

Package ‘hydrogeo’

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Title Groundwater Data Presentation and Interpretation

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Description Contains one function for drawing Piper diagrams (also called Piper-Hill diagrams) of water analyses for major ions.

Depends R (>= 2.6.0)

Imports methods

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hydrogeo

Groundwater data presentation and interpretation.

Description

Contains one function, for drawing Piper (or Piper-Hill) diagrams from water analyses for major ions, and a dataset from Zaporozec

Details

Package: hydrogeo
Type: Package
Version: 0.4-1
Date: 2016-07-09
License: BSD
LazyLoad: yes

Author(s)

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See Also

[piper](#) and [toPercent zaporozec](#)

Examples

```
library(hydrogeo)
data(zaporozec)
zaporozec$CO3 <- rep(0,9) # toPercent expects CO3
zaporozec$Na <- rep(0,9) # toPercent expects Na
z <- toPercent(zaporozec)
pz <- piper(z)
plot(pz,cex=1.5)
```

piper

Create a new piper object

Description

Create a new piper object

Usage

`piper(d, ...)`

Arguments

`d` list passed to class `piper`, [piper](#)
`...` additional arguments, as for [piper](#)

See Also

[piper-class](#) and [toPercent](#)

`piper-class` *Class* `piper`

Description

Objects of this class are plotable as Piper-Hill diagrams. A dataframe of major ions as percentages can be used to initialise a `piper` object.

Usage

```
## S4 method for signature 'piper'
initialize(.Object, l, ..., call = NULL, group = NULL,
          wt.col = NULL, pt.col = NULL)

## S4 method for signature 'piperplot'
labelAxes(x, cex.axis = 0.35, side = -1, ...)

## S4 method for signature 'piper'
plot(x, type = "p", cex = 0.75, ...)

## S4 method for signature 'piper'
show(object)
```

Arguments

`.Object` object of class `piper`
`l` list of data, see 'Examples' below
`...` additional arguments, as for [piper](#)
`call` the call that asked for the new `piper` object
`group` (untested) Object of class `vector`, a second factor to group by, to allow plotting on different plots
`wt.col` (untested) Water type colours
`pt.col` Object of class `vector` of colours for points

<code>x</code>	an object of class <code>piperplot</code>
<code>cex.axis</code>	magnification to be used for axis annotation relative to the current setting of <code>'cex'</code> , see <code>help("par")</code>
<code>side</code>	integer between 1 and 10 specifying which side to label, the default is to label all
<code>type</code>	what type of plot should be drawn, only "p" for <code>*p*oints</code> is useful
<code>cex</code>	magnification to be used for symbols relative to the current setting of <code>'cex'</code> , see <code>help("par")</code>
<code>object</code>	an object of class <code>piper</code>

Methods (by generic)

- `initialize`: Initialiser
- `labelAxes`: Label the axes
- `plot`: Plot an object of class `piper`
- `show`: Show an object of class `piper`

Slots

<code>Ca</code>	Object of class <code>vector</code> — Calcium
<code>Mg</code>	Object of class <code>vector</code> — Magnesium
<code>Cl</code>	Object of class <code>vector</code> — Chloride
<code>SO4</code>	Object of class <code>vector</code> — Sulphate
<code>WaterType</code>	Object of class <code>vector</code> — factor for grouping samples
<code>anion.x</code>	<code>x</code> coordinate of the point on the anion triangle (internal)
<code>anion.y</code>	<code>y</code> coordinate of the point on the anion triangle (internal)
<code>cation.x</code>	<code>x</code> coordinate of the point on the cation triangle (internal)
<code>cation.y</code>	<code>y</code> coordinate of the point on the cation triangle (internal)
<code>diamond.x</code>	<code>x</code> coordinate of the point on the diamond (internal)
<code>diamond.y</code>	<code>y</code> coordinate of the point on the anion diamond (internal)
<code>group</code>	Object of class <code>vector</code> Another way of grouping other than <code>WaterType</code>
<code>IDs</code>	Object of class <code>vector</code> of sample identifiers
<code>pt.col</code>	Object of class <code>vector</code> of colours for points
<code>pt.pch</code>	Object of class <code>vector</code> of symbols for points
<code>call</code>	Object of class <code>character</code> — call that created it

Author(s)

Myles English <myles@rockhead.biz>

References

A. Zaporozec, "Graphical interpretation of water quality data," *Ground Water* 10, no. 2 (1972): 32–43.

Examples

```
showClass("piper")

l <- list( Ca = c(43,10,73,26,32),
          Mg = c(30,50,3,14,12),
          Cl = c(24,10,12,30,43),
          SO4 = c(24,10,12,30,43),
          WaterType = c(2,2,1,2,3),
          IDs = c("A","B","C","D","E") )

lp <- piper(l)
plot( lp, main="Piper-Hill Diagram of Water Quality" )
lp <- piper(l)
plot( lp, main="Piper-Hill Diagram of Water Quality", cex=1.4 )
```

piperPaper

Create a new piperplot object

Description

Create a new piperplot object

Usage

```
piperPaper(size = NULL, ...)
```

Arguments

size	integer related to the size of the plot area
...	additional arguments, as for piperplot

Examples

```
library(hydrogeo)
p = piperPaper(size=1)
plot(p)
```

piperplot-class *Class* piperplot

Description

Objects of this class are plottable as empty (i.e. no points) Piper-Hill diagrams

Usage

```
## S4 method for signature 'piperplot'
Axis(x = NULL)
```

Arguments

x an object of class piperplot

Methods (by generic)

- Axis: Add axes to a piperplot

Slots

size Object of class numeric — Length of the (square) plot area, defaults to 300
 call R call that created it

plot,piperplot-method *Plot the diagram area with two triangles and a diamond*

Description

Plot the diagram area with two triangles and a diamond

Usage

```
## S4 method for signature 'piperplot'
plot(x, axes = TRUE, ...)
```

Arguments

x object of class piperplot
 axes logical saying whether to draw the axes or not, defaults to TRUE
 ... further arguments to plot.default

toPercent

Major ions as a percentage of total major ions

Description

Expects certain column names

Usage

```
toPercent(d)
```

Arguments

d list or data.frame with the following columns: Ca, Mg, Na, K and Cl, SO4, CO3, HCO3

Examples

```
library(hydrogeo)
l <- list( Ca = c(43,10,73,26,32),
          Mg = c(30,50,83,14,62),
          Na = c(54,76,3,14,12),
          K = c(31,22,32,22,11),
          Cl = c(24,10,12,30,43),
          SO4 = c(24,10,12,30,43),
          CO3 = c(24,10,12,30,43),
          HCO3 = c(42,110,12,3,4),
          WaterType = c(2,2,1,2,3),
          IDs = c("A","B","C","D","E") )
d <- toPercent(l)
# check, should add up to 100%
z <- as.data.frame(d)
for(i in 1:length(z[[1]])) { print(sum(z[i,5:8])) }
for(i in 1:length(z[[1]])) { print(sum(z[i,1:4])) }
```

zaporozec

Major ions for groundwaters reported by Zaporozec

Description

This data set contains major ion analyses for three groundwaters.

Format

A data frame with 9 observations on the following 15 variables:

- location a factor with levels Tertiary, Czechoslovakia Upper Cambrian, Wisconsin Upper Cretaceous, Czechoslovakia
- K a numeric vector - potassium
- Mg a numeric vector - magnesium
- Ca a numeric vector - calcium
- Mn a numeric vector - magnesium
- Fe a numeric vector - iron
- Cl a numeric vector - chloride
- NO3 a numeric vector - nitrate
- HCO3 a numeric vector - bicarbonate
- SO4 a numeric vector - sulphate
- sigma a numeric vector - standard deviation
- TDS a numeric vector - total dissolved solids
- tempC a numeric vector - temperature
- pH a numeric vector - pH
- units a factor with levels meq/l meq_pc mg/l

Source

Zaporozec, "Graphical interpretation of water quality data," *Ground Water* 10, no. 2 (1972): pages 32–43.

Examples

```
data(zaporozec)
str(zaporozec)
```

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