

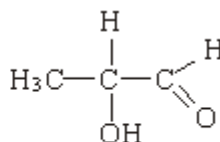
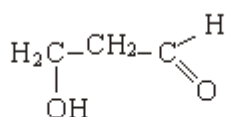
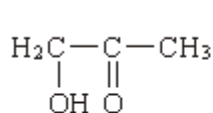
M1. (a) X (O-H) (alcohols)
penalise acid or missing "alcohol"

1

Y C=O

allow carbonyl

1

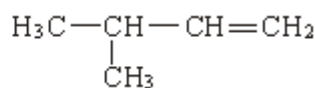
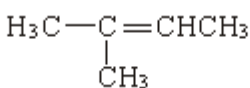
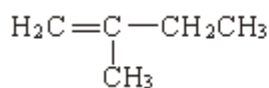


A

NOT acid

4

(b)

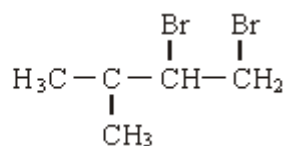
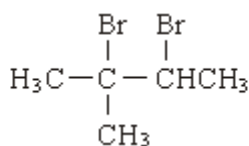
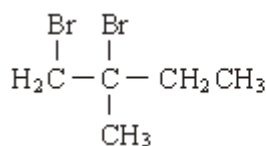


Allow conseq dibromocompounds following incorrect unbranched alkenes

NOT allow dibromocompound consequent on a duplicate alkene

NOT allow monobromocompounds if HBr added

3



3

6:3:1 either next to correct structure or to none

1

Allow a mark for identifying correct dibromocompound with three peaks even if integration ratio is wrong

1

if 6:3:1 missing or wrong, no marks for splitting

Only award a mark for splitting if it is clear which integration number it refers to

6 singlet or drawn

1

3 doublet or drawn	1
1 quartet/quadruplet or drawn	1
	(max 10 marks)

[16]

M2.D

[1]

M3.C

[1]

M4.	(a)	(i)	fractional distillation or fractionation	1
		(ii)	C_9H_{20} <u>only</u>	1
		(iii)	$C_{11}H_{24} + 17O_2 \rightarrow 11CO_2 + 12H_2O$	1
		(iv)	$C_{11}H_{24} + 6O_2 \rightarrow 11C + 12H_2O$	1
	(b)	(i)	$C_{10}H_{22} \rightarrow C_3H_6 + C_7H_{16}$	1
		(ii)	correctly drawn structure of methylpropene (insist on clearly drawn C-C and C=C bonds)	1
	(c)	Any <u>two</u> from		
		o	chemically similar or chemically the same or react in the same way	

- o same functional group
- o same general formula
- o differ by CH₂
(penalise same molecular formula or same empirical formula)

2

[8]

M5.C

[1]

- M6.** (a) (base) elimination
(penalise other words before 'elimination' e.g. nucleophilic)

1

M1: curly arrow from lone pair of electrons on oxygen of hydroxide ion
(insist on a lone pair of electrons on the oxygen atom and a negative charge, but only credit this mark if the attack is to a correct H atom)

1

M2: curly arrow from the middle of the C-H bond to the middle of the C-C bond

1

(only credit this mark if the arrow originates from the correct C-H bond and if an attempt has been made at M1)

M3: curly arrow from the middle of the C-Br bond towards/alongside the Br atom

(credit M3 independently unless the bond breaking is contradicted by an additional arrow)

(penalise curly arrow if the C-Br has a formal positive charge)

(credit full marks for an E1 mechanism, with M2 awarded for a correct curly arrow on the correct carbocation)

(award a maximum of two marks for either an incorrect haloalkane or an incorrect organic product)

(maximum 2 marks for use of 'sticks' for the haloalkane, unless RE from 2(b), when credit can be given)

- (b) (i) **M1:** compounds with the same structural formula 1
- M2:** but the bonds/groups/atoms have different spatial arrangements or orientation or configuration/are arranged differently in space/3D
(ignore reference to the same molecular formula for M1) 1
- (ii) **M1:** correct structural representation for cis-but-2-ene and its name or its identification as the cis isomer 1
- M2:** correct structural representation for trans-but-2-ene and its name or its identification as the trans isomer
(accept representations which are 90° to linear)
(award one mark for two correct structures but either wrong/no names)
(maximum 1 mark for an incorrect alkene) 1
- (iii) geometric(al) or cis-trans 1
- (c) nucleophile or electron pair donor
(penalise 'base') 1
- (d) $\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{Br} + 2\text{NH}_3 \rightarrow \text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{NH}_2 + \text{NH}_4\text{Br}$
(M1 correct product)
(M2 balanced equation using 2NH₃ and leading to NH₄Br)
(penalise M1 for use of C₄H₉NH₂ or for incorrect haloalkane, but allow consequent correct balancing of equation with 2 moles of ammonia) 2
- (1-)butylamine
(credit 1-aminobutane and butyl-1-amine)
(award QoL mark for correct spelling) 1

[13]

M7.B

[1]

M8. (a) M1: $\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{OH}$; 1

M2: $\text{CH}_3\text{CH}(\text{OH})\text{CH}_2\text{CH}_3$;
*(penalise incorrect alcohols in part (a), but mark
consequentially in part (b) and in part (c), if relevant)
(if three alcohols drawn, award MAX. 1 mark)* 1

(b) M1, M2 and M3: Correct structures for butanal, butanone
and butanoic acid;
*(award these structure marks wherever the structures
appear, but insist that the C=O is shown in each structure
and additionally, the C-O in the carboxylic acid* 3

M4: balanced equation for the reaction of butan-1-ol
with [O] to produce butanal and water; 1

M5: balanced equation for the reaction of butan-1-ol
with [O] to produce butanoic acid and water

OR

balanced equation for the reaction of butanal with [O] to
produce butanoic acid; 1

M6: balanced equation for the reaction of butan-2-ol with [O] to
produce butanone and water;
*(Credit condensed structures or molecular formulas in each
equation, provided it is obvious to which reaction the
equation refers) (Insist that whatever formula is used in each
equation that it is a conventional representation of the
compound; for example penalise $\text{CH}_3\text{CH}_2\text{CH}_2\text{COH}$ for
butanal)* 1

- (c) M1: Correct structure for 2-methylpropan-2-ol;
M2: 2-methylpropan-2-ol

1

OR

methylpropan-2-ol;

(penalise on every occasion in parts (a) and (c), structures for the alcohols that are presented with the alcohol functional group as C-H-O)

1

[10]

M9.(a) 2-bromobutane;

1

- (b) Elimination;

(penalise "nucleophilic" OR "electrophilic" before the word "elimination")

1

M1: curly arrow from lone pair on oxygen of hydroxide ion to H atom on correct C-H adjacent to C-Br;

(penalise M1 if KOH shown as covalent with an arrow breaking the bond)

1

M2: curly arrow from single bond of adjacent C-H to adjacent single bond C-C;

(only credit M2 if M1 is being attempted to correct H atom)

1

M3: curly arrow from C-Br bond to side of Br atom;

*(credit M3 independently unless arrows contradict)
(Credit possible repeat error from 2(c)(iii) for M3)*

(If the wrong haloalkane is used OR but-1-ene is produced, award MAX. 2 marks for the mechanism)

(If E1 mechanism is used, give full credit in which M1 and M2 are for correct curly arrows on the correct carbocation)

- (c) (i) (structural) isomers/hydrocarbons/compounds/they have the same molecular formula, but different structural formulas/different structures; 1
(penalise statements which are not expressed in good English and which do not refer clearly to structural isomers i.e. plural)
(penalise statements which refer to “different (spatial) arrangements”)
(credit” different displayed formulas”)
(Q of L mark)
- (ii) Correct structure for but-1-ene; 1

[7]