ Oxidation number separate from name <br> IGNORE <br> state symbols even if incorrect names without oxidation numbers except for $\mathbf{D}$ | Formulae with incomplete or unbalanced charges <br> Incorrect oxidation states even with correct formulae | 6 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 4}$ (b) | (Dilute) sulfuric acid $/ \mathrm{H}_{2} \mathrm{SO}_{4} / \mathrm{H}_{2} \mathrm{SO}_{4}(\mathrm{aq})$ <br> ALLOW concentrated |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 4 ( c ) ( i )}$ | (transition metal / d-block element) complex(es) <br> /complex ion(s) <br> IGNORE <br> ammines | Complex <br> molecules <br> amines, ions, <br> ligands | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 14 (c)(ii) | Copper ion in C has partially filled d orbital(s) / subshell / 3d ${ }^{9}$ <br> ALLOW <br> unpaired d electron <br> d shell <br> Copper ion in $\mathbf{F}$ has (completely) filled $\mathbf{d}$ orbitals / <br> subshell / 3d ${ }^{10}$ <br> Reference to complete / incomplete d orbitals max 1 <br> EITHER <br> Electronic transitions between partially filled (d) orbitals (of different energy) are possible <br> OR <br> Electronic transitions between (completely) filled <br> (d) orbitals (of different energy) are not possible <br> ALLOW <br> Equivalent words for transition e.g. promotion / <br> jump / movement <br> Penalise use of just 'shell' once IGNORE references to electrons returning to lower energy levels and emission of light | d orbitals empty <br> no unpaired electrons (in F) orbital (singular) <br> Splitting impossible because d orbitals full | 3 |


| Question | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| Number |  |  |  |
| $\mathbf{1 4}$ | Copper(I) is oxidized (to copper(II)) | (1) |  |
| (c)(iii) | ALLOW F / it is oxidized <br> By oxygen / air <br> Second mark depends on first <br> IGNORE <br> 'shaking' | (1) |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 4 ( d ) ( i )}$ | (simultaneous) oxidation and reduction (1) <br> OR <br> Simultaneous increase or decrease in oxidation <br> number <br> of an element <br> ALLOW <br> 'Species' 'atoms of the same type' for 'element' <br> Explanation in terms of copper(I) <br> IGNORE <br> Atom / ion / compound / substance / reactant | molecule | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 4}$ | $2 \mathrm{Cu}^{+} \rightarrow \mathrm{Cu}+\mathrm{Cu}^{2+}$ | Non-ionic <br> (d)(ii) <br>  <br>  <br>  <br>  <br>  <br>  <br>  <br>  <br>  <br>  <br>  <br>  <br>  <br> OR <br> $2 \mathrm{CuI}+2 \mathrm{H}^{+} \rightarrow \mathrm{Cu}^{2+}+\mathrm{Cu}^{2+}+2 \mathrm{HI}$ <br> IGNORE state symbols even if incorrect | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{aligned} & 14 \\ & \text { (d) (iii) } \end{aligned}$ | ALLOW <br> The use of cell notation (as in the Data Booklet SEP table) in place of equations $\text { e.g. } \mathrm{Cu}^{+}(\mathrm{aq}) \mid \mathrm{Cu}(\mathrm{~s}) \mathrm{E}^{\ominus}=+0.52(\mathrm{~V})$ <br> (from the data book the equations are) $\begin{align*} & \mathrm{Cu}^{+}(\mathrm{aq})+\mathrm{e}^{-} \rightarrow \mathrm{Cu}(\mathrm{~s}) \mathrm{E}^{\ominus}=+0.52(\mathrm{~V}) \\ & \mathrm{Cu}^{2+}(\mathrm{aq})+\mathrm{e}^{-} \rightarrow \mathrm{Cu}^{+}(\mathrm{aq}) \mathrm{E}^{\ominus}=+0.15(\mathrm{~V})  \tag{1}\\ & \mathrm{So}^{\ominus} \text { cell }=0.52-0.15=+0.37(\mathrm{~V}) \tag{1} \end{align*}$ <br> Correct answer including sign with no working scores full marks <br> TE for second mark for use of $\mathrm{Cu}^{2+} \mathrm{ICu}+0.34(\mathrm{~V})$ which gives $+0.19(\mathrm{~V}) /+0.18(\mathrm{~V})$ <br> No TE on incorrect equation in (d)(ii) | Answer without + sign | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{aligned} & \hline 14 \\ & \text { (d)(iv) } \end{aligned}$ | ALLOW <br> In both schemes the use of cell notation (as in the Data Booklet SEP table) in place of equations e.g. $\mathrm{Cu}^{2+}(\mathrm{aq}) \mid \mathrm{Cu}(\mathrm{s}) \quad \mathrm{E}^{\ominus}=+0.34(\mathrm{~V})$ <br> Penalise omission of electrons from equations and vertical lines from cell diagrams and reversal of equation without reversing sign. once only <br> IGNORE omission of + sign for all $E^{\ominus}$ values <br> Scheme 1 (oxidation of copper) <br> Copper (formed (by disproportionation)) is oxidized (by nitric acid) must be stated in words stand alone mark <br> Relevant half equations are $\begin{align*} & \mathrm{Cu}^{2+}(\mathrm{aq})+2 \mathrm{e}^{-} \rightarrow \mathrm{Cu}(\mathrm{~s}) \quad \mathrm{E}^{\ominus}=+0.34(\mathrm{~V})(\mathbf{1}) \\ & 2 \mathrm{NO}_{3}^{-}(\mathrm{aq})+4 \mathrm{H}^{+}(\mathrm{aq})+2 \mathrm{e}^{-} \rightarrow \mathrm{N}_{2} \mathrm{O}_{4}(\mathrm{~g})+2 \mathrm{H}_{2} \mathrm{O}(\mathrm{I}) \\ & \mathrm{E}^{\ominus}=+0.80(\mathrm{~V}) \\ & \mathrm{OR}^{-} \\ & \mathrm{NO}_{3}^{-}(\mathrm{aq})+3 \mathrm{H}^{+}(\mathrm{aq})+2 \mathrm{e}^{-} \rightarrow \mathrm{HNO}_{2}(\mathrm{aq})+\mathrm{H}_{2} \mathrm{O}(\mathrm{I}) \\ & \mathrm{E}^{\ominus}=+0.94(\mathrm{~V}) \tag{1} \end{align*}$ <br> Correct overall equation scores both marks: $\mathrm{Cu}+2 \mathrm{NO}_{3}^{-}+4 \mathrm{H}^{+} \rightarrow \mathrm{Cu}^{2+}+\mathrm{N}_{2} \mathrm{O}_{4}+2 \mathrm{H}_{2} \mathrm{O}$ <br> OR $\mathrm{Cu}+\mathrm{NO}_{3}^{-}+3 \mathrm{H}^{+} \rightarrow \mathrm{Cu}^{2+}+\mathrm{HNO}_{2}+\mathrm{H}_{2} \mathrm{O}$ <br> So $\mathrm{E}^{\ominus}$ cell is $+0.46(\mathrm{~V}$ ) (or $+0.60(\mathrm{~V}$ ) or just 'positive') <br> Scheme 2 (oxidation of copper(I) <br> Copper(I) iodide / $\mathrm{Cu}^{+}$is oxidized (by nitric acid) must be stated in words <br> stand alone mark $\begin{align*} & \mathrm{Cu}^{2+}(\mathrm{aq})+\mathrm{e}^{-} \rightarrow \mathrm{Cu}^{+}(\mathrm{aq}) \mathrm{E}^{\ominus}=+0.15(\mathrm{~V})  \tag{1}\\ & 2 \mathrm{NO}_{3}^{-}(\mathrm{aq})+4 \mathrm{H}^{+}(\mathrm{aq})+2 \mathrm{e}^{-} \rightarrow \mathrm{N}_{2} \mathrm{O}_{4}(\mathrm{~g})+2 \mathrm{H}_{2} \mathrm{O}(\mathrm{I}) \\ & \mathrm{E}^{\ominus}=+0.80(\mathrm{~V}) \\ & \mathrm{OR} \\ & \mathrm{NO}_{3}^{-}(\mathrm{aq})+3 \mathrm{H}^{+}(\mathrm{aq})+2 \mathrm{e}^{-} \rightarrow \mathrm{HNO}_{2}(\mathrm{aq})+\mathrm{H}_{2} \mathrm{O}(\mathrm{I}) \\ & \mathrm{E}^{\ominus}=+0.94(\mathrm{~V}) \end{align*}$ <br> Correct overall equation scores both marks: |  | 4 |


|  | $2 \mathrm{Cu}^{+}+2 \mathrm{NO}_{3}^{-}+4 \mathrm{H}^{+} \rightarrow 2 \mathrm{Cu}^{2+}+\mathrm{N}_{2} \mathrm{O}_{4}+2 \mathrm{H}_{2} \mathrm{O}$ <br> $2 \mathrm{Cu}^{+}+\mathrm{NO}_{3}^{-}+3 \mathrm{H}^{+} \rightarrow 2 \mathrm{Cu}^{2+}+\mathrm{HNO}_{2}+\mathrm{H}_{2} \mathrm{O}$ <br> So E cell is $+0.65(\mathrm{~V})($ or $+0.79(\mathrm{~V})$ or just (1) <br> 'positive') <br> IGNORE <br> (omission of) state symbols even if incorrect |  |
| :--- | :--- | :--- |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 5} \mathbf{( a ) ( i )}$ | (vitamin C / ascorbic acid ) oxidation / <br> oxidized / oxidised | Redox / oxidation- <br> reduction / <br> reduction-oxidation | $\mathbf{1}$ |
| ALLOW <br> oxidisation |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 5}$ | (very) pale yellow / straw coloured <br> (a)(ii) <br> IGNORE 'just before the end-point' <br> blue-black to colourless (both needed) $\quad$ (1) | Just 'yellow' | Clear |
| Accept (dark) blue or black <br> ALLOW <br> pale yellow / straw coloured to colourless for <br> $1 / 2$ |  | $\mathbf{2}$ |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{align*} & 15  \tag{1}\\ & \text { (a) (iii) } \end{align*}$ | $\begin{align*} & \text { Moles } \mathrm{S}_{2} \mathrm{O}_{3}{ }^{2-}=27.85 \times 10^{-3} \times 0.0631 \\ &\left(=1.757335 \times 10^{-3}\right) \\ & \text { moles of } \mathrm{I}_{2} \text { remaining }=\text { Moles } \mathrm{S}_{2} \mathrm{O}_{3}{ }^{2-} \div 2 \\ & \quad=27.85 \times 10^{-3} \times 0.0631 \div 2 \\ &=8.786675 \times 10^{-4}=8.79 \times 10^{-4} \tag{1} \end{align*}$ <br> Moles ascorbic acid $=$ moles $\mathrm{I}_{2}$ at start moles $\mathrm{I}_{2}$ remaining $\begin{align*} & =2.00 \times 10^{-3}-8.786675 \times 10^{-4} \\ & =1.1213325 \times 10^{-3}=1.12 \times 10^{-3} \tag{1} \end{align*}$ <br> $M_{r}$ (ascorbic acid) $=176$ <br> Mass ascorbic acid in $250 \mathrm{~cm}^{3}=10 \times \mathrm{M}_{\mathrm{r}} \times$ <br> moles ascorbic acid $\begin{align*} & =10 \times 176 \times 1.1213325 \times 10^{-3}  \tag{1}\\ & (=1.97355) \end{align*}$ <br> Percentage ascorbic acid in tablet $100 \times$ mass ascorbic acid in $250 \mathrm{~cm}^{3} \div 2$ $\begin{align*} & =100 \times 10 \times 176 \times 1.1213325 \times 10^{-3} \div 2 \\ & =98.67726=98.7 \% \tag{1} \end{align*}$ <br> IGNORE SF except 1 SF <br> Premature rounding gives $98.5 \%$ (5) <br> Correct answer with no working scores full marks <br> TE at each stage of the calculation. | Answers greater than 100\% | 5 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 5 ( a ) ( i v ) ~}$ | EITHER <br> Using larger mass reduces the percentage <br> error / uncertainty (in weighing) | Just 'reduces the <br> percentage error' | $\mathbf{1}$ |
|  | OR <br> Using larger amount reduces the percentage <br> error / uncertainty in weighing <br> OR <br> Reverse discussion of two tablets | Titration value will <br> be larger (with four <br> tablets) so reduces <br> the percentage <br> error (in volume <br> measurement) | ALLOW <br> using four tablets gives a more <br> representative sample |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 5 ~ ( b ) ( i )}$ |  | 2 |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 5}$ (b)(ii) | First mark <br> Use of (plane-) polarized light (mentioned (1) <br> somewhere) <br> ALLOW <br> Use a polarimeter <br> Second mark <br> Pure optical isomer / enantiomer) rotates <br> the plane of (plane-) polarized light <br> OR <br> racemic mixture has no effect on the plane <br> of (plane-) polarized light <br> (1) |  | $\mathbf{2}$ |
| IGNORE <br> optically active / inactive <br> ALLOW <br> rotates plane-polarized light scores 2 |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 5 ( b ) ( i i i )}$ | (Ester group / vitamin C / it) is hydrolysed <br> ALLOW <br> Vitamin C is oxidized <br> Ester / vitamin C is broken down to form <br> carboxylic acid and alcohol (groups) <br> IGNORE <br> Just 'breaks down' | C=O is broken | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 6 ( a ) ( i )}$ | The delocalization of the ( $\pi$ ) electrons of the ring <br> make benzene more stable (than 1,3,5- (1) <br> cyclohexatriene) |  | $\mathbf{2}$ |
|  | IGNORE bonding in benzene is strong <br> Substitution retains this (stable) arrangement <br> OR <br> Addition removes this (stable) arrangement <br> (1) |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(a)(ii) |  <br> Formation of electrophile (curly arrow, structural formulae not required). Positive charge may be anywhere on the electrophile <br> ALLOW $\mathrm{HCl}+\mathrm{CO}$ for HCOCl <br> ALLOW Non-displayed electrophile <br> Curly arrow from benzene ring to electrophile <br> Wheland structure with gap opposite tetrahedral carbon <br> Curly arrow from $\mathrm{C}-\mathrm{H}$ bond into ring and formation of correct organic product <br> OR <br> Kekulé structures <br> IGNORE <br> Use of $\mathrm{AlCl}_{4}^{-}$to pick off proton <br> Proton product <br> First curly arrow may come from any part of the delocalisation circle <br> Second curly arrow may come from any part of the C-H bond <br> Positive charge on the Wheland structure may be in any part of the horseshoe | $\text { - } \mathrm{COH} /-\mathrm{HCO}$ <br> Positive charge on the tetrahedral carbon | 4 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(a)(iii) | In each step the second mark is dependent on the first <br> Step 2 <br> Potassium dichromate((VI)) / $\mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7}$ / <br> sodium dichromate((VI)) / $\mathrm{Na}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7}$ <br> ALLOW <br> Potassium manganate ((VII)) / $\mathrm{KMnO}_{4}$ <br> Sodium manganate ((VII)) / $\mathrm{NaMnO}_{4}$ <br> Stand alone mark <br> Sulfuric acid / $\mathrm{H}_{2} \mathrm{SO}_{4}$ (ALLOW nitric acid) <br> Ignore 'concentrated' <br> ALLOW <br> Acidified potassium (/ sodium) dichromate((VI)) OR <br> Acid and potassium (/ sodium) dichromate((VI)) <br> $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}$ and $\mathrm{H}^{+} \mathrm{OR}$ acidified dichromate((VI)) <br> (1) <br> Step 3 <br> Lithium tetrahydridoaluminate((III)) / LiAlH ${ }_{4}$ <br> OR <br> Lithium aluminium hydride <br> Stand alone mark <br> (Dry) ether / ethoxyethane / (di)ethyl ether (1) <br> Sodium borohydride / $\mathrm{NaBH}_{4}$ in ethanol, alkali or water scores $1 / 2$ | Incorrect oxidation number <br> Hydrochloric acid <br> Hydrogen and catalyst / Tin and HCl | 4 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 6 ( b )}$ | Marking Point 1 <br> Electron density of the ring increased (1) <br> Stand alone mark <br> Marking Point 2 <br> Due to donation of oxygen / OH group lone pair to <br> the ring <br> (1) <br> Marking Point 3 and 4 <br> Any two from <br> in phenol oxygen / OH group attached directly to <br> ring <br> Oxygen / OH group in phenylmethanol too far away <br> / not attached directly to ring <br> (In phenol) lone pair overlaps with the $\pi$ electrons <br> / delocalised electrons (of the ring) <br> ALLOW p orbital for lone pair for this mark <br> (2) | $\mathbf{4}$ |  |

Total for Q16 = 14 Marks
TOTAL FOR SECTION B = 50

## Section C

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( a ) ( \mathbf { i ) }}$ | There is a barrier to rotation about a (C=C) bond |  |  |
|  | ALLOW restricted / limited / no rotation (1) | Just 'molecule <br> cannot rotate' | $\mathbf{2}$ |
|  | Each carbon atom (in the C=C double bond) has <br> (two) different atoms / groups attached (1) <br> IGNORE reference to priority groups |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( a ) ( i i )}$ | There is a barrier to / restricted rotation about <br> the ring <br> OR <br> The ring behaves like a double bond | Reference to <br> benzene ring | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( a ) ( \text { iii) }}$ |  | Omission of <br> amine $\mathrm{CH}_{2}$ | $\mathbf{1}$ |
|  | Any diagram of the correct molecule showing the <br> groups (attached to the ring) on same side of <br> the ring <br> OR <br> zwitterion <br> ALLOW <br> Amine group in skeletal form |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( a ) ( i v )}$ | Tranexamic acid exists as a zwitterion <br> OR <br> Diagram of zwitterion <br> OR <br> Description of zwitterion formation <br> So the (strongest) intermolecular forces are ionic <br> (strong) <br> ALLOW electrostatic for ionic | $\mathbf{3}$ |  |
|  | IGNORE H bonding in tranexamic acid if either of the <br> first two marks scored. Otherwise... |  |  |
|  | Hydrogen bonding in tranexamic acid scores 1/2 max <br> Undecane has (only) (much weaker) London / <br> dispersion / van der Waals / temporary induced dipole <br> (-induced dipole) forces / interactions |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( b ) ( i )}$ | Phosphorus(v) chloride / $\mathrm{PCl}_{5}$ | HCl | $\mathbf{1}$ |
|  | ALLOW <br> phosphorus pentachloride / <br> phosphorus(III) chloride / $\mathrm{PCl}_{3} /$ phosphorus <br> trichloride <br> Thionyl chloride (sulfur dichloride oxide) $/ \mathrm{SOCl}_{2}$ |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( b ) ( i i )}$ | First mark <br> amide linkage <br> ALLOW CONH for amide linkage <br> Second mark <br> Completion of structure (brackets not required) with <br> displayed or skeletal formula <br> Second mark dependent on first <br> Dimer scores amide linkage mark only | $\mathbf{2}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( b ) ( \text { iii) }}$ | Condensation / addition-elimination <br> (polymerization) | Addition <br> (polymerization) <br> Elimination <br> (polymerization) <br> Polyamide <br> formation | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( b ) ( i v )}$ | Protein / proteins / polypeptide / polypeptides / <br> peptide / peptides <br> ALLOW Enzyme / Enzymes | Nylon <br> Polyamide <br> amino acids | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(c)(i) | Check sequence of letters. Candidates may have labelled the groups of hydrogen atoms with different letters, which is fine. <br> First mark <br> Unique NH (at e) <br> Second mark <br> Unique $\mathrm{CH}_{2}$ (at c) <br> Third mark <br> CH (at d) and CH (at f) with different unique labels <br> Fourth mark <br> (1) <br> $2 \mathrm{CH}_{2}$ (at a) and $2 \mathrm{CH}_{2}$ (at b) with different new labels |  | 4 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( c ) ( \mathbf { i i ) }}$ | $\mathrm{C}=\mathrm{O}$ amide (stretching vibrations are in the region) <br> $1700-1630 \mathrm{~cm}^{-1}$ | Ketone | 2 |
|  | $\mathrm{N}-\mathrm{H}$ amide (stretching vibrations are in the region) <br> $3500-3140 \mathrm{~cm}^{-1}$ | (1) <br> Amine (for <br> amide) |  |
|  | Amide only needs to be mentioned once but... |  |  |
| These answers without mention of amide max 1 |  |  |  |
| Amides have peaks in these regions max 1 |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( c ) ( i i i )}$ | Any two from <br> In the trans isomer the (amine and acid chloride) groups <br> are too far apart to react intramolecularly / to form M <br> OR <br> Because the groups are on opposite sides of the (plane of <br> the) ring <br> OR <br> More likely to polymerize / react with adjacent molecules. <br> Marks may also be scored by a reverse argument: (2) <br> In the cis isomer the (amine and acid chloride) groups <br> are on the same side of the (plane of the) ring <br> So close enough to react intramolecularly / to form M (1) | bond |  |

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