General Certificate of Education (A-level) January 2012

Chemistry

CHEM1

(Specification 2420)

Unit 1: Foundation Chemistry

Final



Mark schemes are prepared by the Principal Examiner and considered, together with the relevant questions, by a panel of subject teachers. This mark scheme includes any amendments made at the standardisation events which all examiners participate in and is the scheme which was used by them in this examination. The standardisation process ensures that the mark scheme covers the candidates' responses to questions and that every examiner understands and applies it in the same correct way. As preparation for standardisation each examiner analyses a number of candidates' scripts: alternative answers not already covered by the mark scheme are discussed and legislated for. If, after the standardisation process, examiners encounter unusual answers which have not been raised they are required to refer these to the Principal Examiner.

It must be stressed that a mark scheme is a working document, in many cases further developed and expanded on the basis of candidates' reactions to a particular paper. Assumptions about future mark schemes on the basis of one year's document should be avoided; whilst the guiding principles of assessment remain constant, details will change, depending on the content of a particular examination paper.

Further copies of this Mark Scheme are available from: aqa.org.uk

Copyright © 2012 AQA and its licensors. All rights reserved.

Copyright

AQA retains the copyright on all its publications. However, registered centres for AQA are permitted to copy material from this booklet for their own internal use, with the following important exception: AQA cannot give permission to centres to photocopy any material that is acknowledged to a third party even for internal use within the centre.

Set and published by the Assessment and Qualifications Alliance.

The Assessment and Qualifications Alliance (AQA) is a company limited by guarantee registered in England and Wales (company number 3644723) and a registered charity (registered charity number 1073334). Registered address: AQA, Devas Street, Manchester M15 6EX.

| Question | Marking Guidance | Mark | Comments |
|----------|--|------|---|
| 1(a) | Covalent | 1 | If not covalent CE = 0/2 |
| | | | If dative covalent CE = 0/2 |
| | | | If blank mark on |
| | | | Ignore polar |
| | | | If number of pairs of electrons specified, must be 3 |
| | Shared <u>pair(</u> s) of electrons / one electron from Br and one | 1 | Not 2 electrons from 1 atom |
| | electron from F | | Not shared pair between ions/molecules |
| 1(b)(i) | F | | |
| | F = P | 1 | $BrF_{3}\ should have 3 \ bp \ and 2 \ lp \ and \ correct \ atoms \ for \ the mark$ |
| | F C | | Penalise Fl |
| | r = 0r F | 1 | Allow 84 - 90° or 120° and ignore 180° |
| | BrF ₃ if trigonal planar shown = 120° or if T shape shown 84 - 90° | | Irrespective of shape drawn |

| 1(b)(ii) | F F F F | 1 | BrF ₄ should have 4 bp and 2 lp and all atoms for the mark (ignore sign) Allow FI |
|----------|---|--------|---|
| | BrF ₄ ⁻ 90° | 1 | Only Ignore 180° |
| 1(c) | Ionic or (forces of) attraction between ions / bonds between ions | 1 | If molecules, IMF, metallic, CE =0 If covalent bonds mentioned, 0/3, unless specified within the BrF_4^- ion and not broken Ignore atoms |
| | Strong (electrostatic) attraction / strong bonds / lots of energy needed to break bonds Between K ⁺ and BrF ₄ ⁻ ions/oppositely charged ions / + and - ions | 1 1 | If ions mentioned they must be correct Strong bonds between + and – ions =3/3 |
| 1(d)(i) | Hydrogen <u>bonds</u> /hydrogen <u>bonding</u> /H <u>bonds</u> /H <u>bonding</u> | 1 | Not just hydrogen |
| 1(d)(ii) | $ \begin{array}{cccccccccccccccccccccccccccccccccccc$ | 3 | One mark for 4 partial charges One mark for 6 lone pairs One mark for H bond from the <u>lone pair to the Hδ+</u> Allow FI If more than 2 molecules are shown they must all be correct. Treat any errors as contradictions within each marking point. CE = 0/3 if incorrect molecules shown. |

| 1(e) | vdw / van der Waals forces between molecules | 1 | QoL |
|------|---|---|--|
| | | | Not vdw between HF molecules, CE = 0/2 |
| | | | vdw between atoms, CE = $0/2$ |
| | | | If covalent, ionic, metallic, CE=0/2 |
| | IMF are weak / need little energy to break IMF / easy to overcome IMF | 1 | |

| Question | Marking Guidance | Mark | Comments |
|----------|--|------|---|
| 2(a) | Lithium / Li | 1 | Penalise obvious capital I (second letter). |
| 2(b)(i) | Increase / gets bigger | 1 | Ignore exceptions to trend here even if wrong |
| 2(b)(ii) | Boron / B | 1 | If not Boron, CE = 0/3 |
| | Electron removed from (2) <u>p</u> orbital /sub-shell / (2)p electrons removed | 1 | If p orbital specified it must be 2p |
| | Which is higher in energy (so more easily lost) / more shielded (so more easily lost) / further from nucleus | 1 | |
| 2(c) | C / carbon | 1 | |
| 2(d) | Below Li 2 (d) Draw a cross on the diagram to show the melting point of nitrogen. | 1 | The cross should be placed on the diagram, on the column for nitrogen, below the level of the cross printed on the diagram for Lithium. |

| 2(e) | Macromolecular / giant molecular / giant atomic | 1 | Allow giant covalent (molecule) = 2 |
|------|---|---|--|
| | Covalent bonds in the structure | 1 | |
| | <u>Strong</u> (covalent) <u>bonds must be broken or overcome</u> / (covalent) <u>bonds need a lot of energy to break</u> | 1 | Ignore weakening / loosening bonds If ionic / metallic/molecular/ dipole dipole/ H bonds/ bonds between molecules, CE = 0/3 Ignore van der Waals forces Ignore hard to break |

| Question | Marking Guidance | Mark | Comments |
|-----------|--|------|--|
| 3(a)(i) | Crude oil / oil / petroleum | 1 | Do not allow 'petrol' |
| 3(a)(ii) | Fractional distillation / fractionation / fractionating | 1 | Not distillation alone |
| 3(b)(i) | 5 | 1 | Allow five / V |
| 3(b)(ii) | Chain (isomerism) | 1 | Allow branched chain / chain branched / side chain (isomerism) |
| | | | Ignore position (isomerism) Do not allow straight chain / geometric / branched / function |
| 3(c)(i) | C ₁₂ H ₂₆ / H ₂₆ C ₁₂ | 1 | Only |
| 3(c)(ii) | Thermal cracking | 1 | If not thermal cracking, CE = 0/2 If blank mark on |
| | High temperature (400°C ≤ T ≤ 900°C) or (650 K ≤ T ≤ 1200 K) | 1 | Allow 'high heat' for 'high temperature' Not 'heat' alone |
| | and High pressure (≥ 10 atm, ≥ 1 MPa, ≥1000 kPa) | | If no T, units must be 650 - 900 |
| 3(c)(iii) | To produce substances which are (more) in demand / produce products with a high value / products worth more | 1 | Ignore 'to make more useful substances' |

| 3(d)(i) | Corrosive or diagram to show this hazard symbol | 1 | Ignore irritant, acidic, toxic, harmful |
|----------|---|---|---|
| 3(d)(ii) | (<u>120.5</u> × 100) (86 + 71) =76.75(%) or 76.8(%) | 1 | Allow answers > 3 sig figs |
| 3(e) | 2,2-dichloro-3-methylpentane C_3H_6CI | 1 | Ignore punctuation Any order |

| Question | Marking Guidance | Mark | Comments |
|----------|--|--------|---|
| 4(a)(i) | $C_8H_{18} + 8\frac{1}{2}O_2 \rightarrow 8CO + 9H_2O$ | 1 | Accept multiples |
| 4(a)(ii) | Not enough oxygen or air (available for complete combustion) /lack of oxygen or air / too much octane | 1 | Ignore poor ventilation, low temp, poor mixing, incomplete combustion |
| 4(b)(i) | $2CO + 2NO \rightarrow 2CO_2 + N_2$ | 1 | Allow multiples |
| 4(b)(ii) | Pt / Pd / Rh / Ir or names Big(ger) surface area / increased reaction rate / removes more of the gases / ensures complete reaction | 1 1 | Apply list principle Allow (ceramic) withstands high temperatures |
| 4(c)(i) | Acid rain | 1 | Allow consequence of acid rain Ignore greenhouse gas / global warming / ozone |
| 4(c)(ii) | CaO/ lime / CaCO $_3$ /limestone Neutralises the gas or words to that effect/it is basic/ SO $_2$ is acidic | 1 | Allow chemical names Allow 'reacts with it' or 'it is alkaline' Ignore 'absorb' |

| Question | Marking Guidance | Mark | Comments |
|----------|---|------|--|
| 5(a) | N ³⁻ / N ⁻³ | 1 | |
| 5(b) | F [−] / fluoride | 1 | Ignore fluorine/F Penalise Fl |
| 5(c) | Li ₃ N / NLi ₃ | 1 | |
| 5(d) | $\frac{81.1}{40.1} \frac{18.9}{14}$ (=2.02 = 1.35) | 1 | M1 for correct fractions |
| | 1.5 1 or 3:2 | 1 | M2 for correct ratio |
| | Ca ₃ N ₂ | 1 | If Ca_3N_2 shown and with no working award 3 marks |
| | | | If Ca_3N_2 obtained by using atomic numbers then lose M1 |
| 5(e) | $3 \text{ Si} + 2 \text{ N}_2 \rightarrow \text{Si}_3 \text{N}_4$ | 1 | Accept multiples |

| Question | Marking Guidance | Mark | Comments |
|-----------|---|------|---|
| 6(a) | Mol Pb = 8.14 / 207(.2) (= 0.0393 mol) | 1 | M1 and M2 are process marks |
| | Mol HNO ₃ = 0.0393 x 8 / 3 =0.105 mol | 1 | Allow mark for M1 x 8/3 or M1 x 2.67 |
| | Vol HNO ₃ = 0.105 / 2 = 0.0524 (dm ³) | 1 | Accept range 0.0520 to 0.0530 |
| | | | No consequential marking for M3 |
| | | | Answer to 3 sig figs required |
| 6(b) | 101000 (Pa) and 638 x 10 ⁻⁶ (m ³) | 1 | |
| | n = pV/RT $(= \frac{101000 \times 638 \times 10^{-6}}{8.31 \times 298})$ | | |
| | (8.31 x 298) | 1 | Can score M2 with incorrect conversion of p and V |
| | | | If T incorrect lose M1 and M3 |
| | <u>0.026</u> (0) (mol) | 1 | If answer correct then award 3 marks |
| | | | Allow answers to 2 sig figs or more |
| | | | 26.02 = 1 |
| | | | If transcription error lose M3 only |
| 6(c)(i) | $2Pb(NO_3)_2(s) \rightarrow 2 PbO(s) + 4NO_2(g) + (1)O_2(g)$ | 1 | Allow multiples |
| | | | Allow fractions |
| 6(c)(ii) | Decomposition not complete / side reactions / by-products / some (NO_2) escapes / not all reacts / impure Pb $(NO_3)_2$ | 1 | Ignore reversible / not heated enough / slow |
| 6(c)(iii) | Hard to separate O_2 from NO_2 / hard to separate the 2 | 1 | Allow mixture of gases |
| | gases | | Not 'all products are gases' |

| Question | Marking Guidance | Mark | Comments |
|----------|---|-------------|---|
| 7(a) | $\frac{(82 \times 2) + (83 \times 2) + (84 \times 10) + (86 \times 3)}{17} (1428) \\ (17) = \underline{84.0}$ | 1 1 1 | M1 for the top line M2 is for division by 17 Not 84 |
| | The <i>A</i> _r in the Periodic table takes account of the <u>other</u> <u>isotopes</u> / <u>different amounts of isotopes</u> (or words to that effect regarding isotopes) | 1 | No consequential marking from M1 or M2 Ignore units Award independently Comparison implied Isotope(s) alone, M4 = 0 |
| 7(b) | (Beam of electrons from) an electron gun / high speed / high energy electrons | 1 | |
| | Knocks out electron(s) (to form a positive ion) $Kr(g) + e^{-} \rightarrow Kr^{+}(g) + 2e(^{-})$ <i>OR</i> $Kr(g) \rightarrow Kr^{+}(g) + e(^{-}) / Kr(g) - e(^{-}) \rightarrow Kr^{+}(g)$ The ⁸⁴ Kr isotope Has 2 electrons knocked out / gets a 2+ charge | 1 1 1 | State symbols must clearly be (g) One mark for identifying the 84 isotope |
| | Thas 2 electrons knocked out / yets a 2+ charge | 1 | One mark for the idea of losing 2 electrons (from this isotope) |

General principles applied to marking CHEM1 papers by CMI+ (January 2012)

It is important to note that the guidance given here is generic and specific variations may be made in the mark scheme.

Basic principles

- Examiners should note that throughout the mark scheme, items that are underlined are required information to gain credit.
- Occasionally a response involves incorrect chemistry and the mark scheme records CE = 0, which means a chemical error has occurred and no credit is given for that section of the clip or for the whole clip.

A. <u>The "List principle" and the use of "ignore" in the mark scheme</u>

If a question requires **one** answer and a candidate gives two answers, no mark is scored if one answer is correct and one answer is incorrect. There is no penalty if both answers are correct.

N.B. Certain answers are designated in the mark scheme as those that the examiner should "Ignore". These answers are not counted as part of the list and should be ignored and will not be penalised.

B. Incorrect case for element symbol

The use of an incorrect case for the symbol of an element should be penalised **once only** within a clip. For example, penalise the use of "h" for hydrogen, "CL" for chlorine or "br" for bromine.

C. Spelling

In general

- The names of organic chemical compounds and functional groups **must be spelled correctly**, when specifically asked for, to gain credit.
- Phonetic spelling may be acceptable for some chemical compounds (eg amonia would be phonetically acceptable. However, ammoniam would be unacceptable since it is ambiguous).

N.B. Some terms may be required to be spelled correctly or an idea needs to be articulated with clarity, as part of the "Quality of Language" (**QoL**) marking. These will be identified in the mark scheme and marks are awarded only if the QoL criterion is satisfied.

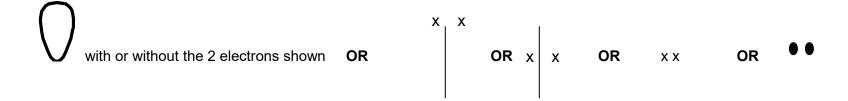
D. Equations

In general

- Equations **must** be balanced.
- State symbols are <u>generally</u> ignored, unless specifically required in the mark scheme.

E. Lone Pairs

The following representations of lone pairs in structures are acceptable.



F. <u>Reagents</u>

The command word "Identify", allows the candidate to choose to use **either** the name or the formula of a reagent in their answer. In some circumstances, the list principle may apply when the name and formula contradict. Specific details will be given in mark schemes.

G. Marking calculations

In general

- A correct answer alone will score **full marks** unless the necessity to show working is specifically required in the question.
- If a candidate has made an arithmetic error or a transcription error deduct one mark, but continue marking (error carried forward).

H. Organic structures

In general

- Displayed formulae must show all of the bonds and all of the atoms in the molecule, but need not show correct bond angles.
- Bonds should be drawn correctly between the relevant atoms.
- Latitude should be given to the representation of C C bonds in structures, given that CH_3 is considered to be interchangeable with H_3C even though the latter would be preferred.
- The following representations are allowed:-

