Cambridge Assessment International Education

# Cambridge International AS & A Level

**CHEMISTRY** **9701/42**

Paper 4 A Level Structured Questions **October/November 2021**

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 2021 series for most Cambridge IGCSE™, Cambridge International A and AS Level components and some Cambridge O Level components.



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**[Turn over**

**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.

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|---|
| GENERIC MARKING PRINCIPLE 5:<br><br>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:<br><br>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. |

**Science-Specific Marking Principles**

| | |
|---|---|
| 1 | Examiners should consider the context and scientific use of any keywords when awarding marks. Although keywords may be present, marks should not be awarded if the keywords are used incorrectly. |
| 2 | The examiner should not choose between contradictory statements given in the same question part, and credit should not be awarded for any correct statement that is contradicted within the same question part. Wrong science that is irrelevant to the question should be ignored. |
| 3 | Although spellings do not have to be correct, spellings of syllabus terms must allow for clear and unambiguous separation from other syllabus terms with which they may be confused (e.g. ethane / ethene, glucagon / glycogen, refraction / reflection). |
| 4 | The error carried forward (ecf) principle should be applied, where appropriate. If an incorrect answer is subsequently used in a scientifically correct way, the candidate should be awarded these subsequent marking points. Further guidance will be included in the mark scheme where necessary and any exceptions to this general principle will be noted. |
| 5 | ‘List rule’ guidance<br><br>For questions that require ***n*** responses (e.g. State **two** reasons …):<br><br>• The response should be read as continuous prose, even when numbered answer spaces are provided.<br>• Any response marked *ignore* in the mark scheme should not count towards ***n***.<br>• Incorrect responses should not be awarded credit but will still count towards ***n***.<br>• Read the entire response to check for any responses that contradict those that would otherwise be credited. Credit should **not** be awarded for any responses that are contradicted within the rest of the response. Where two responses contradict one another, this should be treated as a single incorrect response.<br>• Non-contradictory responses after the first ***n*** responses may be ignored even if they include incorrect science. |

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6 Calculation specific guidance

Correct answers to calculations should be given full credit even if there is no working or incorrect working, **unless** the question states 'show your working'.

For questions in which the number of significant figures required is not stated, credit should be awarded for correct answers when rounded by the examiner to the number of significant figures given in the mark scheme. This may not apply to measured values.

For answers given in standard form (e.g. $a \times 10^n$) in which the convention of restricting the value of the coefficient ($a$) to a value between 1 and 10 is not followed, credit may still be awarded if the answer can be converted to the answer given in the mark scheme.

Unless a separate mark is given for a unit, a missing or incorrect unit will normally mean that the final calculation mark is not awarded. Exceptions to this general principle will be noted in the mark scheme.

7 Guidance for chemical equations

Multiples / fractions of coefficients used in chemical equations are acceptable unless stated otherwise in the mark scheme.

State symbols given in an equation should be ignored unless asked for in the question or stated otherwise in the mark scheme.

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| Question | Answer | Marks |
|---|---|---|
| 1(a) | • enthalpy / energy change / given out / evolved / released<br>• one mole is formed / made [1]<br><br>• of compound / solid / lattice / crystal<br>• (from) **gas**eous **ions** [1] | **2** |
| 1(b)b | $S^-(g) + e^- \rightarrow S^{2-}(g)$ [1] | **1** |
| 1(c)c | (555 + 200 – 532 = 223, 223 × 8 = 1784)<br><br>M1 selecting correct data 555, 200, 532 **only**, (ignore signs and multipliers) [1]<br><br>M2 evaluation to give +223 [1]<br><br>M3 multiplying M2 by 8 and evaluation ans (+) 1784 [1] | **3** |
| 1(d) | (1619 + 555 – 2612 = –438)<br><br>M1 selecting correct data 1619 555 2612 only, (ignore signs and multipliers) [1]<br><br>M2 evaluation to give –438 [1] | **2** |
| 1(e)(i) | ionic radius / size / sum of ionic radii [1]<br><br>ionic charge / product of ionic charges [1] | **2** |
| 1(e)(ii) | M1 (size tends to make $\Delta H^{\ominus}_{latt}$ of radium sulfide) less exothermic since the ions are larger [1]<br><br>M2 (charge tends to make $\Delta H^{\ominus}_{latt}$ of radium sulfide) less exothermic since the ions are more highly charged [1] | **2** |
| 1(e)(iii) | (ionic) charge (since)<br><br>**AND**<br><br>$\Delta H^{\ominus}_{latt}$ of radium sulfide is more exothermic [1] | **1** |

| Question | Answer | Marks |
|---|---|---|
| 2(a) | M1 reference to **equilibrium** not being established / not at **equilibrium** [1]<br><br>M2 concentration in water too high **or** conc in octan-1-ol too low owtte [1] | **2** |
| 2(b)(i) | 1.1 / 1.13 / 1.127 / 1.1272 [1] | **1** |
| 2(b)(ii) | M1 $6.760 = (0.762 - x) / x$ [1]<br><br>M2 $7.76x = 0.762$<br>$x = 0.0982$ g (correct value scores 2) [1]<br><br>M3 0.0982 / 74 = 0.0013 / 0.00133 mol [1] | **3** |
| 2(c)(i) | $1.0 \times 10^{-9}$ [1]<br><br>$mol^3\ dm^{-9}$ [1] | **2** |
| 2(c)(ii) | yellow precipitate / yellow solid / yellow crystals [1]<br><br>common ion effect / $K_{sp}$ has been exceeded [1] | **2** |

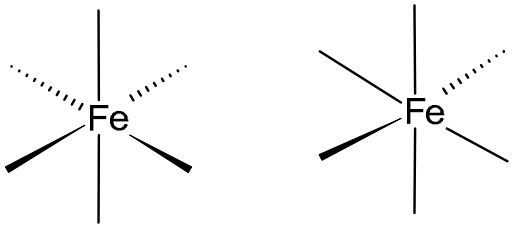
| Question | Answer | Marks |
|---|---|---|
| 3(a) | 0.351 / 24 = 0.015 (mol) [1]<br><br>$0.015 \times 6.02 \times 10^{23} = 9.0 \times 10^{21} / 8.8 \times 10^{21}$ [1] | **2** |
| 3(b) | $1.76 \times 10^{22} / 1.8 \times 10^{22}$ [1] | **1** |
| 3(c) | 2817 / 2816 / 2820 / 2800 C [1] | **1** |
| 3(d) | 15 / 15.1 / 15.05 / 15.15 minutes [1] | **1** |
| 3(e)(i) | –20 [1] | **1** |

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| Question | Answer | Marks |
|---|---|---|
| 3(e)(ii) | states / uses correct Gibbs equation [1]<br>answer = 190 / 191 / 190.0 [1] | **2** |
| 3(e)(iii) | Becomes less feasible / less spontaneous /<br>**AND**<br>because $\Delta S$ is negative / $T\Delta S$ becomes more negative / $-T\Delta S$ becomes more positive [1] | **1** |

| Question | Answer | Marks |
|---|---|---|
| 4(a) | strontium nitrate **AND** because of larger cationic radius [1]<br>$NO_3^-$ / nitrate **ion** / anion is less distorted / polarised<br>**OR**<br>N–O or N=O bond less weakened / distorted [1] | **2** |
| 4(b) | M1 magnesium oxide requires more water to dissolve **AND** because $Sr(OH)_2$ is more soluble [1]<br>M2 lattice and hydration energies less for $Sr(OH)_2$ [1]<br>M3 lattice energy is dominant factor / change in lattice energy is greater [1] | **3** |
| 4(c)(i) | 0.13 [1] | **1** |
| 4(c)(ii) | 26.7 $cm^3$ **must be 3SF** [1] | **1** |

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| Question | Answer | Marks |
|---|---|---|
| 5(a)(i) | donates **one lone pair** / forms **one** dative bond **AND** to **metal atom** / to **metal ion** [1] | **1** |
| 5(a)(ii) | $Mo(CO)_6$ 0 / no charge / neutral / zero<br>$Fe(CN)_6$ 3–<br><br>both formulae [1]<br>both charges [1] | **2** |
| 5(a)(iii) | either of these renditions, with a CN on each bond:<br><br><br><br><br>**AND**<br><br>180° bond angle between two opposite bonds [1] | **1** |
| 5(b)(i) | $[Cu(H_2O)_6]^{2+} + 4NH_3 \rightarrow [Cu(NH_3)_4(H_2O)_2]^{2+} + 4H_2O$<br><br>**OR**<br><br>$Cu^{2+} + 2H_2O + 4NH_3 \rightarrow [Cu(NH_3)_4(H_2O)_2]^{2+}$ [1] | **1** |
| 5(b)(ii) | M1 **d**-orbitals split (could be in diagram) [1]<br><br>M2 electrons transition to higher orbital / electrons promoted or excited [1]<br><br>M3 wavelength / frequency / light / photon / hf / hn absorbed [1]<br><br>M4 colour seen / reflected / transmitted is complement of colour absorbed [1] | **4** |

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| Question | Answer | Marks |
|---|---|---|
| 5(c)(i) | $[Cu(NH_3)_4(H_2O)_2]^{2+} + 4HCl \rightarrow CuCl_4^{2-} + 2H_2O + 4NH_4^+$<br>**OR**<br>$[Cu(NH_3)_4(H_2O)_2]^{2+} + 4Cl^- (+ 4H^+) \rightarrow CuCl_4^{2-} + 2H_2O + 4NH_3 (+ 4H^+)$<br>M1 formula and charge of $CuCl_4^{2-}$ [1]<br>M2 rest of equation [1] | **2** |
| 5(c)(ii) | **ligand** exchange / substitution / replacement / displacement [1] | **1** |
| 5(c)(iii) | tetrahedral [1] | **1** |
| 5(c)(iv) | yellow [1] | **1** |
| 5(c)(v) | **d**-orbital splitting changes / **ΔE** changes [1] | **1** |

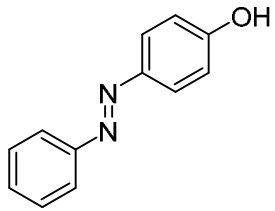
| Question | Answer | Marks |
|---|---|---|
| 6(a) | red or pink to yellow / straw / brown / orange [1] | **1** |
| 6(b) | $[[Co(NH_3)_6]^{2+}] / [[Co(H_2O)_6]^{2+}][NH_3]^6$ [1] | **1** |
| 6(c) | $Co(NH_3)_6^{2+}$ is more stable (than $Co(H_2O)_6^{2+}$) [1] | **1** |
| 6(d)(i) | $E^{\ominus}_{cell}$ = + 0.30 [1] | **1** |
| 6(d)(ii) | $4[Co(NH_3)_6]^{2+} + O_2 + 2H_2O \rightleftharpoons 4[Co(NH_3)_6]^{3+} + 4OH^-$ [1] | **1** |
| 6(d)(iii) | M1 No, because 1.82 V **and** 1.23 V [1]<br>M2 $E^{\ominus}_{cell}$ = –0.59 / 1.23 lower than 1.82 / 1.82 greater than 1.2 [1] | **2** |

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| Question | Answer | Marks |
|---|---|---|
| 7(a) | $C_4H_3O_2$ [1] | **1** |
| 7(b) | benzene-1,3-dicarboxylic acid \| 5 [1]<br>benzene-1,4-dicarboxylic acid \| 3 [1] | **2** |
| 7(c)(i) | benzoic acid [1] | **1** |
| 7(c)(ii) | COOH directs 3 position [1] | **1** |
| 7(c)(iii) | electrophilic substitution [1] | **1** |
| 7(c)(iv) | M1 curly arrow from within hexagon towards $CH_3C^+=O$ [1]<br>M2 correct intermediate [1]<br>M3 curly arrow from C–H bond into hexagon and correct product Q [1] | **3** |
| 7(c)(v) | $MnO_4^-$ / $KMnO_4$ / manganate$\mathrm{VII}$ / permanganate aq / $H^+$ / acidified / $OH^-$ then acid / alkaline then acid heat / boil / reflux / T>50$^o$<br>**OR**<br>alkaline iodine followed by acidification [1] | **1** |

| Question | Answer | Marks |
|---|---|---|
| 8(a) | M1 one diagram correct [1]<br>M2 both diagrams correct 3D and different [1] | **2** |

| Question | Answer | Marks |
|---|---|---|
| 8(b) | alanine because glutamic acid would have more than two / three peaks / absorb$^{ns}$ / proton environments [1]<br><br>reason why alanine has a **doublet** given as one neighbouring proton [1]<br>glutamic acid would have (two) **triplet**(s) **OR** a **multiplet** [1]<br><br>reason why alanine has a **quartet / quadruplet** given as three neighbouring protons [1] | **3** |
| 8(c)(i) | 147 [1] | **1** |
| 8(c)(ii) | 88 = $[H_2NCH(CH_2)CO_2H]+$ [1]<br>131 = $[CH(CH_2CH_2CO_2H)CO_2H]+$ [1] | **2** |
| 8(d)(i) | use of buffer [1] | **1** |
| 8(d)(ii) | • correct circuit including **DC** power supply<br>• paper or gel labelled [1]<br>• sample towards the middle of the paper / gel<br><br>**OR**<br><br>on cathode side [1] | **2** |
| 8(d)(iii) | anode / positive / +<br><br>**AND**<br><br>anode / positive / + [1] | **1** |
| 8(d)(iv) | M1 ala is –1 and glu is –2 [1]<br><br>M2 ala is lighter / has lower $M_r$ [1] | **2** |

<table>
<tr><th>Question</th><th>Answer</th><th>Marks</th></tr>
<tr><td>9(a)</td><td>
<table>
<tr><td>organic starting material</td><td>reagent and conditions</td></tr>
<tr><td>1-butyl halide, e.g. $CH_3CH_2CH_2CH_2Br$</td><td>$NH_3$ under pressure or heated in sealed tube</td></tr>
<tr><td>butanenitrile $CH_3CH_2CH_2CN$</td><td>$H_2$ and Ni or Pt / $LiAlH_4$ / Na + ethanol</td></tr>
<tr><td>butanamide $CH_3CH_2CH_2CONH_2$</td><td>$LiAlH_4$ / Na + ethanol</td></tr>
</table>
[1] + [1]<br>[1] + [1]<br>[1] + [1]
</td><td>6</td></tr>
<tr><td>9(b)</td><td>M1 butylamine > ammonia > phenylamine [1]<br><br>M2 basicity related to ability of lp to accept proton / $H^+$ [1]<br><br>M3 butylamine is stronger because of positive inductive effect of alkyl group / $C_4H_9$ [1]<br><br>M4 phenylamine is weaker because lp on N is delocalised into ring [1]</td><td>4</td></tr>
<tr><td>10(a)</td><td>$C_6H_5O^-$ / $C_6H_5O^-Na^+$ / $C_6H_5ONa$ [1]<br><br>$C_6H_5O^-$ / $C_6H_5O^-Na^+$ / $C_6H_5ONa$ [1]<br><br>$C_6H_5N_2C_6H_4O^-$<br>or<br>$C_6H_5N_2C_6H_4OH$ or <br>[1]<br><br>$C_6H_5OCOC_6H_4CO_2C_6H_5$ [1]</td><td>4</td></tr>
<tr><td>10(b)</td><td>2 and 4 [1]</td><td>1</td></tr>
</table>