



General Certificate of Education

Chemistry 5421

**CHM1 Atomic Structure, Bonding and
Periodicity**

Mark Scheme

2009 examination - June series

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Question	Part	Sub Part	Marking Guidance	Mark	Comments
1	(a)	(i)	$1s^2 2s^2 2p^6 3s^2 3p^6 4s^2 3d^{10} 4p^4$ OR $1s^2 2s^2 2p^6 3s^2 3p^6 3d^{10} 4s^2 4p^4$;	1	Not [He] etc.
1	(a)	(ii)	'p' block; Highest energy / outermost e ⁻ /e ⁻ s in 'p' orbital/sub-level;	1 1	Not sub-orbital/outer orbital is 4p
1	(b)	(i)	Sum of/number of/amount of protons and neutrons	1	accept nucleons in place of p & n
1	(b)	(ii)	⁶⁸ Zn;	element = Zn mass number = 68	1 1 Incorrect atomic number negates 'Zn' Not 'ZN' Not decimal mass number (68.0)
1	(c)	(i)	High energy e ⁻ / high speed e ⁻ / e ⁻ from electron gun; Knock e ⁻ off (atom);	1 1	For M2, allow: $M + e^- \rightarrow M^+ + 2e^-$
1	(c)	(ii)	(Ions) generate/produce a <u>current</u> (on collision at the detector);	1	Not: current detected / current induced / flow of electricity / electric charge / electric impulse
1	(d)		$A_r = \frac{(76 \times 11.2) + (78 \times 23.8) + (80 \times 49.8) + (82 \times 15.2)}{100}$ = 79.4;	1 1	Wrong approach or not dividing by 100 = CE = 0. Answer to 1 d.p. Mark conseq on transcription error.

Question	Part	Sub Part	Marking Guidance	Mark	Comments
2	(a)		$\text{Al}_2\text{Se}_3 + 6\text{H}_2\text{O} \rightarrow 2\text{Al}(\text{OH})_3 + 3\text{H}_2\text{Se}$;	1	Allow fractions and multiples
2	(b)	(i)	H_2Se = 'bent' shape with 2 lone pairs – based on tetrahedral; NH_3 = pyramidal shape with 1 lone pair;	1 1	
2	(b)	(ii)	Bent / V-shaped /triangular/angular;	1	Not bent-linear
2	(b)	(iii)	107° ; Se has 2 lone pairs but N has only 1 lone pair; Lone pairs repel more strongly than bonding pairs / etc;	1 1 1	Allow $106\frac{1}{2}^\circ - 107\frac{1}{2}^\circ$ Allow Se has more lone pairs than N if supported by diagrams
2	(c)		Clear indication that covalent bond is a shared e^- pair; Se supplies one e^- and H supplies one/the other e^- ;	1 1	May be inferred from <u>statement</u> that $1e^-$ used from Se and $1e^-$ from H Allow M2 for a correct dot- <u>and</u> -cross diagram (not if only dots or crosses used)

Question	Part	Sub Part	Marking Guidance	Mark	Comments												
3	(a)	(i)	Moles $\text{BaSO}_4 = 2.10 \div 233.4 = 9.00 \times 10^{-3}$;	1	Accept $8.9 \times 10^{-3} - 9 \times 10^{-3}$												
3	(a)	(ii)	Moles $\text{Al}_2(\text{SO}_4)_3$ in $25 \text{ cm}^3 = 3.00 \times 10^{-3}$; [conseq on error in (i)]	1	Accept $2.9 \times 10^{-3} - 3 \times 10^{-3}$												
			Moles $\text{Al}_2(\text{SO}_4)_3$ in sample = 3.00×10^{-2} ; [conseq on error in 25 cm^3]	1	Accept $2.9 \times 10^{-2} - 3 \times 10^{-3}$												
3	(b)		M_r of $\text{Al}_2(\text{SO}_4)_3 \cdot x\text{H}_2\text{O} = 20.0 \div 3.00 \times 10^{-2}$;	1	If 3.17×10^{-3} used												
			= 666.9;	1	630.9												
			M_r of $\text{Al}_2(\text{SO}_4)_3 = 342(.3)$;	1	342(.3)												
			M_r of $x\text{H}_2\text{O} = 666.9 - 342.3 = 324.6$; [conseq on error(s) in M_r]	1	288.7												
			$X = \underline{18} / \text{Al}_2(\text{SO}_4)_3 \cdot 18\text{H}_2\text{O}$; [conseq on errors above [must be <u>integer</u>]]	1	<u>16</u>												
3	(c)		Mass O = 0.389;	1	At number(s) used M2 & M3 = CE = 0 Conseq on error in mass of O, Not $\text{Ba}(\text{CO}_2)_2 / \text{BrC}_2\text{O}_4$ etc.												
			<table border="0"> <tr> <td></td> <td style="text-align: center;">Ba</td> <td style="text-align: center;">C</td> <td style="text-align: center;">O</td> </tr> <tr> <td></td> <td style="text-align: center;"><u>0.835</u></td> <td style="text-align: center;"><u>1.46</u></td> <td style="text-align: center;"><u>0.389</u></td> </tr> <tr> <td></td> <td style="text-align: center;">137.3</td> <td style="text-align: center;">12</td> <td style="text-align: center;">16</td> </tr> </table>			Ba	C	O		<u>0.835</u>	<u>1.46</u>	<u>0.389</u>		137.3	12	16	1
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<table border="0"> <tr> <td></td> <td style="text-align: center;">0.00608</td> <td style="text-align: center;">0.0122</td> <td style="text-align: center;">0.0243</td> </tr> <tr> <td style="text-align: center;">BaC_2O_4</td> <td style="text-align: center;">1.0</td> <td style="text-align: center;">2.00</td> <td style="text-align: center;">4.00</td> </tr> </table>		0.00608	0.0122	0.0243	BaC_2O_4	1.0	2.00	4.00	1								
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BaC_2O_4	1.0	2.00	4.00														
3	(d)	(i) QoL	(Actual) number of <u>atoms of each element</u> in a molecule/compound;	1													
3	(d)	(ii)	The M_r / relative molecular/formula mass;	1	Not molecular mass												

Question	Part	Sub Part	Marking Guidance	Mark	Comments
4	(a)	(i)	<u>Energy/enthalpy required to remove one electron / energy/enthalpy change</u> when one electron is removed; From a gaseous atom; Enthalpy required/change for $M(g) \rightarrow M^+(g) + e^-$ earns both marks	1 1	If molar def ⁿ give, must be molar throughout Allow state symbol in equation for M2
4	(a)	(ii)	Increase in number of protons/nuclear charge; Decrease in size / e^- closer to nucleus same/similar shielding / same shells;	1 1	Accept 'increased attraction between nucleus and outer electrons' as alternative to <u>either</u> M1 or M2 <u>but tied to</u> M1/M2 or near miss
4	(a)	(iii)	Electron removed from 3p orbital/sub-shell in Al rather than 3s in Mg; Al loses e^- which is of higher energy /is further away from nucleus /is shielded by <u>3s</u> electrons;	1 1	Not just e^- arrangements
4	(b)	(i)	$MgCl_2(aq)$ <u>White</u> ppt/solid formed; $BaCl_2(aq)$ No change/no ppt/clear solution;	1 1	 Not nothing/dissolves
4	(b)	(ii)	$Mg^{2+}(aq) + 2OH^-(aq) \rightarrow Mg(OH)_2(s)$;	1	Penalise 2 equations Allow Ba^{2+} equation conseq on reversed observations in (b)(i)
4	(c)	(i)	(A covalent bond in which) the electron density is unequally shared;	1	Allow e^- cloud is unequally shared Not e^- cloud distorted
4	(c)	(ii)	Greater difference in electronegativity in HCl / Cl more electronegative than I; QoL So electron density more strongly <u>drawn towards the Cl</u> (than towards the I);	1 1	Not HCl/Cl ⁻ is more electronegative Tied to M2 or 'near miss' Not more attracted to Cl but allow e^- cloud more distorted towards Cl

Question	Part	Sub Part	Marking Guidance	Mark	Comments
5	(a)	M1	Macromolecular/giant atomic/giant covalent structure;	1	Allow separate 'giant' and 'covalent' if appropriate Reference to IMF loses M1 Ions/metallic bonds = CE = 0 for M1-M3 Reference to C-C bonds negates M1
		M2	Covalent bonds must be <u>broken/overcome</u> ;	1	Tied to M2 or near miss Not loosened
		M3	Much energy to break bonds/bonds strong/many bonds to break;	1	
		M4	S and Cl both (simple) molecular / S ₈ & Cl ₂ / separate refs to molecules;	1	
		M5	Van der Waals' forces present in both (or implied);	1	Not vdW between atoms
		M6	vdW forces increase with number of e ⁻ /size/M _r etc;	1	
		M7	e ⁻ /size/M _r is smaller for Cl molecules;	1	'molecules' may be implied
5	(b)		Correct bonding in both structures – accept correct descriptions	1	Incorrect bonding loses M1 and either M2 or M3, depending on the bonding error.
			In Na, atoms/ions/layers/particles slide/move when struck (but still metallic bonding);	1	
			Idea that, in NaCl, ions of same charge repel each other;	1	

Question	Part	Sub Part	Marking Guidance	Mark	Comments
6			$pV = nRT;$	1	If $V = p/nRT$ lose M3 and M4
			Moles methanol = $n = \frac{1.04}{32}$ (= 0.0325 mol);	1	
			$V = nRT/p = \frac{0.0325 \times 8.31 \times 350}{101000}$; Conseq on error in moles	1	If <u>1</u> used, & $V = .936 \text{ dm}^3$ allow M3 & M4 101
			= $9.36 \times 10^{-4} \text{ m}^3$; ignore missing units but penalise incorrect units	1	If <u>1</u> used, & $V = .936 \text{ m}^3/\text{cm}^3/\text{no units}$, 101 penalise M3 allow M4 conseq
			= 936 cm^3 ; mark is for correct conversion from their M4 units to cm^3	1	If no units in M4, assume them to be m^3 when awarding M5