Ultrasonic investigation of fluorite-type BaF₂ :Cu and BaF₂ :Ni crystals

N. Yu. Ofitserova¹, M. N. Sarychev¹, I. V. Zhevstovskikh^{1,2}, V. A. Ulanov³, V. T. Surikov⁴, N. S. Averkiev⁵, V. V. Gudkov¹

 ¹Ural Federal University, Ekaterinburg, Russia
²M. N. Miheev Institute of Metal Physics, Ural Branch of Russian Academy of Sciences, Ekaterinburg, Russia
³ E. K. Zavoisky Physical Technical Institute, FRC Kazan Scientific Centre of Russian Academy of Sciences, Kazan, Russia
⁴Institute of Solid State Chemistry, Ural Branch of Russian Academy of Sciences, Ekaterinburg, Russia
⁵A. F. Ioffe Physical Technical Institute of Russian Academy of Sciences, St. Petersburg, Russia

The temperature dependences of the dynamic elastic moduli c_{β} in crystals of BaF₂ doped with Ni and Cu were studied with the use of ultrasonic technique. In BaF₂:Cu, anomalies were found which are typical for the Jahn-Teller effect (JTE) subject to the $T \otimes (e + t_2)$ problem with orthorhombic global minima of the adiabatic potential energy surface similar to what was observed in $CaF_2:Cu^{2+}$ [1]. However, the peak in $\text{Im}[\Delta c_{\beta}(T)/c_{\beta}(T_0)]$ and anomalies in $\text{Re}[\Delta c_{\beta}(T)/c_{\beta}(T_0)]$ (where $T_0 = 4$ K) were much less in BaF₂:Cu. In BaF₂:Ni, there were no visible signs of the JTE although the crystals BaF₂:Cu and BaF₂:Ni were colored meaning the presence of sufficient amount of the 3d ions. EPR studies of BaF₂:Cu [2] indicated the off-centre position of the Cu ion which do not have the JT origin. So, we conclude that only a small amount of the dopants does substitute metal in cation positions in BaF₂:Cu . In BaF₂:Ni, there is no enough such Ni ions to be noticed in an experiment. The technique for evaluation the content of the impurities gives the total amount of the dopant while evaluation of most of the parameters of the JTE using the experimental data requires concentration of the JT complexes. The exception is the temperature dependence of relaxation time which provides information about the parameters of relaxation mechanisms. Namely these parameters can be used for comparison of the properties of the JT complexes in different host crystals and discussed in our poster. The study was supported by the Ministry of Science and Higher Education of the Russian Federation (Ural Federal University Program of Development within the Priority-2030 Program).

[1] M.N. Sarychev et al.; JETP 135 473 (2022).

[2] M.M. Zaripov, V.A. Ulanov; *Fiz. Tverd. Tela* **31** 254 (1989) [*Sov. Phys. Solid State* **31** 1798 (1989)].