M1.	(a)	(i)	1s² 2s² 2p ⁶ 3s² 3p¹ (1) Allow subscripted electron numbers
	(ii)	p (bl	ock) (1)

Allow upper or lower case 's' and 'p' in (a)(i) and (a)(ii)

2

(b) Lattice of metal / +ve ions/ cations / atoms (1) Not +ve nuclei/centres Accept regular array/close packed/tightly packed/uniformly arranged

(Surrounded by) <u>delocalised</u> electrons **(1)** *Note: Description as a 'giant ionic lattice' = CE*

2

(c) Greater nuclear or ionic charge or more protons (1)

Smaller atoms / ions (1)

Accept greater charge density for either M1 or M2

More delocalised electrons / e^- in sea of e^- / free e^- (1)

Stronger attraction between ions and delocalised / free electrons etc. (1)

Max 3 Note: 'intermolecular attraction/ forces' or covalent molecules = CE Accept stronger 'electrostatic attraction' if phrase prescribed elsewhere Ignore references to m/z values If Mg or Na compared to AI, rather than to each other, then: Max 2

Treat description that is effectively one for lonisation Energy as a '**contradiction**'

3

(d) (Delocalised) electrons (1)

Move / flow in a given direction (idea of moving non-randomly) **or** under the influence applied pd *QoL mark* **(1)**

Allow 'flow through metal' Not: 'Carry the charge'; 'along the layers'; 'move through the metal' M2. (a) Elements in the p block have their outer electron(s) in p orbital(s) or levels or sub-shells (1) example of element (1) correct electronic configuration (1)

3

2

(b) Pattern in the change in the properties of a row of elements (1) OR Trend in the properties of elements across a period

Repeated in the next row (1)

OR element underneath (or in same group) has similar properties

atomic radius

decreases across the row (1) CE if trend is wrong

number of protons increases (1) (or nuclear charge increases) more attraction for electrons in the same shell (1)

electronegativity

increases across the row (1) number of protons increases (1) (or nuclear charge) atomic radius decreases (1) (or shielding remains the same or electrons in the same shell) more attraction for <u>bonding</u> or <u>shared</u> electrons (1)

conductivity

decreases row (1) OR significant drop from AI to Si

	Na–Al me	als (1) OR metallic bonding or description of metallic bonding
	Two of Si	- Ar non metals (1) <i>OR molecular or covalent</i>
	EITHER ele OR electro	ectrons free to move (or delocalised) in metals ns unable to move in non-metals (1)
	(a) Ability (or electror in a covale	(or power) of an atom to attract electron density is or - ve charge) (1) nt bond (1)
		or shared pair If remove an electron lose first mark
)	<i>Trend:</i> incl <i>Explanatiol</i> electrons ir	reases (1) n: nuclear charge (number of protons) increases (1) n same shell (1)
		OR similar shielding OR atoms similar size or smaller

OR atoms similar size or smaller OR 1 mol of e

3

2

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(c) Heat / enthalpy / energy for removal of one electron (1) from a gaseous atom (1) can score in an equation must have first mark to score the second

2

(d) (i) 2 (1)

M3.

(b)

- (ii) <u>Two elements</u> (or Na / Mg) before the drop (in energy) to Al (1)
- (iii) ionisation energy of Al < that for Mg (1)
- (iv) fall in energy from P to S (1)

or discontinuity in trend

From Al to P there are 3 additional electrons (1) or three elements For second mark idea of block of 3 elements

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M4.

(a)

(i) <u>Deductions</u>:

lonic (1)
lons not free to move in the solid state (1)
lons free to move when molten or in aqueous solution (1)
ldentity of P: Na₂O or sodium oxide (1) *N.B. If a formula given this must be correct*

Equation: $Na_2O + H_2O \rightarrow 2 NaOH$ (1)

5

(ii) <u>Deductions</u>:

Covalent Intermolecular forces are weak or van der Waals forces, or dipole-dipole *N.B. Any answer including a reference to hydrogen bonding is incorrect*

Identity of **Q**: SO₂ or sulphur dioxide (1)

Equation: $SO_2 + H_2O \rightarrow H_2 SO_3(1)$ NB Allow max one for SO_3

4

(b) (i) Amphoteric (1)

(ii) Equation with NaOH

 $\begin{array}{l} \mathsf{AI}(\mathsf{OH})_{\scriptscriptstyle 3} + \mathsf{Na}\mathsf{OH} \to \mathsf{Na}\mathsf{AI}(\mathsf{OH})_{\scriptscriptstyle 4} \\ & \mathsf{OR}\;\mathsf{AI}(\mathsf{OH})_{\scriptscriptstyle 3}(\mathsf{H}_2\mathsf{O})_{\scriptscriptstyle 3} + \mathsf{OH}^+ \to [\mathsf{AI}(\mathsf{OH})_{\scriptscriptstyle 4}(\mathsf{H}_2\mathsf{O})_{\scriptscriptstyle 2}]^{\scriptscriptstyle +} + \mathsf{H}_2\mathsf{O} \\ & \mathsf{OR}\;\mathsf{AI}(\mathsf{OH})_{\scriptscriptstyle 3} + \mathsf{OH}^- \to [\mathsf{AI}(\mathsf{OH})_{\scriptscriptstyle 4}]^{\scriptscriptstyle +} \end{array}$

R identified as $AI(OH)_3$ or $AI(OH)_3(H_2O)_3$ (1)

A balanced equation (1)

N.B. Allow equation with six co-ordinate Aluminium and up to six OH-ligands N.B. Allow equation mark if $M(OH)_3$ given in a balanced equation

Equation with H_2SO_4

 $2AI(OH)_{\scriptscriptstyle 3} + 3H_{\scriptscriptstyle 2}SO_{\scriptscriptstyle 4} \rightarrow AI_{\scriptscriptstyle 2}(SO_{\scriptscriptstyle 4})_{\scriptscriptstyle 3} + 6H_{\scriptscriptstyle 2}O$

 $OR AI(OH)_{3}(H_{2}O)_{3} + H^{*} \rightarrow [AI(OH)_{2}(H_{2}O)_{4}^{*} + H_{2}O]$

NB Allow equations with six co-ordinate Aluminium and up to six H_2O ligands NB Allow equation mark if $M(OH)_3$ given in a balanced equation

Correct AI species as product (1) A balanced equation (1)

 (iii) Large lattice energy or strong covalent bonds or ΔH_{soln} is very positive or ΔG is positive or sum of hydration energies less than covalent bond energies (1)

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