

## CHEMISTRY

9701/22 March 2017

Paper 2 AS Level Structured Questions MARK SCHEME Maximum Mark: 60

Published

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.

Mark schemes should be read in conjunction with the question paper and the Principal Examiner Report for Teachers.

Cambridge will not enter into discussions about these mark schemes.

Cambridge is publishing the mark schemes for the March 2017 series for most Cambridge IGCSE<sup>®</sup>, Cambridge International A and AS Level components and some Cambridge O Level components.

® IGCSE is a registered trademark.

# Cambridge International AS/A Level – Mark Scheme PUBLISHED

| Question  | Answer  |   |  |                                  |                           | Marks |    |                    |
|-----------|---|---|--|----------------------------------|---------------------------|-------|----|--------------------|
| 1(a)(i)   | max<br>O.N.   | +1  | (+)2                                       | (+)3                             | (+)5                      | (+)6  | +7 | 1                  |
| 1(a)(ii)  | (from Na to   | o Cl) nucle   | ar charge i                                | ncreases                         |                           |       |    | 1                  |
|           | electrons are in the same shell/have same shielding |   |  |                                  | 1                         |       |    |                    |
|           | greater/str   | ronger attra  | ction (of e                                | lectrons to r                    | nucleus)                  |       |    | 1                  |
| 1(a)(iii) | Mg <sup>2+</sup> AND                                | S <sup>2-</sup>   |  |                                  |                           |       |    | 1                  |
|           | ion of Mg/  | Mg <sup>2+</sup> has o                                      | ne fewer s                                 | hell (than io                    | on of S/S <sup>2-</sup> ) | )     |    | 1                  |
| 1(b)(i)   | P <sub>4</sub> + 5O <sub>2</sub>                    | $\rightarrow P_4O_{10}/$                                    | 2P <sub>2</sub> O <sub>5</sub>             |                                  |                           |       |    | 1                  |
| 1(b)(ii)  | any 2 from<br>• yell<br>• whi<br>• whi<br>• soli    | :<br>low/green<br>ite flame<br>ite solid<br>id melts        | colour (of c                               | chlorine gas                     | s) disappea               | rs    |    | 2                  |
| 1(b)(iii) | phosphoric(V) acid                                  |   |  |                                  | 1                         |       |    |                    |
| 1(c)(i)   | € € € € € € € € € € € € € € € € € € €               | • • • • • • • • • • • • • • • • •                           | llar arrang                                | ement of (p<br>ed) electror      | ositive) ion              | S     |    | <b>2</b><br>1<br>1 |
| 1(c)(ii)  | any 2 from<br>higl<br>elec<br>har<br>reta           | :<br>h melting/t<br>ctrical/ther<br>d/rigid<br>ains strengt | ooiling/sub<br>mal insulat<br>h at high te | limation po<br>tor<br>emperature | int<br>/pressure          |       |    | 2                  |

# Cambridge International AS/A Level – Mark Scheme PUBLISHED

| Question  | Answer  | Marks |
|-----------|---|-------|
| 1(c)(iii) | M1<br>% abundance of fourth isotope<br>= 100 - (0.185 + 0.251 + 88.450) = 11.114  | 1     |
|           | $ \frac{M2}{(0.185 \times 135.907) + (0.251 \times 137.906) + (88.450 \times 139.905) + (11.114 \times RIM)}{100} = 140.116 $ $ \therefore (140.116 \times 100) - 12434.35 = 1577.246 = 11.114 \times RIM $ | 1     |
|           | M3<br>RIM = $\frac{1577.246}{11.114}$ = 141.915   | 1     |

| Question  | Answer  | Marks |
|-----------|---|-------|
| 2(a)(i)   | bond in which the centres of positive and negative charges do not coincide <b>OR</b> electron distribution is asymmetric/unequal <b>OR</b> two (bonded) atoms are partially charged | 1     |
| 2(a)(ii)  | HF has the strongest (permanent) dipole–dipole/van der Waals'<br>(forces)/HF has hydrogen bonding   | 1     |
|           | requires more energy to overcome (than weaker (permanent) dipole–dipole/<br>van der Waals' forces between other hydrogen halides)   | 1     |
| 2(a)(iii) | thermal stability of the hydrogen halides decreases down group (17)   | 1     |
|           | larger (halogen) atoms/atomic radius (down group) / increased shielding   | 1     |
|           | bond energies decrease/less energy required to break H–X  | 1     |
| 2(b)(i)   | M1<br>base is $Cl^-$ AND conjugate acid is HC $l$<br>OR<br>base is HSO <sub>4</sub> <sup>-</sup> AND conjugate acid is H <sub>2</sub> SO <sub>4</sub>                               | 1     |
|           | M2<br>$Cl^{-}/HSO_{4}^{-}/base$ is a proton acceptor<br>OR<br>$HCl/H_{2}SO_{4}/(conjugate)$ acid has one more $H^{+}$   | 1     |
| 2(b)(ii)  | H <sub>2</sub> SO <sub>4</sub> is (too strong) an oxidising agent   | 1     |
|           | I <sub>2</sub> would be formed instead  | 1     |

# Cambridge International AS/A Level – Mark Scheme **PUBLISHED**

| Question  | Answer   |                                |   |  |                  |   |
|-----------|--|--------------------------------|---|--|------------------|---|
| 2(c)(i)   | $\Delta_{\rm r} H = \Delta_{\rm r} H \{ \text{products} \} - \Delta_{\rm r} H \{ \text{reactants} \} = 2 \times (-242) - 4 \times (-92)$                               |                                |   |  |                  | 1 |
|           | = –116 (sign <b>AND</b> answer)  |                                |   |  |                  | 1 |
| 2(c)(ii)  | heterogeneous (catalyst)   |                                |   |  |                  | 1 |
|           | provides an alternative reaction pathway of lower activation energy  |                                |   |  |                  | 1 |
| 2(c)(iii) | reaction is exothermic   |                                |   |  |                  | 1 |
|           | (increased temperature) shifts equilibrium to the left <b>AND</b> decreases yield of products ( $Cl_2$ and/or $H_2O$ )/less product formed                             |                                |   |  |                  | 1 |
| 2(c)(iv)  |  | HCl                            | O <sub>2</sub>  | Cl <sub>2</sub>  | H <sub>2</sub> O | 3 |
|           | initial number<br>of moles   | 1.60                           | 0.500   | 0  | 0                |   |
|           | M1 eqm<br>number of<br>moles   | 1.60 -<br>2 × 0.600<br>= 0.400 | 0.500 -<br><sup>1</sup> / <sub>2</sub> × 0.600<br>= 0.200 | 0.600  | 0.600            |   |
|           | M2 mole<br>fraction  |                                |   | <u>0.600</u><br>1.80                                       |                  |   |
|           | M3 partial pressure  |                                |   | $\frac{0.600}{1.80} \times p_{\rm tot} = 5.00 \times 10^4$ |                  |   |
| 2(c)(v)   | $K_{\rm p} = \frac{\left(3.6 \times 10^4\right)^2 \times \left(3.6 \times 10^4\right)^2}{\left(4.8 \times 10^4\right)^4 \times 3.0 \times 10^4} = 1.05 \times 10^{-5}$ |                                |   |  |                  | 1 |
|           | units = Pa <sup>-1</sup>   |                                |   |  |                  | 1 |
| 2(c)(vi)  | K <sub>p</sub> would not change  |                                |   |  |                  | 1 |

| Question | Answer  | Marks |
|----------|---|-------|
| 3(a)(i)  |   | 1     |
| 3(a)(ii) | reaction <b>1</b> = HC <i>l</i> (aq)                    | 1     |
|          | reaction <b>2</b> = (conc.) NaOH/KOH <b>AND</b> ethanol | 1     |

# Cambridge International AS/A Level – Mark Scheme PUBLISHED

| Question  | Answer  | Marks       |
|-----------|---|-------------|
| 3(a)(iii) | $\begin{array}{c} \begin{array}{c} H & C_2H_5 \\ \hline \\ - C & -C \\ H & H \end{array}$ C-C backbone with dangling bonds<br>rest of structure | 2<br>1<br>1 |
| 3(b)      | CH <sub>2</sub>                 | 3<br>1<br>1 |
| 3(c)(i)   | (electrophilic) addition  | 1           |
| 3(c)(ii)  | <b>S</b> has CH <sub>3</sub> CHOH <b>OR</b> methyl/CH <sub>3</sub> group next to CHOH   | 1           |
| 3(c)(iii) | positive inductive effect of more alkyl groups/more alkyl groups donate electron density  | 1           |
|           | secondary carbocation/secondary intermediate is more stable (than primary)  | 1           |
| 3(c)(iv)  | S =   | 1           |
|           | T = HO  | 1           |
|           | U = 0   | 1           |
| 3(c)(v)   | $CH_3CHOHCH_2CH_3 + [O] \rightarrow CH_3COCH_2CH_3 + H_2O$  | 1           |
| 3(d)(i)   | methyl pentanoate   | 1           |
| 3(d)(ii)  | (compound <b>V</b> is) spectrum X   | 1           |
|           | spectra X and Z show a C=O (stretch) at 1730 (cm <sup>-1</sup> )  | 1           |
|           | spectra Y and Z show O–H (stretches) above 2500 (cm <sup>-1</sup> )   | 1           |
|           | V has a C=O (bond) and no O–H (bond)  | 1           |