DiadFit: An Open-SourcePython3 Tool for Peak fitting of Raman Data from silicate melts and CO_2 fluids

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We welcome any feedback – it might not work straight out the box for every Raman set up on the planet, but if you send us some spectra, we will tweak it to ensure its useful for your lab!

Any questions – contact penny_wieser@berkeley.edu

See also our GitHub:

PennyWieser/DiadFit: A python package for fitting Fermi Diads, Ne lines, and other things relating to Raman analyses of CO2 fluids (github.com)

Read the docs page:

https://diadfit.readthedocs.io/