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


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1}$ <br> (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> rust "to quench <br> acid catalyst/stop <br> reaction" |  |


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


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1}$ | 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) | Conical flask <br> Flat bottomed flask | $\mathbf{3}$ |
| Thermometer in correct position with bulb <br> opposite condenser opening (1) | Ignore fractionating column if included <br> between flask and condenser |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| * 1 (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 $\left(\mathrm{CH}_{3}\right)_{3}$ ( 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 <br> 1.3) (1) <br> Fourth mark <br> Singlets/ no splitting as no H on adjacent C 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 <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1}$ | $\mathrm{CH}_{3} \mathrm{COOCH}_{2} \mathrm{CH}\left(\mathrm{CH}_{3}\right)_{2}$ |  | $\mathbf{1}$ |
| (d)(i) | Or correctly displayed |  |  |
|  | Allow $\mathrm{CH}_{3} \mathrm{COOCH}_{2} \mathrm{CH}\left(\mathrm{CH}_{3}\right) \mathrm{CH}_{3}$ |  |  |


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


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


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


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 ( a ) ( i )}$ | (Acid) hydrolysis | substitution | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 ~ ( a ) ( i i ) ~}$ | $\mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{Na}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}$ <br> Potassium dichromate((VI)) / sodium <br> dichromate((VI)) / dichromate((VI)) ions <br> ALLOW manganate((VII)) ions, etc | Just <br> "dichromate" <br> chromates | $\mathbf{1}$ |
| Correct |  |  |  |
| formula with |  |  |  |
| wrong name |  |  |  |
| and vice versa |  |  |  |
| Incorrect |  |  |  |
| oxidation |  |  |  |
| number |  |  |  |\(\quad\left\{\begin{array}{l} \\

\hline\end{array}\right.\)

| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2 (a)(iii) | Lithium tetrahydridoaluminate/ lithium aluminium hydride/ $\mathrm{LiAlH}_{4}$ (in dry ether) | J ust [ $\mathrm{H}^{-}$] | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 2 (a)(iv) | Methyl butanoate (1) <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{CH}_{3} \mathrm{OH} \rightarrow+$ <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{COOCH}_{3}+\mathrm{H}_{2} \mathrm{O}$ (1) <br> ALLOW $\rightleftharpoons$ <br> IGNORE state symbol even if wrong | Methyl <br> butoate | $\mathbf{2}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2 (a)(v) |  <br> Don't penalise undisplayed methyl groups as here. COCl must be displayed as above. | $\begin{aligned} & \mathrm{C}_{3} \mathrm{H}_{7} \text { for } \\ & \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \end{aligned}$ | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 ~ ( b ) ( i ) ~}$ | Nitrogen inert / unreactive / less reactive <br> (than oxygen) <br> OR <br> Oxygen might react with chemicals going <br> through column / sample might oxidise | $\mathbf{1}$ |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2 (b)(ii) | Solubility (in liquid / stationary phase) <br> OR <br> Interaction with liquid / stationary phase <br> OR <br> Interaction between mobile and stationary phase <br> OR <br> Attraction for liquid / stationary phase <br> OR <br> Strength of (named) intermolecular forces <br> OR <br> Adsorption on liquid / stationary phase <br> OR <br> Absorption on liquid / stationary phase | Size of molecule / molar mass <br> Polarity, unless with explanation <br> Boiling point / volatility <br> Viscosity <br> Attraction for carrier gas <br> Just a named intermolecular force <br> J ust 'retention time' <br> Density | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2 (c)(i) |  <br> OR <br> Ester link including $\mathrm{C}=0$ (1) <br> Rest of polymer with oxygens at end correct (1) <br> All H atoms must be shown. <br> PENALISE lack of displayed $\mathrm{C}=0$ once only ACCEPT <br> Without brackets around formula but bonds at end should be shown More than two correct units IGNORE n after brackets |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 2 (c)(ii) | Hydrolysis |  | $\mathbf{1}$ |
|  | OR <br> Splits / breaks ester link <br> OR <br> polymer breaks down to monomers <br> OR <br> equation showing hydrolysis | Just 'breaks <br> polymer down' |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 3(a)(i) | Sodium/potassium dichromate((VI))/potassium manganate ((VII))/ $\mathrm{Na}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{KMnO}_{4}$ <br> IGNORE references to acid | $\begin{aligned} & \text { Just } \mathrm{Cr}_{2} \mathrm{O}_{7}^{2-} \\ & / \mathrm{MnO}_{4}^{-} \end{aligned}$ | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{3 ( a ) ( i i )}$ | (Heat under) reflux (1) |  | $\mathbf{2}$ |
|  | Use excess/sufficient oxidizing agent/reagent <br> named in (a)(i), even if incorrect <br> IGNORE references to (excess) acid <br> Stand alone marks | (1) |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 3(a)(iii) | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CN} / \mathrm{C}_{2} \mathrm{H}_{5} \mathrm{CN}$ <br> ACCEPT displayed or skeletal formulae $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CN}+\mathrm{H}^{+}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{NH}_{4}^{+}$ <br> OR $\begin{equation*} \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CN}+\mathrm{HCl}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{NH}_{4} \mathrm{Cl} \tag{2} \end{equation*}$ <br> If equation is incorrect then presence of $\mathrm{H}^{+}$or acid in equation/or above arrow and water on LHS scores (1) <br> Mark cq on formula of nitrile <br> ALLOW one mark for the following equation without $\mathrm{H}^{+}$. $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CN}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{NH}_{3}$ <br> ALLOW two marks for either of the following with $\mathrm{H}^{+}$above the arrow $\begin{aligned} & \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CN}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{NH}_{3} \\ & \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CN}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{NH}_{4}^{+} \end{aligned}$ <br> ALLOW answers for alkaline hydrolysis followed by acidification $\begin{equation*} \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CN}+\mathrm{OH}^{-}+\mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}+\mathrm{NH}_{3} \tag{1} \end{equation*}$ <br> Then $\begin{equation*} \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}+\mathrm{H}^{+} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH} \tag{1} \end{equation*}$ <br> If propanamide, $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CONH}_{2}$ is given initially then ALLOW the two equation marks for the hydrolysis $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CONH}_{2}+\mathrm{H}^{+}+\mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+$ $\mathrm{NH}_{4}{ }^{+}$ <br> If no acid is used then only one mark $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CONH}_{2}+\mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{NH}_{3}$ | Hydroxynitriles | 3 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 3(b) | Reagent - Propanoyl chloride/ $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COCl}$ <br> Any two from: <br> $\mathrm{C}-\mathrm{Cl}$ bond is weaker (than $\mathrm{C}-0$ ) <br> $\mathrm{Cl}^{-} /$chloride (ion) is a better leaving group <br> Carbonyl carbon is more positive/more $\delta+/$ more attractive to nucleophiles <br> OR <br> Reagent - Propanoic anhydride/( $\left.\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CO}\right)_{2} \mathrm{O}$ <br> $\mathrm{CH}_{3} \mathrm{COO}^{-} /$propanoate (ion) is a better leaving group <br> Carbonyl carbon is more positive/more $\delta+/$ more attractive to nucleophiles <br> IGNORE references to eversible/equilibrium/ catalysts <br> IGNORE bond polarity | Propyl chloride <br> Just Cl is more electronegative | 3 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{3 ( c ) ( \mathbf { i } )}$ | Radio waves/radio frequency | Just radio | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |  |
| :--- | :--- | :--- | :--- | :--- |
| 3(c)(ii) | Any two from: |  | 2 |  |
|  | Protons/nuclei/they <br> have a property called spin/ <br> have a magnetic moment/ <br> have a magnetic field/ <br> are aligned with the external magnetic field | (1) | just dipole <br> moment | starts to spin |
| which flips/changes | (1) <br> align against the external magnetic field (when <br> radiation is absorbed) | polarity flips <br> any reference <br> to electrons or <br> molecules <br> scores zero |  |  |

$\left.\begin{array}{|l|l|l|l}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline \mathbf{3 ( c ) ( i i i )} & \begin{array}{l}\text { Quartet } \\ \text { ALLOW quadruplet/indication of four (peaks) (1) }\end{array} & & \mathbf{2} \\ & \begin{array}{l}\text { Value from 0.1 to } 1.9 \text { (ppm) inclusive } \\ \text { ACCEPT any range within the above range }\end{array} \quad \text { (1) }\end{array}\right) \quad$.

