

15th International Conference on Muon Spin Rotation, Relaxation and Resonance



Contribution ID: 214

Type: Poster

Local electronic structure of dilute hydrogen in gallium oxide

Tuesday, 30 August 2022 18:40 (20 minutes)

Gallium oxide $\beta\text{-Ga}_2\text{O}_3$ is currently drawing much attention as a material for high-voltage power devices because of its large band gap ($E_g \sim 4.9$ eV). We investigated the electronic structures of muon as a pseudo-hydrogen in $\beta\text{-Ga}_2\text{O}_3$ in which the dilute hydrogen is under the central focus as a crucial factor for the bulk conducting properties. We demonstrate by μSR study combined with the first principles calculations that muons in $\beta\text{-Ga}_2\text{O}_3$ have two electronic structures: a state corresponding to hydrogen that acts as an electron donor (Mu_1), bounded to three-coordinated oxygen, and a hydride-like state in rapid motion (Mu_2). Furthermore, we imply from the Hall effect measurements that the fractional yield of Mu_2 exhibits a close link with the mobility and density of the carriers (electrons).

In the presentation, we will compare our results with those of previous study and discuss the electronic structures of Mu_1 and Mu_2 in terms of our recently proposed ambipolar model of muon in oxides.

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Session Classification: Posters

Track Classification: Semiconductors