Oxford Cambridge and RSA

## GCE

## Chemistry A

Unit F321: Atoms, Bonds and Groups
Advanced Subsidiary GCE
Mark Scheme for June 2016

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All examiners are instructed that alternative correct answers and unexpected approaches in candidates' scripts must be given marks that fairly reflect the relevant knowledge and skills demonstrated.

Mark schemes should be read in conjunction with the published question papers and the report on the examination.

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Annotations

| Annotation | Meaning |
| :---: | :--- |
| BOD | Benefit of doubt given |
| CON | Contradiction |
| ECF | Incorrect response |
| I | Error carried forward |
| NAQ | Ignore |
| NBOD | Not answered question |
| POT | Benefit of doubt not given |
| A | Power of 10 error |
| RE | Omission mark |
| $\mathbf{S F}$ | Rounding error |
| $\boldsymbol{N}$ | Error in number of significant figures |
|  | Correct response |

Abbreviations

| Annotation | Meaning |
| :--- | :--- |
| DO NOT ALLOW | Answers which are not worthy of credit |
| IGNORE | Statements which are irrelevant |
| ALLOW | Answers that can be accepted |
| ( ) | Words which are not essential to gain credit |
| - | Underlined words must be present in answer to score a mark |
| ECF | Error carried forward |
| AW | Alternative wording |
| ORA | Or reverse argument |

The following questions should be annotated with ticks $\qquad$ , crosses $x$etc to show where marks have been awarded in the body of the text ignore $2 b i$

| Question |  |  | Answer |  |  |  | Mark | For relative masses GuidanceALLOW $1 / 1800$ to $1 / 2000$ for electron value(0.0005-0.00056)ALLOW 'negligible' for electron valueIGNORE ' + ' in front of correct valuesDO NOT ALLOW '' in fron of $1 / 2000$DO NOT ALLOW 'nil' OR 'zero' for mass of electronFor relative chargesALLOW 1+ and 'neutral' and 1-IGNORE ' '' (ie a dash) for neutronDO NOT ALLOW ''+ 'r ' '' without ' 1 'DO NOT ALLOW '1' without chargeFor position within the atomIGNORE 'middle OR 'centre' for 'nucleus' |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
| 1 | (a) |  | particle <br> proton <br> neutron <br> electron <br> Relative mass <br> Relative charg | relative mass <br> unn $\checkmark$ <br> ND positio | relative charge <br> $+1$ <br> nil/0 <br> - 1 <br> olumns | position <br> within the <br> atom <br> nucleus <br> nucleus <br> shell | 2 |  |
| 1 | (b) | (i) | $\begin{aligned} & \text { s-orbital = sph } \\ & \text { AND } \\ & \text { p-orbital = dum } \end{aligned}$ | al <br> ell shape |  |  | 1 | For s-orbital <br> IGNORE 'circular' <br> For p-orbital <br> ALLOW other words indicating 3-D shape of p-orbital eg 'Peanut-shaped' OR hour glass etc ALLOW 'figure of eight' OR 'figure of 8' IGNORE diagrams |
| 1 |  | (ii) | p-orbitals have (three) p-orbit | ater ener <br> have equa | han s-orb <br> ergy |  | 2 | ALLOW reverse argument <br> ALLOW suitable energy diagram for either part |


|  | uest | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 1 | (c) | ${ }_{\mathrm{x}}^{\mathrm{x}} \mathrm{~N} \underset{\stackrel{i}{\dot{x}}}{\stackrel{\dot{x}}{\dot{x}} \mathrm{x}} \mathrm{~N}$ | 1 | ALLOW all dots or all crosses. |
| 1 | (d) | First check the answer line. <br> If answer $=1.7(0) \times 10^{-3}$ award 2 marks. <br> M1 (Dividing by $6.02 \times 10^{23}$ ) <br> Number of $\mathrm{N}_{2}$ molecules $=\frac{5.117 \times 10^{20}}{6.02 \times 10^{23}}=8.5 . \times 10^{-4}$ <br> OR $0.85 \times 10^{-3}$ OR $0.085 \times 10^{-2}$ OR $0.0085 \times 10^{-1}$ OR 0.00085 <br> M2 (Correct conversion of molecules to atoms + standard form) <br> M1 $\times 2$ and in standard form $\checkmark$ <br> From 0.0085, answer $=2 \times 0.00085=0.00170$ $=1.7(0) \times 10^{-3}$ <br> Alternative method <br> M1 (Correct conversion of molecules to atoms) $=5.117 \times 10^{20} \times 2=1.02(34) \times 10^{21}$ <br> OR $10.2(34) \times 10^{20}$ OR 102.(34) $\times 10^{19}$ etc $\begin{aligned} & \text { M2 (Correct use of } 6.02 \times 10^{23}+\text { standard form) } \\ & \frac{1.02(34) \times 10^{21}}{6.02 \times 10^{23}}=1.7(0) \times 10^{-3} \end{aligned}$ | 2 | ALLOW one mark for <br> $0.17 \times 10^{-2}$ OR $0.017 \times 10^{-1}$ OR 0.0017 (not standard form) <br> ALLOW one mark for <br> $4.25 \times 10^{-4}$ (dividing by 2 in M2 + standard form) <br> ALLOW one mark for <br> $6.16 \times 10^{44}$ (multiplying by $6.02 \times 10^{23}$ in M1 + standard form) |


| Question |  | Answer | Mark | Guidance |
| :--- | :--- | :--- | :--- | :---: | :--- |
| $\mathbf{1}$ | (e) | (i) | $\begin{array}{c}\mathrm{N}_{2} \mathrm{O}_{3}=+3 \\ \mathrm{NO}_{3}=+2\end{array}$ |  |
| $\mathrm{NO}_{2}=+4$ |  |  |  |  |$)$


| Question |  | Answer | Mark | Guidance |
| :--- | :--- | :--- | :--- | :--- | :--- |
| $\mathbf{2}$ | (a) | Simple molecular lattice $\checkmark$ | ALLOW 'simple covalent' OR 'simple molecular' <br> ie 'simple' must be seen. <br> DO NOT ALLOW 'simple covalent bonds' |  |
| $\mathbf{2}$ | (b) | (i)M1 Creating the dipole mark <br> Uneven distribution of electrons $\checkmark$ <br> M2 Type of dipole mark <br> This creates/causes an instantaneous dipole <br> OR temporary dipole $\checkmark$ | IGNORE use of 'atoms' for M1 and M2 <br> ALLOW (random) movement of electrons <br> ALLOW change in electron density |  |


| Question |  | Answer | Mark | Guidance |
| :---: | :---: | :--- | :--- | :--- |
| 2 | (c) |  |  |  |


| Question |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: |
| 3 | (a) | Periodicity $\checkmark$ | 1 |  |
| 3 | (b) | Sodium OR Na Silicon OR Si $\downarrow$ Neon OR Ne $\checkmark$ | 3 |  |
| 3 | (c) | $\mathrm{Ga}^{3+} \checkmark$ | 1 |  |
| 3 | (d) | M1 Number of bonding electrons mark <br> Magnesium has more outer OR bonding electrons <br> M2 Ionic charge mark Magnesium ions have a greater (positive) charge (density) <br> M3 Attraction mark Magnesium has a greater attraction between ions and delocalised electrons | 3 | ALLOW reverse argument throughout <br> ALLOW 'more delocalised electrons' for 'more outer electrons' <br> DO NOT ALLOW 'Magnesium molecules' for M1 <br> ALLOW $\mathrm{Mg}^{2+}$ ion OR Mg ion for 'magnesium ion' ALLOW $\mathrm{Mg}^{2+}$ and $\mathrm{Na}^{+}$for M2 (may be seen in a diagram) IGNORE magnesium has a greater charge but ALLOW magnesium has a greater ionic charge <br> IGNORE nuclear charge <br> DO NOT ALLOW 'atoms' or 'molecules' having a greater charge for M2 <br> ALLOW 'stronger metallic bonds' only when a clear description of metallic bonding is given. Eg 'The attraction of positive (metal) ions to delocalised electrons' <br> QWC 'delocalised/delocalized' spelled correctly at least once in context of M3 (may be seen in M1 but used in M3) <br> 'delocalised' need not be directly next to electrons eg Mg has more delocalised electrons and the ions have a greater attraction to these electrons would secure M3 |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 3 | (e) |  | First check the answer line. <br> If answer $=1200 \mathrm{~cm}^{3}$ award 3 marks. <br> Mol of $\mathrm{Mg}\left(\mathrm{NO}_{3}\right)_{2}=\frac{2.966}{148.3}=2(.00) \times 10^{-2}$ OR 0.02(00) mol V <br> Mol of gas $=2(.00) \times 10^{-2} \times 5 / 2=5(.00) \times 10^{-2}$ <br> OR 0.05(00) mol <br> Vol of Gas $=0.05 \times 24000=1200 \mathrm{~cm}^{3} \checkmark$ | 3 | If answer $=960 \mathrm{~cm}^{3}$ award 2 marks. <br> If answer $=240 \mathrm{~cm}^{3}$ award 2 marks. <br> ALLOW ECF for answers to at least two significant figures up to calculator value, correctly rounded <br> ALLOW separate numbers of mol of each gas for M2 (0.04(00) $\mathrm{mol} \mathrm{NO}_{2}$ and $0.0100 \mathrm{~mol} \mathrm{O}_{2}$ ) <br> ALLOW a second mark if only volume of $\mathrm{O}_{2}\left(240 \mathrm{~cm}^{3}\right)$ OR only volume of $\mathrm{NO}_{2}\left(960 \mathrm{~cm}^{3}\right)$ is calculated |
| 3 | (f) | (i) | $\mathrm{SF}_{6}$ <br> AND <br> Sulfur(VI) fluoride OR Sulfur hexafluoride | 1 | IGNORE sulfur fluoride |
|  |  | (ii) | $2 \mathrm{~F}_{2}+2 \mathrm{NaOH} \rightarrow \mathrm{~F}_{2} \mathrm{O}+2 \mathrm{NaF}+\mathrm{H}_{2} \mathrm{O}$ <br> M1 $\mathrm{F}_{2} \mathrm{O}$ AND $\mathrm{NaF} \checkmark$ <br> M2 Rest of equation (including balance) | 2 | ALLOW multiples <br> IGNORE state symbol <br> ALLOW $\mathrm{OF}_{2}$ for $\mathrm{F}_{2} \mathrm{O}$ AND FNa for NaF <br> ALLOW both marks for alternative equations which have both $\mathrm{F}_{2} \mathrm{O}$ and NaF AND three products <br> $\mathrm{Eg}_{3} \mathrm{~F}_{2}+2 \mathrm{NaOH} \rightarrow 2 \mathrm{~F}_{2} \mathrm{O}+2 \mathrm{NaF}+\mathrm{H}_{2}$ <br> $\mathrm{Eg}_{2} \mathrm{~F}_{2}+\mathrm{NaOH} \rightarrow \mathrm{F}_{2} \mathrm{O}+\mathrm{NaF}+\mathrm{HF}$ |
| 3 | (g) | (i) | $\delta$ - on each F AND $\delta+$ on $\mathrm{O} \checkmark$ | 1 | ALLOW $\delta 2+$ OR $\delta+\delta+$ on O |
|  |  | (ii) | Shape: non-linear <br> AND <br> Bond angle: $104.5^{\circ}$ | 1 | For shape <br> ALLOW alternative words eg 'V-shaped' 'bent' ‘angular'. In the absence of words allow a diagram with a non-linear shape F - $\mathrm{O}-\mathrm{F}$ bond angle $>90^{\circ}$. <br> For bond angle <br> ALLOW 106 $>$ bond angle $\geq 102\left(\right.$ Actual $\left.=102^{\circ}\right)$ |
|  |  | (iii) | $+2 \checkmark$ | 1 | ALLOW 2+ |
|  |  |  | Total | 17 |  |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (a) |  | $1 s^{2} 2 s^{2} 2 p^{6} 3 s^{2} 3 p^{6} 3 d^{10} 4 s^{2} \checkmark$ | 1 | ALLOW 4s ${ }^{2} 3 \mathrm{~d}^{10}$ |
| 4 | (b) | (i) | M1 <br> The (weighted) mean mass of an atom (of an element) <br> M2 <br> Compared with $1 / 12^{\text {th }}$ (the mass) <br> M3 <br> Of (one atom of) carbon-12 | 3 | ALLOW 'average' for 'mean' <br> ALLOW 'mean mass of isotopes' <br> but DO NOT ALLOW 'mean mass of isotope' (singular) <br> DO NOT ALLOW 'mean mass of an element' <br> For M2 and M3 <br> ALLOW compared with the mass of carbon-12 which is 12 <br> ALLOW for three marks <br> Mass of one mole of atoms <br> Compared to $1 / 11^{\text {th }}$ <br> (mass of) one mole OR 12 g of carbon-12 <br> ALLOW for three marks <br> Mass of one mole of atoms <br> $1 / 12^{\text {th }}$ (mass of) one mole OR 12 g of carbon-12 |
| 4 | (b) | (ii) | First check the answer line. <br> If answer $=65.44$ award 2 marks. $\frac{(64 \times 49.0)+(66 \times 27.9)+(67 \times 4.3)+(68 \times 18.8)}{100}$ <br> OR $31.36(0)+18.414+2.881+12.784$ <br> OR <br> 65.439 $=65.44 \checkmark$ | 2 | ALLOW one mark for ECF from transcription error in the first sum provided the final answer is to two decimal places and is between 64 and 68 and is a correct calculation of the transcription |
| 4 | (c) | (i) | Effervescence OR fizzing OR bubbling OR gas produced AND <br> The solid OR zinc carbonate would dissolve OR disappear | 1 | ALLOW 'carbon dioxide produced' DO NOT ALLOW incorrectly named gas eg $\mathrm{H}_{2}$ |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (c) | (ii) | $\mathrm{ZnCO}_{3}+2 \mathrm{HCl} \rightarrow \mathrm{ZnCl}_{2}+\mathrm{CO}_{2}+\mathrm{H}_{2} \mathrm{O} \checkmark$ | 1 | ALLOW multiples IGNORE state symbols |
| 4 | (d) | (i) | Magnesium (atoms) has been oxidised AND <br> Because it has lost two electrons <br> Copper (ions) has been reduced AND <br> Because it has gained two electrons $\checkmark$ | 2 | IGNORE use of oxidation numbers if electron gain/loss is mentioned. <br> Electrons gain/loss could be in half equations In the absence of text look for evidence on the equation ALLOW 'donated' for 'lost' <br> Assume 'Cu' refers to copper in 'CuSO ${ }_{4}$ ' ALLOW one mark two electrons gained and lost for each species but oxidation/reduction is incorrect or is omitted <br> ALLOW one mark for correct oxidation and reduction if electron transfer is omitted and correct changes of oxidation state are shown (ie Mg $0-->(+) 2$ AND Cu (+)2 to 0) <br> ALLOW 'two electrons transferred from magnesium to copper' |
| 4 | (d) | (ii) | $\mathrm{Mg}(\mathrm{~s})+2 \mathrm{H}_{2} \mathrm{O}(\mathrm{l}) \rightarrow \mathrm{Mg}(\mathrm{OH})_{2}(\mathrm{aq})+\mathrm{H}_{2}(\mathrm{~g})$ <br> Correct reactants and products $\checkmark$ <br> Balance and state symbols $\checkmark$ | 2 | ALLOW multiples <br> ALLOW $\mathrm{Mg}(\mathrm{OH})_{2}(\mathrm{~s})$ <br> ALLOW Mg(s) $+\mathrm{H}_{2} \mathrm{O}(\mathrm{g})$ OR $\mathrm{H}_{2} \mathrm{O}(\mathrm{l})$--> $\mathrm{MgO}(\mathrm{s})+\mathrm{H}_{2}(\mathrm{~g})$ <br> including state symbols for one mark |


| Question |  |  | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: | :---: | :---: |
| 4 | (e) |  | First check the answer line. If answer $=0.120$ award 4 marks. <br> M1 Mol of $\mathrm{H}_{2} \mathrm{SO}_{4}=3.00 \times 10^{-2} \times \frac{35.0}{1000}=1.05 \times 10^{-3} \mathrm{~mol}$ <br> M 2 Mol of $\mathrm{Al}_{2}\left(\mathrm{SO}_{4}\right)_{3}=\frac{1.05 \times 10^{-3}}{3}=3.5(0) \times 10^{-4} \mathrm{~mol}$ $\mathrm{M} 3=342.3 \checkmark$ <br> M4 Mass $\mathrm{Al}_{2}\left(\mathrm{SO}_{4}\right)_{3}=3.5(0) \times 10^{-4} \times 342.3$ <br> and $=0.120 \mathrm{~g}$ <br> Answer must be $\mathbf{3}$ sf | 4 | ALLOW ECF <br> ALLOW 0.00105 mol <br> ALLOW 0.00035(0) mol <br> ALLOW 342 <br> DO NOT ALLOW 0.12 |
| 4 | (f) | (i) | $\mathrm{Ca}(\mathrm{OH})_{2}$ OR Calcium hydroxide OR CaO OR Calcium oxide $\checkmark$ | 1 | ALLOW Calcium carbonate $\mathrm{OR} \mathrm{CaCO}_{3}$ |
| 4 | (f) | (ii) | $6 \mathrm{Ca}+\mathrm{P}_{4} \rightarrow 2 \mathrm{Ca}_{3} \mathrm{P}_{2} \checkmark$ | 1 | ALLOW multiples IGNORE state symbols |


| Question | Answer | Mark | Guidance |
| :---: | :---: | :---: | :---: |
| (iii) | $3 x\left[\left[\begin{array}{c} x_{x x} \\ x_{x} \operatorname{Cax}_{x}^{x} \\ x]^{2+} \end{array} \quad 2 x\left[\begin{array}{c} \bullet \bullet \\ \dot{x} P_{0}^{x} \\ 0 x \end{array}\right]^{2}\right.\right.$ <br> Ca with 8 (or no) electrons AND phosphide ion with dot-and-cross outermost octet $\checkmark$ <br> Three Ca ions AND two phosphide ions with correct charges | 2 | For first mark: <br> If 8 electrons are shown on the cation then the extra electron in the anion must match the symbol chosen for the electrons in the cation. <br> IGNORE inner shells <br> IGNORE circles <br> ALLOW one mark if both electron arrangements and charges are correct but only one of each ion is drawn. <br> ALLOW (brackets not required) <br> $3\left[\mathrm{Ca}^{2+}\right] 3[\mathrm{Ca}]^{2+}\left[\mathrm{Ca}^{2+}\right]_{3}$ <br> $2\left[\mathrm{P}^{3-}\right] 2[\mathrm{P}]^{3-}\left[\mathrm{P}^{3-}\right]_{2}$ <br> DO NOT ALLOW <br> $\left[\mathrm{Ca}_{3}\right]^{2+}[3 \mathrm{Ca}]^{2+}[\mathrm{Ca}]_{3}{ }^{2+}$ <br> $\left[\mathrm{P}_{2}\right]^{3-} \quad[2 \mathrm{P}]^{3-}\left[\mathrm{P}_{2}{ }^{3-}\right.$ |
|  | Total | 20 |  |

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