The parameters are an ID of ion, ion column density (N\_ion, [10<sup>18</sup> cm<sup>-2</sup>]), an effective temperature for the Doppler broadening  $kT_{\rm eff}$  [keV], and redshift z. The ion ID is written as (atomic number)+(the number of electrons). For example, Fe xxvI is 2601, that of Fe xxv is 2602 and that of Ni xxvII is 2802. The effective temperature  $kT_{\rm eff}$  is written as the  $2kT_{\rm eff}/m_{\rm atom} = 2kT/m_{\rm atom} + v_{\rm turb}^2$ , where T is the real temperature of the gas and  $v_{\rm turb}$  is the microscopic turbulent velocity of the gas. Thus, the Doppler broadening is calculated as  $\Delta E_D/E = \sqrt{2kT_{\rm eff}/m_{\rm atom}}/c$ . If  $kT_{\rm eff}$  is much higher than the temperature you think, then the microscopic turbulence must be very large. If you use the negative value for this parameter, it returns  $\Delta v_D/c = \Delta E_D/E$  directly.

When you use it, please set the environmental variable "IONABS\_DATA\_PATH" to the directory of data\_ionabs your .zshrc, such as "export IONABS\_DATA\_PATH=/path/to/the/data\_ionabs"