A density model for mantle carbonaterich melts and applications to carbonatite magmatism in the upper mantle

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Volatiles cycle plays a leading role in the evolution and fate of a planet, defining habitability conditions at the surface as well as the physicochemical properties of the inner layers. Conversely, chemical and physical interactions between the inner and outer layers shape the evolution of the volatiles cycle. On Earth, the mantle constitutes a major actor of the volatiles cycle by hosting considerable amounts of carbon and hydrogen. Pressuretemperature-composition conditions prevailing within the Earth's upper mantle allow for stabilizing CO2-rich (carbonatite) melts that are important conveyors for the distribution of volatiles CO₂-H₂O. Their efficiency in transporting such volatiles resides in their mobility properties. Yet, modelling their mobility is still critically prevented by the lack of fundamental constraints on their density (and viscosity). Taking advantage of new high pressure-high temperature data recently acquired by both experimental and computational methods, we propose here a model for the density of mantle-derived carbonate-rich melts in the system MgO-CaO-Na₂O-K₂O-Li₂O-H₂O-CO₂.

Calibrated over a comprehensive database comprising >880 density data points, our model is conceived to be continuous in pressure-, temperature- and composition-space, and allows for predictions of the carbonate melt density up to 15 GPa and 2000 °C. It reasonably assumes a linear dependence on composition of the multicomponent carbonate melt volumes, and employs a simple Murnaghan equation of state in order to calculate the volumetric properties of multicomponent carbonate melts in the system MgO-CaO-Na₂O-K₂O-Li₂O-H₂O-CO₂. Combined with reasonable assumptions on volumetric properties for SiO₂ and FeO, we simulate the density and seismic P-wave velocity V_P of carbonate-rich melts experimentally equilibrated with various mantle lithologies (e.g., peridotite, eclogite, pelite), and discuss about implications.