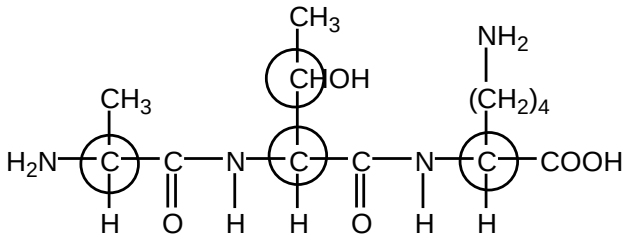
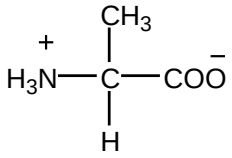
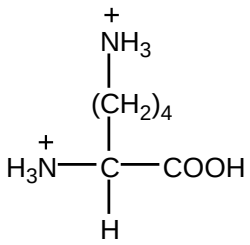
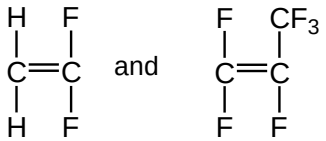
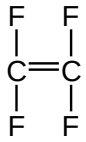
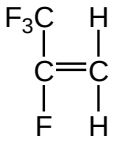
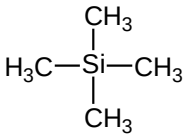
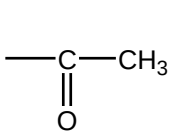
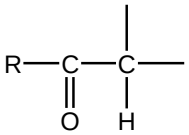
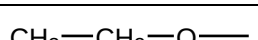
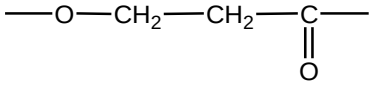
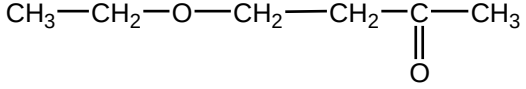
Question 1: N/A

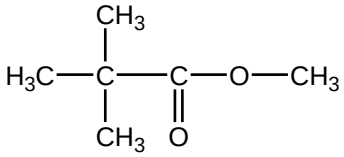
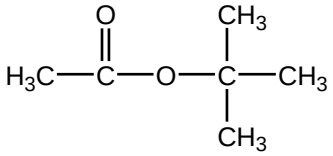
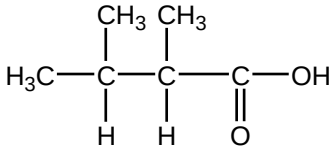
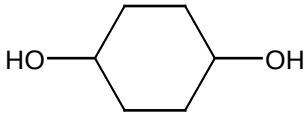
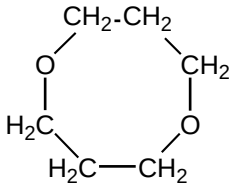
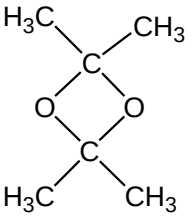
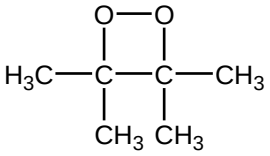
Question 2: N/A

Question	Marking Guidance	Mark	Comments
3(a)(i)	$\underline{3}CH_3OH$ $HOCH_2CH(OH)CH_2OH$	1 1	Not molecular formula
3(a)(ii)	$\rightarrow$ $19CO_2$ + $19H_2O$ $C_{17}H_{35}COOCH_3$ + 27½ or 55/2 O_2	1 1	Or doubled Consequential on correct right-hand side
3(b)(i)	A 0.7 Ethanol 6.4 Water 3.6	1 1 1	
3(b)(ii)	No effect Equal moles on each side of equation ***OR*** V cancels	1 1	If wrong, CE= 0 Ignore moles of gas
3(b)(iii)	M1 $K_c = \frac{[DEM][H_2O]^2}{[A][C_2H_5OH]^2}$	1	Must have all brackets but allow ( )
3(b)(iv)	M2 $\frac{2.1 \times (3.4)^2}{0.85 \times (7.2)^2}$ M3 0.55 (min 2dp) M4 No units	1 1 1	If K_c wrong can only score M4 for units consequential to their K_c working in (b)(iv)

Question	Marking Guidance	Mark	Comments
4(a)(i)		1	These four only
4(a)(ii)		1	Allow $-NH_3^+$ and $^+NH_3-$
4(a)(iii)	<u>2-amino-3-hydroxybutanoic acid</u> Do not penalise commas or missing hyphens	1	Ignore 1 in butan-1-oic acid Penalise other numbers
4(a)(iv)		1	Allow $-NH_3^+$ and $^+NH_3-$

4(b)(i)	Condensation	1	Allow polyester
4(b)(ii)	propan**e**-1,3-diol	1	Must have e Allow 1,3-propan**e**diol
4(c)(i)	Addition	1	Not additional
4(c)(ii)	 and ***OR***  and 	1 for each structure within each pair	Allow monomers drawn either way round Allow bond to F in CF_3
4(d)	c C-C or C-F bonds too strong	1 1	If wrong, CE = 0

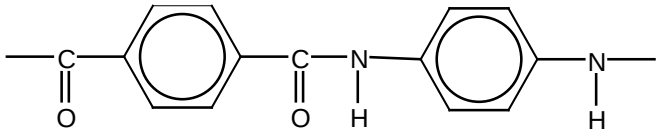
Question	Marking Guidance	Mark	Comments
5(a)(i)	Single/one (intense) peak/signal ***OR*** all H or all C in same environment ***OR*** 12 equiv H or 4 equiv C ***OR*** Upfield / to the right of (all) other peaks ***OR*** well away from others ***OR*** doesn't interfere with other peaks ***OR*** Low bp ***OR*** volatile ***OR*** can easily be removed	2	**Do not allow non-toxic or inert (both given in Q)** Any 2 from three Ignore peak at zero Ignore cheap Ignore non-polar Ignore mention of solubility
5(a)(ii)		1	Allow $Si(CH_3)_4$
5(b)(i)	 or with sticks or 	1	Ignore any group joined on other side of CO Ignore missing trailing bond Ignore charges
5(b)(ii)	 or with sticks	1	Ignore any group joined on other side of –O– Ignore missing trailing bond Ignore charges as if MS fragment
5(b)(iii)	 or with sticks	1	Ignore missing trailing bond Ignore charges as if MS fragment
5(b)(iv)		1	

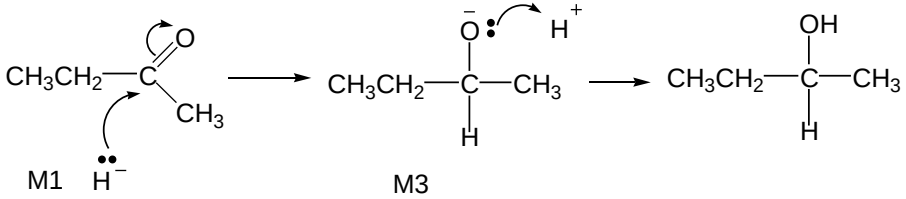
5(c)(i)	Check structure has 6 carbons  	1 1	Allow $(CH_3)_3CCOOCH_3$ or $(CH_3)_3CCO_2CH_3$ Allow $CH_3COOC(CH_3)_3$ or $CH_3CO_2C(CH_3)_3$
5(c)(ii)	Check structure has 6 carbons 	1	Allow $(CH_3)_2CHCH(CH_3)COOH$ or $(CH_3)_2CHCH(CH_3)CO_2H$ Penalise C_3H_7
5(c)(iii)	Check structure has 6 carbons  ***OR*** 	1	Allow  OR 

Question	Marking Guidance	Mark	Comments
6	In each section • If wrong or no reagent given, no marks for any observations; • Penalise incomplete reagent or incorrect formula – but mark observations • Mark each observation independently • Allow *no reaction* for no change / no observable reaction in all three parts, but not *none* or *nothing* • Q says **one test.** If two tests are given, score zero		

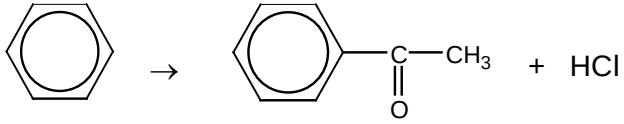
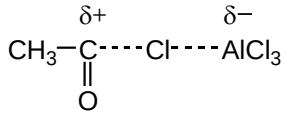
6(a)		$K_2Cr_2O_7$ / H^+	$KMnO_4$ / H^+	Lucas test ($ZnCl_2$ / HCl)	1	Allow acidified potassium manganate and acidified potassium dichromate without oxidation numbers			
	R Primary alcohol	(Orange) goes green Penalise wrong starting colour	(purple) goes colourless / decolourises allow goes brown	No cloudiness	1				
	S Tertiary alcohol	no change / no observable reaction	no change / no observable reaction	Rapid cloudiness	1				
6(b)		Na_2CO_3 / $NaHCO_3$ named carbonate	metal eg Mg	named indicator	1	PCl_5 PCl_3 $SOCl_2$	Named alcohol + HCl / H_2SO_4		
	T ester	no change / no observable reaction	no change / no observable reaction	no effect	1	no change / no observable reaction	no change / no observable reaction		
	U Acid	Effervescence or (CO_2) gas formed	Effervescence or (H_2) gas formed	acid colour	1	Fumes / (HCl) gas formed	Sweet smell		

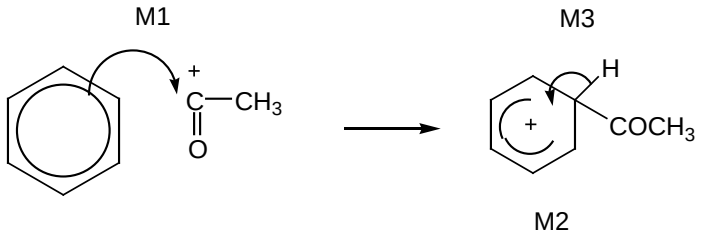
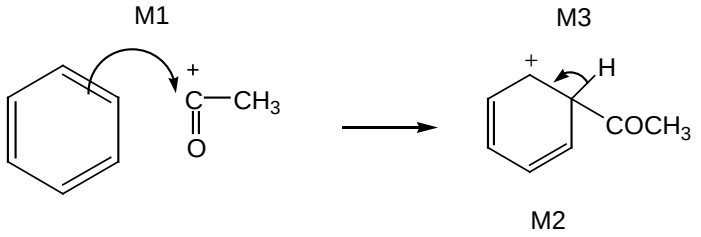
Question	Marking Guidance				Mark	Comments		
	In each section • If wrong or no reagent given, no marks for any observations; • Penalise incomplete reagent or incorrect formula – but mark observations • Mark each observation independently • Allow *no reaction* for no change / no observable reaction in all three parts, but not *none* or *nothing* • Q says **one test.** If two tests are given, score zero							
6(c)		Fehling's / Benedict's	Tollens' / $[Ag(NH_3)_2]^+$	$K_2Cr_2O_7$ / H^+	1	I_2 / NaOH		
	V Ketone	no change / no observable reaction	no change / no observable reaction	no change / no observable reaction	1	Yellow ppt		
	W aldehyde	Red <u>ppt</u>	Silver mirror	(Orange) goes green Penalise wrong starting colour	1	no change / no observable reaction		

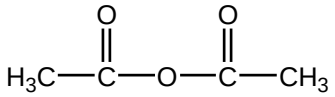
Question	Marking Guidance	Mark	Comments
7(a)	Sn / HCl ***OR*** Fe / HCl not conc H_2SO_4 nor any HNO_3 Ignore subsequent use of NaOH	1	Ignore reference to Sn as a catalyst with the acid Allow H_2 (Ni / Pt) but penalise wrong metal But NOT $NaBH_4$ $LiAlH_4$ Na / C_2H_5OH
	Equation must use molecular formulae $C_6H_4N_2O_4 + 12\,[H]$	1	12[H] and $4H_2O$ without correct molecular formula scores 1 out of 2
	$\rightarrow C_6H_8N_2 + 4H_2O$	1	Allow ….. + $6H_2$ if H_2 / Ni used
		2	Allow –CONH— or —COHN— or —C_6H_4—
			Mark two halves separately : lose 1 each for • error in diamine part • error in diacid part • error in peptide link • missing trailing bonds at one or both ends • either or both of H or OH on ends
			Ignore *n*
7(b)	H_2 (Ni / Pt) but penalise wrong metal	1	NOT Sn / HCl, $NaBH_4$ etc.
	CH_2	1	
	In benzene 120^o	1	
	In cyclohexane $109^o\,28'$ or $109½^o$	1	Allow 108^o - 110^o
	If only one angle stated without correct qualification, no mark awarded		

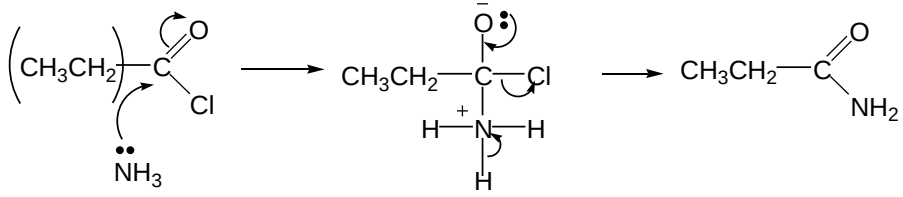
7(c)(i)	Nucleophilic addition M2 M4 for lp, arrow and H+  	1 4	• M2 not allowed independent of M1, but allow M1 for correct attack on C+ • + rather than δ+ on C=O loses M2 • M3 is for correct structure including minus sign but lone pair is part of M4 • Allow C_2H_5 • M1 and M4 include lp and curly arrow • Allow M4 arrow to <u>H</u> in H_2O (ignore further arrows)

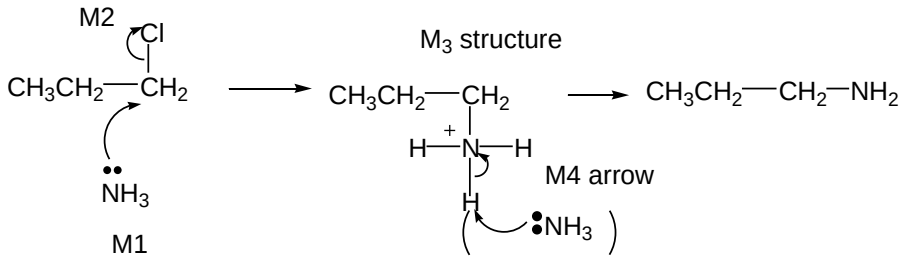
7(c)(ii)	M1 Planar C=O (bond/group) M2 Attack (equally likely) from either side M3 (about product): Racemic mixture formed ***OR*** 50:50 mixture or each enantiomer equally likely	1 1 1	Not just planar molecule Not just planar bond without reference to carbonyl

Question	Marking Guidance	Mark	Comments
8(a)(i)	$CH_3COCl + C_6H_6 \rightarrow C_6H_5COCH_3 + HCl$ ***OR*** $CH_3COCl +$ 	1	Not molecular formulae Not allow $C_6H_5CH_3CO$
	phenylethanone	1	Ignore number 1 in name but penalise other numbers
	$AlCl_3$ can be scored in equation	1	
	$CH_3COCl + AlCl_3 \longrightarrow CH_3\overset{+}{C}O + AlCl_4^-$	1	Allow RHS as  Allow + on C or O in equation but + must be on C in mechanism below Ignore curly arrows in balanced equation even if wrong

8(a)(ii)	Electrophilic substitution  ***OR***  	1 3	• M1 arrow from within hexagon to C or to + on C • + must be on C of CH_3CO in mechanism • + in intermediate not too close to C1 • Gap in horseshoe must be centred approximately around C1 • M3 arrow into hexagon unless Kekule • Allow M3 arrow independent of M2 structure, ie + on H in intermediate loses M2 not M3 • Ignore base removing H for M3

8(b)	Electron pair donor or lone pair donor  (acid) anhydride	1 1 1	Allow donator Allow lone pair used in description of (dative) bond formation Allow $(CH_3CO)_2O$ Allow ethanoic anhydride but not any other anhydride

Question	Marking Guidance	Mark	Comments
10(a)	(Nucleophilic) addition-elimination M2 M3  M1 M4 for 3 arrows and lp propanamide (Ignore -1-)	1 4 1	• Minus sign on NH_3 loses M1(but not M4 also) • M2 not allowed independent of M1, but • allow M1 for correct attack on C+ • + rather than δ+ on C=O loses M2 • **If Cl lost with C=O breaking, max1 for M1** • **M3** for correct structure <u>with charges</u> but lp on O is part of **M4** • only allow **M4** after correct/very close **M3** • For M4, ignore NH_3 removing H^+ but lose M4 for Cl^- removing H^+ in mechanism, • but ignore HCl shown as a product penalise other numbers penalise propaneamide and N-propanamide

10(b)	Nucleophilic substitution  Propylamine (ignore number 1) or propan-<u>1</u>-amine or <u>1</u>-aminopropane (<u>number 1 needed</u>)	1 4 1	• Minus sign on NH_3 loses M1 (not M4 also) • + rather than δ+ on C=O loses M2 • ALLOW SN1 so allow M2 for loss of Cl^- before attack of NH_3 on C+ for M1 • only allow M4 after correct/very close M3 • For M4, ignore NH_3 removing H^+ but lose M4 for Cl^- removing H^+ in mechanism, • but ignore HCl shown as a product penalise other numbers allow <u>1</u>-propanamine
10(c)	electron rich ring or benzene or pi cloud <u>repels</u> nucleophile/ammonia	1 max	Allow • C–Cl bond is short/stronger than in haloalkane • C–Cl is less polar than in haloalkane • resonance stabilisation between ring and Cl

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.

Basic principles

- **Examiners should note that throughout the mark scheme, items that are underlined are <u>required information</u> to gain credit.**
- **Occasionally an answer involves incorrect chemistry and the mark scheme records CE = 0, which means a chemical error has occurred and no credit is given for that section of the clip or for the whole clip.**
-

A. <u>The "List principle" and the use of "ignore" in the mark scheme</u>

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 if both answers are correct.

N.B. Certain answers are designated in the mark scheme as those which the examiner should "Ignore". These answers are not counted as part of the list and should be ignored and will not be penalised.

B. <u>Incorrect case for element symbol</u>

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.

C. <u>Spelling</u>

In general

- The names of chemical compounds and functional groups **must be spelled correctly** to gain credit.
- Phonetic spelling may be acceptable for some chemical terminology.

N.B. Some terms may be required to be spelled correctly or an idea needs to be articulated with clarity, as part of the "Quality of Language" (**QoL**) marking. These will be identified in the mark scheme and marks are awarded only if the QoL criterion is satisfied.

D. Equations

In general

- Equations **must** be balanced.
- When an equation is worth two marks, one of the marks in the mark scheme will be allocated to one or more of the reactants or products. This is independent of the equation balancing.
- State symbols are generally ignored, unless specifically required in the mark scheme.

E. Reagents

The command word "Identify", allows the candidate to choose to use **either** the name or the formula of a reagent in their answer. In some circumstances, the list principle may apply when both the name and the formula are used. Specific details will be given in mark schemes.

The guiding principle is that a reagent is a chemical which can be taken out of a bottle or container. Failure to identify complete reagents **will be penalised**, but follow-on marks (e.g. for a subsequent equation or observation) can be scored from an incorrect attempt (possibly an incomplete reagent) at the correct reagent. Specific details will be given in mark schemes.
For example, **no credit** would be given for

- the cyanide ion or CN^- when the reagent should be potassium cyanide or KCN;
- the hydroxide ion or OH^- when the reagent should be sodium hydroxide or NaOH;
- the $Ag(NH_3)_2^+$ ion when the reagent should be Tollens' reagent (or ammoniacal silver nitrate). In this example, no credit is given for the ion, but credit could be given for a correct observation following on from the use of the ion. Specific details will be given in mark schemes.

In the event that a candidate provides, for example, **both** KCN and cyanide ion, it would be usual to ignore the reference to the cyanide ion (because this is not contradictory) and credit the KCN. Specific details will be given in mark schemes.

F. Oxidation states

In general, the sign for an oxidation state will be assumed to be positive unless specifically shown to be negative.

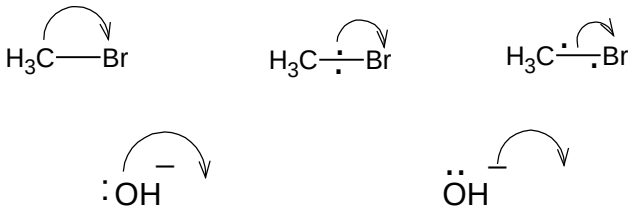
G. Marking calculations

In general

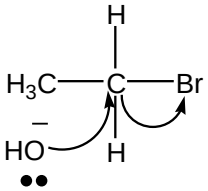
- A correct answer alone will score **full marks** unless the necessity to show working is specifically required in the question.
- An arithmetic error may result in a one mark penalty if further working is correct.
- A chemical error will usually result in a two mark penalty.

H. Organic reaction mechanisms

Curly arrows should originate either from a lone pair of electrons or from a bond.
The following representations should not gain credit **and will be penalised each time** within a clip.



For example, the following would score zero marks



When the curly arrow is showing the formation of a bond to an atom, the arrow can go directly to the relevant atom, alongside the relevant atom or **more than half-way** towards the relevant atom.

In free-radical substitution

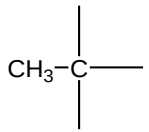
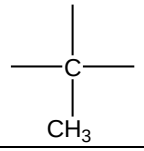
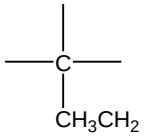
- The absence of a radical dot should be penalised **once only** within a clip.
- The use of double-headed arrows or the incorrect use of half-headed arrows in free-radical mechanisms should be penalised **once only** within a clip

In mass spectrometry fragmentation equations, the absence of a radical dot on the molecular ion and on the free-radical fragment would be considered to be two independent errors and both would be penalised if they occurred within the same clip.

I. Organic structures

In general

- Displayed formulae must show all of the bonds and all of the atoms in the molecule, but need not show correct bond angles.
- Bonds should be drawn correctly between the relevant atoms. This principle applies in all cases where the attached functional group contains a carbon atom, e.g nitrile, carboxylic acid, aldehyde and acid chloride. The carbon-carbon bond should be clearly shown. Wrongly bonded atoms will be penalised **on every occasion**. (see the examples below)
- The same principle should also be applied to the structure of alcohols. For example, if candidates show the alcohol functional group as C ─ HO, they should be penalised **on every occasion**.
- Latitude should be given to the representation of C ─ C bonds in alkyl groups, given that CH_3─ is considered to be interchangeable with H_3C─ even though the latter would be preferred.
- Similar latitude should be given to the representation of amines where NH_2─ C will be allowed, although H_2N─ C would be preferred.
- Poor presentation of vertical C ─ CH_3 bonds or vertical C ─ NH_2 bonds should **not** be penalised. For other functional groups, such as ─ OH and ─ CN, the limit of tolerance is the half-way position between the vertical bond and the relevant atoms in the attached group.
  By way of illustration, the following would apply.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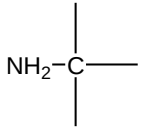	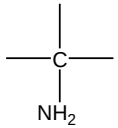	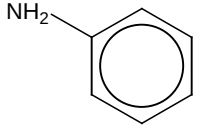	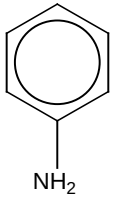	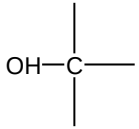	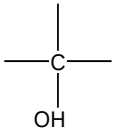
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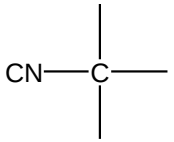	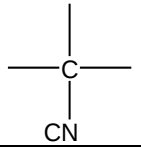	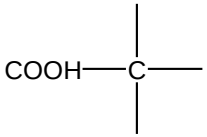	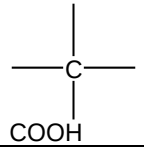	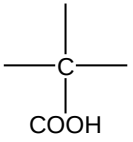	
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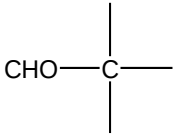	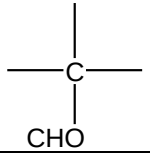	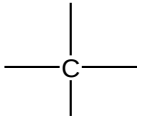	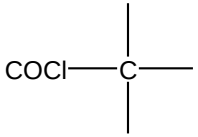	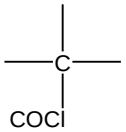	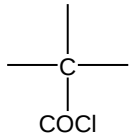
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- In most cases, the use of "sticks" to represent C ─ H bonds in a structure should **not** be penalised. The exceptions will include structures in mechanisms when the C ─ H bond is essential (e.g. elimination reactions in haloalkanes) and when a displayed formula is required.
- Some examples are given here of **structures** for specific compounds that should **not** gain credit

CH_3COH	for	ethanal
CH_3CH_2HO	for	ethanol
$OHCH_2CH_3$	for	ethanol
C_2H_6O	for	ethanol
CH_2CH_2	for	ethene
$CH_2.CH_2$	for	ethene
$CH_2:CH_2$	for	ethane

N.B. Exceptions <u>may</u> be made in the context of balancing equations

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

$CH_2 = CH_2$	for	ethene, $H_2C=CH_2$
$CH_3CHOHCH_3$	for	propan-2-ol, $CH_3CH(OH)CH_3$

J. Organic 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	should be **butan-2-ol**
2-hydroxybutane	should be **butan-2-ol**
butane-2-ol	should be **butan-2-ol**
2-butanol	should be **butan-2-ol**
2-methpropan-2-ol	should be **2-methylpropan-2-ol**
2-methylbutan-3-ol	should be **3-methylbutan-2-ol**
3-methylpentan	should be **3-methylpentane**
3-mythylpentane	should be **3-methylpentane**
3-methypentane	should be **3-methylpentane**
propanitrile	should be **propanenitrile**
aminethane	should be **ethylamine** (although aminoethane can gain credit)
2-methyl-3-bromobutane	should be **2-bromo-3-methylbutane**
3-bromo-2-methylbutane	should be **2-bromo-3-methylbutane**
3-methyl-2-bromobutane	should be **2-bromo-3-methylbutane**
2-methylbut-3-ene	should be **3-methylbut-1-ene**
difluorodichloromethane	should be **dichlorodifluoromethane**