Poster-1-3

Charge order above room temperature in kagome superconductor LaRu3Si2

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The present work focuses on LaRu3Si2, which has one of the highest critical temperatures (Tc = 7 K) among kagome superconductors. The structural motif of this material represents Rubased kagome layers placed between LaSi2 hexagonal nets. Our current single-crystal X-ray diffraction study of the pristine and Fe-doped LaRu3Si2 reveals a cascade of structural phase transitions with temperature, consistent with the previously predicted charge order instabilities. Above 600K, the material crystallizes in an undistorted P6/mmm structure (HT-hex), which doubles along the c-direction upon cooling (LT-hex). Below 400K, additional charge order (CO-I) reflections with $q1 = (1/4\ 0\ 0)$ evolve, followed by the second charge order transition into CO-II with $q2 = (1/6\ 0\ 0)$ below 180K. Apart from detailing the transition in reciprocal space, I will also describe the corresponding structural transformations observed in real space and discuss their potential implications on superconductivity.