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## Alkanes: Formulae, Reactions \& Structure

## Mark Scheme 2

| Level | International A Level |
| :--- | :--- |
| Subject | Chemistry |
| Exam Board | Edexcel |
| Topic | 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* | A | B | C | D | E | U |
| :--- | :--- | :--- | :--- | :--- | :--- | :--- |
| $>85 \%$ | $77.5 \%$ | $70 \%$ | $62.5 \%$ | $57.5 \%$ | $45 \%$ | $<45 \%$ |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 ( a )}$ | Fractional distillation |  | $\mathbf{1}$ |
|  | Both words needed |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 ( b ) ( i )}$ | $\mathrm{C}_{9} \mathrm{H}_{20}$ |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 1(b)(ii) | Correct skeletal formula <br> Correct name for the structure drawn providing that the structure is a branched-chain isomer of $\mathrm{C}_{9} \mathrm{H}_{20}$ <br> NO TE for name if skeletal formula is incorrect <br> ALLOW <br> Correct name, even if structural or displayed formula has been drawn <br> EXAMPLES of correct skeletal formulae and names <br> 2-methyloctane <br> 3-methyloctane <br> 4-methyloctane | Structural or displayed formula | 2 |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 ( c ) ( i )}$ | $\mathrm{C}_{15} \mathrm{H}_{32} \rightarrow \mathrm{C}_{13} \mathrm{H}_{28}+\mathrm{C}_{2} \mathrm{H}_{4}$ <br> IGNORE <br> State symbols, even if incorrect <br> ALLOW <br> Correct structural OR displayed OR <br> skeletal OR mixture of these (as long as <br> unambiguous) | $\mathbf{1}$ |  |


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

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 ( d ) ( \mathbf { i ) }}$Two double bonds anywhere <br> on the RING (allow them to be adjacent). <br> e.g. | If any other <br> incorrect structure <br> is included with <br> the final answer | $\mathbf{1}$ |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 1(d)(ii) | NOTE <br> The answer must relate to combustion or burning <br> To promote efficient combustion <br> OR <br> To increase octane number <br> OR <br> To reduce knocking <br> OR <br> Pre-ignition less likely <br> ALLOW <br> To allow smoother burning <br> OR <br> More efficient fuels <br> OR <br> Better burning / fuels easier to burn <br> OR <br> Combust more easily <br> OR <br> Improves combustion <br> ALLOW <br> Reverse argument for straight-chain hydrocarbons <br> IGNORE <br> References to: <br> 'less pollution' / 'burning more cleanly' / <br> 'better fuels' / 'to form alkenes' / 'to form more useful products' /'branched chains form' / <br> boiling point / volatility / 'to form $\mathrm{H}_{2}{ }^{\prime}$ | 'Ignition less likely' <br> (0) | 1 |

(Total for Question 1 = 8 marks)

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2(a)(i) | Curly arrow from double bond towards iodine atom AND curly arrow from the $\mathrm{I}-\mathrm{Cl}$ bond to the chlorine atom <br> Carbocation intermediate <br> Curly arrow from the chloride ion to the correct $\mathrm{C}^{+}$in the intermediate <br> NOTE <br> Curly arrow can originate from anywhere on the $\mathrm{Cl}^{-}$ion in the final step. <br> Do not have to have a lone pair of $\mathrm{e}^{-}$on the $\mathrm{Cl}^{-}$ion | $\delta^{+} \text {for }+$ $\delta^{-} \text {for - on } \mathrm{Cl}^{-}$ | 3 |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2(a)(ii) | Electrophilic <br> Addition <br> ALLOW answers in either order <br> IGNORE <br> 'heterolytic' |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 2(a)(iii) | ALLOW <br> Correct structural OR displayed OR <br> skeletal formula OR mixture of these (so <br> long as unambiguous) <br> Eg CH2CICHICH |  |  |
| IGNORE <br> Any name given, even if incorrect | $\mathbf{1}$ |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{2 ( b ) ( i )}$ | Ultraviolet / UV | OR |  |
|  | Sun (light) <br> OR <br> Light | Mention of a <br> 'catalyst' | $\mathbf{1}$ |
|  | ALLOW <br> High temperature / $300^{\circ} \mathrm{C}$ (minimum) <br> IGNORE <br> Just heat / just radiation / rays |  |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 2(b)*(ii) | First mark: |  | 7 |
|  | (Free) radical substitution (1) |  |  |
|  | Second mark: |  |  |
|  | Homolytic (fission) (1) | Heterolytic |  |
|  | Third mark: |  |  |
|  | Initiation |  |  |
|  | AND |  |  |
|  | $\mathrm{ICl} \rightarrow \mathrm{I} \bullet+\mathrm{Cl} \bullet$ (1) |  |  |
|  | BOTH needed for the 3rd mark |  |  |
|  | Fourth mark: |  |  |
|  | Propagation (1) |  |  |
|  | Fifth and sixth marks: |  |  |
|  | $\mathrm{CH}_{4}+\mathrm{Cl} \bullet \rightarrow \mathrm{CH}_{3} \bullet+\mathrm{HCl}$ (1) | $\mathrm{H} \cdot$ (the fifth and |  |
|  | $\mathrm{CH}_{3} \bullet+\mathrm{ICl} \rightarrow \mathrm{CH}_{3} \mathrm{I}+\mathrm{Cl} \bullet$ (1) | cannot be awarded if $\mathrm{H} \bullet$ |  |
|  | IGNORE $\mathrm{CH}_{4}+\mathrm{I} \bullet \rightarrow \mathrm{CH}_{3} \bullet+\mathrm{HI}$ | appears in either propagation step) |  |
|  | $\mathrm{CH}_{3} \bullet+\mathrm{ICl} \rightarrow \mathrm{CH}_{3} \mathrm{Cl}+\mathrm{I} \bullet$ |  |  |
|  | Seventh mark: $\mathrm{CH}_{3} \bullet+\mathrm{I} \bullet \rightarrow \mathrm{CH}_{3} \mathrm{I}$ |  |  |
|  | OR |  |  |
|  | $\mathrm{Cl} \bullet+\mathrm{Cl} \bullet \rightarrow \mathrm{Cl}_{2}$ |  |  |
|  | $\mathrm{CH}_{3} \bullet+\mathrm{Cl} \bullet \rightarrow \mathrm{CH}_{3} \mathrm{Cl}$ |  |  |
|  | OR |  |  |
|  | $\mathrm{CH}_{3} \bullet+\mathrm{CH}_{3} \bullet \rightarrow \mathrm{C}_{2} \mathrm{H}_{6}$ |  |  |
|  | OR |  |  |
|  | $\mathrm{I} \bullet+\mathrm{Cl} \bullet \rightarrow \mathrm{ICl}$ |  |  |
|  | ALLOW |  |  |
|  | $\mathrm{I} \bullet+\mathrm{I} \bullet \rightarrow \mathrm{I}_{2}$ |  |  |
|  | IGNORE |  |  |
|  | Any INCORRECT termination step(s) |  |  |
|  | IGNORE <br> State symbols, even if incorrect |  |  |
|  | Curly arrows / half curly arrows, even if |  |  |

(Total for Question 2 = 14 marks)

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


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{3}$ (a)(ii) | $\mathrm{Br}_{2} \rightarrow \mathrm{Br}^{+}+\mathrm{Br}^{-}$ <br> Ignore state symbols and curly arrows <br> even if incorrect | $\mathbf{\delta}^{+} / \mathbf{\delta}^{-}$for the + or - | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{array}{\|l\|} \hline 3 \\ (a)(i i i) \\ \hline \end{array}$ | (free radical) Br• | Br | 2 |
|  | NOTE: |  |  |
|  | No TE, except Cl• |  |  |
|  | (1) |  |  |
|  | Penalise omission of the dot only once in (a)(i) and (a)(iii) |  |  |
|  | (electrophile) $\mathbf{B r}^{+}$ |  |  |
|  | NOTE: |  |  |
|  | No TE, except $\mathrm{Cl}^{+}$ |  |  |
|  | (1) |  |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 3 (b)(i) |    <br> Isomers can be in any order <br> ALLOW skeletal or structural formulae | Any branched-chain isomers | 3 |


| Question | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| Number | Corrosive / toxic / poisonous | Flammable / 'naked flames' | $\mathbf{1}$ |
| $\mathbf{3}$ (b)(ii) | Allow correct symbols for corrosive or <br> toxic / poisonous |  |  |
|  | IGNORE <br> harmful / dangerous / irritant / acidic <br> lvolatile / any references to state of <br> HBr <br> IGNORE <br> Any precautions taken, EXCEPT those <br> related to flammability |  |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{aligned} & 3 \\ & \text { (b)(iii) } \end{aligned}$ | First mark |  | 2 |
|  | Calculation of the $\mathrm{C}_{6} \mathrm{H}_{13} \mathrm{Br} \mathrm{M}_{\mathrm{r}}$ value and the total of the product Mr |  |  |
|  | EXPECTED |  |  |
|  | 164.9 AND 245.8 |  |  |
|  | ALLOW |  |  |
|  | 165 AND 246 |  |  |
|  | Second mark |  |  |
|  | EXPECTED |  |  |
|  | 164.9 ( $\times 100 \%$ ) |  |  |
|  | 245.8 |  |  |
|  | $=67.08706265(\%)$ |  |  |
|  | $=67.1(\%)$ to 3 s.f. |  |  |
|  | ALLOW |  |  |
|  | $\frac{165}{24}$ ( $\times 100 \%$ ) |  |  |
|  | 246 |  |  |
|  | $=67.07317073$ (\%) |  |  |
|  | $=67.1(\%)$ to 3 s.f. |  |  |
|  | ALLOW |  |  |
|  | TE from any incorrect $M_{r}$ value(s) |  |  |
|  | provided answer is not greater than 100\% |  |  |
|  | (1) |  |  |
|  | Answer MUST be rounded correctly to 3 s.f. for the second mark |  |  |
|  | Correct answer with no working |  |  |
|  | (2) |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{3 ( c ) ( i )}$ | $\mathrm{CH}_{4}+\mathrm{F}_{2} \rightarrow \mathrm{CH}_{3} \mathrm{~F}+\mathrm{HF}$ <br> IGNORE state symbols, even if <br> incorrect | $\mathrm{Cl}_{2}$ | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 3 (c)(ii) | NOTE |  | 2 |
|  | Allow reverse argument throughout |  |  |
|  | $1^{\text {st }}$ Mark |  |  |
|  | Fluorine / $F$ (atom is) smaller (than a Cl atom) | $\mathrm{F}_{2}$ / 'fluorine molecule' |  |
|  | (1) |  |  |
|  | $\mathbf{2}^{\text {nd }}$ Mark |  |  |
|  | Any ONE of:- |  |  |
|  | (so expect) F-F bond to be shorter (than the $\mathrm{Cl}-\mathrm{Cl}$ bond) | Mention of 'Intermolecular forces' (no $\mathbf{2}^{\text {nd }}$ mark) |  |
|  | OR |  |  |
|  | F-F bonding electrons / bond pair / / shared pair closer to (both) nuclei |  |  |
|  | OR <br> (so) attraction between nuclei and bonding electrons / bond pair / shared pair expected to be stronger |  |  |
|  | IGNORE <br> Any references to the strengths of the $\mathrm{F}-\mathrm{F}$ and/or $\mathrm{Cl}-\mathrm{Cl}$ bonds |  |  |
|  | Any references to the 'repulsion between nuclei' |  |  |
|  | Any references to 'shielding' / 'Charge density' / 'Electronegativity' / outer electrons |  |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $\begin{array}{\|l\|} \hline 3 \\ \text { (c)(iii) } \end{array}$ | Shared pair of electrons shown <br> The remaining six electrons on each $F$ atom <br> NOTE <br> Can be dots or crosses - only total number of electrons matters <br> Circles not required <br> IGNORE <br> Two inner-shell electrons <br> ALLOW <br> ' Fl ' or F symbol missing |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{3}$ | 'Repulsion between electrons' scores <br> (c)(iv) | Just repulsion between <br> bonding / shared electrons | $\mathbf{2}$ |
| BUT | 'Repulsion between lone pairs (of <br> electrons)' scores (2) <br> ALLOW <br> 'Non-bonding electrons' for lone pairs |  |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 3 (c)(v) | UV (light) / (sun) light / heat / energy required to break $\mathrm{Cl}-\mathrm{Cl}$ bond <br> OR <br> UV (light) / (sun) light / heat / energy required to form $\mathrm{Cl} \cdot$ <br> OR <br> $\mathrm{F}-\mathrm{F}$ requires less energy to break OR <br> $F-F$ requires less energy to form $F$. <br> IGNORE <br> Just $\mathrm{F}_{2}$ more reactive (than $\mathrm{Cl}_{2}$ ) <br> Just $\mathrm{F}-\mathrm{F}$ bond is weaker (than $\mathrm{Cl}-\mathrm{Cl}$ ) <br> Just $\mathrm{F}-\mathrm{F}$ bond energy is lower (than $\mathrm{Cl}-\mathrm{Cl})$ |  | 1 |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 3 (d) | Mark independently <br> First mark: <br> For both arrows in initial step <br> Allow upper arrow as in diagram or directly to Br atom <br> Second mark: <br> Carbocation intermediate <br> Third mark: <br> Arrow from anywhere on the bromide ion to the C or to the + sign on the intermediate <br> Lone pair(s) on $\mathrm{Br}^{-}$not required | Half-arrow(s) <br> Incorrect polarities <br> Full-charges on $\mathrm{Br}_{2}$ <br> Half-arrow(s) <br> $\boldsymbol{\delta}^{-}$instead of the full - sign on the $\mathrm{Br}^{-}$ | 3 |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 3(e)(i) |  |  |  |

(Total for Question 3 = 23 marks)

