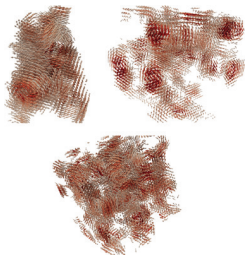

First-principles atomistic simulations of polarization patterns showing swirling structures in nanocomposites of dielectric $BaZrO_3$ nanoregions randomly distributed within a ferroelectric $BaTiO_3$ matrix. [R. Machado *et al.*, Phys. Rev. Lett. **136**, 126201 (2026)]

NEWSPAPER

PHYSICAL REVIEW LETTERS

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