# **A-Level Chemistry**

# Paper 4

**Unsolved Topical** 

**Past Papers with Marking Schemes** 

**All Variants** 

2014-2021

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#### A-LEVEL TOPICAL CHEMISTRY PAPER 4

**Published by** 

MS Books (042-35774780)

**Legal Advisor** 

Ashir Najeeb Khan (Advocate) **AKBAR LAW CHAMBERS** 

39-40, 1<sup>st</sup> Floor, Sadiq Plaza, The Mall, Lahore.

0307-4299886, 042-36314839

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**MS Books** 

83-B Ghalib Market, Gulberg III Lahore

(042-35774780),(03334504507),(03334548651)

**Price** 

Rs.

#### **PREFACE**

Excellence in learning cannot be claimed without application of concepts in a dexterous way. In this regard one of the logical approach is to start in chunks; like chapter wise learning and applying the concept on exam based questions.

This booklet provides an opportunity to candidates to practice topic wise questions from previous years to the latest. Extensive working of Team MS Books has tried to take this booklet to perfection by collaborating with top of the line teachers.

We have added answer key / marks scheme at the end of each topic for the candidate to compare the his/her answer to the best.

MS Books strives to maintain actual spacing between consecutive questions and within options as per CAIE format which gives students a more realistic feel of attempting question.

Review, feedback and contribution in this booklet by various competent teachers of a subject belonging to renowned school chains make it most valuable resource and tool for both teachers and students.

With all belief in strength of this resource material I can confidently claim that it is worth in achieving brilliance.

Our sincere thanks and gratification to Mr. Waqar Ahmad who took out special time to help compile and manage this booklet. We would also like to appreciate chemistry faculty for reviewing and indorsing it.

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## **CHEMISTRY PAPER 4**

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# **Enthalpy and Born-Haber Cycles**

Q2(d)/42/M/J/14

- **1 (d) (i)** Write an equation to represent the lattice energy of PbC  $l_2$ . Show state symbols.
  - (ii) Use the following data, together with appropriate data from the *Data Booklet*, to calculate a value for the lattice energy of  $PbCl_2$ .

electron affinity of chlorine =  $-349 \,\mathrm{kJ} \,\mathrm{mol}^{-1}$ enthalpy change of atomisation of lead =  $+195 \,\mathrm{kJ} \,\mathrm{mol}^{-1}$ enthalpy change of formation of PbC  $l_2(\mathrm{s})$  =  $-359 \,\mathrm{kJ} \,\mathrm{mol}^{-1}$ 

	lattice energy =kJ mol <sup>-1</sup>
iii)	How might the lattice energy of $PbCl_2$ compare to that of $PbBr_2$ ? Explain your answer.

Q2(c)/43/M/J/14

2 (c) Copper(I) oxide and copper(II) oxide can both be used in the ceramic industry to give blue, green or red tints to glasses, glazes and enamels.

The table lists the  $\Delta H_f^e$  values for some compounds.

compound	$\Delta H_{\rm f}^{\rm e}/{\rm kJmol^{-1}}$
Cu <sub>2</sub> O(s)	-168.6
CuO(s)	-157.3
Cu(NO <sub>3</sub> ) <sub>2</sub> (s)	-302.9
NO <sub>2</sub> (g)	+33.2

(i) Copper(II) oxide can be produced in a pure form by heating copper(II) nitrate. Use suitable  $\Delta H_f^{\bullet}$  values from the table to calculate the  $\Delta H^{\bullet}$  for this reaction.

$$Cu(NO_3)_2(s) \rightarrow CuO(s) + 2NO_2(g) + \frac{1}{2}O_2(g)$$

 $\Delta H^{\circ} = \dots kJ \, \text{mol}^{-1}$ 

- (ii) Copper(I) oxide can be produced from copper(II) oxide.
  - Use suitable  $\Delta H_f^{e}$  values from the table to calculate  $\Delta H^{e}$  for the reaction.

$$2CuO(s) \iff Cu_2O(s) + \frac{1}{2}O_2(g)$$

•	Hence suggest whether a low or a high temperature of oxidation would favour the production of copper(I) oxide. Explain your reasoning.

Q3(	(م	/42	10	/N/	14
QU,	С,	// <del>+</del> _	$\cup$	1 1/	14

3	(e)	(i)	What is meant by the term lattice energy?
		(ii)	Explain why the lattice energy of calcium phosphate is <b>less</b> exothermic than that of magnesium phosphate.
			[3]

Q1(b(ii)/43/O/N/14

**4 (b) (ii)** Use the following data, together with relevant data from the *Data Booklet*, to calculate a value for the lattice energy of strontium chloride. You may find it helpful to construct a Born-Haber cycle.

electron affinity per mole of chlorine atoms	−349 kJ mol <sup>-1</sup>
standard enthalpy of atomisation of Sr(s)	+164 kJ mol <sup>-1</sup>
standard enthalpy of formation of SrCl <sub>2</sub> (s)	-830 kJ mol⁻¹

	(0.0)	lattice energy =kJ mol-
Q1(d,e)/42		
5 (d)	(i)	What is meant by the term standard enthalpy change of hydration, $\Delta H_{\text{hyd}}^{\bullet}$ ?
		[2]

Q2(d)/42/M/J/14

**1** (d) (i) 
$$Pb^{2+}(g) + 2Cl^{-}(g) \longrightarrow PbCl_{2}(s)$$
 [1]

(ii) 
$$\Delta H_f = \Delta H_{at} + E(Cl - Cl) + 1^{st} IE + 2^{nd} IE + 2 \times E_A(Cl) + LE$$
  
 $-359 = 195 + 242 + 716 + 1450 - 2 \times 349 + LE$   
 $LE = 2 \times 349 - 359 - 195 - 242 - 716 - 1450$   
 $LE = -2264 \text{ (kJ mol}^{-1})$  [3]

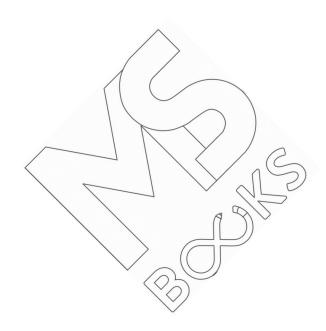
(iii) 
$$LE(PbCl_2) > LE(PbBr_2)$$
 or more exothermic or stronger lattice [1]

[6]

Q2(c)/43/M/J/14

**2** (c) (i) 
$$\Delta H^{\text{e}} = +2 \times 33.2 - 157.3 + 302.9 = (+) 212 \text{ kJ mol}^{-1} \text{ ecf}$$
 [2]

(ii) 
$$\Delta H^{\text{e}} = -168.6 + 2 \times 157.3 = (+)146 \text{ kJ mol}^{-1}$$
 allow ecf from (c)(i) [1] high T/temperature since  $\Delta H$  is positive/endothermic [1]



Q3(e)/42/O/N/14 Q 3

(e) (i)	(enthalpy change) when 1 mole of an ionic compound is formed from its gaseous ions		
(ii)	${\rm Mg^{2^+}}$ has a smaller (ionic) radii than ${\rm Ca^{2^+}}$ <b>OR</b> ${\rm Mg^{2^+}}$ is smaller than ${\rm Ca^{2^+}}$	1	[3]

Q1(b(ii)/43/O/N/14 Q4

[6]
111
(ii) Lattice energy = $\Delta H_{\rm f}({\rm SrC} I_2) - (\Delta H_{\rm atom}({\rm Sr}) + \Delta H_{\rm i1}({\rm Sr}) + \Delta H_{\rm i2}({\rm Sr}) + \Delta H_{\rm atom}({\rm C} I) + 2\Delta H_{\rm ea}({\rm C} I))$ = +(-830) - (+ 164 + 548 + 1060 + 242 + (2 × -349)) = -2146 (kJ mol <sup>-1</sup> )

Q1(d,e)/42/O/N/15 Q 5

(i) (b)	(energy change when) 1 mole of ions	2
	gaseous (ions) dissolve in water (to form an infinitely dilute solution) or gaseous (ions) form an aqueous solution	
(II)	(ii) $\Delta H_{\text{latt}}^{\text{fatt}} \text{Ca(NO}_3)_2 + \Delta H_{\text{sol}}^{\text{Ca(NO}_3)_2} = \Delta H_{\text{hyd}}^{\text{hyd}} \text{Ca}^{2^+} + 2\Delta H_{\text{hyd}}^{\text{hyd}} \text{NO}_3^-$ $\Delta H_{\text{latt}}^{\text{latt}} - 19 = -1650 + (2x - 314)$ $-2259 \text{ kJ mol}^{-1}$	က
(a)	Ca <sup>(2+)</sup> is a smaller(lion) <b>or</b> Ca <sup>(2+)</sup> has a larger charge density Ca <sup>(2+)</sup> has a stronger attraction/bond to H <sub>2</sub> O	7