
CHEMISTRY

9701/43

Paper 4 A Level Structured Questions

October/November 2019

MARK SCHEME

Maximum Mark: 100

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 International will not enter into discussions about these mark schemes.

Cambridge International is publishing the mark schemes for the October/November 2019 series for most Cambridge IGCSE™, Cambridge International A and AS Level components and some Cambridge O Level components.

This document consists of **14** printed pages.

PUBLISHED**Generic Marking Principles**

These general marking principles must be applied by all examiners when marking candidate answers. They should be applied alongside the specific content of the mark scheme or generic level descriptors for a question. Each question paper and mark scheme will also comply with these marking principles.

GENERIC MARKING PRINCIPLE 1:

Marks must be awarded in line with:

- the specific content of the mark scheme or the generic level descriptors for the question
- the specific skills defined in the mark scheme or in the generic level descriptors for the question
- the standard of response required by a candidate as exemplified by the standardisation scripts.

GENERIC MARKING PRINCIPLE 2:

Marks awarded are always **whole marks** (not half marks, or other fractions).

GENERIC MARKING PRINCIPLE 3:

Marks must be awarded **positively**:

- marks are awarded for correct/valid answers, as defined in the mark scheme. However, credit is given for valid answers which go beyond the scope of the syllabus and mark scheme, referring to your Team Leader as appropriate
- marks are awarded when candidates clearly demonstrate what they know and can do
- marks are not deducted for errors
- marks are not deducted for omissions
- answers should only be judged on the quality of spelling, punctuation and grammar when these features are specifically assessed by the question as indicated by the mark scheme. The meaning, however, should be unambiguous.

GENERIC MARKING PRINCIPLE 4:

Rules must be applied consistently e.g. in situations where candidates have not followed instructions or in the application of generic level descriptors.

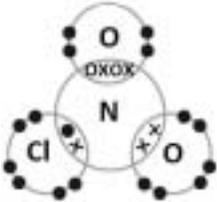
GENERIC MARKING PRINCIPLE 5:

Marks should be awarded using the full range of marks defined in the mark scheme for the question (however; the use of the full mark range may be limited according to the quality of the candidate responses seen).

GENERIC MARKING PRINCIPLE 6:

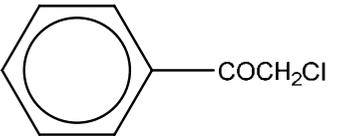
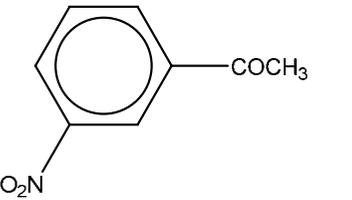
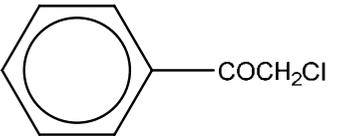
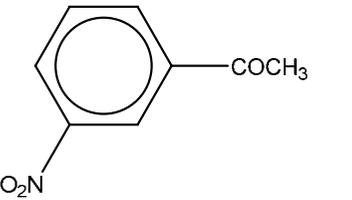
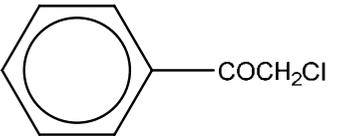
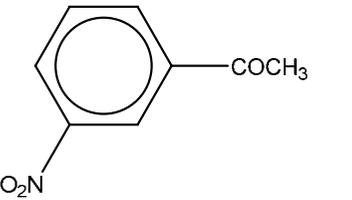
Marks awarded are based solely on the requirements as defined in the mark scheme. Marks should not be awarded with grade thresholds or grade descriptors in mind.

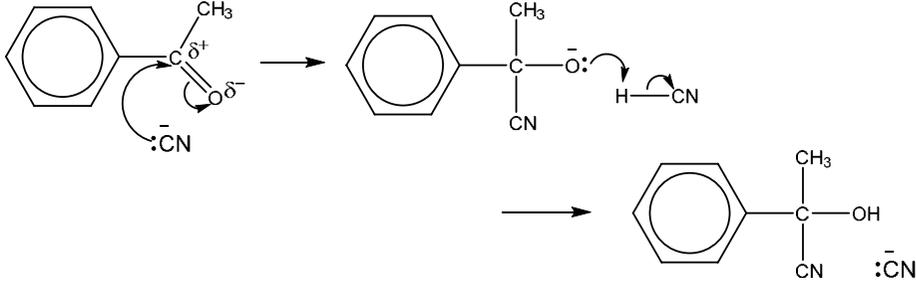
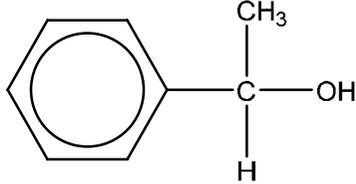
| Question | Answer | Marks |
|-----------|---|----------|
| 1(a) | Platinum / Pt Aluminium / Al BOTH | 1 |
| 1(b)(i) | M1: use of or quoting a valid Nernst equation $E = E^{\circ} + 0.0590 / z \log [\text{ox}] / [\text{red}]$ OR $E = 0.15 + (0.0590 / 2)\log 2$ M2: $E = (+)0.16$ (0.159) V minimum 2 sig. fig. correct answer scores 2 marks | 2 |
| 1(b)(ii) | $E_{\text{cell}} = 0.16 - (-1.66) = +1.82 \text{ V}$ minimum 3 sig. fig. | 1 |
| 1(b)(iii) | $2\text{Al} + 3\text{Sn}^{4+} \rightarrow 2\text{Al}^{3+} + 3\text{Sn}^{2+}$ M1: species M2: balancing | 2 |
| 1(c) | M1: number of C (= $300\,000 \times 60 \times 60 \times 24$) = 2.59×10^{10} (C) M2: number of F (= $2.592 \times 10^{10} / 9.65 \times 10^4$) = 2.69×10^5 (moles of electrons) M3: moles of Al (= $2.69 \times 10^5 / 3$) = 8.95×10^4 M4: mass of Al (= $8.95 \times 10^4 \times 27$) = 2420 kg correct answer scores 4 marks | 4 |
| 1(d) | M1: ($\text{Cr}^{2+} + 2\text{e}^{-} \rightleftharpoons \text{Cr}$) $E^{\circ} = -0.91$ and ($2\text{H}^{+} + 2\text{e}^{-} \rightleftharpoons \text{H}_2$) $E^{\circ} = 0.00$ seen M2: hydrogen formed instead / hydrogen (ions) easier to reduce / hydrogen has more positive E° | 2 |

| Question | Answer | Marks |
|-----------|---|-------|
| 2(a) |  <p>M1: eight electrons around N atom [N=O, N-O, N-Cl with N-O as dative] M2: all other electrons correct</p> | 2 |
| 2(b)(i) | (rate =) $k[\text{ClNO}_2][\text{NO}]$ | 1 |
| 2(b)(ii) | $\text{mol}^{-1} \text{dm}^3 \text{s}^{-1}$ | 1 |
| 2(b)(iii) | Yes AND number of moles of reactants in overall equation is the same as order in rate equation | 1 |
| 2(c)(i) | <ul style="list-style-type: none"> • straight line with a negative gradient • starting at 2.0×10^{-4} • reaches at 1.8×10^{-4} at 0.2 seconds <p>Award 1 mark for two points, award 2 marks for all three points</p> | 2 |
| 2(c)(ii) | $2 \times 10^{-5} \text{ (mol dm}^{-3}\text{)}$ | 1 |
| 2(c)(iii) | The reaction has reached equilibrium | 1 |

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| Question | Answer | Marks |
|-----------------|---|--------------|
| 3(a) | a measure / degree of disorder / randomness of a system | 1 |
| 3(b) | M1: negative – molecules have less energy in the system M2: positive – solid being converted into an aqueous solution M3: negative – gaseous ions being converted into a solid | 3 |
| 3(c)(i) | (standard) Gibbs free energy <u>change</u> | 1 |
| 3(c)(ii) | M1: $(\Delta)G = \Delta H - T\Delta S$ M2: description of calculating the minimum value of T for which ΔG is zero / becomes negative OR $T = \Delta H / \Delta S$ [1] | 2 |

| Question | Answer | | | Marks | | | | | | | | | | | | |
|-------------------|---|----------------------------|--|----------|-----------------|-------------------|----|--|---------------------------|-------------------|--|----------------------------|----|--------------------------------|--|----------|
| 4(a) | M1: CH_3COCl or ethanoyl chloride M2: AlCl_3 catalyst | | | 2 | | | | | | | | | | | | |
| 4(b) | <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th data-bbox="338 347 562 413">reagent</th> <th data-bbox="562 347 958 413">organic product</th> <th data-bbox="958 347 1256 413">name of mechanism</th> </tr> </thead> <tbody> <tr> <td data-bbox="338 413 562 651">Cl</td> <td data-bbox="562 413 958 651">  chlorine atom(s) in side chain only </td> <td data-bbox="958 413 1256 651">free radical substitution</td> </tr> <tr> <td data-bbox="338 651 562 954">nitric / sulfuric</td> <td data-bbox="562 651 958 954">  one only $-\text{NO}_2$ group added at 3 position </td> <td data-bbox="958 651 1256 954">electrophilic substitution</td> </tr> <tr> <td data-bbox="338 954 562 1023">Br</td> <td data-bbox="562 954 958 1023">no reaction with Br_2</td> <td data-bbox="958 954 1256 1023" style="background-color: #cccccc;"></td> </tr> </tbody> </table> <p data-bbox="338 1054 1016 1091">Award 1 mark for each correct entry to the table [5]</p> | | | reagent | organic product | name of mechanism | Cl |  chlorine atom(s) in side chain only | free radical substitution | nitric / sulfuric |  one only $-\text{NO}_2$ group added at 3 position | electrophilic substitution | Br | no reaction with Br_2 | | 5 |
| reagent | organic product | name of mechanism | | | | | | | | | | | | | | |
| Cl |  chlorine atom(s) in side chain only | free radical substitution | | | | | | | | | | | | | | |
| nitric / sulfuric |  one only $-\text{NO}_2$ group added at 3 position | electrophilic substitution | | | | | | | | | | | | | | |
| Br | no reaction with Br_2 | | | | | | | | | | | | | | | |
| 4(c)(i) | nucleophilic addition | | | 1 | | | | | | | | | | | | |

| Question | Answer | Marks |
|-----------|---|-------|
| 4(c)(ii) |  <p>M1 and M2: first structure – award one mark for two points correct, award two marks for four points correct.</p> <ul style="list-style-type: none"> • correct polarity shown on C=O bond • curly arrow from double bond to O of the C=O • lone pair shown on C of CN⁻ ion • curly arrow from C (on CN⁻ ion) to C of C=O <p>M3: middle structure</p> <ul style="list-style-type: none"> • correct intermediate AND curly arrow from lone pair on O⁻ to H⁺ or H of H₂O or HCN <p>M4: third structure</p> <ul style="list-style-type: none"> • correct product AND either CN⁻ reformed (if HCN seen in step 2) OR curly arrow on H-CN bond towards CN in step 2 | 4 |
| 4(d)(i) |  | 1 |
| 4(d)(ii) | LiAlH ₄ or NaBH ₄ | 1 |
| 4(d)(iii) | conc H ₂ SO ₄ / conc H ₃ PO ₄ / Al ₂ O ₃ | 1 |
| 4(d)(iv) | 6 | 1 |

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| Question | Answer | Marks |
|----------|--|----------|
| 4(d)(v) | <ul style="list-style-type: none"> • 25–50 • 110–160 • 190–220 <p>Award 1 mark for two points, award 2 marks for three points</p> | 2 |

| Question | Answer | Marks |
|-----------|--|----------|
| 5(a)(i) | $K_{sp} = [Ag^+]^2[S^{2-}]$ | 1 |
| 5(a)(ii) | <ul style="list-style-type: none"> • $[S^{2-}] = 1.16 \times 10^{-17}$ • $[Ag^+] = 2.32 \times 10^{-17}$ • $K_{sp} = \mathbf{6.2(4)} \times 10^{-51}$ minimum 2 sig. fig. <p>correct answer scores 2 marks Award 1 mark for two points, award 2 marks for three points</p> | 2 |
| 5(a)(iii) | <p>M1: moles $Ag_2S = 1 / 247.9 = 0.00403$ moles [1] 2sf min</p> <p>M2: $1.16 \times 10^{-17} = 0.0040 / V$ so $V = \mathbf{3.5 \times 10^{14}}$(dm³) [1] 2sf min ecf on M1</p> <p>correct answer scores 2 marks</p> | 2 |
| 5(b)(i) | <p>M1: $[H^+] = \sqrt{2.0 \times 10^{-9} \times 0.20}$ $[H^+] = 2.0 \times 10^{-5}$ (1.9976×10^{-5})</p> <p>M2: pH = 4.7 (4.699) minimum 2 sig. fig. min</p> <p>correct answer scores 2 marks</p> | 2 |

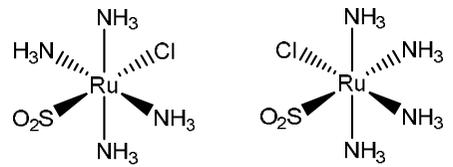
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| Question | Answer | Marks |
|----------|--|----------|
| 5(b)(ii) | <p>M1: Both equilibria correctly stated moles KOH = $0.005 \times 0.2 = 1 \times 10^{-3}$ moles HOBr(initial) = $0.020 \times 0.2 = 4 \times 10^{-3}$ moles HOBr(eqm) = $4 \times 10^{-3} - 1 \times 10^{-3} = 3 \times 10^{-3}$ moles BrO⁻(eqm) = 1×10^{-3}</p> <p>M2: ratio [OBr⁻]/[HOBr] = 1/3 [H⁺] = $3 \times 2.0 \times 10^{-9} = 6 \times 10^{-9}$ pH = 8.2(2)</p> <p>correct answer scores 2 marks</p> | 2 |

| Question | Answer | Marks |
|-----------|--|----------|
| 6(a)(i) | $\text{Mg(g)} \rightarrow \text{Mg}^{\text{+}}(\text{g}) + \text{e}^{-}$ | 1 |
| 6(a)(ii) | $\text{Sr(s)} + 2\text{H}_2\text{O(l)} \rightarrow \text{Sr(OH)}_2(\text{aq}) + \text{H}_2(\text{g})$ | 1 |
| 6(a)(iii) | more reactive and easier to ionise down the group OR more reactive and ionisation energies decrease down the group | 1 |
| 6(b)(i) | brown gas and white solid | 1 |
| 6(b)(ii) | $2\text{Ca(NO}_3)_2 \rightarrow 2\text{CaO} + 4\text{NO}_2 + \text{O}_2$ | 1 |
| 6(b)(iii) | <p>M1: more stable (down the group)</p> <p>M2: cationic radius increases / charge density of M²⁺ decreases (down the group)</p> <p>M3: NO₃⁻ anion is less polarised / distorted</p> | 3 |

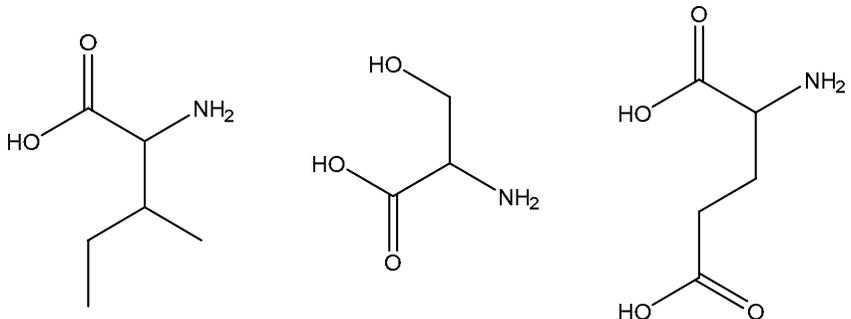
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| Question | Answer | Marks |
|----------|--|----------|
| 6(c) | <p>M1: less soluble / decreases (down the group)</p> <p>M2: ΔH_{latt} and ΔH_{hyd} both decrease / less exothermic down the group</p> <p>M3: ΔH_{hyd} decreases by more (than ΔH_{latt})</p> <p>M4: ΔH_{sol} becomes more endothermic / less exothermic</p> | 4 |

| Question | Answer | Marks |
|-----------|---|----------|
| 7(a) | forms one or more stable ions with incomplete / partially filled d-orbitals / d-subshell | 1 |
| 7(b)(i) | purple to pale pink / colourless AND orange to green | 1 |
| 7(b)(ii) | $3\text{Sn}^{2+} + \text{Cr}_2\text{O}_7^{2-} + 14\text{H}^+ \rightarrow 3\text{Sn}^{4+} + 2\text{Cr}^{3+} + 7\text{H}_2\text{O}$ | 1 |
| 7(c)(i) | <ul style="list-style-type: none"> • six • coordinate bonds / dative bonds / lone pairs donated • to the (central) metal ion <p>award 1 mark for all three points</p> | 1 |
| 7(c)(ii) | $[\text{Ru}(\text{NH}_3)_4\text{Cl}(\text{SO}_2)]^+$ | 1 |
| 7(c)(iii) |  <p>M1: six correct ligands around Ru, bonds are shown from S of SO_2</p> <p>M2: 3-D bonds used correctly for octahedral</p> <p>M3: cis and trans isomers shown</p> | 3 |

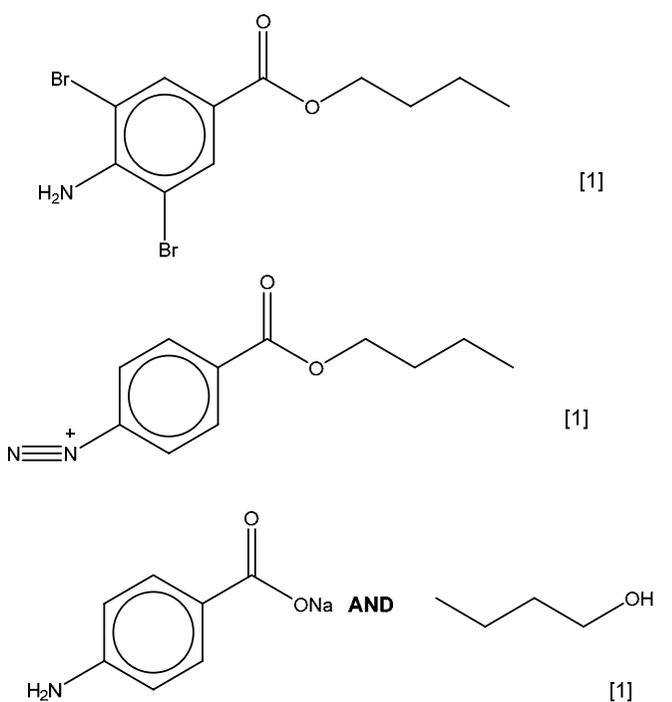
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| Question | Answer | Marks |
|----------|---|-------|
| 7(c)(iv) | cis-trans or geometric(al) [1] | 1 |
| 7(c)(v) | <p>M1: complexes have two sets of d orbital(s) of different energy OR d-orbitals splits into two sets (of orbitals)</p> <p>M2: visible light absorbed (and complementary colour observed)</p> <p>M3: electron(s) promoted / excited OR electron(s) moves to higher (d–) orbital</p> | 3 |

| Question | Answer | Marks |
|----------|---|-------|
| 8(a)(i) | <p>any one from:</p> <ul style="list-style-type: none"> • OH⁻ / NaOH; aqueous / dilute; heat under reflux • H⁺ / HCl / H₂SO₄, aqueous / dilute; heat under reflux • protease or named protease; water; T = 30° – 40°C <p>all three points in each bullet [1]</p> | 1 |
| 8(a)(ii) |  <p>M1: three amino acids in any ionic / non-ionic states [1]</p> <p>M2: three amino acids in the correct ionic state for their conditions [1]</p> | 2 |

| Question | Answer | Marks |
|----------|---|----------|
| 8(b) | <ul style="list-style-type: none"> permanent dipole-dipole one group that will have a δ^+ and another with δ^- e.g. CO, NH, COOH, OH BOTH [1] hydrogen bonds one group that will have a $H^{\delta+}$, e.g. NH, OH and another with lone pair, e.g. NH, COOH, OH, CONH₂ BOTH [1] ionic bonding NH₃⁺ and COO⁻ BOTH [1] <p>ALLOW</p> <ul style="list-style-type: none"> London forces C₄H₉ groups or parts of these alkyl groups | 3 |
| 8(c)(i) | any structure containing one COOH / COCl and NH ₂ groups in the same molecule [1] | 1 |
| 8(c)(ii) | HOCH ₂ CH ₂ OH [1] ethan(e)-1,2-diol [1] ecf for diols HO ₂ CCO ₂ H or ClOCCOCl [1] ethan(e)dioic acid or ethan(e)diol chloride [1] ecf for diacids / diacyl chlorides | 4 |

| Question | Answer | Marks |
|----------|--|----------|
| 9(a)(i) | RNH ₂ + H ⁺ → RNH ₃ ⁺ OR RNH ₂ + HCl → RNH ₃ Cl [1] | 1 |
| 9(a)(ii) | weaker AND lone pair of N delocalised into benzene ring [1] | 1 |

| Question | Answer | Marks |
|-----------|--|-------|
| 9(b) |  <p>[1]</p> <p>[1]</p> <p>[1]</p> | 3 |
| 9(c)(i) | 2 [1] | 1 |
| 9(c)(ii) | CH ₂ next to ester and terminal CH ₃ are circled [1] | 1 |
| 9(c)(iii) | <ul style="list-style-type: none"> • one less peak • the lost peak is NH₂ / aryl amine • protons exchange with D OR protons are labile OR valid equation <ul style="list-style-type: none"> • ✓✓ for two marks [2] | 2 |
| 9(d) | C ₆ H ₄ NH ₂ ⁺ and CH ₃ CH ₂ CH ₂ CH ₂ ⁺ [1] | 1 |