

Session #9: Homework Problems

Problem #1

Estimate the ionic radius of Cs^+ . The lattice energy of CsCl is 633 kJ/mol . For CsCl the Madelung constant, M , is 1.763 , and the Born exponent, n , is 10.7 . The ionic radius of Cl^- is known to be 1.81 \AA .

Problem #2

- CFCs have been implicated in ozone depletion. Show that when Freon 12 (CCl_2F_2) is exposed to ultraviolet radiation, the compound decomposes to produce chlorine.
- Draw the Lewis structure of Freon 12 and indicate the polarities of each bond within this compound.
- Determine the percent ionic character of the C--Cl and C--F bonds.

DATA:

Average Bond Energies (kJ/mol)

Single Bonds

H-H	435
F-F	155
Cl-Cl	242
C-C	347

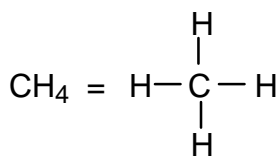
Multiple Bonds

C = C	610
C \equiv C	836

Problem #3

The compound trichloroethylene has the composition $\text{C}_2\text{Cl}_3\text{H}$.

- Formulate this compound in "Lewis" notation. Example:



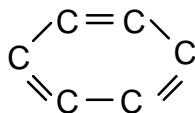
- List the types of atomic orbitals which, upon overlap, lead to the formation of the bonds indicated in (a).
- List all bonds (number and type) involved in the formation of this compound.

Problem #4

Frequently the statement is made that the crystal energy (ΔE_{cryst}) is predominantly given by the attractive energy term. How can you account for this approximation?

Problem #5

For benzene (C_6H_6), which has a schematic *carbon* arrangement?



- Determine the total number and types of covalent bonds.
- For each bonding type, determine the bonding orbitals which, by overlap, lead to their formation.

Problem #6

Why is double bonding encountered in some carbon compounds, but not in germanium compounds – although both exhibit sp^3 hybridization?

Problem #7

List the individual steps with the corresponding chemical equations used in constructing a Born-Haber cycle for the formation of $CaBr_2$ from the element and identify those which you expect to be exothermic.

Problem #8

- Given the ionic radii, $Cs^+ = 1.67\text{\AA}$, $Cl^- = 1.81\text{\AA}$, and the Madelung constant $M(CsCl) = 1.763$, determine to the best of your ability the molar Crystal energy (ΔE_{cryst}) for $CsCl$.
- Not given additional data, do you expect the value obtained to be larger or smaller than theoretical, and by how many percent do you anticipate to be off?

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