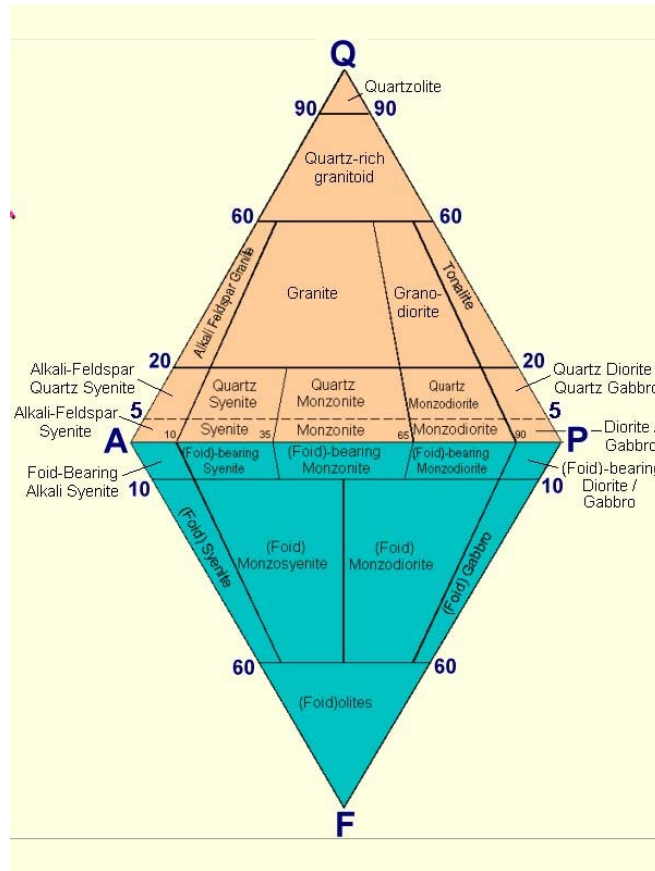


Normative Classification (Phaneritic):



5. Other buttons:

New clears all elemental information and results for a new analysis.

Open retrieves previously saved compositions and results.

Print prints a copy of the the Magma window to your default pringer

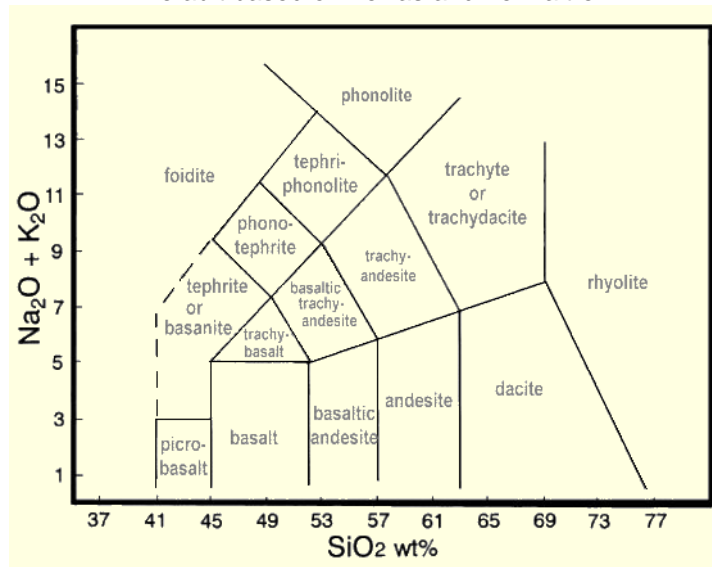
TAS Plot displays an total alkali-silica (TAS) diagram with your analyses plotted on the classification diagrams of *Le Bas & Streckheisen (1991)* and *Le Maitre et al (1989)*, or *Cox et al. (1979)*. The plot window allows clearing of specified or all analysis and printing via a print window.

Save saves the displayed results to a ascii file for later printing.

Exit ends the application.

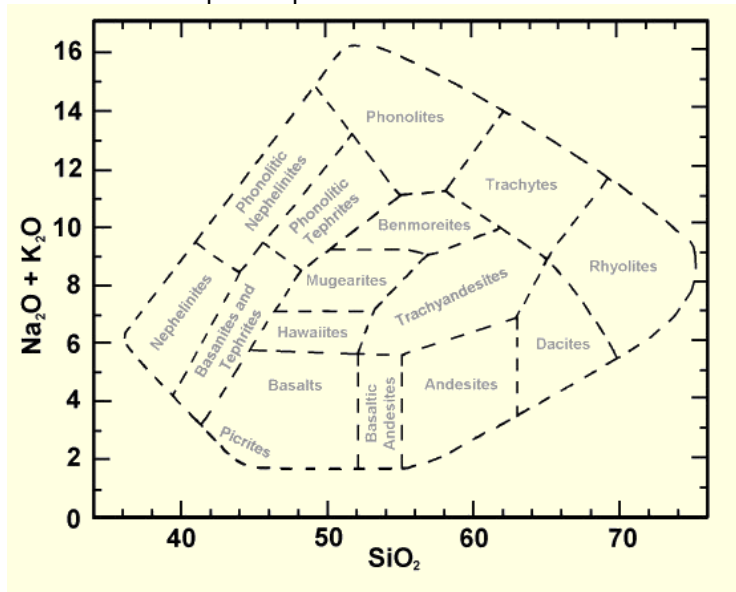
Total Alkali-Silica Classification:

Default based on LeBas and Le Maitre



Note: tephrite when OL < 10%, trachyte when Q < 20%

Optional plot based on Cox et al.



NOTES:

1. The Fe_3+/Fe_{Tot} ratio depends on the oxidation state of the magma, often a function of silica content. The present version of magma will suggest a ratio to use if you have not entered a value for either Fe_2O_3 or FeO . This value ranges between 0.3 for basalts to 0.60 for rhyolites, with a value near 0.50 reasonable for all magmas.
2. The CIPW norm calculation is performed according to *Myron G. Best; Igneous and Metamorphic Petrology, Freeman, 1982, Appendix E, pp. 616-619.*
3. The Felsic mineral designation are values calculated to use for the IUGG classification tetrahedron. The distribution of Ab between K-feldspar and plagioclase for IUGS classification is somewhat complex, depending upon the overall composition of the magma. In order to make the classification as robust as possible for standard compositions, a 3rd order polynomial equation is used to determine the distribution coefficient, XPLAG, as follows.

$$\text{FELD} = \text{AN} - ((\text{ORT} + \text{LC}) / (\text{AB} + \text{NE})) + (\text{NE} * (\text{AN} / (\text{AN} + \text{AB})))$$

$$\text{XPLAG} = 0.02108 + 0.01716 * (\text{FELD}) + 0.00234 * (\text{FELD}^2) - 0.00006 * (\text{FELD}^3)$$

This method seems to provide the best overall IUGS classification for volcanic rocks that I have found. If you have a suggestion to improve the method, please let me know.

- Densities (*Bottinga and Weil, 1972, Amer. J. Sci., 269: 169-182*) take into account normative crystal content the amount of water in the sample. Bulk density is calculated with respect to vesicles that may be present if water is excess of saturation at the calculated pressure. Particle density does not include vesicle volumes. The texture designation is dependent upon bulk density; if vesicle volumes are in excess of 0.75, the rock is arbitrarily assigned to be "Tephra."

- The water solubility model used is a simple Henry's Law model where

$$n = sP^\beta$$

- for basalts: $s = 0.0215$ and the exponent $\beta = 0.5$
- for rhyolites: $s = 0.130$ and the exponent $\beta = 0.7$
- for intermediate compositions a simple linear mixing model between these endpoint compositions is employed

- The liquidus temperature calculation follows *Sisson and Grove (1993) Contrib. Mineral. Petrol., 113, 143 - 166*:

$$\text{Tliq} = 969 - 33.1 * \text{Hx} + 742.7 * \text{AlNum} - 138 * \text{NaKNum} + 125.3 * \text{MgNum}$$

where $\text{AlNum} = \text{ALx} / (\text{Six} + \text{ALx})$, $\text{NaKNum} = (\text{Kx} + \text{NAx}) / (\text{Kx} + \text{NAx} + \text{CAx})$, $\text{MgNum} = \text{MGx} / (\text{MGx} + \text{F2x})$, and variables denoted with small x are oxide weight fractions (H is water).

- The viscosity values are given in log units of Pa s (10 Poise). The methods for calculation are given in

Bottinga, Y. A., and Weil, D. F., 1972. The viscosity of magmatic silicate liquids: a model for calculation. Amer. J. Sci., 272: 438-473.

Shaw, H. R., 1972. Viscosities of magmatic silicate liquids: an empirical method of prediction. Amer. Jour. Sci. 272: 870-893.

Effective viscosities are calculated to estimate the effects of crystal and vesicles on viscosity. This calculation involves the Einstein-Roscoe equation from *McBirney and Murase, Ann. Rev. Earth Planet. Sci. 12: 337-357, 1984.*

$$\eta' = \eta (1 - R \phi)^{-2.5} \quad (\text{for the Bottinga-Weil estimation})$$

and

$$\ln \eta' = \ln \eta + \alpha D / [(\phi'/\phi)^{1/3} - 1] - 0.15 \quad (\text{Sherman's modification for the Shaw viscosity}).$$

where η' is the effective, viscosity, η is the liquid viscosity, ϕ is volume fraction of crystals (and vesicles), R is volumetric ratio of solids at maximum packing that is similar to ϕ' , the concentration of crystals (and vesicles) at maximum packing, and α is a constant that varies with mean diameter, D . Following *McBirney and Murase (1984)*, this program assumes for a range in compositions the best fits to data are obtained using a value for $\alpha = 0.011$ and $R(\phi')$ at 1.0. Certainly there is much room for debate about these constants, so one should be careful to recognize the limitations of these estimations.

- Option check boxes activate convenient features:

Add Water: Checking this box forces Magma to add water to make the total = 100 wt%

Auto Pressure: If a composition has a water content in excess of saturation (calculated in 5 above), the pressure will be automatically increased to minimize the excess water

Adjust FE Ratio: If Fe is present in the magma composition but is expressed as a total of FeO and

Fe₂O₃ so that one or the other is entered as zero, Magma will prompt the user with a suitable ratio to be applied.

9. Results are models and must be viewed as such!
(There are more recent models for solubility and viscosity functions that may be more appropriate for your work.)

Help: Please send email to KWare (KWare@lanl.gov) for further support.



Software by KWare: <http://www-geo.lanl.gov/Wohletz/KWare.htm>

Magma (Igneous rock norms, classification, and physical properties)

This Windows Help file was written by Ken Wohletz
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