

**Assessment Schedule – 2025****Chemistry: Demonstrate understanding of the properties of selected organic compounds (91165)****Evidence**

Q	Evidence			Achievement	Merit	Excellence
ONE (a)	Compound	Structure	Name	<ul style="list-style-type: none"> <li>THREE of six correct.</li> </ul>	<ul style="list-style-type: none"> <li>SIX of six correct.</li> </ul>	
	A	$  \begin{array}{c}  \text{H} \quad \text{H} \\    \quad   \\  \text{H}-\text{C}-\text{C}-\text{N} \\    \quad   \\  \text{H} \quad \text{H}  \end{array}  $	ethanamine			
	B	$  \begin{array}{c}  \text{Br} \\    \\  \text{Br}-\text{C}-\text{H} \\    \\  \text{H}  \end{array}  $	dibromomethane			
	C	$  \begin{array}{c}  \text{H} \quad \text{CH}_3 \text{H} \\    \quad   \\  \text{H}-\text{C}-\text{C}-\text{C}-\text{C} \\    \quad   \quad   \\  \text{H} \quad \text{CH}_3 \text{H} \quad \text{OH}  \end{array}  $	3,3-dimethylbutanoic acid			
	D	$  \begin{array}{ccccccccc}  \text{H} & \text{H} & \text{Br} & \text{H} & \text{H} & \text{H} \\    &   &   &   &   &   \\  \text{H}-\text{C}-\text{C}-\text{C}-\text{C}=\text{C}-\text{C}-\text{H} \\    &   &   & & &   \\  \text{H} & \text{H} & \text{H} & & & \text{H}  \end{array}  $	4-bromohex-2-ene			
	E	$  \begin{array}{ccccc}  \text{H} & \text{H} & \text{H} \\    &   &   \\  \text{H}-\text{C}-\text{C}-\text{C}-\text{C}\equiv\text{C}-\text{H} \\    &   &   \\  \text{H} & \text{H} & \text{H}  \end{array}  $	1-pentyne			
	F	$  \begin{array}{c}  \text{H} \quad \text{CH}_3 \text{H} \\    \quad   \\  \text{H}-\text{C}-\text{C}-\text{C}-\text{H} \\    \quad   \quad   \\  \text{H} \quad \text{OH} \quad \text{H}  \end{array}  $	2-methyl propan-2-ol			

(b)	<p>Acidified dichromate will react with the alcohol group in Compound H, but will not react with G. This is an oxidation reaction, producing a carboxylic acid. This reaction must be heated to occur. The dichromate solution will be orange when mixed with the alcohol and turn green as the alcohol is oxidised.</p> $\begin{array}{c} \text{O} \\ \text{  } \\ \text{C} - \text{CH}_2 - \text{CH} - \text{CH}_3 \\   \\ \text{HO} \\ \text{CH}_3 \end{array}$	<ul style="list-style-type: none"> <li>Identifies type of reaction.</li> <li>OR</li> <li>Colour change.</li> <li>OR</li> <li>Compound H chosen.</li> </ul>	<ul style="list-style-type: none"> <li>Full explanation missing one aspect (reaction type, conditions, structure of product or observation).</li> </ul>	<ul style="list-style-type: none"> <li>Full explanation with product drawn correctly and reference to G not reacting.</li> </ul>				
(c)(i)	<p>Compound G: Primary Compound J: Primary Compound I: Tertiary Compound K: Secondary</p>	<ul style="list-style-type: none"> <li>TWO of four correct.</li> </ul>	<ul style="list-style-type: none"> <li>Both structures.</li> <li>AND</li> <li>Major / minor correct.</li> </ul>					
(ii)	<table border="1" data-bbox="233 509 990 695"> <tr> <td data-bbox="233 509 608 636"> <math display="block">\begin{array}{c} \text{H}_3\text{C} \\ \diagdown \\ \text{C} = \text{C} \\ \diagup \\ \text{H} \quad \text{CH}_3 \\ \diagup \\ \text{CH}_3 \end{array}</math> </td><td data-bbox="608 509 990 636"> <math display="block">\begin{array}{c} \text{H} \\ \diagdown \\ \text{C} = \text{C} \\ \diagup \\ \text{H} \quad \text{H} \\ \diagup \\ \text{CH}(\text{CH}_3)_2 \end{array}</math> </td></tr> <tr> <td data-bbox="233 636 608 695">Major</td><td data-bbox="608 636 990 695">Minor</td></tr> </table>	$\begin{array}{c} \text{H}_3\text{C} \\ \diagdown \\ \text{C} = \text{C} \\ \diagup \\ \text{H} \quad \text{CH}_3 \\ \diagup \\ \text{CH}_3 \end{array}$	$\begin{array}{c} \text{H} \\ \diagdown \\ \text{C} = \text{C} \\ \diagup \\ \text{H} \quad \text{H} \\ \diagup \\ \text{CH}(\text{CH}_3)_2 \end{array}$	Major	Minor	<ul style="list-style-type: none"> <li>Correctly identifies alkene is made.</li> </ul>	<ul style="list-style-type: none"> <li>Explains why two products are formed.</li> <li>OR</li> <li>States how major / minor products can be identified.</li> </ul>	<ul style="list-style-type: none"> <li>Justifies why two products are formed and how to identify major / minor products, with reference to the structure of Compound K.</li> </ul>
$\begin{array}{c} \text{H}_3\text{C} \\ \diagdown \\ \text{C} = \text{C} \\ \diagup \\ \text{H} \quad \text{CH}_3 \\ \diagup \\ \text{CH}_3 \end{array}$	$\begin{array}{c} \text{H} \\ \diagdown \\ \text{C} = \text{C} \\ \diagup \\ \text{H} \quad \text{H} \\ \diagup \\ \text{CH}(\text{CH}_3)_2 \end{array}$							
Major	Minor							
(iii)	<p>Compound K is an asymmetrical haloalkane, meaning two products can be formed depending upon which neighbouring carbon the hydrogen atom is eliminated from. The major product forms when the Cl group is eliminated, along with a H atom from the neighbouring carbon with the least number of H atoms attached. In Compound K, C1 has three H atoms, while C3 only has one; therefore the major product is when the hydrogen is eliminated from C3, forming 2-methylbut-2-ene (and the minor product is 3-methylbut-1-ene).</p>							

N0	N1	N2	A3	A4	M5	M6	E7	E8
No response; no relevant evidence.	1a	2a	3a	4a	3m	4m	2e minor error	2e

Q	Evidence			Achievement	Merit	Excellence
TWO (a)(i)	Compound B  $\begin{array}{ccccccc} & H & H & CH_3 & H & H & \\ &   &   &   &   &   & \\ H & - C & - C & - C & - C & - C & - H \\ &   &   &   &   &   & \\ & H & H & H & NH_2 & H & \end{array}$	Compound C  $\begin{array}{ccccccc} & H & H & CH_3 & H & H & \\ &   &   &   &   &   & \\ H & - C & - C & - C & - C & - C & - H \\ &   &   &   &   &   & \\ & H & H & H & OH & H & \end{array}$	Compound D  $\begin{array}{ccccccc} & H & H & CH_3 & H & H & \\ &   &   &   &   &   & \\ H & - C & - C & - C & = C & - C & - H \\ &   &   &   & &   & \\ & H & H & H & & H & \end{array}$	• FOUR of eight correct.	• SIX of eight correct.	• EIGHT of eight correct.
(ii)	Reaction type W: substitution Reaction type X: elimination Reaction type Y: addition Reaction type Z: substitution					
(iii)	Reagent 4: HCl					

(b)(i)	$\left[ \begin{array}{ccccccc} \text{H} & \text{H} & \text{H} & \text{H} & \text{H} & \text{H} \\   &   &   &   &   &   \\ \text{C} - \text{C} \\   &   &   &   &   &   \\ \text{C}_4\text{H}_9 & \text{H} & \text{C}_4\text{H}_9 & \text{H} & \text{C}_4\text{H}_9 & \text{H} & \text{C}_4\text{H}_9 \end{array} \right]_n$	<ul style="list-style-type: none"> <li>• Correctly drawn polymer.</li> </ul>		
(ii)	<p>C=C, carbon to carbon double bond</p> <p>In an addition polymerisation reaction, C=C double bonds are broken in order for new single bonds to form between monomers, joining into long repeating chains called polymers. The monomers have an unsaturated C=C double bond, which is much more reactive than the resulting C–C saturated single bond in the polymer structure. This is why the monomer can undergo addition polymerisation, while the polymer with only C–C single bonds, cannot.</p>	<ul style="list-style-type: none"> <li>• Double bond is broken.</li> </ul> <p>OR</p> <p>Identifies polymer is saturated (C–C) or monomer is unsaturated (C=C).</p>	<ul style="list-style-type: none"> <li>• Links saturated / unsaturated nature of polymer / monomer to materials reactivity.</li> <li>• Explains addition polymerisation.</li> </ul>	<ul style="list-style-type: none"> <li>• Explains the difference of monomer and polymer in terms of structure, reactivity.</li> </ul>
(iii)	$\begin{array}{c} \text{F} \quad \text{F} \\ \diagdown \quad \diagup \\ \text{C} = \text{C} \\ \diagup \quad \diagdown \\ \text{F} \quad \text{F} \end{array}$	<ul style="list-style-type: none"> <li>• Correctly drawn monomer.</li> </ul>		<p>AND</p>
(iv)	<p>Teflon does not melt at high temperatures. The polymer Teflon is not (chemically) reactive.</p>	<ul style="list-style-type: none"> <li>• States Teflon has a high melting point.</li> </ul>	<ul style="list-style-type: none"> <li>• Links Teflon does not melt at high temperatures to use.</li> </ul>	<p>OR</p>
		<p>OR</p>	<p>States that Teflon is not (chemically) reactive.</p>	<p>Links Teflon is not (chemically) reactive to use.</p>
(c)	$\begin{array}{c} \text{Cl} \\   \\ \text{CH}_2 - \text{CH}_2 - \text{CH}_2 - \text{CH}_2 - \text{CH}_3 \end{array}$	<ul style="list-style-type: none"> <li>• THREE correct.</li> </ul>		
	$\begin{array}{c} \text{Cl} \\   \\ \text{CH}_3 - \text{CH} - \text{CH}_2 - \text{CH}_2 - \text{CH}_3 \end{array}$			
	$\begin{array}{c} \text{Cl} \\   \\ \text{CH}_3 - \text{CH}_2 - \text{CH} - \text{CH}_2 - \text{CH}_3 \end{array}$			

<b>NØ</b>	<b>N1</b>	<b>N2</b>	<b>A3</b>	<b>A4</b>	<b>M5</b>	<b>M6</b>	<b>E7</b>	<b>E8</b>
No response; no relevant evidence.	1a	2a	3a	4a	3m	4m	2e minor error	2e

Q	Evidence	Achievement	Merit	Excellence
THREE (a)(i)	Carboxylic acid.	<ul style="list-style-type: none"> <li>Correct group.</li> </ul>		
(ii)	1-heptanamine and 1-propanamine both contain a polar amine group, which is soluble in water, as the forces of attraction between the water and the amine group (solvent to solute) are sufficient to overcome the solute-solute interactions.	<ul style="list-style-type: none"> <li>Gives a basic method (water / heating).</li> </ul>	<ul style="list-style-type: none"> <li>Explains that chain length will determine which is soluble in water.</li> </ul>	<ul style="list-style-type: none"> <li>Fully explains either method, including associated observation, for correct identification and explains relevant carbon chain length effect.</li> </ul>
	However, 1-propanamine has a short chain of 3 carbons, so will be water soluble, while 1-heptanamine has a longer chain with 7 carbons; therefore it is not fully soluble in water.	OR		
	Identify by mixing a sample of each with water. 1-heptanamine will form two layers, but 1-propanamine will not.	Identifies a difference in properties (solubility / boiling point).	OR	Explains that chain length will determine which has lower/higher boiling point.
	The longer carbon chain also means 1-heptanamine has a higher melting and boiling point. Therefore, heating each compound to observe the temperature at which they boil will allow identification. 1-propanamine will boil first at a lower temperature, due to the shorter carbon chain length.			
(iii)	Gallic acid contains a carboxylic acid group which will turn damp blue litmus red as it is acidic.	<ul style="list-style-type: none"> <li>Correct observations.</li> </ul>		
	1-heptanamine is an amine, which is basic. Therefore, the sample will turn damp red litmus blue.			
(iv)	L2 acid reaction.	<ul style="list-style-type: none"> <li>A correct method is described.</li> </ul>	<ul style="list-style-type: none"> <li>A correct method is explained with observations.</li> </ul>	
	The acid reacts with a metal carbonate in an acid-base reaction, forming a metal salt, carbon dioxide, and water. The CO <sub>2</sub> will form as a gas, so bubbles will be observed in the solution.			

(b)(i)	(ii)	<p>Group B.</p> <p>B forms cis / trans isomers, as it has both requirements for geometric isomerism. It contains a C=C double bond, which prevents rotation of atoms about the bond. A also has this. However, the second requirement is that each carbon of the double bond is bonded to two different atoms or groups. A doesn't meet this requirement, as one of the carbons has two CH<sub>3</sub> groups attached. In B, one of the carbons of the double bond has a hydrogen atom and an R<sub>3</sub> group, while the other has a CH<sub>3</sub> group and a different R<sub>2</sub> group. This means that B can be arranged differently in space, forming geometric isomers.</p>	<ul style="list-style-type: none"> <li>Identifies B.</li> <li>Identifies C=C double bond.</li> </ul> <p>OR</p> <p>Identifies two different atoms / groups need to be attached to the carbons in the double bond.</p>	<ul style="list-style-type: none"> <li>Explains why the double bond is required for geometric isomerism.</li> <li>Explains why both C atoms on the double bond must have two different atoms / groups.</li> </ul> <ul style="list-style-type: none"> <li>Justifies why B forms geometric isomers, while A does not.</li> </ul>
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NØ	N1	N2	A3	A4	M5	M6	E7	E8
No response; no relevant evidence.	1a	2a	3a	4a	2m	3m	2e minor error	2e

**Cut Scores**

Not Achieved	Achievement	Achievement with Merit	Achievement with Excellence
0 – 7	8 – 13	14 – 18	19 – 24