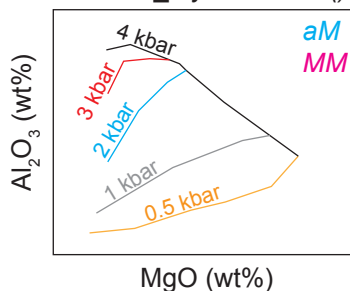


## a) Crystallisation calcs

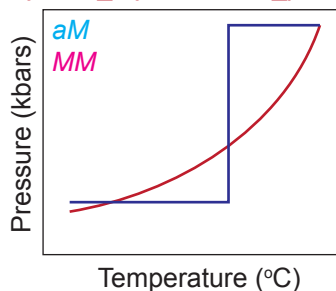
aM - compatible with alphaMELTS for Python  
MM - compatible with MAGEMin\_C

*isobaric\_crystallisation()*

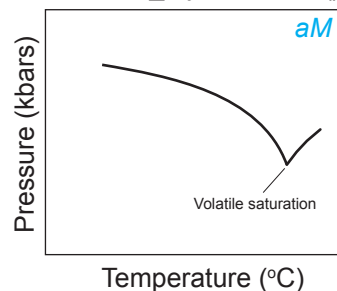


*polybaric\_crystallisation\_onestep()*

*polybaric\_crystallisation\_path()*



*isochoric\_crystallisation()*

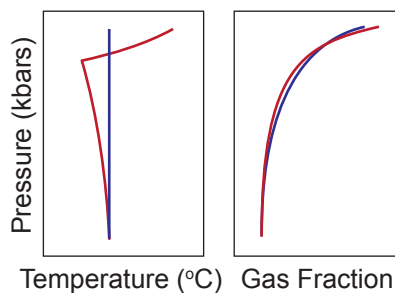


## b) Decompression calcs

aM

*isothermal\_decompression()*

*isentropic\_decompression()*

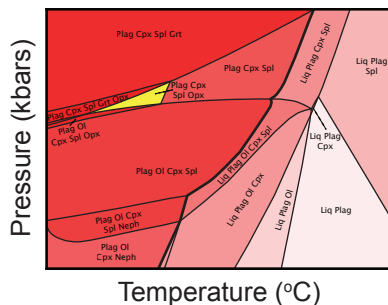


## c) Phase Diagram calcs

aM

MM

*phaseDiagram\_calc()*

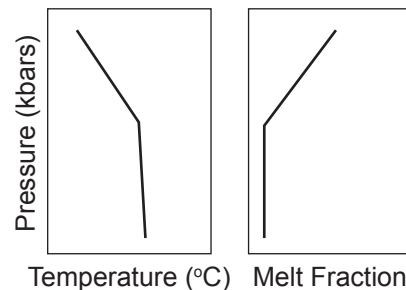


## d) Melting calcs

aM

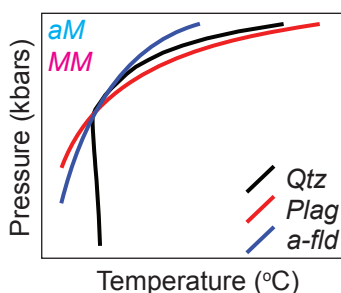
MM

*adiabatic\_melting()*

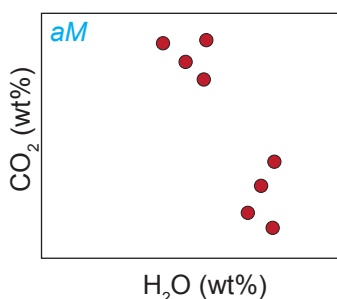


## e) Barometry calcs

*mineral\_cosaturation()*



*saturationP\_calc()*



## e) Additional core functions

*findLiq\_multi()* - load in an Excel spreadsheets of liquids and calculate their liquidus temperatures

*findCO2\_multi()* - given a particular melt composition, pressure and H2O content, what is the maximum CO2 content of the melt phase (aM only)

*equilibrate\_multi()* - takes multiple input variables and returns data for phase composition, abundance, and thermodynamic properties