

# Molecular Structure and Thermal Analysis of Potassium Hydrogenphthalate Monohydrate

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Single crystals of potassium hydrogenphthalate monohydrate, KHPht( $\text{H}_2\text{O}$ ), were grown from aqueous solution using a slow cooling method and the structure was determined by X-ray diffraction analysis. The crystals belong to the monoclinic space group  $P2_1/c$ . Its unit cell parameters are as follows:  $a = 1.1235(1)$ ,  $b = 0.6689(1)$ ,  $c = 1.1998(2)$  nm,  $\beta = 98.85^\circ$ ,  $V = 0.8909(1)$  nm<sup>3</sup>,  $D_c = 1.657$  g/cm<sup>3</sup>,  $Z = 4$ ,  $F(000) = 456$ . The thermal decomposition of the complex was studied using differential scanning calorimetry (DSC), thermogravimetry-derivative thermogravimetry (TG-GTG) and FT-IR techniques. With a linear heat rate, the diagrams show three endothermic processes. RCOOK and  $\text{K}_2\text{CO}_3$  are produced at 330 and 467 °C, respectively, according to FT-IR analysis.

*Key words:* Potassium Hydrogenphthalate Monohydrate, Molecular Structure, Thermal Analysis