# Mark Scheme (Results) June 2010 

## GCE

## GCE Chemistry (6CH04/ 01)

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Section A (multiple choice)

| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 1 (a) | D | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 1 (b) | D | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 1 (c) | A | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 2 | B | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 3 | C | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 4 | D | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 5 | B | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 6 | A | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 7 (a) | C | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 7 (b) | B | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 7 (c) | D | 1 |


| Question <br> Number | Correct Answer | Mark |
| :--- | :--- | :--- |
| 8 | B | 1 |


| Question Number | Correct Answer | Mark |
| :---: | :---: | :---: |
| 9 | D | 1 |
| Question 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 |
| :--- | :--- | :--- | :--- |
| 17 (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 $6 \times 10^{-5}$ ) | 1 |  |


| Question <br> Number | Accept able Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 17 (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] const ant )/ rate and <br> concentration increase in proportion (1) <br> ALLOW'time halves' instead of 'rate doubles' | 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> [C449Br] constant ) (1) | ALLOW' doubling in concentration of OH' <br> instead of 'increase in concentration' | ALLOWtime increases by the same factor as <br> increase in hydroxide concentration (5/ 3) | May refer to experiment number rather than <br> concentrations |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 17 (a)(iii) | Rate $=\mathrm{k}\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]$ <br> OR Rate $=\mathrm{k}\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]^{1}[\mathrm{OH}]^{0}$ <br> ALLOWk 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 | 1 |  |


| Question Number | Acceptable Answers | Rej ect | 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> ALLOWk $=1.68 \times 10^{-3}$ <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 $=k\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]^{2}$ gives $\mathrm{k}=0.10036 \mathrm{dm}^{3} \mathrm{~mol}^{-1} \mathrm{~s}^{-1}$ <br> Rate $=\mathrm{k}\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right][\mathrm{OH}]$ 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 $=k\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right][\mathrm{OH}]^{2}$ gives $\mathrm{k}=1184.6$ $\mathrm{dm}^{6} \mathrm{~mol}^{-2} \mathrm{~s}^{-1}$ <br> Rate $=k\left[\mathrm{C}_{4} \mathrm{H}_{9} \mathrm{Br}\right]^{2}[\mathrm{OH}]$ gives $\mathrm{k}=83.62$ $\mathrm{dm}^{6} \mathrm{~mol}^{-2} \mathrm{~s}^{-1}$ |  | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 17 (b) | $[\mathrm{OH}]$ 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 OH- is in a step other than <br> rate determining step) | 1 |  |


| Question Number | Acceptable Answers | Mark |
| :---: | :---: | :---: |
| 17 (c) | First mark <br> Choice of bromoalkane must be consistent with rate equation in (a)(iii). <br> If $[\mathrm{OH}]$ is not in rate equation, secondary/tertiary bromoalkane. <br> If $[\mathrm{OH}]$ is in rate equation, primary/ secondary bromoalkane. <br> Second and third marks <br> Ether SN1 or SN2 mechanism can score 2 marks regardless of choice of bromoalkane. <br> Lone pairs not required <br> Curly arrow from $\mathrm{C}-\mathrm{Br}$ bond to Br (making Br ) (1) <br> Curly arrow from anywhere on $\mathrm{OH}^{-} / \mathrm{HO}^{-}$to $\mathrm{C}^{+}$in correct intermediate (making alcohol) (1) <br> OR <br> Both curly arrows from $\mathrm{OH}^{-}$and from $\mathrm{C}-\mathrm{Br}$ bond to Br (may both be shown at start) (1) <br> Transition state including minus charge (and product) (1) <br> Do not penalise if $\mathrm{C}_{2} \mathrm{H}_{5}$ shown instead of $\mathrm{C}_{3} \mathrm{H}_{7}$. <br> Bonds in transition state can be dotted. <br> Do not penalise the missing H atoms in alkyl groups in mechanism. | 3 |


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


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 18 (a)(i) | (Acid) hydrolysis | substitution | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 18 (a)(ii) | $\mathrm{K}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{Na}_{2} \mathrm{Cr}_{2} \mathrm{O}_{7} / \mathrm{Cr}_{2} \mathrm{O}_{7}^{2-}$ <br> Potassium dichromate((V)) / sodium <br> dichromate((VI)) / dichromate((VI)) ions <br> ALLOW manganate((VII)) ions, etc | Just <br> "dichromate" <br> chromates | 1 |
| Correct <br> formula with <br> wrong name <br> and vice versa <br> Incorrect | loxidation <br> oxumber |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 18 (a)(iii) | Lithium tetrahydridoaluminate/ lithium <br> aluminium hydride/ LiAlH (in dry ether) |  |  |\(~\left(\begin{array}{ll}Just[H] \& 1 <br>

\hline\end{array}\right.\)
$\left.\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Reject } & \text { Mark } \\ \hline 18 \text { (a)(iv) } & \begin{array}{l}\text { Methyl butanoate (1) } \\ \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{COOH}^{2}+\mathrm{CH}_{3} \mathrm{OH} \rightarrow+ \\ \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{COOCH}_{3}+\mathrm{H}_{2} \mathrm{O} \text { (1) }\end{array} & \begin{array}{l}\text { Methyl } \\ \text { butoate }\end{array} & 2\end{array}\right\}$

| Question Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 18 (a) (v) |  <br> Don't penalise undisplayed methyl groups as here. COCl must be displayed as above. | $\begin{aligned} & \mathrm{C}_{3} \mathrm{H}_{7} \text { for } \\ & \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \end{aligned}$ | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 18 (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 | 1 |  |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 18 (b)(ii) | Solubility (in liquid / stationary phase) <br> OR <br> Interaction with liquid / stationary phase OR <br> Interaction between mobile and stationary phase <br> OR <br> Attraction for liquid / stationary phase <br> OR <br> Strength of (named) intermolecular forces 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> Just a named intermolecular force <br> Just 'retention time' <br> Density | 1 |


| Question Number | Acceptable Answers | Rej ect | 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> PENALISElack of displayed $\mathrm{C}=\mathrm{O}$ 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 | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 18 (c)(ii) | Hydrolysis <br> OR <br> Splits/breaks ester link <br> OR <br> polymer breaks down to monomers <br> OR <br> equation showing hydrolysis | Just 'breaks <br> polymer down' | 1 |


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


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 19 (a)(ii) | $\mathrm{P} \mathrm{CH} 3 \mathrm{OH}=4.9(\operatorname{atm})(1)$ <br> $\mathrm{P} \mathrm{CO}=4.9(\operatorname{atm})(1)$ <br> 1 mark for recognition that pressures are equal <br> IGNORE units | 2 |  |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| 19 (a)(iii) | $K_{p}=\left((22.2) /(4.9)^{2}\right)$ <br> $=0.925(1)$ <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 | 2 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 19 (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 | 2 |  |

$\left.\begin{array}{|l|l|l|l|}\hline \begin{array}{l}\text { Question } \\ \text { Number }\end{array} & \text { Acceptable Answers } & \text { Rej ect } & \text { Mark } \\ \hline 19 \text { (b)(ii) } & \left(\frac{46.8 \times 32}{53.2}\right)=28.2 / 28.1504 \text { (atm) } & 28.1 & 1 \\ & \text { IGNORE sf except } 1 \\ \text { Value }=28.16 \text { if mol fraction rounded } \\ \text { ALLOW TE from (b)(i) } & \frac{46.8 \times 32}{50}= & 29.95 \text { (atm) }\end{array}\right]$

| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 19 (b)(iii) | exothermic as yield / pp of ethanoic acid / <br> conversion of reactants/ K is higher at lower <br> temperature / as equilibrium moves (right) at <br> lower temperature | 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 Number | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: |
| 19 (c)(i) | No effect <br> and <br> other concentrations change to keep $\mathrm{K}_{\mathrm{p}}$ constant / $K_{p}$ is only affected by temperature/ as equilibrium moves (right) to keep $\mathrm{K}_{\mathrm{p}}$ constant / change in pressure does not change $\mathrm{K}_{\mathrm{p}}$ | As $K_{p}$ is a constant | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 19 (c)(ii) | Yield increased to restore fraction / quotient / <br> partial pressure ratio back to K |  | 1 |
|  | ALLOW (equilibrium moves) to use up the <br> methanol / answers based on entropy or Le <br> Chatelier <br> Correct prediction in (c)(i) and (c)(ii) with <br> inadequate explanations scores 1 mark in <br> (c)(ii) | Just <br> equilibrium <br> moves to the <br> right' |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 19 (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 |  |
| Enables use of an alternative process with <br> higher atom economy (1) <br> fewer raw materials needed / less natural <br> resources used (1) | 2 |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 20 (a)(i) | Correct answer with or without working scores <br> 2 marks <br> $\left[\mathrm{H}^{+}\right]=\left(1.00 \times 10^{-14} / 0.250\right)=4 \times 10^{-14}(1)$ <br> $\mathrm{pH}=(13.39794=) 13.4(1)$ <br> OR <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 | 2 |  |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 20 (a)(ii) | $\begin{equation*} \left(K_{\mathrm{a}}=\right) \frac{\left[\mathrm{CH}_{3} \mathrm{COO}-\right]\left[\mathrm{H}^{+}\right]}{\left[\mathrm{CH}_{3} \mathrm{COOH}\right]} \tag{1} \end{equation*}$ <br> ALLOW <br> $\mathrm{H}_{3} \mathrm{O}^{+}$instead of $\mathrm{H}^{+}$ <br> [ $\left.\mathrm{A}^{-}\right]\left[\mathrm{H}^{+}\right]$if key to symbols given [HA] <br> IGNOREstate symbols | $\frac{\left[\mathrm{H}^{+}\right]^{2}}{\left[\mathrm{CH}_{3} \mathrm{COOH}\right]}$ | 1 |


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


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 20 (a)(iv) | $\begin{aligned} & \mathrm{pH}=4.8 / 4.77(1) \\ & \mathrm{pH}=\mathrm{p} K_{\mathrm{a}} /\left[\mathrm{H}^{+}\right]=K_{\mathrm{a}}(\text { when acid is half } \\ & \text { neutralized) (1) } \end{aligned}$ | $\mathrm{H}^{+}=K_{a}$ | 2 |


| Question <br> Number | Acceptable Answers | Rej ect | Mark |
| :--- | :--- | :--- | :--- |
| 20 (a)(v) | Sgmoid 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. Soping 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 | 3 |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 20 (a)(vi) | 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) | 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 |  |  |  |
| 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 |  |
| 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 | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 20 (b) | Sodium ethanoate/ $\mathrm{CH}_{3} \mathrm{COONa}$ <br> Potassium ethanoate $/ \mathrm{CH}_{3} \mathrm{COOK}$ <br> ALLOW <br> other cations as alternatives to sodium | Use of sodium <br> hydroxide <br> (because it's <br> in food) | 1 |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 21 (a)(i) | $\begin{align*} & \Delta S_{\text {system }}=109.2+(6 \times 69.9)-343(1) \\ & =(+) 185.6\left(\mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right) /(+) 186\left(\mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right) \tag{1} \end{align*}$ <br> OR $(+) 0.186\left(\mathrm{~kJ} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ <br> IGNORE units even if incorrect correct answer with no working scores 2 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 |
| :--- | :--- | :--- | :--- |
| 21 (a)(ii) | Yes as (solid and) liquid forms (from solid) / <br> number of moles increases | Disorder <br> increases, <br> with no ref to <br> liquid or <br> number of <br> moles | 1 |
|  | If $\Delta S_{\text {ssstem in (i) is negative the sign is not as }}^{\text {expected as liquid forms from solid / number }}$ <br> of moles increases |  |  |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| 21 (a)(iii) | First mark $\begin{equation*} \Delta \mathcal{S}_{\text {surroundings }}=\frac{-88.1 \times(1000)}{298} \tag{1} \end{equation*}$ <br> Second mark $=-295.6375$ $=-295.6 \mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}$ <br> correct units must be shown but order not important <br> OR <br> $-0.2956 \mathrm{~kJ} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}(1)$ <br> 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 <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 21 (a)(iv) | $(185.6-295.6)$ <br> $=-110\left(\mathrm{~J} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ <br> OR <br> $-0.110\left(\mathrm{~kJ} \mathrm{~mol}^{-1} \mathrm{~K}^{-1}\right)$ <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 $6 \mathrm{H}_{2} \mathrm{O}(\mathrm{g})$ was <br> 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 <br> used in (a) (i) | Answers where <br> values in J are <br> added to kJ | 1 |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 21 (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> l decomposition not thermodynamically <br> feasible / (hydrated cobalt chloride) is <br> thermodynamically stable | 1 |  |
| TE if answer to (a)(iv) is positive showing <br> decomposition (at 298 K) may occur <br> OR <br> Oositive total entropy change doesn't indicate <br> rate of reaction |  |  |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 21 (b)(i) | First mark <br> Thermometer (1) <br> Second mark (dependent on first) <br> depends on choosing thermometer <br> as temperat ure change is small / <br> (9/ 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' | 2 |  |


| Question <br> Number | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- |
| 21 (b)(ii) | Use more cobalt chloride / less water (1) <br> To increase temperature rise (1) <br> Mark independently | Just 'use more <br> reactants' <br> Use more <br> cobalt <br> chloride and <br> more water <br> repeat expt <br> add a lid or <br> extra <br> insulation to <br> beaker | 2 |
| use distilled <br> water |  |  |  |


| Question Number | Acceptable Answers | Rej ect | Mark |
| :---: | :---: | :---: | :---: |
| $21 \text { (c)(i) }$ <br> QWC | Radius (of cation) increases (down group) <br> OR any two values of radius: <br> $\mathrm{Mg}^{2+}=0.072, \mathrm{Ca}^{2+}=0.100 / \mathrm{Sr}^{2+}=0.113(\mathrm{~nm})$ <br> data may be shown beside the table (1) <br> Radius $\mathrm{Co}^{2+}=0.065 \mathrm{~nm}$ <br> $\mathrm{OR} \mathrm{Co}^{2+}$ 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 (kJ mol ${ }^{-1}$ ) <br> (1) <br> IGNORE sign | Atomic radii unless ionic radii also given <br> Radius of cobalt chloride <br> Polarising power decreases | 4 |


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| :--- | :--- | :--- | :--- |
| 21 (c)(ii) | First mark <br> Reference to enthalpy of hydration (may be in <br> equation $\Delta H_{\text {solution }}=-L E+\Delta H_{\text {hydration }}$ (1) <br> QWC <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 | 3 |  |
| (If $\Delta H_{\text {solution }}$ is positive,) may / will dissolve if |  |  |  |
| $\Delta$ Sotal $_{\text {is positive }}$ |  |  |  |
| $A C C E P T$ solvation instead of hydration |  |  |  |


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

