## Section A (multiple choice)

Question 1: N/A
Question 2: N/A
Question 3: N/A

| 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 |
| ---: | :--- | :--- | :--- |
| (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 |
| ---: | :--- | :--- | :--- |
| $\mathbf{1 3 ( a )}$ | D |  | $\mathbf{1}$ |
| (b) | B |  | $\mathbf{1}$ |
| (c) | A |  | $\mathbf{1}$ |

TOTAL FOR SECTION = 20 MARKS

## Section B

Question 14: N/A

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


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


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{aligned} & 15 \\ & (a)(i i i) \end{aligned}$ | $\begin{align*} & \text { Moles } \mathrm{S}_{2} \mathrm{O}_{3}{ }^{2-}=27.85 \times 10^{-3} \times 0.0631 \\ & \left(=1.757335 \times 10^{-3}\right) \\ & \text { moles of } \mathrm{I}_{2} \text { remaining }=\text { Moles } \mathrm{S}_{2} \mathrm{O}_{3}{ }^{2-} \div 2 \\ & \quad=27.85 \times 10^{-3} \times 0.0631 \div 2 \\ & \quad=8.786675 \times 10^{-4}=8.79 \times 10^{-4} \tag{1} \end{align*}$ <br> Moles ascorbic acid $=$ moles $I_{2}$ 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 <br> $100 \times$ 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 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) <br> OR <br> Using larger amount reduces the percentage <br> error / uncertainty in weighing <br> OR <br> Reverse discussion of two tablets <br> ALLOW <br> using four tablets gives a more <br> representative sampleJust 'reduces the <br> percentage error' | $\mathbf{1}$ |  |
| Titration value will <br> be larger (with four <br> tablets) so reduces <br> the percentage <br> error (in volume <br> measurement) |  |  |  |


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


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 15 (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 |  | $\mathbf{2}$ |
| IGNORE <br> IGtically active / inactive |  |  |  |
| 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' <br> C=O is broken | 1 |  |

Total for Q15 = 14 Marks

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


| 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 | $\begin{equation*} -\mathrm{COH} /-\mathrm{HCO} \tag{1} \end{equation*}$ <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> Step 3 <br> Lithium tetrahydridoaluminate((III)) / $\mathrm{LiAlH}_{4}$ <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> Marking Point 3 and 4 | 4 |  |
| Any two from |  |  |  |
| in phenol oxygen / OH group attached directly to |  |  |  |
| ring |  |  |  |
| Oxygen / OH group in phenylmethanol too far away |  |  |  |
| / not attached directly to ring |  |  |  |
| (In phenol) lone pair overlaps with the $\pi$ electrons |  |  |  |
| / delocalised electrons (of the ring) |  |  |  |
| ALLOW p orbital for lone pair for this mark |  |  |  |
| (2) |  |  |  |$\quad$| (1) |
| :--- |

Total for Q16 = 14 Marks

## Section C

| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(a)(i) | There is a barrier to rotation about a $(\mathrm{C}=\mathrm{C})$ bond <br> ALLOW restricted / limited / no rotation <br> Each carbon atom (in the $\mathrm{C}=\mathrm{C}$ double bond) has (two) different atoms / groups attached <br> IGNORE reference to priority groups | Just `molecule cannot rotate' & 2 \\ \hline Question Number & Acceptable Answers & Reject & Mark \\ \hline 17(a)(ii) & \begin{tabular}{l} There is a barrier to / restricted rotation about the ring \\ OR \\ The ring behaves like a double bond \end{tabular} & \begin{tabular}{l} Reference to benzene ring \\ Just `molecule cannot rotate' |  | \& 1 <br>

\hline
\end{tabular}

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( a ) ( i i i )}$ |  | 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 | $(\mathbf{1 )}$ |  |  |


| 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 |
| :--- | :--- | :--- | :--- |
| Number | Acceptable Answers $\quad$ 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 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( b ) ( i i i ) ~}$ | 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 |
| :--- | :--- | :--- | :--- |
| 17(b)(iv) | 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> $\mathbf{C H}$ (at d) and $\mathbf{C H}$ (at f) with different unique labels <br> Fourth mark <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 |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( c ) ( i i )}$ | $\mathrm{C}=\mathrm{O}$ amide (stretching vibrations are in the region) (1) <br> $1700-1630 \mathrm{~cm}^{-1}$ <br> $\mathrm{~N}-\mathrm{H}$ amide (stretching vibrations are in the region) (1) <br> $3500-3140 \mathrm{~cm}^{-1}$ <br> Amide only needs to be mentioned once but... | Amine (for <br> amide) | 2 |
|  | 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>  <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 |  |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 8}$ | B | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :---: |
| $\mathbf{1 9}$ | A | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Mark |
| :--- | :---: | :---: |
| $\mathbf{2 0}$ | D | $\mathbf{1}$ |

Section B

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 1}$ | Addition (1) <br> Nucleophilic (1) <br> Either order | SN1 <br> SN2 | $\mathbf{2}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| (a)(ii) | Hydrogen cyanide / HCN (1) <br> Potassium cyanide / KCN/ sodium cyanide/ <br> NaCN (1) <br> OR <br> Potassium cyanide / KCN (1) <br> With hydrochloric acid / sulfuric acid (to <br> generate HCN) (1) <br> Ignore concentration of acids <br> Mark for HCl etc is consequential on KCN <br> OR <br> Hydrogen cyanide / HCN (1) <br> With sodium hydroxide / other base (to make <br> cyanide ions) (1) <br> Mark for NaOH etc is consequential on HCN | Just CN- <br> Just CN ${ }^{-}$ <br> Just acid/ $\mathrm{H}^{+}$ any weak acid <br> Just $\mathrm{OH}^{-}$ | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{aligned} & 21 \\ & (\mathrm{a})(\mathrm{iii}) \end{aligned}$ |   <br> (1) <br> (1) <br> (1) <br> Both arrows in first step of mechanism above correctly drawn <br> (1) <br> Correct intermediate with charge <br> Both arrows in second step with correct organic product ( $\mathrm{CN}^{-}$is not required) (1) <br> Use of HCN for first step max 2 marks <br> Allow omission of lone pair on $\mathrm{CN}^{-}$and $\mathrm{O}^{-}$ Allow curly arrow from negative charge or elsewhere on cyanide ion <br> Allow arrow from $\mathrm{O}^{-}$in $2^{\text {nd }}$ step to $\mathrm{H}^{+}$(no other product or only one product) or $\mathrm{H}_{2} \mathrm{O}$ ( with $\mathrm{OH}^{-}$ formed) | $\mathrm{C}=\mathrm{O}$ breaking before attack by $\mathrm{CN}^{-}$ <br> Arrows from atoms when they should be from bonds and vice versa | 3 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| *21 <br> (a)(iv) | Attack (by nucleophile on the C) is from both <br> sides (equally)/ above and below (at the <br> planar reaction site in the aldehyde group) <br> (1) | Attack on <br> intermediate in <br> reaction <br> mechanism is <br> from both sides <br> Attack from both <br> ends/two angles | $\mathbf{2}$ |
|  | So a mixture of two <br> enantiomers/(optical)isomers in equal <br> proportions forms <br> OR <br> racemic mixture forms (1) <br> First and second marks are independent | Just "both <br> enantiomers form" |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (b) | Any named (aqueous) strong acid or its <br> formula. | Water | $\mathbf{1}$ |
|  | Allow <br> (aqueous) sodium hydroxide followed by <br> named acid or formula <br> Ignore references to concentration | Potassium <br> dichromate + <br> sulfuric acid <br> Carboxylic acids | $\mathbf{l}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (c)(i) | 2-hydroxypropanoic acid | 2- <br> hydroxylpropanoic <br> acid <br> 2- <br> hydroxopropanoic <br> acid <br> 2-hydroxypropan- <br> 1 -oic acid | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{gathered} 21 \\ (\mathrm{c})(\mathrm{ii}) \end{gathered}$ |  <br> OR <br> All bonds in ester link must be shown More than 2 units may be shown but structure shown should be a repeat unit Ignore brackets/n | A dimer <br> Missing H atoms <br> Missing bonds at ends | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (c)(iii) | Ester (link/bond) in PLA can be <br> hydrolysed/broken down (by enzymes) <br> OR Ester (link/bond) in PLA can be broken <br> down | Just "it can be <br> hydrolysed" | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (c)(iv) | Ethene is (from crude oil so) non-renewable/ <br> milk is from a renewable source/ <br> energy required to make ethene is high/ <br> high temperatures needed to make ethene/ <br> energy requirements for process from sour <br> milk less/ <br> process from milk doesn't use toxic <br> chemicals / process from milk doesn't use <br> cyanide | Milk is more readily <br> available <br> Greater atom <br> economy | $\mathbf{1}$ |
| Allow other chemicals <br> needed in process <br> from milk <br> process from ethene requires many steps so <br> expensive/so loss of material occurs at each <br> step /so more reagents needed | Just "process from <br> ethene requires <br> many steps" | Just "cheaper" |  |

## Section C

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- |
| $\mathbf{2 2 ( a )}$ | Alcohol; (2)-methylpropan-2-ol (1) | Formula of alcohol | $\mathbf{2}$ |
|  | Catalyst: sulfuric acid OR any named strong <br> acid Ignore concentration of acid (1) <br> Accept formula for acid | Just acid $/ \mathrm{H}^{+}$for <br> catalyst |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (b)(i) | Tap funnel / separating funnel | Buchner funnel <br> Filter funnel | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (b)(ii) | To neutralize / remove/ react with (excess) <br> acid | To purify it | $\mathbf{1}$ |
|  | Allow <br> To neutralize / remove / react with (excess) <br> $\mathrm{H}^{+}$ <br> To remove acidic impurities <br> To remove ethanoic acid <br> To remove the acid (used as a) catalyst <br> Ignore additional comments on quenching or <br> reaction stopping | To remove excess <br> acid and alcohol <br> Just "to quench <br> acid catalyst/stop <br> reaction" |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (b)(iii) | Add (anhydrous) calcium chloride/ sodium <br> sulfate/ magnesium sulfate/ <br> Allow silica gel <br> Allow formulae of drying agents | Conc. sulfuric acid <br> Anhydrous copper <br> sulphate <br> Just "silica" | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| (b)(iv) | Round bottomed or pear-shaped flask + still <br> head with stopper or thermometer + heat <br> source (1) <br> This mark cannot be given if apparatus is <br> completely sealed /large gaps between <br> components <br> Downwards sloping condenser (with correct <br> water flow) + collection vessel (1) <br> Thermometer in correct position with bulb <br> opposite condenser opening (1) | Conical flask <br> Flat bottomed flask | $\mathbf{3}$ |
| Ignore fractionating column if included <br> between flask and condenser |  |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| *22 (c) | First mark <br> (Two signals so) two hydrogen environments <br> (1) <br> This mark may be gained by a description of the only two environments, but reference to hydrogen must be made. <br> Second mark <br> (Numbers of hydrogen in each environment are/ are predicted to be) in ratio $3: 9$ or $1: 3$ <br> OR <br> Peak due to $\left(\mathrm{CH}_{3}\right)_{3}$ is $3 x$ higher than peak due to $\mathrm{CH}_{3}$ (1) <br> Third mark <br> Environments are $\mathrm{CH}_{3} \mathrm{COO}$ and $\left(\mathrm{CH}_{3}\right)_{3}$ <br> ( H may have been specified in first marking point) <br> These may be shown on a diagram of the formula of the molecule <br> OR <br> $\mathrm{H}-\mathrm{C}-\mathrm{C}=\mathrm{O}$ (peak at 2.1) and $\mathrm{H}-\mathrm{C}-\mathrm{C}$ (peak at 1.3) (1) <br> Fourth mark <br> Singlets/ no splitting as no H on adjacent C <br> OR <br> Singlets as the hydrogen environments are not adjacent to other H environments <br> Allow <br> "only one peak" for no splitting (1) | Just "the peaks are due to $\left(\mathrm{CH}_{3}\right)_{3}$ and $\mathrm{CH}_{3}$ | 4 |
| Question Number | Acceptable Answers | Reject | Mark |
| (d)(i) | $\mathrm{CH}_{3} \mathrm{COOCH}_{2} \mathrm{CH}\left(\mathrm{CH}_{3}\right)_{2}$ <br> Or correctly displayed <br> Allow $\mathrm{CH}_{3} \mathrm{COOCH}_{2} \mathrm{CH}\left(\mathrm{CH}_{3}\right) \mathrm{CH}_{3}$ |  | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 2}$ <br> (d)(ii) | The H on the $\mathrm{CH}_{3} \mathrm{COO}$ <br> Accept circle round all of first methyl group <br> of molecule is incorrect | Circle round C of <br> first methyl group | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| (e)(i) | Any acid with $6 \mathrm{C}(5 \mathrm{C}+\mathrm{COOH})$ which is | Infrared indicates O-H <br> Infrared indicates alkyl group | 5 |
|  | chiral, so will have a branched chain |  |  |
|  | $\mathrm{C}_{3} \mathrm{H}_{7} \mathrm{CH}\left(\mathrm{CH}_{3}\right) \mathrm{COOH}$ |  |  |
|  | OR |  |  |
|  | $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{CH}\left(\mathrm{CH}_{3}\right) \mathrm{CH}_{2} \mathrm{COOH}$ |  |  |
|  | $\begin{align*} & \mathrm{OR} \\ & \left(\mathrm{CH}_{3}\right)_{2} \mathrm{CHCH}\left(\mathrm{CH}_{3}\right) \mathrm{COOH} \tag{1} \end{align*}$ |  |  |
|  | Infrared indicates ( $\mathrm{O}-\mathrm{H}$ present in a) carboxylic acid (1) |  |  |
|  | High boiling temperature due to hydrogen bonding (between atoms in OH groups so not an ester.) Hydrogen bonds must be possible for structure shown |  |  |
|  | Allow acids can form dimers. |  |  |
|  | Allow TE from formula of straight chain molecule with explanation that London forces are higher in a linear molecule (1) |  |  |
|  | (Optically active so) contains chiral C/ C bonded to four different groups The formula suggested must contain a chiral carbon to score this mark |  |  |
|  | This may be shown by a chiral carbon being labelled in the formula (1) |  |  |
|  | Carbonyl compound/ Carbonyl group/ Aldehyde and ketone absent (as no reaction with 2,4-dinitrophenylhydrazine)/ Allow carboxylic acids do not react with 2,4dinitrophenylhydrazine/ <br> (1) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 2}$ | No because the isomers (which are <br> (earboxylic acids) contain same bonds / <br> groups (C=O, C-O, C-H etc) (1) | (1) <br> OR <br> Yes because could be distinguished by <br> infrared fingerprint (1) | Yes because <br> spectrum is unique |

