Version 1.2



# General Certificate of Education June 2010 

## Chemistry

## CHEM5

Energetics, Redox and Inorganic Chemistry

## Mark Scheme

Mark schemes are prepared by the Principal Examiner and considered, together with the relevant questions, by a panel of subject teachers. This mark scheme includes any amendments made at the standardisation meeting attended by all examiners and is the scheme which was used by them in this examination. The standardisation meeting ensures that the mark scheme covers the candidates' responses to questions and that every examiner understands and applies it in the same correct way. As preparation for the standardisation meeting each examiner analyses a number of candidates' scripts: alternative answers not already covered by the mark scheme are discussed at the meeting and legislated for. If, after this meeting, examiners encounter unusual answers which have not been discussed at the meeting they are required to refer these to the Principal Examiner.

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| Q | Part | Sub Part | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) |  | $\mathrm{CaF}_{2}(\mathrm{~s}) \rightarrow \mathrm{Ca}^{2+}(\mathrm{g})+2 \mathrm{~F}^{-}(\mathrm{g})$ | 1 |  |
| 1 | (b) | (i) | Enthalpy change for formation of 1 mol of substance <br> From its elements <br> Reactants and products/all substances in their standard states | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | Allow heat energy change, NOT energy <br> Or normal states at $298 \mathrm{~K}, 1$ bar (100 kPa ) |
| 1 | (b) | (ii) | $\mathrm{Ca}(\mathrm{s})+\mathrm{F}_{2}(\mathrm{~g}) \rightarrow \mathrm{CaF}_{2}(\mathrm{~s})$ | 1 |  |
| 1 | (b) | (iii) | $\begin{aligned} & \Delta H_{f}\left(\mathrm{CaF}_{2}\right)=\Delta H_{\mathrm{a}}(\mathrm{Ca})+1 \mathrm{st} \mathrm{IE}(\mathrm{Ca})+2^{\text {nd }} \mathrm{IE}(\mathrm{Ca})+\mathrm{BE}\left(\mathrm{~F}_{2}\right)+2 \text { xEA }(\mathrm{F})- \\ & \Delta H_{\mathrm{L}}\left(\mathrm{CaF}_{2}\right) \\ & =193+590+1150+158+(2 \mathrm{x}-348)-2602 \\ & =-1207 \mathrm{~kJ} \mathrm{~mol}^{-1} \end{aligned}$ | 1 1 1 | Or labelled diagram <br> Correct answer scores 3 -842 scores 2 (transfer error) -859 scores 1 only (using one E.A.) Units not required, wrong units lose 1 mark |
| 1 | (c) |  | Electrostatic attraction stronger/ionic bonding stronger/attraction between ions stronger/more energy to separate ions <br> Because fluoride (ion) smaller than chloride | 1 1 | Molecular attraction /atoms/intermolecular forces CE=0 <br> Do not allow $F$ or fluorine |
| 1 | (d) | (i) | $\begin{aligned} & \Delta H=\Delta H_{\mathrm{L}}+\Sigma \Delta H_{\mathrm{hyd}}=2237-1650+(2 \mathrm{x}-364) \\ & =-141 \mathrm{~kJ} \mathrm{~mol}^{-1} \end{aligned}$ | 1 1 | Can be on cycle/diagram <br> Correct answer scores 2 <br> Units not required, wrong units lose 1 mark |


| 1 | (d) | (ii) | Decreases <br> Reaction exothermic/ $\Delta \mathrm{H}$-ve <br> (Equilibrium )shifts to left/backwards (as temperature rises)/ equilibrium opposes the change | 1 1 1 | If ans to (d)(i) positive allow increases If (d)(i) +ve allow endothermic/ $\Delta H+$ ve <br> If (d) (i) +ve allow shifts to right/forwards / equilibrium opposes the change <br> If no answer to (d) (i) assume -ve $\Delta H$ used <br> If effect deduced incorrectly from any $\Delta H C E=0$ for these 3 marks |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (e) |  | u.v. absorbed: electrons/they move to higher energy (levels)/ electrons excited visible light given out: electrons/they fall back down/move to lower energy (levels) |  | Must refer to absorbing u.v. NOT visible light or this must be implied. |


| Q Part Sub <br> Part Marking Guidance Mark Comments <br> 2 (a)  Macromolecular 1 Or giant molecule <br> Or giant covalent (also gains M 2$)$ <br> Do <br> not allow giant atomic <br> Ionic/metallic $\mathrm{CE}=0$ for all 3 marks      <br> Do NOT allow if between molecules      |
| :--- |


| 2 | (f) | (ii) | $\mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{NaOH}+3 \mathrm{H}_{2} \mathrm{O} \rightarrow 2 \mathrm{NaAl}(\mathrm{OH})_{4}$ | 1 | Other equations with $\mathrm{Al}_{2} \mathrm{O}_{3}$ are possible e.g. $\begin{aligned} & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{OH}^{-}+3 \mathrm{H}_{2} \mathrm{O} \rightarrow 2\left[\mathrm{Al}(\mathrm{OH})_{4}\right]^{-} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{OH}^{-}+7 \mathrm{H}_{2} \mathrm{O} \rightarrow \\ & 2\left[\mathrm{Al}^{\left.\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}(\mathrm{OH})_{4}\right]^{-}}\right. \end{aligned}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | (g) |  | $\mathrm{SiO}_{2}$ acidic/Lewis acid/electron pair acceptor $\mathrm{SiO}_{2}+2 \mathrm{NaOH} \rightarrow \mathrm{Na}_{2} \mathrm{SiO}_{3}+\mathrm{H}_{2} \mathrm{O}$ | 1 1 | Allow $\mathrm{SiO}_{2}$ not amphoteric Do NOT allow BL acid <br> Other equations with $\mathrm{SiO}_{2}$ are possible e.g. $\begin{aligned} & \mathrm{SiO}_{2}+2 \mathrm{OH}^{-} \rightarrow \mathrm{SiO}_{3}{ }^{2-}+\mathrm{H}_{2} \mathrm{O} \\ & \mathrm{SiO}_{2}+2 \mathrm{OH}^{-}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{Si}(\mathrm{OH})_{6}{ }^{2-} \end{aligned}$ |


| Q | Part | $\begin{array}{\|l} \hline \text { Sub } \\ \text { Part } \\ \hline \end{array}$ | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) |  | Same phase/state | 1 |  |
| 3 | (b) |  | Because only exist in one oxidation state | 1 | Allow do not have variable oxidation states |
| 3 | (c) |  | $21^{-}+\mathrm{S}_{2} \mathrm{O}_{8}{ }^{2-} \rightarrow \mathrm{I}_{2}+2 \mathrm{SO}_{4}{ }^{2-}$ | 1 | Ignore state symbols Allow multiples |
| 3 | (d) |  | Both (ions)have a negative charge | 1 | Or both have the same charge Or (ions) repel each other Do not allow both molecules have the same charge (contradiction) |
| 3 | (e) |  | $\begin{aligned} & 2 \mathrm{Fe}^{2+}+\mathrm{S}_{2} \mathrm{O}_{8}^{2-} \rightarrow 2 \mathrm{Fe}^{3+}+2 \mathrm{SO}_{4}{ }^{2-} \\ & 2 \mathrm{Fe}^{3+}+2 \mathrm{I}^{-} \rightarrow 2 \mathrm{Fe}^{2+}+\mathrm{I}_{2} \end{aligned}$ <br> Positive and negative (ions)/oppositely charged (ions) | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | Equations can be in any order <br> Mark independently |
| 3 | (f) |  | Equations 1 and 2 can occur in any order | 1 | Allow idea of $\mathrm{Fe}^{3+}$ converted to $\mathrm{Fe}^{2+}$ then $\mathrm{Fe}^{2+}$ converted back to $\mathrm{Fe}^{3+}$ |


| Q | Part | Sub Part | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (a) |  | Partially filled/incomplete d sub-shell/orbital/shell | 1 | Ignore reference to forbitals <br> Do not allow d block <br> Do not allow half-filled d orbitals |
| 4 | (b) |  | Has ligand(s) <br> linked by co-ordinate bonds | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | Allow molecules/ions with lone pairs <br> Allow dative/donation of lone pair |
| 4 | (c) |  | (Blue) light is absorbed (from incident white light) <br> Due to electrons moving to higher levels / electrons excited <br> Red light (that) remains (is transmitted) / light that remains (transmitted light) is the colour observed | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | Allow d $\rightarrow$ d transitions <br> Allow red light reflected |
| 4 | (d) | (i) | Circle round any $\mathrm{O}^{-}$ <br> Circle round either N | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | List principle |
| 4 | (d) | (ii) | $\mathrm{EDTA}^{4-}+\left[\mathrm{Co}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+} \rightarrow[\mathrm{CoEDTA}]^{2-}+6 \mathrm{H}_{2} \mathrm{O}$ | 1 | Allow missing square brackets Ignore state symbols |
| 4 | (d) | (iii) | Increase in entropy/ $\Delta S$ positive <br> Because 2 mol (of particles/molecules/species/entities) form 7 mol | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | Or increase in disorder <br> Allow 'increase in number' as stated in words or as shown by any numbers deduced correctly from an incorrect equation <br> Do not allow increase in ions/atoms |

\begin{tabular}{|c|c|c|c|c|c|}
\hline 4 \& (e) \& (i) \& \begin{tabular}{l}
Co-ordinate/dative/dative covalent bond \\
Covalent bond
\end{tabular} \& 1

1 \& | Allow pair of electrons donated by nitrogen/ligand |
| :--- |
| Do not allow pair of electrons donated from Iron/Fe |
| Shared electron pair | <br>

\hline 4 \& (e) \& (ii) \& Transport of oxygen/ $\mathrm{O}_{2}$ \& 1 \& | Allow any statement that implies oxygen carried (around the body) |
| :--- |
| Do not allow transport of carbon dioxide $\left(\mathrm{CO}_{2}\right)$. This also contradicts the mark (list principle) | <br>


\hline 4 \& (e) \& (iii) \& | Because it bonds to the iron/haemoglobin |
| :--- |
| Displaces oxygen | \& 1

1 \& | Allow blocks site /CO has greater affinity for haemoglobin /carboxyhaemoglobin more stable than oxyhaemoglobin |
| :--- |
| Or prevents transport of oxygen QoL | <br>

\hline
\end{tabular}

| Q | Part | $\begin{aligned} & \hline \text { Sub } \\ & \text { Part } \end{aligned}$ | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (a) |  | W is $\mathrm{CuCl}_{4}{ }^{2-}$ <br> Yellow-green/yellow/green $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}+4 \mathrm{Cl}^{-} \rightarrow \mathrm{CuCl}_{4}{ }^{2-}+6 \mathrm{H}_{2} \mathrm{O}$ | $1$ <br> 1 $1$ | Not necessary to indicate solution Do not allow precipitate/solid $\text { Allow }+4 \mathrm{HCl} \rightarrow 4 \mathrm{H}^{+}$ |
| 5 | (b) |  | X is $\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}(\mathrm{OH})_{2}$ <br> Blue precipitate/solid $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}+2 \mathrm{NH}_{3} \rightarrow \mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}(\mathrm{OH})_{2}+2 \mathrm{NH}_{4}^{+}$ |  | Allow $\mathrm{Cu}(\mathrm{OH})_{2} /$ copper hydroxide Ignore shades <br> Allow any balanced equation/equations leading to this hydroxide or $\mathrm{Cu}(\mathrm{OH})_{2}$ But must use ammonia |


| 5 | (c) |  | Y is $\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}\right]^{2+}$ <br> Deep/dark/royal blue solution $\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{4}(\mathrm{OH})_{2}+4 \mathrm{NH}_{3} \rightarrow\left[\mathrm{Cu}\left(\mathrm{NH}_{3}\right)_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}\right]^{2+}+2 \mathrm{H}_{2} \mathrm{O}+2 \mathrm{OH}^{-}$ | $1$ <br> 1 <br> 1 | QoL <br> Accept equation for formation from $\mathrm{Cu}(\mathrm{OH})_{2}$ |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (d) |  | Z is $\mathrm{CuCO}_{3}$ <br> Green solid/precipitate $\left[\mathrm{Cu}\left(\mathrm{H}_{2} \mathrm{O}\right)_{6}\right]^{2+}+\mathrm{CO}_{3}^{2-} \rightarrow \mathrm{CuCO}_{3}+6 \mathrm{H}_{2} \mathrm{O}$ | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | Allow copper carbonate <br> Allow blue-green precipitate |
| 5 | (e) | (i) | $\mathrm{Cu}^{2+}(\mathrm{aq})+\mathrm{Fe}(\mathrm{~s}) \rightarrow \mathrm{Cu}(\mathrm{~s})+\mathrm{Fe}^{2+}(\mathrm{aq})$ <br> Blue <br> Green | 1 1 | Allow hydrated ions State symbols not essential but penalise if wrong <br> Do not allow description of solids <br> Allow yellow/(red-)brown/orange |


| 5 | (e) | (ii) | Any two correct points about copper extraction from two of these three <br> categories: <br> Any relevant mention of lower energy consumption <br> Any relevant mention of benefits of less mining (of copper ore) <br> Less release of $\mathrm{CO}_{2}$ (or CO ) into the atmosphere | Max 2 <br> Do not allow reference to electricity <br> alone or to temperature alone. |
| :--- | :--- | :--- | :--- | :--- | :--- |
| Allow avoids depletion of (copper ore) |  |  |  |  |
| resources |  |  |  |  |


| Q | Part | Sub <br> Part | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 6 | (a) |  | $\begin{aligned} & \Delta H=\Sigma \Delta H_{f}(\text { products })-\Sigma \Delta H_{f}(\text { reactants }) \\ & =-201-242-(-394) \\ & =-49 \mathrm{~kJ} \mathrm{~mol}^{-1} \end{aligned}$ | $1$ $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | $+49 \mathrm{~kJ} \mathrm{~mol}^{-1}=1$ mark units not required, wrong units lose 1 mark |
| 6 | (b) |  | $\begin{aligned} & \Delta S=\Sigma S(\text { products })-\Sigma S(\text { reactants }) \\ & =238+189-(214+3 \times 131) \\ & =-180 \mathrm{~J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1} \end{aligned}$ | $\begin{aligned} & 1 \\ & 1 \\ & 1 \end{aligned}$ | $+180=1 \text { mark }$ <br> units not required, wrong units lose 1 mark |
| 6 | (c) |  | $\Delta G=\Delta H-T \Delta S$ <br> ( $\Delta S$ is negative so) at high temp $-T \Delta S$ (is positive and) greater than $\Delta H /$ large <br> So $\Delta G>0$ <br> (Limiting condition $\Delta G=0$ so) $T=\Delta H / \Delta S$ $=272 \mathrm{~K}$ <br> Reaction is too slow at this temperature/to speed up the reaction | 1 <br> 1 <br> 1 <br> 1 <br> 1 <br> 1 | If use $G$ not $\Delta G$ penalise M1 but not M2 and M3 <br> Do not award M2 or M3 if positive $\Delta S$ value used <br> Independent mark unless positive $\Delta S$ value used <br> Allow 297-298 if used given values. Do not award M5 if T -ve or if M4 should give T -ve |



| Q | Part | Sub <br> Part | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 7 | (a) |  | Hydrogen $/ \mathrm{H}_{2}$ gas/bubbles <br> $1.0 \mathrm{~mol} \mathrm{dm}^{-3} \mathrm{HCl} / \mathrm{H}^{+}$ <br> At 298 K and 100 kPa <br> Pt (electrode) | $\begin{aligned} & 1 \\ & 1 \\ & 1 \\ & 1 \end{aligned}$ | Allow 1 bar instead of 100 kPa Do not allow 1 atm |
| 7 | (b) |  | $\begin{aligned} & \mathrm{Li}^{+}+\mathrm{MnO}_{2}+\mathrm{e}^{-} \rightarrow \mathrm{LiMnO}_{2} \\ & -0.13(\mathrm{~V}) \end{aligned}$ | $\begin{aligned} & 1 \\ & 1 \\ & \hline \end{aligned}$ | Ignore state symbols |
| 7 | (c) |  | $\begin{aligned} & \mathrm{Fe}^{3+} \text { ions reduced to } \mathrm{Fe}^{2+} \\ & \text { Because } E\left(\mathrm{Fe}^{3+}\left(/ \mathrm{Fe}^{2+}\right)\right)>E\left(\mathrm{H}^{+} / \mathrm{H}_{2}\right) / E(\text { hydrogen }) \end{aligned}$ | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | Can score from equation/scheme <br> Allow emf $/ E_{\text {cell }}+\mathrm{ve} / 0.77 \mathrm{~V}$ <br> Allow $\mathrm{Fe}^{3+}$ better oxidising agent than $\mathrm{H}^{+}$ <br> Allow $\mathrm{H}_{2}$ better reducing agent than $\mathrm{Fe}^{2+}$ <br> Only award this explanation mark if previous mark given |



