

## High pressure study of langasite $\text{Ba}_3\text{TaFe}_3\text{Si}_2\text{O}_{14}$ up to 60 GPa and low temperatures

Yu.A. Nikiforova<sup>1,2\*</sup>, I.S. Lyubutin<sup>1</sup>, A.G. Gavriiliuk<sup>1,2</sup>, I.A. Troyan<sup>1,2</sup>, A.G. Ivanova<sup>1,2</sup>, S.S. Starchikov<sup>1,2</sup>, S.N. Sul'yanov<sup>1</sup> and S.N. Axenov<sup>2</sup>

<sup>1</sup>Shubnikov Institute of Crystallography, FRSC «Crystallography and Photonics»,

Russian Academy of Sciences, Leninsky av., Moscow, 119333, Russia

<sup>2</sup>Institute for Nuclear Research, Russian Academy of Sciences, Phisicheskaya str., Troitsk, Moscow, 117312, Russia

Keywords: high pressure phase transitions, Synchrotron X-ray diffraction, langasites.

\*e-mail: juliadavudova@gmail.com

Langasite  $\text{La}_3\text{Ga}_5\text{SiO}_{14}$  gave the name to the big family of crystals, which have the unique piezoelectric properties in combination with luminescent, laser, and nonlinear optical properties. Recently, a strong interest was attracted to the langasite-type compounds containing magnetic ions [1]. Coexistence of electric and magnetic order parameters in these materials provide a new type of multiferroics.

In this work the high-pressure properties of a new multiferroic of the langasite family  $\text{Ba}_3\text{TaFe}_3\text{Si}_2\text{O}_{14}$  up to 60 GPa were investigated in diamond-anvil cells (DAC) by synchrotron X-ray diffraction (XRD). A preliminary X-ray analysis of the initial sample revealed that value of cell parameters is  $a = 8.538(2)$  Å and  $c = 5.237(2)$  Å. Two structural phase transitions at  $P \approx 5.5$  and 20 GPa were observed [2].

For the Mössbauer studies, iron in the samples was enriched in the  $^{57}\text{Fe}$  isotope up to 50%. Synchrotron Mossbauer source (SMS) technique were used in the temperature range of 4.2–295 K. Strong enhancement of the Néel temperature  $T_N$  was observed at pressures above 20 GPa and associated with the structural transformation. The highest value of  $T_N$  is about 130 K which is almost five times larger than the value at ambient pressure (about 27 K) [3]. Raman spectroscopy data also reveal anomalies in phonon and electronic properties of these compounds at about 5 and 20 GPa.

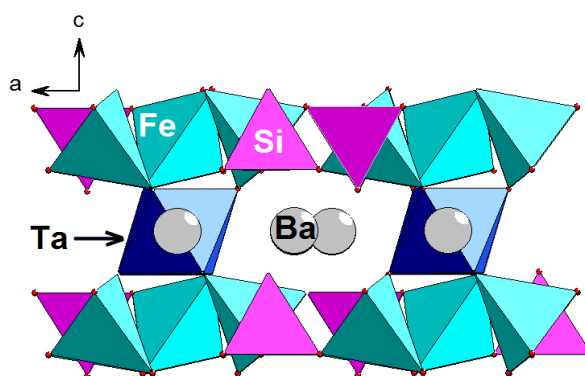


Figure 1. The unit cell of the iron langasite (sp. gr.  $P321$ ,  $Z=1$ ). All  $\text{Fe}^{3+}$  ions are in the tetrahedral oxygen sites

**Acknowledgments:** This work was supported by RFBR grant #17-02-00766 in part of Mossbauer spectroscopy measurements and spectra analysis and by the Ministry of Science and Higher Education within the State assignment FSRC “Crystallography and Photonics” RAS in part of preparation of the DAC. The SMS measurements were performed at the Nuclear Resonance beamline ID18 of ESRF. We are grateful to Dr. P. Liermann and Dr. K. Glazyrin for assistance at the P02.2 beamline of DESY (PETRA-III, Hamburg, Germany)

- [1] K. Marty, V. Simonet, E. Ressouche, R. Ballou, P. Lejay, P. Bordet, *Phys. Rev. Lett.* 2008, **101**, 247201.
- [2] Lyubutin, I. S.; Gavriilyuk, A. G.; Davydova, Yu. A.; Ivanova, A. G.; Troyan, I. A.; Sul'yanov, S. N.; Starchikov, S. S.; Aksenov, S. N.; Glazyrin, K. V., *Jetp Lett. (JETP Letters)*, 2015, **100**, 798
- [3] Lyubutin, I. S.; Starchikov, S. S.; Gavriiliuk, A. G.; Troyan, I. A.; Nikiforova, Yu. A.; Ivanova, A. G.; Chumakov, A. I.; Ruffer, R., *Jetp Lett. (JETP Letters)*, 2017, **105**, 26.