## **PCP5-3**

## Study of µSR in Iron-Based Superconductor LaFeAs<sub>1-x</sub>P<sub>x</sub>O<sub>0.9</sub>F<sub>0.1</sub>

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In the iron-based superconductors  $LaFeAs_{1-x}P_xO_{1-y}F_y$ , the electron doping level and the local crystal structure can be controlled by the F substitution for O and P substitution for As. With these chemical substitutions, Fermi surface (FS) topology changes giving three different superconducting (SC) phases [1]. For example, at y=0.1, the As-rich compounds are in the first superconducting phase (SC1), while the P-rich compounds are in the second superconducting phase (SC2) [2]. The theoretical study by Kuroki and coworkers has indicated that the different nesting in LaFeAsO-type and LaFePO-type FSs induces the different SC gap symmetries, i.e., full and nodal gaps [3].

In the present work, we have investigated the difference between SC gap symmetry in SC1 and SC2 using  $\mu$ SR measurement in LaFeAs<sub>1-x</sub>P<sub>x</sub>O<sub>0.9</sub>F<sub>0.1</sub> (x=0.0~0.8). The  $\mu$ SR measurement were performed at TRIUMF in Canada and Research Center for Nuclear Physics (RCNP), Osaka University in Japan using a He gas-flow cryostat in a magnetic field of 250G. At x=0, the temperature (*T*) dependence of the muon spin relaxation rate  $\sigma$  shows a rapid increase with decreasing *T* below *T*<sub>c</sub> and a saturation at low temperatures, indicating the s-wave behavior. In contrast, LaFeAs<sub>1-y</sub>P<sub>y</sub>O<sub>0.9</sub>F<sub>0.1</sub> (y=0.2~0.8) show the slightly different *T* dependence of the relaxation rate  $\sigma$ . In these P doping compounds, the *T* dependence of the relaxation rate  $\sigma$  does not show a clear saturation at low temperatures and cannot be fitted by the simple s-wave model. These results suggest that the P-doped compounds have several SC gaps with different gap sizes or a nodal SC gap, and the SC gap symmetries in the SC1 and SC2 phases may be different.

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