Lattice energy - 2021

(b)

- 1. Nov/2020/Paper_41/No.2
 - (a) The lattice energies of three ionic compounds are given.

| compound | lattice energy/kJ mol ⁻¹ |
|----------|-------------------------------------|
| LiF(s) | -1022 |
| CaO(s) | -3513 |
| SrO(s) | -3310 |

| Define the term lattice energy. |
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| |
| [2 |
| Explain why the lattice energy of CaO is more exothermic than the lattice energy of LiF. |
| |
| [1 |
| Use the data in the table to estimate approximate values for the lattice energies of magnesium oxide and barium oxide. |
| $\Delta H_{\text{latt}} MgO(s) = \dots kJ mol$ |
| $\Delta H_{\text{latt}} \text{BaO}(\text{s}) = \dots \text{kJ mol-}$ |
| Write an equation for the reaction between BaO and $\rm H_2O$. Include state symbols. |
| |

| (ii) | State and explain how the solubilities of the hydroxides of the Group 2 elements vary down the group. |
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| (c) | | the following data and relevant data from the <i>Data Booklet</i> to calculate a value for the ce energy of magnesium fluoride, MgF ₂ (s). |
|-----|------|--|
| | | might find it helpful to construct an energy cycle. ow your working. |
| | | electron affinity of F(g) = $-348 \text{kJ} \text{mol}^{-1}$ enthalpy change of atomisation of Mg(s) = $+147 \text{kJ} \text{mol}^{-1}$ enthalpy change of formation of MgF ₂ (s) = $-1102 \text{kJ} \text{mol}^{-1}$ |
| | | $\Delta H_{\text{latt}} \text{MgF}_2(\mathbf{s}) =$ |
| (d) | (i) | $\Delta H_{\text{latt}} \text{MgF}_2(\mathbf{s}) =$ |
| | | [2] |
| | (ii) | The electron affinity of carbon, C(g), is -120 kJ mol ⁻¹ . |
| | | Suggest an explanation for the difference between the electron affinity of fluorine and the electron affinity of carbon. |
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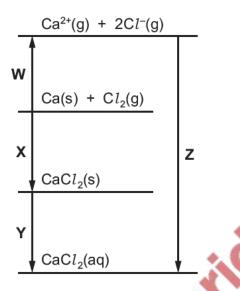
......[1]

[Total: 15]

2. Nov/2020/Paper_42/No.3

(a) The energy cycle shown can be used, along with suitable data, to calculate the enthalpy change of hydration of Ca²⁺(g).

Each arrow indicates a transformation, **W**, **X**, **Y** and **Z**. Each transformation consists of one or more steps.



The following data and data from the Data Booklet should be used.

electron affinity of Cl(g) = $-349 \,\mathrm{kJ}\,\mathrm{mol}^{-1}$ enthalpy change of atomisation of Ca(s) = $+193 \,\mathrm{kJ}\,\mathrm{mol}^{-1}$ enthalpy change of formation of $CaCl_2(s)$ = $-795 \,\mathrm{kJ}\,\mathrm{mol}^{-1}$ enthalpy change of solution of $CaCl_2(s)$ = $-83 \,\mathrm{kJ}\,\mathrm{mol}^{-1}$ enthalpy change of hydration of $Cl^-(g)$ = $-364 \,\mathrm{kJ}\,\mathrm{mol}^{-1}$

(i) Calculate the value of the enthalpy change corresponding to transformation **W**. Show your working.



enthalpy change $W = \dots kJ \text{ mol}^{-1}$ [2]

| (ii) | Use your answer to (a)(i) and other data to calculate the value of the enthalpy change corresponding to transformation Z . |
|-------|---|
| | |
| | enthalpy change Z = kJ mol ⁻¹ [2] |
| (iii) | Use your answer to (a)(ii) to calculate the enthalpy change of hydration of Ca ²⁺ (g). |
| | |
| | enthalpy change of hydration of Ca ²⁺ (g) =kJ mol ⁻¹ [2] |
| (iv) | Write an expression, in terms of \mathbf{W} , \mathbf{X} , \mathbf{Y} and/or \mathbf{Z} , to show how the enthalpy changes of \mathbf{two} of the transformations can be used to calculate the lattice energy of $\mathrm{CaC} l_2(\mathbf{s})$. |
| | lattice energy of $CaCl_2(s) =$ [1] |
| (v) | State whether the lattice energy of $CaCl_2(s)$ is more or less exothermic than the lattice energy of $MgF_2(s)$. |
| | Explain your answer. |
| | |
| | [1] |
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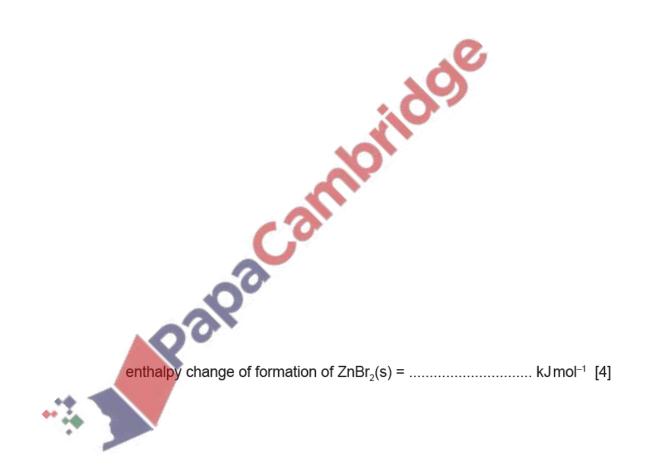
| (b) | The | e sulfates of the Group 2 elements vary in solubility down Group 2. |
|-----|------|---|
| | (i) | Give the names of two solutions that could be mixed to form barium sulfate. |
| | | [1] |
| | (ii) | State and explain how the solubilities of the sulfates of the Group 2 elements vary down Group 2. |
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| | | [4] |
| | | [Total: 13] |
| | | |

3. June/2020/Paper_42/No.7b,7c

(b) Use the following data and relevant data from the *Data Booklet* to calculate a value for the enthalpy change of formation of zinc bromide, ZnBr₂(s).

You might find it helpful to construct an energy cycle.

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electron affinity of Br(g) = -325 \,\mathrm{kJ} \,\mathrm{mol}^{-1}
enthalpy change of atomisation of Zn(s) = +131 \,\mathrm{kJ} \,\mathrm{mol}^{-1}
enthalpy change of vaporisation of Br<sub>2</sub>(l) = +31 \,\mathrm{kJ} \,\mathrm{mol}^{-1}
lattice energy of ZnBr<sub>2</sub>(s) = -2678 \,\mathrm{kJ} \,\mathrm{mol}^{-1}
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(c) The lattice energies of ${\rm ZnBr_2}$, ${\rm ZnC}\,l_2$ and ${\rm ZnO}$ are shown.

| compound | lattice energy/kJ mol ⁻¹ |
|-------------------|-------------------------------------|
| ZnBr ₂ | -2678 |
| ZnCl ₂ | -2734 |
| ZnO | -3971 |

| (i) | Explain why there is a difference between the lattice energies of ${\sf ZnBr_2}$ and ${\sf ZnCl_2}$. | |
|------|--|-------|
| | | . [1] |
| (ii) | Explain why there is a difference between the lattice energies of ${\rm ZnC}\it{l}_{2}$ and ${\rm ZnO}.$ | |
| | | . [1] |
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| | Palpa | |
| | | |