

# Study of atomic motion in rubidium borate glasses

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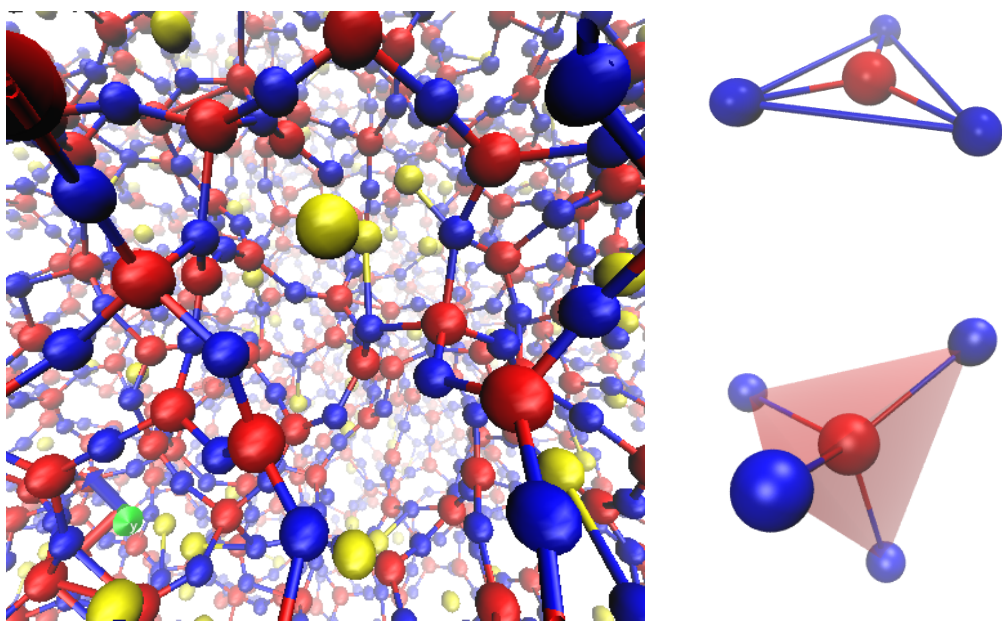
## Introduction

Atomic Scale X-ray Photon Correlation Spectroscopy (aXPCS) uses coherent X-rays to probe the dynamics of materials on an atomic scale. It was applied to study atomic diffusion in crystals [1], but its application was

recently extended to glasses as well [2]. Results of dynamics and Small-Angle X-ray Scattering (SAXS) studies on high- and low-alkali content rubidium borate glasses are presented here.

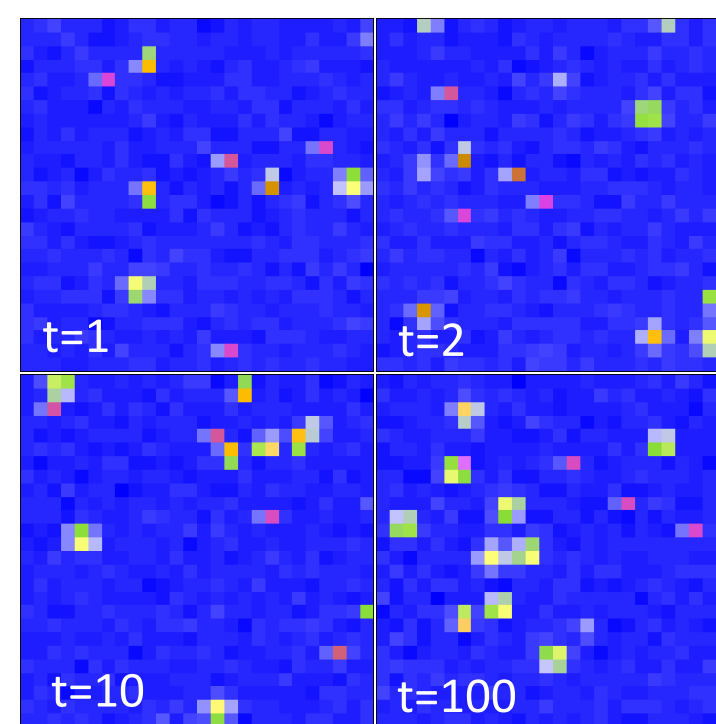
## Theory

Real space structure: changes in time results in different scattering patterns



Space and time dependence described by the Van Hove Pair Correlation Function

Sequence of scattering patterns  
→ Time correlation

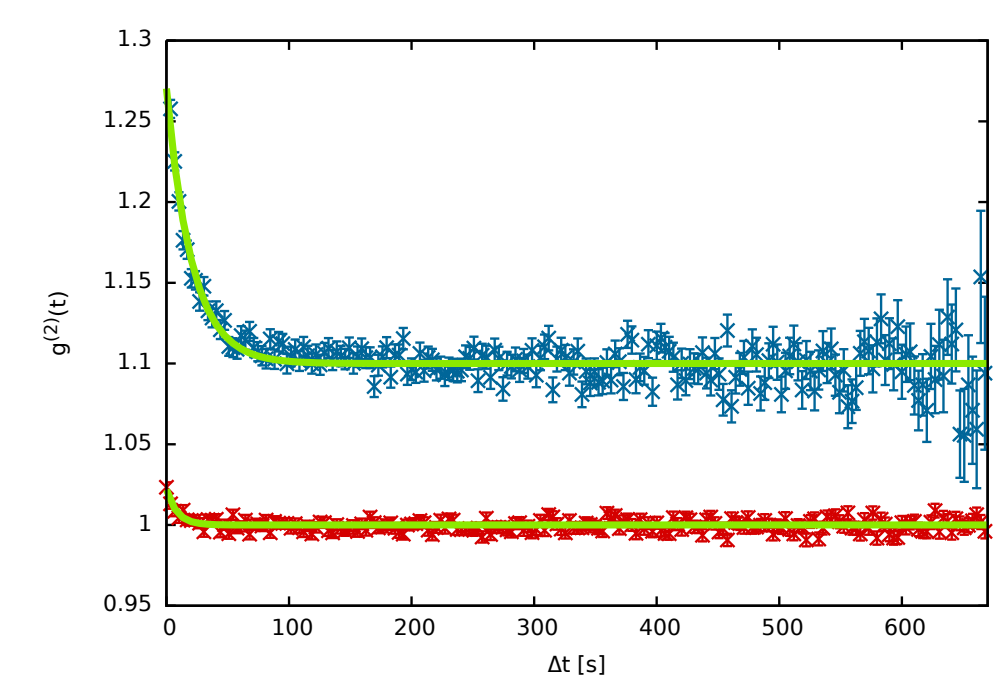


Intensity Auto Correlation Function (ACF)

$$g^{(2)}(\vec{q}, \Delta t) := \frac{\langle I(\vec{q}, t) I(\vec{q}, t + \Delta t) \rangle}{\langle I(\vec{q}, t) \rangle^2}$$

Functional form of the ACF:

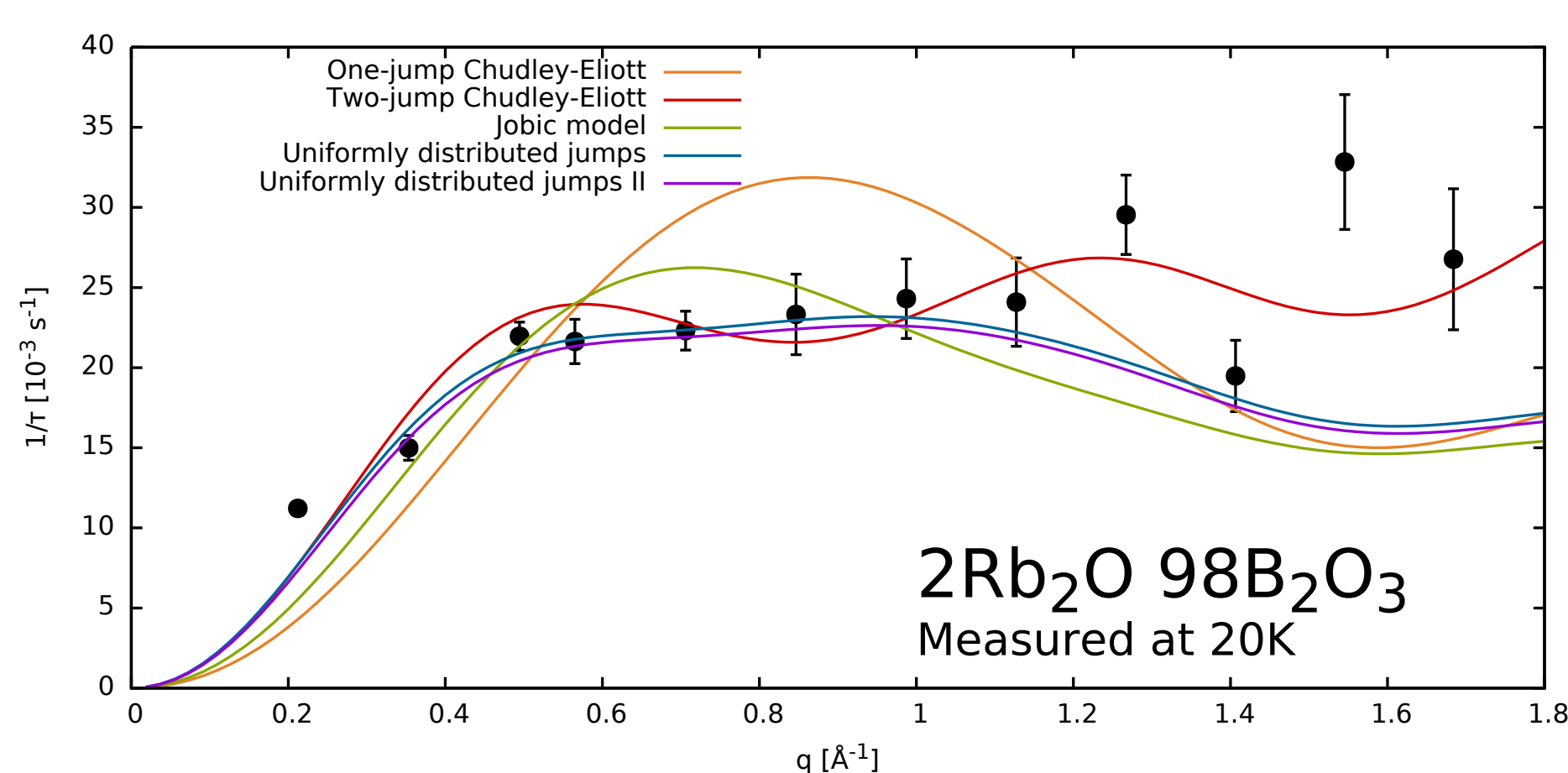
$$g^{(2)}(q) = 1 + \beta e^{-\left(\frac{2\Delta t}{\tau}\right)^\alpha}$$



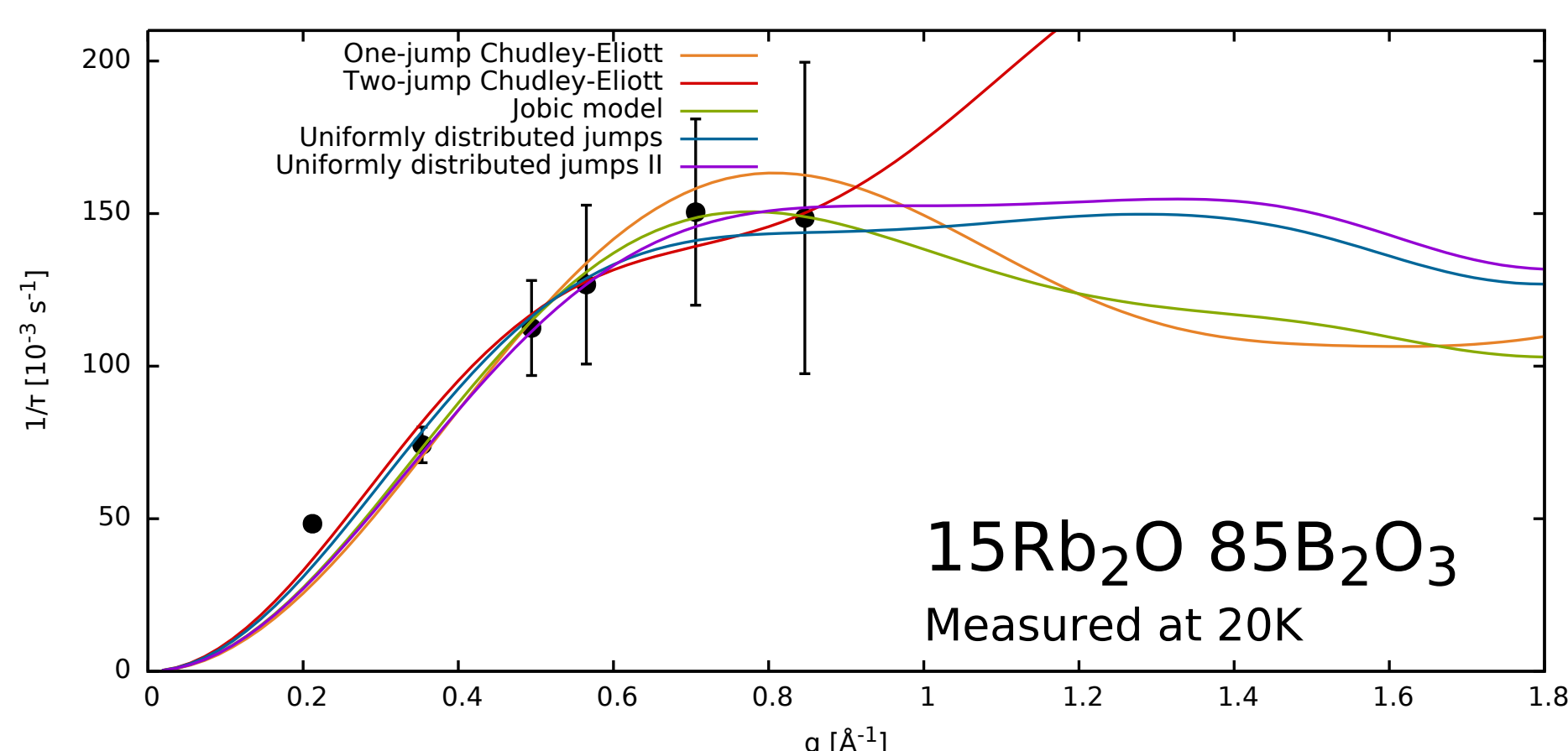
Normalization by the total structure factor

$$\tau_{coh}(\vec{q})^{-1} = \frac{\tau_{inc}(\vec{q})^{-1}}{S_{total}(q)}$$

## Results: fit to jump models

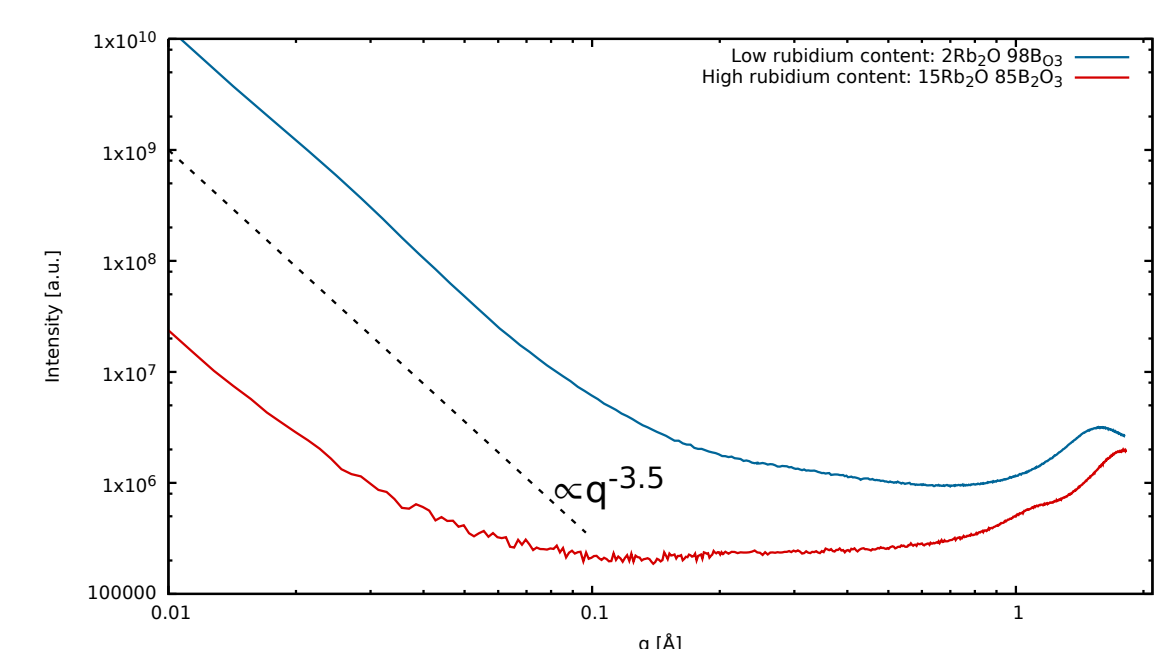


Best fit:  
Two-jump  
distances  
Chudley-Elliott  
Model with  
jump lengths  
 $d_1 = 1.0 \text{ \AA}$  and  
 $d_2 = 8.5 \text{ \AA}$ , with  
3/4 of  
short-distance  
jumps



**top:** low alkali  
content glass,  
**bottom:** High  
alkali content  
glass. Several  
models can be  
fitted at small  
q-range. Fits  
are equally  
good in the  
hydrodynamic  
limit.

## SAXS



SAXS intensities as a function of momentum transfer

## Conclusion

- Diffusion in low rubidium content borate glasses best described by two-jump model
- Large difference in diffusivity between low and high alkali glass
- Diffusion at 20K still too fast to be measured in high alkali glass
- Inhomogeneities are much larger in high alkali sample