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Synthesis and study of Lithium Triuranate $\text{Li}_2\text{U}_3\text{O}_{10}\cdot 6\text{H}_2\text{O}$

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In this work, a method of synthesis of lithium triuranate hexahydrate $\text{Li}_2\text{U}_3\text{O}_{10}\cdot 6\text{H}_2\text{O}$ is described. The chemical and functional composition of this compound has been investigated; its crystallographic characteristics have been determined; the state of H_2O and its role in the formation of the structure have been studied.

Synthesis of the investigated compound is a reaction of schoepite $\text{UO}_3\cdot 2.25\text{H}_2\text{O}$ with aqueous solution of lithium nitrate under hydrothermal conditions at 200°C. The synthesized lithium triuranate hexahydrate is an easily reproducible individual crystalline compound. The X-ray diffraction picture contains a series of reflections from planes with indices which, in combination with an intense reflection peak at $2\theta = 12.04^\circ$, indicate a typical layered structure of the triuranate.

For evaluation of its functional composition of $\text{Li}_2\text{U}_3\text{O}_{10}\cdot 6\text{H}_2\text{O}$, we have performed the IR spectroscopic research. The spectrum contains two groups of vibrations associated with H_2O and uranyl group. The vibrations of H_2O are very characteristic. The band of $\delta(\text{H}_2\text{O})$ vibrations at 1620 cm^{-1} is not split. Due to the participation of H_2O molecules in the formation of the branched system of H-bonds, the bands of vibrations ν_s and ν_a are represented in the spectra by a broad and intense band with faint maxima at 3511 and 3414 cm^{-1} . On the whole, all H_2O molecules in the IR spectrum of $\text{Li}_2\text{U}_3\text{O}_{10}\cdot 6\text{H}_2\text{O}$ retain their vibrational identity. The vibrations of the uranyl group are represented in the spectrum by the only band ν_3 at 917 cm^{-1} , which is typical for the seven-fold coordination of uranium(VI) in its uranium-oxygen polyhedron. The absence of the band allows us to consider the uranyl group as having a linear and the equal-shoulder configuration.

To specify the state of H_2O in $\text{Li}_2\text{U}_3\text{O}_{10}\cdot 6\text{H}_2\text{O}$ and to estimate its position in the structure, we have performed thermographic study. According to the first effect in the DTA curve at 162°C , the elimination of four H_2O molecules per $\text{Li}_2\text{U}_3\text{O}_{10}\cdot 6\text{H}_2\text{O}$ formula unit proceeds in a single stage. The elimination of the two additional H_2O molecules also proceeds in a single stage, but at a higher temperature 393°C . The dehydration process is completed at 393°C by the total destruction of the crystal structure and the transition into the amorphous state. The crystallization of $\text{Li}_2\text{U}_3\text{O}_{10}$ over wide time and temperature ranges occurs above 393°C .

Primary author: Ms KOSTROVA, Elena (Leonidovna)

Co-authors: Ms CHAPLIEVA, Kseniya (Alexsandrovna); Prof. CHERNORUKOV, Nicholay (Georgievich); Ms NIPRUK, Oxana (Valentinovna)

Presenter: Ms KOSTROVA, Elena (Leonidovna)

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