# Mark Scheme (Results) J une 2010 

## GCE

## GCE Chemistry (6CH04/ 01)

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## Section A (multiple choice)

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


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


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


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


| Question | Correct Answer | Mark |
| :--- | :--- | :--- |
| Number |  |  |
| $\mathbf{4}$ | D | $\mathbf{1}$ |


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


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


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{7 ~ ( a )}$ | $\mathbf{C}$ | $\mathbf{1}$ |


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


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| $\mathbf{7}$ (c) | D | $\mathbf{1}$ |


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


| Question Number | Correct Answer | Mark |
| :---: | :---: | :---: |
| 9 | D | 1 |
| Question <br> Number | Correct Answer | Mark |
| 10 | D | 1 |
| Question Number | Correct Answer | Mark |
| 11 | B | 1 |
| Question Number | Correct Answer | Mark |
| 12 | A | 1 |
| Question Number | Correct Answer | Mark |
| 13 | B | 1 |
| Question Number | Correct Answer | Mark |
| 14 | C | 1 |
| Question Number | Correct Answer | Mark |
| 15 | C | 1 |
| Question Number | Correct Answer | Mark |
| 16 | A | 1 |

## Section B

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7}$ (a)(i) | $5.7 \times 10^{-5} / 5.71 \times 10^{-5} / 5.714 \times 10^{-5} / 0.000057$ <br> IGNORE SF except 1 (ie don't accept $6 \times 10^{-5}$ ) | $\mathbf{1}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7}$ (a)(ii) | $\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}$ : first order / 1 (1) <br> (going from first to second experiment) <br> rate doubles when concentration / number of <br> moles doubles (and [OH-] constant )/ rate and <br> concentration increase in proportion (1) <br> ALLOW 'time halves' instead of 'rate doubles' | $\mathbf{3}$ |  |
| OH'$^{-}$: zero order / 0 <br> and <br> (going from second to third expt) as increase in <br> concentration does not affect rate (and <br> [C4H9Br] constant ) (1) | ALLOW 'doubling in concentration of OH' <br> instead of 'increase in concentration' | ALLOW time increases by the same factor as <br> increase in hydroxide concentration (5/3) <br> May refer to experiment number rather than <br> concentrations |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7}$ (a)(iii) | Rate $=\mathrm{k}\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]$ <br> OR Rate $=\mathrm{k}\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]^{1}\left[\mathrm{OH}^{-}\right]^{0}$ <br> ALLOW k in lower or upper case <br> Rate equation must be consistent with orders in <br> (a)(ii) <br> If no order is given for hydroxide in (ii) mark <br> cannot be given | $\mathbf{1}$ |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17 (a)(iv) | $\begin{aligned} \mathrm{k} & =\frac{2.9 \times 10^{-5}}{0.017} \\ & =1.7 \times 10^{-3} / 1.71 \times 10^{-3} / 1.706 \times 10^{-3} \mathrm{~s}^{-1} \end{aligned}$ <br> ALLOW k=1. $68 \times 10^{-3}$ <br> (value obtained from experiment 2 or 3 ) <br> value of $k$ (1) <br> units (1) stand alone mark <br> ALLOW TE from (a)(iii) <br> IGNORE SF except 1 <br> Rate $=k\left[C_{4} \mathrm{H}_{9} B r\right]^{2}$ gives $k=0.10036 \mathrm{dm}^{3} \mathrm{~mol}^{-1} \mathrm{~s}^{-1}$ <br> Rate $=\mathrm{k}\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]\left[\mathrm{OH}^{-}\right]$gives $\mathrm{k}=1.42 \mathrm{dm}^{3} \mathrm{~mol}^{-1} \mathrm{~s}^{-1}$ <br> ALLOW $\mathrm{k}=1.39 \mathrm{dm}^{3} \mathrm{~mol}^{-1} \mathrm{~s}^{-1}$ <br> (value obtained from experiment 2 or 3 ) <br> Rate $=k\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]\left[\mathrm{OH}^{-}\right]^{2}$ gives $\mathrm{k}=1184.6$ $\mathrm{dm}^{6} \mathrm{~mol}^{-2} \mathrm{~s}^{-1}$ <br> Rate $=\mathrm{k}\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]^{2}\left[\mathrm{OH}^{-}\right]$gives $\mathrm{k}=83.62$ $\mathrm{dm}^{6} \mathrm{~mol}^{-2} \mathrm{~s}^{-1}$ |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 7 ( b )}$ | $\left[\mathrm{OH}^{-}\right]$is (in chemical equation but) not in rate <br> equation / not in rate determining step (so is in <br> a step other than rate determining step) <br> OR <br> Only $\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}$ is in rate equation / rate <br> determining step (so $\mathrm{OH}^{-}$is in a step other than <br> rate determining step) | $\mathbf{1}$ |  |



| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 17 (d) <br> QWC | (Primary and tertiary) carbocation intermediates have different stabilities (1) as (inductive effects of) alkyl groups stabilise tertiary carbocation (1) <br> OR <br> Steric hindrance differs for attack on primary and tertiary carbon (in the molecule) / less space available for attack by $\mathrm{OH}^{-}$on tertiary carbon / more space for attack by $\mathrm{OH}^{-}$on primary carbon (1) <br> as bulky / three alkyl groups obstruct attack (1) | "Tertiary bromoalkanes react by SN1" without further explanation <br> carbocation intermediates have different reactivity <br> steric hindrance in carbocation | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 8}$ (a)(i) | (Acid) hydrolysis | substitution | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 8 ( a ) ( i i )}$ | $\mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{Na}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{Cr}_{2} \mathrm{O}_{7}^{2-}$ <br> Potassium dichromate $((\mathrm{VI})) /$ sodium <br> dichromate((VI)) / dichromate((VI)) ions | Just <br> "dichromate" | $\mathbf{1}$ |
| ALLOW manganate((VII)) ions, etc | Correct <br> formula with <br> wrong name <br> and vice versa <br> Incorrect <br> oxidation <br> number |  |  |

$\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline \mathbf{1 8} \text { (a)(iii) } & \begin{array}{l}\text { Lithium tetrahydridoaluminate/ lithium } \\ \text { aluminium hydride/ LiAlH }\end{array} \text { (in dry ether) }\end{array} \quad$ Just [H $\left.\mathrm{H}^{-}\right] \quad \mathbf{1}$.

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 8}$ (a)(iv) | Methyl butanoate (1) <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{COOH}^{2}+\mathrm{CH}_{3} \mathrm{OH} \rightarrow$ <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{COOCH}_{3}+\mathrm{H}_{2} \mathrm{O}$ (1) <br> ALLOW $\rightleftharpoons$ <br> IGNORE state symbols even if wrong | Methyl <br> butoate | $\mathbf{2}$ |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 8 ( a ) ( v )}$ | $\mathrm{CH}_{3}-\mathrm{CH}_{2}-\mathrm{CH}_{2}-\mathrm{C}^{-}=\mathrm{O}$ <br> Don't penalise undisplayed methyl groups as <br> here. <br> COCl must be displayed as above. | $\mathrm{C}_{3} \mathrm{H}_{7} \mathrm{for}$ <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2}$ | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 8}$ (b)(i) | Nitrogen inert / unreactive / less reactive <br> (than oxygen) <br> OR <br> Oxygen might react with chemicals going <br> through column / sample might oxidise | $\mathbf{1}$ |  |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 18 (b)(ii) | Solubility (in liquid / stationary phase) <br> OR <br> Interaction with liquid / stationary phase <br> OR <br> Interaction between mobile and stationary phase <br> OR <br> Attraction for liquid / stationary phase <br> OR <br> Strength of (named) intermolecular forces <br> OR <br> Adsorption on liquid / stationary phase <br> OR <br> Absorption on liquid / stationary phase | Size of molecule / molar mass <br> Polarity, unless with explanation <br> Boiling point / volatility <br> Viscosity <br> Attraction for carrier gas <br> $J$ ust a named intermolecular force <br> J ust 'retention time' <br> Density | 1 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18 (c)(i) |  <br> OR <br> Ester link including C=O (1) <br> Rest of polymer with oxygens at end correct (1) <br> All H atoms must be shown. <br> PENALISE lack of displayed $\mathrm{C}=0$ once only ACCEPT <br> Without brackets around formula but bonds at end should be shown More than two correct units IGNORE n after brackets |  | 2 |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 8}$ (c)(ii) | Hydrolysis |  | $\mathbf{1}$ |
|  | OR <br> Splits / breaks ester link <br> OR <br> polymer breaks down to monomers <br> OR <br> equation showing hydrolysis | Jolymer down' |  |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 19 (a)(i) | $\left(\mathrm{K}_{\mathrm{p}}=\frac{\mathrm{pCH}_{3} \mathrm{CO}_{2}-\frac{\mathrm{H}}{\mathrm{pCH}_{3}} \mathrm{OH}^{(\mathrm{x})} \mathrm{pCO}}{}\right.$ <br> Partial pressure symbol can be shown in various ways, eg pp, $p_{\mathrm{co}}$, (CO)p, etc <br> ALLOW $p$ in upper or lower case, round brackets <br> IGNORE units | [ ] <br> State symbols given as (I) <br> +in bottom line | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9}$ (a)(ii) | $\mathrm{P} \mathrm{CH} 3 \mathrm{OH}=4.9(\mathrm{~atm})(\mathbf{1})$ <br> $\mathrm{P} \mathrm{CO}=4.9(\mathrm{~atm})(\mathbf{1})$ <br> 1 mark for recognition that pressures are equal <br> IGNORE units | $\mathbf{2}$ |  |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9}$ (a)(iii) | $\mathrm{K}_{\mathrm{p}}=\left((22.2) /(4.9)^{2}\right)$ <br> $=0.925(\mathbf{1})$ <br> atm $^{-1}$ (1) stand alone mark but must match <br> expression used in (a)(iii) <br> OR <br> $9.25 \times 10^{4} \mathrm{~Pa}^{-1} / 92.5 \mathrm{kPa}^{-1}$ (2) <br> ALLOW TE from (a)(i) if inverted and/ or (a)(ii) | Answers to <br> other than 3 <br> significant <br> figures | $\mathbf{2}$ |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9}$ (b)(i) | $\mathrm{CH}_{3} \mathrm{OH}: 3.2$ <br> $\mathrm{CO}: 3.2$ (1) for both values <br> $\mathrm{CH}_{3} \mathrm{CO}_{2} \mathrm{H}: 46.8$ (1) <br> ALLOW TE for moles of ethanoic acid based on <br> numbers of methanol and carbon monoxide <br> used, as long as moles of methanol and carbon <br> monoxide are equal and moles ethanoic acid + <br> moles methanol =50 | $\mathbf{2}$ |  |

$\left.\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline \mathbf{1 9} \text { (b)(ii) } & \left(\frac{46.8 \times 32}{53.2}\right)=28.2 / 28.1504 \text { (atm) } & 28.1 & \mathbf{1} \\ & \text { IGNORE sf except } 1 \\ \text { Value = 28.16 if mol fraction rounded } \\ \text { ALLOW TE from (b)(i) } & \frac{46.8 \times 32}{50}= & 29.95 \text { (atm) }\end{array}\right]$

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9}$ (b)(iii) | exothermic as yield / pp of ethanoic acid / <br> conversion of reactants/ Kp is higher at lower <br> temperature / as equilibrium moves (right) at <br> lower temperature | $\mathbf{1}$ |  |
| ALLOW <br> if partial pressure of ethanoic acid <22.2 atm <br> in(b)(ii), endothermic as yield / pp of ethanoic <br> acid / conversion of reactants/ Kp is lower at <br> lower temperature |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9}$ (c)(i) | No effect <br> and <br> other concentrations change to keep $K_{p}$ <br> constant / K $K_{p}$ is only affected by temperature/ <br> as equilibrium moves (right) to keep K <br> constant / change in pressure does not change <br> $\mathrm{K}_{\mathrm{p}}$ | As K is a <br> constant | $\mathbf{1}$ |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9}$ (c)(ii) | Yield increased to restore fraction / quotient / <br> partial pressure ratio back to K |  |  |
|  | ALLOW (equilibrium moves) to use up the <br> methanol / answers based on entropy or Le <br> Chatelier | Just <br> Coquilibrium <br> moves to the <br> right' |  |
| inadequate explanations scores 1 mark in <br> (c)(ii) |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{1 9}$ (d) | Mark independently <br> Reaction can occur at lower temperature / has <br> lower activation energy / requires less energy <br> (1) <br> less fuel needed / fewer emissions (from fuels) <br> / fewer raw materials needed / less natural <br> resources used (1) <br> OR | Answer based <br> on car exhaust <br> emissions |  |$土$ 2 $\quad$| Enables use of an alternative process with |
| :--- |
| higher atom economy (1) |
| fewer raw materials needed / less natural |
| resources used (1) |$\quad$|  |
| :--- |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ( a ) ( i )}$ | Correct answer with or without working scores <br> $\mathbf{2}$ marks <br> $\left[\mathrm{H}^{+}\right]=\left(1.00 \times 10^{-14} / 0.250\right)=4 \times 10^{-14} \quad$ (1) <br> $\mathrm{pH}=(13.39794=) 13.4$ (1) <br> $\mathbf{O R}$ <br> $\mathrm{pOH}=-\log 0.250=0.602$ (1) <br> $\mathrm{pH}=(13.39794=) 13.4$ (1) <br> ALLOW <br> TE in second mark if error in $\left[\mathrm{H}^{+}\right]$calculation <br> gives pH more than 7 <br> 3 or more sf <br> IGNORE rounding errors e.g. accept 13.39 | $\mathbf{2}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0}$ (a)(ii) | $\left(\mathrm{K}_{\mathrm{a}}=\right) \frac{\left[\mathrm{CH}_{3} \mathrm{COO}^{-}\right]\left[\mathrm{H}^{+}\right]}{\left[\mathrm{CH}_{3} \mathrm{COOH}\right]}$ (1) <br>  $\mathrm{ALLOW}^{\mathrm{H}_{3} \mathrm{O}^{+} \text {instead of } \mathrm{H}^{+}}$ <br> $\frac{\left[A^{-}\right]\left[\mathrm{H}^{+}\right]}{[\mathrm{HA}]}$ if key to symbols given <br> IGNORE state symbols <br> $\left[\mathrm{CH}_{3} \mathrm{COOH}\right]$  | $\mathbf{1}$ |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 20 (a)(iii) | Correct answer with or without working scores 2 marks $\begin{align*} & 1.7 \times 10^{-5}=\frac{\left[\mathrm{H}^{+}\right]^{2}}{0.125}  \tag{1}\\ & \\ & {\left[\mathrm{H}^{+}\right]=1.46 \times 10^{-3}} \\ & \mathrm{pH}=2.84 / 2.8(1) \end{align*}$ <br> no TE from an incorrect $\left[\mathrm{H}^{+}\right]$ |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0}$ (a)(iv) | $\mathrm{pH}=4.8 / 4.77$ (1) <br> $\mathrm{pH}=\mathrm{pK}_{\mathrm{a}} /\left[\mathrm{H}^{+}\right]=\mathrm{K}_{\mathrm{a}}$ (when acid is half <br> neutralized) (1) | $\mathrm{H}^{+}=\mathrm{K}_{\mathrm{a}}$ |  |$\quad$ 2 $\quad$|  |
| :--- |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0}$ (a)(v) | Sigmoid curve starting between pH 2 and 4 <br> (2.8), ending between pH 12 and 14 inclusive <br> (1) <br> with steep rise (may be vertical or gently <br> sloping) of between 3-7 units between pH 6 <br> and 12. Sloping section should not extend over <br> more than $5 \mathrm{~cm}^{3} .(1)$ <br> When 12.5 $\mathrm{cm}^{3}, \mathrm{NaOH}$ added. (1) <br> ALLOW tolerance for grid <br> Reverse curves lose first mark | $\mathbf{3}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0 ~ ( a ) ( v i ) ~}$ | First mark <br> Thymolphthalein more suitable as it changes <br> (from colourless to blue) in steep region of <br> titration (pH 8.3 to 10.6)/ at the equivalence <br> point / at the end point <br> OR <br> thymolphthalein has pH range in steep region <br> of titration (1) | $\mathbf{2}$ |  |
| Second mark <br> Methyl yellow changes (from red to yellow at <br> pH 2.9 to 4) before equivalence point / before <br> the end point / doesn't change in steep section <br> OR <br> Methyl yellow has pH range before / outside <br> steep region of titration (1) | ALLOW 'Thymolphthalein more suitable as it <br> changes at the equivalence point but methyl <br> yellow does not.' This scores 2 marks | OR <br> OR <br> First mark <br> pK in $\pm 1$ must lie within vertical region on <br> titration curve (1) <br> Second mark <br> hence thymolphthalein is suitable and methyl <br> yellow is not (1) |  |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 0}$ (b) | Sodium ethanoate/ $\mathrm{CH}_{3} \mathrm{COONa}$ <br> Potassium ethanoate $/ \mathrm{CH}_{3} \mathrm{COOK}$ | Use of sodium <br> hydroxide <br> (because it's <br> in food) | $\mathbf{1}$ |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 21 (a)(i) | $\begin{align*} & \Delta S_{\text {system }}^{9}=109.2+(6 x 69.9)-343 \text { (1) } \\ & =(+) 185.6\left(\left(\mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right) /(+) 186(\mathrm{~J} \mathrm{~mol}\right. \tag{1} \end{align*}$ <br> OR $(+) 0.186\left(\mathrm{~kJ} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ <br> IGNORE units even if incorrect <br> correct answer with no working scores 2 <br> Value using 1 for $\mathrm{H}_{2} \mathrm{O}=-163.9$ scores 1 <br> Use of value for $\mathrm{H}_{2} \mathrm{O}(\mathrm{g})$ (188.7) gives $898.4\left(\mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ (1) <br> correct value with incorrect sign scores 1 | 185 | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 1}$ (a)(ii) | Yes as (solid and) liquid forms (from solid) / <br> number of moles increases | Disorder <br> increases, <br> with no ref to <br> ORuid or <br> number of <br> moles | $\mathbf{1}$ |
|  | If $\Delta \mathbf{S}_{\text {system }}$ in (i) is negative the sign is not as <br> expected as liquid forms from solid / number <br> of moles increases |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 21 (a)(iii) | First mark $\begin{equation*} \Delta S_{\text {surroundings }}^{s}=\frac{-88.1 \times(1000)}{298} \tag{1} \end{equation*}$ <br> Second mark $\begin{aligned} & =-295.6375 \\ & =-295.6 \mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1} \end{aligned}$ <br> correct units must be shown but order not important <br> OR <br> $-0.2956 \mathrm{~kJ} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}(\mathbf{1})$ correct units must be shown but order not important <br> correct answer with or without working and correct units scores (2) <br> ignore sf except 1 <br> correct value with positive sign scores 1 |  | 2 |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 21 (a)(iv) | $\begin{aligned} & (185.6-295.6) \\ & =-110(\mathrm{~J} \mathrm{~mol} \\ & \left.\mathrm{m}^{-1}\right) \end{aligned}$ <br> OR $-0.110\left(\mathrm{~kJ} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ <br> could use 186 or 296 etc <br> TE from (a)(i) and (iii) <br> $(+) 602.8\left(\mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ if value for $\mathbf{6 H} \mathbf{~} \mathbf{O}(\mathrm{g})$ was used in (a) (i) <br> $-459.5\left(\mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ if value for one $\mathrm{H}_{2} \mathrm{O}$ was used in (a) (i) | Answers where values in J are added to kJ | 1 |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 1 ( a ) ( v )}$ | Decomposition (at 298 K) will not occur as <br> $\Delta S_{\text {total }}$ is negative / Reactions are only <br> spontaneous if total entropy change is positive <br> ldecomposition not thermodynamically <br> feasible / (hydrated cobalt chloride) is <br> thermodynamically stable | $\mathbf{1}$ |  |
| TE if answer to (a)(iv) is positive showing <br> decomposition (at 298 K) may occur <br> OR | Positive total entropy change doesn't indicate <br> rate of reaction |  |  |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 1}$ (b)(i) | First mark <br> Thermometer (1) <br> Second mark (dependent on first) <br> depends on choosing thermometer <br> as temperature change is small / <br> (\%) error in balance smaller than for <br> temperature reading <br> (\%) error in pipette smaller than for <br> temperature reading <br> (can be shown by calculation) / <br> as scale with greater degree of precision <br> needed / scale with more graduations needed <br> (1) <br> IGNORE any references to 'accurate <br> thermometer' | $\mathbf{2}$ |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| $\mathbf{2 1 ( b ) ( i i ) ~}$ | Use more cobalt chloride / less water (1) <br> To increase temperature rise (1) <br> Mark independently | J ust 'use more <br> reactants' <br> Use more <br> cobalt <br> chloride and <br> more water | $\mathbf{2}$ |
| repeat expt |  |  |  |
| add a lid or |  |  |  |
| extra |  |  |  |
| insulation to |  |  |  |
| beaker |  |  |  |$\quad$| use distilled |
| :--- |
| water |$\quad$.


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| $21 \text { (c)(i) }$ <br> QWC | Radius (of cation) increases (down group) OR any two values of radius: <br> $\mathrm{Mg}^{2+}=0.072, \mathrm{Ca}^{2+}=0.100 / \mathrm{Sr}^{2+}=0.113(\mathrm{~nm})$ data may be shown beside the table (1) <br> Radius $\mathrm{Co}^{2+}=0.065 \mathrm{~nm}$ <br> OR $\mathrm{Co}^{2+}$ radius smaller than other ions (1) <br> Data on EITHER $\mathrm{Co}^{2+}$ OR data showing increase in radius down Group II required for BOTH of first two marks <br> Force of attraction between ions decreases (as radius of ions increases) / charge density of ions decreases / negative ion can come closer to nucleus of positive ion (1) <br> ALLOW "weaker ionic bonds" <br> Predict lattice energy -2550 to $-2900\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> (1) <br> IGNORE sign | Atomic radii unless ionic radii also given <br> Radius of cobalt chloride <br> Polarising power decreases | 4 |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| 21 (c)(ii) | First mark <br> Qeference to enthalpy of hydration (may be in <br> equation $\Delta \mathrm{H}_{\text {solution }}=-$ LE $+\Delta H_{\text {hydration }}$ (1) <br> Second mark <br> Solubility depends on relative size of lattice <br> energy and enthalpy of hydration (1) <br> Third mark <br> EITHER <br> Solubility more likely if $\Delta H_{\text {solution }}$ is negative <br> OR | $\mathbf{3}$ |  |
| (If $\Delta H_{\text {solution }}$ is positive,) may / will dissolve if |  |  |  |
| $\Delta \mathrm{S}_{\text {total }}$ is positive |  |  |  |
| ACCEPT solvation instead of hydration |  |  |  |


| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 21 (d) QWC | First mark <br> Third ionization energy high(er) for $\mathrm{Mg} / \mathrm{Mg}=$ $7733 \mathrm{~kJ} \mathrm{~mol}^{-1}$, (third ionization energy for $\mathrm{Co}=$ $3232 \mathrm{~kJ} \mathrm{~mol}^{-1}$ ) (1) <br> Second mark <br> (Third ionization energy for Mg is high) because the electron is being removed from an inner shell / full shell / $2 p$ level / $2 p$ orbital (1) <br> OR <br> Not compensated by higher lattice energy for $\mathrm{Mg}^{3+}$ (and so $\Delta \mathrm{H}_{\text {formation }}$ of $\mathrm{MgCl}_{3}$ would be highly endothermic) (1) |  | 2 |

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