| 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 $\mathrm{C}_{n} \mathrm{H}_{2(n-1)}$ |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 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 $(2 \times 415)+(837)+(2.5 \times 498)$ <br> ALLOW $(4 \times-805)+(2 \times-464)$ <br> OR $(4 \times 805)+(2 \times 464)$ <br> ALLOW ECF for calculation of enthalpy of combustion ALLOW 2 marks for +1236 with no working out |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 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 $=-(-60)-(2-286) /(+)$ $632 \checkmark$ <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 |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 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 $1.50 \times 10^{4} \mathrm{etc}$. |
|  |  | 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 |  |


| Question |  | Expected Answers | Marks | Additional Guidance |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\mathbf{2}$ | Sideways overlap of two porbitals on each carbon atom $\checkmark$ |  |  |  |
| forms m-orbital or m-bond above and below plane of |  |  |  |  |
| molecule $\checkmark$ |  |  |  |  |

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

p-orbitals

p-orbitals

p-orbitals


p-orbital

Each of the diagrams on its own scores no mark

p-orbitals

| Question |  | Expected Answers | Marks | Additional Guidance |  |
| :--- | :--- | :--- | :--- | :---: | :--- |
| $\mathbf{2}$ | b | 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 $\pi$ 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 $\pi$-bond is <br> bonded to different groups or atoms $\checkmark$ |



| Question |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 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 n.b. the bracket is part of the answer DO NOT ALLOW polybutene |
| f | i | (Lots of) OH group present <br> Can form hydrogen bonds with water | 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 <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 |  |


| Question |  |  | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) |  | $\mathrm{C}_{n} \mathrm{H}_{2 n+2} \checkmark$ | 1 | ALLOW $\mathrm{C}_{n} \mathrm{H}_{2(n+1)} \checkmark$ <br> IGNORE size of subscripts |
|  | (b) | (i) | $\mathrm{C}_{8} \mathrm{H}_{18}+8{ }^{1 / 2} \mathrm{O}_{2} \rightarrow 8 \mathrm{CO}+9 \mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 | ALLOW any correct multiples IGNORE state symbols |
|  |  | (ii) | limited supply of air OR not enough $\mathrm{O}_{2} \checkmark$ | 1 | ALLOW use of air or oxygen <br> IGNORE it is not completely oxidised |
|  | (c) | (i) | $2 \mathrm{CO}+2 \mathrm{NO} \rightarrow 2 \mathrm{CO}_{2}+\mathrm{N}_{2} \checkmark$ | 1 | ALLOW any correct multiples including fractions IGNORE state symbols |
|  | (c) | (ii) | CO and NO are adsorbed (onto surface) OR reactants are adsorbed (onto surface) <br> weakening of bonds OR lowers activation energy <br> $\mathrm{CO}_{2}$ and $\mathrm{N}_{2}$ desorbs (from the surface) OR products desorbs (from the surface) | 3 | ALLOW CO and NO stick onto surface OR CO and NO form weak attractions to the surface OR gases are adsorbed onto surface <br> NOT absorb but allow ecf for deabsorb later on <br> IGNORE alternative pathway <br> Requires less energy is not sufficient <br> ALLOW products leave the surface OR products diffuse away from surface OR weak attraction to surface is broken <br> ALLOW deadsorb |
|  | (d) |  | skeletal formula of a branched isomer of $\mathrm{C}_{8} \mathrm{H}_{18} \checkmark$ <br> skeletal formula of a cyclic hydrocarbon OR skeletal formula of substituted arene of $\mathrm{C}_{8} \mathrm{H}_{10}$ v | 2 | ALLOW any ring between $\mathrm{C}_{3}$ and $\mathrm{C}_{8}$ with 8 carbon atoms per molecule <br> IGNORE wrong names <br> If two correct structural or displayed formulae drawn award one mark |


| Questi | Expected Answers | Marks | Additional Guidance |
| :---: | :---: | :---: | :---: |
| (e) | Any TWO from: atmospheric concentration ability to absorb infrared radiation residence time | 2 | ALLOW the amount of the gas OR abundance of gas <br> ALLOW how much IR it absorbs OR ability to absorb heat <br> IGNORE global warming potential / heat reflected / how much is produced <br> ALLOW how long it stays in the atmosphere |
|  | Any TWO from: <br> deep in the oceans OR on the sea-bed <br> storage in geological formations OR under the sea-bed $\checkmark$ by reaction (with metal oxides) to form carbonates $\checkmark$ | 2 | ALLOW piped into disused or partially filled oil wells <br> ALLOW stored as a carbonate OR equation to show formation of suitable carbonate from an oxide IGNORE mineral storage <br> IGNORE reforestation |
|  | Total | 13 |  |


| Question |  | Expected Answers | Marks | Rationale |
| :---: | :---: | :---: | :---: | :---: |
| 4 | (a) | one mark for each correct structure $\checkmark \checkmark \checkmark \checkmark$ | 4 | ALLOW skeletal formula OR displayed formulae IGNORE molecular formulae <br> IF two answers given e.g. name and structure then both must be correct to be given a mark <br> ALLOW methylpropane OR $\left(\mathrm{CH}_{3}\right)_{3} \mathrm{CH} \checkmark$ <br> ALLOW 1,2-dibromo-methylpropane OR $\mathrm{CH}_{2} \mathrm{BrCBr}\left(\mathrm{CH}_{3}\right)_{2} \checkmark$ <br> ALLOW 1-bromo-methylpropane $\mathbf{O R} \mathrm{CH}_{2} \mathrm{BrCH}\left(\mathrm{CH}_{3}\right)_{2}$ <br> ALLOW 2-bromo-methylpropane OR $\mathrm{CH}_{3} \mathrm{CBr}\left(\mathrm{CH}_{3}\right)_{2}$ <br> ALLOW ecf if wrong carbon skeleton is used in all of the structures mark first structure wrong and then apply ecf for the rest |
|  | (b) | curly arrow from double bond to $\mathrm{Br}^{\delta+}$ and curly arrow from $\mathrm{Br}-$ Br bond pair to $\mathrm{Br}^{\overline{ }-}$ in 1st step <br> curly arrow in 2nd step from bromide ion $\checkmark$ <br> correct dipole shown on $\mathrm{Br}_{2}$ <br> correct carbocation shown $\checkmark$ | 4 | Curly arrow must start from the double bond and not a carbon atom, other curly arrow must start from $\mathrm{Br}-\mathrm{Br}$ bond <br> ALLOW curly arrow from any part of bromide ion The bromide ion does not need to show a lone pair <br> Dipole must be partial charge and not full charge Carbocation needs a full charge and not a partial charge (charges do not need to be surrounded by a circle) <br> ALLOW carbocation on carbon 1 where electrophile attacks carbon 2 i.e. ${ }^{+} \mathrm{CH}_{2} \mathrm{CBr}\left(\mathrm{CH}_{3}\right)_{2}$ |


| Question |  | Expected Answers | Marks | Rationale |
| :---: | :---: | :---: | :---: | :---: |
| (c) | (i) | $\mathrm{C}_{6} \mathrm{H}_{10} \checkmark$ | 1 |  |
|  | (ii) | ```Mr(cyclohexanol)}=100 amount of cyclohexanol = 0.0765 mol percentage yield =35.0% ``` | 3 | ALLOW full marks for correct answer with no or limited working out <br> ALLOW ecf from wrong molar mass i.e. $7.65 \div$ molar mass <br> ALLOW ecf from wrong amount in moles i.e. [0.0268 $\div$ moles] $\times 100$ <br> ALLOW 35\% <br> ALLOW two marks for $0.35 \%$ <br> If $M_{r}$ of 82 is used then $\%$ yield will be 28.7 or 29 and this is worth two marks |
| (d) | (i) | (sum of) the molecular masses of the desired product :sum of molecular masses of all products $\times 100 \checkmark$ | 1 | ALLOW (sum of) the molecular masses of the desired product $\div$ <br> sum of molecular masses of all reactants $\times 100 \checkmark$ |
|  | (ii) | this preparation is addition OR has 100\% atom economy OR there is only one product $\checkmark$ <br> preparation from cyclohexanol has less than $100 \%$ atom economy OR $\mathrm{H}_{2} \mathrm{O}$ is produced as well OR calculated atom economy $=82 \% \checkmark$ | 2 | ALLOW no by products formed <br> ALLOW other substances formed OR cyclohexene is not the only product |
|  |  | Total | 15 |  |

