\begin{tabular}{|c|c|c|c|c|c|}
\hline \multicolumn{3}{|r|}{Question} \& Answer \& Mark \& Guidance \\
\hline 1 \& (a) \& \& \begin{tabular}{l}
FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer \(=8.3 \times 10^{4}\) OR 83333 award 2 marks THEN IF units are \(\mathbf{d m}^{6} \mathrm{~mol}^{-2} \mathbf{s}^{-1}\), award 1 further mark
\[
\begin{aligned}
\& k=\frac{r a t e}{\left[\mathrm{H}_{2}(\mathrm{~g})\right][\mathrm{NO}(\mathrm{~g})]^{2}} \text { OR } \frac{3.6 \times 10^{-2}}{\left(1.2 \times 10^{-2}\right) \times\left(6.0 \times 10^{-3}\right)^{2}} \\
\& \checkmark \\
\& =8.3 \times 10^{4} \mathrm{OR} 83000 \text { OR } 83333 \mathrm{r}
\end{aligned}
\] \\
units: \(\mathrm{dm}^{6} \mathrm{~mol}^{-2} \mathrm{~s}^{-1} \checkmark\)
\end{tabular} \& 2

1 \& | ALLOW 1 mark for $8.3 \times 10^{x}$ with no working (power of 10 is error) |
| :--- |
| ALLOW 2 SF up to calculator value of $8.33333333 \times 10^{4}$ correctly rounded |
| ALLOW ECF for calculated answer from incorrectly rearranged $k$ expression but not for units (Marked independently see below) |
| ALLOW dm ${ }^{6}, \mathrm{~mol}^{-2}$ and $\mathrm{s}^{-1}$ in any order, eg $\mathrm{mol}^{-2} \mathrm{dm}^{6} \mathrm{~s}^{-1}$ DO NOT ALLOW other units (Rate equation supplied on paper - not derived from data ) | \\

\hline \& (b) \& (i) \& effect on rate $\times 2 \checkmark$ \& 1 \& ALLOW 'doubles' OR rate $=7.2 \times 10^{-2}\left(\mathrm{~mol} \mathrm{dm}^{-3} \mathrm{~s}^{-1}\right)$ \\

\hline \& \& (ii) \& effect on rate $\times 1 / 4 \mathrm{OR} \times 0.25 \checkmark$ \& 1 \& | ALLOW 'a quarter' OR decrease by $1 / 4$ OR decrease by 0.25 OR rate decreases by 4 OR decrease by $75 \%$ OR rate $=0.9 \times 10^{-2}\left(\mathrm{~mol} \mathrm{dm}^{-3} \mathrm{~s}^{-1}\right)$ |
| :--- |
| DO NOT ALLOW just $0.5^{2}$ of rate OR rate decreases by $2^{2}$ | \\


\hline \& \& (iii) \& effect on rate $\times 64 \checkmark$ \& 1 \& | ALLOW rate $=2.3(04)\left(\mathrm{mol} \mathrm{dm}^{-3} \mathrm{~s}^{-1}\right)$ |
| :--- |
| DO NOT ALLOW just 'increases by 4 and then by $16 / 4^{2}$ OR increases by $4^{3}$ | \\

\hline
\end{tabular}

| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (c) | (i) | (initial) rate increases <br> AND <br> more frequent collisions OR more collisions per second/time | 1 | BOTH points required for mark <br> ALLOW rate increases AND concentration increases <br> For concentration increases, ALLOW particles closer together OR less space between particles <br> DO NOT ALLOW just more collisions OR collisions more likely |
|  |  | (ii) | rate constant does not change $\checkmark$ | 1 |  |
|  | (d) |  | $\text { step 1: } \mathrm{H}_{2}(\mathrm{~g})+2 \mathrm{NO}(\mathrm{~g}) \longrightarrow \mathrm{N}_{2} \mathrm{O}(\mathrm{~g})+\mathrm{H}_{2} \mathrm{O}(\mathrm{~g})$ <br> LHS of step one <br> step 2: $\mathrm{H}_{2}(\mathrm{~g})+\mathrm{N}_{2} \mathrm{O}(\mathrm{g}) \longrightarrow \mathrm{N}_{2}(\mathrm{~g})+\mathrm{H}_{2} \mathrm{O}(\mathrm{g})$ rest of equations for step 1 AND step $2 \checkmark$ | 2 | State symbols NOT required <br> For 'rest of equations', <br> This mark can only be awarded if 1 st mark can be awarded <br> ALLOW other combinations of two steps that together give the overall equation (shown above part in scoris window), eg step 1: $\longrightarrow \mathrm{N}_{2}(\mathrm{~g})+1 / 2 \mathrm{O}_{2}(\mathrm{~g})+\mathrm{H}_{2} \mathrm{O}(\mathrm{g})$ step 2: $\mathrm{H}_{2}(\mathrm{~g})+1 / 2 \mathrm{O}_{2}(\mathrm{~g}) \longrightarrow \mathrm{H}_{2} \mathrm{O}(\mathrm{g})$ <br> step 1: $\longrightarrow \mathrm{H}_{2} \mathrm{O}_{2}(\mathrm{~g})+\mathrm{N}_{2}(\mathrm{~g})$ <br> step 2: $\mathrm{H}_{2}(\mathrm{~g})+\mathrm{H}_{2} \mathrm{O}_{2}(\mathrm{~g}) \longrightarrow 2 \mathrm{H}_{2} \mathrm{O}(\mathrm{I})$ <br> There may be others with species, such as $\mathrm{H}_{2} \mathrm{~N}_{2} \mathrm{O}_{2}$ and HNO . Provided the two steps add up to give the overall equation AND charges balance, the 2nd mark can be awarded |
|  |  |  | Total | 10 |  |


| Question |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 2 | (a) | $\begin{aligned} & \text { Fe: } \quad\left(1 s^{2} 2 s^{2} 2 p^{6}\right) 3 s^{2} 3 p^{6} 3 d^{6} 4 s^{2} \checkmark \\ & \text { Fe}^{2+}:\left(1 s^{2} 2 s^{2} 2 p^{6}\right) 3 s^{2} 3 p^{6} 3 d^{6} \checkmark \end{aligned}$ | 2 | ALLOW $4 s$ before 3 d, i.e. $\left(1 s^{2} 2 s^{2} 2 p^{6}\right) 3 s^{2} 3 p^{6} 4 s^{2} 3 d^{6}$ <br> ALLOW $4 s^{0}$ <br> ALLOW subscripts <br> IGNORE $1 s^{2} 2 s^{2} 2 p^{6}$ is written out a second time |
|  | (b) | coloured (compound/complex/precipitate/ions) OR <br> catalyst $\checkmark$ | 1 | IGNORE 'variable oxidation states' .... but ALLOW the idea that $\mathrm{Fe}^{2+}$ can react to form an ion with a different charge/oxidation state. 'ion' is essential: 'atom' or 'metal' is not sufficient <br> IGNORE partially filled d sub-shell/d orbital (question refers to property of $\mathrm{Fe}^{2+}$ ) |
|  | (c) | Fe oxidised from +2 to $+3 \checkmark$ Cr reduced from +6 to $+3 \checkmark$ | 2 | CHECK and credit oxidation numbers on equation <br> ALLOW $\mathrm{Fe}^{2+}$ oxidised to $\mathrm{Fe}^{3+}$ <br> ALLOW Cr ${ }^{6+}$ reduced to $\mathrm{Cr}^{3+}$ <br> ALLOW + sign after number in oxidation number, ie 2+, etc <br> ALLOW 1 mark only if oxidation numbers given with no identification of which species has been oxidised or reduced, ie <br> Fe goes from +2 to +3 AND Cr goes from +6 to +3 <br> Fe reduced from +2 to +3 AND Cr oxidised from +6 to +3 (oxidation and reduction the wrong way around) <br> DO NOT ALLOW just ' Fe is oxidised and Cr reduced' <br> IGNORE other oxidations numbers (even if wrong) <br> IGNORE any references to electrons |



| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | (e) | (ii) | (For complex) with CO , stability constant is greater (than with complex in $\mathrm{O}_{2}$ ) <br> OR with CO, stability constant is high $\checkmark$ <br> (Coordinate) bond with CO is stronger (than $\mathrm{O}_{2}$ ) OR bond with CO is strong $\checkmark$ | 2 | ANNOTATE WITH TICKS AND CROSSES, etc <br> Comparison of CO and $\mathrm{O}_{2}$ is NOT required ALLOW stability constant with/of CO is greater IGNORE (complex with) CO is more stable <br> ALLOW bond with CO is less likely to break OR bond with CO more likely to form OR 'CO cannot be removed' OR idea that attachment of CO is irreversible OR CO is a stronger ligand (than $\mathrm{O}_{2}$ ) OR CO has greater affinity for ion/metal/haemoglobin (than $\mathrm{O}_{2}$ ) <br> IGNORE CO bonds more easily |
|  | (f) | (i) | $\mathrm{Pt}^{2+} / \mathrm{Pt}$ is $+2 / 2+, 2 \times \mathrm{Cl}^{-}-2 \checkmark$ | 1 | DO NOT ALLOW response in terms of $\mathrm{Cl}_{2}$ rather than $\mathrm{Cl}^{-}$ DO NOT ALLOW 'charges cancel' without the charges involved being stated |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | (f) | (ii) | OR <br> $\checkmark \checkmark$ For each structure <br> Ligand donates an electron pair to metal (ion) $/ \mathrm{Pt}^{2+} / \mathrm{Pt}$ <br> OR forms a coordinate bond to the metal (ion)/ $/ \mathrm{Pt}^{2+} / \mathrm{Pt} \checkmark$ | 3 | IGNORE any charge, ie $\mathrm{Pt}^{2+} \mathrm{ORCl}^{-}$, even if wrong IGNORE any angle, even if wrong <br> ACCEPT bonds to $\mathrm{H}_{3} \mathrm{~N}$ (does not need to go to ' N ') <br> Assume that a solid line is in plane of paper <br> Each structure must contain 2 'out wedges' AND 2 'in wedges' or dotted lines <br> OR 4 solid lines at right angles (all in plane of paper) <br> DO NOT ALLOW any structure that cannot be in one plane DO NOT ALLOW any structure with $\mathrm{Cl}_{2}$ as a ligand DO NOT apply ECF from one structure to the other <br> ALLOW coordinate bonds shown on diagrams provide that they start from a lone pair <br> ALLOW 'dative covalent bond' or 'dative bond' as alternative for 'coordinate bond <br> IGNORE cis and trans labels (even if incorrect) IGNORE incorrect connectivity to $\mathrm{NH}_{3}$, ie ALLOW NH3— |
|  |  | (iii) | platin binds to DNA (of cancer cells) OR platin stops (cancer) cells dividing/replicating $\checkmark$ | 1 |  |


| Question |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 2 | (g) | 1,1-cyclobutanedicarboxylate ion <br> Correct charge required (could also be 2- outside square brackets) <br> carboplatin (cis isomer shown below) | 2 | Must show cyclobutane ring with both $\mathrm{COO}^{-}$groups bonded to same carbon <br> ALLOW $\mathrm{COO}^{-} \mathrm{OR} \mathrm{CO}_{2}^{-}$for each carboxylate ion ALLOW structures showing $\mathrm{CH}_{2}$ or C atoms provided it is clear that C skeleton is shown, <br> Note: H atoms are not required if C atoms shown, ie <br> DO NOT ALLOW circle inside cyclobutane ring <br> Two bonds from Pt to O atoms <br> Any bonds from ligand MUST come from O OR from atom with lone pair <br> IGNORE any charge shown <br> Note: H atoms are not required if C atoms shown, (see ion in 1st structure) <br> ALLOW ECF from 1st structure provided that the attached atoms are capable of forming coordinate bonds (ie they contain a lone pair of electrons) <br> Example if 1st structure is as below, then ALLOW 1 mark ECF  <br> X  <br> ECF. $\checkmark$ |
|  |  | Total | 18 |  |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (a) | (i) | $\mathrm{HOCH}_{2} \mathrm{COOH}+\mathrm{NaOH} \rightarrow \mathrm{HOCH}_{2} \mathrm{COONa}+\mathrm{H}_{2} \mathrm{O}$ | 1 | ALLOW: $\mathrm{HOCH}_{2} \mathrm{COOH}+\mathrm{OH}^{-} \rightarrow \mathrm{HOCH}_{2} \mathrm{COO}^{-}+\mathrm{H}_{2} \mathrm{O}$ <br> ALLOW: $\mathrm{H}^{+}+\mathrm{OH}^{-} \rightarrow \mathrm{H}_{2} \mathrm{O}$ <br> DO NOT ALLOW molecular formulae <br> (cannot see which OH has reacted) |
|  |  | (ii) | FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer $=0.142\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$, award 2 marks $\begin{aligned} & \text { amount of } \mathrm{HOCH}_{2} \mathrm{COOH}=0.125 \times \frac{25.0}{1000} \\ & =0.003125(\mathrm{~mol}) \checkmark \\ & \text { concentration } \mathrm{NaOH}=0.003125 \times \frac{1000}{22.00} \\ & =0.142\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)^{\checkmark} \end{aligned}$ | 2 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below <br> ANNOTATE WITH TICKS AND CROSSES, etc <br> ALLOW $3.125 \times 10^{-3} \mathrm{~mol}$ <br> ALLOW ECF: answer above $\times \frac{1000}{22.00}$ <br> ALLOW 2 SF: 0.14 to calculator value: 0.142045454 <br> If candidate has written in (a)(i): $\mathrm{HOCH}_{2} \mathrm{COOH}+\mathbf{2 N a O H}$, mark by ECF: <br> concentration $\mathrm{NaOH}=2 \times 0.003125 \times \frac{1000}{22.00}$ $=0.284\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ |
|  |  | (iii) | Vertical section matches the ( pH ) range (of the indicator) <br> OR colour change (of the indicator) <br> OR end point (of the indicator) $\checkmark$ | 1 | ALLOW stated pH range for vertical section at about 7-10, 6-10, etc <br> ie ALLOW ' pH range must be about 7-10' <br> ALLOW 'pH changes rapidly' for vertical section ALLOW 'equivalence point' for vertical section, ie ALLOW equivalence point matches the $(\mathrm{pH})$ range, etc <br> DO NOT ALLOW just 'end point matches $(\mathrm{pH})$ range' DO NOT ALLOW just 'indicator matches vertical section' <br> Response must link either the pH range or colour change or end point with the vertical section / pH range $\sim 7-10$ |

\begin{tabular}{|c|c|c|c|c|c|}
\hline \multicolumn{3}{|r|}{Question} \& Answer \& Mark \& Guidance \\
\hline 3 \& (b) \& (i) \& \[
\left(K_{\mathrm{a}}=\right) \frac{\left[\mathrm{H}^{+}\right]\left[\mathrm{HOCH}_{2} \mathrm{COO}^{-}\right]}{\left[\mathrm{HOCH}_{2} \mathrm{COOH}\right]} \checkmark
\] \& 1 \& IGNORE state symbols IGNORE \(\qquad\) \(\frac{\left[\mathrm{H}^{+}\right]^{2}}{\left[\mathrm{HOCH}_{2} \mathrm{COOH}\right]}\) in (i) but ALLOW in (ii) \\
\hline \& \& (ii) \& \begin{tabular}{l}
FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer \(=1.46 \times 10^{-4}\), award 2 marks THEN IF units are \(\mathrm{mol} \mathrm{dm}^{-3}\), award 1 further mark
\[
\begin{aligned}
\& {\left[\mathrm{H}^{+}\right]=10^{-2.37}=0.00427\left(\mathrm{~mol} \mathrm{dm}^{-3}\right) \checkmark} \\
\& K_{\mathrm{a}}=\frac{0.00427^{2}}{0.125}=1.46 \times 10^{-4} \checkmark
\end{aligned}
\] \\
units: \(\mathrm{mol} \mathrm{dm}^{-3} \checkmark\)
\end{tabular} \& 2

1 \& | IF there is an alternative answer, check to see if there is any ECF credit possible using working below UNITS can be credited with no numerical answer |
| :--- |
| ANNOTATE WITH TICKS AND CROSSES, etc |
| ALLOW $4.27 \times 10^{-3}$ (mol) |
| ALLOW 2 SF: 0.0043 up to 0.004265795188 (calc value) |
| IF candidate has rounded to $0.00427\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ in 1 st response, credit |
| EITHER |
| 2 SF: $1.5 \times 10^{-4}$ up to $1.458632 \times 10^{-4}$ (from 0.00427 ) |
| OR |
| 2 SF: $1.5 \times 10^{-4}$ up to $1.455760687 \times 10^{-4}$ (from unrounded calculator value of 0.004265795188 ) |
| ALLOW calculation based on equilibrium conc of glycolic acid as $0.125-\left[\mathrm{H}^{+}\right]$: |
| Using $\left[\mathrm{H}^{+}\right]=0.00427, K_{\mathrm{a}}=\frac{0.00427^{2}}{0.125-0.00427}=1.51 \times 10^{-4}$ |
| For UNITS this is the ONLY correct answer | \\

\hline \& \& (iii) \& | $\% \text { dissociation }=\frac{0.00427}{0.125} \times 100=3.4(\%) \checkmark$ |
| :--- |
| Assume working from EITHER from a rounded $\left[\mathrm{H}^{+}\right]$ OR unrounded calculator value of $\mathbf{b}$ (ii) $\left[\mathrm{H}^{+}\right]$ | \& 1 \& | ALLOW ECF using calculated $\left[\mathrm{H}^{+}\right]$from $\mathbf{b}$ (ii), ALLOW 2 SF: 3.4 \% up to calculator value |
| :--- |
| Note: $\left[\mathrm{H}^{+}\right]$from b(ii) displayed at top of answer window DO NOT MARK THIS TWICE! | \\

\hline
\end{tabular}



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| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (e) |  | $\begin{aligned} & 2 \mathrm{HSCH}_{2} \mathrm{COO}^{-}+\mathrm{R}-\mathrm{S}-\mathrm{S}-\mathrm{R} \\ & \\ & \longrightarrow \mathrm{OOCCH}_{2} \mathrm{~S}-\mathrm{SCH}_{2} \mathrm{COO}^{-}+2 \mathrm{R}-\mathrm{SH} \checkmark \\ & 2 \mathrm{R}-\mathrm{SH}+\mathrm{H}_{2} \mathrm{O}_{2} \longrightarrow \mathrm{R}-\mathrm{S}-\mathrm{S}-\mathrm{R}+2 \mathrm{H}_{2} \mathrm{O} \end{aligned}$ | 2 | ALLOW $\left(\mathrm{SCH}_{2} \mathrm{COO}^{-}\right)_{2}$ <br> ALLOW equation with ammonium salt, ie: $\begin{aligned} & 2 \mathrm{HSCH}_{2} \mathrm{COONH}_{4}+\ldots \ldots . . \\ & \longrightarrow \mathrm{H}_{4} \mathrm{NOOCCH}_{2} \mathrm{~S}-\mathrm{SCH}_{2} \mathrm{COONH}_{4}+\ldots \ldots . . \end{aligned}$ |
|  |  |  | Total | 20 |  |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (a) | (i) | Complete circuit with electrodes to voltmeter AND salt bridge between solutions <br> $\mathrm{Sn}^{4+} / \mathrm{Sn}^{2+}$ half cell with Pt electrode AND both solutions labelled as $1 \mathrm{~mol} \mathrm{dm}^{-3} / 1 \mathrm{M}$ <br> $\mathrm{H}^{+} / \mathrm{H}_{2}$ half cell with Pt electrode AND $\mathrm{H}^{+}$solution labelled as $1 \mathrm{~mol} \mathrm{dm}^{-3} / 1 \mathrm{M} \checkmark$ | 3 | ANNOTATE WITH TICKS AND CROSSES, etc circuit shown must be complete, ie must be capable of working salt bridge must be labelled and must dip into both solutions <br> ALLOW concentration label of 'equimolar' or similar wording for $\mathrm{Sn}^{4+} / \mathrm{Sn}^{2+}$ half cell <br> ALLOW any strong acid <br> IF both half cells are correct with no concentrations, ALLOW 1 out of the 2 marks available for the 2 half cells <br> IGNORE any stated temperature or pressure, even if wrong |
|  |  | (ii) | $\begin{array}{ll} 2 \mathrm{Cr}+3 \mathrm{Sn}^{4+} & \rightarrow \quad 2 \mathrm{Cr}^{3+}+3 \mathrm{Sn}^{2+} \\ \mathrm{Cr}+3 \mathrm{Cu}^{+} \rightarrow & \mathrm{Cr}^{3+}+3 \mathrm{Cur} \\ \mathrm{Sn}^{2+}+2 \mathrm{Cu}^{+} & \rightarrow \quad \mathrm{Sn}^{4+}+2 \mathrm{Cu} \checkmark \end{array}$ <br> Conditions not standard OR concentrations not $1 \mathrm{~mol} \mathrm{dm}^{-3} \checkmark$ <br> High activation energy OR slow rate $\checkmark$ | 5 | ANNOTATE WITH TICKS AND CROSSES, etc Correct species AND balancing needed for each mark ALLOW equations as shown with equilibrium sign ALLOW multiples but electrons must not be shown <br> IF three equations have correct species but no balancing, AWARD 1 mark <br> ALLOW not favoured kinetically |
|  | (b) | (i) | $\mathrm{CH}_{3} \mathrm{OH}+11 / 2 \mathrm{O}_{2} \rightarrow \mathrm{CO}_{2}+2 \mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 | Correct species AND balancing needed ALLOW multiple, ie $2 \mathrm{CH}_{3} \mathrm{OH}+3 \mathrm{O}_{2} \quad \rightarrow \quad 2 \mathrm{CO}_{2}+4 \mathrm{H}_{2} \mathrm{O}$ ALLOW CH ${ }_{4} \mathrm{O}$ for formula of methanol |
|  |  | (ii) | $\mathrm{CH}_{3} \mathrm{OH}+\mathrm{H}_{2} \mathrm{O} \rightarrow 6 \mathrm{H}^{+}+6 \mathrm{e}^{-}+\mathrm{CO}_{2} \checkmark$ | 1 |  |
|  |  | (iii) | less $\mathrm{CO}_{2}$ OR less greenhouse gases $\checkmark$ greater efficiency $\checkmark$ | 2 | ALLOW no $\mathrm{CO}_{2}$ OR no greenhouse gases ALLOW (very) efficient <br> IGNORE less pollution OR 'renewable fuels' |
|  |  | (iv) | methanol is a liquid <br> AND <br> methanol is easier to store/transport $\checkmark$ | 1 | Both points required for mark Response MUST state that methanol is a liquid IGNORE methanol has a higher boiling point Assume that 'it' refers to methanol IGNORE safety issues, eg $\mathrm{H}_{2}$ leakage, flammability, explosive |
|  |  |  | Total | 13 |  |

\begin{tabular}{|c|c|c|c|c|}
\hline \multicolumn{2}{|l|}{Question} \& Answer \& Mark \& Guidance \\
\hline 5 \& (a) \& \begin{tabular}{l}
A: forms fewer moles/molecules of gas \(\checkmark\) \\
B: forms gas from a liquid \(\checkmark\) \\
C: forms liquid from gases \(\checkmark\) \\
D: forms more moles/molecules of gas
\end{tabular} \& 4 \& \begin{tabular}{l}
Note: Responses must imply the key difference between the sides of the equation \\
IGNORE comments about C(s)
\end{tabular} \\
\hline \& (b) \& \begin{tabular}{l}
\[
\begin{aligned}
\& \Delta S=\Sigma S \text { (products) }-\Sigma S(\text { reactants) } \\
\& =40+214-89=165\left(\mathrm{~J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1}\right) \\
\& =0.165\left(\mathrm{~kJ} \mathrm{~K}^{-1} \mathrm{~mol}^{-1}\right) \checkmark
\end{aligned}
\] \\
At \(25^{\circ} \mathrm{C}, \Delta \mathrm{G}=+178-298 \times 0.165\)
\[
\begin{array}{ll}
=(+) 129 \checkmark \& \text { units: } \mathrm{kJ} \mathrm{~mol}^{-1} \checkmark \\
\text { OR }(+) 129,000 \checkmark \& \text { units: } \mathrm{J} \mathrm{~mol}^{-1} \checkmark
\end{array}
\] \\
As \(\Delta G>0\), reaction is not feasible OR as \(\Delta G>0, \mathrm{CaCO}_{3}\) is stable \(\checkmark\) \\
Minimum temperature for feasibility when
\[
\begin{aligned}
\& 0=\Delta H-T \Delta S \text { OR } \Delta H=T \Delta S \text { OR } T=\frac{\Delta H}{\Delta S} \\
\& =\frac{178}{0.165}=1079 \mathrm{~K} \mathrm{OR} 806^{\circ} \mathrm{C}
\end{aligned}
\] \\
The units must be with the stated temperature
\end{tabular} \& 4

2 \& | ANNOTATE WITH TICKS AND CROSSES, etc |
| :--- |
| Mark is for the working line: $40+214-89=165$ |
| UNITS have a separate mark |
| ALLOW 129 to calculator value of 128.83 |
| DO NOT ALLOW 128 (incorrect rounding) |
| IF $25^{\circ} \mathrm{C}$ used rather than 298 K , credit by ECF, calculated $\Delta G$ |
| $=174$ to calculator value of 173.875 |
| ENTROPY APPROACH- |
| ALLOW At $25^{\circ} \mathrm{C}, \Delta S_{\text {total }}=0.165-\frac{178}{298} \checkmark$ $=-0.432 \checkmark \mathrm{~kJ} \mathrm{~K}^{-1} \mathrm{~mol}^{-1} \checkmark$ $\text { OR -432 } \checkmark \mathrm{J} \mathrm{~K}^{-1} \mathrm{~mol}^{-1} \checkmark$ |
| As $\Delta S<0$, reaction is not feasible |
| ENTROPY APPROACH- |
| Minimum temperature for feasibility when $0=\Delta S_{\text {system }}+\Delta S_{\text {surroundings }} \text { OR } \quad \Delta S_{\text {system }}=\frac{\Delta H}{T}$ |
| ALLOW 1080 K up to calculator value of 1078.787879 , correctly rounded, eg 1078.79 is correct value to 6SF DO NOT ALLOW 1078 (incorrect rounding) |
| IF 1079 K is given and additional temperature in ${ }^{\circ} \mathrm{C}$ is incorrect, IGNORE ${ }^{\circ} \mathrm{C}$ temperature (and vice versa) | \\

\hline \& \& Total \& 11 \& \\
\hline
\end{tabular}

| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 6 | (a) | (i) | $\left(K_{\mathrm{w}}=\right.$ ) $\left[\mathrm{H}^{+}(\mathrm{aq})\right]\left[\mathrm{OH}^{-}(\mathrm{aq})\right] \checkmark$ | 1 | IGNORE state symbols ALLOW $\left[\mathrm{H}_{3} \mathrm{O}^{+}(\mathrm{aq})\right]\left[\mathrm{OH}^{-}(\mathrm{aq})\right]$ |
|  |  | (ii) | FIRST, CHECK THE ANSWER ON ANSWER LINE IF answer $=2.3 \times 10^{-10}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$, award 2 marks IF answer $=2.34 \times 10^{-10}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$, award 1 mark $\qquad$ $\begin{aligned} & {\left[\mathrm{H}^{+}\right]=10^{-\mathrm{pH}}=4.27 \times 10^{-5}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right) \checkmark} \\ & {\left[\mathrm{OH}^{-}\right]=\frac{1.0 \times 10^{-14}}{4.27 \times 10^{-5}}} \\ & =2.34 \times 10^{-10} \\ & =2.3 \times 10^{-10}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)^{\checkmark} \end{aligned}$ | 2 | IF there is an alternative answer, check to see if there is any ECF credit possible using working below ANNOTATE WITH TICKS AND CROSSES, etc <br> ALLOW $4.3 \times 10^{-5}$ up to calculator: $4.265795188 \times 10^{-5}$ ALLOW 0.0000427 <br> Answer MUST be to 2 SF (in question) ALLOW $=2.3 \times 10^{-x}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$ for 1 mark (must be a negative power) <br> ALLOW alternative approach based on pOH : $\begin{aligned} & \mathrm{pOH}=14-4.27=9.63 \checkmark(\text { DO NOT ALLOW 9.6 }) \\ & {\left[\mathrm{OH}^{-}\right]=10^{-\mathrm{pOH}}=10^{-9.63}=2.3 \times 10^{-10}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)^{\checkmark}} \end{aligned}$ |
|  | (b) | (i) | Endothermic because $K_{w}$ increases with temperature | 1 | Endothermic AND reason required for the mark ALLOW Endothermic because increasing temperature shifts equilibrium/reaction to the right |
|  |  | (ii) | $K_{w}$ value from graph from 2.2 to $2.6 \times 10^{-14}\left(\mathrm{~mol}^{2}\right.$ $\left.\mathrm{dm}^{-6}\right)^{\vee}$ <br> Using $2.4 \times 10^{-14}$, <br> $\left[\mathrm{H}^{+}\right]=\sqrt{2.4 \times 10^{-14}}$ OR $1.55 \times 10^{-7} \checkmark$ $\begin{aligned} & \mathrm{pH}=-\log \left(1.55 \times 10^{-7}\right)=6.81 \\ & \left(\text { using } K_{\mathrm{w}}=2.4 \times 10^{-14}\right)^{\checkmark} \end{aligned}$ | 3 | ANNOTATE WITH TICKS AND CROSSES, etc <br> Actual $K_{w}=2.38 \times 10^{-14} \mathrm{~mol}^{2} \mathrm{dm}^{-6}$ <br> For this mark, candidate must use a value between 2.0 and $3.0 \times 10^{-14}\left(\mathrm{~mol}^{2} \mathrm{dm}^{-6}\right)$, ie from the approximately correct region of the graph, <br> ALLOW 6.8 up to calculator value <br> Note: You will need to calculate the pH value from the candidate's estimate of $K_{\mathrm{w}}$ at $37^{\circ} \mathrm{C}$ before awarding the 3 rd marking point <br> ONLY award an ECF pH mark if candidate has generated a value of $\left[\mathrm{H}^{+}\right]$by attempting to take a square root of a value between 2.0 and $3.0 \times 10^{-14}$ |



| Question |  | Answer | Mark | Guidance |
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| 6 | (d) | Ionic radius <br> Potassium ion OR K ${ }^{+}$OR K ion is smaller <br> OR $\mathrm{K}^{+}$has greater charge density $\checkmark$ <br> Lattice enthalpy <br> Lattice enthalpy of KF is more negative than RbF $\checkmark$ OR <br> $\mathrm{K}^{+}$has greater attraction for $\mathrm{F}^{-}$ <br> Hydration enthalpy <br> $\Delta H$ (hydration) of $\mathrm{K}^{+}$is more negative than $\mathrm{Rb}^{+}$ OR <br> $\mathrm{K}^{+}$has greater attraction for $\mathrm{H}_{2} \mathrm{O}$ <br> Enthalpy change of solution Idea that $\Delta H$ (solution) is affected more by lattice enthalpy than by hydration enthalpy $\checkmark$ | 4 | ANNOTATE WITH TICKS AND CROSSES, etc <br> Throughout question, ORA in terms of $\mathrm{Rb}^{+}$ <br> Throughout question, ALLOW energy for enthalpy <br> DO NOT ALLOW potassium OR K OR reference to atoms (ie reference to ions is required throughout a response) <br> ALLOW lattice enthalpy of KF > lattice enthalpy of RbF <br> ALLOW more energy needed to separate $\mathrm{K}^{+}$AND $\mathrm{F}^{-}$ IGNORE KF has stronger bonds <br> ALLOW $\Delta H$ (hydration) of $\mathrm{K}^{+}>\Delta H$ (hydration) of $\mathrm{Rb}^{+}$ <br> ALLOW more energy needed to separate $\mathrm{K}^{+}$AND $\mathrm{H}_{2} \mathrm{O}$ IGNORE $\mathrm{K}^{+}$has a stronger bond to $\mathrm{H}_{2} \mathrm{O}$ <br> ALLOW a correct attempt to link the contribution of lattice enthalpy and hydration enthalpy to $\Delta H$ (solution), ie lattice enthalpy is a more important factor than hydration enthalpy |
|  | (e) | (During dissolving,) entropy/disorder increases OR disorder increases <br> $T \Delta S>\Delta H$ <br> OR $T \Delta S$ is more positive than $\Delta H$ OR $\Delta H-T \Delta S$ is negative $\checkmark$ | 2 | ALLOW entropy change is positive OR $\Delta S$ is positive OR $T \Delta S$ is positive <br> ALLOW $\Delta \mathrm{S}$ (system) $>\Delta H / T$ <br> ALLOW $\Delta S$ (system) is more positive than $\Delta H / T \checkmark$ <br> ALLOW $\Delta S$ (system) $+\Delta S$ (surroundings) is positive <br> ALLOW Energy contribution from increase in entropy is greater than decrease in energy from enthalpy change OR entropy change outweighs enthalpy change <br> IGNORE $\Delta G$ is negative |
|  |  | Total | 20 |  |


| Question |  | Answer | Mark | Guidance |
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| 7 | (a) | (i) | amount $\mathbf{S}_{2} \mathbf{O}_{3}{ }^{2-}$ used <br> $=0.00100 \times \frac{24.6}{1000}=2.46 \times 10^{-5} \mathrm{~mol} \checkmark$ <br> amount $\mathbf{O}_{2}$ in $25 \mathrm{~cm}^{3}$ sample <br> $=\frac{2.46 \times 10^{-5}}{4}=6.15 \times 10^{-6} \mathrm{~mol} \checkmark$ <br> Concentration of $\mathbf{O}_{2}$ in sample <br> $=6.15 \times 10^{-6} \times \frac{1000}{25}=2.46 \times 10^{-4}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right) \checkmark$ <br> mass concentration of $\mathbf{O}_{2}$ in $\mathrm{mg} \mathrm{dm}^{-3}$ <br> $=2.46 \times 10^{-4} \times 32 \mathrm{~g}=7.872 \times 10^{-3}\left(\mathrm{~g} \mathrm{dm}^{-3}\right)$ <br> $=7.872\left(\mathrm{mg} \mathrm{dm}^{-3}\right) \checkmark$ | ANNOTATE WITH TICKS AND CROSSES, etc <br> ALLOW $0.0000246(\mathrm{~mol})$ |


| Question |  |  | Answer | Mark | Guidance |
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| 7 | (b) | (ii) | $\begin{aligned} & 2 \mathrm{H}_{2} \mathrm{O}+2 \mathrm{I}^{-}+2 \mathrm{NO}_{2}^{-} \longrightarrow 2 \mathrm{NO}+\mathrm{I}_{2}+4 \mathrm{OH}^{-} \\ & \begin{array}{l} \mathrm{OR} 2 \mathrm{H}^{+}+2 \mathrm{I}^{-}+2 \mathrm{NO}_{2}^{-} \longrightarrow 2 \mathrm{NO}+\mathrm{I}_{2}+2 \mathrm{OH}^{-} \\ \text {species } \checkmark \\ \text { balance } \checkmark \end{array} \end{aligned}$ | 2 | IGNORE state symbols <br> ALLOW multiples <br> For species ONLY, IGNORE any extra $\mathrm{H}_{2} \mathrm{O}$ or $\mathrm{e}^{-}$on either <br> side of the equation <br> ALLOW on LHS: $2 \mathrm{HI}+2 \mathrm{NO}_{2}{ }^{-}$OR $2 \mathrm{I}^{-}+2 \mathrm{HNO}_{2}$ <br> ALLOW species and equation involving $\mathrm{N}_{2} \mathrm{H}_{2}$ : $\begin{aligned} & \quad 6 \mathrm{H}_{2} \mathrm{O}+8 \mathrm{I}^{-}+2 \mathrm{NO}_{2}^{-} \longrightarrow \mathrm{N}_{2} \mathrm{H}_{2}+4 \mathrm{I}_{2}+10 \mathrm{OH}^{-} \\ & \text {OR } 6 \mathrm{H}^{+}+8 \mathrm{I}^{-}+2 \mathrm{NO}_{2}^{-} \longrightarrow \mathrm{N}_{2} \mathrm{H}_{2}+4 \mathrm{I}_{2}+4 \mathrm{OH}^{-} \\ & \text {species } \checkmark \\ & \text { balance } \checkmark \end{aligned}$ |
|  |  |  | Total | 8 |  |

