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Detailed mark scheme

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Time allowed 394 Minutes

Score

/328

Percentage

%

CHEMISTRY

Mark Scheme

OCR AS & A LEVEL

Module 4: Core organic chemistry

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- **1.** (i) 120–130 (**1**)
 - (ii) boiling point increases with increase in *M*r/molecular formula/number of carbon atoms/chain length (1) more intermolecular forces/electrons/surface area/surface interactions/van der Waal forces (1) □
- [3]

2

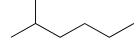
2. $C_{13}H_{28}$

[1]

3. $C_9H_{20} \rightarrow C_7H_{16} + C_2H_4$ (1)

[1]

4. (i) Any branched isomer of heptane with correct name, e.g.



2-methylhexane (1)

2

2

(ii)
$$+ H_2$$

[4]

5. (i) species with an unpaired electron (1)

1

(ii) uv (light)/high temperature/min of 400° C/sunlight (1)

1

(iii) homolytic (fission) (1)

1

2

(iv) $C_4H_{10} + Cl \cdot (1) \rightarrow C_4H_9 \cdot + HCl (1)$

$$C_4H_9^{\bullet} + Cl_2(1) \rightarrow C_4H_9Cl + Cl^{\bullet}(1)$$

[5]

6. (i) 8.72/136.9 = 0.0637 mol (1)

1

(ii) M_r butan-1-ol = 74(.0) (1)

moles =
$$4.28/74.0 = 0.0578 \text{ mol } (1)$$

2

(iii) $0.0578/0.0637 \times 100 = 90.7\%$ (1)

[4]



7. Availability of starting materials:

availability

sugar is renewable because it can be grown (1) ethane is finite because it is obtained by processing of crude oil (1)

energy:

fermentation: energy is required for distillation/ hydration: energy is required to generate steam (1)

atom economy and waste products:

atom economy for fermentation < atom economy hydration (1) In fermentation, CO₂ is produced in addition to ethanol/ethanol is not the only product (1)

In hydration, ethanol is the only product/hydration is an addition reaction (1)

Atom economy of fermentation could be increased by finding a use $CO_2(1)$



Atom economy linked to a chemical equation to show that hydration has 100% atom economy/fermentation has 51% atom economy (1) 7max

[7]

[5]

8. $M_{\rm r} \, {\rm C_7 H_{16}} = 100 \, (1)$ (i) amount = 2000/100 = 20 mol (1)

2

energy saved = $20 \times 4817 = 9634 \text{ kJ}$ (1) (ii)

1

(iii) moles $CO_2 = 7 \times 20 = 140 \text{ mol } (1)$

2

decrease in $CO_2 = 140 \times 24 = 3360 \text{ dm}^3$ (1)

9. structural isomerism:

> structural isomers: same molecular formula, different structural formula (1) structural isomers of but-1-ene: but-2-ene (1) and methylpropene (1)

geometric isomerism

C=C prevents rotation of the double bond (1)

each C in the C=C double bond bonded to 2 different atoms or groups (1)

a clear statement that links non-rotation of the double bond to the idea of groups being trapped on one side of the double bond (1)

cis but-2-ene clearly identified (1) trans but-2-ene clearly identified (1)

[7]



10. 1st bullet

product: CH₃CH₂CHBrCH₂Br (1)

equation: $CH_3CH_2CH=CH_2 + Br_2 \rightarrow CH_3CH_2CHBrCH_2Br$ (1)

products: CH₃CH₂CHBrCH₃ and CH₃CH₂CH₂CH₂Br (1)

(or statement that 2-bromo- is formed)

equation: $CH_3CH=CHCH_3 + HBr \rightarrow CH_3CH_2CHBrCH_3$ (1)

(i.e. for one product)

products: CH₃CH₂CHOHCH₃ and CH3CH2CH2CH2OH (1)

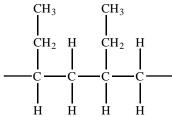
(or statement that 2-ol is formed)

equation: $CH_3CH=CHCH_3 + H_2O \rightarrow CH_3CH_2CHOHCH_3$ (1)

(i.e. for one product)

6

2nd bullet



1 mark for skeleton with two repeat units (1)

1 mark for correct groups on side chains (1)

2

3rd bullet

two (1) (1) from energy from incineration development of biodegradable polymers

cracking of waste polymers

2

[10]

11. separation by (differences in) boiling point

 $C_7H_{16} \rightarrow C_4H_{10} + C_3H_6$

1

1

(i) Any of

1



(ii)
$$C_7H_{16} \rightarrow C_7H_{14} + H_2$$

(or by structural formula)

1

(i) 2,2-dimethylpentane



- (ii) 3-methylhexane, 3,3 dimethylpentane or (3)-ethylpentane in any unambiguous form.
- (iii) 2,2,3-trimethylbutane 1
- (iv) if branched, difficult to pack/less surface interaction/less points of contact 1 less van der Waals' forces/ less intermolecular bonds/less energy needed to boil 1

[10]

2

1

1

- **12.** (a) (i) phosphoric acid/H⁺/sulphuric acid
 - (ii) lone/electron pair of electrons acceptor
 - (b) (i)

- Step 1curly arrow from π-bond to H^+ 1Step 2curly arrow from lone pair on the O^{δ^-} to C+1Step 3curly arrow from O—H bond to O+1
- (ii) catalyst ... no marks because it is **not** consumed/used up in the reaction/owtte 1

[6]

13. (a) 3-chloro(-2-)methylprop-1-ene/1-chloro(-2-)methylprop-2-ene



Backbone of 4 carbons and a reasonable attempt gets 1 mark.

2 **[3]**

14. (a) (i) uv/sunlight/high temperature (range
$$400 - 700$$
 °C) 1

(ii) $Cl_2 \rightarrow 2Cl$ 1

 $C_4H_{10} + Cl \rightarrow HCl + C_4H_9/C_4H_9 \rightarrow 1$

• $C_4H_9/C_4H_9 \rightarrow Cl_2 \rightarrow C_4H_9Cl + Cl \rightarrow 1$

(iii) any two free radicals from (a) (ii) 1

(iv) homolytic (fission) 1

(b) (i) 2,3-dichlorobutane 1
(ii) 1

(iii) any dichlorobutane **except** 2,3-dichlorobutane.

[9]

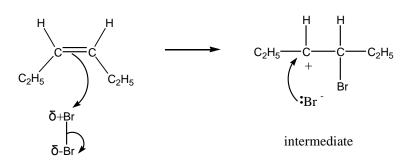


15. Bonding: π -bond formed by overlap of (adjacent) p-orbitals/ π -bond labelled on diagram 1 diagram to show formation of the π -bond 1 minimum allowed for diagram mark **Shape/bond angles:** tetrahedral around the CH₃ 1 bond angle = $109^{\circ}28/(109-110^{\circ})$ 1 trigonal planar around each C in the C=C 1 bond angle = 120° (118-122°) 1 **Cis-trans** cis & trans correctly labelled eg but-2-ene 1 require a double bond because it restricts rotation 1 each C in the C=C double bond must be bonded to two different atoms or groups 1 **QWC** Allow mark for well constructed answer and use of **three** terms like: orbital, tetrahedral, trigonal, planar, rotation, spatial, stereoisomers, geometric 1 [10] 16. (i) (free radical) substitution 1 1-bromohexane, 2-bromohexane and 3-bromohexane 3 (ii)

[4]



17. (a)



curly dipoles shown correctly on the Br–Br and curly arrow from the Br–Br bond towards the Br^{δ} correct intermediate shown 1 curly arrow from the lone pair or the negative charge on the Br $^-$ to the C+

- (b) (i) Hs are diagonal to each other in the *trans/* difference clearly shown in a diagram
 - (ii) (the product is saturated hence) there is no restricted rotation/single bonds allow rotation/because C=C prevents rotation

[6]

18. Recognises that either a catalyst or high temperature (heat is not sufficient) is required

crackingsuitable balanced equation1reformingequation or statement indicating formation of a ring/cycliccompoundsuitable balanced equation with H_2 1(balanced equation showing formation of a ring scores both marks)1

isomerisation suitable balanced equation
The processed products are: 1

The **processed products** are:

• used in fuels/used in petrol

- better /more efficient fuels/increase octane number/rating
- alkenes (from cracking) produce polymers/alcohols
- H₂ used for Haber process/fuels/hydrogenation of oils

QWC SPAG – look for two complete sentence that present a coherent argument

[9]

3



- **19.** (i) C_6H_{10}
 - (ii) C_3H_5 / ecf to (i)
 - (iii) M_r of cyclohexene = 82

$$\% C = (72/82) \times 100 = 88\%$$

87.8% gets 1 mark

ecf to (i) and (ii) for both marks

Alternative calculation based on empirical formula:

Mass of empirical unit = 41, % $C = (36/41) \times 100 = 88\%$

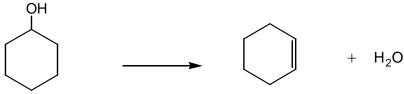
- **20.** H₂
 - Ni/Pt/Pd (catalyst) 1 [2]

[4]

21. (a) (i)



- (ii) $H_2SO_4/Al_2O_3/(hot)$ pumice/ H_3PO_4 1 $(H_2SO_4(aq) \text{ or dil } H_2SO_4 \text{ loses the mark})$
- (iii) 1



 $C_6H_{11}OH \ / \ C_6H_{12}O \rightarrow C_6H_{10} + H_2O$



(b) (i)

OH

also allow

OH

CI

CI-alcohol

from the diol allow from the Cl-alcohol allow

[6]

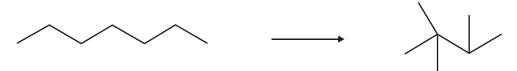
- 22. (a) (i) compound/molecule containing hydrogen and carbon only 1 (ii) $C_{10}H_{22}$ 1 $C_5H_{11} \left\{ ecf \ from \ (ii) \right\}$ (iii) 1 (b) (i) (a particle that) contains/has a single/unpaired electron 1 (ii) UV (light) /sunlight/high temp 1 (iii) homolytic (fission)/ homolysis (iv) $C_{12}H_{26} + Cl \bullet \rightarrow \bullet C_{12}H_{25} + HCl$ 1 (the dot for the free radical does not have to be on the C) $\bullet \mathrm{C}_{12}\mathrm{H}_{25} + \mathrm{C}l_2 \to \mathrm{C}_{12}\mathrm{H}_{25}\mathrm{C}l + \mathrm{C}l \bullet$ 1 (v) 1
 - (c) (i) $C_{12}H_{26} \rightarrow 2C_2H_4 + 1C_8H_{18}$ 2

 (1 mark for correct formula of octane or ethene)

 (ii) octane/ ecf from (c) (i) 1



(d) (i)



1 mark for correct reagent and 1 mark for correct product.

1 mark for any unambiguous formula of cyclohexane 1 (ii)

1 mark for 1H₂ but check that formula of heptane is correct/equation balanced.

 $C_7H_{16} \longrightarrow C_7H_{14} + H_2$ gets 1 marks

23. 1 (i) alkene (a) bromine 1 decolourises 1

> (ii) 3-methylhex-2-en-1-ol/ 1-hydroxy-3-methylhex-2-ene

[4]

1

[16]

2



24. margarine

Ni catalyst 1
hydrogen/ hydrogenated 1
unsaturated vegetable oil/fat 1

poly(propene)

equation

two repeat units

(Ziegler) catalyst / high temp/heat/use of an initiator

Problems with disposal

non-biodegradable/don't decompose/not broken down by bacteria etc 1
when burnt produces toxic fumes 1

Future methods of disposal

recycling (to produce new polymers)

incineration for energy (production)

1

cracking/owtte (to produce useful organic molecules)

use gas scrubbers to reduce toxic fumes

any two

max = 9

QWC

Answer is well organised/structure and using at least three of:

catalyst, hydrogenation, addition polymerisation, Ziegler, incineration, feedstock, recycling, non-biodegradable, initiator, monomer, unsaturated.

in the correct context.

[10]

1



- 25. (a) octane, 400 +/- 5 1 hexadecane. 545 +/- 5 1 if °C penalise once.
 - (b) fractional distillation 1
 - (c) (i)
 - (ii) 2-methylpentane 1
 - (iv) the more branching/the shorter the chain... the lower the boiling point/
 less energy needed to separate the molecules 1
 long chain have greater surface area/surface interactions/more VdW forces
 or converse argument about short/branched chains. 1
 - (d) (i)

(iii)

C, B and A

$$\operatorname{not} \operatorname{just} C_6 H_{12} \qquad \operatorname{or} \quad \operatorname{H}_2 C \subset \operatorname{C}_{H_2} \qquad \operatorname{or} \quad \operatorname{H} \subset \operatorname{C}_{H_2} \subset \operatorname{H}$$

- (ii) $C_6H_{14} \rightarrow C=H_{12} + H_2$
- (iii) <u>better fuels/more volatile/lower boiling point/reduces knocking/increases octane rating/used as (petrol) additives</u>

 1
- (e) (i) M_r of $(CH_3)_3COH = 74$ 1 % oxygen = $(16/74) \times 100 = 21.6$ %
 - (ii) $(CH_3)_3COH + 6O_2 \rightarrow 4CO_2 + 5H_2O$ 1 mark for CO_2 and H_2O only

[16]

2



26. (a) (i) C_5H_8

(i) C_5H_8 (ii) C_5H_8

1

1

1

[6]

[4]

[4]

(b) (i) Ni/Pt/Pd 1

(ii) 1 mark for C_5H_{12} 1 1 mark for correct balancing 1

(iii)

27. (i) electron/lone pair acceptor 1

(ii)

curly arrow from
$$\pi$$
-bond to $\operatorname{Br}^{\delta+}$

Dipoles on the Br–Br bond and curly arrow from Br–Br bond to $\operatorname{Br}^{\delta-}$

C2H₅

Br δ -

Curly arrow from π -bond to $\operatorname{Br}^{\delta-}$

Curly arrow from π -bond to $\operatorname{Br}^{\delta-}$

Curly arrow from π -bond to $\operatorname{Br}^{\delta-}$

1 1 1

28. (i) M_r of 2-methylpropan-1-ol = 74

moles = 4.44/74 = 0.06

(ii) moles = 5.48/137 = 0.04

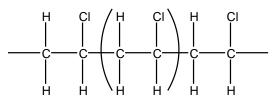
(iii) 66.7%



29. (i) correctly shows three repeat units with 'end bonds'

1
correctly identifies the repeat unit

1



- (ii) harmful/toxic fumes are produced 1
- (iii) recycle/remove HCl by using gas scrubbers or wtte/crack polymers/used a feedstock/ source of fuel (in an incinerator)/developing biodegradable alternatives.

[5]

2

- **30.** (i) $Cl_2 \rightarrow 2Cl \bullet$
 - (ii) uv (light)/high temperature/min of 400 C/sunlight 1
 - (iii) $Cl \bullet + C_6H_{12} \longrightarrow C_6H_{11} \bullet + HCl$ $C_6H_{11} \bullet + Cl_2 \longrightarrow C_6H_{11}Cl + Cl \bullet$
 - (iv) react with each other/suitable equation
 solvent W = water/aqueous/aqueous ethanol
 solvent X = ethanol/alcohol

[5]

31. Structural/chain/positional isomers have the same molecular formula, different structure

but-1-ene/ but-2-ene/ methylpropene / cyclobutane/ methylcyclopropane (any three or two with correct structures and names) 3

4 marks for structural isomerism

Cis-trans / geometric isomerism 1

cis & trans but-2-ene clearly identified 1

C=C prevents rotation 1

each C in the C=C double bond must be bonded to two different atoms or groups 1

4 marks for cis-trans isomerism

QWC: Well organised answer making use of correct terminology to include any **three** from: structural, geometric, cis-trans, molecular formula, restricted, rotation, stereoisomerism, stereoisomers, chain isomerism, positional isomerism, if all isomers are correctly named

[9]

1

32. (a) (i) 24.7/12: 2.1/1: 73.2/35.5



2.06: 2.1: 2.06 1

CHCl 1

(ii) (CHCl = 12 + 1 + 35.5 =) 48.5 1 $48.5 \times 3 = 145.5$ 1

(b) (i) Any two from

CI H C=C H CH₂CI H C=C CH₂CI CH₂CI CH₂CI

1,1,3 -trichloro

3,3,3 -trichloro

2,3,3 -trichloro

1,3,3 -trichloro

2

(ii) 1, 2,3-trichloropropene

(trichloropropene scores 1 mark ✓)

3 marking points:

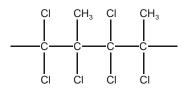
- correct numbers 1, 2,3
- trichloro
- propene/prop-1-ene

any two gets 1 mark

2

2

(c) (i)



1 mark if backbone contains 4 carbons with 'endbonds' and a reasonable attempt has been made e.g used the wrong isomer.... max = 1 mark

(ii) non-biodegradable

1

toxic fumes evolved when burnt

1

1

 $\mathrm{HC}l$ or $\mathrm{C}lullet$ or chlorinated organic compounds such as $\mathrm{COC}l_2$ also evolved when burnt

[13]



33. (i) decolourises 1

(ii)

QWC

$$\begin{array}{c} \text{CH}_3\text{CH}_2 & \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \begin{array}{c} \text{CH}_3\text{CH}_2 & \text{CH}_2\text{CH}_2\text{OH} \\ \text{Br} & \end{array} \\ \text{H} & \begin{array}{c} \text{CH}_3\text{CH}_2 & \text{CH}_2\text{CH}_2\text{OH} \\ \text{Br} & \end{array} \\ \text{H} & \begin{array}{c} \text{CH}_3\text{CH}_2 & \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{H} & \begin{array}{c} \text{CH}_3\text{CH}_2 & \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{Br} & \begin{array}{c} \text{CH}_3\text{CH}_2 & \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{H} & \begin{array}{c} \text{CH}_3\text{CH}_2 & \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{H} & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{OH} \\ \text{CH}_2\text{CH}_2\text{OH} \\ \text{CH}_3\text{CH}_2 & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{CH}_2\text{OH} \\ \text{CH}_3\text{CH}_2 & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{CH}_2\text{OH} \\ \text{CH}_3\text{CH}_2 & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{CH}_2\text{OH} \\ \text{CH}_3\text{CH}_2 & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{CH}_2\text{OH} \\ \text{CH}_3\text{CH}_2 & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2\text{CH}_2\text{OH} \\ \text{CH}_3\text{CH}_2 & \end{array} \\ \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_3\text{CH}_2 & \\ \text{CH}_3\text{CH}_2 & \end{array} \\ \text{CH}_3\text{CH}_3 & \begin{array}{c} \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_3\text{CH}_2 & \\ \text{CH}_3\text{CH}_3 & \end{array} \\ \text{CH}_3\text{CH}_3 & \begin{array}{c} \text{CH}_3\text{CH}_2 & \begin{array}{c} \text{CH}_3\text{CH}_3 & \\ \text{CH}_3\text{CH}_3 & \\$$

curly arrow from C=C bond to bromine 1 dipoles on Br_2 or curly arrow to show movement of bonded pair of electrons 1 intermediate carbonium ion/carbocation 1 curly arrow from lone pair on the Br- ion to carbonium ion (Br^{δ} - loses 1 mark) 1

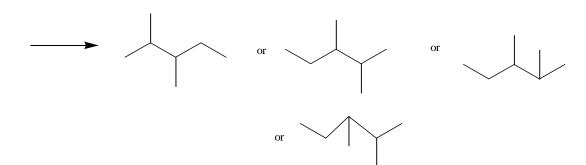
[5]

- 34. identifies the three process as cracking, reforming, isomerisation 1 recognises the need for high temperature or a catalyst 1 equation for cracking equation for isomerisation state that reforming converts chains into rings/cyclic compounds equation for reforming (balanced with H₂ could score two marks) oil is finite/non-renewable ethanol is renewable/sustainable 1 from plants/crops/sugar cane/sugar beet/glucose/sugar/ fermentation 1 $C_2H_5OH + 3O_2 \rightarrow 2CO_2 + 3H_2O$ 1
 - organise relevant information clearly and coherently, using specialist vocabulary when appropriate (minimum of 4 from cracking/isomerisation/reforming/renewable/feedstock/finite/fermentation/non-renewable/sustainable/zeolite/bimetallic catayst/etc)
 - reasonable spelling, punctuation and grammar throughout

[11]



- **35.** (a) C_6H_{14}
 - (b) (i) boiling point increases with increase in M_R /molecular formula/ N° of carbon atoms/chain length 1
 - (ii) more intermolecular forces/electrons/surface area/surface interactions/van der Waal forces1
 - (iii) 120 130 °C 1 [4]
- **36.** (i) $C_9H_{20} \longrightarrow C_7H_{16} + C_2H_4$
 - (ii) $C_2H_4 + H_2O \longrightarrow C_2H_5OH$ 1 temperature > 100 °C/ steam 1 phosphoric acid (catalyst) 1
- **37.** (a) (i)



(ii) 85 –98 °C 1



(b) $C_7H_{16} \longrightarrow C_6H_{11}CH_3 / \downarrow + H_2$ He as a product)

$$\left\{ \begin{array}{c} H_2 \text{ as a product} \\ C_7H_{16} \\ \\ \end{array} \right\} \begin{array}{c} \text{either of these} \\ \text{scores 1 mark} \end{array}$$

(c) more efficient fuel/better fuel/ higher octane number/reduces knocking/more volatile/lower boiling points/burn better/burn more easily/quicker ✓

[5]

1

1

- **38.** (a) (i) reaction 1
 - (ii) reaction 4
 - (iii) reaction 3
 - (b) (i) lone pair/electron pair donor 1

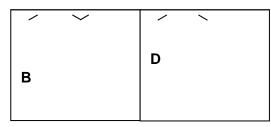
Correct dipole 1

Curly arrow from the O in the OH^- to C in the CH_2

Curly arrow to show movement of bonded pair in the C-Cl bond 1

CI as a product 1

- (c) (i) same molecular formula, different structure/arrangement of atoms. 2 (same formula, different structure.)
 - (ii) 2



(d) (i) addition, (not additional)



- (ii) poly(propene)/ polypropene/ polypro-1-ene, polypropylene 1

 (iii)

 H CH₃ H CH₃

 -C C C C C
 H H H H H
- **39.** (i) decolourises/not clear/not discolours
 - (ii)

$$H_3C-C = CH_2$$
 $H_3C-C + CH_2Br$
 Or
 $H_3C-CHBr-CH_2Br$
 $H_3C-CHBr-CH_2Br$

- curly arrow from C=C to $Br^{\delta+}$
- dipole on Br-Br **and** curly arrow showing movement of bonded pair of electrons
- correct intermediate/carbonium ion/carbocation **and** curly arrow from Br to C+
- 1, 2-dibromopropane as product 1
- 40. CH₃CBr₂CH₃ 1
 CH₃CHBrCH₂Br 1
 - CH₃CH₂CHBr₂
 (CH₃CHBrCH₂Br has a chiral centre, hence optical isomers of
 - 1, 2-dibromopropane are acceptable but must be drawn with 'wedge-shape' bonds and be non-superimposable mirror images)

1

[5]

[15]



41. (i) unsaturated contains a double/multiple/π bond ✓
hydrocarbon contains hydrogen and carbon only. ✓
(ii) angle a 109 −110° ✓
angle b 117 −120° ✓
1

(iii)



Diagram to show a minimum of 2 carbons, each with a $\sigma\text{-bond}$ and p-orbitals \checkmark

Overlap of adjacent p-orbitals (in words or in diagram) ✓ 2

42. (i) *electrophile*: lone pair (of electrons) acceptor. ✓

(ii)

essential mark intermediate carbocation/carbonium ion, accept primary

/"triangular"/ ✓

essential mark product 🗸

curly arrow from double bond to $Br_2 \checkmark$

curly arrow showing movement of electrons in the Br-Br bond \mathbf{or} the dipole in the Br-Br \checkmark

curly arrow from lone pair of electrons in Br[−] to intermediate ✓

mark any errors first

5 max

[5]

[6]



43. (i) Addition (not additional) ✓

1

(ii) 🗸

1

(iii) ✓

- or but-1-ene
- (iv) Poly(but-1-ene) ✓

[4]

44. (a) (i) alkene ✓

1 1

1

alcohol/hydroxy/hydroxyl ✔

1

(b) (i) I = alkene & II = alcohol... both are needed \checkmark

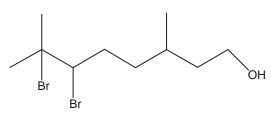
-

(ii) decolourised / colourless ✓

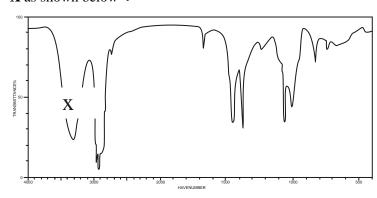
1

(iii) ✓

.



(iv) \mathbf{X} as shown below \checkmark

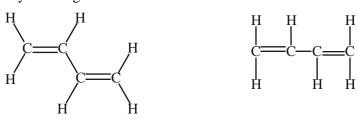




- (c) (i) Ni/Pt/Rh/Pd ✓
 - (ii) compound **B** is $C_{10}H_{22}O$ \checkmark
 - (iii) $C_{10}H_{20}O + H_2 \rightarrow C_{10}H_{22}O \checkmark$

[9]

- **45.** (a) (i) C_4H_{10}
 - (ii) $C_2H_5O \checkmark$
 - (iii) B and E \checkmark 1
 - (iv) A and F \checkmark
 - (b) $(C_4H_9OH \to) C_4H_8 + H_2O \checkmark$
 - (c) any unambiguous formula: ✓



CH₂CHCHCH₂

buta-1,3-diene ✓ 1

name ecf to the structure only if structure above has formula C_4H_6

[7]

46.

1 mark is available if the backbone consists of 4 C atoms and a reasonable attempt has been made $\checkmark \checkmark$

[2]



47. Same molecular formula, different structure /displayed formula/ (a) arrangement of atoms/bonds ✓✓

(Same <u>formula</u>, different structure/displayed formula/arrangement of atoms ✓

- 3-methylbut-1-ene and 2-methylbut-2-ene (b) (any unambiguous structure/formula is acceptable) ✓✓ 2
 - 2-methylbut-1-ene/2-methyl-1-butene ✓ 1
 - (iii) 1



2

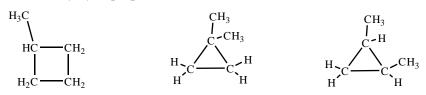
2

[6]

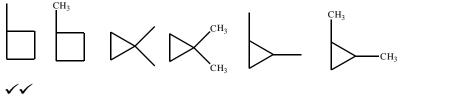
[4]

[4]

48. any two from methylcyclobutane, 1,1-dimethylcyclopropane and (i) 1,2-dimethylcyclopropane



allow



- cyclopentane ✓ 1 (ii)
- (iii) 1



49. (i) homolytic 🗸 1

- $Cl_2 \rightarrow 2Cl \bullet (need \bullet on the Cl... penalise only once in the 3 equations) \checkmark$ 1
- $(C_5H_{10}) + \underline{Cl} \bullet \rightarrow (\bullet C_5H_9) + \underline{HCl} \checkmark$ (iii) 1

 $(\bullet C_5 H_9) + \underline{Cl_2} \rightarrow \underline{C_5 H_9 Cl} + \underline{Cl} \bullet \checkmark$ II 1



50.	(a)	(i)	Alkene/C=C ✓		1
			Alcohol/ROH/hy	droxy/hydroxyl/OH (not OH⁻ or hydroxide) ✓	1
		(ii)	i) One of the C in both C=C is joined to two atoms or groups that are the same ✓		1
	(b)	Obse	rvation	decolourisation (of Br_2) \checkmark	1
		Mole	cular formula	$C_{10}H_{18}OBr_4 \checkmark \checkmark$	2
				$C_{10}H_{18}OBr_2$ gets 1 mark	
	(c)	reage	ent	CH₃COOH ✓	1
		catal	yst	$\mathrm{H_2SO_4/H^+/HC}\mathit{l}$ (aq) or dilute loses the mark \checkmark	1
	(d)	(i)	$C_{10}H_{18}O + 2[O] \rightarrow C_{10}H_{16}O_2 + H_2O \checkmark \checkmark$		2
			1 mark for H ₂ O a	and 1 mark for 2[O]	
		(ii)	The infra-red spectrum was of compound \mathbf{Y}		
	because absorption between 1680 − 1750 cm ⁻¹ indicates a C=O ✓				1
	and the absence of a peak between $2500 - 3300 \text{ cm}^{-1}$ shows the absence				
			of the OH hydrog	gen bonded in a carboxylic acid ✓	1 [12]
51.	Vari	ation ir	n boiling points.	(max = 4 marks)	
J1.	As chain length increases, boiling point increases ✓				
	due to increased number of electrons/ surface area/ more van der Waals forces / intermolecular forces/ more surface interactions 🗸				
	As branching increases, boiling point decreases ✓				
	straight chains can pack closer together/ straight chains have greater surface area/				
	more van der Waals forces /more intermolecular forces/ more surface interactions				
	Ison				
				(produces) branched chain alkanes ✓	1
				equation to illustrate any isomerisation (of octane) \checkmark	1
	<u> </u>	<u> </u>	into a	any one of or or	
				or any other branched isomer of octane	1



Branched chains are better/more efficient fuels/used as additives ✓	1	
because they are more volatile/easier to ignite/burn more easily/higher octane number(rating)/lower boiling points/reduces knocking (pinking) ✓	1	
QWC mark		
• use of suitable chemical terms such as van der Waals, intermolecular forces/intermolecular bonds/volatile/knocking/pinking/pre-ignition		
• reasonable spelling, punctuation and grammar throughout ✓	1	[9]