

Molecular dynamics simulations of creatine kinase and adenine nucleotide translocase in mitochondrial membrane patch

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Molecular dynamics simulations on PGLa using NMR orientational constraints

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Molecular dynamics study of bending deformation in fivefold twinned silver nanowires

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Simulation of oriented NMR spectra : combining molecular dynamics and chemical shift tensor calculations

Sternberg, Ulrich; **Witter, Raiker** Magnetic resonance in chemistry 2024 / p. 125-144 <https://doi.org/10.1002/mrc.5403>

Transformations of gold nanocontacts studied by molecular dynamics simulations

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