

## Mark schemes

**1** B [1]

**2** A [1]

**3** (a) M1 have the same molecular formula  
or are  $C_3H_6O$   
or both have the same number/amount of each type of atom or same amount of each  
element  
or are isomers

*Not just the same atoms;*

1

M2 identical / exactly the same / same precise (relative) molecular  
mass / formula mass /  $M_r$

*Same (relative) molecular mass / formula mass /  $M_r$  is NOT enough  
got score M2*

*Allow same accurate (relative) molecular mass / formula mass /  $M_r$*

*Ignore reference to number of decimal places*

1

(b) M1 prop-2-en-1-ol

*Must refer to this compound clearly by name or structure (not to  
alcohol alone); ignore minor slips in name/structure*

1

M2 O(-)H (alcohol) and 3230–3550 ( $cm^{-1}$ ), or  
C=C and 1620–1680 ( $cm^{-1}$ )

*Marked independently from M1*

*Could score from bond labelled on correct signal on spectrum*

*Allow any value within these ranges*

*If additional incorrect signals given penalise M2*

*Ignore signals below 1500  $cm^{-1}$  and C-H signals*

1

- (c) (i) Determine the level by looking at the chemical content. (**NB** - If there is clear breakage of covalent bonds then max level 2 (max 3 marks).
- (ii) The mark within that level is then determined by looking at how coherent and logical the answer is and by use of terminology; start at the higher mark and penalise poor terminology/explanation; examples of terminology that would reduce the mark to the lower one:
- reference to van der Waals 'bonds' or dipole-dipole 'bonds' in relevant compounds that are being credited
  - uncertainty about whether hydrogen bonds are the O-H bonds within or are forces/bonds between molecules (if the alcohol is being credited)
  - use of 'vdw' or 'dip-dip' unless these terms 'van der Waals' for 'dipole-dipole' have been used elsewhere in answer (note that IMF and H-bond would not be penalised)
- (iii) If the answer does not achieve level 1, then 1 mark maximum could be scored for any correct point from the list of indicative content

### Level 3

- **Relative order** of boiling points of **all three** compounds
- Strongest intermolecular force of **all three** compounds identified
- Answer explains this coherently and logically and uses correct terminology for all **three** compounds

5-6 marks

### Level 2

- **Relative** boiling points of **two** compounds correctly compared
- Strongest intermolecular force for these **two** compounds correctly identified
- Answer explains this coherently and logically and uses correct terminology for **these two** compounds

3-4 marks

### Level 1

- **One** compound with the **highest** or **lowest** boiling point is correctly identified
- Strongest intermolecular force for that **one** compound identified
- Answer explains this coherently and logically and uses correct terminology for **this one** compound
- Allow 1 mark for individual correct point from indicative content on the right if no other mark scored

1-2 marks

### Level 0

None of the indicative chemistry content given.

0 marks

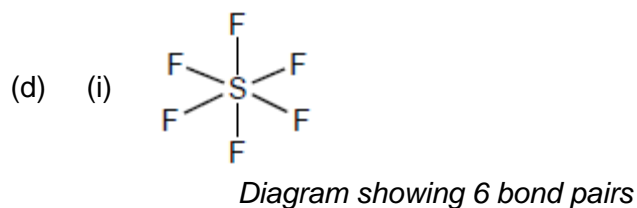
**Indicative chemistry content:**

- Correct order (highest to lowest) = prop-2-en-1-ol > propanal > butane
  - Prop-2-en-1-ol has hydrogen bonds
  - Propanal has (permanent) dipole-dipole forces
  - Butane has van der Waals' forces
  - Strength of intermolecular forces:  
hydrogen bonds > dipole-dipole > van der Waals
- (Note - actual values for reference are prop-2-en-1-ol 97°C, propanal 46°C and butane -1°C)

[10]

4

- (a) Electron movement in first molecule / temporary dipole  
*Allow description* 1
- Induces a dipole in another molecule  
*Allow description* 1
- (Induced-temporary) attraction or  $\delta^+$  attracts  $\delta^-$  in different/adjacent molecules  
*M3 dependent on M1 and M2*  
*Allow electrostatic attraction*  
*M3 could be scored in diagram*  
*If other type of force / metallic / ionic / polar bonds / permanent dipoles / difference in electronegativity mentioned CE = 0* 1
- (b) (i) (Methanol) H-bonds / hydrogen bonding 1
- (Methanethiol) dipole-dipole forces or van der Waals 1
- H-bonds are a stronger / are the strongest IMF 1
- Allow H-bonds require more energy to overcome*  
*If M1 and M2 not scored, allow 1 for methanol has stronger IMFs*  
*If breaking covalent bonds then CE=0*
- (ii) (Fractional) distillation 1
- Allow description*  
*Do not allow heating unqualified*
- (c) (Methaneselenol is a) bigger molecule / larger Mr / larger no of electrons / Se bigger atom 1
- With stronger/more vdw forces between molecules  
*If breaking covalent bonds then CE=0* 1



1

(Bond angle)  $90^\circ$  for  $SF_6$

Ignore  $180^\circ$

1

Octahedral

1

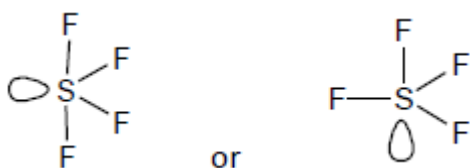


Diagram showing 4 bond pairs and 1 lone pair

1

(Bond angles) for  $SF_4$

If shape of  $SF_4$  is not based on 4 bond pairs and 1 lone pair cannot score M4 or M5

Any **two** from:

Allow  $85 - 89^\circ$

Do not allow  $90^\circ$

Allow  $100 - 119^\circ$

Do not allow  $120^\circ$

Allow  $170 - 179^\circ$

Do not allow  $180^\circ$

2

(ii) NaCl (as product in any equation)

1



Allow multiples

Ignore states

1

[17]

5

(a) General increase

If not increase then CE

1

Greater nuclear charge / more protons

1

Same shielding / electrons added to same shell

Allow similar

1

Stronger attraction (from nucleus) for outer electron(s)

Allow electron in outer shell

1

- (b) Aluminium / Al (lower than Mg)  
*CE if not Al or S* 1
- (Outer) electron in (3)p orbital / sub-shell (level)  
*If 2p or 4p orbital lose M2 and M3* 1
- (3p) higher in energy  
*Allow more shielded or weaker nuclear attraction*  
*M3 is dependent on M2* 1
- or
- Sulfur / S (lower than P)  
 (Outer) electrons in (3)p orbital begin to pair  
 Repel  
*If 2p or 4p orbital lose M2 and M3*  
*Allow 2 electrons in (3)p*  
*M3 is dependent on M2* 1
- (c) Sulfur / S  
*CE if not S* 1
- Large jump after 6<sup>th</sup> or between 6<sup>th</sup> and 7<sup>th</sup>  
*Do not allow M2 if atom/ion is removed* 1
- (d) Silicon  
*CE if not Si* 1
- Giant covalent structure / macromolecule 1
- Covalent (bonds)  
*Giant covalent scores M2 and M3* 1
- Many / strong (covalent bonds) or  
 (covalent bonds) need lots of energy to break  
*CE for M2-M4 if molecules / metallic / ionic / IMFs mentioned* 1
- [13]
- 6** (a) OH AND alcohol  
**IGNORE** hydroxy(l) 1

(b) **A** = butan-2-ol /  $\text{CH}_3\text{CH}(\text{OH})\text{CH}_2\text{CH}_3$

*If formulae given then must be unambiguous*

*If both formula and name given then formula must match name for mark to be awarded*

1

**B** = butan-1-ol /  $\text{CH}_3\text{CH}_2\text{CH}_2\text{CH}_2\text{OH}$

1

Product from **A** / **P** is a ketone

**AND**

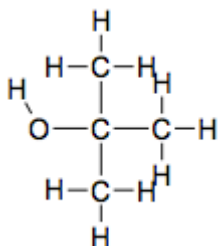
Product from **B** / **Q** is an aldehyde

*Penalise reference to incorrect class of alcohol*

1

(c) Type of Bond: C=C

1



*Must show all bonds in Isomer C including O-H bond*

1

Reagent: conc.  $\text{H}_2\text{SO}_4$  / conc.  $\text{H}_3\text{PO}_4$

*If incorrect attempt at correct reagent, mark on*

*Apply list principle for reagents and conditions marks*

*Conc required - may appear on conditions line*

*NOT (aq) For M3 even if seen on conditions line*

*ALLOW*

*Reagent =  $\text{Al}_2\text{O}_3$*

*Condition = 'passing vapour over hot solid' owtte*

1

Conditions: 180 °C / High temp / Hot / Reflux /

*ALLOW stated temp in range 100-300 °C/373-573 K*

*IGNORE 'heat'*

*M4 dependent on correct reagent in M3*

1

(d) (i) **S** = aldehyde/CHO **AND** **T** = carboxylic/COOH/CO<sub>2</sub>H

1

T forms hydrogen bonds

1

(Which are) stronger than / need more energy to break than forces between molecules/IMFs in S ora (or reverse argument)

*If implication of breaking covalent bonds max M1 only*

1

(ii) (No oxidation has occurred as..)

(Still) contains peak at 3230–3550 cm<sup>-1</sup> due to O–H/alcohol

Does not contain peak at 2500–3000 cm<sup>-1</sup> due to  
O–H/carboxylic acid

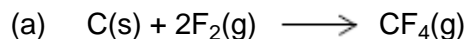
Does not contain peak at 1680–1750 cm<sup>-1</sup> due to C=O

*Must have wavenumber range (or value within range) and bond or  
functional group to score each mark.*

Any 2

[13]

7



*State symbols essential*

1

(b) Around carbon there are 4 bonding pairs of electrons (and no lone pairs)

1

Therefore, these repel equally and spread as far apart as possible

1

(c)  $\Delta H = \sum \Delta_f H \text{ products} - \sum \Delta_f H \text{ reactants}$  or a correct cycle

1

$$\text{Hence } = (2 \times -680) + (6 \times -269) - (x) = -2889$$

1

$$x = 2889 - 1360 - 1614 = -85 \text{ (kJ mol}^{-1}\text{)}$$

1

*Score 1 mark only for +85 (kJ mol<sup>-1</sup>)*

(d) Bonds broken = 4(C–H) + 4(F–F) = 4 × 412 + 4 × F–F

Bonds formed = 4(C–F) + 4(H–F) = 4 × 484 + 4 × 562

*Both required*

1

$$-1904 = [4 \times 412 + 4(F-F)] - [4 \times 484 + 4 \times 562]$$

$$4(F-F) = -1904 - 4 \times 412 + [4 \times 484 + 4 \times 562] = 632$$

1

$$F-F = 632 / 4 = 158 \text{ (kJ mol}^{-1}\text{)}$$

1

The student is correct because the F–F bond energy is much less than the C–H or  
other covalent bonds, therefore the F–F bond is weak / easily broken

*Relevant comment comparing to other bonds*

*(Low activation energy needed to break the F–F bond)*

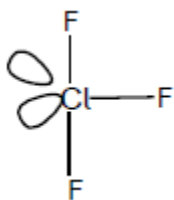
1

[10]

8	B		[1]
9	A		[1]
10	C		[1]
11	C		[1]
12	D		[1]
13	B		[1]
14	C		[1]
15	C		[1]
16	(a) Silicon / Si		
		<i>If not silicon then CE = 0 / 3</i>	1
		<u>covalent</u> (bonds)	
		<i>M3 dependent on correct M2</i>	1
		Strong or many of the (covalent) bonds need to be <u>broken</u> / needs a lot of energy to <u>break</u> the (covalent) bonds	
		<i>Ignore hard to break</i>	1
	(b) Argon / Ar		
		<i>If not argon then CE = 0 / 3. But if Kr chosen, lose M1 and allow M2+M3</i>	1
		Large(st) number of protons / large(st) nuclear charge	
		<i>Ignore smallest atomic radius</i>	1
		Same amount of shielding / same number of shells / same number of energy levels	
		<i>Allow similar shielding</i>	1
	(c) Chlorine / Cl		
		<i>Not Cl<sub>2</sub>, Not CL, Not Cl<sup>2</sup></i>	1

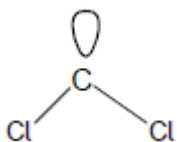


(d) (i)



Or any structure with 3 bonds and 2 lone pairs  
Ignore any angles shown

1



Or a structure with 2 bonds and 1 lone pair

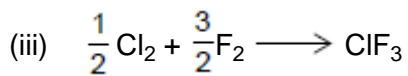
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(ii) Bent / v shape

Ignore non-linear, angular and triangular

Apply list principle

1



No multiples

Ignore state symbols

1

[11]

17

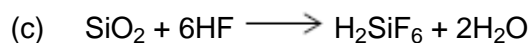
(a) Macromolecular / giant covalent / giant molecule

Not giant atomic

1

(b) No delocalised electrons / no free ions / no free charged particles

1



Accept multiples

1

[3]

18

(a) 94–105.5°

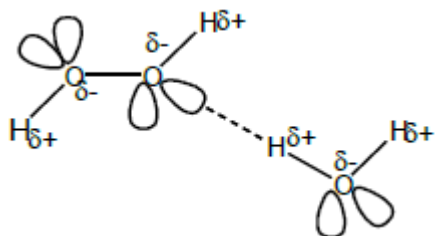
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(b) (i) Hydrogen bond(ing) / H bonding / H bonds

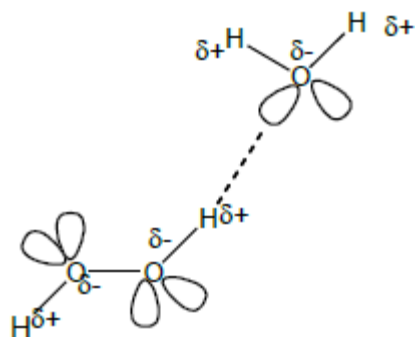
Not just hydrogen

1

(ii)



OR



1 mark for all lone pairs

1 mark for partial charges on the O and the H that are involved in H bonding

1 mark for the H-bond, from  $H\delta^+$  on one molecule to lone pair on O of other molecule

3

- (c) Electronegativity of S lower than O or electronegativity difference between H and S is lower

Mark independently

1

No hydrogen bonding between  $H_2S_2$  molecules

Or only van der Waals / only dipole-dipole forces between  $H_2S_2$  molecules

If breaking covalent bonds CE = 0

1

[7]