

“Validation of inter-atomic potential for WS₂ and WSe₂ crystals through assessment of thermal transport properties”. Mobaraki A, Kandemir A, Yapicioglu H, Gulseren O, Sevik C, Computational materials science **144**, 92 (2018).
<http://doi.org/10.1016/J.COMMATSCI.2017.12.005>