1.	Cu c Cu h	as no d electrons as an ion / Cu has d electrons; ompounds are coloured / Sc compounds are colourless; as more than one oxidation state / Sc has only one oxidation state; ompounds can act as catalysts / Sc cannot act as catalysts;	3 max	[3]
; 2.	(i)	the amount of energy required to remove one (mole of) electron(s); from (one mole of) an atom(s) in the <u>gaseous</u> state;	2	
	(ii)	greater positive charge on nucleus / greater number of protons / greater core charge; greater attraction by Mg nucleus for electrons (in the same shell) / smaller atomic radius;	2	[4]
	Cu c	ompounds can act as catalysts / Sc cannot act as catalysts;	3 max	[3]
3.	A			[1]
4.	meta radii nucle Dow incre	Ilic bonding gets weaker; /atoms get bigger / delocalized electrons shielded/screened from eus by filled shells; /a group 7: eased M_r of halogen molecules / OWTTE; molecular/van der Waals/London/dispersion forces increase;	4	[4]
5.	(a) (b)	as (cat)ion becomes more positive / Na ⁺ , Mg ²⁺ , Al ³⁺ / size/radius decreases / charge density increases; Do not allow increasing number of protons or increasing nuclear charge. attraction for mobile/valence/delocalized/sea of electrons increases; Do not accept "cloud of electrons". larger molecule / higher $M_{\rm I}/M$ / greater number of electrons; Do not accept "larger/higher/greater mass".	2	

IB Questionbank Chemistry

greater van der Waals'/dispersion/London forces;	

(c) Si: giant/network/macromolecular/3-D covalent bonding;
No mark for strong bonding without reference to covalent and network.
No mark for molecular.

Ar: (simple) atomic / (only weak) van der Waals'/dispersion/London forces; *No mark for (simple) molecular.*

[6]

6. (i) atomic number / Z; Accept nuclear charge / number of protons.

1

2

2

	(ii)	Across period 3: increasing number of protons / atomic number / Z / nuclear charge; (atomic) radius/size decreases / same shell/energy level / similar shielding/screening (from inner electrons); No mark for shielding/screening or shielding/screening increases.		
		Noble gases: do not form bonds (easily) / have a full/stable octet/shell/energy level / cannot attract more electrons; Do not accept "inert" or "unreactive" without reference to limited ability/ inability to form bonds or attract electrons.	3	[4]
7.	(i)	first ionization energy: $M(g) \rightarrow M^+(g) + e^-/e$ / the (minimum) energy (in kJ mol ⁻¹) to remove one electron from a <u>gaseous</u> atom / the energy required to remove one mole of electrons from one mole of <u>gaseous</u> atoms; periodicity: <u>repeating</u> pattern of (physical and chemical) properties;	2	
	(ii)	2.8.8/sp version;		
		Accept any two of the following: the outer energy level/shell is full; the increased charge on the nucleus; great(est) attraction for electrons;	3 max	
	(iii)	17 p in Cl nucleus attract the outer level more than 11 p in Na nucleus / greater nuclear charge attracts outer level more; <i>Allow converse for Na</i> .		
		Do not accept larger nucleus.	1	
	(iv)	S^{2-} has one proton less / smaller nuclear charge so outer level held less strongly / <i>OWTTE</i> ; <i>Allow converse for chloride</i> .		
		Do not accept larger nucleus.	1	
	(v)	the radii of the metal atoms increase (from Li \rightarrow Cs) (so the forces of attraction are less between them) / <i>OWTTE</i> ;		
		the forces of attraction between halogen molecules are van der Waals; forces increase with increasing mass/number of electrons;	3	[10]
8.	A			[1]

 ${\it Silicon\ dioxide:}\ strong/covalent\ bonds\ in\ network/giant\ structure/macromolecule;$

IB Questionbank Chemistry

9.

	Carb	on dioxide: weak/van der Waals'/dispersion/London forces between molecules	; 2	[2]
10.	one e coord	(covalent) bond; electron pair donated by oxygen to carbon atom / dative (covalent)/ dinate (covalent) bond; ed [1 max] for representation of $C \equiv O$. Ed [2] if CO shown with dative covalent bond.	2	[2]
→ M	(in k. requi	re ⁻ /e / the (minimum) energy I mol ⁻¹) to remove one electron from a <u>gaseous</u> atom / the energy red to remove one mole of electrons from one mole of <u>gaseous</u> atoms; dicity: <u>repeating</u> pattern of (physical and chemical) properties;	2	
	(ii)	2.8.8/sp version; Accept any two of the following: the outer energy level/shell is full; the increased charge on the nucleus; great(est) attraction for electrons;	3 max	
	(iii)	17 p in Cl nucleus attract the outer level more than 11 p in Na nucleus / greater nuclear charge attracts outer level more; Allow converse for Na. Do not accept larger nucleus.	1	
	(iv)	S ²⁻ has one proton less / smaller nuclear charge so outer level held less strongly / <i>OWTTE</i> ; Allow converse for chloride. Do not accept larger nucleus.	1	
	(v)	the radii of the metal atoms increase (from Li \rightarrow Cs) (so the forces of attraction are less between them) / $OWTTE$; the forces of attraction between halogen molecules are van der Waals; forces increase with increasing mass/number of electrons;	3	[10]

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11. A [1] **12.** D [1] **13.** D [1] **14.** Α [1] **15.** В [1] **16.** (a) (i) $\begin{matrix} & H \\ H & \overset{\times}{N} & \overset{\times}{N} & H \end{matrix};$ *Accept* ×'s, dots or lines for electron pairs. 2 *H–C–H:* (ii) any angle between 118° and 122°; due to three negative charge centres/electron domains/electron pairs; H–N–H: any angle between 104° and 108°; due to four negative charge centres/electron domains/electron pairs; extra repulsion due to lone electron pairs; 5 Do not allow ECF for wrong Lewis structures. (b) (relative) measure of an atoms attraction for electrons; (i) in a covalent bond / shared pair; 2

		(11)	as N is more electronegative / difference in electronegativity is greater for N-H than C-H;		
		(iii)	bond polarities cancel in C_2H_4 / OWTTE;	1	
	(c)	stron	ter van der Waals'/London/dispersion/intermolecular forces in ethene; ger (intermolecular) hydrogen bonding in hydrazine; comparison between strengths then [1 max].	2	[13]
17.	A				[1]
18.	$M_{\rm r}$ /n $Acce$ great NH_3	umber <i>pt elec</i> ter disp /ammo	nts increase going down the group (from PH ₃ to AsH ₃ to SbH ₃); of electrons/molecular size increases down the group; tron cloud increases down the group for the second marking point. persion/London/van der Waals' forces; nia has a higher boiling point than expected due to the hydrogen tween the molecules;		
		_	ept hydrogen bonding alone.	4	[4]
19.	(i)	centr electr	number of electron pairs/charge centres in (valence shell of) al atom; ron pairs/charge centres (in valence shell) of central atom repel other;		
		to po pairs	one of the following: sitions of minimum energy/repulsion / maximum stability; forming a double or triple bond act as a single bond; bonding pairs repel more than bonding pairs / OWTTE;	3 max	
			Do not accept repulsion between bonds or atoms.		

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(ii) 6

Species	Lewis (electron-dot) structure	Shape	Bond angle(s)
$\mathrm{H_3O}^{^+}$	$\left[\begin{array}{c} \bar{O} \\ H \end{array}\right]^+;$	Trigonal/triangular pyramidal;	Allow values in the range 106° to 109.5°;
C_2H_4	$H \rightarrow H$;	Trigonal/triangular planar;	Allow values of approximately 120°;

Accept crosses and dots for electrons in Lewis structures also. As the Lewis structures were asked for, and not 3D representations, do not penalize incorrectly drawn geometries. Do not accept structure of hydronium cation without lone pair on oxygen.

No penalty for missing charge.

(iii) H₃O⁺: is polar and explanation either using a diagram or in words, involving the net dipole moment;

e.g. the three individual O-H bond dipole moments add as vectors to give a net dipole moment.

C₂H₄: is non-polar and explanation either using a diagram or in words, involving no net dipole moment;

e.g. the vector sum of the individual bond dipole moments is zero.

For simple answers such as bond polarities do not cancel for H_3O^+ and do cancel for C_2H_4 , Award [1], only for the last two marking points.

(iv) O-H is most polar;

O-H has greatest difference between electronegativities $\!\!\!/$ calculation showing values of 1.4, 0.5 and 0.9 respectively;

[13]

2

2