# Mark Scheme (Results) January 2008 

## GCE

## GCE Chemistry (6244) Paper 1

## General Marking Guidance

- All candidates must receive the same treatment. Examiners must mark the first candidate in exactly the same way as they mark the last.
- Mark schemes should be applied positively. Candidates must be rewarded for what they have shown they can do rather than penalised for omissions.
- Examiners should mark according to the mark scheme not according to their perception of where the grade boundaries may lie.
- There is no ceiling on achievement. All marks on the mark scheme should be used appropriately.
- $\quad$ All the marks on the mark scheme are designed to be awarded. Examiners should always award full marks if deserved, i.e. if the answer matches the mark scheme. Examiners should also be prepared to award zero marks if the candidate's response is not worthy of credit according to the mark scheme.
- Where some judgement is required, mark schemes will provide the principles by which marks will be awarded and exemplification may be limited.
- When examiners are in doubt regarding the application of the mark scheme to a candidate's response, the team leader must be consulted.
- Crossed out work should be marked UNLESS the candidate has replaced it with an alternative response.


## Using the mark scheme

1 / means that the responses are alternatives and either answer should receive full credit.
2 ( ) means that a phrase/word is not essential for the award of the mark, but helps the examiner to get the sense of the expected answer.
3 [ ] words inside square brackets are instructions or guidance for examiners.
4 Phrases/words in bold indicate that the meaning of the phrase or the actual word is essential to the answer.
$5 \mathrm{ecf} / \mathrm{TE} / \mathrm{cq}$ (error carried forward) means that a wrong answer given in an earlier part of a question is used correctly in answer to a later part of the same question.

| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 1.(a)(i) | Ionic | Giant ionic or <br> electrovalent |  | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 1.(a)(ii) | Covalent | Giant covalent | Convalent | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 1.(b)(i) | Basic | Base or alkali <br> or <br> alkaline |  | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 1.(b)(ii) | Acidic | Acid <br> Weakly acidic <br> Weak acid |  | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :---: | :---: | :---: |
| 1.(c)(i) | $3 \mathrm{Na}_{2} \mathrm{O}+2 \mathrm{H}_{3} \mathrm{PO}_{4} \rightarrow 2 \mathrm{Na}_{3} \mathrm{PO}_{4}+3 \mathrm{H}_{2} \mathrm{O}$ <br> OR <br> $\mathrm{Na}_{2} \mathrm{O}+\mathrm{H}_{3} \mathrm{PO}_{4} \rightarrow \mathrm{Na}_{2} \mathrm{HPO}_{4}+\mathrm{H}_{2} \mathrm{O}$ <br> OR <br> $\mathrm{Na}_{2} \mathrm{O}+2 \mathrm{H}_{3} \mathrm{PO}_{4} \rightarrow 2 \mathrm{NaH}_{2} \mathrm{PO}_{4}+\mathrm{H}_{2} \mathrm{O} \quad$ (1) <br> Ignore state symbols |  |  | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 1.(c)(ii) | $\mathrm{SiO}_{2}+2 \mathrm{NaOH} \rightarrow \mathrm{Na}_{2} \mathrm{SiO}_{3}+\mathrm{H}_{2} \mathrm{O}$ (1) <br> lgnore state symbols | $\mathrm{SiO}_{2}+2 \mathrm{OH}^{-} \rightarrow \mathrm{SiO}_{3}^{-2}+\mathrm{H}_{2} \mathrm{O}$ |  | 1 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 1.(d) | First mark: <br> $\mathrm{Al}_{2} \mathrm{O}_{3(\mathrm{~s})}+6 \mathrm{H}^{+}{ }_{(\mathrm{aq})} \rightarrow 2 \mathrm{Al}^{3+}{ }_{(\mathrm{aq})}+3 \mathrm{H}_{2} \mathrm{O}_{(\mathrm{l})}$ <br> This mark is for correct species and balancing <br> Second mark: $\mathrm{Al}_{2} \mathrm{O}_{3(\mathrm{~s})}+2 \mathrm{OH}_{(\mathrm{aq})}^{-}+3 \mathrm{H}_{2} \mathrm{O}_{(\mathrm{l})} \rightarrow 2 \mathrm{Al}(\mathrm{OH})_{4(\mathrm{aq})}^{-}$ <br> OR $\mathrm{Al}_{2} \mathrm{O}_{3(\mathrm{~s})}+6 \mathrm{OH}_{(\mathrm{aq})}^{-}+3 \mathrm{H}_{2} \mathrm{O}_{(\mathrm{l})} \rightarrow 2 \mathrm{Al}(\mathrm{OH})_{6(\mathrm{aq})}^{3-}$ <br> OR $\mathrm{Al}_{2} \mathrm{O}_{3(\mathrm{~s})}+2 \mathrm{OH}_{(\mathrm{aq})} \rightarrow 2 \mathrm{AlO}_{2(\mathrm{aq})}+\mathrm{H}_{2} \mathrm{O}(\mathrm{l})$ <br> This mark is for correct species and balancing <br> Third mark is for the state symbols <br> Correct state symbols in either equation, but all species must be correct. <br> This mark may be awarded from an unbalanced equation. | Two correct 'molecular’ equations with correct state symbols scores (2) |  | 3 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 1.(e) | $\mathrm{PbO}_{2}+4 \mathrm{HCl} \rightarrow \mathrm{PbCl}_{2}+\mathrm{Cl}_{2}+2 \mathrm{H}_{2} \mathrm{O}$ <br> Ignore state symbols | $\mathrm{PbO}_{2}+6 \mathrm{HCl} \rightarrow \mathrm{H}_{2} \mathrm{PbCl}_{4}$ <br> $\mathrm{Cl}_{2}+2 \mathrm{H}_{2} \mathrm{O}$ |  | 1 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 1.(f) | First mark: |  |  | 2 |
|  | Tin more stable in the +4 oxidation state (than the +2 oxidation state) whereas lead more stable in the +2 oxidation state (than in the +4 oxidation state) <br> OR <br> +2 oxidation state becomes more stable relative to +4 oxidation state as group descended. (1) <br> Second Mark: <br> (So) $I_{2}$ reduced to $I^{-}$(by $\mathrm{Sn}^{2+}$ ) <br> OR <br> $\mathrm{Sn}^{2+}+\mathrm{I}_{2} \rightarrow \mathrm{Sn}^{4+}+2 \mathrm{I}^{-}$ <br> OR <br> Therefore $\operatorname{tin}$ (II) is a strong(er) reducing agent (than lead(II)) (1) | redox reaction between $\mathrm{Sn}^{2+}$ and $\mathrm{I}_{2}$ OR <br> $\mathrm{Sn}^{2+}$ oxidised (to $\mathrm{Sn}^{4+}$ ) OR <br> Sn (II) acts as (a strong) reducing agent | $\mathrm{Sn}^{2+}$ ions less stable than $\mathrm{Pb}^{2+}$ <br> OR <br> Pb (II) is more stable than Sn (II) |  |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 2.(a) | IGNORE 'alkane' in any answer <br> X: ester (1) <br> Y: both <br> alkene <br> and <br> alcohol or hydroxyl (1) | carbon-carbon double <br> bond <br> "hydroxy" | carbonyl |  |
| Z: both alcohol or hydroxyl <br> and <br> aldehyde (1) | "hydroxide" <br> OH $^{-}$or <br> "hydroxide" or <br> "carbonyl" <br> Just the formula |  |  |  |



| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 2. (c)(i) |  <br> (1) <br> Allow $\mathrm{C}_{3} \mathrm{H}_{7} / \mathrm{C}_{2} \mathrm{H}_{5} \mathrm{CH}_{2}$ <br> (1) |  <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{COO}^{-} /$ <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CO}_{2} \mathrm{Na} /$ <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{CH}_{2} \mathrm{CO}_{2}^{-} \mathrm{Na}^{+}$ | Carboxylic acid <br> Or $. . . \mathrm{O}^{-}-\mathrm{Na}^{+}$ | 2 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 2. (c)(ii) |  <br> (1) <br> / $\mathrm{CH}_{2} \mathrm{CHCOONa} / \mathrm{CH}_{2} \mathrm{CHCO}_{2} \mathrm{Na} /$ $/ \mathrm{CH}_{2} \mathrm{CHCOO}^{-} \mathrm{Na}^{+} / \mathrm{CH}_{2} \mathrm{CHCO}_{2}^{-} \mathrm{Na}^{+}$ $/ \mathrm{CH}_{2}=\mathrm{CHCOONa} / \mathrm{CH}_{2}=\mathrm{CHCO}_{2} \mathrm{Na}$ $/ \mathrm{CH}_{2}=\mathrm{CHCOO}^{-} \mathrm{Na}^{+} / \mathrm{CH}_{2}=\mathrm{CHCO}_{2}{ }^{-} \mathrm{Na}^{+}$(1) | $\mathrm{CH}_{2} \mathrm{CHCOO}^{-}$ <br> Allow carboxylic acid as product e.g. $\mathrm{CH}_{2} \mathrm{CHCOOH}$ |  | 2 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 2. (c)(iii) |  |  |  | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| $\mathbf{3 . ( a ) ( i )}$ | To slow down the reaction/to stop the <br> reaction <br> OR to quench the reaction <br> OR to freeze the (position of) <br> equilibrium OWTTE (1) <br> so that the (equilibrium) <br> concentrations/amounts do not change <br> (1) | To stop equilibrium <br> shifting to the left | 2 |  |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 3.(a)(ii) | First mark: $\left[H_{2(g)}\right]=\left[I_{2(g)}\right]$ <br> OR <br> Use of $\left(5.0 \times 10^{-4}\right)^{2}$ <br> Second mark: $\left[H I_{(g)}\right]^{2}=\frac{\left(5.0 \times 10^{-4}\right)^{2}}{0.019}$ <br> OR <br> $\left.0.019=\frac{\left(5.0 \times 10^{-4}\right.}{\left[\mathrm{HI}_{(\mathrm{g})}\right)^{2}}\right)^{2}$ <br> OR $\begin{equation*} [\mathrm{HI}(\mathrm{~g})]=\int\left(\frac{\left(5.0 \times 10^{-4}\right)^{2}}{0.019}\right) \tag{1} \end{equation*}$ <br> Third mark: $\begin{equation*} \left[H I_{(g)}\right]=3.6 \times 10^{-3}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right) \tag{1} \end{equation*}$ <br> Correct answer scores 3 marks. Ignore state symbols. Ignore units unless wrong. Ignore s.f. | If [HI] not squared, first mark only. | If first mark not awarded, total (0). | 3 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 3.(b)(i) | $K_{p}=\frac{p_{H I}^{2}}{p_{H_{2}} \times p_{I_{2}}}$ |  |  |  |
|  |  |  |  | 1 |
|  | Ignore position of any ( ) scores (0) |  |  |  |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 3. (b)(ii) | Each step of this calculation must be looked at. <br> $1^{\text {st }}$ mark is for calculating equilibrium moles $\begin{array}{\|l} \mathrm{H}_{2}=0.2 \\ \mathrm{I}_{2}=0.2 \\ \mathrm{HI}=1.6 \tag{1} \end{array}$ <br> $2^{\text {nd }}$ mark is for dividing these by 2 (to get mole fractions) $\begin{align*} & x_{H_{2}}=\frac{0.2}{2.0}=0.1 \\ & x_{I_{2}}=\frac{0.2}{2.0}=0.1 \\ & x_{H I}=\frac{1.6}{2.0}=0.8 \tag{1} \end{align*}$ <br> $3^{\text {rd }}$ mark is for multiplying by 1.1 (to get partial pressures) $\begin{align*} \mathrm{P}_{\mathrm{H}_{2}} & =\frac{0.2}{2.0} \times 1.1 \\ & =0.11(\mathrm{~atm}) \\ \mathrm{P}_{\mathrm{I}_{2}}= & \frac{0.2}{2.0} \times 1.1 \\ & =0.11(\mathrm{~atm}) \\ \mathrm{P}_{\mathrm{HI}} & =\frac{1.6}{2.0} \times 1.1 \\ & =0.88(\mathrm{~atm}) \tag{1} \end{align*}$ <br> $4^{\text {th }}$ mark is for substituting into their expression and calculating $K_{p}$ $\begin{align*} K_{P}= & \frac{(0.88)^{2}}{(0.11) \times(0.11)} \\ = & 64 \tag{1} \end{align*}$ <br> Ignore s.f. <br> Correct answer with no working scores (1) | Mark consequentially <br> Mark consequentially <br> Mark consequentially <br> If moles HI given as $0.8, K_{\mathrm{p}}=16 \max (3)$ |  | 4 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 3.(b)(iii) | Same number of moles on each side | 'Powers cancel' | 'Partial pressures | 1 |
|  | OR | OR | cancel' |  |
|  | (Total) pressure cancels | OR cancel' | OR | 'mol dm ${ }^{-3}$ cancel' |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 4.(a)(i) | $\Delta H_{6}$ |  |  |  |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :--- | :--- |
| 4.(a)(ii) | $\frac{\Delta H_{5}}{2} O R \frac{1}{2} \Delta H_{5}$ |  | $\Delta H_{5}$ | 1 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 4.(b) | Either $\Delta H f=\Delta H_{2}+\Delta H_{3}+\Delta H_{4}+\Delta H_{5}+\Delta H_{6}$ <br> OR $\begin{aligned} & \Delta H f=(+178)+(1735)+2 \times(+218)+2 \times \\ & (-73)+(-2389) \end{aligned}$ $=-186\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$ <br> Correct answer with no working (2) Ignore kJ | [First mark only if doubles both $\Delta H_{a t}$ and electron affinity for hydrogen] <br> [2nd mark is only consequential on failure to multiply either $\Delta$ Hat or electron affinity or both giving: $-404 /-113 /-331$ (kJ mol-1)] | +186 scores (0) $\begin{aligned} & +404 /+113 /+331 \\ & \text { scores (0) } \end{aligned}$ | 2 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 4. (c) | EITHER <br> First mark: <br> Magnesium/Mg ion smaller (radius) <br> than calcium/ Ca ion <br> Or <br> the sum of the ionic radii in $\mathrm{MgH}_{2}$ <br> smaller (than in $\mathrm{CaH}_{2}$ ) (1) <br> Second mark: <br> but charges the same (1) <br> Third mark: <br> (so) stronger (forces of) attraction <br> between ions (in $\mathrm{MgH}_{2}$ ) <br> [Correct reverse arguments can score both marks] <br> OR <br> First and second mark combined: $\mathrm{Mg}^{2+}$ (ion) or $\mathrm{Mg}^{2+}$ (cation) smaller (radius) than $\mathrm{Ca}^{2+}$ <br> Third mark: <br> (so) stronger (forces of) attraction <br> between ions (in $\mathrm{MgH}_{2}$ ) <br> [Correct reverse arguments can score both marks] <br> Ignore references to polarisation of the hydride ion or "covalent character" in the hydrides. <br> Ignore references to "energy required to separate ions/break bonds" | Magnesium ion has greater charge density than calcium ion for first mark. <br> "stronger ionic bonding" for $3^{\text {rd }}$ mark in either case. | Reference to 'atoms’ or 'molecules' or ' $\mathrm{H}_{2}$ ' scores zero overall. <br> If " $\mathrm{H}^{+}$ions" or "hydrogen ions" referred to, $3^{\text {rd }}$ mark cannot be awarded in either case <br> If just "stronger bonding in $\mathrm{MgH}_{2}$ ", $3^{\text {rd }}$ mark cannot be awarded in either case | 3 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 4.(d)(i) | Enthalpy/energy/heat change when 1 <br> mol of gaseous ions (1) | Heat released.... <br> $\mathrm{X}^{+}(\mathrm{g})+\mathrm{aq} \rightarrow \mathrm{X}^{+}(\mathrm{aq})$ and <br> statement of energy <br> change per mole for <br> first mark. | Any implication of <br> endothermic, do not <br> award $1^{\text {st }}$ mark | 2 |
|  | Is dissolved in (a large) excess of water <br> Or <br> Is dissolved until further dilution <br> causes no further heat change (1) | "Added to water" or <br> "reacts with water" <br> instead of "dissolved" <br> "Ignore any reference to "standard <br> conditions" <br> Mark independently | "Ininitely dilute <br> solution" | "Dissolves <br> completely" <br> "Is completely <br> surrounded by water <br> molecules" |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 4. (d)(ii) | $\delta^{\circ} 0$ (in water) attracted to positive ions/cations (1) <br> $\delta^{+} \mathrm{H}$ (in water) attracted to negative ions/anions (1) | 'forms (dative) bonds' instead of 'attracted' <br> Just "attraction between water (molecules) and ions" (1 max) | Reference to full charges on water molecules scores zero overall <br> "energy required" or implication of an endothermic process scores (0) overall. <br> Dipole-dipole attractions and/or "polarisation" scores zero overall | 2 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 5.(a)(i) | One acid: $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}^{2}(\mathrm{aq})$ <br> Conjugate base: $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}(\mathrm{aq)} \mathrm{(1)}$ <br> Other acid: $\mathrm{H}_{3} \mathrm{O}^{+}(\mathrm{aq})$ <br> Conjugate base: $\mathrm{H}_{2} \mathrm{O}(\mathrm{l}) \quad$ (1) <br> Ignore state symbolsAccept correct acids <br> with conjugate bases in <br> either order | 2 |  |  |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 5.(a)(ii) | WEAK: dissociates/ionises to a small <br> extent (1) OWTTE | 'Few molecules <br> dissociate' <br> 'Incomplete' or <br> 'partial' dissociation <br> "Does not fully <br> dissociate" | "ions partially <br> dissociate" | 2 |
|  | ACID: proton donor (1) | Produces $\mathrm{H}_{3} \mathrm{O}^{+} /$ <br> hydrogen / $\mathrm{H}^{+}$ions | Just "contains $\mathrm{H}_{3} \mathrm{O}^{+}$ <br> ..." |  |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :---: | :---: | :---: | :--- |
| 5.(b)(i) | $\mathrm{Ka}=\frac{\left[\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}\right]\left[\mathrm{H}_{3} \mathrm{O}^{+}\right]}{\left[\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}\right]}$ | $\left[\mathrm{H}^{+}\right]$instead of $\left[\mathrm{H}_{3} \mathrm{O}^{+}\right]$ | Any expression <br> containing $\left[\mathrm{H}_{2} \mathrm{O}\right]$ | 1 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 5. (b)(ii) | $\left(\left[\mathrm{H}^{+}\right]=3.63 \times 10^{-4}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)\right.$ (1) Or $10^{-3.44}$ $\left[\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}\right]=\frac{\left[\mathrm{H}^{+}\right]^{2}}{1.30 \times 10^{-5}}$ <br> Or $\begin{equation*} \left[\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}\right]=\frac{\left(3.63 \times 10^{-4}\right)^{2}}{1.30 \times 10^{-5}} \tag{1} \end{equation*}$ <br> ASSUMPTIONS: <br> First assumption mark: <br> negligible $\left[\mathrm{H}^{+}\right]$from ionisation of water $\begin{equation*} \operatorname{Or}\left[\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}\right]=\left[\mathrm{H}^{+}\right] \tag{1} \end{equation*}$ <br> Second assumption mark: Ionisation of the (weak) acid is negligible <br> Or $\mathrm{x}-\left[\mathrm{H}^{+}\right] \approx \mathrm{x}$ where x is initial concentration of $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}$ Or $[\mathrm{H}+] \ll[\mathrm{HA}]$ | If $K_{\mathrm{a}}$ expression incorrect in (b)(i) or [ $\mathrm{H}^{+}$] not squared, only $1^{\text {st }}$ mark available <br> "No other source of $\mathrm{H}^{+}$ ions" <br> "Very slight ionisation ..." <br> "the initial [HA] = equilibrium [HA]" | Just " $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}=$ $\mathrm{H}^{+"}$ (ie no square brackets) <br> Any mention of nonstandard conditions or 'temperature not at $298 \mathrm{~K}^{\prime}$ | 5 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 5.(c) | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}+\mathrm{H}_{2} \mathrm{O} \rightleftharpoons / \rightarrow \mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}$ <br> $+\mathrm{OH}^{-}$ <br> Or <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COONa}+\mathrm{H}_{2} \mathrm{O} \rightleftharpoons / \rightarrow$ <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}+\mathrm{NaOH} \quad(1)$ | $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COO}^{-}+\mathrm{H}^{+} \rightleftharpoons$ <br> $\mathrm{CH}_{3} \mathrm{CH}_{2} \mathrm{COOH}$ <br> and causes the <br> following eqm to shift <br> to the right <br> $\mathrm{H}_{2} \mathrm{O} \rightleftharpoons \mathrm{H}^{+}+\mathrm{OH}^{-}$ |  | 2 |
|  | $\mathrm{OH}^{-}$ions produced cause the solution <br> to be alkaline (1) <br> Mark independently | Causing an excess of <br> $\mathrm{OH}^{-}$ions (1) | "OH- ions from <br> water" |  |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 5.(d)(i) | Ignore "A solution of known pH <br> which...." <br> maintains nearly constant pH <br> OR <br> resists change in pH (1) OWTTE <br> on adding small amounts of acid or <br> alkali (1) <br> Mark independently |  | 2 |  |



| Question <br> Number | Correct Answer | Receptable Answers | Rark |  |
| :--- | :---: | :--- | :--- | :--- |
| 6.(a)(i) |  | Delocalised <br> carboxylate group with <br> a negative charge <br> shown | Compressed <br> structural formula | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| $\mathbf{6 . ( a ) ( i i )}$ | (H ${ }^{+}$from) COOH (group) protonates the <br> $-\mathrm{NH}_{2}$ (group) | Transfer of $\mathrm{H}^{+}$from <br> COOH to $\mathrm{NH}_{2}$ <br> Or <br> "self-protonation" | Just "protonation" <br> Just "acid-base <br> reaction" | 1 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 6.(a)(iii) | Read the whole answer! <br> High energy needed to overcome (strong) ionic attractions (1) <br> between zwitterions (1) <br> Ignore reference to "molecules" if clearly used in the context of attraction between ions | "ionic bonds" or "ionic lattice" instead of "ionic attractions" <br> between adjacent species | J ust "intermolecular forces" <br> Or H bonding <br> Or van der Waals’ <br> forces etc <br> award zero overall | 2 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 6.(b)(i) | $\begin{aligned} & { }^{+} \mathrm{NH}_{3} \mathrm{CH}_{2} \mathrm{COOH} /{ }^{+} \mathrm{H}_{3} \mathrm{NCH}_{2} \mathrm{COOH} / \\ & { }^{+} \mathrm{H}_{3} \mathrm{NCH}_{2} \mathrm{COOH} \end{aligned}$ <br> OR written right to left | $-\mathrm{CO}_{2} \mathrm{H}$ <br> OR <br> $-\mathrm{NH}_{3}{ }^{+} \mathrm{Cl}^{-}$ <br> Or $-\mathrm{NH}_{3} \mathrm{Cl}$ | Molecular formula | 1 |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 6. (b)(ii) | $\mathrm{NH}_{2} \mathrm{CH}_{2} \mathrm{COO}^{-} / \mathrm{NH}_{2} \mathrm{CH}_{2} \mathrm{CO}_{2}^{-} /$  | $-\mathrm{COONa}$ <br> or $-\mathrm{COO}^{-} \mathrm{Na}^{+}$ | Molecular formula | 1 |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :--- | :--- | :--- | :--- |
| 6.(b)(iii) | $\mathrm{CH}_{3} \mathrm{CONHCH}_{2} \mathrm{COOH} /$ | $\mathrm{CH}_{3} \mathrm{CONHCH}_{2} \mathrm{CO}_{2} \mathrm{H}$ | Molecular formula | 1 |
|  | OR |  |  |  |
| 'no reaction' (1) |  |  |  |  |
|  |  |  |  |  |
|  |  |  |  |  |
|  |  |  |  |  |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 6.(b)(iv) |  | $\mathrm{NH}_{2} \mathrm{CH}_{2} \mathrm{CO}_{2} \mathrm{CH}_{3}$ |  | 1 |


| Question <br> Number | Correct Answer | Receptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| 6.(c)(i) | (Glutamic acid molecule) has four <br> different groups attached to a C <br> (atom) <br> Or <br> (Glutamic acid molecule) has four <br> different groups attached to a chiral <br> centre | Contains an <br> asymmetric carbon <br> (atom) <br> Or <br> molecule has no plane <br> of symmetry <br> OR <br> has mirror images which are not <br> superimposable | Just "has a chiral <br> centre" | Or ust "the molecule <br> is asymmetrical" |


| Question <br> Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :--- | :--- | :--- | :--- | :--- |
| $\mathbf{6 . ( c ) ( i i )}$ | (the isomers) rotate the plane (or <br> polarisation) of (plane-) polarised light <br> (1) <br> in opposite directions (1) <br> Ignore any reference to polarimeter | "...rotate plane <br> polarised light" | Just "in different <br> directions" |  |


| Question Number | Correct Answer | Acceptable Answers | Reject | Mark |
| :---: | :---: | :---: | :---: | :---: |
| 6.(d) | $\mathrm{H}_{2} \mathrm{~N}\left(\mathrm{CH}_{2}\right)_{6} \mathrm{NH}_{2}(1)$ $\mathrm{ClOC}\left(\mathrm{CH}_{2}\right)_{4} \mathrm{COCl} /$  <br> [Monomers can be given in either order] |  |  | 2 |

