

Toward a unified theory for dynamics of glassy materials

April 22 2024



A scaling relation between the new theoretical order parameter $IS D_{min}$ and the relaxation damping phase angle (also known as internal friction in materials sciences). The data cover 5 orders of magnitude in timescale. Credit: Science China Press



In the realm of disorder and amorphous systems, such as oxide glasses utilized in display technologies and the cryogenic preservation of biological materials, there exists a substantial body of contemporary scientific and technological exploration.

A distinguishing feature of disordered materials is the presence of intricate dynamic behaviors, known as relaxation processes, which span from <u>atomic vibrations</u> on the picosecond timescale to aging and densification processes that can extend over thousands of years. These relaxation processes play a pivotal role in shaping the diverse properties of glassy materials.

Recent research in the field of glass science has brought to light a variety of specific dynamic phenomena within glassy materials, prompting researchers to seek a unifying principle that can elucidate these processes across a wide spectrum of materials.

Hai-Bin Yu from Huazhong University of China and Konrad Samwer from the University of Gottingen recognized the absence of a comprehensive theoretical framework for understanding relaxation dissipation in disordered systems. Their research is <u>published</u> in the journal *National Science Review*.

Rising to the challenge, they proposed a novel perspective to tackle this issue. While previous studies typically delved into the relaxation dynamics of individual particles within glassy materials, Yu and Samwer opted to view the system as a whole, focusing on the overarching patterns of inherent structures.

This novel approach sheds light on the complex challenges in the field. Embracing this concept, they introduced a global order parameter,



termed the inherent structure minimal displacement (IS D_{min}), to measure the variability of configurations using a pattern-matching methodology.

By conducting atomic simulations on seven model glass-forming liquids, they were able to unify the impacts of temperature, pressure, and perturbation time on relaxation dissipation through a scaling law linking the mechanical damping factor to IS D_{min} . They elucidated that this scaling law is a reflection of the curvature of the local potential energy landscape.

Consequently, they successfully identified a universal foundation for glassy relaxation, proposing that the variability of configurations, as quantified by IS D_{min} uniquely determines the <u>relaxation</u> damping.

This work not only presents an innovative approach to studying disordered systems but also serves as an inspiration, showcasing the potential of advanced pattern-matching techniques as potent tools for analyzing complex systems.

More information: Hai-Bin Yu et al, Universal origin of glassy relaxation as recognized by configuration pattern matching, *National Science Review* (2024). DOI: 10.1093/nsr/nwae091

Provided by Science China Press

Citation: Toward a unified theory for dynamics of glassy materials (2024, April 22) retrieved 23 June 2024 from <u>https://phys.org/news/2024-04-theory-dynamics-glassy-materials.html</u>



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