

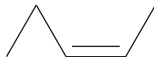
A Level AQA Chemistry

Chapter 14 – answers

| Question | Answers | Extra information | Mark | AO Spec reference |
|----------|---|---|-------------|-------------------------------|
| 01.1 | A <u>compound</u> made up of hydrogen and carbon <u>only</u> . | The underlined terms are essential. | 1 | AO1 3.3.2 |
| 01.2 | Its molecular formula is C ₅ H ₁₂ . Relative molecular mass = 5 × 12 + 12 × 1 = 72 | | 1 1 | AO2 3.1.2.4 |
| 01.3 | Isomers have the same molecular formula but different structural formulae. | | 1 | AO1 3.3.1.3 |
| 01.4 | CH ₃ CH ₂ CH ₂ CH ₂ CH ₃ pentane CH ₃ CH(CH ₃)CH ₂ CH ₃ 2-methylbutane C(CH ₃) ₄ or (CH ₃) ₄ C 2,2-dimethylpropane | Both formula and name required for each mark. Any order will do. For the 2 nd and 3 rd isomers = make sure the commas and dashes are correct. | 1 1 1 | AO2 3.3.1.3 |
| 01.5 | 0.36 g = 0.36/72 mol = 5 × 10 ⁻³ mol Number of molecules = 5 × 10 ⁻³ × 6.022 × 10 ²³ = 3.01 × 10 ²¹ In each molecule there are 17 atoms. ∴ Total number of atoms = 3.01 × 10 ²¹ × 17 = 5.12 × 10 ²² | Just the answer with no reasoning shown is 1 mark only. | 1 1 1 | AO2 and AO3 MS1.4; 3.1.2.2 |

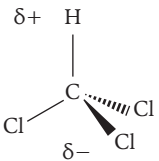
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| 02.1 | I pent-1-ene II 2-methylbut-2-ene III pent-2-ene | Do not allow pentene | 1 1 1 | AO2 3.3.1.3 |
| 02.2 | Stereoisomers have the same structural formula Have different arrangement of bonds in space | For 2 nd mark allow have different displayed formulae | 1 1 | AO1 3.3.1.3 |
| 02.3 | Isomer III  | | 1 1 | AO2 3.3.1.3 |
| 02.4 | There is no free rotation about the C=C double bond. On each carbon of the C=C bond, there are 2 different atoms or groups. | | 1 1 | AO1 3.3.1.3 |
| 02.5 | Methylcyclobutane | | 1 | AO2 3.3.1.1 |
| 03.1 | Use the ideal gas equation $n = \frac{PV}{RT}$ $n = \frac{1.01 \times 10^5 \times 85 \times 10^{-6}}{8.31 \times 450}$ $= 2.2957... \times 10^{-3} \text{ mol}$ $M_r = m/n = 0.20/2.2957... \times 10^{-3} = 87 \text{ g mol}^{-1}$ | Some evidence of using the equation is required answer to 2 s.f. | 1 1 1 | AO2 3.1.2.3 |

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| 03.2 | The last (highest value) significant peak | | 1 | AO1 3.3.6.2 |
| 03.3 | Error = $100\% \times 88-87 / 88$ 1.14% | | 1 1 | AO2 PS2.3 |
| 03.4 | $a = 109.5 / 109^\circ$ $b = 104.5^\circ$ | Allow 109° | 1 1 | AO1 MS4.1; 3.1.3.5 |
| 04.1 | Chlorine is more <u>electronegative</u> than carbon electrons are not shared equally / The electrons in the covalent bond are attracted to more strongly to the chlorine atom electron cloud distorted towards chlorine | | 1 1 1 | AO2 3.1.3.6 |
| 04.2 | The molecule is (tetrahedral and) symmetrical (The electrons (electron clouds) are symmetrically distributed) This so the dipoles cancel out | | 1 1 | AO2 3.1.3.6 |
| 04.3 | δ^+ H  δ^- Cl | The diagram would be reversed showing δ^- end of CHCl_3 pointing towards the δ^+ end of propanone | 2 | AO3 3.1.3.6 |
| 04.4 | CHCl_3 – it is more polar It interacts more strongly/is attracted more strongly to the polar stationary phase and takes longer to leave the column | | 1 1 | AO3 3.1.3.6; 3.3.16 |

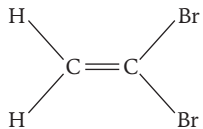
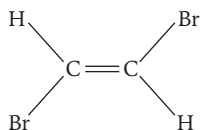
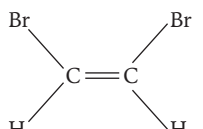
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| 04.5 | <p>Both compounds have permanent dipoles Diagram shows the orientation of the molecules as shown below with dipole-dipole forces shown in-between molecules</p> | The dipoles must be shown correctly | 1 2 | AO3 3.1.3.7 |
| 05.1 | <p>CH₃CH₂CH₂CH₂OH CH₃CH₂CH₂CH₂CH₃</p> | | 1 1 | AO1 3.3.1.3 |
| 05.2 | <p>Butan-1-ol has an OH group, which can hydrogen bond to other OH groups on the butan-1-ol molecules. Pentane has only very weak instantaneous dipole-instantaneous dipole forces between the molecules which are transient (temporary). (Butan-1-ol also has these forces.) The hydrogen bonds are stronger than the instantaneous dipole-instantaneous dipole forces between the molecules.</p> | | 1 1 1 | AO2 3.1.3.6 and 3.1.3.7 |
| 05.3 | <p>Fill both burettes with the two liquids and rub the nylon rod with the silk rag to produce a charged rod. Run each liquid <u>slowly</u> from the burette (into a beaker). Observe any deviation of the liquid stream. The butan-1-ol will be deviated a lot because it is a polar liquid. The pentane stream will not be deviated because it is non-polar and will not be affected by the electrical field.</p> | | 1 1 1 1 | AO1 and AO3 PS1.2; 3.1.3.7 |

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| 06.1 | $\text{CH}_2\text{BrCH}_2\text{Br}$ 1, 2-dibromoethane CH_3CHBr_2 1, 1-dibromoethane | Both correct name and structure required for each mark | 1 1 | AO2 3.3.1.3 |
| 06.2 |  <p>1, 1-dibromoethene</p>  <p><i>E</i>-1, 2-dibromoethene</p>  <p><i>Z</i>-1, 2-dibromoethene</p> | Both correct name and structure required for each mark | 1 1 1 | AO2 3.3.1.3 |
| 06.3 | 1, 1-dibromoethene and <i>Z</i> -1, 2-dibromoethene are polar. The more electronegative bromine atoms are at one end or on one side of the molecule. This leads to an asymmetric distribution of electrons/charge and therefore the molecule is polar. | | 1 1 1 | AO3 3.1.3.6; 3.3.1.3 |
| 06.4 | In $\text{C}_2\text{H}_4\text{Br}_2$ the bond angle is 109.5° In $\text{C}_2\text{H}_2\text{Br}_2$ the angle is 120° | | 1 1 | AO2 3.1.3.5 |

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Skills Box answers:

- $\frac{0.1}{12} \times 100 = 0.83\%$
- $\frac{0.1}{45} \times 100 = 0.22\%$
- $\Delta V = 36.75 - 12.50 = 24.25 \text{ cm}^3$
 $\frac{2 \times 0.05}{24.25} \times 100 = 0.41\%$
- $\Delta T = 45.0 - 22.5 = 22.5$
 $\frac{2 \times 0.05}{22.5} \times 100 = 0.44\%$