

Cambridge Assessment International Education

Cambridge International Advanced Subsidiary and Advanced Level

CHEMISTRY 9701/41

Paper 4 A Level Structured Questions

October/November 2018

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.

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Cambridge International AS/A Level – Mark Scheme

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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.

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GENERIC MARKING PRINCIPLE 5:

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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.

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| Question | | | Answ | er | Marks | | |
|----------|-----------------------------------|---|--|--|-------|--|--|
| 1(a)(i) | peak organic compound explanation | | | | | | |
| | х | alkane | London forces only OR no hydrogen bonding | | | | |
| | Y | aldehyde | (Permanent dipole-dipole and London forces) | | | | |
| | Z | carboxylic acid | (contains) hydrogen bonding | | | | |
| | | assignments [1] nation of Z OR X [1] | | | | | |
| 1(a)(ii) | % of Z = | 47/98 = 48% | | | 1 | | |
| 1(b)(i) | ³⁷ C <i>l</i> and | ⁸¹ Br | | | 1 | | |
| 1(b)(ii) | two corre | ${ m CH_2}^{35}{ m C}l^{79}{ m Br}$ k ${ m CH_2}^{37}{ m C}l^{79}{ m Br}$ ${ m OR}$ Cl k ${ m CH_2}^{37}{ m C}l^{81}{ m Br}$ ect scores 1 mark ect scores 2 marks | H ₂ ³⁵ C <i>l</i> ⁸¹ Br | | 2 | | |
| 1(c)(i) | H ₂ C | CH ₃ CH ₃ S+ δ- H Br | H ₃ C C+ CH ₃ CH ₃ | H ₃ C — C CH ₃ CH ₃ | 3 | | |
| | M2 two o | | any correct curly arrow [1] ws AND lone pair required on Br | ⁻ [1] | | | |

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| Question | Answer | Marks |
|----------|---|-------|
| 1(c)(ii) | (major product is) formed via the most stable tertiary carbocation / intermediate OR tertiary halogenoalkane formed via more stable carbocation / intermediate | 1 |
| 1(d)(i) | M1 ratio of the concentrations of solute in two (immiscible) solvents [1] M2 at equilibrium [1] | 2 |
| 1(d)(ii) | $K_{\text{partition}} = (x/10)/(1.25-x/50)$ [1] 4.75(1.25-x) = 5x x = 5.9375/9.75 = 0.61 g [1] correct answer scores [2] | 2 |

| Question | Answer | Marks |
|----------|---|-------|
| 2(a) | species that forms dative bond(s) to a (central) metal atom / ion | 1 |
| 2(b) | $\begin{array}{c} OH_2 \\ OH_2 \\ OH_2 \\ OH_2 \\ \end{array}$ $\begin{array}{c} OH_2 \\ OH_2 \\ OH_2 \\ \end{array}$ $\begin{array}{c} OH_2 \\ OH_2 \\ OH_2 \\ \end{array}$ $\begin{array}{c} OH_2 \\ OH_2 \\ \end{array}$ | 2 |
| 2(c)(i) | $K_{sp} = [Ca^{2+}][C_2O_4^{2-}][1]$ units mol ² dm ⁻⁶ [1] | 2 |
| 2(c)(ii) | $[Ca^{2+}] = [C_2O_4^{2-}] = 6.65 \times 10^{-3}/128.1 = 5.19 \times 10^{-5} \text{ mol dm}^{-3} [1]$ $K_{sp} = (5.19 \times 10^{-5})^2 = 2.7 \times 10^{-9} \text{ mol}^2 \text{ dm}^{-6} [1]$ | 2 |

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| Question | | | | Answer | Marks | | |
|-----------|---|--|----------------------------------|---|-------|--|--|
| 3(a) | [1] for each column | [1] for each column | | | | | |
| | | | of unpaired rons in | | | | |
| | element | 3d | 4s | | | | |
| | Cr | 5 | 1 | | | | |
| | Mn | 5 | 0 | | | | |
| | Fe | 4 | 0 | | | | |
| 3(b) | $2KMnO_4 \rightarrow K_2MnO_4$ formulae of K_2MnO_4 rest of the equation | and O ₂ [1] | nO ₂ | | 2 | | |
| 3(c) | M1 d orbitals split in M2 visible light is ab M3 electron(s) pron | sorbed and t | he complemer | per orbitals [1] htary colour observed [1] | 3 | | |
| 3(d)(i) | precipitate A [Cu(H ₂ solution B [Cu(NH ₃). solution C Cu(CH ₃ C | ₄ (H ₂ O) ₂] ²⁺ [1] | R Cu(OH) ₂ [1] | | 3 | | |
| 3(d)(ii) | Na ₂ CO ₃ or CO ₃ ²⁻ | | | | 1 | | |
| 3(d)(iii) | $CuCO_3 + 2CH_3CO_2H \rightarrow Cu(CH_3CO_2)_2 + CO_2 + H_2O$ | | | | | | |
| 3(d)(iv) | any two for one ma fizzing / bubbles solid disappea green / blue solu | / effervescen 's | | | 1 | | |

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| Question | Answer | Marks |
|----------|--|-------|
| 3(e) | sum of the charges of the (four) ligands equals the oxidation number / charge of Pt OR a calculation Pt +2, NH ₃ neutral / no charge, both Cl^{-1} s –1 (so no overall charge) | 1 |
| 3(f)(i) | CI Pt NH_3 CI I I I I I I I I I | 2 |
| 3(f)(ii) | M1 this can bond / bind with DNA [1] M2 which prevents replication of the DNA / strand OR prevents cell division [1] | 2 |
| 3(g) | H_3N Pt O H_3N O | 1 |

| Question | Answer | Marks |
|----------|--|-------|
| 4(a) | M1 solubility decreases (down the Group) [1] M2 because lattice energy and hydration energy decreases OR lattice energy and hydration energy become less exothermic / more endothermic [1] M3 because hydration energy decreases to a greater extent (than does ΔH_{Latt}) [1] | 3 |
| 4(b)(i) | $(K_{w} =) [H^{+}][OH^{-}]$ | 1 |

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| Question | | | | Answer | | Marks | |
|----------|---|-----------------------------------|---|-------------------------------|---|-------|--|
| 4(b)(ii) | [1] or each correct tick | | | | | | |
| | effect of increasing temperature | decreases | stay the same | increase | | | |
| | рН | ✓ | | | | | |
| | ratio of [H ⁺]:[OH ⁻] | | ✓ | | | | |
| 4(c) | $[H^{+}] = 10^{-13.25} = 5.62 \times 1$ $[OH^{-}] = K_{w}/[H^{+}] = 1.0 \times 1$ $[OH^{-}] = 0.18 (0.178) (mag)$ | $10^{-14}/5.62 \times 10^{-14}$ | r ^{–14} of correct answer s | cores [2] | | 2 | |
| 4(d) | $HCO_3^- + H^+ \rightarrow H_2CO_3$ OR $HCO_3^- + H^+ \rightarrow CO_2 + H_2O$ [1] $H_2CO_3 + OH^- \rightarrow HCO_3^- + H_2O$ [1] | | | | | 2 | |
| 4(e)(i) | CH_3COOH + H_2O \rightleftharpoons acid + base \rightleftharpoons base | • | H ₃ O ⁺ [1] | | | 2 | |
| 4(e)(ii) | M1 moles NaOH = 0.15 | × 20/1000 = 0 | 0.0030 AND initial | moles CH ₃ C0 | OOH = 0.25 × 30/1000 OR 0.0075 [1] | 4 | |
| | M2 equilibrium moles C | H ₃ COOH = 0. 0 | 0045 AND equilibri | um moles CH | I ₃ COONa = 0.0030 [1] | | |
| | M3 [CH ₃ COOH] = 0.004 [H ⁺] = K_a | | O AND [CH3COON /[CH3COONa] = 2 | | | | |
| | M4 pH = $-\log[H^+]$ = 4.6 | [1] correct an | swer scores [4] | | | | |
| 4(f)(i) | end point = 28 cm ³ | | | | | 1 | |
| 4(f)(ii) | M1 reaction M bromoth | ıymol (blue)/b | romocresol (green |) AND reaction | on N bromothymol (blue) / thymolphthalein [1] | 2 | |
| | M2 (both indicators hav | e) a pH range | / colour change wi | thin / in end-ր | point / vertical region / sharp fall of the graph [1] | | |

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| Question | | | | Answer | | Marks | | |
|----------|--------------------------------------|--------------|-------------------|----------------------------------|--|-------|--|--|
| 5(a)(i) | [1] for each correct answer | | | | | | | |
| | number of p | eaks | | | | | | |
| | F 3 | | | | | | | |
| | G 6 | | | | | | | |
| 5(a)(ii) | one amide bond displayed in full [1] | | | | | 2 | | |
| | rest of the structure – c | | y [1] | | | | | |
| 5(b) | | | | | | 2 | | |
| | | σ-bonds only | π -bonds only | both σ - and π -bonds | | | | |
| | bonds broken | | ✓ | | | | | |
| | bonds formed | ✓ | | | | | | |

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| Question | Answer | Marks |
|----------|---|-------|
| 5(c) | C ₆ H ₅ H CH ₃ H CH ₄ C C C C C C C C C C C C C C C C C C C | 2 |
| 5(d)(i) | C-C bonds are non-polar / have no dipole so cannot be hydrolysed [1] | 1 |
| 5(d)(ii) | M1 Hydrolysis using acid / base / alkali / enzymes [1] M2 action of UV light [1] | 2 |

| Question | Answer | Marks |
|----------|--|-------|
| 6(a) | $\begin{array}{c ccccccccccccccccccccccccccccccccccc$ | 2 |
| | M1 amide bond displayed [1] M2 rest of the structure [1] | |

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| Question | Answer | Marks |
|----------|--|-------|
| 6(b) | mixture applied here M1 valine on the cross [1] M2 Val-Lys and Lys on the right of the cross (in any order) [1] M3 relative order of Val-Lys and Lys (on the same side of the cross) [1] Explanation Val does not move as it is a zwitterion / neutral (at pH6) OR Lys / Val-Lys move towards negative (pole) as they are positively charged Lys moves the furthest as it has the lowest M _r (with the same positive charge) OR Val-Lys moves the least as it has the largest M _r (with the same positive charge) [1] × 2 | 5 |

| Question | Answer | Marks |
|----------|---|-------|
| 7(a) | M1 C-X/C-C l /C-O bond is stronger (in chlorobenzene/phenol) [1] M2 p-orbital/lone pair on C l /O(H)/X (in chlorobenzene/phenol) [1] M3 electrons of the (C l /O/electronegative atom) AND overlap/delocalise with π -electron cloud/delocalise into ring [1] | 3 |

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| Question | Answer | | | |
|-----------|---|---|--|--|
| 7(b) | OH 2-bromophenol structure and name correct [1] | 2 | | |
| 7(c)(i) | step 1 conc . HNO ₃ + H ₂ SO ₄ (and temperare 50–55 °C) [1] step 2 Sn + HC <i>l</i> AND one of conc.HC <i>l</i> + heat [1] step 4 H ₂ O warm / heat [1] | 3 | | |
| 7(c)(ii) | N+ CIT OR CIT | 1 | | |
| 7(c)(iii) | step 1 electrophilic substitution | 1 | | |
| 7(c)(iv) | $C_6H_5NO_2 + 6[H] \rightarrow C_6H_5NH_2 + 2H_2O$ | 1 | | |

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| Question | Answer | | | | |
|----------|--|------------------------------|-------------------------|---|---|
| 8(a) | M1 continuous increase in S | rature/K S from 0–300 K (| | [1] inuous increase in <i>S</i> after m.p. [1] | 2 |
| 8(b) | [1] for each correct tick | | | 1 | |
| | | negative ∆S ^e | positive ΔS^{e} | | |
| | solid dissolving in water | | ✓ | | |
| | water boiling to steam | | ✓ | | |
| 8(c) | $\Delta H^{\text{e}} = (2 \times \text{C=O}) + (3 \times \text{H-H}) - (3 \times \text{C-H}) - (\text{C-O}) - (3\text{xO-H})$ $\Delta H^{\text{e}} = (2 \times 805) + (3 \times 436) - (3 \times 410) - (1 \times 360) - (3 \times 460)$ [1] $\Delta H^{\text{e}} = 1610 + 1308 - 1230 - 360 - 1380 = -52$ (kJ mol ⁻¹) [1] ecf correct answer scores [2] | | | | |
| 8(d)(i) | $\Delta S^{0} = 127 + 70 - (214 + 3 \times 131) [1]$ = -410 (J K ⁻¹ mol ⁻¹) [1] ecf correct answer scores [2] | | | | |
| 8(d)(ii) | $\Delta G^{\circ} = \Delta H^{\circ} - T\Delta S^{\circ}$ [1] $\Delta G^{\circ} = -131 - (298 \times -0.41) = -8.8(2)$ (kJ mol ⁻¹) [1] correct answer scores [2] | | | | |

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| Question | Answer | Marks |
|-----------|--|-------|
| 8(d)(iii) | (as temperature increases) feasibility decreases | 1 |
| 8(e)(i) | $2CH_3OH + 3O_2 \Rightarrow 2CO_2 + 4H_2O$ OR $2CH_3OH + 3O_2 \Rightarrow 2CO_2 + 4H^+ + 4OH^-$ | 1 |
| 8(e)(ii) | $E_{\text{cell}}^{\circ} = 1.23 - 0.02 = 1.21 \text{ V}$ | 1 |

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