

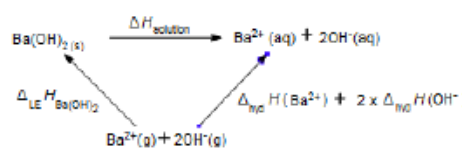
## Lattice Energy - Mark Scheme

Q1.

Question number	Answer	Additional guidance	Mark
(a)(i)	<ul style="list-style-type: none"> <li>• <math>R = O^{2-}(g)</math> (1)</li> <li>• <math>W =</math> first electron affinity <math>O(g)</math> (1)</li> </ul>	Allow alternative ways to express electron affinity, e.g. EA State required Do not allow $O_2/O^-$	2

Question number	Answer	Additional guidance	Mark
(a)(ii)	<ul style="list-style-type: none"> <li>• correct application of cycle (1)</li> <li>• correct value (1)</li> <li>• correct sign and units (1)</li> </ul>	Example of calculation: $\Delta_f H (BaO(s)) = \Delta_{at} H(Ba(s)) + \Delta_{at} H(\frac{1}{2}O_2(g)) + I^{st} IE (Ba)(g) + 2^{nd} IE (Ba(g)) + 2^{nd} EA (O(g)) + 1^{st} EA (O(g)) + \Delta_{LE} H (BaO(s))$ or Correct numbers $= 180.0 + 249.2 + 503 + 965 + 798 - 141.1 - 3054$ $= -499.9 / (-)500 (kJ mol^{-1})$ Allow TE from incorrect application of cycle Allow TE for incorrect numbers Correct answer with no working scores 3	3

Question number	Answer	Additional guidance	Mark
(a)(iii)	<ul style="list-style-type: none"> <li>• ionic radius of <math>Ba^{2+} \gg</math> ionic radius of <math>Mg^{2+}</math> / (have) lower charge density <b>and</b> <math>Ba^{2+}</math> (ions are) less polarising / (have) lower charge density (1)</li> <li>• iodide ions / <math>I^-</math> are large <b>and</b> their electron clouds are easily distorted / polarised (by Group 2 cations) or oxide ions / <math>O^{2-}</math> are small(er) <b>and</b> their electron clouds are less easily distorted / polarised (1)</li> <li>• more distortion / covalency leads to greater difference between theoretical and experimental values</li> </ul>	Allow reverse argument	3

Question number	Answer	Additional guidance	Mark
(b)(i)	 <ul style="list-style-type: none"> <li>all arrows in the correct direction (1)</li> <li>correct formulae at each corner and enthalpies of hydration, and solution and LE correctly identified (1)</li> <li>correct expression or correct substitution of values (1)</li> <li>correct evaluation (1)</li> </ul>	<p>Do not allow energy profile or energy level diagrams</p> <p>Species at each corner must be approximately correct</p> <p>Allow missing minor detail: brackets, position of subscripts, etc. but not absence of subscripts</p> <p>Example of calculation:  <math>\Delta_{\text{sol}}H = (\Delta_{\text{hyd}}H(\text{Ba}^{2+}) + 2\Delta_{\text{hyd}}H(\text{OH}^-)) - \text{LE}(\text{Ba}(\text{OH})_2)</math>  or  <math>= (-1360 + (2 \times -460)) - (-2230)</math>  <math>= -50 \text{ (kJ mol}^{-1}\text{)}</math>  Allow TE from their cycle if <math>\Delta_{\text{hyd}}H(\text{OH}^-)</math> is not doubled</p>	4

Question number	Answer	Additional guidance	Mark
b(ii)	<ul style="list-style-type: none"> <li>entropy (change) of system / <math>\Delta S_{\text{system}}</math> / <math>\Delta S_{\text{dissolving}}</math> is large and positive (and outweighs negative <math>\Delta S_{\text{surroundings}}</math> (<math>-\Delta H/T</math>)) (1)</li> <li>overall entropy change / <math>\Delta S_{\text{total}}</math> is positive (1)</li> <li>or</li> <li>use of <math>\Delta S_{\text{total}} = \Delta S_{\text{surroundings}} + \Delta S_{\text{system}}</math> (1)</li> <li><math>\Delta S_{\text{total}} = \Delta S_{\text{surroundings}} + (-\Delta H/T)</math> (1)</li> </ul>	Allow use of $\Delta G$	2