| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) | (i) |  <br> all correct 2 marks $\checkmark \checkmark$ <br> amide link only $\checkmark$ | 2 | ALLOW <br> - any correct structural formula including skeletal <br> - mixtures of structural and skeletal <br> - -CONH- <br> - Cs and Hs on ring <br> IGNORE brackets and n etc |
|  |  | (ii) | -CONH- circled $\checkmark$ | 1 | ALLOW adjacent C atoms in circle |
|  |  | (iii) | 1,6-diamino $\checkmark$ hexane $\checkmark$ OR hexane $\checkmark$-1,6-diamine $\checkmark$ | 2 | If butane ALLOW 1,4-diamino for ecf mark IGNORE commas and dashes ALLOW 1,6-hexanediamine |
|  | (b) | (i) |  | 1 | ALLOW any formula that makes structure clear ALLOW cyclic amide |


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| :--- | :--- | :--- | :--- | :--- | :--- |
| (c) | (i) | condensation <br> AND water is eliminated/formed/lost $\checkmark$ <br> IF cyclic amide in 1b(i) THEN addition AND hydrolysis (of ring) / <br> water added | 'addition' or 'addition polymerisation' is a CON <br> ALLOW 'small molecule' instead of water <br> any other named small molecule is a CON |  |
| (phains/molecules are (highly) ordered/aligned AW $\checkmark$ | 1 | ALLOW labelled diagram that shows alignment of <br> chains by using parallel lines |  |  |
| AW means other suitable phrases |  |  |  |  |
| eg |  |  |  |  |
| 'stacked closely and neatly' |  |  |  |  |
| 'arranged regularly'/'regularity of chains' |  |  |  |  |


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| (d) | (i) | $\begin{array}{ll}\text { 1. } & \text { elimination } \checkmark \\ \text { 2. } & \text { addition } \checkmark \\ \text { 3. } & \text { addition } \checkmark\end{array}$ | 3 | IGNORE 'nucleophilic' |
|  | (ii) | can be used in step 3 AW $\checkmark$ | 1 | NOT step 2 <br> ALLOW <br> - can be sold <br> - can be used as a source of energy <br> - can be recycled |
| (e) |  | acidified dichromate $\checkmark$ | 1 | IGNORE <br> - metal cation <br> - name of acid <br> - wrong formulae if name given <br> - oxidation state of 'dichromate' <br> ALLOW $\mathrm{H}_{2} \mathrm{SO}_{4} / \mathrm{H}^{+}$and $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}$ DO NOT ALLOW just dichromate |
|  |  | Total | 17 |  |


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| 2 | (a) |  | $-\mathrm{COOH} /$ carboxyl / carboxylic acid group $\checkmark$ is a proton $/ \mathrm{H}^{+}$donor / loses $\mathrm{H}^{+} \checkmark$ | 2 | correct equation showing dissociation gains both marks ALLOW 'gives $\mathrm{H}^{+}$ <br> Mark separately |
|  | (b) | (i) | $\mathrm{CH}_{2} \mathrm{OHCOOH}+\mathrm{NaOH} \rightarrow \mathrm{CH}_{2} \mathrm{OHCOONa}+\mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 | ```ALLOW CH2OHCOO- Na+, Na+CH2OHCOO- OR Na(CH2OHCOO) IGNORE state symbols NOT CH2OHNaCOO``` |
|  |  | (ii) | 1. moles of NaOH used in titration $=16.00 / 1000 \times 0.250 \checkmark=0.00400$ <br> 2. moles of glycolic acid used in titration $=$ answer from 1, scaled by ratio in equation in 2bi $\checkmark$ 0.00400 <br> CHECK equation in 2(b)(i) <br> 3A. moles of glycolic acid in $250 \mathrm{~cm}^{3}$ $=(\text { answer from 2) } \times 10 \checkmark=0.0400 \mathrm{~mol}$ <br> OR <br> 3B. mass of glycolic in $25 \mathrm{~cm}^{3}$ $=\left(\right.$ answer from 2) $\times \boldsymbol{M}_{\mathrm{r}}$ of glycolic acid $\checkmark 0.304 \mathrm{~g}$ <br> OR <br> 3C. concentration of glycolic acid $=\left(\right.$ answer from 2) $\times 1000 / 25 \checkmark 0.16 \mathrm{~mol} \mathrm{dm}^{-3}$ <br> 4. $\quad M_{\mathrm{r}}$ of glycolic acid $=76$.(0) $\checkmark$ | 6 | There are several possible routes through this question after point 2, the 'mole route' $A$, the 'mass route' $B$ and the 'concentration route' $C$ <br> If final answer is incorrect please annotate with ticks where the marks are awarded <br> 2. ALLOW by implication if 0.004 used subsequently <br> The marks are awarded for the working out given in bold OR the correctly calculated answer to that working (but no mark if calculated answer is shown and is wrong) |


| Question |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
|  |  | ```5A. moles of glycolic acid in \(100 \mathrm{~cm}^{3}\) \(=(\) answer from 3A) \(\times 100 / 14 \checkmark \quad 0.286 \mathrm{~mol}\) OR 5Bi. mass of glycolic acid in \(250 \mathrm{~cm}^{3}\) \(=\left(\right.\) answer from 3A) \(\times \boldsymbol{M}_{\mathbf{r}}\) of glycolic acid \(\checkmark 3.04 \mathrm{~g}\) OR 5Bii. mass of glycolic in \(250 \mathrm{~cm}^{3}\) \(=(\) answer from 3B \() \times 10=\checkmark 3.04 \mathrm{~g}\) OR 5C. concentration of undiluted glycolic acid \(=\left(\right.\) answer from 3C) \(\times 250 / 14 \checkmark 2.86 \mathrm{~mol} \mathrm{dm}^{-3}\) 6. mass of glycolic acid in \(100 \mathrm{~cm}^{3}\) undiluted \(=\left(\right.\) answer from 5A) \(\times \boldsymbol{M}_{r}\) of glycolic acid OR = (answer from 5B) \(\times\) 100/14 OR = (answer from 5C/10) \(\times M_{r}\) of glycolic acid \(=21.7\) (3 sf) AND correct comment``` |  | ALLOW ecf for incorrect equation AND between each step <br> ALLOW 'Acnegone' for 'glycolic acid' <br> ALLOW answers in standard form <br> The following on the answer line with correct corresponding comment, score as follows, irrespective of working or lack of it: <br> 21.7 scores 6 <br> 1.22 scores 5 (error in 5C) <br> 12.2 scores 5 (error in 6C) <br> 2.17 scores 5 (error in 3A or 5Bii) <br> 3.04 scores 5 (error in 6) <br> these to other sf OR with incorrect comment score one mark less <br> If one of the answers above applies place correct number of ticks by answer |
| (c) | (i) |  | 2 | IGNORE where the circle cuts the bond as long COO is included ALLOW adjacent $C$ atoms in circle <br> ALLOW butylethanoate without gap |
|  | (ii) | butan-1-ol <br> concentrated sulfuric acid/hydrochloric acid $\checkmark$ | 2 | MUST HAVE number 1 <br> DO NOT ALLOW ecf for alcohol in 2(c)(i) <br> ALLOW formula for acid ONLY <br> IGNORE spelling of name for $\mathrm{H}_{2} \mathrm{SO}_{4}$ as long as it is clear |


| Question |  | Answer | Marks | Guidance |
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|  | (iii) | ethanoic acid: hydrogen bonding $\checkmark$ IGNORE pd-pd compound D: pd-pd $\checkmark$ hydrogen bonding is CON instantaneous dipole - induced dipole / id-id in both $\checkmark$ | 3 | ALLOW 'pd-pd' / id-id abbreviations here ALLOW Van der Waals for id-id |
| (d) | (i) | The answer requires a comment for a carboxylic acid AND an alcohol so award ONE mark from each section below: <br> CARBOXYLIC ACID <br> (strong) peak at $1743\left(\mathrm{~cm}^{-1}\right)$ shows $\mathrm{C}=\mathrm{O}$ in ester not acid OR no peak $1700-1725\left(\mathrm{~cm}^{-1}\right)$ shows no $\mathrm{C}=\mathrm{O}$ in acid OR no broad peak at 2500-3200 (cm ${ }^{-1}$ ) shows no O-H in acid $\checkmark$ <br> ALCOHOL <br> no peak greater than 3000 / in range $3200-3600$ (allow 3640) so no O-H in alcohol | 2 | answers may be given on spectrum <br> to score each point, range, bond and group in which it is found must be given <br> ALLOW carboxyl or carboxylic acid or COOH or ethanoic acid (or formula) for 'acid' <br> FOR O-H ALLOW OH / hydroxyl FOR C=O ALLOW carbonyl NOT CO |
|  | (ii) | peak at $\boldsymbol{m l z} 73$ : $\mathrm{CH}_{3} \mathrm{COOCH}_{2} / \mathrm{C}_{3} \mathrm{H}_{5} \mathrm{O}_{2}$ <br> positive charge on any formula $\checkmark$ <br> species lost: $\mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3} / \mathrm{C}_{3} \mathrm{H}_{7}(\mathrm{NO} \text { charge) } \checkmark$ | 3 | ALLOW any correct structural or molecular formula for both answers <br> ALLOW $\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{O}^{+}$ <br> IF $\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{O}^{+}$given above <br> THEN species lost must be $\mathrm{C}_{2} \mathrm{H}_{3} \mathrm{O} / \mathrm{CH}_{3} \mathrm{CO}$ |
| (e) | (i) | nucleophilic $\checkmark$ addition $\checkmark$ | 2 |  |


| Question |  | Answer | Marks | Guidance |
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| (ii) |  |  |  |  |



| Question |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| (c) | (i) |  <br> ANY TWO of the THREE CORRECT chiral Cs ONLY $\checkmark$ | 1 |  |
|  | (ii) | arrows as in (d) (i) $\checkmark$ | 1 | IGNORE extra arrows pointing to $\mathrm{C}-\mathrm{N}$ bonds at ends of chain. <br> Any other arrows are a CON |
|  | (iii) |  <br> correct amino acid (see opposite) $\checkmark$ <br> correct formula of ion $\checkmark$ | 2 | IGNORE any cations <br> ALLOW any correct structural formula <br> IGNORE species formed from amino acids to right and left <br> IGNORE <br> - added $\mathrm{H}^{+}$to $-\mathrm{NH}_{2}$ <br> - negative charge on alcohol i.e. $-\mathrm{O}^{-}$ <br> - -COOH instead <br> NO ecf <br> IF anion formed at $\mathrm{CH}_{2} \mathrm{O}^{-}$then 1 mark max for amino acid |
|  | (iv) | secondary: <br> folding/twisting of polypeptide/amino acid chains/primary structure <br> tertiary: <br> further/final folding OR 3D shape/structure | 2 | ALLOW (alpha)-helix/coiled OR (beta-)sheets / pleated sheets <br> ALLOW overall/global structure |
|  |  | Total | 14 |  |


| Question |  |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (a) | (i) | correct inside oval (around peroxy Os) rest correct $\checkmark$ | 2 | ALLOW another symbol for S electrons <br> second mark depends on first |
|  |  | (ii) | $\left(\mathrm{NH}_{4}\right)_{2} \mathrm{~S}_{2} \mathrm{O}_{8} \checkmark$ | 1 | ALLOW <br> - with correct charges <br> - $\mathrm{S}_{2} \mathrm{O}_{8}\left(\mathrm{NH}_{4}\right)_{2}$ |
|  | (b) |  | $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}$ (is the stronger oxidising agent) because <br> $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}$ has a more positive $E^{9}$ value <br> indicates a greater tendency AW to gain/attract/accept electrons/to be reduced <br> ORA ie: <br> $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}$ (is the stronger oxidising agent) because $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}$ has a less positive / more negative $E^{\theta}$ value $\checkmark$ <br> indicates a smaller tendency AW to gain/attract/accept electrons/to be reduced <br> OR indicates a greater tendency AW to lose electrons/to be oxidised $\checkmark$ | 2 | without reference to oxidising strength of $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}$ only scores 1 as it doesn't answer the question <br> NOT 'higher' for 'more positive' both statements MUST BE comparative |
|  | (c) | (i) | $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}+2 \mathrm{I}^{-} \rightarrow 2 \mathrm{SO}_{4}{ }^{2-}+\mathrm{I}_{2}$ <br> correct equation balanced | 1 | IGNORE state symbols |


| Question | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: |
| (ii) | 1. use filter of complementary colour to iodine/solution $\checkmark$ | 6 | Using starch ANYWHERE is a CON and CANNOT gain marking point 3 ALLOW <br> - filter/wavelength giving maximum absorption/absorbance <br> - green/blue filter <br> - suitable filter |
|  | 2. zero colorimeter with water $\checkmark$ <br> 3. measure absorbance readings of standard solutions/solutions of known concentration (of iodine) |  | orange/yellow/brown is CON NOT 'solvent' instead of 'water' |
|  | 4. plot calibration graph $\checkmark$ <br> 5. take absorbance readings of the reacting mixture at known/certain times AW (must refer to time - may state units of time) |  |  |
|  | 6. convert absorbance readings to iodine concentrations using the calibration curve <br> QWC: <br> In order to gain the mark for point 3,5 or 6, absorbance must be used AND spelled correctly at least once |  | IGNORE references to quenching procedures |


| Question |  |  | Marks | Guidance |
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| (iii) |  | gnswer <br> decreasing from left to right (ANY decreasing <br> curve will do) |  |  |


| Question |  | Answer | Marks | Guidance |
| :---: | :---: | :---: | :---: | :---: |
|  | (iii) | 1. $\mathrm{Fe}^{3+}$ reacts with/oxidises $\mathrm{I}^{-}$ORA $\checkmark$ $2 \mathrm{Fe}^{3+}+2 \mathrm{I}^{-} \rightarrow 2 \mathrm{Fe}^{2+}+\mathrm{I}_{2} \checkmark$ <br> explanation: <br> the $E^{\circ}$ of $\mathrm{Fe}^{3+} / \mathrm{Fe}^{2+}$ half-cell is more positive / less negative than that of the $\mathrm{I}^{-1} \mathrm{I}_{2}$ half-cell ORA $\checkmark$ AW <br> 2. $\mathrm{Fe}^{2+}$ reacts with/reduces $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}$ ORA $\checkmark$ $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}+2 \mathrm{Fe}^{2+} \rightarrow 2 \mathrm{SO}_{4}{ }^{2-}+2 \mathrm{Fe}^{3+} \checkmark$ <br> explanation: <br> the $E^{\circ}$ of $\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-} / \mathrm{SO}_{4}{ }^{2-}$ half-cell is more positive than that of the $\mathrm{Fe}^{3+} / \mathrm{Fe}^{2+}$ half-cell ORA $\checkmark$ AW <br> QWC: to gain the explanation mark for either 1 or 2, the data has to be linked correctly to the reaction | 6 | IGNORE references to activation enthalpy OR variable valency/oxidation states <br> ALLOW $E^{\circ}$ of $\mathrm{Fe}^{3+}$ is more positive etc <br> ALLOW $E^{\circ}$ of $\mathbf{S}_{2} \mathrm{O}_{8}{ }^{2-}$ is more positive etc <br> NOTE <br> If answer starts with $\mathrm{Fe}^{2+}$ rather than $\mathrm{Fe}^{3+}$ lose $1^{\text {st }}$ mark but ecf since not answering question so can get 5 marks |
| (e) | (i) | Rate $=\mathrm{kx}\left[\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-}\right] \checkmark \times\left[\mathrm{I}^{-}\right] \checkmark$ | 2 | IGNORE state symbols |
|  | (ii) | uses one of the 3 sets of results: <br> 1. $\mathrm{k}=2.0 \times 10^{-5} /(0.075 \times 0.040)$ <br> 2. $k=4.0 \times 10^{-5} /(0.150 \times 0.040)$ <br> 3. $k=1.0 \times 10^{-5} /(0.075 \times 0.020) \checkmark$ $k=0.0067 \checkmark$ <br> $\mathrm{mol}^{-1} \mathrm{dm}^{3} \mathrm{~s}^{-1} \checkmark$ | 3 | ecf from part e(i) provided working is shown ALLOW standard form (eg $6.7 \times 10^{-3}$ ) Correct numerical answer without working scores 2 for calculation part <br> NO ecf from first mark to second <br> ALLOW any number of sig figs e.g. 0.007, 0.00667 NOT 0.006 NOR 0.0066 <br> ALLOW units in any order, e.g. $\mathrm{dm}^{3} \mathrm{~s}^{-1} \mathrm{~mol}^{-1}$ |
|  | (iii) | $\begin{aligned} & 2.0 \times 10^{-5} \checkmark \\ & \mathrm{~mol} \mathrm{dm}^{-3} \mathrm{~s}^{-1} \end{aligned}$ | 2 | Mark separately |
|  |  | Total | 31 |  |

