



**General Certificate of Education**

**Chemistry 5421**

**CHM3/W Introduction to  
Organic Chemistry**

**Mark Scheme**

*2008 examination - June series*

Mark schemes are prepared by the Principal Examiner and considered, together with the relevant questions, by a panel of subject teachers. This mark scheme includes any amendments made at the standardisation meeting attended by all examiners and is the scheme which was used by them in this examination. The standardisation meeting ensures that the mark scheme covers the candidates' responses to questions and that every examiner understands and applies it in the same correct way. As preparation for the standardisation meeting each examiner analyses a number of candidates' scripts: alternative answers not already covered by the mark scheme are discussed at the meeting and legislated for. If, after this meeting, examiners encounter unusual answers which have not been discussed at the meeting they are required to refer these to the Principal Examiner.

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## CHM3/W

## Question 1

- |     |       |   |   |   |
|-----|-------|---|---|---|
| (a) | (i)   | A compound OR molecule containing OR consisting of hydrogen and carbon <u>only</u>  | 1 |   |
| (a) | (ii)  | $\text{C}_9\text{H}_{20} + 9\frac{1}{2}\text{O}_2 \longrightarrow 9\text{CO} + 10\text{H}_2\text{O}$  | 1 | Or multiple   |
| (a) | (iii) | <b>M1</b><br><br>High temperature<br><br>OR spark<br><br>OR $2500^\circ\text{C} \leq T \leq 4000^\circ\text{C}$<br><br><b>M2</b><br><br>$\frac{1}{2}\text{N}_2 + \frac{1}{2}\text{O}_2 \longrightarrow \text{NO}$ | 2 | Ignore "very" when referring to high temperature<br><br><br><br><br><br><br>Or multiple |
| (b) | (i)   | Platinum OR Pt<br><br>OR<br><br>Palladium OR Pd<br><br>OR<br><br>Rhodium OR Rh  | 1 | Credit phonetic spelling  |
| (b) | (ii)  | <b>25 NO    8 CO<sub>2</sub>    12½ N<sub>2</sub>    9 H<sub>2</sub>O</b>   | 1 | Credit whole equation doubled   |
| (c) |       | Releases OR is a source of OR provides a useable form of <u>heat OR energy</u> (when burned)  | 1 | Ignore "enthalpy"<br><br>Ignore "exothermic"  |
| (d) |       | $\text{H}_2\text{S} + 1\frac{1}{2}\text{O}_2 \longrightarrow \text{SO}_2 + \text{H}_2\text{O}$  | 1 | Or multiple   |

**Question 2**

- |           |  |   |  |
|-----------|--|---|--|
| (a)       | <b>M1</b><br>propanone   | 2 | <i>Credit propan-2-one as the <u>only</u> other acceptable response.</i>   |
|           | <b>M2</b><br>CH <sub>3</sub> CH <sub>2</sub> CHO<br>OR correct aldehyde with functional group drawn out.   |   | <i>Penalise CH=O for the aldehyde group</i>  |
| (b)       | <u>Functional group</u> (isomerism) ONLY   | 1 | <i>Both words needed</i>   |
| (c)       | The formula which shows the (actual) <u>number of each type of atom</u> OR <u>atoms of each element</u> in the molecule / compound.  | 1 | <i>Penalise reference to <u>ratio</u> of atoms</i>   |
| (d)       | C <sub>3</sub> H <sub>6</sub> O ONLY   | 1 | <i>Elements in any order</i>   |
| (e) (i)   | Redox OR Oxidation   | 1 | <i>Either word ONLY</i>  |
| (e) (ii)  | The mark is for the <u>structure</u> of propanoic acid<br><br>$\text{H}_3\text{C}-\text{CH}_2-\text{C} \begin{array}{l} \nearrow \text{O} \\ \searrow \text{OH} \end{array}$                                       | 1 | <i>Carboxylic acid group must show both the double and the single bonds to oxygen.</i><br><br><i>Formula could show C<sub>2</sub>H<sub>5</sub> OR CH<sub>3</sub>CH<sub>2</sub></i> |
| (e) (iii) | <i>If <u>Tollens'</u> is chosen, then</i><br><br><u>silver mirror / coating</u><br><br><u>OR black precipitate / solid</u><br><br><i>If <u>Fehling's</u> is chosen, then</i><br><br><u>red precipitate / solid</u> | 1 | <i><u>No mark for choice of reagent</u>, but do not award the mark if no reagent given.</i><br><br><i>Insist on Tollens' OR Fehling's as the ONLY reagents</i>                     |

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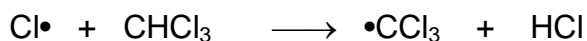
(f)	(i)	Contains a C=C OR double bond	1	
(f)	(ii)	1° OR Primary (alcohol)	1	
(f)	(iii)	<u>Structure</u> for		<i>Structure must be clear</i>
			1	
		CH <sub>2</sub> BrCHBrCH <sub>2</sub> OH		<i>Accept the alternatives</i>
				CH <sub>2</sub> BrCHBrCHO
		as shown or drawn out		<i>and</i>
				CH <sub>2</sub> BrCHBrCOOH

**Question 3**

(a) (i) (free-) radical substitution 1 *Both words needed*

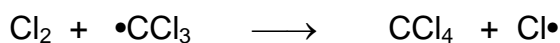
(a) (ii) *Initiation:*  $\text{Cl}_2 \longrightarrow 2\text{Cl}\cdot$  4 *Ignore reference to uv/sunlight/heat.*

*First propagation:*



*The dot can be anywhere on the  $\text{CCl}_3$*

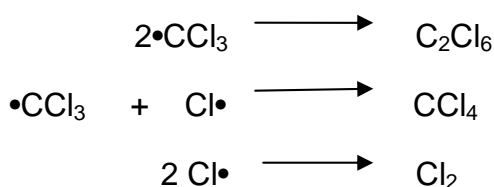
*Second propagation:*



*Penalise the absence of a dot once only*

*Termination:*

*One from*



*If half-arrows used, they must be correct.*

(b) (i) dichlorodifluoromethane 1 *ONLY these*

OR

1,1-dichloro-1,1-difluoromethane

**QoL**

*Penalise "cloro" and penalise "flouro"*

(b) (ii) (free- OR chlorine- ) radical 1 *Penalise reference to "substitution"*

OR chlorine atoms

*Ignore "alkyl"*

(b) (iii) propagation (step) 1 *Ignore "first" or "second"*

**Question 4**(a) (i) Polar C-Br (bond)

1

*The C-Br bond needs to be drawn out or stated*

OR

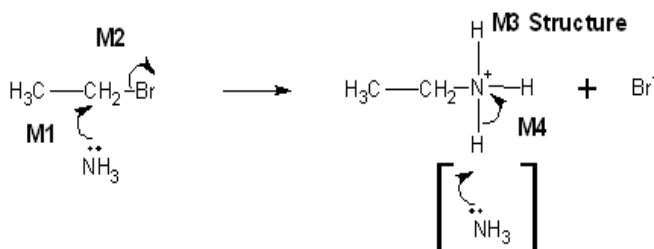
 $\delta+$  ( $\delta-$ )C — Br

OR

partially positive carbon atom on C-Br (bond)

(a) (ii)

4

*Penalise M1 if negative charge on ammonia**Penalise M2 for formal charge on C of C-Br or incorrect partial charges***M1** must show an arrow from the lone pair of electrons on the nitrogen atom of the ammonia to the C atom of the C-Br bond.*Penalise once only for a line and two dots to show a bond.***M2** must show the movement of a pair of electrons from the C-Br bond to the Br atom.

Mark M2 independently.

*Max 3 marks for wrong reactant or "sticks"***M3** is for the structure of the alkylammonium ion and could be obtained from  $\text{CH}_3\text{CH}_2\text{NH}_3^+$ **M4** is for an arrow from the N-H bond to the N atom.The second mole of ammonia is not essential for full credit, therefore ignore this part even if other species are used.Award full marks for an  $\text{S}_{\text{N}}1$  mechanism in which M1 is the attack of the ammonia on the intermediate carbocation.

*ONLY these names*

1

(c)	(i)	<u>Electron pair donor</u>	1	<i>Insist on both “electron pair” <u>and</u> an action by the electron pair.</i>
		OR		

Accept "lone pair"

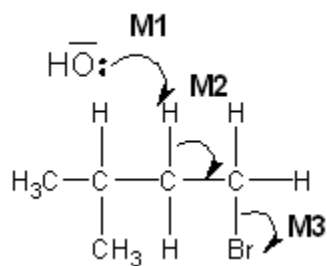
(c)	(ii)	1	<i>The minus sign can be anywhere on the OH</i>
	<u>Hydroxide ion</u> OR <u>OH<sup>-</sup></u> ONLY		<i>The electron pair (if drawn) MUST be on the oxygen atom</i>



## Question 5

- (a) **3-methylbut-1-ene** 1 **ONLY**
- (b) Elimination 1 *Credit "base elimination" but NOT "nucleophilic elimination"*

*No other prefix.*



3

*Penalise M1 if covalent KOH*

*Penalise M3 for formal charge on C or incorrect partial charges or extra arrow from Br to e.g. K+*

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

*Penalise once only for a line and two dots to show a bond.*

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

*Max 2 marks for the mechanism for wrong reactant or "sticks"*

**M3** is independent.

*Ignore incorrect organic product*

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

- (c) (i) *Structure OR name* 1 *If name is given it must be correct with "bromo" before "methyl"*

**2-bromo-3-methylbutane**

OR

$(\text{CH}_3)_2\text{CHCHBrCH}_3$

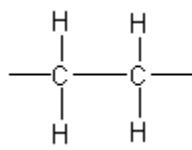
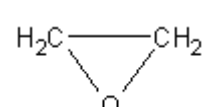
*Apply list principle if both structure and name given*

- (c) (ii) Electrophilic addition ONLY 1 *Both words*

- (d) **M1** Structure of pent-2-ene ONLY: 2 *M1 and M2 should be marked independently*
- $\text{CH}_3\text{CH}_2\text{CH}=\text{CHCH}_3$
- Credit M1 with a structure which is either linear or cis or trans (or both)*
- Structure MUST show a double bond
- M2** Type: Geometric(al) OR cis-trans  
(OR E/Z)

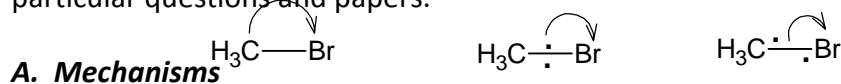
**Question 6**

- (a) **M1 (Separation based on boiling point)** 4 **Maximum 2 marks if candidate refers to bond breaking or cracking or using a Blast furnace or adding oxygen/air.**
- Separation depends on boiling point. (Ignore reference to melting point or density)
- M2 (Link between boiling point and size/forces)**
- Boiling point depends on  $M_r$  OR molecular size OR chain length OR intermolecular / Van der Waals forces OR candidate links boiling point with heavier/lighter fractions.
- M3 (Specified temperature gradient)**
- Specified temperature gradient OR difference stated on column/tower OR explained e.g. hotter at bottom.
- (If numbers used, accept up to 400°C at the base and down to 25°C at top)
- M4 (Position of molecules/fractions on column)**
- Lower  $M_r$  OR lower boiling point OR shorter chains OR smaller molecules OR more volatile/gaseous molecules OR lighter fractions (condense) at top
- (OR converse at bottom)
- M2** *Ignore references to smaller/larger fractions.*
- QoL for M2**
- (b) **M1** 3 **NOT just “hot” for M2**
- Thermal cracking
- M2**
- Any T (or range) in the range 400°C to 900°C OR
- High temperature.
- M3**
- (Free- OR alkyl-) radical
- Mark M1, M2 and M3 independently
- Penalise M2 for any reference to catalyst.**
- Penalise M3 for reference to “substitution”**

(c)	<p><b>M1</b></p> <p>Yeast or suitable enzyme (zymase)</p> <p><b>M2</b></p> <p>Fermentation</p> <p><b>M3</b></p> $\text{C}_6\text{H}_{12}\text{O}_6 \longrightarrow 2\text{CH}_3\text{CH}_2\text{OH} + 2\text{CO}_2$ <p style="text-align: center;">(or <math>2\text{C}_2\text{H}_5\text{OH}</math>)</p> <p>If <math>\text{O}_2</math> is in the equation, penalise the equation (M3) but NOT M1</p> <p>Mark M1, M2 and M3 independently</p>	<p>3</p> <p><i>Penalise M1 if oxygen/air added to process .</i></p> <p><i>Ignore reference to temperature and water.</i></p> <p><i>Penalise <math>\text{C}_2\text{H}_6\text{O}</math></i></p>
(d)	<p><b>M1</b></p> <p>(fractional) distillation OR fractionation</p> <p><b>M2</b></p> <p><u>c(oncentrated)</u> <math>\text{H}_3\text{PO}_4</math> OR <u>c(oncentrated)</u> <math>\text{H}_2\text{SO}_4</math> OR <math>\text{Al}_2\text{O}_3</math> OR alumina OR porous pot OR pumice</p> <p><b>M3</b></p> <p>(acid-catalysed) <u>dehydration</u> OR <u>elimination</u></p> <p>Mark M1, M2 and M3 independently</p>	<p>3</p> <p><i>There are two processes and therefore assume that the order of answers is the order of marking unless annotated by the candidate</i></p> <p><i>Penalise M3 for any other prefix such as "base"</i></p>
(e)	<p><b>M1</b></p> <div style="text-align: center;">        OR    <math>\text{—CH}_2\text{—CH}_2\text{—}</math> </div> <p><b>M2</b></p> <div style="text-align: center;">  </div>	<p>2</p> <p><i>Accept structures in either order</i></p> <p><i>For M1, insist on bonds either side of <math>\text{CH}_2</math> groups</i></p> <p><i>For M1 ignore "n" and brackets</i></p> <p><i>Insist on C-O bonds in M2 structure.</i></p>

**General principles applied to marking CHM3/W papers**

It is important to note that the guidance given here is generic and specific variations may be made at individual standardising meetings in the context of particular questions and papers.



Curly arrows should originate either from a lone pair of electrons or from a bond. Each of the following representations **should not gain credit**.

1. The absence of a radical dot in a free radical substitution should be penalised **once only** within a clip.
2. The use of double-headed arrows or the incorrect use of half-headed arrows in free-radical mechanisms should be penalised **once only** within a clip.

**B. Structures**

1. Bonds should be drawn correctly between the relevant atoms. For example, if candidates show the alcohol functional group as C-H-O, they should be penalised **on every occasion**.
2. Some latitude should be given to the representation of C-C bonds in structures, given that CH<sub>3</sub>— is considered to be interchangeable with H<sub>3</sub>C— even though the latter would be preferred.
3. Poor presentation of vertical C — C bonds should **not** gain credit.
4. The use of 'sticks' in structures should **not** gain credit. The occasions that this applies will be indicated in the mark scheme.
5. Some examples of formulae for specific compounds which should **not** gain credit are given here

CH<sub>3</sub>COH for ethanal

CH<sub>2</sub>OCH<sub>2</sub> or CH<sub>2</sub>CH<sub>2</sub>O for epoxyethane

CH<sub>3</sub>CH<sub>2</sub>HO for ethanol

OHCH<sub>2</sub>CH<sub>3</sub>

C<sub>2</sub>H<sub>6</sub>O (except when specifically indicated in the mark scheme)

CH<sub>2</sub>CH<sub>2</sub> for ethene

CH<sub>2</sub>.CH<sub>2</sub>

CH<sub>2</sub>:CH<sub>2</sub>

(N.B. Exceptions may be made in the context of balancing equations)

**C. Names**

As a general principle, non-IUPAC names or incorrect spelling or incomplete names should **not** gain credit. Some illustrations are given here.

but-2-ol	
2-hydroxybutane	all should be <b>butan-2-ol</b>
butane-2-ol	
2-butanol	
2-methopropan-2-ol	should be <b>2-methylpropan-2-ol</b>
2-methylbutan-3-ol	should be <b>3-methylbutan-2-ol</b>
3-methylpentan	all should be <b>3-methylpentane</b>
3-mythylpentane	
3-methypentane	
propanitrile	should be <b>propanenitrile</b>
aminethane	should be <b>ethylamine</b> (although aminoethane can gain credit)
2-methyl-3-bromobutane	all should be <b>2-bromo-3-methylbutane</b>
3-bromo-2-methylbutane	
3-methyl-2-bromobutane	
2-methylbut-3-ene	should be <b>3-methylbut-1-ene</b>
difluorodichloromethane	should be <b>dichlorodifluoromethane</b>

**D. Reagents**

The guiding principle is that a reagent is a chemical which can be taken out of a bottle or container. Failure to identify whole reagents **will be penalised**.

cyanide (ion)	should be e.g. potassium cyanide
hydroxide (ion)	should be e.g. sodium hydroxide

**E. Some general guidance on organic structures**

Each of the following **should gain credit** as alternatives to correct representations of the structures.

$\text{CH}_2 = \text{CH}_2$	for	ethene,	$\text{H}_2\text{C}=\text{CH}_2$
$\text{CH}_3\text{CHOHCH}_3$	for	propan-2-ol,	$\text{CH}_3\text{CH}(\text{OH})\text{CH}_3$
$\text{CH}_2\text{OHCH}_2\text{OH}$	for	ethane-1,2-diol	

**F. Incorrect case for element symbol**

The use of an incorrect case for the symbol of an element should be penalised **once only** within a clip. For example, penalise the use of “h” for hydrogen, “CL” for chlorine or “br” for bromine.

**G. The “List principle”**

If a question requires one answer and a candidate gives two answers, no mark is scored if one answer is correct and one answer is incorrect.

There is no penalty, however, if both answers are correct.

N.B. Certain answers are designated in the mark scheme as those which the examiner should “Ignore” and these answers are not counted as part of the list.