

Supporting Information

Syntheses and Structures of Two Chiral Zincophosphite Compounds

[Zn(C₈H₈N₂)(HPO₃)] and (C₆H₁₃N₂) [Zn₃(C₆H₁₂N₂)(HPO₃)₃(H₂PO₃)]

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Fig. 1 IR spectra for compound **1(a)** and **2(b)**.

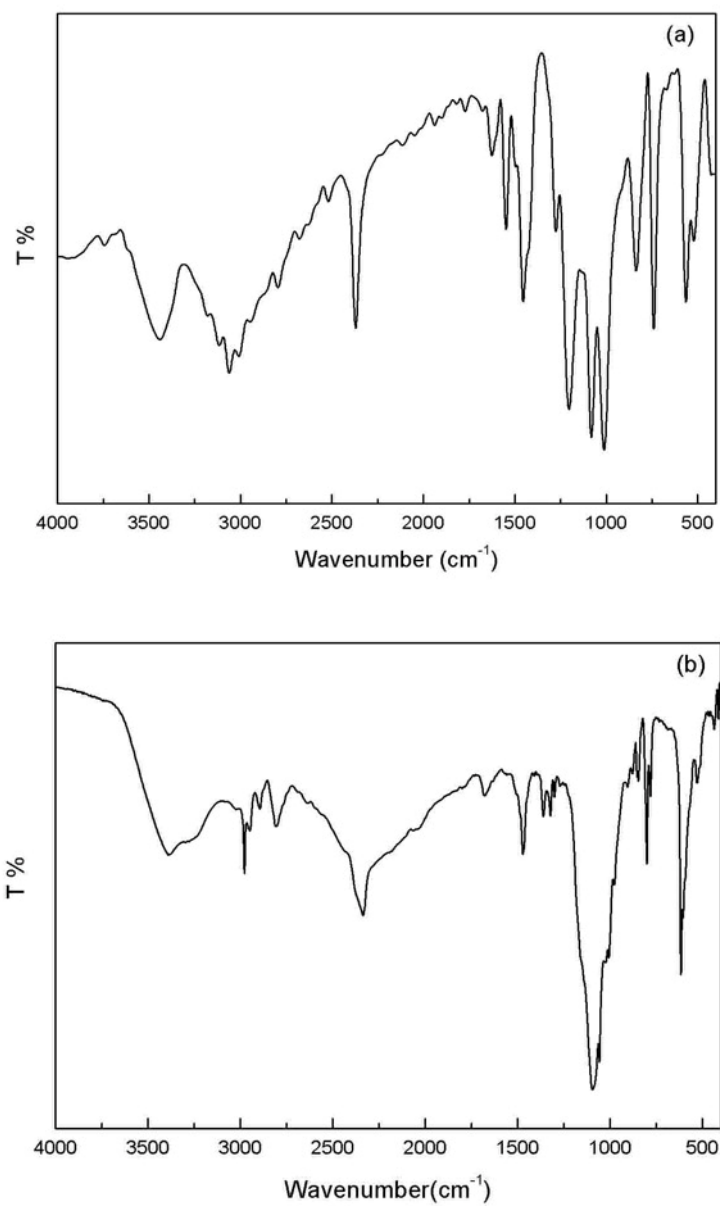


Fig. 2 TG curves for compound **1(a)** and **2(b)**.

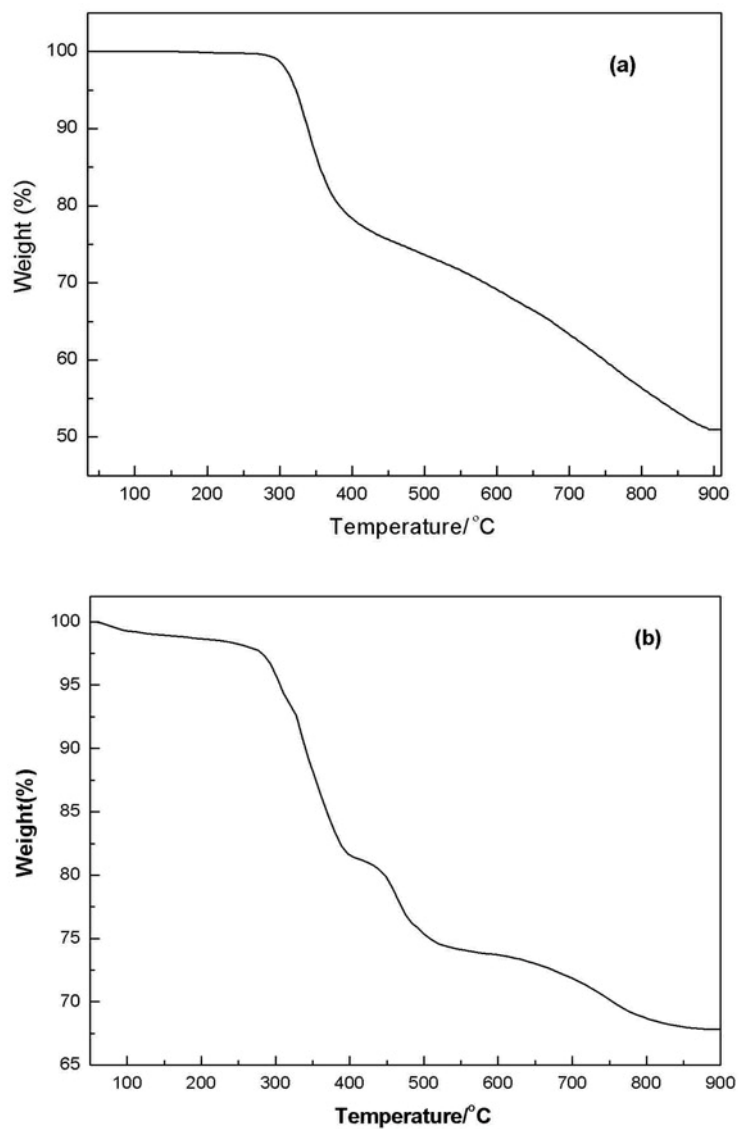
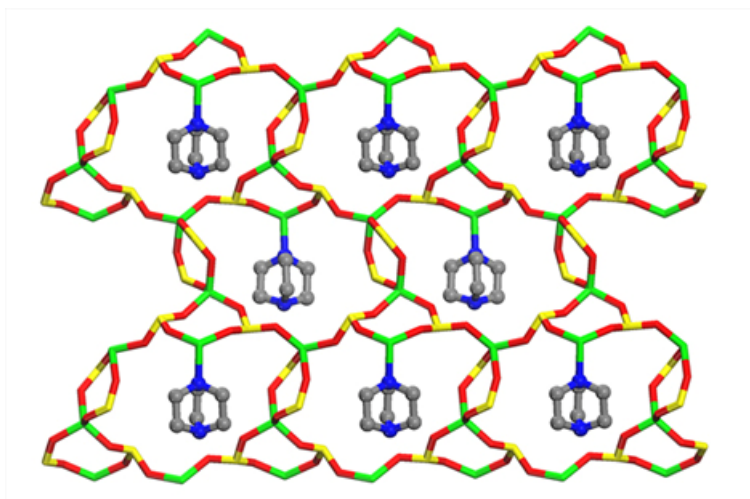
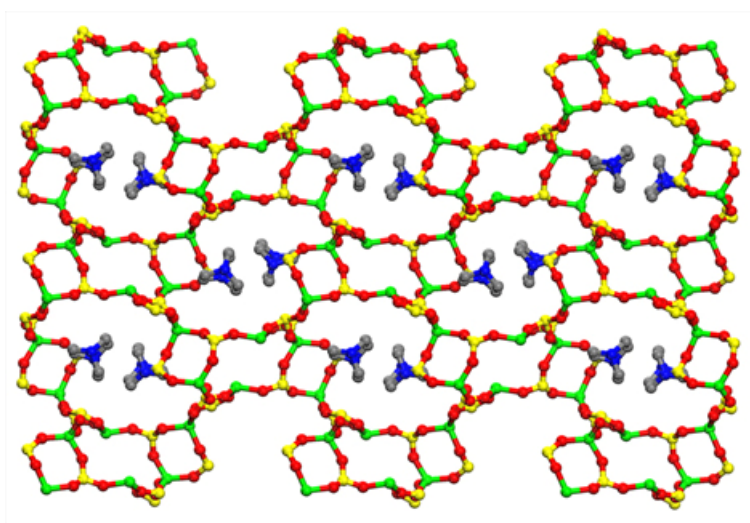


Fig. 3 In compound **2**, **(a)** bonded DABCO molecule extending from Zn(3) atoms occupies the 12-ring window, **(b)** free monoprotonated DABCO molecule resides in the 16-ring channel and compensates the charge of framework. (All the H atoms are omitted for clarity).



(a)



(b)

Fig. 4 The 3D framework structure of **2** viewed along the *b* axis. (All the H atoms are omitted for clarity).

