

Multicomponent diffusion in natural silicate melts: Toward a universal eigenvector matrix

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Multicomponent diffusion in natural silicate melts is fundamental to understanding igneous processes such as magma mixing. In N -component silicate melts, multicomponent diffusion is characterized by an $N-1$ square matrix, termed the diffusivity matrix $[D]$. Eigenvectors of $[D]$ define $N-1$ eigen-components that diffuse independently of each other. Diffusion eigenvectors in basalt appear to be roughly temperature independent [1]. In this study, we hypothesize that diffusion eigenvectors in natural silicate melts are also roughly compositionally independent. We use a single eigenvector matrix, $[P]$, to simultaneously fit concentration profiles from 27 diffusion couple experiments at three temperatures from [1-2], and one additional experiment conducted in this work. The goodness of our fitting further supports the temperature independence of $[P]$.

Oxide concentrations were converted into eigen-component "concentrations", Z_i , and plotted against distance for both the 28 diffusion experiments in basalt and those from literature in rhyolitic to basaltic melts. If $[P]$ is compositionally invariant, all Z_i profiles should be monotonic. Approximately 96% of >1000 plots are monotonic, showing that we are getting close to a universal eigenvector matrix. Notably, all Z_i profiles during quartz dissolution in basalt/rhyolite and olivine dissolution in basalt/andesite are monotonic. About 2% (23 out of 1190) show clear uphill diffusion, including one with Z_3 in our diffusion-couple experiment, and 22 with Z_1 , Z_2 , Z_4 and Z_5 in diopside/rutile dissolution in andesite, diopside/plagioclase dissolution in basalt, and diffusion couple experiments in haplo-basalt, implying the need to further constrain $[P]$. The other 2% show complexities but it is less clear whether they reflect uphill diffusion.

We propose to study multicomponent diffusion in eigen-component space through Z_i vs x plots. This approach combines the full rigor of multicomponent treatment with the simplicity of effective binary diffusion treatment. For example, a few well-designed experiments will suffice to extract all eigenvalues at desired melt composition. Moreover, this approach can treat the compositional dependence of eigenvalues. Lastly, its applicability is self-testable: If there are non-monotonic eigen-component profiles, for example, the approach should not be used.

2. Guo and Zhang (2018). *GCA*, 228, 190-204.

1. Guo and Zhang (2020). *CG*, 549, 119700.