

Research at BU — Computer Simulations of SiO₂

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Department of Physics & Astronomy

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(Summer) Research Students:

Laura Owens	Robin Bjorkquist
Sean G. McMahon	Landon M. Chambers
Christopher H. Gorman	Jakob A. Roman
& benefits in classroom	

Collaborations Beyond Bucknell:

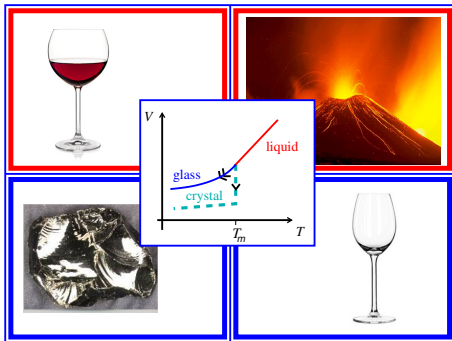
Annette Zippelius	Jürgen Horbach
Horacio E. Castillo	Karina Avila

Computers:

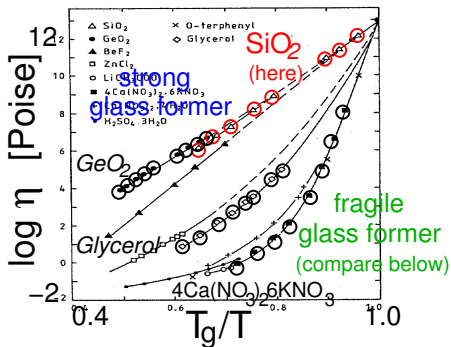
J. Dreese, M. Harvey, B. Temelso, G. P. Shrivastav

Department of Physics & Astronomy

Computer Simulations of SiO_2 Glass



Viscosity vs $1/\text{Temperature}$:



[C.A. Angell et al. 1976]

Computer Simulations of SiO₂ Glass

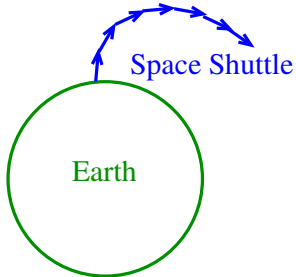
Hidden Figures:

Katherine G. Johnson

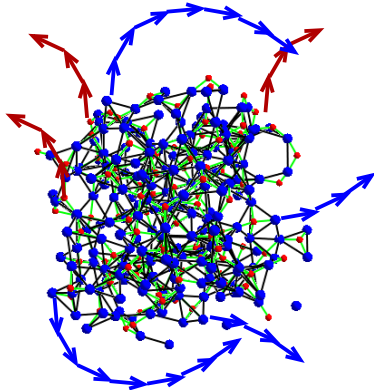
Euler's Method

Space Shuttle

Earth

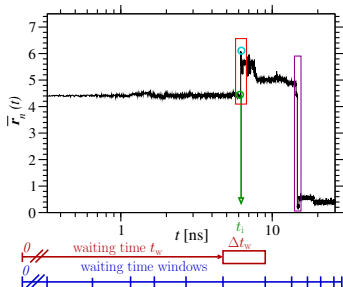


Molecular Dynamics Simulation:
SiO₂ glass:



Results: Computer Simulations of SiO₂ Glass

Single Particle Jumps:

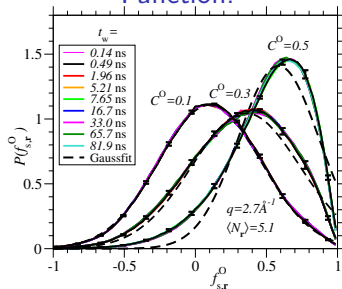


- ▶ t_w -dep.: no. jump. part., not size, not time in between jumps

Phys.Rev.Lett. 110, 017801 (2013)

Similar results for fragile & strong glass formers!

Local Incoh. Intern. Scattering Function:



- ▶ scaling collapse

JCP Editor's Choice: J.Chem.Phys. 144, 234510 (2016)