



| Question | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: |
|  | Repeat units $n=10000 / 72=139$ <br> Equation <br> (1 mark) <br> Balanced equation for formation of P from $\mathrm{N} \checkmark$ e. |  | MUST be a whole number. <br> ALLOW 138 OR140 <br> For equation, ALLOW molecular OR structural OR skeletal OR displayed formulae OR mixture of the above e.g. ALLOW $n \mathrm{C}_{3} \mathrm{H}_{4} \mathrm{O}_{2} \longrightarrow\left(\mathrm{C}_{3} \mathrm{H}_{4} \mathrm{O}_{2}\right)_{n}$ <br> $n$ on LHS can be at any height to the left of formula AND $n$ on the RHS must be a subscript (essentially below the side link if displayed/skeletal formula is used) <br> ALLOW use of calculated value for $n$ in equation $\text { e.g. } 139 \mathrm{C}_{3} \mathrm{H}_{4} \mathrm{O}_{2} \longrightarrow\left(\mathrm{C}_{3} \mathrm{H}_{4} \mathrm{O}_{2}\right)_{139}$ |
|  | Total | 12 |  |






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| (d) | mole ratio C : H : O <br> $\frac{66.7}{12.0}: \frac{11.1}{1.0}: \frac{22.2}{16.0}$ OR $5.56: 11.1: 1.39$ $4: 8: 1 \mathrm{OR} \mathrm{C}_{4} \mathrm{H}_{8} \mathrm{O} \checkmark$ <br> contains a $\mathrm{C}=\mathrm{O}$ or carbonyl because of absorbance at about $1710 \mathrm{~cm}^{-1} \checkmark$ <br> Any two from: | 5 | PLEASE LOOK AT THE SPECTRA AND ABOVE THE SPECTRA FOR POSSIBLE ANSWERS <br> ALLOW two marks for $\begin{aligned} & 72 \times 66.7 / 100=48 / 12=4(\mathrm{C}) \\ & 72 \times 11.1 / 100=8=8(\mathrm{H}) \\ & 72 \times 22.2 / 100=16=1(\mathrm{O}) \end{aligned}$ <br> ALLOW C=O or carbonyl since has absorbance within the range 1640 to $1750 \mathrm{~cm}^{-1}$ <br> ALLOW ketone OR aldehyde linked to correct absorbance ALLOW 'could be aldehyde, ketone, carboxylic acid, ester (or amide) because of absorbance between range 1640 to $1750 \mathrm{~cm}^{-1}$, (ie direct quote from the data book) <br> DO NOT ALLOW reference to $M$ being a carboxylic acid, ester or amide unless they are included in a list with aldehyde/ketone in which case IGNORE carboxylic acid/ester/amide <br> IGNORE reference to $\mathrm{C}-\mathrm{O} /$ absence of $\mathrm{O}-\mathrm{H}$ DO NOT ALLOW has O—H <br> ALLOW correct structural OR displayed OR skeletal formula OR mixture of the above (as long as unambiguous) <br> eg $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CHO}, \mathrm{CH}_{3} \mathrm{COCH}_{2} \mathrm{CH}_{3}$ OR $\left(\mathrm{CH}_{3}\right)_{2} \mathrm{CHCHO}$ <br> DO NOT ALLOW $\mathrm{C}_{3} \mathrm{H}_{7} \mathrm{CHO}$ <br> IGNORE incorrect name <br> correct name on its own is not sufficient |
|  | Total | 11 |  |



| Question | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: |
|  | Compound Y <br> QWC Y contains O-H because of absorption between 3100 and $3500 \mathrm{~cm}^{-1}$ <br> QWC Mass spec of $\mathbf{Y}$ has molecular ion, $m / z=46$ so $M_{r}$ is 46 <br> Correct identification of one fragment from a $\mathrm{m} / \mathrm{z}$ value e.g. $\mathrm{m} / \mathrm{z}=31$ is $\mathrm{CH}_{2} \mathrm{OH}^{+} ; \mathrm{m} / \mathrm{z}=29$ is $\mathrm{C}_{2} \mathrm{H}_{5}^{+} ; \mathrm{m} / \mathrm{z}=15$ is $\mathrm{CH}_{3}{ }^{+}$ |  | ANNOTATE ANSWER WITH TICKS AND CROSSES <br> ALLOW $\mathbf{Y}$ is an alcohol (or phenol) because of absorption between 3200 and $3550 \mathrm{~cm}^{-1}$ <br> ALLOW Y contains $\mathrm{C}-\mathrm{O}, \mathrm{C}-\mathrm{H}$ and $\mathrm{O}-\mathrm{H}$ bonds because of absorptions at approximately 1030, 2950 and $3350 \mathrm{~cm}^{-1}$ <br> ALLOW $m / z=46$ so $M_{r}$ is 46 <br> OR mass spectrum has a peak at 46 which is the $M_{r}$ <br> OR $M_{\mathrm{r}}$ is 46 because of $m / z$ peak shown on the actual spectra <br> $M_{\mathrm{r}}=46$ on its own is not sufficient <br> $\mathrm{m} / \mathrm{z}=46$ on its own is not sufficient <br> ALLOW $m / z=31$ shows $\mathrm{CH}_{2} \mathrm{OH}$ (fragment); <br> $\mathrm{m} / \mathrm{z}=29$ shows $\mathrm{C}_{2} \mathrm{H}_{5}$ (fragment); <br> $\mathrm{m} / \mathrm{z}=15$ is $\mathrm{CH}_{3}$ (fragment) |
|  | Identification of compounds <br> So $\mathbf{X}$ must be $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}$ OR propanoic acid $\checkmark$ <br> So $\mathbf{Y}$ is ethanol $\mathrm{OR} \mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OH}$ OR $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH} \checkmark$ <br> Z is $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOC}_{2} \mathrm{H}_{5}$ OR ethyl propanoate $\checkmark$ | 3 | Note: an incorrect name CONs a correct structure ALLOW skeletal OR displayed formula throughout <br> DO NOT ALLOW propanoic acid with wrong structure or incorrect molecular formula <br> DO NOT ALLOW ethanol with wrong structure or incorrect molecular formula <br> DO NOT ALLOW ethyl propanoate with wrong structure or incorrect molecular formula <br> ALLOW ECF for identification of $\mathbf{Z}$ from incorrect $\mathbf{X}$ and $\mathbf{Y}$. DO NOT ALLOW this ECF if name and structures of $\mathbf{X}$ or $\mathbf{Y}$ do not match |
|  | Total | 10 |  |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | a | i | Any two from: <br> Any value between 1000-1300 $\checkmark$ <br> Any value between 2850-3100 $\checkmark$ <br> Any value between 3200-3550 $\checkmark$ | 2 |  |
|  |  | ii | Orange to green or blue $\checkmark$ | 1 |  |
|  |  | iii | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH}+[\mathrm{O}] \rightarrow \mathrm{CH}_{3} \mathrm{CHO}+\mathrm{H}_{2} \mathrm{O}$ <br> OR $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH}+2[\mathrm{O}] \rightarrow \mathrm{CH}_{3} \mathrm{COOH}+\mathrm{H}_{2} \mathrm{O}$ <br> Correct organic product $\checkmark$ Balanced equation $\checkmark$ | 2 | IGNORE any state symbols <br> ALLOW $\mathrm{CH}_{3} \mathrm{COH}$ in equation but not for the structure <br> ALLOW equations with molecular formulae but not the product mark |
|  | b | i | Absorption around 2850-3100 ( $\mathrm{cm}^{-1}$ ) so contains C-H bonds <br> No other important absorptions present / no other characteristic absorptions $\checkmark$ | 2 | Answer must have a reference to infrared spectrum i.e. use of $\mathrm{cm}^{-1}$ or data from the infrared spectrum <br> 'Has no other peaks so no functional groups present' is not sufficient <br> BUT <br> There are no peaks due to functional groups is sufficient <br> ALLOW peaks instead of absorption ALLOW no absorption due to $\mathrm{C}=\mathrm{O}$ and $\mathrm{O}-\mathrm{H} /$ no absorption due to carbonyl and hydroxyl |
|  |  | ii | Peak furthest to right hand side is 58 / molecular ion peak is 58 / peak at highest mass $\checkmark$ | 1 | ALLOW peak at $m / z 58$ marked on the mass spectrum / M peak is 58 / peak at 58 linked to the molecular mass <br> DO NOT ALLOW highest peak but ALLOW 58 is the highest peak |


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| b | iii |   <br> BOTH isomers correct $\checkmark$ | 1 | If three structures are drawn then do not award mark ALLOW skeletal formulae / structural formulae IGNORE incorrect names |
|  | iv | $\begin{aligned} & \mathrm{CH}_{3}^{+} \checkmark \\ & \mathrm{C}_{2} \mathrm{H}_{5}^{+} \downarrow \\ & \mathrm{C}_{3} \mathrm{H}_{7}^{+} / \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2}^{+} /\left(\mathrm{CH}_{3}\right)_{2} \mathrm{CH}^{+} \downarrow \end{aligned}$ | 3 | Essentially marks are allocated as positive ions <br> Formula of two fragments correct (ignore charge) $\checkmark$ <br> BUT <br> formulae of all three fragments correct (ignore charge) |
|  | V | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3}$ because there is a peak at $m / z=29 \checkmark$ | 1 | ALLOW name, displayed or skeletal structure ALLOW butane because there is a $\mathrm{C}_{2} \mathrm{H}_{5}$ fragment ALLOW butane because it gives all three fragments listed in (iv) |
|  |  | Total | 13 |  |

