## CAMBRIDGE INTERNATIONAL EXAMINATIONS

## MARK SCHEME for the October／November 2015 series

## 9701 CHEMISTRY

9701／22
Paper 2 （AS Structured Questions），maximum raw mark 60

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) | name of isotope | type of particle | charge | symbol | electron configuration | [5] | [5] |
|  | carbon-13 | atom | 0 | ${ }_{6}^{13} \mathrm{C}$ | $1 s^{2} 2 s^{2} 2 p^{2}$ |  |  |
|  | chloride(-37) | anion | $1-$ | Cl | $1 s^{2} 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{6}$ |  |  |
|  | sulfur-34 | atom | 0 | ${ }_{16}^{34} \mathrm{~S}$ | $1 s^{2} 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{4}$ |  |  |
|  | iron-54 | cation | 2+ | ${ }_{26}^{54} \mathrm{Fe}^{(2+)}$ | $1 s^{2} 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{6} 3 d^{6}$ |  |  |
| (b) (i) | ability/tendency/power of an atom/nucleus to attract/pull electron(s) in a covalent bond/shared pair of electrons/bonding pair of electrons |  |  |  |  | [1] <br> [1] | [2] |
| (ii) | Covalent overlap of orbitals OR shared pair(s) (of electrons) <br> OR <br> metallic positive ions/cations surrounded by delocalised electrons |  |  |  |  | ${ }^{[11}$ <br> [1] <br> [1] <br> [1] | [2] |
| (iii) | Ionic/electrovalent (electrostatic) Attraction between oppositely charged/+ve and -ve ions |  |  |  |  | $\begin{aligned} & {[1]} \\ & {[1]} \end{aligned}$ | [2] |
| (c) (i) | similar strength/amount/number of intermolecular forces/induced dipole/van der Waals'/VdW/London forces/LDF/dispersion forces <br> therefore similar energy needed |  |  |  |  | [1] <br> [1] | [2] |


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| (ii) | M1 HCl polar/has a dipole AND $F_{2}$ non-polar/has no dipole OR (permanent) dipole (-dipole) attractions/forces between HCl (molecules) AND induced dipole (-induced dipole) attractions/forces/LDFs between $\mathrm{F}_{2}$ (molecules) <br> M2 more energy needed for HCl than $\mathrm{F}_{2}$ <br> OR <br> pd-pd forces stronger than id-id forces <br> OR <br> IMFs/VdWs in HCl stronger than in $\mathrm{F}_{2}$ | [1] <br> [1] | [2] |
| (iii) | Hydrogen bonding (between methanol molecules) <br> Stronger than IMFs/van der Waals' in other three/is the strongest intermolecular force | [1] <br> [1] | [2] |
|  |  |  | [17] |
| 2 (a) | M1 Heat (energy) change (or $\mathrm{H}_{\text {prod }}-\mathrm{H}_{\text {react }}$ ) measured at constant pressure OR enthalpy change when the amount/moles of reactants as shown in a (reaction) equation react together to give products <br> M2 measured at standard conditions | [1] <br> [1] | [2] |
| (b) (i) | $\mathrm{q}=2125.53$ | [1] | [1] |
| (ii) | amount $=0.025(0)$ | [1] | [1] |
| (iii) | -85.(0) | [1] | [1] |


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| (iv) |  | [1] | [1] |
| (v) | $\begin{aligned} & \Delta H+9.6=-85.0 \\ & \Delta H=-85.0-9.6=-94.6\left(\mathrm{~kJ} \mathrm{~mol}^{-1}\right) \end{aligned}$ | [1] | [1] |
|  |  |  | [7] |
| 3 (a) (i) | $\begin{aligned} & \mathrm{Na}_{2} \mathrm{O} \text { or } \mathrm{Na}_{2} \mathrm{O}_{2} ; \mathrm{MgO} ; \\ & \mathrm{P}_{4} \mathrm{O}_{10} \text { or } \mathrm{P}_{4} \mathrm{O}_{6} ; \mathrm{SO}_{2} \end{aligned}$ | $\begin{aligned} & {[1]} \\ & {[1]} \end{aligned}$ | [2] |
| (ii) | Na: Yellow/orange/gold flame/white solid/powder/smoke $4 \mathrm{Na}+\mathrm{O}_{2} \rightarrow 2 \mathrm{Na}_{2} \mathrm{O} \text { or } 2 \mathrm{Na}+\mathrm{O}_{2} \rightarrow \mathrm{Na}_{2} \mathrm{O}_{2}$ <br> S: Blue flame/(yellow) solid melts/turns red/amber/white fumes $\mathrm{S}+\mathrm{O}_{2} \rightarrow \mathrm{SO}_{2}$ | $\begin{aligned} & {[1]} \\ & {[1]} \\ & {[1]} \\ & {[1]} \end{aligned}$ | [4] |
| (b) (i) | acidic P and S <br> amphoteric Al and basic Na and Mg | $\begin{gathered} {[1]} \\ {[1]} \end{gathered}$ | [2] |
| (ii) | acidic: covalent (bonding) <br> basic: ionic (bonding) | [1] <br> [1] | [2] |


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| (iii) | $\begin{aligned} & \mathrm{Al}_{2} \mathrm{O}_{3}+6 \mathrm{HCl} \rightarrow 2 \mathrm{AlCl}_{3}+3 \mathrm{H}_{2} \mathrm{O} \mathrm{OR} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+6 \mathrm{H}^{+} \rightarrow 22 \mathrm{Al}^{3+}+3 \mathrm{H}_{2} \mathrm{O} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{NaOH}+7 \mathrm{H}_{2} \mathrm{O} \rightarrow 2 \mathrm{NaAl}(\mathrm{OH})_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2} \mathrm{OR} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{NaOH}+3 \mathrm{H}_{2} \mathrm{O} \rightarrow 2 \mathrm{NaAl}(\mathrm{OH})_{4} \mathrm{OR} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{NaOH}_{2} \mathrm{NaAlO}_{2}+\mathrm{H}_{2} \mathrm{OOR} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{OH}^{-}+7 \mathrm{H}_{2} \mathrm{O} \rightarrow 2\left[\mathrm{Al}(\mathrm{OH})_{4}\left(\mathrm{H}_{2} \mathrm{O}\right)_{2}\right]^{-} \mathrm{OR} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{OH}^{-}+3 \mathrm{H}_{2} \mathrm{O} \rightarrow 2\left[\mathrm{Al}(\mathrm{OH})_{4}\right]^{\mathrm{OR}} \\ & \mathrm{Al}_{2} \mathrm{O}_{3}+2 \mathrm{OH}^{-} \rightarrow 2 \mathrm{AlO}_{2}^{-}+\mathrm{H}_{2} \mathrm{O} \end{aligned}$ | [1] <br> [1] | [2] |
| (c) | ```sulfur forms \(\mathrm{SO}_{2} / \mathrm{SO}_{2}+/\) mixes \(\mathrm{H}_{2} \mathrm{O} \rightarrow \mathrm{H}_{2} \mathrm{SO}_{3}\) or in words OR \(\mathrm{SO}_{2}+/\) mixes \(\mathrm{H}_{2} \mathrm{O}(\rightarrow\) acid \() /\) or in words OR \(\mathrm{SO}_{2}+/\) mixes \(\mathrm{H}_{2} \mathrm{O}+\left(1 / 2 \mathrm{O}_{2}\right) \rightarrow \mathrm{H}_{2} \mathrm{SO}_{4} /\) or in words``` | $\begin{gathered} {[1]} \\ {[1]} \end{gathered}$ | [2] |
|  |  |  | [14] |
| $4 \quad$ (a) (i) | Nucleophilic Substitution | [1] | [1] |
| (ii) | Has a chiral centre/carbon OR has a carbon/C attached to 4 different groups/atoms/chains OR <br> has no plane/line of symmetry | [1] | [1] |
| (iii) |  | [1+1] | [2] |
| (iv) | Elimination | [1] | [1] |


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| (v) |   <br> cis-but-2-ene <br> trans-but-2-ene |  | [1] <br> [1] | [2] |
| (vi) | But-1-ene <br> 2 Hs on one of the double-bonded Cs OR does not have 2 different groups on both atoms/each atom in $\mathrm{C}=\mathrm{C}$ |  | [1] <br> [1] | [2] |
| (b) (i) | ammonia/ $\mathrm{NH}_{3}$ |  | [1] | [1] |
| (ii) | propanoyl chloride $/ \mathrm{C}_{2} \mathrm{H}_{5} \mathrm{COCl}$ |  | [1] | [1] |
| (iii) | $\mathrm{CH}_{3} \mathrm{CH}\left(\mathrm{NHCOC}_{2} \mathrm{H}_{5}\right) \mathrm{CH}_{3}$ |  | [1] | [1] |
| (iv) | Reduction <br> $\mathrm{LiAlH}_{4}$ / lithium aluminium hydride / lithium tetrahydridoaluminate | (1) | $\begin{gathered} {[1]} \\ {[1]} \end{gathered}$ | [2] |
| (v) | aluminium oxide |  | [1] | [1] |


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| (vi) | M 1 = correct structure of Y and curly arrow from double bond to H <br> M 2 = dipole and curly arrow from $\mathrm{H}-\mathrm{Br}$ bond to Br <br> M3 = correct intermediate <br> $\mathrm{M} 4=\mathrm{Br}^{-}$with lone pair and curly arrow from lone pair to $\mathrm{C}(+)$ | $\begin{gathered} {[1]} \\ {[1]} \\ {[1]} \\ {[1]} \end{gathered}$ | [4] |
| (vii) | electrophilic addition | [1] | [1] |
| (viii) | secondary carbocation more stable than primary due to electron releasing character/(positive) inductive effect of alkyl groups | $\begin{aligned} & {[1]} \\ & {[1]} \end{aligned}$ | [2] |
|  |  |  | [22] |

