

Figure 1. Structure of cronstedtite- $6T_o$, side view, projection close to a, For sake of clarity, only a small part of every OD packet (1:1 layer) is displayed: one ring of tetrahedra and three adjacent octahedra. Delimitations of packets $(P_o, P_1, P_2, ...)$ are indicated on the right side.

Keywords: Cronstedtite, 1:1 layer silicate, polytypism, non-MDO polytype 6T2, crystal structure

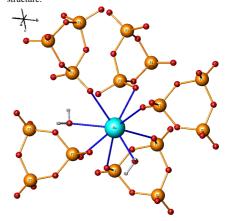
$\frac{\text{MS15-P12}}{\text{Chemical preparation,}} \label{eq:ms15-P12} \text{Chemical preparation,} \\ \text{crystallographic characterization and} \\ \text{vibrational study of condensed phosphates} \\ \text{associated to Barium-Cesium} \\ \text{BaCs}(P_3O_9)_2.2H_2O$

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Methods of chemical preparation and XRD data are reported for the new condensed phosphates associated to Barium-Cesium BaCs(P_Q_0)₂,2H,O. BaCs(P_Q_0)₂,2H,O was prepared by the method of ion-exchange resin. This salt crystallizes in the monoclinic system, space group P21/n a = 7.6992(2) Å b = 12.3237(3) Å c = 11.8023(3) Å, β =101.181°(3)., M(20) = 1313,35; F (20) = 1004,53 and V = 333,95(2) (Å3), the vibrational study by IR absorption spectroscopy of the title compound reveals the presence of three bands and confirm the existence of non-equivalent positions of water molecules in the structure.



 $\begin{array}{llll} \textbf{Figure} & \textbf{1.} & \text{Projection} & \text{of} & \text{the} & \text{structure Barium-Cesium} \\ \text{BaCs}(P_3O_0)_2.2H_2O & & & \\ \end{array}$

Keywords: condensed phosphates, ion-exchange resin, vibrational study