

Single-crystal thermometric calibration of Fe-Mg order-disorder in pigeonites

DARIA PASQUAL,¹ GIANMARIO MOLIN,^{1,*} AND MARIO TRIBAUDINO²

¹Dipartimento di Mineralogia e Petrologia, Università di Padova, I-35127 Padova, Italy

²Dipartimento di Scienze Mineralogiche e Petrologiche, Università di Torino, I-10125 Torino, Italy

ABSTRACT

The single-crystal X-ray technique was used to calibrate a new intracrystalline geothermometer based on equilibrium Mg-Fe* fractionation ($\text{Fe}^* = \text{Fe}^{2+} + \text{Mn}^{2+}$) between M1 and M2 sites of natural $P2_1/c$ pigeonite. Suitable crystals free of exsolution textures and sharp diffraction maxima were selected by careful TEM and XRD investigations from a large number of samples. Two single crystals, PCA82506-3 ($\text{Wo}_6\text{En}_{76}\text{Fs}_{18}$) from the Pecora Escarpment 82506 Antarctic ureilite, and BTS308-2 ($\text{Wo}_{10}\text{En}_{47}\text{Fs}_{43}$) from the BTS308 Paranà rhyodacite, were annealed at temperatures ranging from 600 to 1000 °C. The TEM investigation, carried out on fragments of selected single crystals both before and after thermal treatment, shows heating-induced texture modifications preliminary to spinodal decomposition in both crystals and a size increase in antiphase domains in BTS308-2. The two geothermometric equations calculated by linear regression of $\ln K_B^*$ vs. $1/T$ are:

$$\ln K_B^* = -3291(\pm 269)/T(\text{K}) + 0.971(\pm 0.253); (r^2 = 0.974)$$

$$\ln K_B^* = -2816(\pm 83)/T(\text{K}) + 0.542(\pm 0.083); (r^2 = 0.995)$$

respectively for PCA82506-3 ($X_{\text{Fe}^*} = 0.20$) and BTS308-2 ($X_{\text{Fe}^*} = 0.49$) [$X_{\text{Fe}^*} = \text{Fe}^*/(\text{Fe}^* + \text{Mg})$]. These results imply negligible compositional effects on Mg-Fe* site partitioning within the range of compositions encompassed by the samples. Comparison of the intracrystalline fractionation data of pigeonites with those of *Pbca* orthopyroxenes shows a similar degree of ordering for both at a given temperature. This result suggests only a small effect of Ca on Fe*-Mg ordering in pyroxene with Ca content up to Wo_{10} .