## A Level OCR Chemistry

## Chapter 10 - answers

| Question | Answers | Extra information | Mark | AO Spec reference |
| :---: | :---: | :---: | :---: | :---: |
| 1(a) | A compound made up of hydrogen and carbon only. |  | 1 | $\begin{gathered} \text { AO1 } \\ \text { 4.1.1a } \end{gathered}$ |
| 1(b) | Because it is an alkane it has the general formula $\mathrm{C}_{\mathrm{n}} \mathrm{H}_{2 \mathrm{n}+2}$. Its molecular formula is $\mathrm{C}_{5} \mathrm{H}_{12}$. <br> Relative molecular mass $=5 \times 12+12 \times 1=72$ |  | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO2 } \\ \text { 2.1.1e } \end{gathered}$ |
| 1(c) | Isomers have the same molecular formula but different structural formulae. |  | 1 | $\begin{gathered} \text { AO1 } \\ \text { 4.1.1e } \end{gathered}$ |
| 1(d) | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3}$ pentane $\mathrm{CH}_{3} \mathrm{CH}\left(\mathrm{CH}_{3}\right) \mathrm{CH}_{2} \mathrm{CH}_{3}$ 2-methylbutane $\mathrm{C}\left(\mathrm{CH}_{3}\right)_{4}$ or $\left(\mathrm{CH}_{3}\right)_{4} \mathrm{C}$ 2,2-dimethylpropane | Both formula and name required for each mark. | $1$ <br> 1 <br> 1 | $\begin{gathered} \text { AO2 } \\ \text { 4.1.1a } \end{gathered}$ |
| 1(e) | $0.36 \mathrm{~g}=0.36 / 72 \mathrm{~mol}=5 \times 10^{-3} \mathrm{~mol}$ <br> Number of molecules $=5 \times 10^{-3} \times 6.022 \times 10^{23}=3.01 \times 10^{21}$ <br> In each molecule there are 17 atoms. Total number of atoms $=3.01 \times 10^{21} \times 17=5.12 \times 10^{22}$ | Just the answer with no reasoning shown is $\mathbf{1}$ mark | 1 <br> 1 <br> 1 | $\begin{gathered} \text { AO2 } \\ \text { 2.3.1a } \\ \text { 2.3.1a } \\ \text { AO3 } \\ \text { 2.3.1a } \end{gathered}$ |
| 2(a) | I pent-1-ene <br> II 2-methylbut-2-ene <br> III pent-2-ene | Do not allow pentene | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO2 } \\ \text { 4.1.1a } \end{gathered}$ |
| 2(b)(i) | Stereoisomers have the same structural formula Have different arrangement of bonds in space |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO1 } \\ 4.1 .3 \mathrm{C} \end{gathered}$ |

## A Level OCR Chemistry

## Chapter 10 - answers

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| 2(b)(ii) | Isomer III |  | $1$ <br> 1 | $\begin{gathered} \text { AO2 } \\ 4.1 .3 \mathrm{C} \end{gathered}$ |
| 2(b)(iii) | There is no free rotation about the $\mathrm{C}=\mathrm{C}$ double bond. On each carbon of the $\mathrm{C}=\mathrm{C}$ bond there are 2 different atoms or groups |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO1 } \\ 4.1 .3 \mathrm{C} \end{gathered}$ |
| 2(c) | Methylcyclobutane |  | 1 | $\begin{gathered} \mathrm{AO} 2 \\ 4.1 .1 \mathrm{c}(\mathrm{v}) \end{gathered}$ |
| 3(a) | Use the ideal gas equation $\begin{aligned} & n=\frac{p V}{R T} \\ & n=\frac{1.01 \times 10^{5} \times 85 \times 10^{-6}}{8.31 \times 450} \\ & =2.30 \times 10^{-3} \mathrm{~mol} \\ & M_{\mathrm{r}}=\mathrm{m} / \mathrm{n}=0.20 / 2.30 \times 10^{-3}=87(.1) \mathrm{g} \mathrm{~mol}^{-1} \end{aligned}$ | Some evidence of using the equation is required | 1 <br> 1 <br> 1 $1$ | $\begin{gathered} \mathrm{AO2} \\ 4.2 .4 \mathrm{f} \\ 2.1 .3 \mathrm{f} \end{gathered}$ |
| 3(b) | The last /highest value significant peak |  | 1 | $\begin{aligned} & \text { AO1 } \\ & \text { 4.2.4f } \end{aligned}$ |
| 3(c) | $\begin{aligned} & \text { Error }=100 \% \times\|88-87\| / 88 \\ & 1.14 \% \end{aligned}$ |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \mathrm{AO2} \\ \text { 1.1.4d } \end{gathered}$ |

## A Level OCR Chemistry

## Chapter 10 - answers

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| 3(d)(i) | 2 marks for $1^{\text {st }}$ (left-hand) diagram <br> 1 mark for $2^{\text {nd }}$ diagram | Give both marks if as shown Lose 1 mark if give $\mathrm{CH}_{3}$ on CH but does not show all hydrogens <br> All bonds must be shown | $2$ <br> 1 | $\begin{gathered} \text { AO3 } \\ \text { 4.1.1b } \end{gathered}$ |
| 3(d)ii) | $\begin{aligned} & a=109.5 \\ & b=104.5^{\circ} \end{aligned}$ |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO1 } \\ \text { 4.1.1b } \end{gathered}$ |
| 4(a) | Chlorine is more electronegative than carbon electrons are not shared equally / the electrons are attracted more by the chlorine atom electron cloud distorted towards chlorine |  | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO2 } \\ \text { 2.2.2j } \end{gathered}$ |
| 4(b)(i) | The molecule is tetrahedral and symmetrical The electrons (electron clouds) are symmetrically distributed so the polar bonds cancelout OR not asymmetrically distributed |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO2 } \\ \text { 2.2.2j } \end{gathered}$ |

## A Level OCR Chemistry

## Chapter 10 - answers

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| 4(b)(ii) | 1 mark for atoms and bonds 1 for dipole <br> $\delta-$ |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO3 } \\ \text { 2.2.2j } \end{gathered}$ |
| 4(c) | intermolecular bond is formed <br> Diagram shows the orientation of the molecules as shown below with dipoledipole forces shown in-between molecules <br> $\delta-$ <br> мимим <br> $\delta$ |  | 1 <br> 1 <br> 1 | AO3 2.2.2k |
| 5(a) | $\begin{aligned} & \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{OH} \\ & \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CH}_{3} \end{aligned}$ |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO1 } \\ \text { 4.1.1b } \end{gathered}$ |
| 5(b) | Butan-1-ol has an OH group which can hydrogen bond to other - OH groups on the butan-1-ol molecules. <br> Pentane has (weak) instantaneous dipole-instantaneous dipole forces between the molecules which are transient (temporary). (Butan-1-ol also has these forces.) <br> The hydrogen bonds are stronger than the instantaneous dipole-instantaneous dipole forces between the molecules. |  | $1$ <br> 1 $1$ | $\begin{gathered} \mathrm{AO2} \\ \text { 2.2.2l/4.2.1a } \end{gathered}$ |

## A Level OCR Chemistry

## Chapter 10 - answers

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| 5(c) | Fill both burettes with the two liquids and rub the nylon rod with the silk rag to produce a charged rod <br> Run each liquid slowly from the burette (into a beaker) <br> Observe any deviation of the liquid stream <br> The butan-1-ol will be deviated a lot because it is a polar liquid. <br> The pentane stream will not be deviated because it is non-polar and will not be affected by the electrical field. |  | $\begin{aligned} & 1 \\ & 1 \\ & 1 \\ & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO3 } \\ \text { 1.2.1a } \end{gathered}$ |
| 6(a) | $\mathrm{CH}_{2} \mathrm{BrCH}_{2} \mathrm{Br}$ 1, 2-dibromoethane $\mathrm{CH}_{3} \mathrm{CHBr}_{2}$ 1, 1-dibromoethane |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{gathered} \text { AO2 } \\ \text { 4.1.1e } \end{gathered}$ |
| 6(b)(i) |  <br> 1, 1-dibromoethene <br> E-1, 2-dibromoethene <br> Z-1, 2-dibromoethene |  | 1 <br> 1 <br> 1 | $\begin{gathered} \text { AO2 } \\ 4.1 .3 \mathrm{c} \end{gathered}$ |
| 6(b)(ii) | 1, 1-dibromoethene and Z-1, 2-dibromoethene are polar <br> The more electronegative bromine atoms are at one end or on one side of the molecule <br> This leads to an asymmetric distribution of electrons/charge and therefore the molecule is polar |  | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | $\begin{aligned} & \text { AO3 } \\ & 4.1 .3 \mathrm{C} \end{aligned}$ |
| 6(b)(iii) | In $\mathrm{C}_{2} \mathrm{H}_{4} \mathrm{Br}_{2}$ the bond angle is $109.5^{\circ}$ In $\mathrm{C}_{2} \mathrm{H}_{2} \mathrm{Br}_{2}$ the angle is $120^{\circ}$ |  | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $\begin{aligned} & \mathrm{AO} 2 \\ & \text { 2.2.2g/4.1.3b } \end{aligned}$ |

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## Chapter 10 - answers

## Answers:

1 (a) False.
$\pi$ and 10 never changes/they are not variables so they can't be proportional to each other
(b) False. $\pi^{2}=9.8696 \ldots$ this is not the same as 10 (you could write: $\pi^{2} \neq 10$ )
(c) True. $\pi^{2}=9.8696$ which is very close to 10
(d) True.
(e) False e to a negative power is less than 1. The more negative it is, the smaller it gets. $\mathrm{e}^{-100}=3.74 \times 10^{-44}$
(f) False.

For a zero order reaction, rate does not vary with concentration
(g) False.

Rate is proportional to the rate constant, $k . k=A e^{-\frac{E_{a}}{R T}}$ (Arrhenius equation). So, rate does not increase linearly with temperature even though it does increase with temperature. (Actually rate $\propto \mathrm{e}^{-\frac{1}{T}}$.
(h) False.

The enthalpy change of combustion of propane is higher than ethane because there are more bonds made for the combustion of propane. However, it is incorrect to use the $\gg$ symbol, because $\Delta_{c} H^{\circ}\left(\mathrm{C}_{3} \mathrm{H}_{8}\right)$ is of the same order of magnitude as $\Delta_{c} H^{\circ}\left(\mathrm{C}_{2} \mathrm{H}_{6}\right)$. Both enthalpy changes are approximately in the $1000 \mathrm{skJ} \mathrm{mol}{ }^{-1}$
2 (a) >
$A_{\mathrm{r}}(\mathrm{Se})=34, A_{\mathrm{r}}(\mathrm{Se})=16$. It would be wrong to use the $\gg$ symbol as they are both the same order of magnitude. Both in the tens.
(b) $\gg$

The mass of a diamond is likely to be comparable to the mass of plastic bag, but a polymer has millions of molecules in it. So, there would be millions more carbon atoms in diamond than polymer chains in a plastic bag. It's right to use $\gg$ and not >, because it is many orders of magnitude higher.
(c) $\sim o r \approx$.

Chemistry A level has a even gender split - it is the most gender balanced of the science subjects (at time of writing). It would not be right to use = as it is very, very unlikely to be perfectly balanced.

