Erratum to “Lattice distortion and thermoelectric property for Zn_{1-(x+y)}Ga_xIn_yO system (x + y = 0.007, 0 ≤ x ≤ 0.007, 0 ≤ y ≤ 0.007)” [Transactions of the Materials Research Society of Japan 37 [3] 399-403 (2012)]

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Unfortunately, we found some errors in Fig. 5 and Fig. 6 which were displayed in the original contribution. We have wrongly put the values of the full width at half maximum (FWHM) in Fig 5. Now we have corrected the mistakes and the corrected Fig. 5 and Fig. 6 are shown below. The authors apologize for the inconvenience that this may have caused.

Fig.5 β – tanθ plot of Zn_{1-(x+y)}Ga_xIn_yO system by Rietveld analysis (x+y=0.007, 0≤x≤0.007, 0≤y≤0.007) for r_{av.} = 0.048 nm (●), 0.051 nm (■), 0.054 nm (●), 0.057 nm (▲), 0.060 nm (▼), Zn_{9.993}Ga_{0.007}O (△), Zn_{9.993}In_{0.007}O (○).

Fig.6 Relationship between mean ionic radius and lattice distortion (∆d/d) for Zn_{1-(x+y)}Ga_xIn_yO system (x+y=0.007, 0≤x≤0.007, 0≤y≤0.007).

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