

Mark Scheme

| Question | | Answer | Marks | Guidance |
|----------|-----|--|-------|---|
| 1 | (a) | $\text{Ba(OH)}_2 + 2\text{HCl} \rightarrow \text{BaCl}_2 + 2\text{H}_2\text{O} \checkmark$ | 1 | ALLOW multiples IGNORE state symbols (even if wrong) |
| | (b) | <p><i>Increasing size:</i> Atomic radius increases OR more shells OR more (electron) shielding \checkmark</p> <p><i>Attraction</i> Nuclear attraction decreases OR (outer) electron(s) experience less attraction \checkmark</p> <p><i>Ionisation energy</i> Ionisation energy decreases OR less energy needed to remove electron(s) \checkmark</p> | 3 | <p>FULL ANNOTATIONS WITH TICKS, CROSSES, CON, etc MUST BE USED</p> <p>IGNORE more orbitals OR more sub-shells <i>Alternative must refer to shells</i></p> <p>ALLOW Energy levels for shells</p> <p>ALLOW more electron repulsion between shells IGNORE just 'shielding' (<i>more/greater needed</i>) IGNORE 'nuclear shielding'</p> <p>IGNORE 'pull' for attraction IGNORE 'electrons less tightly held' IGNORE 'nuclear charge' for 'nuclear attraction'</p> <p>IGNORE 'easier to remove electron' <i>Energy is required</i></p> <p>ALLOW less energy to oxidise</p> |

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| (c) | (i) | <p><i>Disproportionation:</i> oxidation and reduction of the same element ✓</p> <p><i>Redox:</i> Cl is oxidised from +5 (in KClO_3) to +7 (in KClO_4) ✓</p> <p>Cl is reduced from +5 (in KClO_3) to -1 (in KCl) ✓</p> | 3 | <p>ALLOW 'chlorine' OR 'Cl' for same element IGNORE 'species' for 'element'</p> <p>ALLOW after number, e.g. 5+ IGNORE ionic charges, e.g. Cl^{5+}</p> <p>IGNORE '5' (signs required)</p> <p>IGNORE any reference to electron loss/gain (even if wrong)</p> <p>ALLOW one redox mark if oxidation numbers are correct but reduction/oxidation is incorrectly assigned</p> |
| | (ii) | potassium chlorate(VII) ✓ | 1 | Brackets required |
| (d) | (i) | <p><i>Equation</i> $\text{Ba}(\text{NO}_3)_2(\text{aq}) + \text{Na}_2\text{SO}_4(\text{aq}) \rightarrow \text{BaSO}_4(\text{s}) + 2\text{NaNO}_3(\text{aq})$ ✓</p> <p><i>Entropy change and explanation</i> entropy decreases OR entropy change negative AND (BaSO_4) solid/ppt has less disorder/ more order/ fewer ways of arranging energy/ less freedom/ less random particles/dispersal of energy ✓</p> | 2 | <p>ALLOW multiples</p> <p>M2 is dependent on $\text{BaSO}_4(\text{s})$ (even if formula is incorrect – eg $\text{Ba}(\text{SO}_4)_2(\text{s})$ seen as a product in the attempted equation as long as reactants are not solid. BaSO_4 solid/ppt may be assumed from $\text{BaSO}_4(\text{s})$ seen in the attempted equation.</p> |

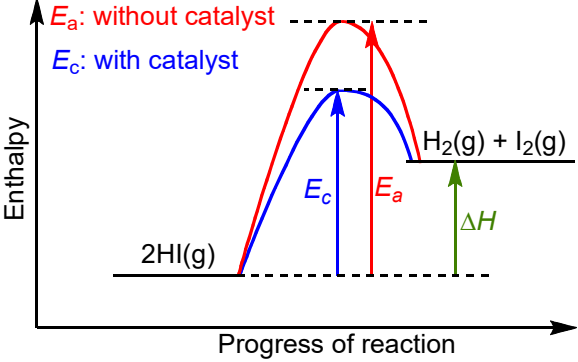
Mark Scheme

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| | (ii) | <p><i>Equation</i> $\frac{1}{2}\text{I}_2(\text{s}) \rightarrow \text{I}(\text{g}) \checkmark$ <i>state symbols required</i></p> <p><i>Entropy change and explanation</i> entropy increases OR entropy change positive AND gas has more disorder/ less order/ more ways of arranging energy/ more freedom/ more random particles / more dispersal of energy \checkmark</p> | 2 | <p>DO NOT ALLOW $\text{I}_2(\text{s}) \rightarrow 2\text{I}(\text{g})$</p> <p>DEPENDENT on $\frac{1}{2}\text{I}_2(\text{s}) \rightarrow \text{I}(\text{g})$ OR $\text{I}_2(\text{s}) \rightarrow 2\text{I}(\text{g})$</p> |
| | | Total | 12 | |

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| 2 | (a) | <p>ASSUME trend is down the group (unless stated otherwise)</p> <p>Forces London forces increase OR induced dipole(-dipole) interactions increase ✓</p> <p>Reason (Number of) electrons increases ✓</p> <p>Link to energy and particles More energy to break intermolecular forces OR to break London forces OR to break induced dipole(-dipole) interactions ✓</p> | 3 | <p>FULL ANNOTATIONS MUST BE USED ----- ALLOW reverse argument throughout</p> <p>IGNORE van der Waals'/vdW forces DO NOT ALLOW hydrogen bonds OR permanent dipole(-dipole) interactions for first and third marking points</p> <p>ALLOW more (electron) shells</p> <p>DO NOT ALLOW covalent bonds break</p> |

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| (b) |  <p>2HI(g) on LHS AND H₂(g) + I₂(g) on RHS ✓</p> <p>ΔH labelled with product above reactant AND arrow upwards ✓</p> <p>E_a AND E_c correctly labelled with E_c below E_a ✓</p> | 3 | <p>FULL ANNOTATIONS MUST BE USED</p> <p>Mark each point independently</p> <p>IGNORE state symbols.</p> <p>ΔH: DO NOT ALLOW -ΔH. ALLOW ΔH arrow even with a gap at the top and bottom, i.e. does not quite reach reactant or product line</p> <p>E_a: ALLOW no arrowhead or arrowheads at both ends of E_a line E_a line must reach (near or not too far beyond) maximums regardless of position</p> <p>ALLOW AE or EA for E_a</p> <p>Exothermic diagram can access the first and third marks</p> |

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| (c) | <p>FIRST CHECK THE ANSWER ON THE ANSWER LINE IF $M = 183$ AND Formula = Cl_2O_7 award 4 marks IF $M = 183$ award 3 marks</p> <hr/> <p>Use of data and unit conversions</p> <ul style="list-style-type: none"> • (R = 8.314) • T in K: 373K • V in m^3: 76.0×10^{-6} • (p in Pa: 1.00×10^5) ✓ <p>Calculation of n</p> $n = \frac{(1.00 \times 10^5) \times (76.0 \times 10^{-6})}{8.314 \times 373}$ $n = 2.45 \times 10^{-3} \text{ (mol) } \checkmark$ <p>Molar mass</p> $M = \frac{m}{n} = \frac{0.4485}{2.45 \times 10^{-3}} = 183 \text{ (g mol}^{-1}\text{)} \checkmark$ <p>Molecular formula</p> <p>Cl_2O_7 ✓</p> | 4 | <p>If there is an alternative answer, check to see if there is any ECF credit possible using working below</p> <p>Correct value of n subsumes first mark</p> <p>ALLOW ECF from incorrectly calculated n</p> <p>ALLOW ECF from incorrect M if formula of Cl_xO_y is the closest to the with calculated value of M</p> <p>IGNORE use of $24\,000 \text{ cm}^3$ for calculation of n BUT then Mark molar mass and Molecular formula by ECF for two marks maximum.</p> $n = \frac{76.0}{24000} = 3.17 \times 10^{-3} \text{ (mol)}$ $M = \frac{0.4485}{3.17 \times 10^{-3}} = 141.6/141.5 \text{ (g mol}^{-1}\text{)} \checkmark$ <p>Molecular formula = Cl_3O_2 ✓</p> |

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|----------|-------|---|----------|---|
| (d) | (i) | Titres correct and ALL recorded to 2 decimal places Titre: 24.00 23.40 23.75 23.85 ✓ mean titre = 23.80 (cm ³) ✓ | 2 | ALLOW 23.8 cm ³ |
| (d) | (ii) | Percentage uncertainty = $\frac{0.05 \times 2}{23.40} \times 100 = 0.43 (\%)$ ✓ | 1 | ALLOW ECF from incorrect subtraction in (i) or incorrect mean ALLOW 0.42% from titre values 2, 3 or 4 or mean titre or trial titre. 2 DP required |
| (d) | (iii) | Add starch (near the end point) ✓ Blue to colourless ✓ | 2 | ALLOW blue/black OR black OR purple for colour of mixture ALLOW blue colour disappears (to colourless) IGNORE 'clear' IGNORE 'colorimetry' |

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| (d) (iv) | <p>FIRST CHECK THE ANSWER ON THE ANSWER LINE IF B = RbIO₃ AND relative formula mass = 260.5 award 5 marks IF relative formula mass = 260.5 award 4 marks</p> <p>-----</p> <p>$n(\text{S}_2\text{O}_3^{2-})$ in titration $= \frac{0.150 \times 23.80}{1000} = 3.57 \times 10^{-3} \text{ (mol) } \checkmark$</p> <p>$n(\text{IO}_3^-)$ in titration $= \frac{3.57 \times 10^{-3}}{6} = 5.95 \times 10^{-4} \text{ (mol) } \checkmark$</p> <p>$n(\text{IO}_3^-)$ in original 250 cm³ $= 10 \times 5.95 \times 10^{-4} = 5.95 \times 10^{-3} \text{ (mol) } \checkmark$</p> <p>Relative formula mass of B $= \frac{1.55}{5.95 \times 10^{-3}} = 260.5 \text{ (g mol}^{-1}\text{) } \checkmark$</p> <p>Formula of B (must be derived from relative formula mass) iodate of Group 1 metal that most closely matches calculated molar mass of B</p> <p>Formula from 260.5 = RbIO₃ ✓</p> | 5 | <p>ALLOW ECF from incorrect mean titre in (a)(i)</p> <p>ECF from $n(\text{S}_2\text{O}_3^{2-})$ in titration ALLOW a two-step calculation $n(\text{I}_2) = n(\text{S}_2\text{O}_3^{2-}) \div 2$ and $n(\text{IO}_3^-) = n(\text{I}_2) \div 3$</p> <p>ECF from $n(\text{IO}_3^-)$ in titration</p> <p>ECF from $n(\text{IO}_3^-)$ in original 250 cm³ IF scaling $\times 10$ is omitted, ALLOW ECF from $n(\text{IO}_3^-)$ in titration</p> <p>ALLOW ECF from incorrect RFM of B provided metal is from Group 1 ALLOW RbIO₃⁻ DO NOT ALLOW RbIO₃ without relative formula mass value. DO NOT ALLOW 260.4 (without working) and RbIO₃ IF B = RbIO₃ AND relative formula mass = 261 award 5 marks</p> |
| | Total | 20 | |

Mark Schemes

| Question | Answer | Marks | AO element | Guidance |
|----------|--------|-------|------------|----------|
| 3 | C | 1 | AO1.2 | |
| 4 | B | 1 | AO2.6 | |

Mark Scheme

| Question | | Answer | Marks | AO element | Guidance |
|----------|-----|---|-------|--|---|
| 5 | (a) | <p>Interpretation of Results Orange contains bromine AND no reaction AND violet contains iodine ✓</p> <p>Ionic equation $\text{Br}_2 + 2\text{I}^- \rightarrow 2\text{Br}^- + \text{I}_2$ ✓</p> <p>Reactivity (down the group) Reactivity decreases AND oxidising power decreases OR gains electrons less easily OR forms negative ion/1- ion less easily OR less energy released when electron gained ✓ OR more negative electron affinity</p> <p>Size/shells/shielding (down the group) Greater atomic radius OR more shells OR more shielding ✓</p> <p>Attraction (down the group) Less nuclear attraction down the group ✓</p> | 5 | <p>2.3×1</p> <p>2.6×1</p> <p>1.1×3</p> | <p>Results can be interpreted anywhere in answer.</p> <p>ALLOW multiples, e.g. $\frac{1}{2}\text{Br}_2 + \text{I}^- \rightarrow \text{Br}^- + \frac{1}{2}\text{I}_2$ IGNORE other halogen/halide equations</p> <p>IGNORE state symbols</p> <p>ALLOW ORA</p> <p>DO NOT ALLOW idea of losing electrons/ionisation energy</p> <p>IGNORE chlorine is the most electronegative</p> <p>IGNORE explanations in terms of displacement</p> |

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|--------------|-----|---|----------|------------------|--|
| | (b) | <p><i>Benefit AND risk required for ONE mark</i></p> <p>Benefit: kills bacteria ✓ AND Risk: toxic/poisonous OR forms chlorinated hydrocarbons OR forms carcinogens/toxic compounds ✓</p> | 1 | 1.1 | <p>ALLOW kills micro-organisms OR kills pathogens OR kills viruses OR sterilises/disinfects water</p> <p>IGNORE antiseptic, reduces risk of disease, cleans water</p> <p>IGNORE 'harmful/'dangerous'</p> <p>IGNORE chlorine is carcinogenic/ dangerous for health/causes breathing problems</p> |
| | (c) | $n(\text{A}) = \frac{0.209}{29} = 0.00721 \text{ (mol)} \checkmark$ $M_r = \frac{1.26}{0.00721} = 174.8 \checkmark$ <p>Molecular formula = BrF₅ ✓ Formula is dependent on M_r</p> | 3 | 2.2×2 3.2 | <p>ALLOW ECF</p> <p>ALLOW 2SF 0.0072 up to calculator value 0.0072068965517</p> <p>ALLOW 175 up to calculator value 174.8325359</p> <p>ALLOW F₅Br</p> <p>ALLOW ECF that matches calculated M_r</p> |
| Total | | | 9 | | |