Alkanes: Formulae, Reactions & Structure

Mark Scheme 2

Level	International A Level
Subject	Chemistry
Exam Board	Edexcel
Торіс	The Core Principles of Chemistry
Sub Topic	Alkanes: Formulae, Reactions & Structure
Booklet	Mark Scheme 2

Time Allowed:	54 minutes
Score:	/45
Percentage:	/100

Grade Boundaries:

A*	А	В	С	D	E	U
>85%	'77.5%	70%	62.5%	57.5%	45%	<45%

Question Number	Acceptable Answers	Reject	Mark
1(a)	Fractional distillation		1
	Both words needed		

Question Number	Acceptable Answers	Reject	Mark
1(b)(i)	C ₉ H ₂₀		1

Question Number	Acceptable Answers	Reject	Mark
1(b)(ii)	Correct skeletal formula (1) Correct name for the structure drawn providing that the structure is a branched-chain isomer of C ₉ H ₂₀	Structural or displayed formula	2
	ALLOW Correct name, even if structural or displayed formula has been drawn (1)		
	and names 2-methyloctane 3-methyloctane		
	4-methyloctane		

Question Number	Acceptable Answers	Reject	Mark
1(c)(i)	$C_{15}H_{32} \rightarrow C_{13}H_{28} + C_2H_4$ IGNORE State symbols, even if incorrect ALLOW Correct structural OR displayed OR skeletal OR mixture of these (as long as unambiguous)		1
	5,		

Question Number	Acceptable Answers	Reject	Mark
1(c)(ii)	Any carbon-carbon bond (in the chain) can break OR The carbon chain can break/split in different places OR Carbon chain is cracked in many places / different places OR $C_{13}H_{28}$ / product will break down further IGNORE 'Molecule can break anywhere' / 'It breaks into smaller molecules' / 'large number of C atoms' / 'bonds break randomly' / 'hydrocarbon chain is long'		1

Question Number	Acceptable Answers	Reject	Mark
1(d)(i)	Two double bonds anywhere on the RING (allow them to be adjacent). e.g.	If any other incorrect structure is included with the final answer	1
		Any 5-valent C atom in structure scores (0)	
	ALLOW One triple bond (instead of two double bonds) BUT not adjacent to a methyl group		
	ALLOW: (ie double bond(s) on side-chain)	If the methyl groups are joined by a bond (0)	
		Benzene ring (0)	

Question Number	Acceptable Answers	Reject	Mark
Question Number 1(d)(ii)	Acceptable Answers NOTE The answer must relate to combustion or burning To promote efficient combustion OR To increase octane number OR To reduce knocking OR Pre-ignition less likely ALLOW To allow smoother burning OR More efficient fuels OR Better burning / fuels easier to burn OR Combust more easily OR Improves combustion ALLOW Reverse argument for straight-chain hydrocarbons IGNORE References to: 'less pollution' / 'burning more cleanly' / 'better fuels' / 'to form alkenes' / 'to form more useful products' /'branched abaias form' /	Reject `Ignition less likely' (0)	1 1
	boiling point / volatility / 'to form H_2'		

(Total for Question 1 = 8 marks)

Question Number	Acceptable Answers	Reject	Mark
2(a)(i)	Curly arrow from double bond towards iodine atom AND curly arrow from the I-Cl bond to the chlorine atom (1) $H \xrightarrow[\delta^+]{\delta^+}_{H} \xrightarrow[\delta^-]{C}_{H}$		3
	Carbocation intermediate (1)		
	$H = CH_{3}$ $H = C = C = C = H$ $H = C = C = H$ $H = C = C$ $H = C$	δ^+ for +	
	correct C ⁺ in the intermediate (1)		
	H CH ₃ ₊ H C C C H H C C C I)CI [−]	δ⁻ for – on Cl⁻	
	NOTE Curly arrow can originate from anywhere on the CI^- ion in the final step. Do not have to have a lone pair of e^- on the CI^- ion		

Question Number	Acceptable Answers		Reject	Mark
2(a)(ii)	Electrophilic Addition ALLOW answers in either order IGNORE 'heterolytic'	(1) (1)		2

Question Number	Acceptable Answers	Reject	Mark
2(a)(iii)	$H = CH_{3}$ $H = C = C = H$ $H = C = C$ $H =$		1

Question Number	Acceptable Answers	Reject	Mark
2(b)(i)	Ultraviolet / UV OR Sun (light) OR Light	Mention of a `catalyst'	1
	ALLOW High temperature / 300°C (minimum)		
	IGNORE Just heat / just radiation / rays		

Question Number	Acceptable Answers		Reject	Mark
2(b)*(ii)	First mark:			7
	(Free) radical substitution	(1)		
	Second mark:			
	Homolytic (fission)	(1)	Heterolytic	
	Third mark:		(fission)	
	Initiation			
	AND			
	$ICI \rightarrow I \bullet + CI \bullet$	(1)		
	BOTH needed for the 3rd mark			
	Fourth mark:			
	Propagation	(1)		
	Fifth and sixth marks:			
	$CH_4 + CI \bullet \rightarrow CH_3 \bullet + HCI$	(1)	$H \bullet$ (the fifth and	
	$CH_3 \bullet + ICI \rightarrow CH_3I + CI \bullet$	(1)	sixth marks cannot be	
	IGNORE $CH_4 + I \bullet \rightarrow CH_3 \bullet + HI$		appears in either propagation step)	
	$CH_3 \bullet + ICI \rightarrow CH_3CI + I \bullet$			
	Seventh mark: $CH_3 \bullet + I \bullet \rightarrow CH_3I$ OR $CI \bullet + CI \bullet \rightarrow CI_2$ OR $CH_3 \bullet + CI \bullet \rightarrow CH_3CI$ OR $CH_3 \bullet + CH_3 \bullet \rightarrow C_2H_6$ OR $I \bullet + CI \bullet \rightarrow ICI$			
	$\begin{array}{llllllllllllllllllllllllllllllllllll$	(1)		
	IGNORE Any INCORRECT termination step(s) IGNORE State symbols, even if incorrect Curly arrows / half curly arrows, even incorrect	ı if		

(Total for Question 2 = 14 marks)

Question Number	Acceptable Answers	Reject	Mark
3 (a)(i)	Penalise use of chlorine once only in Q21(a)(i), (ii) and (iii) IGNORE lone pairs of electrons, even if incorrect in Q21(a)(i), (ii) and (iii) ALLOW one slip in the formula of the element if it is correctly given elsewhere in the answer e.g B for Br $Br_2 \rightarrow Br \bullet + Br \bullet / Br_2 \rightarrow 2Br \bullet$ Ignore position of dot Ignore state symbols and curly arrows even if incorrect	Br	1

Question Number	Acceptable Answers	Reject	Mark
3 (a)(ii)	$Br_2 \rightarrow Br^+ + Br^-$ Ignore state symbols and curly arrows even if incorrect	$\mathbf{\delta}^+$ / $\mathbf{\delta}^-$ for the + or –	1

Question Number	Acceptable Answers	Reject	Mark
3 (a)(iii)	(free radical) Br•	Br	2
	NOTE:		
	No TE, except CI•		
	(1)		
	Penalise omission of the dot only once in (a)(i) and (a)(ii)		
	(electrophile) Br ⁺		
	NOTE:		
	No TE, except CI ⁺ (1)		

Question Number	Acceptable Answers Reject	Mark
3 (b)(i)	H H H H H Any branched-chain isomers $ $ $ $ $ $ $ $ $ $ $ $ $ $ $ $ H $ H$ $ H$ $ H$ $ H$ $ H$ $ H$ $ -$ <th>3</th>	3
	(1)	
	$ \begin{array}{c ccccccccccccccccccccccccccccccccccc$	
	(1) Н Н Н Н Н Н	
	 H H Br H H H	
	(1) Isomers can be in any order	
	ALLOW skeletal or structural formulae	

Question Number	Acceptable Answers	Reject	Mark
3 (b)(ii)	Corrosive / toxic / poisonous Allow correct symbols for corrosive or toxic / poisonous	Flammable / `naked flames'	1
	IGNORE harmful / dangerous / irritant / acidic / volatile / any references to state of HBr IGNORE Any precautions taken, EXCEPT those related to flammability		

Question	Acceptable Answers	Reject	Mark
Number	First mark		2
(b)(iii)	Calculation of the $C_6H_{13}Br M_r$ value		-
	and the total of the product Mr		
	EXPECTED		
	164.9 AND 245.8		
	ALLOW		
	165 AND 246		
	(1)		
	Second mark		
	Second mark		
	EXPECTED		
	<u>164.9</u> (x 100%)		
	245.8		
	= 67.08706265(%)		
	-671(0()) to 2.55		
	= 67.1(%) to 3 s.r.		
	ALLOW		
	165 (x 100%)		
	246		
	= 67.07317073 (%)		
	= 67.1(%) to 3 s.f.		
	ALLOW		
	TE from any incorrect M _r value(s)		
	provided answer is not greater than		
	(1)		
	Answer MUST he rounded correctly to		
	3 s.f. for the second mark		
	Correct answer with no working (2)		
	(=)		

Question	Acceptable Answers	Reject	Mark
Number			
3 (c)(i)	$CH_4 + F_2 \rightarrow CH_3F + HF$ IGNORE state symbols, even if	Cl ₂	1
	incorrect	"FI" if used more than once	

Question Number	Acceptable Answers	Reject	Mark
3 (c)(ii)	NOTE Allow reverse argument throughout 1 st Mark Fluorine / F (atom is) smaller (than a Cl atom) (1)	F ₂ / `fluorine molecule '	2
	2 nd Mark		
	(so expect) F—F bond to be shorter (than the CI—CI bond)	Mention of `Intermolecular forces' (no 2nd mark)	
	OR F—F bonding electrons / bond pair / / shared pair closer to (both) nuclei		
	OR (so) attraction between nuclei and bonding electrons / bond pair / shared pair expected to be stronger (1)		
	IGNORE Any references to the strengths of the F-F and/or CI-CI bonds		
	Any references to the `repulsion between nuclei'		
	Any references to 'shielding' / 'Charge density' / 'Electronegativity' / outer electrons		

Question Number	Acceptable Answers	Reject	Mark
3 (c)(iii)			2
	Shared pair of electrons shown (1)		
	The remaining six electrons on each F		
	(1)		
	NOTE Can be dots or crosses – only total number of electrons matters		
	Circles not required		
	IGNORE Two inner-shell electrons		
	ALLOW `FI' or F symbol missing		

Question Number	Acceptable Answers	Reject	Mark
3 (c)(iv)	'Repulsion between electrons' scores (1) BUT	Just repulsion between bonding / shared electrons	2
	'Repulsion between lone pairs (of electrons)' scores (2) ALLOW		

Question Number	Acceptable Answers	Reject	Mark
3 (c)(v)	UV (light) / (sun) light / heat / energy required to break CI—Cl bond OR UV (light) / (sun) light / heat / energy required to form CI•		1
	OR F—F requires less energy to break OR F—F requires less energy to form F•		
	IGNORE Just F_2 more reactive (than Cl_2)		
	Just F—F bond is weaker (than CI—CI)		
	Just F—F bond energy is lower (than CI—CI)		

Question Number	Acceptable Answers	Reject	Mark
3 (d)	Mark independently H ₅ C ₂ H ₅ C		3
	For both arrows in initial step	Half-arrow(s)	
	Allow upper arrow as in diagram or directly to Br atom (1)	Incorrect polarities Full-charges on Br ₂	
	Second mark:		
	Carbocation intermediate (1)		
	Third mark:	Half-arrow(s) δ⁻ instead of the full – sign	
	Arrow from anywhere on the bromide ion to the C or to the + sign on the intermediate (1)	on the Br ⁻	
	Lone pair(s) on Br ⁻ not required		

Question Number	Acceptable Answers	Reject	Mark
3(e)(i)	$H \qquad C_2H_5 \qquad C_2H_5 \qquad C_2H_5 \qquad OR \qquad C_2H_5 \qquad OR \qquad C_2H_5 \qquad C_2H_$		1
	Diagram clearly shows that H atoms are diagonal to each other in the <i>E</i> - isomer/correct relative positions of hydrogen atoms and ethyl groups ALLOW Skeletal or displayed formula		

Question Number	Acceptable Answers	Reject	Mark
3(e)(ii)	EITHER		1
	Rotation around C—C bond (in product molecule)		
	OR Double bond is broken so rotation (is now possible)		
	ALLOW Same carbocation / intermediate formed (so product is the same)		
	IGNORE Comments about optical isomerism		

(Total for Question 3 = 23 marks)