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## Hess's Law

## Mark Scheme 1

| Level | International A Level |
| :--- | :--- |
| Subject | Chemistry |
| Exam Board | Edexcel |
| Topic | The Core Principles of Chemistry |
| Sub Topic | Hess's Law |
| Booklet | Mark Scheme 1 |


| Time Allowed: | $\mathbf{7 6}$ minutes |
| :--- | :--- |
| Score: | $/ 63$ |
| 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 | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1}$ | A |  | 1 |
|  | Incorrect Answers: <br> B - The enthalpy changes are added <br> and not subtracted <br> C - The enthalpy changes are <br> incorrectly doubled <br> D- T enthalpy changes are doubled <br> and added both incorrectly |  |  |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{2}$ | C |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{3}$ | B |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{4 ( a )}$ | C |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 4(b) | D |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 4(c) | C |  | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{5}$ | C |  | $\mathbf{1}$ |

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| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{6}$ | A | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{7}$ | B | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{8}$ | B | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{9}$ | D | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 0}$ | B | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{1 1}$ | A | $\mathbf{1}$ |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 2}$ | B |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 3}$ | B |  | 1 |


| Question <br> Number | Correct Answer | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 4}$ | A |  | 1 |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 15(a) | Diagram similar to: |  | 3 |
|  | Marking point 1 <br> Arrow upwards for first ionisation energy of sodium and correct label on arrow (from correct entities) <br> (1) <br> Marking point 2 <br> Arrow downwards for electron affinity of iodine and correct label on arrow (from correct entities) <br> (1) <br> Marking point 3 <br> Correct entities with states (on horizontal line) <br> I gnore missing electron <br> (1) <br> ALLOW <br> Numerical values for labels on arrows <br> Recognisable symbols for labels on arrows, such as $\Delta H_{I E}, \Delta H_{E A}$ |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :---: | :---: |
| $\mathbf{1 5}$ (b) | $(\mathrm{LE}=107+107+496+288-295=)-703 \mathrm{~kJ} \mathrm{~mol}^{-1}$ |  | 1 |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 5}$ (c) | Energy is required to break bonds <br> (1) <br> In sodium these are metallic <br> bonds/(electrostatic) attractions between <br> metallic cations and the sea of delocalised <br> electrons (1) | 3 |  |
| In iodine these are covalent bonds <br> (between the iodine atoms and London <br> forces) (1) <br> Mark independently |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| 15(d)(i) | (Sodium iodide has) some covalent character <br> / some covalency/some polarisation <br> ALLOW <br> the electron cloud of the iodide ion is distorted <br> Ignore references to <br> Nal being not 100\% ionic/ <br> Nal being just 'covalent' <br> (1) <br> which results in stronger bonding (than purely <br> ionic) <br> (1) <br> Ignore <br> References to standard conditions/expt. error | 2 |  |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 5 ( d ) ( i i )}$ | Diagram with distorted electron density cloud <br> towards the sodium ion | lodine <br> contour line <br> overlaps <br> with sodium <br> contour line | 1 |
|  | Example |  |  |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i )}$ |  | + | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 6 ( a ) ( i i )}$ | From red-brown / red / brown to colourless | Clear/white <br> Orange/yellow/ <br> Orange-brown | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(b)(i) | (Bonds broken =) $612+193=(+) 805$ <br> (Bonds made=) $347+(290 \times 2)=(-) 927$ <br> (1) <br> Enthalpy of reaction $=(805-927=)-122(\mathrm{~kJ}$ $\mathrm{mol}^{-1}$ ) <br> Correct answer with no working scores two marks <br> ALLOW <br> (All bonds broken $=$ ) +4803 <br> (All bonds made $=$ ) -4925 <br> (1) <br> Enthalpy of reaction $=(+4803-4925=)-122(\mathrm{~kJ}$ $\mathrm{mol}^{-1}$ ) (1) <br> Award one mark for ( + ) $122\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ Award one mark for a correct subtraction using one of the correct values above, example $4538-4925=-387\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :---: | :---: |
| $\mathbf{1 6 ( b ) ( i i )}$ | Bond enthalpies are for gaseous compounds <br> and <br> bromine is a liquid / 1,2 dibromobutane is a <br> liquid <br> IGNORE <br> Reference to just 'different states' | 1 |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(b)(iii) | Mechanism drawn similar to <br> Marking point 1 <br> Curly arrow from double bond to Br and curly arrow from $\mathrm{Br}-\mathrm{Br}$ bond to the Br (dipoles not required) <br> (1) <br> Marking point 2 <br> Correct carbocation structure <br> (1) <br> Marking point 3 <br> Curly arrow from anywhere on the bromide ion (including the minus sign) towards the carbocation and the correct product ALLOW TE on primary carbocation <br> (1) <br> Note the bromide ion must have a full negative charge but the lone pair of electrons need not be shown | Incorrect dipole ${ }^{\delta-}-\mathrm{Br}$ | 3 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 16(b)(iv) | 1-bromobutan-2-ol / $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CHOHCH}_{2} \mathrm{Br} /$ <br> ALLOW <br> 2-bromobutan-1-ol/ $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CHBrCH}_{2} \mathrm{OH} /$ <br> ALLOW 2-bromo-1-butanol <br> ALLOW skeletal or structural formulae <br> Penalise contradictory names/formulae | Missing H's | 1 |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( a ) ( i )}$ | $\Delta \mathrm{H}_{2}$ <br> ALLOW $\Delta \mathrm{H}_{2}=\ldots \ldots .$. |  | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( a ) ( i i )}$ | $\Delta H_{5}$ <br> ALLOW $\Delta H_{5}=\ldots \ldots .$. | $\frac{\Delta H_{5}}{2}$ | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( a ) ( \text { iii) }}$ | $\frac{\Delta \mathrm{H}_{6}}{2}$ <br> $\mathrm{OR} \Delta \mathrm{H}_{6} / 2$ OR $\Delta \mathrm{H}_{6} \div 2 \mathrm{OR} 0.5 \Delta \mathrm{H}_{6}$ | $\Delta \mathrm{H}_{6}$ | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 7 ( a ) ( i v )}$ | $\Delta H_{1}$ <br> ALLOW $\Delta H_{1}=\ldots \ldots \ldots$. | $\Delta H_{\mathbf{7}}$ | $\mathbf{1}$ |

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| Question | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(b)(i) | (The energy change / enthalpy change that accompanies / energy released / enthalpy released) the formation of one mole of a( n ionic) compound <br> ALLOW as alternative for compound: lattice /crystal / substance / solid / product <br> from its gaseous ions <br> NOTE <br> 'one mole of gaseous ions' scores $\max$ (1) <br> (ie 2nd mark only available) <br> IGNORE <br> References to 'constituent elements' References to 'standard conditions' <br> ALTERNATIVE RESPONSE <br> If no mark(s) already awarded from above, can answer by giving:- <br> energy change / enthalpy change per mole $\begin{equation*} \mathrm{Sr}^{2+}(\mathrm{g})+2 \mathrm{Cl}^{-}(\mathrm{g}) \rightarrow \mathrm{SrCl}_{2}(\mathrm{~s}) \tag{1} \end{equation*}$ <br> ALLOW <br> Any correct 'generic' equation with state symbols included | 'Energy / enthalpy required' / 'used' <br> 'molecule' no $\mathbf{1}^{\text {st }}$ mark <br> 'gaseous atoms' no $2^{\text {nd }}$ mark | 2 |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17(b)(ii) | [FIRST, check the answer on the answer line IF answer $=\mathbf{- 2 1 5 3}\left(\mathrm{kJ} \mathrm{mol}^{-1}\right)$ then award (2) marks, with or without working] <br> 1st Mark: $\begin{aligned} & \Delta \mathrm{H}_{1}=\Delta \mathrm{H}_{2}+\Delta \mathrm{H}_{3}+\Delta \mathrm{H}_{4}+\Delta \mathrm{H}_{5}+\Delta \mathrm{H}_{6}+\Delta \mathrm{H}_{7} \\ & \mathrm{OR} \\ & \Delta \mathrm{H}_{7}=\Delta \mathrm{H}_{1}-\left[\Delta \mathrm{H}_{2}+\Delta \mathrm{H}_{3}+\Delta \mathrm{H}_{4}+\Delta \mathrm{H}_{5}+\right. \\ & \left.\Delta \mathrm{H}_{6}\right] \end{aligned}$ <br> OR $\begin{align*} \Delta H_{7}=-829- & {[164+550+1064+} \\ & (122 \times 2)+(2 \times-349)] \tag{1} \end{align*}$ <br> 2nd Mark: $\begin{equation*} \Delta \mathrm{H}_{7}=-2153\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right) \tag{1} \end{equation*}$ <br> NOTE: <br> The following answers score (1) mark with or without working <br> $+2153\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> $-2031\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> $-2502\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> $-2380\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> NO OTHER TEs are allowed on an incorrect expression involving $\Delta \mathrm{H}_{7}$ |  | 2 |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17*(c) | (Lattice energy of $\mathrm{MgF}_{2}$ more exothermic than that of NaF because) <br> 1st mark: <br> $\mathrm{Mg}^{2+}$ is smaller (than $\mathrm{Na}^{+}$) <br> ALLOW <br> "Magnesium / Mg is smaller (than sodium / Na )" <br> 2nd mark: <br> $\mathrm{Mg}^{2+}$ higher charge / higher charge density (than $\mathrm{Na}^{+}$) <br> ALLOW <br> Any reference to $\mathrm{Mg}^{2+}$ and $\mathrm{Na}^{+}$in answer for the $2^{\text {nd }}$ mark, unless nuclear charge mentioned <br> 3rd mark: <br> (So electrostatic forces of) attraction between ions stronger in $\mathrm{MgF}_{2}$ (than in NaF ) <br> ALLOW <br> Stronger ionic bonds in $\mathrm{MgF}_{2}$ / stronger ionic bonding in $\mathrm{MgF}_{2}$ <br> OR reverse arguments | No $1^{\text {st }}$ mark if only mention Mg atom or atomic radius <br> " $\mathrm{Mg}^{2+}$ higher nuclear charge" | 3 |

(Total for Question 17 = 11 marks)

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( a )}$ | $\mathrm{C}_{\mathrm{n}} \mathrm{H}_{2 \mathrm{n}}$ |  | $\mathbf{1}$ |
|  | ALLOW <br> Letters other than n |  |  |

ALLOW: (partially) displayed or skeletal formulae throughout Q18(b)
IGNORE: additional incorrect non-organic products

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( b ) ( i )}$ | $\mathrm{CH}_{3} \mathrm{CH}_{3}$ | $\mathrm{C}_{2} \mathrm{H}_{6}$ | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( b ) ( i i ) ~}$ | $\mathrm{ClCH}_{2} \mathrm{CH}_{2} \mathrm{Cl} / \mathrm{CH}_{2} \mathrm{ClCH}_{2} \mathrm{Cl}$ | $\mathrm{C}_{2} \mathrm{H}_{4} \mathrm{Cl}_{2}$ | $\mathbf{1}$ |

ONLY PENALISE ONCE ONLY in (b)(iii) \& (b)(iv) THE CONNECTIVITY BETWEEN C and OH if CLEARLY a $\mathbf{C}$ to $\mathbf{H}$ covalent bond has been drawn

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :---: | :--- | :---: |
| $\mathbf{1 8 ( b ) ( i i i ) ~}$ | $\mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{OH} / \mathrm{CH}_{2} \mathrm{OHCH}_{2} \mathrm{OH}$ | $\mathrm{C}_{2} \mathrm{H}_{6} \mathrm{O}_{2} /$ <br> $\mathbf{O H C H}_{2} \mathrm{CH}_{2} \mathrm{OH}$ | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( b ) ( i v ) ~}$ | $\mathrm{HOCH}_{2} \mathrm{CH}_{2} \mathrm{Br} / \mathrm{CH}_{2} \mathrm{OHCH}_{2} \mathrm{Br}$ | $\mathrm{BrCH}_{2} \mathrm{CH}_{2} \mathrm{Br} /$ <br> $\mathrm{C}_{2} \mathrm{H}_{5} \mathrm{OBr} /$ <br> $\mathrm{C}_{2} \mathrm{H}_{4} \mathrm{Br}_{2}$ | $\mathbf{1}$ |

## PENALISE USE OF Br instead of $\mathbf{C l}$ once only in parts (c)(i) \& <br> (c)(ii) <br> PENALISE missing $H$ atoms from displayed formulae once only in parts (c)(i) \& (c)(ii)

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 8 ( c ) ( i ) ~}$ |  |  |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(c)(ii) |   <br> attack of chloride ion (1) <br> 1st mark: <br> Curly arrow from $\mathrm{C}=\mathrm{C}$ to H (in $\mathrm{H}-\mathrm{Cl}$ ) AND curly arrow from bond in $\mathrm{H}-\mathrm{Cl}$ to the Cl (dipole not reqd) Curly arrows must start from the bonds NOT the atoms | Full + and charges on HCl <br> Incorrect polarity on HCl | 3 |
|  | 2nd mark: <br> Structure of correct secondary carbocation <br> 3rd mark: <br> Curly arrow from anywhere on the chloride ion (including the minus sign) towards the $\mathrm{C}+$ on the carbocation <br> NOTE: The chloride ion must have a full negative charge, but the lone pair of electrons on the $\mathrm{Cl}^{-}$ need not be shown <br> ALLOW: TE on major product given in (c)(i) <br> Skeletal formulae can be used <br> Mark the three points independently | Extra / <br> spare <br> bond <br> dangling <br> from <br> the C+ <br> carbon <br> $\delta$ - on chloride ion instead of $\mathrm{Cl}^{-}$ |  |

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| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(d)(i) |  <br> TWO ' $n$ ' in the equation and a correct formula (molecular or structural) for propene on the left-hand side of the equation <br> One correct repeating unit, with the methyl branch shown <br> ALLOW <br> $\mathrm{CH}_{3}$ fully displayed or just as $\mathrm{CH}_{3}$ <br> BOTH continuation bonds (with or without bracket shown) <br> If $\mathrm{C}=\mathrm{C}$ bond left in polymer on righthand side, then max (1) <br> Mark the three points independently |  | 3 |

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| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :---: |
| $\mathbf{1 8 ( d ) ( i i )}$ | Non-biodegradable |  |  |
|  | IGNORE <br> References to toxicity of <br> poly(propene) / flammability <br> IGNORE <br> Litter / pollution / waste of resources <br> / costs |  |  |
|  | ALLOw <br> People are reluctant to recycle <br> OR <br> Harmful to marine life / harmful to <br> wildlife <br> OR <br> References to 'landfill' <br> OR <br> References to 'incineration' producing <br> toxic fumes/toxic gases / CO / <br> Greenhouse gases <br> OR <br> References to use of energy/fuel <br> used in transport (of waste) <br> OR <br> It takes a long time to degrade |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(e)(i) | Both arrows in the correct direction <br> AND <br> $3 \mathrm{CO}_{2}$ and $3 \mathrm{H}_{2} \mathrm{O}$ in lowest box <br> IGNORE state symbols, even if incorrect <br> IGNORE extra $\mathrm{O}_{2}$ molecules in box or alongside arrows |  | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18(e)(ii) | $\begin{aligned} & \mathbf{1}^{\text {st }} \text { mark } \\ & (-394 \times 3)+(-286 \times 3) \end{aligned}$ <br> OR $\begin{equation*} =-2040\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right) \tag{1} \end{equation*}$ <br> 2nd mark: <br> $\Delta H_{f} \quad=-2040-(-2058)$ $\begin{equation*} =(+) 18\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right) \tag{1} \end{equation*}$ <br> NOTE: <br> The following answers score (1) mark <br> with or without working <br> $-18\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> $(+) 1378\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> $(+) 806\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> $(+) 590\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> -4098 (kJ mol ${ }^{-1}$ ) <br> IGNORE units even if incorrect |  | 2 |

