

Mapping ...

... the protein conformational energy landscape is possible by using a combination of high-resolution nuclear magnetic resonance spectroscopy and accelerated molecular dynamics simulation. In their Communication on page 6103 ff., M. Blackledge and coworkers provide a detailed description of the different conformational sub-states of the interaction site of the SH3C domain of CD2 associated protein that are sampled on nanosecond to millisecond timescales.

