
CHEMISTRY

9701/43

Paper 4 A Level Structured Questions

May/June 2017

MARK SCHEME

Maximum Mark: 100

Published

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

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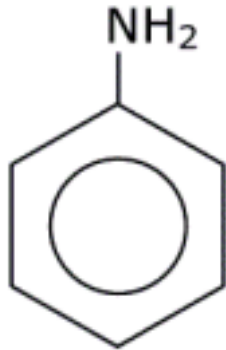
| Question | Answer | Marks |
|-----------|---|-------|
| 1(a) | solubility increases down the group | 1 |
| | ΔH_{latt} and ΔH_{hyd} both decrease or ΔH_{latt} and ΔH_{hyd} both become less exothermic / more endothermic | 1 |
| | ΔH_{latt} decreases / changes more (than ΔH_{hyd} as OH^- being smaller than M^{2+}) | 1 |
| | ΔH_{sol} becomes more exothermic / more negative / less endothermic / less positive | 1 |
| 1(b)(i) | $\Delta H_{\text{r1}} - (538 + 2 \times 230 + 394) = -(1216 + 286)$ $\Delta H_{\text{r1}} - 1392 = -1502$ | 1 |
| | $\Delta H_{\text{r1}} = \mathbf{-110}$ | 1 |
| | | |
| 1(b)(ii) | let $\Delta H_{\text{f}}(\text{HCO}_3^-(\text{aq})) = y$ $2y - 538 = -1216 - 394 - 286 - 26$ | 1 |
| | $y = \mathbf{-692}$ | 1 |
| 1(b)(iii) | $\Delta H_{\text{r3}} - 538 - 2(230 + 394) = -538 - 2(692)$ $\Delta H_{\text{r3}} = \mathbf{-136}$ | 1 |
| | | |
| 1(b)(iv) | ΔH_{r3} will be identical to ΔH_{r4} , / unchanged | 1 |
| | as the reaction is the same, or: $2\text{OH}^-(\text{aq}) + 2\text{CO}_2(\text{g}) \longrightarrow 2\text{HCO}_3^-(\text{aq})$ or metal ions stay in solution/metal ions are unchanged / are spectators | 1 |

| Question | Answer | Marks |
|----------|--|-----------|
| 1(c) | more gaseous moles are being consumed (in reaction 3) or more CO₂ moles are being consumed (in reaction 3) | 1 |
| | ΔS is therefore expected to be more negative/less positive for reaction 3. | 1 |
| | Total: | 13 |

| Question | Answer | Marks |
|-----------|---|--------------|
| 2(a)(i) |  | 1 + 1 |
| | 16 electrons on each diagram | 1 |
| 2(a)(ii) | HNC = 115–125° AND NCO = 180° | 1 |
| 2(a)(iii) | cyanic acid, because it's a stronger / higher bond enthalpy / triple / C≡N / more electrons involved bond | 1 |
| 2(b)(i) | $[H^+] = \sqrt{([HNCN]K_a)} = \sqrt{(0.1 \times 1.2 \times 10^{-4})}$ or 3.46×10^{-3} | 1 |
| | pH = log $[H^+] = 2.5$ (2.46) | 1 |
| 2(b)(ii) | $Na_2CO_3 + 2(NH_2)_2CO \longrightarrow 2NaNCO + CO_2 + 2NH_3 + H_2O$ | 1 |
| 2(c)(i) | $n(OH^-)$ at start = $(2 \times 0.1 \times 30) / 1000 = 6 \times 10^{-3}$ mol $n(OH^-)$ reacted = $(0.1 \times 20) / 1000 = 2 \times 10^{-3}$ mol $n(OH^-)$ remaining = $(6-2) \times 10^{-3} = 4 \times 10^{-3}$ mol, (in 50 cm ³) | 1 |
| | so $[OH^-]_{end} = (4 \times 10^{-3} \times 1000) / 50 = 0.08 \text{ mol dm}^{-3}$ | 1 |

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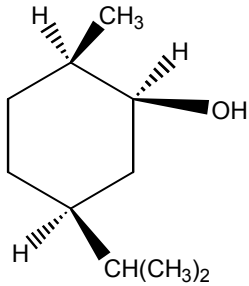
| Question | Answer | Marks |
|-----------|--|----------|
| 2(c)(ii) | $[H^+] = K_w / [OH^-] = (1 \times 10^{-14}) / 0.08 = 1.25 \times 10^{-13} \text{ mol dm}^{-3}$ | 1 |
| | so $\text{pH} = -\log(1.25 \times 10^{-13}) = \mathbf{12.9}$ | 1 |
| 2(c)(iii) | curve starts at 2.46 / 2.5 | 1 |
| | vertical portion (end point) at vol added = 10.0 cm ³ | 1 |
| | finishes at pH = 12.9 | 1 |
| 2(d)(i) | <i>monodentate</i> : (a species that) forms one dative / coordinate bond | 1 |
| | <i>ligand</i> : a species that uses a lone pair of electrons to form a dative / coordinate bond to a metal atom / metal ion | 1 |
| 2(d)(ii) | $[Ag(NCO)_2]^-$ or $[Ag(OCN)_2]^-$ correct formula | 1 |
| | correct charge | 1 |
| 2(e)(i) | $n(\text{BaCO}_3) = 1.66 / 197.3 = 8.4(1) \times 10^{-3} \text{ mol}$ | 1 |
| 2(e)(ii) | $n(\text{RNCO}) = 8.41 \times 10^{-3} \text{ mol}$, so $M_r = 1 / (8.41 \times 10^{-3}) = \mathbf{119}$ | 1 |
| 2(e)(iii) | molecular formula = C ₇ H ₅ NO | 1 |

| Question | Answer | Marks |
|----------|---|-----------|
| 2(e)(iv) |  | 1 |
| | Total: | 23 |

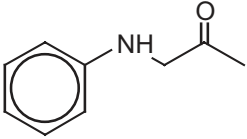
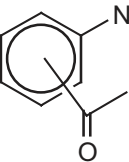
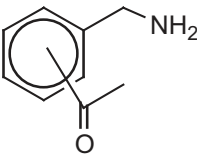
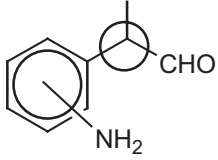
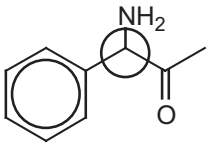
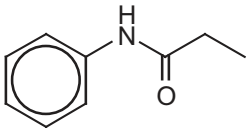
| Question | Answer | Marks |
|----------|---|-------|
| 3(a)(i) | +3 or Co^{3+} | 1 |
| 3(a)(ii) | oxidation | 1 |
| | ligand displacement / replacement / exchange / substitution | 1 |

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| Question | Answer | Marks |
|-----------|--|--------------|
| 4(a)(i) | optical, because it contains a / one chiral C-atom or chiral C-atoms or chiral atom / centre or C* indicated or C with 4 different groups | 1 |
| 4(a)(ii) | $\text{C}_{10}\text{H}_{14}\text{O} + 3\text{H}_2 \longrightarrow \text{C}_{10}\text{H}_{20}\text{O}$ correct formulae | 1 |
| | balancing | 1 |
| 4(b)(i) | electrophilic substitution | 1 |
| 4(b)(ii) | step 3 reduction | 1 |
| | step 5 substitution / hydrolysis | 1 |
| 4(b)(iii) | step 1 $(\text{CH}_3)_2\text{CHCl} + \text{AlCl}_3 / \text{AlBr}_3 / \text{FeCl}_3 / \text{FeBr}_3$ | 1 + 1 |
| | step 2 $\text{HNO}_3 + \text{H}_2\text{SO}_4$ conc (T < 55 °C) | 1 |
| | step 3 $\text{Sn} + \text{HCl}$ | 1 |
| | step 4 HNO_2 (or $\text{NaNO}_2 + \text{HCl}$) (at T < 10 °C) | 1 |
| | the two temperatures for steps 2 and 4 | 1 |
| 4(c)(i) | $\text{H}_2 + \text{Pt}$ or $\text{H}_2 + \text{Ni}$ + heat or pressure | 1 |

| Question | Answer | Marks |
|----------|--|-----------|
| 4(c)(ii) |  <p>(CH₃)₂CH, CH₃ and OH on the correct ring atoms i.e. structure is correct</p> | 1 |
| | all Hs on the same side of the ring | 1 |
| | Total: | 15 |

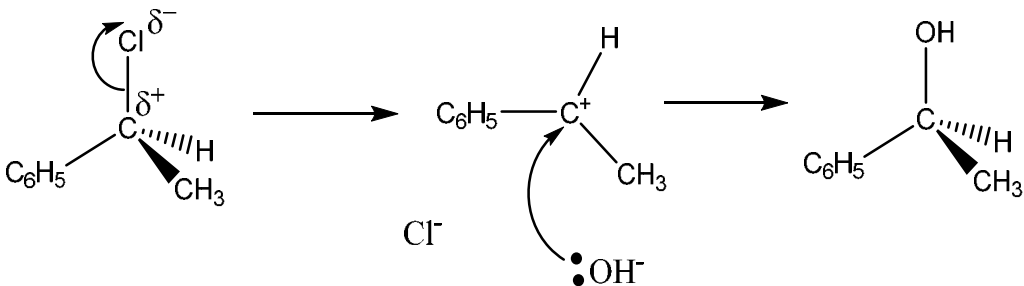
| Question | Answer | | | | Marks | | | | | | | | |
|-----------|---|----------------------------|------------------------|-------|-------|---|---|---|------------------------|----------------------------|------------------------|-------|--|
| 5(a) | <table><tr><td>J</td><td>K</td><td>L</td><td>M</td></tr><tr><td>amine methyl ketone</td><td>aromatic amine aldehyde</td><td>amine methyl ketone</td><td>amide</td></tr></table> | | | | J | K | L | M | amine methyl ketone | aromatic amine aldehyde | amine methyl ketone | amide | |
| | J | K | L | M | | | | | | | | | |
| | amine methyl ketone | aromatic amine aldehyde | amine methyl ketone | amide | | | | | | | | | |
| | J and L correct | | | | 1 + 1 | | | | | | | | |
| K correct | | | | 1 + 1 | | | | | | | | | |
| M correct | | | | 1 | | | | | | | | | |
| 5(b)(i) | hydrolysis | | | | 1 | | | | | | | | |
| 5(b)(ii) | P is C ₆ H ₅ NH ₂ | | | | 1 | | | | | | | | |
| | Q is CH ₃ CH ₂ CO ₂ Na | | | | 1 | | | | | | | | |

| Question | Answer | Marks |
|----------|---|-----------|
| 5(c) | J is  <i>or</i>  <i>or</i>  | 1 |
| | K is  | 1 |
| | L is  | 1 |
| | M is  | 1 |
| | K&L only: two chiral atoms shown | 1 |
| 5(d) | W is C ₆ H ₅ CO ₂ Na | 1 |
| | Total: | 14 |

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| Question | Answer | Marks |
|-----------|---|----------|
| 6(a) | <p>Any of the three methods possible. Any 4 of the 5 points for each method available for maximum 4 marks.</p> <p>Method 1</p> <ol style="list-style-type: none"> 1 Ensure both solutions (A and B) at 40 °C before mixing 2 mix known volumes of A and B and start the clock 3 at known time take out a sample / X and add it to ice-cold solvent 4 titrate against HCl 5 repeat at time at known time intervals <p>Method 2</p> <ol style="list-style-type: none"> 1 Ensure both solutions (A and B) at 40 °C before mixing 2 mix known volumes of A and B and start the clock 3 at known time pour into ice-cold solvent or pour ice-cold solvent in 4 titrate against HCl 5 repeat with different concentrations of either A or B, or repeat using different times <p>Method 3</p> <ol style="list-style-type: none"> 1 Ensure both solutions (A and B) at 40 °C before mixing 2 mix known volumes of A and B and start the clock and add pH meter 3 at a known time 4 record the pH 5 repeat pH readings at known time intervals | 4 |
| 6(b)(i) | from 1 and 3: when $[\text{RCI}]$ is trebled, so is rate, so order w.r.t. $[\text{RCI}] = 1$ | 1 |
| | from 1 and 2: when both concentrations are doubled, rate doubles so $[\text{OH}^-]$ has no effect on rate, so order w.r.t. $[\text{OH}^-] = 0$ | 1 |
| 6(b)(ii) | rate = $k[\text{RCI}]$ AND units: $\text{sec}^{-1} \text{ 1 / s}$ | 1 |
| 6(b)(iii) | relative rate = 2.0 | 1 |

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| Question | Answer | Marks |
|----------|---|----------------------------|
| 6(c)(i) |  <p>C-Cl dipole and first curly arrow</p> <p>intermediate cation</p> <p>OH⁻ with lone pair and curly arrow</p> | <p>1</p> <p>1</p> <p>1</p> |
| 6(c)(ii) | <p>Beginning with candidate's mechanism in (c)(i):</p> <p>If S_N1: racemate / mixture of / two optical isomers will be formed, because: the intermediate is planar / has a plane of symmetry / OH⁻ can approach from top or bottom or from any direction</p> <p>If S_N2: one optical isomer because attack always from fixed direction / from same side / the "configuration" always inverts / there is an asymmetric transition state</p> | 1 |

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| Question | Answer | | | | | Marks | | | | | | | | | | | | | | | | | | | | |
|--|---|-------------------|--------------------------|-----------|------------------------------|----------------|-------------------|-------|-----------|------------------------------|-----|---|--------------------------|---------|--------------|-----|---|-------------------------|---------|-----------------|-----|---|----|---------|--------------|--|
| 6(d)(i) | <table><tr><td>δ value</td><td>number of H atoms</td><td>group</td><td>splitting</td><td>result with D₂O</td></tr><tr><td>1.4</td><td>3</td><td>CH₃ / methyl</td><td>doublet</td><td>peak remains</td></tr><tr><td>2.7</td><td>1</td><td>OH / hydroxyl / alcohol</td><td>singlet</td><td>peak disappears</td></tr><tr><td>4.0</td><td>1</td><td>CH</td><td>quartet</td><td>peak remains</td></tr></table> | | | | | δ value | number of H atoms | group | splitting | result with D ₂ O | 1.4 | 3 | CH ₃ / methyl | doublet | peak remains | 2.7 | 1 | OH / hydroxyl / alcohol | singlet | peak disappears | 4.0 | 1 | CH | quartet | peak remains | |
| | δ value | number of H atoms | group | splitting | result with D ₂ O | | | | | | | | | | | | | | | | | | | | | |
| | 1.4 | 3 | CH ₃ / methyl | doublet | peak remains | | | | | | | | | | | | | | | | | | | | | |
| | 2.7 | 1 | OH / hydroxyl / alcohol | singlet | peak disappears | | | | | | | | | | | | | | | | | | | | | |
| | 4.0 | 1 | CH | quartet | peak remains | | | | | | | | | | | | | | | | | | | | | |
| | the three groups are in their correct places wrt the δ values | | | | | 1 | | | | | | | | | | | | | | | | | | | | |
| no. of H atoms for each peak agrees with group column | | | | | 1 | | | | | | | | | | | | | | | | | | | | | |
| splitting patterns doublet, singlet and quartet are assigned to correct groups | | | | | 1 | | | | | | | | | | | | | | | | | | | | | |
| peak identified as OH disappears with D ₂ O, no other peak disappears | | | | | 1 | | | | | | | | | | | | | | | | | | | | | |
| | Total: | | | | | 16 | | | | | | | | | | | | | | | | | | | | |