# General Certificate of Education 

## Chemistry 1421

## CHEM1 Foundation Chemistry

Mark Scheme

2010 examination - January series

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[^0]| Question | Part | $\begin{aligned} & \hline \text { Sub } \\ & \text { Part } \end{aligned}$ | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) |  | $2 s^{2} 2 p^{6} 3 s^{1}$ | 1 | $1 \mathrm{~s}^{2}$ can be rewritten <br> Allow $2 s^{2} 2 p_{x}{ }^{2} 2 p_{y}{ }^{2} 2 p_{z}^{2} 3 s^{1}$ <br> Allow subscripts and capitals |
| 1 | (b) | (i) | Energy/enthalpy (needed) to remove one mole of electrons from one mole of atoms/compounds/molecules/elements <br> OR <br> Energy to form one mole of positive ions from one mole of atoms <br> OR <br> Energy/enthalpy to remove one electron from one atom <br> In the gaseous state (to form 1 mol of gaseous ions) |  | Energy given out loses M1 <br> M 2 is dependent on a reasonable attempt at M1 <br> Energy needed for this change $X(\mathrm{~g}) \rightarrow \mathrm{X}^{+}(\mathrm{g})+\mathrm{e}^{(-)}=2$ marks This equation alone scores one mark |
| 1 | (b) | (ii) | $\begin{aligned} & \mathrm{Mg}^{+}(\mathrm{g}) \rightarrow \mathrm{Mg}^{2+}(\mathrm{g})+\mathrm{e}^{(-)} \\ & \mathrm{Mg}^{+}(\mathrm{g})+\mathrm{e}^{(-)} \rightarrow \mathrm{Mg}^{2+}(\mathrm{g})+2 \mathrm{e}^{(-)} \\ & \mathrm{Mg}^{+}(\mathrm{g})-\mathrm{e}^{(-)} \rightarrow \mathrm{Mg}^{2+}(\mathrm{g}) \\ & \hline \end{aligned}$ | 1 | Do not penalise MG Not equation with X |
| 1 | (b) | (iii) | Electron being removed from a positive ion (therefore need more energy)/ electron being removed is closer to the nucleus $/ \mathrm{Mg}^{+}$ smaller (than Mg )/ $\mathrm{Mg}^{+}$more positive than Mg | 1 | Allow from a + particle/ species Not electron from a higher energy level/or higher sub-level More protons $=0$ |
| 1 | (b) | (iv) | Range from 5000 to $9000 \mathrm{~kJ} \mathrm{~mol}^{-1}$ | 1 |  |
| 1 | (c) |  | Increase <br> Bigger nuclear charge (from Na to Cl )/more protons <br> electron (taken) from same (sub)shell/ similar or same shielding/ electron closer to the nucleus/smaller atomic radius | $\begin{array}{\|l\|} \hline 1 \\ 1 \\ 1 \end{array}$ | If decrease CE = 0/3 <br> If blank mark on QWC <br> If no shielding $=0$ <br> Smaller ionic radius $=0$ |

\(\left.$$
\begin{array}{|l|l|l|l|l|l|}\hline 1 & \text { (d) } & & \text { Lower } & 1 & \begin{array}{l}\text { If not lower CE }=0 / 3 \\
\text { If blank mark on }\end{array}
$$ <br>
Allow does not increase <br>

Not 2 p\end{array}\right]\) M3 dependent upon a reasonable | Two/pair of electrons in (3)p orbital or implied |
| :--- |
| repel (each other) |


| Question | Part | Sub Part | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 2 | (a) | (i) | $\begin{aligned} & M_{\mathrm{r}}=132.1 \\ & 0.0238 \end{aligned}$ | $1$ | 132 <br> Allow 0.024 <br> Allow 0.0237 <br> Penalise less than 2 sig fig once in <br> (a) |
| 2 | (a) | (ii) | 0.0476 | 1 | $\begin{aligned} & 0.0474-0.0476 \\ & \text { Allow (a) (i) } \times 2 \\ & \hline \end{aligned}$ |
| 2 | (a) | (iii) | 1.21 | 1 | Allow consequential from (a) (ii) ie allow (a) (ii) $\times 1000$ / 39.30 Ignore units even if wrong |
| 2 | (b) |  | $\frac{34 \times 100}{212.1}$ $=16.0(3) \%$ |  | Allow mass or Mr of desired product times one hundred divided by total mass or Mr of reactants/products <br> If $34 / 212.1$ seen correctly award M1 Allow 16\% <br> 16 scores 2 marks |
| 2 | (c) |  | 100(\%) | 1 | Ignore all working |
| 2 | (d) |  | $\begin{aligned} \mathrm{PV} & =\mathrm{nRT} \text { or } \mathrm{n}=\frac{\mathrm{PV}}{\mathrm{RT}} \\ \mathrm{n} & =\frac{100000 \times 1.53 \times 10^{-2}}{8.31 \times 310} \\ & =0.59(4) \end{aligned}$ | $1$ <br> 1 | If rearranged incorrectly lose M1 and M3 <br> M2 for mark for converting P and T into correct units in any expression <br> Allow 0.593 <br> M3 consequential on transcription error only not on incorrect $P$ and $T$ |


| 2 | (e) | $\left(\mathrm{Na}_{2} \mathrm{SO}_{4}\right)$ | $\mathrm{H}_{2} \mathrm{O}$ |  |  |
| :---: | :---: | :---: | :---: | :---: | :---: |
|  |  | (44.1\%) | 55.9\% | 1 | M 1 is for 55.9 |
|  |  | $\begin{aligned} & 44.1 / 142.1 \\ & 0.310 \\ & =1 \end{aligned}$ | $\begin{aligned} & 55.9 / 18 \\ & 3.11 \\ & =10 \end{aligned}$ | 1 | Alternative method gives180 for water part =2 marks |
|  |  | $x=10$ |  | 1 | $\begin{aligned} & X=10=3 \text { marks } \\ & 10.02=2 \text { marks } \end{aligned}$ |


| Question | Part | Sub <br> Part | Marking Guidance | Mark | Comments |
| :--- | :--- | :--- | :--- | :--- | :--- |
| 3 | (a) |  | Hydrogen/H bonds <br> van der Waals/vdw/ dipole-dipole/London/temporarily induced <br> dipole/dispersion forces | 1 | 1 |



| 3 | (e) | Dative (covalent)/ coordinate <br> (Lone) pair/both electrons/two electrons on $\mathrm{O}\left(\mathrm{H}_{2}\right)$ donated (to $\mathrm{H}^{+}$) OR pair/both electrons come from $\mathrm{O}\left(\mathrm{H}_{2}\right)$ | 1 1 | If not dative/coordinate CE = 0/2 If covalent or blank read on <br> Explanation of a coordinate bond specific to oxygen or water required Not just $\mathrm{H}+$ attracted to lone pair since that is nearer to a H bond |
| :---: | :---: | :---: | :---: | :---: |
| 3 | (f) | ionic <br> oppositely charged ions /+ and - ions or particles <br> ions attract strongly OR strong/many (ionic) bonds must be broken | 1 | ```if not ionic CE = 0 atoms or molecules loses M2 and M3 S loses M2 Reference to IMF loses M2 and M3``` |


| Question | Part | Sub <br> Part | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (a) | (i) | single (C-C) bonds only/ no double ( $\mathrm{C}=\mathrm{C}$ ) bonds C and H (atoms) only/purely/solely/entirely | 1 1 | Allow all carbon atoms bonded to four other atoms <br> Single C-H bonds only $=0$ <br> $\mathrm{C}=\mathrm{H} C E$ <br> Not consists or comprises <br> Not completely filled with hydrogen <br> CH molecules $=\mathrm{CE}$ <br> Element containing C and $\mathrm{H}=\mathrm{CE}$ |
| 4 | (a) | (ii) | $\mathrm{C}_{\mathrm{n}} \mathrm{H}_{2 \mathrm{n}+2}$ | 1 | $\begin{aligned} & \text { Formula only } \\ & \mathrm{C}_{\mathrm{x}} \mathrm{H}_{2 x+2} \end{aligned}$ |
| 4 | (b) | (i) | $\mathrm{C}_{5} \mathrm{H}_{12}+8 \mathrm{O}_{2} \rightarrow 5 \mathrm{CO}_{2}+6 \mathrm{H}_{2} \mathrm{O}$ | 1 | Accept multiples Ignore state symbols |
| 4 | (b) | (ii) | gases produced are greenhouse gases/contribute to Global warming/effect of global warming/climate change | 1 | Allow $\mathrm{CO}_{2}$ or water is greenhouse gas/causes global warming Acid rain/ozone $\mathrm{CE}=0$ |
| 4 | (c) |  | carbon | 1 | Allow C Allow soot |
| 4 | (d) | (i) | $\mathrm{C}_{9} \mathrm{H}_{20} \rightarrow \mathrm{C}_{5} \mathrm{H}_{12}+\mathrm{C}_{4} \mathrm{H}_{8}$ <br> OR $\mathrm{C}_{9} \mathrm{H}_{20} \rightarrow \mathrm{C}_{5} \mathrm{H}_{12}+2 \mathrm{C}_{2} \mathrm{H}_{4}$ | 1 | Accept multiples |
| 4 | (d) | (ii) | Plastics, polymers | 1 | Accept any polyalkene / haloalkanes / alcohols |
| 4 | (d) | (iii) | so the bonds break OR because the bonds are strong | 1 | IMF mentioned = 0 |
| 4 | (e) | (i) | 1,4-dibromo-1-chloropentane / 1-chloro-1,4-dibromopentane | 1 | Ignore punctuation |
| 4 | (e) | (ii) | Chain/position/positional | 1 | Not structural or branched alone |


| Question | Part | $\begin{aligned} & \text { Sub } \\ & \text { Part } \end{aligned}$ | Marking Guidance | Mark | Comments |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 5 | (a) |  | Average/mean mass of (1) atom(s) (of an element) <br> $1 / 12$ mass of one atom of ${ }^{12} \mathrm{C}$ <br> OR <br> (Average) mass of one mole of atoms <br> $1 / 12$ mass of one mole of ${ }^{12} \mathrm{C}$ <br> OR <br> (Weighted) average mass of all the isotopes <br> $1 / 12$ mass of one atom of ${ }^{12} \mathrm{C}$ <br> OR <br> Average mass of an atom/isotope compared to C-12 on a scale in which an atom of C-12 has a mass of 12 | $\begin{aligned} & 1 \\ & 1 \end{aligned}$ | If moles and atoms mixes Max = 1 <br> This expression = 2 marks |
| 5 | (b) |  | d block <br> [Ar] $3 d^{2} 4 s^{2}$ <br> 27 | 1 1 1 | Allow 3d/D <br> Other numbers lose M1 <br> Ignore transition metals <br> Can be written in full <br> Allow subscripts <br> $3 d^{2}$ and $4 s^{2}$ can be in either order |



\begin{tabular}{|c|c|c|c|c|c|}
\hline Question \& Part \& \[
\begin{aligned}
\& \hline \text { Sub } \\
\& \text { Part } \\
\& \hline
\end{aligned}
\] \& Marking Guidance \& Mark \& Comments \\
\hline \multirow[t]{7}{*}{6} \& \multirow[t]{7}{*}{} \& \multirow[t]{7}{*}{} \& \multirow[t]{7}{*}{\begin{tabular}{l}
 \\
trigonal / triangular bipyramid(al) \\
Bent / V shape / non-linear / triangular / angular
\[
104^{\circ}-106^{\circ}
\] \\
(For candidates who thought this was \(\mathrm{CIF}_{2}{ }^{+}\)which contained iodine allow \\
Trigonal / triangular planar \\
\(120^{\circ}\)
\end{tabular}} \& \multirow[t]{2}{*}{1

1} \& Mark M1 - M5 independently M1 for 5 bond pairs around As Do not penalise A for As or FI for F <br>
\hline \& \& \& \& \& Allow trigonal dipyramid <br>

\hline \& \& \& \& 1 \& | M3 for 2 bond pairs to $F$ and 2 lone pairs |
| :--- |
| Lone pairs can be shown as lobes with or without electrons or as $x x$ or $\qquad$ X |
| X | <br>

\hline \& \& \& \& 1
1 \& Bent-linear = contradiction Do not allow trigonal <br>
\hline \& \& \& \& \& <br>
\hline \& \& \& \& \& Not just triangular <br>
\hline \& \& \& \& \& <br>
\hline
\end{tabular}


[^0]:    Set and published by the Assessment and Qualifications Alliance.

