

C153 Dynamics in Ethylammonium Nitrate: The Effects of Dilution with Methanol

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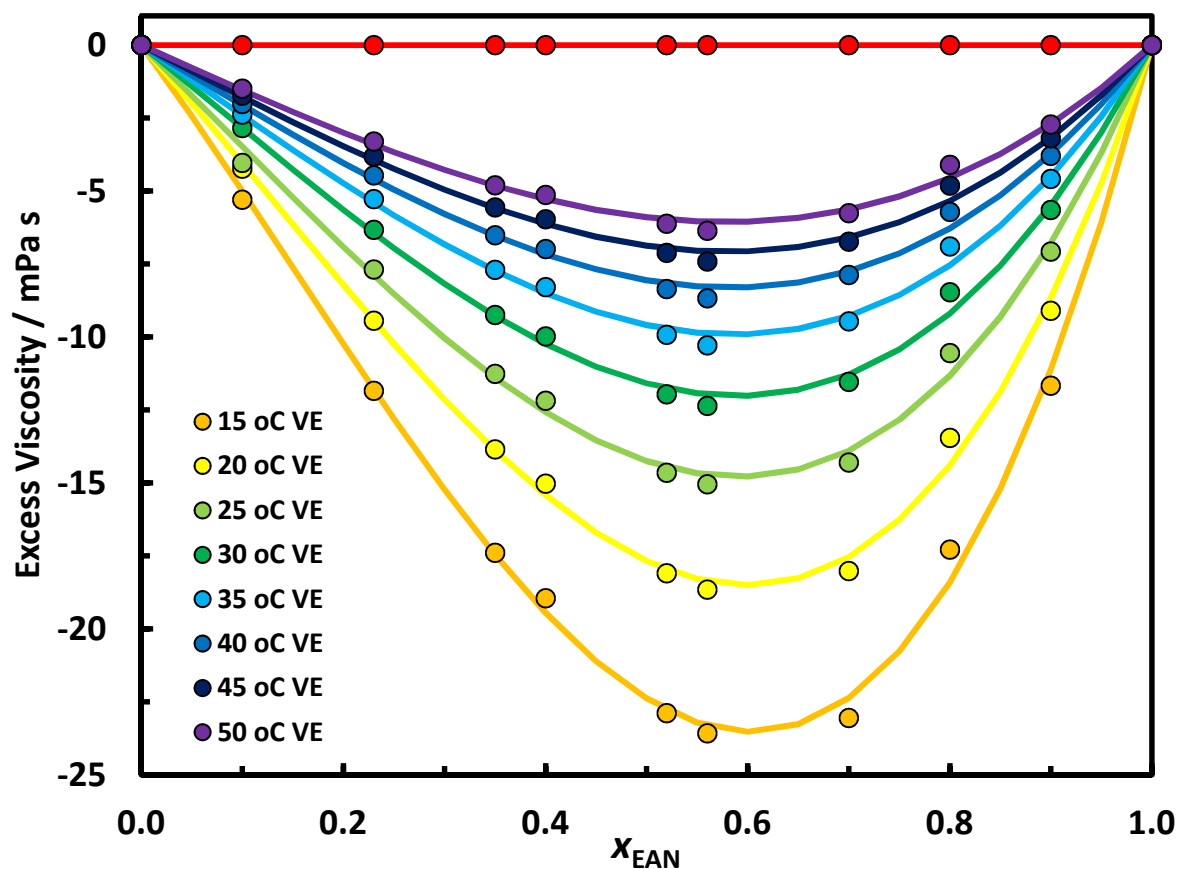
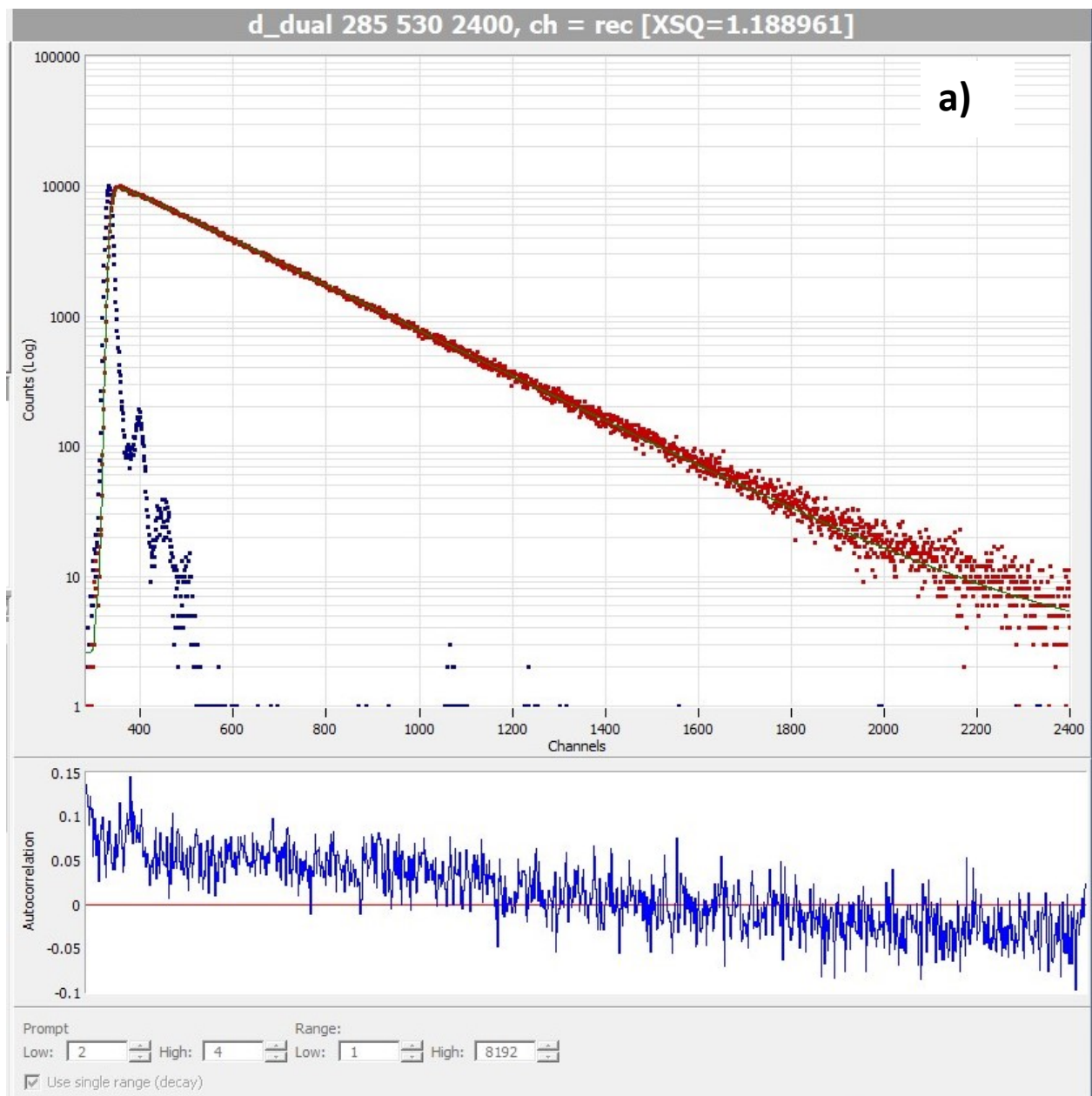


Figure S1. Excess viscosities (symbols) for EAN/MeOH as a function of EAN mole fraction for temperatures from 278 – 323 K. Lines are fits to a two component Redlich-Kister polynomial of functional form:¹.

$$\eta^E = x_{EAN} (1 - x_{EAN}) \sum_{i=1}^N A_i [2x_{EAN} - 1]^{i-1}$$



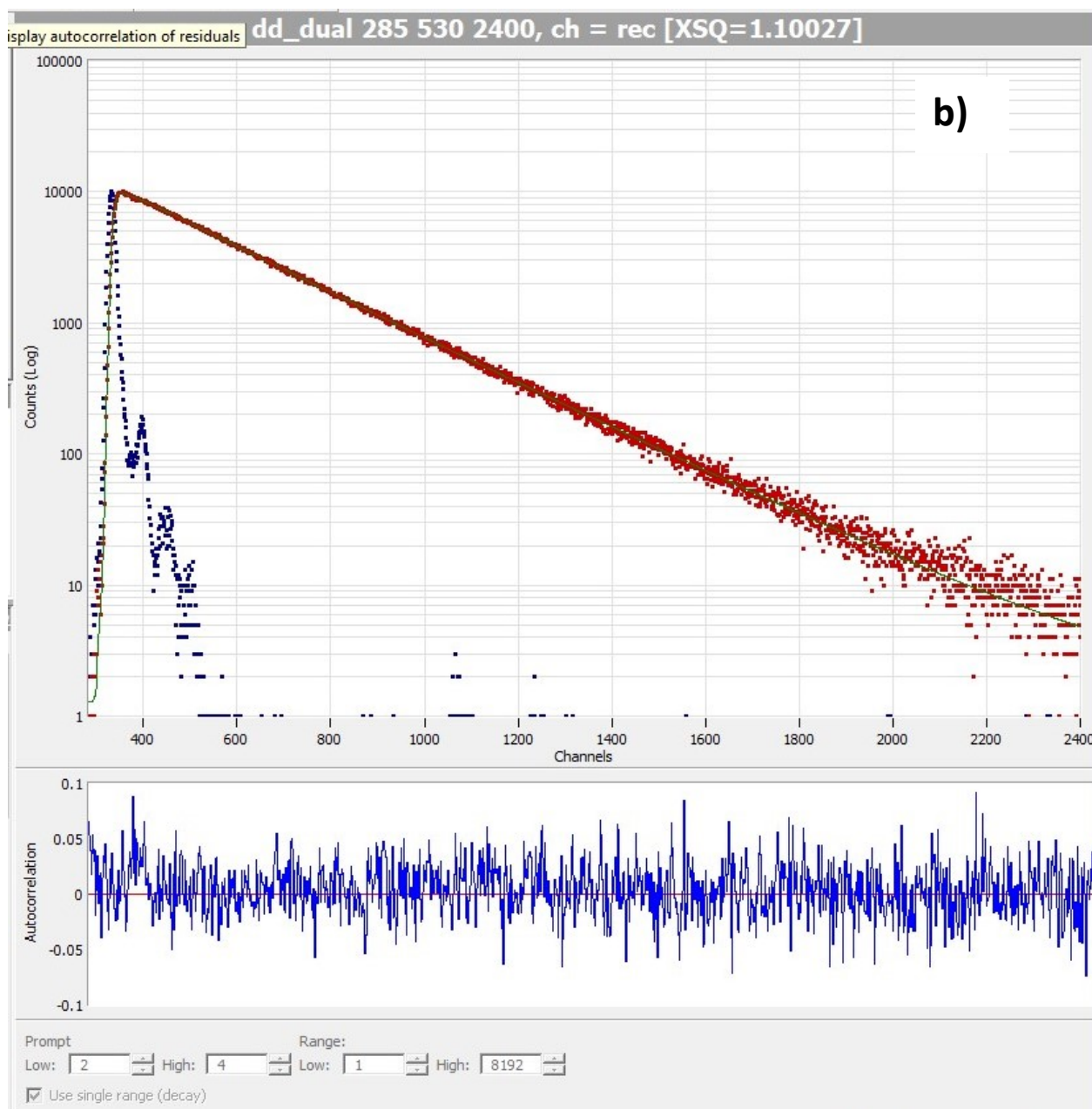


Figure S2. a) Autocorrelation of the residuals for $x_{IL} = 0.2$ EAN/MeOH at 295 K fit to a single exponential model. Note the $\chi^2 = 1.18$, which by consideration is an “acceptable fit”. Although the residuals appear convincing that this model is acceptable, the autocorrelation shows a very clear bias across the entire fitting range. b) Autocorrelation of the residuals for $x_{IL} = 0.2$ EAN/MeOH at 295 K fit to a double exponential model. Note the small improvement in $\chi^2 = 1.10$ but much more significant is the randomness of the autocorrelation function. This argues for inclusion of a second time constant to properly describe the intensity decay.

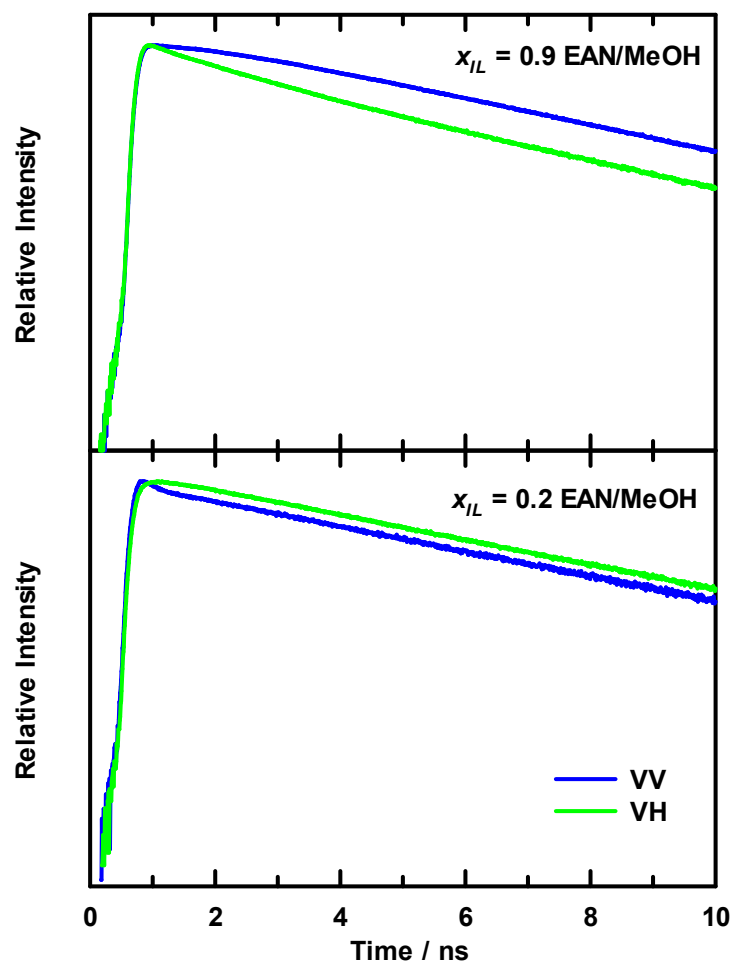


Figure S3. Parallel and perpendicular intensity decays for EAN/MeOH at 295 K. Lower panel shows polarized intensity decays for $x_{IL} = 0.2$ and upper panel is $x_{IL} = 0.9$. The traces intensities were normalized to 1. Note that more time is required for the decays to tail-match, which is the criterion to compute the anisotropy decay.

Reference

1. McAtee, Z. P.; Heitz, M. P. Density, Viscosity and Excess Properties in the Trihexyltetradecylphosphonium Chloride Ionic Liquid/Methanol Cosolvent System. *J. Chem. Thermodyn.*, **2016**, *93*, 34-44.