## Bi<sub>2</sub>Te<sub>2</sub>Se – promising material for topological insulators.

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Among  $Bi_2X_3$  family compounds  $Bi_2Te_2Se$  crystals are the most promising material for topological insulators. These ternary crystals have a layered structure with the atoms layers sequence of: Te - Bi - Se - Bi - Te. They are characterized by high resistivity at low temperatures around 10K  $\rho > 1 \Omega cm$ , [1,2] much higher than crystals of binary compounds as  $Bi_2Se_3$  and  $Bi_2Te_3 - (\rho \sim 10^{-3} \Omega cm)$ .

 $Bi_2Te_2Se$  is obtained by phase transformation in the solid state, which occurs as a result of long time heating the mixture of binary compounds  $Bi_2Se_3 + Bi_2Te_3$  at a specific temperature [3,4,5].

The mixture is formed by  $Bi_2Te_3 - p - type'$  material which properties are determined by high concentration of acceptor defects  $Bi_{Te}$  (p = 2 x 10<sup>19</sup> cm<sup>-3</sup>) and  $Bi_2Se_3 - n$ - type' material which properties resulting from the high concentration of donor defects  $V_{Se}$  (n = 2 x 10<sup>19</sup> cm<sup>-3</sup>).

The resulting new structure quintuples has the following advantages over the binary compounds  $Bi_2Te_3$  and  $Bi_2Se_3$  [1].

- Generation of V<sub>Se</sub> defects is suppressed because Se atom trapped between Bi atoms is less exposed to evaporation due to its stronger chemical binding with Bi,
- Formation of the anti-site defects is also suppressed because of the stronger bond between the Se - Bi than Te - Bi,
- Nature of bonds minimizes the additional disorder caused by random impact of Se / Te.

In this work we applied Vertical Bridgman technique (VB) for crystal growth. Crystallization was carried out for the charges compositions of a different ratio of Se / Te : Bi<sub>2</sub> Te<sub>1.95</sub> Se<sub>1.05</sub> to Bi<sub>2</sub> Te<sub>2.2</sub> Se<sub>0.8</sub>. As the optimum composition Se  $\approx 1.00 - 1.05$  was determined. Specified conditions of heat treatment, resulting in the obtaining of crystals with resistivity  $\rho \sim 7.5 \Omega$ cm at 10 K. After annealing under conditions of 500 °C/520h, we observe transition of conductivity type from *n* to *p*. Measurements of resistivity vs T have proved semiconductor behavior from the temperature T  $\leq 200$  K. The crystals were characterized by the following methods: Hall measurements in the temperature range 300K – 8K, Scanning Electron Microscopy (SEM), Energy Dispersive X– ray Spectroscopy (EDS), Secondary Ion Mass Spectroscopy (SIMS).X-ray Difractometry (XRD)

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