# MARK SCHEME for the May/June 2010 question paper for the guidance of teachers 

## 9701 CHEMISTRY

9701/42
Paper 4 (A2 Structured Questions), maximum raw mark 100

This mark scheme is published as an aid to teachers and candidates, to indicate the requirements of the examination. It shows the basis on which Examiners were instructed to award marks. It does not indicate the details of the discussions that took place at an Examiners' meeting before marking began, which would have considered the acceptability of alternative answers.

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1
(a) $\mathrm{C}_{6} \mathrm{H}_{5}-\mathrm{COCH}_{2} \mathrm{OH}$ or $\mathrm{C}_{8} \mathrm{H}_{8} \mathrm{O}_{2}$ and NaCl or $\mathrm{Cl}^{-}$
$(1)+(1)$
[2]
(b) (i) the exponent / power to which a concentration is raised in the rate equation (or in an equation, e.g. "a" in the equ: rate $=k[A]^{a}$ )
(ii) from 1 and 2: rate increases by $50 \%$ as does $[R C l]$, so rate $\propto[R C l]^{1}$
from 1 and 3: rate $\propto[\mathrm{NaOH}]^{1}$
(iii) (rate =) k[RCl][OH $\left.{ }^{-}\right]$
(iv)

marking points:

- $\quad(+)$ or ${ }^{\delta+}$ on C and (-) or ${ }^{\delta-}$ on Cl
- lone pair and charge on: $\mathrm{OH}^{-}$
- curly arrow from OH (lone pair) to ${ }^{(\delta+)} \mathrm{C}$, and either a curly arrow breaking $\mathrm{C}-\mathrm{Cl}$ bond or 5 -valent transition state (ignore charge)
- $\mathrm{S}_{\mathrm{N}} 1$ alternative for last mark (only award mark if candidate's rate equation shows first order reaction): curly arrow breaking C-Cl bond and carbocation intermediate.
(c) (i) (add $\mathrm{RCl} / \mathrm{RCOCl}$ to) (aq) $\mathrm{Ag}^{+} / \mathrm{AgNO}_{3}$ or named indicator (e.g. MeOr) or use pH probe

White ppt appears (faster with RCOCl ) or turns acidic colour (e.g. red) or shows pH decrease
if water is the only reagent, and no pH meter used: award only the second mark, for "steamy / white fumes"
(ii) ( $\mathrm{C}=\mathrm{O}$ is polarised $/$ ) carbon is more $\delta+$ than in $\mathrm{R}-\mathrm{Cl}$ or carbon is positive or RCOCl can react via addition-elimination
(mention of electronegativity on its own is not enough for the mark)

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2 (a) less soluble down group
lattice energy and hydration energies both decrease (i.e. become less negative)
but H.E. decreases more (than L.E.) or change in H.E. outweighs L.E.
so $\Delta \mathrm{H}_{\text {sol }}$ becomes more endothermic / less exothermic
(b) (i) for Mg: $\Delta \mathrm{H}=2993-1890-(2 \times 550)=(+) 3\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$
for Sr: $\Delta \mathrm{H}=2467-1414-(2 \times 550)=-47\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right)$
(ii) $\operatorname{Sr}(\mathrm{OH})_{2}$ should be more soluble in water, and $\Delta \mathrm{H}$ is more exothermic / negative

Assuming "other factors" (e.g. $\Delta \mathrm{S}$, or temperature etc.) are the same
(iii) $\operatorname{Sr}(\mathrm{OH})_{2}$ should be less soluble in hot water, because $\Delta \mathrm{H}$ is negative / exothermic
(c) (i) $\mathrm{K}_{\mathrm{sp}}=\left[\mathrm{Ca}^{2+}\right]\left[\mathrm{OH}^{-}\right]^{2}$ (needs the charges) units: $\mathrm{mol}^{3} \mathrm{dm}^{-9}$
(ii) $\mathrm{n}\left(\mathrm{H}^{+}\right)=\mathrm{n}\left(\mathrm{OH}^{-}\right)=0.05 \times 21 / 1000=1.05 \times 10^{-3} \mathrm{~mol}$ in $25 \mathrm{~cm}^{3}$
$\left[\mathrm{OH}^{-}\right]=1.05 \times 1000 / 25=4.2 \times 10^{-2}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$
$\left[\mathrm{Ca}^{2+}\right]=2.1 \times 10^{-2}\left(\mathrm{~mol} \mathrm{dm}^{-3}\right)$
$\mathrm{K}_{\mathrm{sp}}=2.1 \times 10^{-2} \times\left(4.2 \times 10^{-2}\right)^{2}=3.7 \times 10^{-5}$
(iii) less soluble in NaOH due to the common ion effect or equilibrium is shifted to the I.h.s. by high $\left[\mathrm{OH}^{-}\right]$(NOT just a mention of Le Chat' on its own)

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3 (a) $\mathrm{SiF}_{4}$ is symmetrical or tetrahedral or bonds are at $109^{\circ}$ or has no lone pair or 4 electron pairs shared equally or all Si-F dipoles cancel out, or $\mathrm{SF}_{4}$ has a lone pair (on S).
(b)

| compound | molecule has <br> an overall dipole | molecule does not have <br> an overall dipole |
| :---: | :---: | :---: |
| $\mathrm{BCl}_{3}$ |  | $\checkmark$ |
| $\mathrm{PC} l_{3}$ | $\checkmark$ |  |
| $\mathrm{CCl} l_{4}$ |  | $\checkmark$ |
| $\mathrm{SF}_{6}$ |  | $\checkmark$ |

mark row-by-row,
(c) (i) Si and B have empty / available / low-lying orbitals or C does not have available orbitals (allow "B is electron deficient" but not mention or implication of d-orbital on B)
(ii) $\mathrm{BCl}_{3}+3 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{H}_{3} \mathrm{BO}_{3}+3 \mathrm{HCl}$ or $2 \mathrm{BCl}_{3}+3 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{B}_{2} \mathrm{O}_{3}+6 \mathrm{HCl}$
$\mathrm{SiCl}_{4}+2 \mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{SiO}_{2}+4 \mathrm{HCl}$ etc., e.g. $\rightarrow \mathrm{Si}(\mathrm{OH})_{4}, \mathrm{H}_{2} \mathrm{SiO}_{3}$
(d) (i) $\mathrm{Si}_{3} \mathrm{Cl}_{8} \mathrm{O}_{2}$ (this has $\mathrm{M}_{\mathrm{r}}=84+280+32=396$ ) or $\mathrm{Si4Cl}_{4} \mathrm{O}_{9}$ or $\mathrm{Si}_{8} \mathrm{Cl}_{4} \mathrm{O}_{2}$
(ii)

| mass number | structure |
| :---: | :---: |
| 133 | $\mathrm{Cl}_{3} \mathrm{Si}$ |
| 247 | $\mathrm{Cl}_{3} \mathrm{Si}_{\mathrm{Si}}-\mathrm{SiCl} l_{2}$ |
| 263 | $\mathrm{Cl}_{3} \mathrm{Si}-\mathrm{O}-\mathrm{SiCl}_{2}-\mathrm{O}$ |

(if correct structures are not given for last 2 rows, you can award (1) mark for two correct molecular formulae:
either $\mathrm{Si}_{2} \mathrm{Cl}_{5} \mathrm{O}+\mathrm{Si}_{2} \mathrm{Cl}_{5} \mathrm{O}_{2}$ or $\mathrm{Si}_{3} \mathrm{ClO}_{8}+\mathrm{Si}_{3} \mathrm{ClO}_{9}$ or $\mathrm{Si}_{7} \mathrm{ClO}+\mathrm{Si}_{7} \mathrm{ClO}_{2}$ )
(iii)

allow ecf on the structure drawn in the third row of the table in (ii)
but any credited structure must show correct valencies for $\mathrm{Si}, \mathrm{Cl}$ and O .

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4 (a) $\mathrm{Cr}^{3+}$ : $1 \mathrm{~s}^{2} 2 \mathrm{~s}^{2} 2 \mathrm{p}^{6} \ldots 3 \mathrm{~s}^{2} 3 p^{6} 3 d^{3}$
$M n^{2+}: \quad 1 s^{2} 2 s^{2} 2 p^{6} \ldots 3 s^{2} 3 p^{6} 3 d^{5}$
(allow (1) out of (2) for $3 s^{2} 3 p^{6} 4 s^{2} 3 d^{1}$ and $3 s^{2} 3 p^{6} 4 s^{2} 3 d^{3}$ )
(b) (i) any three of the following points:

- initial (pale) green (solution)
- fades to (almost) colourless (allow yellow)
- then (permanent faint) pink
- finally (deep) purple
(ii) $\mathrm{MnO}_{4}^{-}+8 \mathrm{H}^{+}+5 \mathrm{Fe}^{2+}\left(+5 \mathrm{e}^{-}\right) \rightarrow \mathrm{Mn}^{2+}+4 \mathrm{H}_{2} \mathrm{O}+5 \mathrm{Fe}^{3+}\left(+5 \mathrm{e}^{-}\right)$
(c) $E^{\ominus}$ values:

$$
\begin{equation*}
\mathrm{O}_{2}+2 \mathrm{H}_{2} \mathrm{O} / 4 \mathrm{OH}^{-}=+0.40 \mathrm{~V} \quad \mathrm{Fe}(\mathrm{OH})_{3} / \mathrm{Fe}(\mathrm{OH})_{2}=-0.56 \mathrm{~V} \tag{2}
\end{equation*}
$$

$E^{\ominus}{ }_{\text {cell }}=+0.46 \mathrm{~V}$ (allow -0.37 ) in acid, but +0.96 V in alkali or $\mathrm{E}^{\ominus}\left(\mathrm{OH}^{-}\right)>\mathrm{E}^{\ominus}\left(\mathrm{H}^{+}\right)$
If $\mathbf{E}_{\text {cell }}$ is more positive it means a greater likelihood of reaction
(d)

(1) and
$\mathrm{CH}_{3} \mathrm{CO}_{2} \mathrm{H}$


(1)
(1)
(e) (i) $\left(\mathrm{CH}_{3}\right)_{2} \mathrm{C}(\mathrm{OH})-\mathrm{CH}_{2} \mathrm{OH}$
(ii) reaction I: (cold dilute) $\mathrm{KMnO}_{4}$ ("cold" not needed, but "hot" or "warm" negates) reaction II: $\mathrm{Cr}_{2} \mathrm{O}_{7}{ }^{2-}+\mathrm{H}^{+}+$distil
[Total: 18 max 17]

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5 (a) (i) because the carbons are $\mathrm{sp}^{2} /$ trigonal planar / bonded at $120^{\circ}$ or are joined by $\pi$ bonds / orbitals
(ii) because the $\pi$ electrons / double bonds are delocalised / in resonance or electrons are evenly distributed / spread out
(b) (i) $\mathrm{HNO}_{3}+2 \mathrm{H}_{2} \mathrm{SO}_{4} \rightarrow \mathrm{NO}_{2}^{+}+\mathrm{H}_{3} \mathrm{O}^{+}+2 \mathrm{HSO}_{4}^{-}$
or $\mathrm{HNO}_{3}+\mathrm{H}_{2} \mathrm{SO}_{4} \rightarrow \mathrm{H}_{2} \mathrm{NO}_{3}^{+}+\mathrm{HSO}_{4}^{-}$or $\rightarrow \mathrm{H}_{2} \mathrm{O}+\mathrm{NO}_{2}^{+}+\mathrm{HSO}_{4}^{-}$
(ii) electrophilic substitution mechanism:

curly arrows from benzene to $\mathrm{NO}_{2}{ }^{+}$, and showing loss of $\mathrm{H}^{+}$ correct intermediate (with " + " in the 'horse-shoe')
(c) $\mathrm{Cl}_{2}+\mathrm{AlCl}_{3} / \mathrm{FeCl}_{3} / \mathrm{Fe} / \mathrm{Al} / \mathrm{I}_{2}$ (aq or light negates this mark)
(d) (i) $\mathbf{Y}$ is chlorobenzene (1) $\mathbf{Z}$ is 4-chloronitrobenzene (1)
(ii) $\mathrm{Sn} / \mathrm{Fe}+($ conc) HCl

HCl is conc, and second step is to add $\mathrm{NaOH}(\mathrm{aq})$
(iii)

ignore orientation


| no reaction |
| :---: |
| $\mathbf{C}$ |

allow $\mathrm{NHOCCH}_{3}$, but not $\mathrm{NHCH}_{3} \mathrm{CO}$ or $\mathrm{NHCH}_{3} \mathrm{OC}$


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6 (a) (i) Primary - the amino acid sequence / order / chain or diag. e.g. NH-C-CO-NH-C-CO
or amino acids bonded by covalent / amide / peptide bonds
(ii) Tertiary - the coiling / folding of the protein / polypeptide chain due to interactions between side-chains on the amino acids or the structure which gives the protein its 3-D / globular shape
(b) (i) Diagram:

Minimum is $\mathrm{CH}_{2} \mathrm{~S}-\mathrm{SCH}_{2}$
(ii) Oxidation / dehydrogenation / redox
(iii) Hydrogen / H bonds; ionic interactions / bonds or ion-dipole or salt bridges; van der Waals' or id-id or induced / instantaneous dipole forces (ignore hydrophobic interactions)
(2)
(c) (i) Hydrogen bonds
(ii) Correct new strand present (see below) Diagram showing $\mathrm{C}=\mathrm{O}$ bonding to $\mathrm{N}-\mathrm{H}$ in new strand... ...and $\mathrm{N}-\mathrm{H}$ bonding to $\mathrm{C}=\mathrm{O}$ in new strand e.g.


New strand must contain a minimum of two amino acid residues in a single chain. Deduct a penalty of -(1) for any wrong H-bond only if (2) marks have already been scored.
(d) There are bonds or S-S bridges / linkages between the layers / sheets (in $\beta$-keratin) (but only van der Waals interactions between the layers in silk)

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7 (a) The amino acid is uncharged / neutral / a zwitterion or charges balance / are equal (NOT "is non-polar")

It is equally attracted by the anode / + and the cathode / - or attracted by neither
The pH of the buffer is at the isoelectric point/IEP of the amino acid any two $\checkmark \checkmark$
(2) [2]
(b) (at pH 10), $\mathrm{H}_{2} \mathrm{NCH}_{2} \mathrm{CO}_{2}^{-}$or $\mathrm{NH}_{2} \mathrm{CH}_{2} \mathrm{COO}^{-}$
(1) [1]
(c)

| amino acid | relative size | charge |
| :---: | :--- | :---: |
| A | small(est) (1) | -ve |
| B | large(st) (3) | - ve |
| C | middle (2) | +ve |

(numbers are OK to show relative sizes)
Mark each row
(d) (i) lys - val - ser - ala - gly - ala - gly - asp
(ii) gly - ala - gly
(iii) aspartic acid (or lysine)
[Total: 10]

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8 (a) Reaction II - since electrons are used up / required / gained / received (from external circuit)
(b) $\left(\mathrm{Pb}^{2+}+2 \mathrm{e}^{-} \rightarrow \mathrm{Pb}\right)$

$$
\left(\mathrm{PbO}_{2}+4 \mathrm{H}^{+}+2 \mathrm{e}^{-} \rightarrow \mathrm{Pb}^{2+}+2 \mathrm{H}_{2} \mathrm{O}\right)
$$

$$
\begin{align*}
& \mathrm{E}^{\ominus}=-0.13 \mathrm{~V} \\
& \mathrm{E}^{\ominus}=+1.47 \mathrm{~V} \\
& \text { two correct } E^{\ominus} \text { values } \tag{1}
\end{align*}
$$

Cell voltage is $1.6(0)(\mathrm{V})$
(c) (i) $3(+)$
(ii) They are less heavy / poisonous / toxic / polluting or are safer due to no (conc) $\mathrm{H}_{2} \mathrm{SO}_{4}$ within them
(d) (i) Platinum or graphite / carbon
(ii) They need large quantities of compressed gases which take up space or the hydrogen would need to be liquefied or the reactant is (highly) flammable / explosive / combustible
(e) Glass: saves energy - the raw materials are easily accessible / cheap or making glass is energy-intensive

Steel: saves energy - extracting iron from the ore
or mining the ore is energy intensive
or saves a resource - iron ore (NOT just "iron") is becoming scarce either one (1)

Plastics: saves a valuable / scarce resource: (crude) oil / petroleum

