
CHEMISTRY

9701/21

Paper 2 AS Level Structured Questions

May/June 2016

MARK SCHEME

Maximum Mark: 60

Published

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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.

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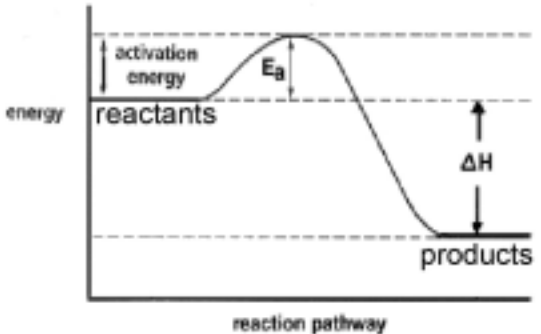
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|-----------------|---|-------------------|----------------|-----------------|------------------|-----------------|------------------|----------------|---------|---|---|---|---|---|----|--------|----|---|---|---|----|----|------|----|----|----|----|----|----|----------|----|----|----|----|----|---|--------------------------|-----|
| 1 (a) | <table border="1"> <thead> <tr> <th>name of element</th> <th>nucleon no.</th> <th>atomic no.</th> <th>no. of protons</th> <th>no. of neutrons</th> <th>no. of electrons</th> <th>overall charge</th> </tr> </thead> <tbody> <tr> <td>lithium</td> <td>6</td> <td>3</td> <td>3</td> <td>3</td> <td>2</td> <td>+1</td> </tr> <tr> <td>oxygen</td> <td>17</td> <td>8</td> <td>8</td> <td>9</td> <td>10</td> <td>-2</td> </tr> <tr> <td>iron</td> <td>54</td> <td>26</td> <td>26</td> <td>28</td> <td>24</td> <td>+2</td> </tr> <tr> <td>chlorine</td> <td>35</td> <td>17</td> <td>17</td> <td>18</td> <td>17</td> <td>0</td> </tr> </tbody> </table> | name of element | nucleon no. | atomic no. | no. of protons | no. of neutrons | no. of electrons | overall charge | lithium | 6 | 3 | 3 | 3 | 2 | +1 | oxygen | 17 | 8 | 8 | 9 | 10 | -2 | iron | 54 | 26 | 26 | 28 | 24 | +2 | chlorine | 35 | 17 | 17 | 18 | 17 | 0 | [1] [1] [1] [1] | [4] |
| name of element | nucleon no. | atomic no. | no. of protons | no. of neutrons | no. of electrons | overall charge | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| lithium | 6 | 3 | 3 | 3 | 2 | +1 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| oxygen | 17 | 8 | 8 | 9 | 10 | -2 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| iron | 54 | 26 | 26 | 28 | 24 | +2 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| chlorine | 35 | 17 | 17 | 18 | 17 | 0 | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| (b) | line straight on labelled 'neutrons' line (curving) up labelled 'protons' proton line clearly shows less (overall) deflection than electron curve | [1] [1] [1] | [3] | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| (c) (i) | Group 16/6/VI AND Big (owtte) increase/big difference/big gap/big jump/jump in increase/jump in difference after 6th IE | [1] | [1] | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| (ii) | increases (across period) due to increasing attraction (of nucleus for electrons) due to increasing nuclear charge/atomic/proton number AND constant/similar shielding/same (outer/number of) shell/energy level | [1] [1] | [2] | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| (iii) | electron (pair) repulsion (Y has a) pair of electrons in a (3)p orbital/a (3)p orbital is full ORA | [1] [1] | [2] | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| (iv) | $(1s^2)2s^22p^63s^23p^5$ | [1] | [1] | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
| (d) (i) | 0.56(%) | [1] | [1] | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |

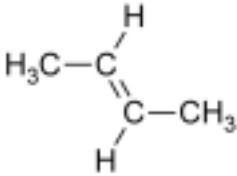
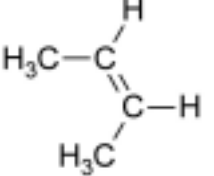
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| (ii) | $\frac{(A \times 0.56) + (86 \times 9.86) + (87 \times 7.00) + (88 \times 82.58)}{100} = 87.71$ <p>A = 84</p> | [1] [1] | [2] |
| | | | [16] |
| 2 (a) | D = Ga G = Se | [1] | [1] |
| (b) (i) | $\text{D}_2\text{O}_3 + 6\text{HCl} \rightarrow 2\text{DCl}_3 + 3\text{H}_2\text{O}$ <p>M1 = species; M2 = balancing</p> | [1] [1] | [2] |
| (ii) | $\text{D}_2\text{O}_3 + 2\text{NaOH} + 7\text{H}_2\text{O} \rightarrow 2\text{NaD}(\text{OH})_4(\text{H}_2\text{O})_2 \text{ OR}$ $\text{D}_2\text{O}_3 + 2\text{NaOH} + 3\text{H}_2\text{O} \rightarrow 2\text{NaD}(\text{OH})_4 \text{ OR}$ $\text{D}_2\text{O}_3 + 2\text{NaOH} \rightarrow 2\text{NaDO}_2 + \text{H}_2\text{O} \text{ OR}$ $\text{D}_2\text{O}_3 + 2\text{OH}^- + 7\text{H}_2\text{O} \rightarrow 2[\text{D}(\text{OH})_4(\text{H}_2\text{O})_2]^- \text{ OR}$ $\text{D}_2\text{O}_3 + 2\text{OH}^- + 3\text{H}_2\text{O} \rightarrow 2[\text{D}(\text{OH})_4]^- \text{ OR}$ $\text{D}_2\text{O}_3 + 2\text{OH}^- \rightarrow 2\text{DO}_2^- + \text{H}_2\text{O}$ <p>M1 = species; M2 = balancing</p> | [1] [1] | [2] |
| (c) | giant ionic / ionic lattice | [1] | [1] |
| (d) | $\text{GO}_2 + \text{H}_2\text{O} \rightarrow \text{H}_2\text{GO}_3$ | [1] | [1] |
| | | | [7] |

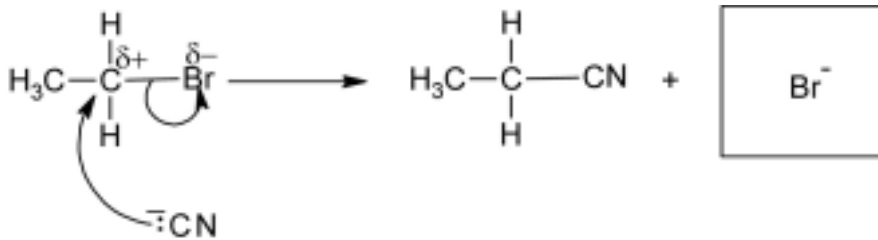
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| 3 (a) (i) | bubbles / effervescence / fizzing calcium gets smaller / disappears water turns cloudy / milky calcium sinks | [1] [1] [1] [1] | max [3] |
| (ii) | $\text{Ca} + 2\text{H}_2\text{O} \rightarrow \text{Ca}(\text{OH})_2 + \text{H}_2$ | [1] | [1] |
| (iii) | faster bubbling / disappearance of Ba OR no / less precipitate forms (owtte) | [1] | [1] |
| (b) (i) |  <p>M1 – general layout with products below reactants AND both labelled</p> <p>M2 – E_a and ΔH / energy change / released labelled with vertical lines</p> | [1] [1] | [2] |
| (ii) | activation energy is high so few / no particles with $E \geq E_a$ | [1] [1] | [2] |

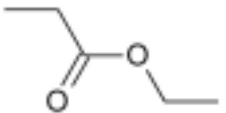
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| (iii) | high melting/boiling point strong forces (of attraction / between oppositely charged ions) / strong (ionic) bonding | [1] [1] | [2] |
| (iv) | MgO is basic / reacts with acid | [1] | [1] |
| (c) (i) | increases (down the group) | [1] | [1] |
| (ii) | $\text{MgCO}_3 \rightarrow \text{MgO} + \text{CO}_2$ | [1] | [1] |
| (iii) | $2\text{Ca}(\text{NO}_3)_2 \rightarrow 2\text{CaO} + 4\text{NO}_2 + \text{O}_2$ | [1] | [1] |
| | | | [15] |
| 4 (a) | $\text{CH}_2=\text{CHCH}_2\text{CH}_3 / \text{CH}_2\text{CHCH}_2\text{CH}_3$ AND $\text{CH}_3\text{CH}=\text{CHCH}_3 / \text{CH}_3\text{CHCHCH}_3$ | [1] | [1] |
| (b) | $\text{CH}_2=\text{CHCH}_2\text{CH}_3 / \text{CH}_2\text{CHCH}_2\text{CH}_3$ AND $(\text{CH}_3)_2\text{C}=\text{CH}_2 / (\text{CH}_3)_2\text{CCH}_2$ | [1] | [1] |
| (c) |   <i>trans-but-2-ene (or E)</i> <i>cis-but-2-ene (or Z)</i> | [1] [1] | [2] |

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| (d) | B is CH ₂ =CHCH ₂ CH ₃ OR CH ₃ CH=CHCH ₃ OR (CH ₃) ₂ C=CH ₂ distinguished by addition of bromine brown/red/orange/yellow to colourless/decolourises with B (but not A) | [1] [1] [1] | [3] |
| | | | [7] |
| 5 (a) |  <p>M1 = lone pair on C of CN⁻ AND curly arrow from lone pair to C of C—Br M2 = correct dipole on C—Br, curly arrow from C—Br bond to Br AND Br⁻</p> | [1] [1] | [2] |
| (b) (i) | reduction | [1] | [1] |
| (ii) | disappearance of peak / dip / trough / absorption at 1680–1730 due to (loss of) C=O OR peak at 3200–3650 due to (alcohol) O—H (formation) | [1] [1] [1] [1] | [2] |
| (c) (i) | sodium/potassium hydroxide aqueous | [1] [1] | [2] |

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| (ii) | ethanol | [1] | [1] |
| (d) (i) | (conc) H ⁺ / (conc) acid / (conc)H ₂ SO ₄ / (conc)H ₃ PO ₄ | [1] | [1] |
| (ii) |  | [1] | [1] |
| (iii) | ethyl propanoate | [1] | [1] |
| (e) (i) | V = CH ₃ CH ₂ CHCHCH ₂ CH ₃ / CH ₃ CH ₂ CH=CHCH ₂ CH ₃ T = CH ₃ CH ₂ CH(OH)CH(OH)CH ₂ CH ₃ | [1] [1] | [2] |
| (ii) | V = geometric(al) / <i>cis-trans</i> / <i>E-Z</i> T = optical | [1] [1] | [2] |
| | | | [15] |