

# **TOPIC 9 HW MS**

1. Compare spectrum of aldehyde with known spectrum of pentanal *Must be a specific reference to a comparison.* 

#### Exact match

Allow 'fingerprint regions match exactly'.

1

1

[2] **2.** (a) If 2 stage test for one compound, award no marks for that compound, eg no mark for ROH or RX to alkene then Br<sub>2</sub> test. If reagent is wrong or missing, no mark for that test; if wrong but close/incomplete, lose reagent mark but can award for correct observation. In each test, penalise each example of wrong chemistry, eg AgClr<sub>2</sub> propan-1-ol acidifiedpotassiumdichromate cture sodium Named acid + conc  $H_2SO_4$ named acyl chloride M1 1 (orange) turns green effervescence Sweet smell Sweet smell /misty fumes Misty fumes M2 1 propanal add Tollens or Fehlings / Benedicts acidifiedpctassiumdichromate Bradys or 2,4-dnph if dichromate used for alcohol cannot be used for aldehyde M3 1 Tollens: silver mirror or Fehlings/ Benedicts: red ppt (orange) turns green Yellow or orange ppt M4 1



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	propanoic acid Named carbonate/ hydrogencarbonate water and UI (paper) Named alcohol + conc H₂SO₄ sodium or magnesium		
	if sodium used for alcohol cannot be used for acid M5 effervescence	1	
	orange/red Sweet smell effervescence Misty fumes <i>if PCI₅ used for alcohol cannot be used for acid</i>		
	Mo 1-chloro propane NaOH then acidified AgNO <sub>3</sub> AgNO <sub>3</sub> If acidification missed after NaOH,no mark here	1	
	but allow mark for observation M7 white ppt white ppt	1	
	M8	1	[8]
(a)	(i) C	1	
	(ii) A	1	
	(iii) D	1	
	(iv) B	1	

3.



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(b) M1 Br<sub>2</sub> OR bromine (water) OR bromine (in CCl<sub>4</sub> / organic solvent)

If M1, has no reagent or an incorrect reagent,

CE=0

Ignore "acidified"

For M1 penalise Br (or incorrect formula of other correct reagent), but mark on

For M1, it must be a whole reagent and/or correct formulae

## Either order

If oxidation state given in name, it must be correct.

M2 cyclohexane OR A or the alkane: remains orange / red/vellow / brown / the same OR no reaction OR reference to colour going to cyclohexane layer

For M2 credit "no change"

Ignore "nothing" Ignore "nothing happens"

Ignore "no observation"

M3 cyclohexene OR D or the alkener decolourised / goes colourless / loses its colour

For M3, ignore "goes clear"

Alternatives : potassium manganate(VII)

M1 KMnO₄ in acid M2 purple M3 colourless

M1 KMnO in alkali / neutral M2 purple M3 brown solid

Give appropriate credit for the use of iodine and observations No credit for combustion observations

3

- (c) M1 <u>acidified potassium or sodium dichromate</u> For M1, it must be a whole reagent and/or correct formulae
  - **OR** eg  $\underline{H_2SO_4} / \underline{K_2Cr_2O_7}$  **OR**  $\underline{H_+} / \underline{K_2Cr_2O_7}$
  - *OR* correct combination of formula and name If oxidation state given in name, it must be correct.

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#### M2 oxidation OR oxidised OR redox

Do not penalise incorrect attempt at formula if name is correct or vice versa

M3 secondary / 2° (alcohol)

Credit acidified potassium chromate(VI) /  $\underline{H_2SO_4}$ +  $K_2CrO_4$ 

3



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	(d)	M1	(free-) <u>radical substitution</u> (mechanism) M1 both words required		
		M2	$Br_2$ 2 $Br_2$		
			Penalise absence of dot once only.		
		M3	$Br\bullet + CH_4 \qquad \bullet CH_3 + HBr$		
			Penalise + or – charges every time		
		M4	$Br_2 + \bullet CH_3$ $CH_3Br + Br \bullet$		
			Accept dot anywhere on methyl radical		
			Accept a <u>correct</u> termination step for 1 mark if neither M3 nor M4 are scored; otherwise ignore		
			termination steps		
			Mark independently		
			NB If Cl <sub>2</sub> is used, penalise every time ( this may be for M2, M3 and M4)		
			If cyclohexane is used, penalise every time (this may be for M3 and M4)		
		M5	Condition		
			ultra-violet / uv / sun light		
		OR	<u>high</u> temperature		
		OR	125 °C <sup>≤</sup> T <sup>≤</sup> 600 °C		
		OR	400 K T 800 K		
			For M5 gnore "heat"		
			5		[15]
			No.		
4.	(a)	Fractio	onal distillation (under reduced pressure)	1	
	(b)	Pas	Q incoluble / remove by filtration		
	(U)	Dao	Do not allow answers which refer to reaction		
			rate		
				1	
	(c)	Both	n contain OH group		
			Allow OH stretch in ir spectrum of each		
			compound Do not allow 'same bonds'		
			5		

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[3]

1



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1

1

2

1

3

1

1

[5]

- 5. (a) Secondary OR 2° (alcohol);
  - (b) Spectrum is for **butanone (or formula) or butan-2-one** <u>The explanation marks depend on correctly</u> <u>identifying butanone.</u>

If butanone is correctly identified, award any two from

- (Strong) absorption / peak at approximately 1700 (cm<sup>-1</sup>) / 1710 (cm<sup>-1</sup>) / in the range 1680 – 1750 (cm<sup>-1</sup>) This needs to be stated.
- (Characteristic) absorption / peak for C=O (may be shown on the spectrum in the correct place).
- No absorption / peak in range 3230 to 3550 cm<sup>-1</sup>
  - No absorption / peak for an OH group. Look at the spectrum to see if anything is written on it that might gain credit. Allow the words "dip" OR "spike" OR "low transmittance" as alternatives for absorption.

(c) <u>Displayed structure</u> for 2-methylpropan-2-ol

 (a) Allow 1 mark each for any correctly drawn primary, secondary and tertiary alcohol of molecular formula C₄H<sub>8</sub>O

Must have all bonds drawn out but ignore the

Tertiary alcohol cannot be oxidised

bond angles

(b) Region 1500–400 cm<sup>-1</sup>





exact match to spectrum of known compound

1



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Reacting Tollens' / Fehling's with alcohols directly is incorrect and gains no M2 or M3. Detailed observations relating to the reactions are not needed but should be penalised where incorrect.

[3]

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9.	Tes	t	bromine (water) / iodine Accept 'Br₂' or 'bromine in a named solvent'. Do not accept 'Br' Use of UV light, CE (lose next mark as well)	1	
Obse	ervati	on oi	range / yellow / (red-)brown to colourless Must have correct reagent to score this mark. For I <sub>2</sub> , allow red-brown / purple to colourless.	1	[2]
10.	(a)	(i)	Green Ignore shades of green.	1	
		(ii)	Excess acidified potassium dichromate(VI)	1	
			Reflux (for some time)	1	
			In the diagram credit should be given for • a vertical condenser	1	
			<ul> <li>an apparatus which would clearly work</li> <li>Do not allow this mak for a flask drawn on its own.</li> <li>Penalise diagrams where the apparatus is sealed.</li> </ul>	1	
		(iii)	Distillation	1	
			Immediately the reagents are mixed)	1	
	(b)	Kee	p away from naked flames Allow heat with water-bath or heating mantle. If a list is given ignore eye protection, otherwise lose this mark	1	
	(c)	(i)	Tollens' or Fehling's reagents Incorrect reagent(s) loses <b>both</b> marks. Accept mis-spellings if meaning is clear.	1	
			Silver mirror / red ppt. formed Accept 'blue to red' but not 'red' alone.	1	
		(ii)	Sodium carbonate (solution) / Group II metal Allow indicator solutions with appropriate colours. Accept any named carbonate or hydrogen	1	

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carbonate.

Effervescence / evolves a gas Accept 'fizzes'. 1

1



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(d) Propanoic acid

If this mark is lost allow one mark if there is reference to stronger intermolecular forces in the named compound. Lose M1 and M3.

Contains hydrogen bonding

Some comparison with other compounds explaining that the intermolecular forces are stronger in propanoic acid

[15]

1

1

1

1

## 11. (a) (i) <u>Electron pair donor</u>

OR

Species which uses a pair of electrons to form a co-ordinate/covalent bond.

Credit "lone pair" as alternative wording Credit "electron pair donator"

(ii) Replacement of the halogen (atom) (by the nucleophile)

OR

The <u>carbon-halogen bond/C-X</u> breaks and a bond forms with the nucleophile or between the carbon and the nucleophile

They must describe the idea of substitution in a haloalkane.

Accept the idea that a nucleophile replaces the halogen which becomes a halide ion

Penalise reference to "halogen molecule" and penalise the idea that the haloalkane contains a halide

(iii) <u>Splitting molecules using/by water</u>

OR

breaking/splitting/dissociating (C<sub>i</sub>VX) bond(s)/using/by water NOT simply the reaction with water or simply the addition of water. Ignore "compound"

1

1

(iv) (Heat) energy/enthalpy required/needed/absorbed (at constant

13



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pressure) to break/split it/the (carbon-halogen) bond

OR

(Heat) <u>energy/enthalpy required/needed/absorbed</u> (at constant pressure) for <u>homolysis</u> of <u>the</u> (C–X/the carbon-halogen) <u>bond</u> *Ignore bond formation Ignore "average"* 



- M1 must show an arrow from the lone pair of electrons on the oxygen atom of the negatively charged hydroxide ion to the central C atom.
- M2 must show the movement of a pair of electrons from the C-Br bond to the Br atom. Mark M2 independently.

Award full marks for an  $S_{\mathbb{N}}$ 1 mechanism in which M1 is the attack of the hydroxide ion on the intermediate carbocation.

Penalise M1 if covalent KOH is used Penalise M2 for formal charge on C or incorrect partial charges Penalise once only for a line and two dots to show a bond. Max 1 mark for the wrong reactant Accept the correct use of "sticks"

2

1



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M1 must show an arrow from the lone pair on oxygen of a negatively charged hydroxide ion to the correct H atom

M2 must show an arrow from the correct C-H bond to the C-C bond and should only be awarded if an attempt has been made at M1

M3 is independent provided it is from the original molecule

Award full marks for an E1 mechanism in which M2 is on the correct carbocation.

Penalise M1 if covalent KOH Penalise M3 for formal charge on C or incorrect partial charges Penalise once only for a line and two dots to show a bond Max 2 marks for wrong reactant

3

Accept the correct use of "sticks" for the molecule except for the C-H being attacked

M1 Stated that the spectrum has an <u>absorption/absorbance/</u> peak in the range 1620 cm<sup>-1</sup> to 1680 (cm<sup>-1</sup>) or specified correctly in this range from the spectrum

(ii)

- M2 depends on correct <u>range or wavenumber being</u> <u>specified</u>
- M2 (Infrared absorption) <u>due to C=C OR carbon-carbon double</u> bond

**QoL for correct M1 statement which** *includes both the word absorption (or alternative)* <u>*and*</u> *the correct range or wavenumber* 

15

Allow "peak" OR "dip" OR "spike" OR "trough" OR "low transmittance" as alternatives for



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absorption. For M2 it is not sufficient simply to state that an alkene has C=C M2 could be on the spectrum Ignore reference to other absorptions

[11]

2



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- **12.** B
- 13. D 14. D
  - [1]

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