

Isobar_Comparison_Plotter

September 28, 2020

```
[2]: import sys
      sys.path.append('../..../')

      import VESical as v
      import pandas as pd
```

```
[6]: alkbasalt = {'SiO2': 49.0,
                  'TiO2': 1.27,
                  'Al2O3': 19.7,
                  'Fe2O3': 3.74,
                  'FeO': 5.33,
                  'MnO': 0.17,
                  'MgO': 4.82,
                  'CaO': 8.85,
                  'Na2O': 4.23,
                  'K2O': 1.00,
                  'P2O5': 0.37,
                  'H2O': 4.51,
                  'CO2': 0.25}

      rhyolite = {'SiO2': 77.19,
                  'TiO2': 0.06,
                  'Al2O3': 12.80,
                  'FeO': 0.94,
                  'MgO': 0.03,
                  'CaO': 0.53,
                  'Na2O': 3.98,
                  'K2O': 4.65,
                  'CO2': 0.05,
                  'H2O': 0.26}

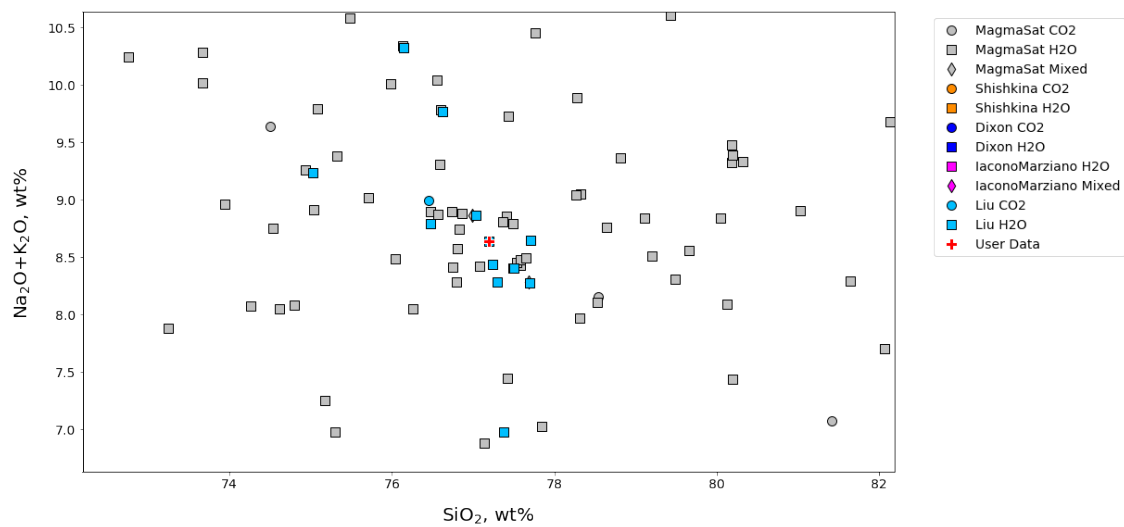
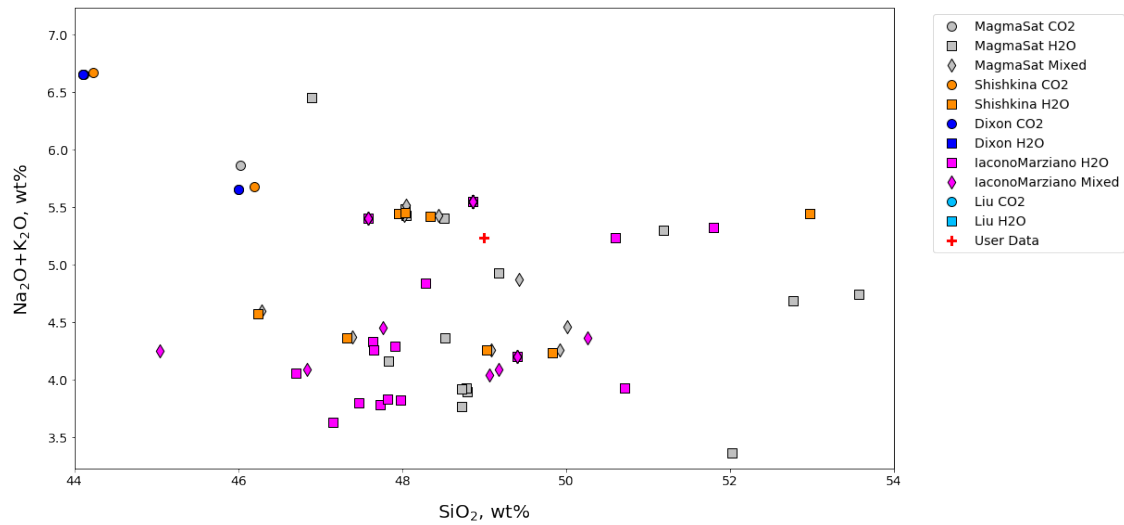
      sample_table = pd.DataFrame([alkbasalt, rhyolite], index=["Alkali Basalt",
      ↪ "Rhyolite"])
      sample_table
```

```
[6]:           SiO2  TiO2  Al2O3  Fe2O3  FeO  MnO  MgO  CaO  Na2O  K2O  \
Alkali Basalt  49.00  1.27   19.7   3.74  5.33  0.17  4.82  8.85  4.23  1.00
```

Rhyolite	77.19	0.06	12.8	NaN	0.94	NaN	0.03	0.53	3.98	4.65
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	P205	H2O	CO2
Alkali Basalt	0.37	4.51	0.25
Rhyolite	NaN	0.26	0.05

```
[15]: #check calibration
v.calib_plot(user_data=alkbasalt, model='mixed', zoom='user_data')
v.calib_plot(user_data=rhyolite, model='mixed', zoom='user_data')
```



```
[5]: alkbasalt_isobars, alkbasalt_isopleths = v.
      ↪calculate_isobars_and_isopleths(sample=alkbasalt, temperature=1200,
      ↪pressure_list=[500, 1000, 2000], isopleth_list=[0.5], print_status=True).
      ↪result

      rhyolite_isobars, rhyolite_isopleths = v.
      ↪calculate_isobars_and_isopleths(sample=rhyolite, temperature=800,
      ↪pressure_list=[500, 1000, 2000], isopleth_list=[0.5]).result
```

```
Calculating isobar at 500 bars
Calculating isopleth at 0
Calculating isopleth at 0.5
Calculating isopleth at 1
Calculating isobar at 1000 bars
Calculating isopleth at 0
Calculating isopleth at 0.5
Calculating isopleth at 1
Calculating isobar at 2000 bars
Calculating isopleth at 0
Calculating isopleth at 0.5
Calculating isopleth at 1
Done!

Calculating isobar at 500 bars
Calculating isopleth at 0
Calculating isopleth at 0.5
Calculating isopleth at 1
Calculating isobar at 1000 bars
Calculating isopleth at 0
Calculating isopleth at 0.5
Calculating isopleth at 1
Calculating isobar at 2000 bars
Calculating isopleth at 0
Calculating isopleth at 0.5
Calculating isopleth at 1
Done!
```

```
[6]: Iac_alkbasalt_isobars, Iac_alkbasalt_isopleths = v.
      ↪calculate_isobars_and_isopleths(sample=alkbasalt, temperature=1200,
      ↪pressure_list=[500, 1000, 2000], isopleth_list=[0.5],
      ↪model="IaconoMarziano").result
      Dixon_alkbasalt_isobars, Dixon_alkbasalt_isopleths = v.
      ↪calculate_isobars_and_isopleths(sample=alkbasalt, temperature=1200,
      ↪pressure_list=[500, 1000, 2000], isopleth_list=[0.5], model="Dixon").result
      Shish_alkbasalt_isobars, Shish_alkbasalt_isopleths = v.
      ↪calculate_isobars_and_isopleths(sample=alkbasalt, temperature=1200,
      ↪pressure_list=[500, 1000, 2000], isopleth_list=[0.5], model="Shishkina").
      ↪result
```

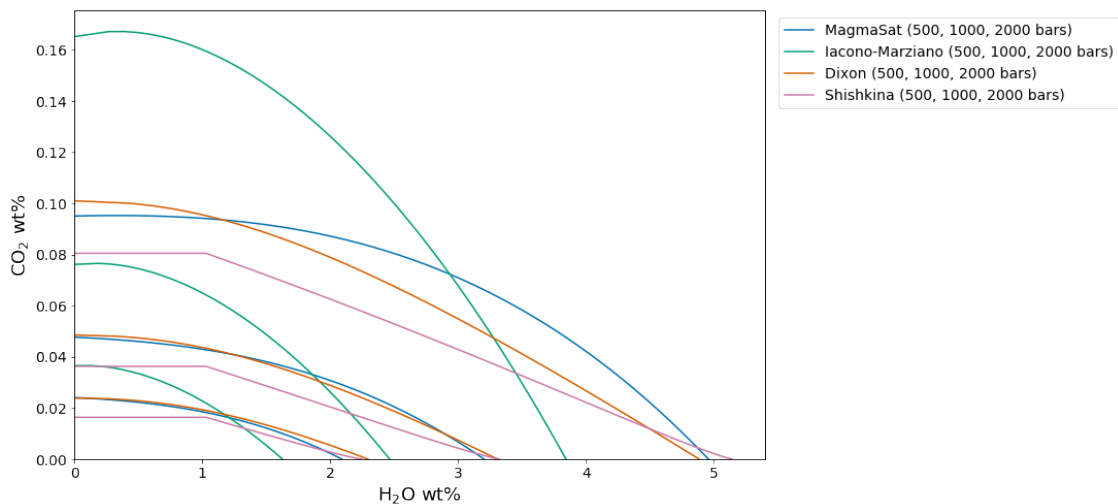
```
Liu_rhyolite_isobars, Liu_rhyolite_isopleths = v.
    ↪calculate_isobars_and_isopleths(sample=rhyolite, temperature=800,
    ↪pressure_list=[500, 1000, 2000], isopleth_list=[0.5], model="Liu").result
```

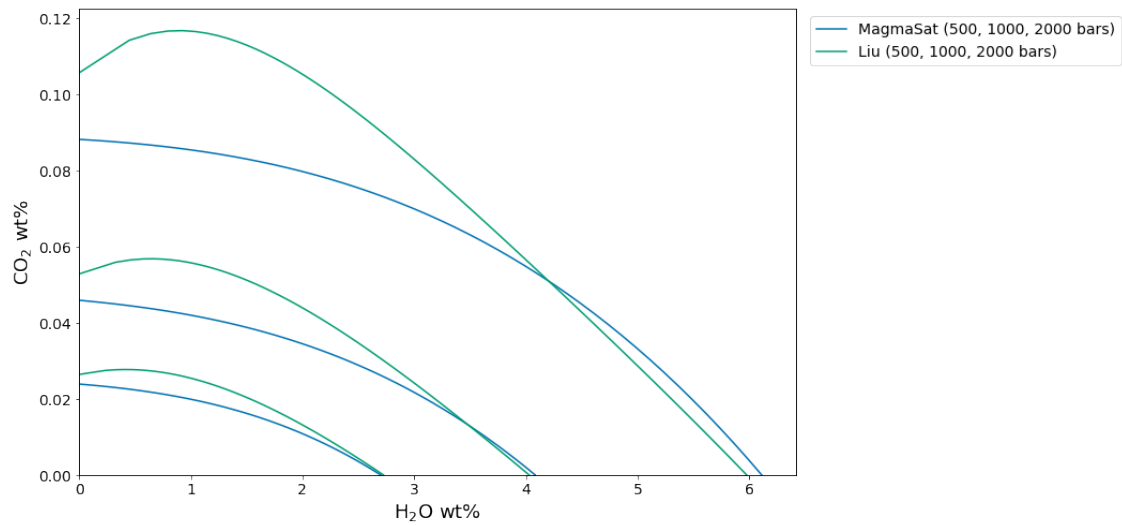
```
../../../../VESIcal.py:1719: RuntimeWarning: pressure exceeds 1000 bar, which
Iacono-Marziano et al. (2012) suggest as an upper calibration limit of the Dixon
(1997, Pi-SiO2 simpl.) Model,
```

```
    w.warn(self.calib_check,RuntimeWarning)
../../../../VESIcal.py:1719: RuntimeWarning: These calibration limits were
selected based on the minimum and maximum values of these oxides (+5%) in the
calibration dataset. As the Liu et al. model incorporates no term for
compositional dependence, users must take extreme care when extrapolating this
model to compositions which differ significantly from the haplogranites and
rhyolites in the calibration dataset. These warnings are simply a guide; we
suggest that users carefully compare their major element data to the calibration
dataset to check for suitability
```

```
    w.warn(self.calib_check,RuntimeWarning)
```

```
[9]: v.plot(isobars=[alkbasalt_isobars, Iac_alkbasalt_isobars,
    ↪Dixon_alkbasalt_isobars, Shish_alkbasalt_isobars],
    ↪isobar_labels=["MagmaSat", "Iacono-Marziano", "Dixon", "Shishkina"])
v.plot(isobars=[rhyolite_isobars, Liu_rhyolite_isobars],
    ↪isobar_labels=["MagmaSat", "Liu"])
```





```
[ ]: v.plot(isobars=Shish_alkbasalt_isobars)
```

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[ ]:
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