

## Poster Presentation

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*New intermetallic La<sub>117</sub>Ru<sub>57</sub>Sn<sub>112</sub>.*

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The new ternary compound with giant unit cell La<sub>117</sub>Ru<sub>57</sub>Sn<sub>112</sub> has been yielded by interaction between pure components lanthanum, ruthenium and tin and investigated by X-ray diffraction and scanning electron microscopy (SEM) in combination with energy dispersive X-ray spectroscopy (EDX) means. Previously, intermetallics with cobalt of similar composition and structure were found in the ternary systems and Dy-Co-Sn and Pr-Co-Sn – Dy<sub>117</sub>Co<sub>57</sub>Sn<sub>112</sub> and Pr<sub>117</sub>Co<sub>57</sub>Sn<sub>112</sub> respectively [1, 2]. The intermetallic La<sub>117</sub>Ru<sub>57</sub>Sn<sub>112</sub> crystallizes in a cubic Dy<sub>117</sub>Co<sub>57</sub>Sn<sub>112</sub> type structure with space group Fm-3m (No. 225) and lattice parameter  $a = 31.529(5)$ . A single-crystal suitable for the X-ray measurements was isolated from the of the equiatomic alloy La<sub>33.3</sub>Ru<sub>33.3</sub>Sn<sub>33.4</sub>. The structure was solved by means of direct methods and refined using the SHELXS-97 and SHELXL-97 programs ( $R_1 = 0.033$  for  $1042 F_o > 4\sigma(F_o)$  and  $0.061$  for all collected data). The additional sample of the La<sub>40.2</sub>Ru<sub>19.9</sub>Sn<sub>39.9</sub> composition was prepared and investigated X-ray by powder diffraction technique (CuK $\alpha$ 1 radiation,  $5 < 2\theta < 90^\circ$ ). The collected intensities data were refined by the Rietveld method using the La<sub>117</sub>Ru<sub>57</sub>Sn<sub>112</sub> single crystal model and FullProf program. Based on a minimum of differences between the observed and calculated theoretically intensities one can judge the good convergence results.

[1] P. Salamakha et al., *Journal of Alloys and compounds* 314 (2001) 177-180., [2] W. He et al., *Journal of Alloys and compounds* 491 (2010) 49-52.

**Keywords:** X-ray single-crystal diffraction, intermetallics