



Cambridge International AS & A Level

CHEMISTRY

9701/43

Paper 4 A Level Structured Questions

October/November 2020

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.

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This document consists of **15** 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.

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
 For questions that require *n* responses (e.g. State **two** reasons ...):
 - The response should be read as continuous prose, even when numbered answer spaces are provided.
 - Any response marked *ignore* in the mark scheme should not count towards *n*.
 - Incorrect responses should not be awarded credit but will still count towards *n*.
 - 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.
 - Non-contradictory responses after the first *n* responses may be ignored even if they include incorrect science.

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.

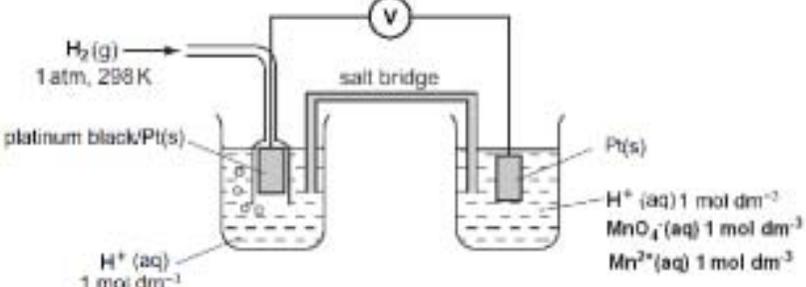
7 Guidance for chemical equations

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.

| Question | Answer | Marks | | | | | | |
|---|---|--|---|---|---|-------------------------------|---|----------|
| 1(a) | <table border="1" data-bbox="763 215 1514 411"> <tr> <td>the order of reaction with respect to [NO]</td> <td>2</td> </tr> <tr> <td>the order of reaction with respect to [O₂]</td> <td>1</td> </tr> <tr> <td>the overall order of reaction</td> <td>3</td> </tr> </table> <p>ALL CORRECT [1]</p> | the order of reaction with respect to [NO] | 2 | the order of reaction with respect to [O ₂] | 1 | the overall order of reaction | 3 | 1 |
| the order of reaction with respect to [NO] | 2 | | | | | | | |
| the order of reaction with respect to [O ₂] | 1 | | | | | | | |
| the overall order of reaction | 3 | | | | | | | |
| 1(b)(i) | $k = (1.51 \times 10^{-4}) / (0.003^2 \times 0.00200)$ $k = \mathbf{8389}$ [1] min 2sf $\text{mol}^{-2} \text{dm}^6 \text{s}^{-1}$ [1] | 2 | | | | | | |
| 1(b)(ii) | $8400 = (6.05 \times 10^{-5}) / (x^2 \times 0.005)$ $x = \sqrt{(6.05 \times 10^{-5}) / (8400 \times 0.005)}$ $x = \mathbf{0.00120 / 1.20 \times 10^{-3}}$ [1] min 2sf ecf from Q1bi | 1 | | | | | | |
| 1(c) | slow(est) [1] | 1 | | | | | | |
| 1(d)(i) | <p>correct RDS identified as step 1 with only one S₂O₈²⁻ and one I⁻ [1]</p> <p>overall mechanism adds up to chemical equation and no cancellable species on LHS / RHS in each of the equations [1]</p> <p>M2 DEP on one S₂O₈²⁻ and one I⁻ in step 1</p> <p>e.g. step 1 S₂O₈²⁻ + I⁻ → SO₄²⁻ + SO₄I⁻ RDS = step 1</p> <p>step 2 SO₄I⁻ + I⁻ → SO₄²⁻ + I₂</p> | 2 | | | | | | |
| 1(d)(ii) | no. of $t_{1/2} = 192 / 48 = 4$ $[I^-] = 0.0078 / 16 = \mathbf{4.9 \times 10^{-4}}$ [1] min 2sf | 1 | | | | | | |

| Question | Answer | Marks |
|-----------|--|----------|
| 2(a)(i) | M1 energy released when 1 mole of an ionic compound is formed [1] M2 from gaseous ions (under standard conditions) [1] | 2 |
| 2(a)(ii) | Ca ²⁺ & O ²⁻ have a higher charge / charge density (than Li ⁺ and F ⁻) [1] | 1 |
| 2(a)(iii) | MgO –3600 or more negative AND BaO –3200 or less negative BOTH [1] | 1 |
| 2(b)(i) | BaO(s) + H ₂ O(l) → Ba(OH) ₂ (aq) [1] | 1 |
| 2(b)(ii) | M1 (solubility) increases (down the group) [1] M2 both ΔH_{latt} and ΔH_{hyd} become less exothermic / less negative [1] M3 ΔH_{latt} changes more / is dominant factor [1] M4 ΔH_{sol} becomes more negative / more exothermic [1] | 4 |
| 2(c) | M1: Use of 2×-348 (EA F) and +158 (bond energy of F ₂) [1] M2: Use of +147 (at Mg) and +736 and +1450 (IEs of Mg) [1] M3: evaluation and calculation of their answer (–1102 – (147 + 158 + 736 + 1450 – 696)) = –2897 (kJ mol ^{–1}) [1] ecf | 3 |
| 2(d)(i) | <ul style="list-style-type: none"> • (energy change) when an / one electron is added to • each atom / ion in one mole of • gaseous atoms / ions mark as • ✓ ✓ [2] | 2 |
| 2(d)(ii) | F has greater nuclear charge / more protons AND greater attraction between F atom / nucleus and the electrons • ✓ BOTH [1] | 1 |

| Question | Answer | Marks |
|----------|---|----------|
| 3(a) | (anode =) oxygen / O ₂ AND (cathode =) hydrogen/H ₂ BOTH [1] | 1 |
| 3(b) | <p>M1: $Q = 1.5 \times 60 \times 60 \times 4.5 = 24300$ (C) [1]</p> <p>M2: no. of F / moles of e⁻ = $24300 / 96500 = 0.25(1813)$ [1] ecf</p> <p>M3: volume of Cl₂ = $24 \times 0.252 / 2 = 3.02$ dm³ [1] ecf min 2sf</p> <p>M4: mass of Na = $0.252 \times 23 = 5.79$ (5.7917) g Na [1] ecf min 2sf</p> | 4 |
| 3(c)(i) | <div style="text-align: center;">  </div> <ul style="list-style-type: none"> • MnO₄⁻, H⁺, Mn²⁺ in same beaker AND H⁺ in other beaker • both electrodes Pt(s) (ALLOW graphite) • one solute clearly identified as 1M / 1 mol dm⁻³ • 298 K OR 1 atm • voltmeter / potentiometer labelled (or circled V) • salt bridge labelled (must touch the solution) • a good delivery system for H₂(g) • H₂(g) <p style="text-align: center;">mark as two correct points = 1 mark [4]</p> | 4 |

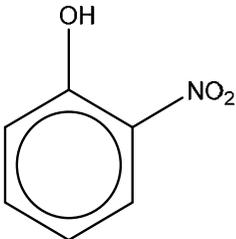
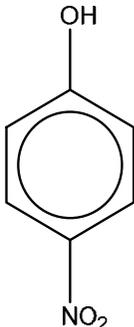
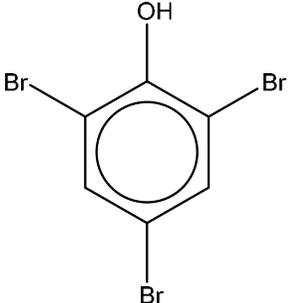
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| Question | Answer | Marks |
|----------|--|-------|
| 3(c)(ii) | <p>F_2 OR $S_2O_8^{2-}$ OR H_2O_2 OR $HOC\ell$ OR Co^{3+} OR Pb^{4+} [1]</p> <p>$2Mn^{2+} + 8H_2O + 5F_2 \rightarrow 2MnO_4^- + 16H^+ + 10F^-$ [1] OR $2Mn^{2+} + 5S_2O_8^{2-} + 8H_2O \rightarrow 2MnO_4^- + 16H^+ + 10SO_4^{2-}$ OR $Mn^{2+} + 4H_2O + 5Co^{3+} \rightarrow MnO_4^- + 8H^+ + 5Co^{2+}$ OR $2Mn^{2+} + 8H_2O + 5Pb^{4+} \rightarrow 2MnO_4^- + 16H^+ + 5Pb^{2+}$ OR $2Mn^{2+} + 5H_2O_2 \rightarrow 2MnO_4^- + 6H^+ + 2H_2O$ OR $2Mn^{2+} + 10HOC\ell \rightarrow 2MnO_4^- + 6H^+ + 5Cl_2 + 2H_2O$</p> | 2 |

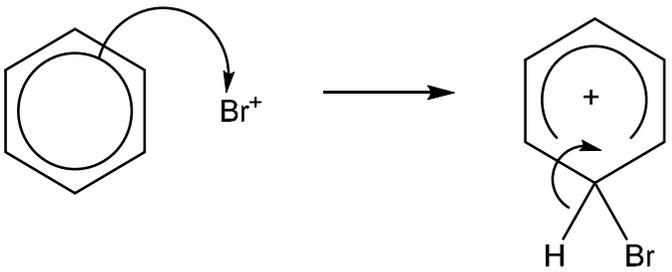
| Question | Answer | Marks |
|----------|---|-------|
| 4(a)(i) | <p>(pH =) $-\log[H^+]$ OR $-\lg[H^+]$ [1]</p> <p>(K_w =) $[H^+][OH^-]$ [1]</p> | 2 |
| 4(a)(ii) | <p>$[H^+] = 1 \times 10^{-14} / 0.027 = 3.7037 \times 10^{-13}$</p> <p>pH = $-\log(3.7037 \times 10^{-13}) = 12.4$ [1] min 3sf</p> | 1 |
| 4(b) | <p>$[H^+] = \sqrt{3.72 \times 10^{-8} \times 0.010} = 1.9287 \times 10^{-5}$</p> <p>pH = $-\log(1.9287 \times 10^{-5}) = 4.7$ [1] min 2sf</p> | 1 |
| 4(c)(i) | <p>$K_{pc} = (0.935 / 50) / (0.065 / 50)$</p> <p>$K_{pc} = 14.4$ (14.38) [1] min 3sf</p> | 1 |
| 4(c)(ii) | <p>M1: $14.4 = ((0.935 - x) / 50) / (x / 100)$ [1] ecf from 4(c)(i)</p> <p>M2: $x = 0.114$ g [1] min 2sf ecf from M1</p> | 2 |

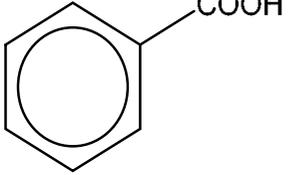
| Question | Answer | Marks | | | | | | | | | | | | |
|---------------------|---|--|---|---|-------------------|--------|------------------|---------------------|-------------|------------|--------------------|--|--|---|
| 5(a)(i) | $K_{stab} = \frac{[(Cu(NH_3)_4]^{2+}]}{[(Cu(H_2O)_6]^{2+}] [NH_3]^4}$ [1] | 1 | | | | | | | | | | | | |
| 5(a)(ii) | deep / dark / royal blue [1] | 1 | | | | | | | | | | | | |
| 5(b) | $[Cu(NH_3)_4]^{2+} + 2H_2O \rightarrow Cu(OH)_2 + 2NH_4^+ + 2NH_3$ [1] OR $[Cu(NH_3)_4]^{2+} + 2H_2O \rightarrow Cu(OH)_2 + 2H^+ + 4NH_3$ | 1 | | | | | | | | | | | | |
| 5(c) | $Cu(OH)_2 + 4HCl \rightarrow [CuCl_4]^{2-} + 2H_2O + 2H^+$ OR $Cu(OH)_2 + 4Cl^- + 2H^+ \rightarrow [CuCl_4]^{2-} + 2H_2O$ [CuCl ₄] ²⁻ complex including charge [1] rest of equation fully correct [1] | 2 | | | | | | | | | | | | |
| 5(d) | <table border="1" style="margin-left: auto; margin-right: auto;"> <thead> <tr> <th></th> <th>Y</th> <th>Z</th> </tr> </thead> <tbody> <tr> <td>colour of complex</td> <td>yellow</td> <td>blue / pale blue</td> </tr> <tr> <td>geometry of complex</td> <td>tetrahedral</td> <td>octahedral</td> </tr> <tr> <td>formula of complex</td> <td style="background-color: #cccccc;"></td> <td>[Cu(H₂O)₆]²⁺</td> </tr> </tbody> </table> <p>one mark for any three cells [1] ••✓ two marks for all five cells [2] ••✓•✓</p> | | Y | Z | colour of complex | yellow | blue / pale blue | geometry of complex | tetrahedral | octahedral | formula of complex | | [Cu(H ₂ O) ₆] ²⁺ | 2 |
| | Y | Z | | | | | | | | | | | | |
| colour of complex | yellow | blue / pale blue | | | | | | | | | | | | |
| geometry of complex | tetrahedral | octahedral | | | | | | | | | | | | |
| formula of complex | | [Cu(H ₂ O) ₆] ²⁺ | | | | | | | | | | | | |
| 5(e) | M1: d orbitals splits into two sets of energy levels of different energy [1] M2: wavelength / frequency / light / photon / hv absorbed [1] M3: electron(s) promoted / excited [1] M4: colour seen is complementary (to colour absorbed) [1] M5: d–d energy gap / ΔE is different for Y and Z AND so different frequency / wavelength of light absorbed • ✓ [1] | 5 | | | | | | | | | | | | |

| Question | Answer | Marks |
|-----------|--|-------|
| 6(a)(i) | (a species) that donates <u>one</u> lone pair [1] to form a dative / coordinate to a central metal atom / metal ion [1] | 2 |
| 6(a)(ii) | $[\text{Ag}(\text{S}_2\text{O}_3)_2]^{3-}$ [1] | 1 |
| 6(b)(i) | $[\text{Ag}(\text{S}_2\text{O}_3)_2]^{3-} + 2\text{CN}^- \rightarrow [\text{Ag}(\text{CN})_2]^- + 2\text{S}_2\text{O}_3^{2-}$ [1] OR $\text{Ag}(\text{S}_2\text{O}_3)_2^{3-} + 2\text{NaCN} \rightarrow [\text{Ag}(\text{CN})_2]^- + \text{Na}_2\text{S}_2\text{O}_3 + \text{S}_2\text{O}_3^{2-}$ | 1 |
| 6(b)(ii) | Q is more stable / has a larger K_{stab} than P [1] | 1 |
| 6(b)(iii) | ligand exchange / displacement / substitution | 1 |
| 6(c)(i) |  <p>both correct [1]</p> | 1 |
| 6(c)(ii) | square planar [1] | 1 |
| 6(c)(iii) | cis-trans OR geometric(al) [1] | 1 |

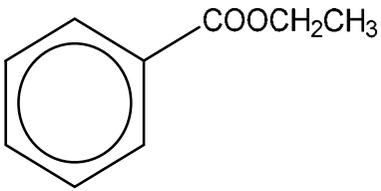
| Question | Answer | Marks |
|----------|--|----------|
| 7(a) | M1: HNO_2 OR $\text{NaNO}_2 + \text{HCl}$ [1] M2: $T \geq 10^\circ\text{C}$ / warm AND water [1] | 2 |
| 7(b) | <div style="display: flex; justify-content: space-around; align-items: center;"> <div style="text-align: center;">  <p>2-nitrophenol</p> </div> <div style="text-align: center;">  <p>4-nitrophenol</p> </div> </div> <p>2 × [1]</p> | 2 |
| 7(c)(i) | <div style="text-align: center;">  <p>2,4,6-tribromophenol</p> </div> <p>✓ ✓ [2]</p> | 2 |
| 7(c)(ii) | bromine is decolourised AND white precipitate is formed BOTH [1] | 1 |
| 7(d) | $\text{C}_6\text{H}_5\text{OH} + \text{NaOH} \rightarrow \text{C}_6\text{H}_5\text{ONa} + \text{H}_2\text{O}$ [1] ALLOW any equation for phenol acting as an acid | 1 |

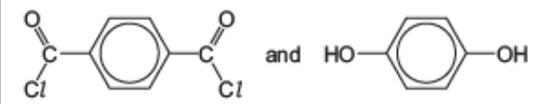
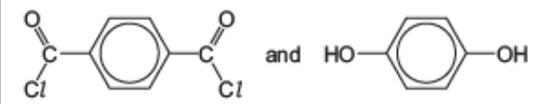
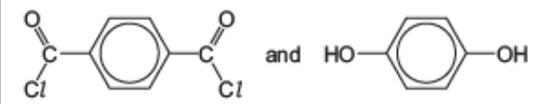
| Question | Answer | Marks |
|----------|--|-------|
| 7(e) | phenol>water>ethanol [1] <ul style="list-style-type: none"> (phenol:) lone pair on oxygen is delocalised into the benzene ring (ethanol:) positive inductive effect / electron donating effect of alkyl / ethyl group correct statement about stabilisation of anion/ conjugate base OR weakening of O-H bonds once <i>in the context of phenol / ethanol</i> correct statement about ease of proton/H⁺ donation <i>in the context of phenol / ethanol</i> [2] Two correct statements = 1 mark | 3 |

| Question | Answer | Marks |
|-----------|--|-------|
| 8(a)(i) | HBr / hydrogen bromide [1] | 1 |
| 8(a)(ii) |  <p>M1 curly arrow to Br⁺ AND curly arrow from C–H bond as shown [1] M2 correct intermediate [1]</p> | 2 |
| 8(a)(iii) | electrophilic substitution [1] | 1 |
| 8(b)(i) | reagent: chloroethane / bromoethane / iodoethane OR formula [1] catalyst: FeCl ₃ / AlCl ₃ etc. [1] | 2 |

| Question | Answer | Marks |
|-----------|---|-------|
| 8(b)(ii) |  <p>[1] ALLOW C₆H₅COONa</p> | 1 |
| 8(b)(iii) | step 3 = LiAlH ₄ [1] step 4 = Pt AND H ₂ [1] | 2 |
| 8(b)(iv) | 5 / five [1] | 1 |

| Question | Answer | Marks |
|----------|---|-------|
| 9(a) | (because CDCl ₃ / it) does not give a peak [1] OR because CHCl ₃ does give a peak | 1 |
| 9(b) | as a standard / reference for (chemical shift measurements) [1] | 1 |
| 9(c) | ester [1] | 1 |
| 9(d)(i) | <ul style="list-style-type: none"> • (δ = 1.4) triplet • (δ = 1.4) two H on neighbouring C atom • (δ = 4.3) quartet / quadruplet • (δ = 4.3) three H on neighbouring C atom mark as • ✓ • ✓ [2] | 2 |
| 9(d)(ii) | aryl group / arene / phenyl [1] | 1 |

| Question | Answer | Marks |
|-----------|--|-------|
| 9(d)(iii) |  OR $C_6H_5CO_2C_2H_5$ [1] | 1 |
| 9(e) | $CH_3CH_2^+ / C_2H_5^+$ [1] $C_6H_5^+$ [1] | 2 |

| Question | Answer | Marks | | | | | | | | |
|---|--|------------------|------------------------|---|--------------|---|--------------|---|----------|---|
| 10(a) | <table border="1" style="width: 100%; border-collapse: collapse;"> <thead> <tr> <th style="width: 50%;">pair of monomers</th> <th style="width: 50%;">type of polymerisation</th> </tr> </thead> <tbody> <tr> <td>HOCH₂CH₂OH and HO₂CCH₂CO₂H</td> <td>condensation</td> </tr> <tr> <td>  </td> <td>condensation</td> </tr> <tr> <td>CH₃CHCF₂ and CH₃CHCH₂</td> <td>addition</td> </tr> </tbody> </table> ALL correct [1] | pair of monomers | type of polymerisation | HOCH ₂ CH ₂ OH and HO ₂ CCH ₂ CO ₂ H | condensation |  | condensation | CH ₃ CHCF ₂ and CH ₃ CHCH ₂ | addition | 1 |
| pair of monomers | type of polymerisation | | | | | | | | | |
| HOCH ₂ CH ₂ OH and HO ₂ CCH ₂ CO ₂ H | condensation | | | | | | | | | |
|  | condensation | | | | | | | | | |
| CH ₃ CHCF ₂ and CH ₃ CHCH ₂ | addition | | | | | | | | | |

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| Question | Answer | Marks |
|-----------|--|----------|
| 10(b)(i) | <p> <ul style="list-style-type: none"> • amide links displayed correct (NHCO OR CONH) • three monomers of Ala only • one repeat unit correctly identified • continuation bonds (with a polypeptide involving Ala only) mark as • ✓ ✓ ✓ [3] </p> | 3 |
| 10(b)(ii) | <p>3D, tetrahedral, both isomers of 2-aminopropanoic acid [1]</p> <p>optical [1]</p> | 2 |
| 10(c)(i) | epoxy resin [1] ALLOW Super Glues | 1 |
| 10(c)(ii) | compound with two amine groups per molecule, amine groups must not be on the same carbon atom [1] e.g. $\text{H}_2\text{NCH}_2\text{CH}_2\text{NH}_2$ | 1 |