

“Native point defects in $\text{CuIn}_{1-x}\text{Ga}_x\text{Se}_2$: hybrid density functional calculations predict the origin of p- and n-type conductivity”. Bekaert J, Saniz R, Partoens B, Lamoen D, Physical chemistry, chemical physics **16**, 22299 (2014). <http://doi.org/10.1039/c4cp02870h>