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**Half-Metallic Ferromagnetism in LaAlO<sub>3</sub>/SrMnO<sub>3</sub> Nanosheet Superlattices** FANG HOU, TIAN-YI CAI, SHENG JU, MING-RONG SHEN, Department of Physics, Soochow University, Suzhou, PR China — Based on first-principle density-functional theory, we have revealed a robust half-metallic ferromagnetism in LaAlO<sub>3</sub>/SrMnO<sub>3</sub> nanosheet superlattices. Interface electronic reconstruction, where electrons transfer from the (LaO)<sup>+</sup> layer to the adjacent (MnO<sub>2</sub>)<sup>0</sup> layer, is found to lead to the partially occupied e<sub>g</sub> orbitals at the Mn sites and the half-metallic state in nn-type superlattice via the Zener double-exchange mechanism. On the other hand, holes transfer from (AlO<sub>2</sub>)<sup>-</sup> layer to (SrO)<sup>0</sup> layer and reside mainly at oxygen sites in SrMnO<sub>3</sub>, leading to either the preserved G-type AFM ordering in pp-type superlattices or complex magnetic ordering in np-type superlattices. When these systems transit to ferromagnetic ordering by an external magnetic field, an obvious change of electronic states at the Fermi level is found, suggesting a large magnetoresistive effect therein.

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