

First-principles atomistic simulations of polarization patterns showing swirling structures in nanocomposites of dielectric BaZrO<sub>3</sub> nanoregions randomly distributed within a ferroelectric BaTiO<sub>3</sub> matrix. [R. Machado *et al.*, Phys. Rev. Lett. **136**, 126201 (2026)]

# NEWSPAPER

## PHYSICAL REVIEW LETTERS

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