## GCE

## Mark Scheme

| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | a | i | Series having same functional group and a general formula | 1 | ALLOW same functional group and members vary by $\mathrm{CH}_{2}$ <br> ALLOW organic compounds with the same functional group that differ in length of their hydrocarbon chain |
|  |  | ii | More surface contact OR bigger molecules <br> More van der Waals' forces | 2 | BOTH answers need to be comparisons <br> ALLOW higher relative formula mass OR has more electrons OR longer chain length OR more carbon atoms <br> IGNORE surface area / bigger compounds <br> ALLOW stronger van der Waals' forces / stronger induced dipoles <br> VDW forces is not sufficient <br> More intermolecular forces is not sufficient <br> DO NOT ALLOW breaking bonds within the chain / breaking covalent bonds <br> IGNORE reference to bonds if not linked to covalent bonds |
|  | b | i | Pent-1-yne OR pent-2-yne $\checkmark$ | 1 | ALLOW pentyne <br> Look for answer in the table if not on answer line but answer line takes precedence |
|  |  | ii | $\mathrm{C}_{n} \mathrm{H}_{2 n-2}{ }^{\checkmark}$ | 1 | ALLOW $\mathrm{C}_{n} \mathrm{H}_{2(n-1)}$ |


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| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | b | iii | Correct displayed formula $\checkmark$ | 1 |  |
|  |  | iv | Correct skeletal formula of cyclic hydrocarbon with formula $\mathrm{C}_{6} \mathrm{H}_{10} \checkmark$ | 1 |  |
|  | c |  | Energy required to break bonds $=(+) 2912 \checkmark$ <br> Energy released to make bonds $=(-) 4148 \checkmark$ <br> Enthalpy of combustion $=-1236 \checkmark$ | 3 | ALLOW full marks for correct answer with no working out <br> ALLOW $(2 \times 415)+(837)+(2.5 \times 498)$ $\begin{aligned} & \text { ALLOW }(4 \times-805)+(2 \times-464) \\ & \text { OR }(4 \times 805)+(2 \times 464) \end{aligned}$ <br> ALLOW ECF for calculation of enthalpy of combustion ALLOW 2 marks for +1236 with no working out |

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| 1 | d | i | (Enthalpy change) when one mole of a compound $\checkmark$ <br> is made from its elements (in their standard states) <br> (Standard conditions are) 298 K and 100 kPa | 3 | IGNORE energy required / energy released ALLOW (energy change) when one mole of a substance <br> DO NOT ALLOW enthalpy change for one mole of products <br> ALLOW 1 atmosphere pressure / $101 \mathrm{kPa} / 10^{5} \mathrm{~Pa} /$ $1.01 \times 10^{5} \mathrm{Nm}^{-2} / 1000$ millibars / $25^{\circ} \mathrm{C} /$ any stated temperature in words <br> IGNORE $1 \mathrm{~mol} \mathrm{dm}^{-3}$ for solutions |
|  |  | ii | From energy cycle <br> Enthalpy change to get elements $=-(-60)-(2-286) /(+)$ <br> $632 \checkmark$ <br> Enthalpy change from elements $=-987+(+227) /(-) 760 \checkmark$ <br> Enthalpy change $=-128 \checkmark$ | 3 | ALLOW full marks for -128 with no working out <br> ALLOW ECF from errors in calculation <br> ALLOW two marks for answer of $-414 /+128 /-1392$ / +1392 <br> ALLOW one mark for answer of +414 |
|  | e | i | $\begin{aligned} & \frac{26.0}{100.1} \times 100 \\ & 26.0 \% \checkmark \end{aligned}$ | 2 | First mark for 100.1 OR (64.1 + 36.0) OR (74.1 + 26.0) at bottom of fraction with or without $\times 100$ <br> ALLOW full marks for 26.0 or $26 \%$ with no working out <br> ALLOW from two significant figures up to calculator value <br> ALLOW 25.97 / 26\% <br> NO ECF for this part from incorrect numbers in first expression |


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| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | e | ii | $1.56 \times 10^{4}$ OR 15600 OR $15601 \checkmark$ | 1 | ALLOW calculator value of 15600.62402 and any rounded value to a minimum of three significant figures |
|  |  | iii | $1.5 \times 10^{4}$ OR $15000 \checkmark$ | 1 | ALLOW $1.50 \times 10^{4}$ etc. |
|  |  | iv | $96.2 \checkmark$ | 1 | ALLOW ECF from (iii) - (ii) ALLOW calculator value 96.1538461 and any rounded value to a minimum of two significant figures ALLOW 96.14768284 if 15601 is used <br> ALLOW any value between 88 to 89 if answer to (iii) was calculated by dividing by 26 |
|  |  | v | Any two from: <br> Low atom economy gives a poor sustainability OR low atom economy means lots of waste <br> A use for the aqueous calcium hydroxide needs to be developed to increase atom economy <br> Alternative process needs to be developed with high atom economy | 2 | ANNOTATE WITH TICKS AND CROSSES IGNORE comments about percentage yield <br> ALLOW ECF from (i) e.g. high atom economy will have good sustainability <br> ALLOW find a use for the waste to increase atom economy |
|  |  |  | Total | 23 |  |


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| :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | a | i | Branched chain alkane of formula $\mathrm{C}_{5} \mathrm{H}_{12}$ to $\mathrm{C}_{9} \mathrm{H}_{20}$ e.g. 2-methylpentane, 3-methyloctane | 1 | Must have position number but ALLOW methylbutane DO NOT ALLOW 1-methylpentane or 2-ethylpentane etc <br> DO NOT ALLOW incorrect nomenclature e.g. <br> 2-methypentane etc |
|  | b | i | Vibrate (more) $\checkmark$ | 1 | ALLOW bend / stretch / oscillate IGNORE rotate NOT break / molecules vibrate |
|  |  | ii | Incomplete combustion $\checkmark$ | 1 | ALLOW not enough oxygen |
|  |  | iii | NO for photochemical smog OR low level ozone $\checkmark$ <br> CO is toxic | 2 | ALLOW NO can (eventually) cause acid rain OR can result in respiratory irritation OR can (eventually) depletes high level ozone OR depletes ozone layer IGNORE greenhouse gas <br> ALLOW poisonous OR kills OR lethal ALLOW CO reduces the capacity of blood to carry oxygen Oxygen combines with haemoglobin is insufficient <br> IGNORE CO is harmful / suffocates / greenhouse gas |
|  | C | i | Makes nitrogen AND carbon dioxide $\checkmark$ $2 \mathrm{CO}+2 \mathrm{NO} \rightarrow \mathrm{~N}_{2}+2 \mathrm{CO}_{2} \checkmark$ | 2 | ALLOW any correct multiples IGNORE state symbols |

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| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | b | i | Nucleophilic substitution $\checkmark$ <br> Heterolytic <br> Dipole shown on $\mathrm{C}-\mathrm{I}$ bond, $\mathrm{C}^{\delta+}$ and $\mathrm{I}^{\delta-}$ <br> Curly arrow from $\mathrm{OH}^{-}$to carbon atom of $\mathrm{C}-\mathrm{I}$ bond $\checkmark$ <br> Curly arrow from C-I bond to the iodine atom $\checkmark$ | 5 | ANNOTATE WITH TICKS AND CROSSES <br> DO NOT ALLOW fish hooks <br> No need to show lone pair on $\mathrm{OH}^{-}$or $\mathrm{I}^{-}$ Curly arrow must come from the negative sign or lone pair on the oxygen of the hydroxide ion <br> ALLOW $\mathrm{S}_{\mathrm{N}} 1$ mechanism <br> dipole shown on $\mathrm{C}-\mathrm{I}$ bond, $\mathrm{C}^{\delta+}$ and $\mathrm{I}^{\delta-} \checkmark$ <br> curly arrow from C-I bond to the iodine atom $\checkmark$ <br> curly arrow from $\mathrm{OH}^{-}$to correct carbonium ion $\checkmark$ |
|  |  | ii | Use reflux OR heat for more than 20 minutes $\checkmark$ <br> $\mathrm{C}-\mathrm{Cl}$ stronger bond (than $\mathrm{C}-\mathrm{I}$ bond) OR $\mathrm{C}-\mathrm{Cl}$ shorter bond (than C-I bond) OR C-Cl bond is harder to break OR needs more energy to break $\mathrm{C}-\mathrm{Cl}$ bond OR ora $\checkmark$ | 2 | ALLOW heat stronger OR heat for longer OR heat at a higher temperature OR more heat <br> Answer must refer to the $\mathrm{C}-\mathrm{Cl}$ bond or $\mathrm{C}-\mathrm{I}$ bonds |
|  |  |  | Total | 11 |  |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | 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 <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{OH}+2[\mathrm{O}] \rightarrow \mathrm{CH}_{3} \mathrm{COOH}+\mathrm{H}_{2} \mathrm{O}$ <br> Correct organic product $\checkmark$ <br> 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 $\checkmark$ <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 |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | 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 | $\begin{aligned} & \mathrm{CH}_{3}^{+} \checkmark \\ & \mathrm{C}_{2} \mathrm{H}_{5}^{+} \checkmark \\ & \mathrm{C}_{3} \mathrm{H}_{7}^{+} / \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2}^{+} /\left(\mathrm{CH}_{3}\right)_{2} \mathrm{CH}^{+} \checkmark \end{aligned}$ | 3 | Essentially marks are allocated as positive ions $\checkmark$ <br> Formula of two fragments correct (ignore charge) $\checkmark$ BUT <br> formulae of all three fragments correct (ignore charge) $\checkmark \checkmark$ |
|  |  | V | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3}$ because there is a peak at $\mathrm{m} / \mathrm{z}=29 \checkmark$ | 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 |  |

Mark Scheme

| Question |  | Expected Answers | Marks | Additional Guidance <br> $\mathbf{5}$ $\mathbf{a}$ | Sideways overlap of two $p$ orbitals on each carbon atom $\checkmark$ <br> forms m-orbital or m-bond above and below plane of <br> molecule $\checkmark$ |
| :--- | :--- | :--- | :--- | :--- | :--- |
| see additional page with typical diagrams you might |  |  |  |  |  |
| see |  |  |  |  |  |

Each of the following diagrams is worth one mark. The words p-orbitals must be present to score the mark


One
p-orbitals

p-orbitals



p-orbitals

Each of the diagrams on its own scores no mark

p-orbitals

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| :--- | :--- | :--- | :--- | :---: | :---: |
| $\mathbf{5}$ | $\mathbf{b}$ | $\mathbf{i}$ | Double bond does not rotate / restricted rotation of the <br> double bond $\checkmark$ <br> Each carbon atom of double bond is bonded to (two) <br> different groups $\checkmark$ | 2 | ALLOW m bond does not rotate |
|  |  | ii | C and E $\checkmark$ | 1 | ALLOW each carbon atom of double bond is bonded <br> to (two) different atoms / each end of the $\pi-b o n d ~ i s ~$ <br> bonded to different groups or atoms $\checkmark$ |

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| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | e | i |  | 1 | Must have at least two repeat units and the free bonds at the end <br> All carbon-carbon bonds in the polymer chain must be shown <br> ALLOW bond to ethyl group to any part of ethyl group <br> IGNORE any brackets drawn |
|  |  | ii | Poly(but-1-ene) $\checkmark$ | 1 | ALLOW polybut-1-ene <br> n.b. the bracket is part of the answer <br> DO NOT ALLOW polybutene |
|  | f | i | (Lots of) OH group present <br> Can form hydrogen bonds with water $\checkmark$ | 2 | ALLOW hydroxyl group present / hydroxy group Alcohol group is not sufficient |
|  |  | ii | Any two from: <br> Incineration to produce energy OR combustion to produce energy $\checkmark$ <br> Sorting and recycling OR sorting and remoulding $\checkmark$ <br> Cracked (to give monomers) OR as an organic feedstock | 2 | Used as a fuel is not sufficient <br> IGNORE use photodegradable or biodegradable polymers |
|  |  |  | Total | 21 |  |

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| :---: | :---: | :---: | :---: | :---: | :---: |
| 6 | a |  | Low pressure because more (gas) molecules on right hand side of equation OR low pressure because $\Delta V=$ positive $\checkmark$ <br> Low temperature because the (forward) reaction is exothermic | 2 | ALLOW low pressure because more (gas) moles on right hand side of equation |
|  | b |  | Increased pressure speeds up reaction / ora $\checkmark$ $900^{\circ} \mathrm{C}$ increases the rate OR increased temperature speeds up reaction / ora <br> Idea that high enough temperature without compromising yield OR idea that high enough pressure without compromising yield | 3 | ANNOTATE WITH TICKS AND CROSSES ALLOW 'pushes gases through system' |
|  | C | i | $5.68 \times 10^{7} / 5.7 \times 10^{7} \checkmark$ | 1 | ALLOW two or more significant figures Calculator answer is $5.6812500 \times 10^{7}$ |
|  |  | ii | Used to heat the incoming gases $\checkmark$ | 1 | ALLOW used to heat rest of factory OR sold to the national grid <br> Provide energy to create conditions is not sufficient because one condition is pressure |
|  |  |  | Total | 7 |  |



