



**Experiment title: Connection between the nematic fluctuations and local structure in iron pnictide superconductors**

**Experiment number: HC-1021**

<b>Beamline:</b> BM-26A	<b>Date of experiment:</b> from: 29-01-2014 to: 04-02-2014	<b>Date of report:</b>
<b>Shifts:</b> 18	<b>Local contact(s):</b> Dipanjan Banerjee (email: banerjee@esrf.fr)	<i>Received at ESRF:</i>
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## Report:

During this beamtime, we have studied the local structure of a  $\text{Ba}(\text{Fe}_{1-x}\text{Co}_x)_2\text{As}_2$  superconductor using temperature dependent extended x-ray absorption fine structure (EXAFS) measurements. Polarized EXAFS at the Fe K-edge on an optimally doped ( $x = 0.06$ ) single crystal has permitted us to determine atomic displacements across the superconducting transition temperature ( $T_c$ ). The Fe–As bondlength is found to have hardly any change with temperature; however, the Fe–Fe bonds revealed a significant drop in mean square relative displacements across  $T_c$ . We have also found a large atomic disorder around the substituted Co, revealed by polarized Co K-edge EXAFS measurements. The Co–Fe/Co bonds are more flexible than the Fe–Fe bonds with the As-height in Co-containing tetrahedra being larger than the one in  $\text{FeAs}_4$ . The results suggest that the local Fe–Fe bondlength fluctuations and the atomic disorder in this sub-lattice should have some important role in the superconductivity of  $\text{Ba}(\text{Fe}_{1-x}\text{Co}_x)_2\text{As}_2$  pnictides.

**Publication: A Study of temperature dependent local atomic displacements in a  $\text{Ba}(\text{Fe}_{1-x}\text{Co}_x)_2\text{As}_2$**

**superconductor:** M. Y. Haciasalihoglu, E. Paris, B. Joseph, L. Simonelli, T. J. Sato, T. Mizokawa, N. L. Saini, in

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