

# Package ‘provenance’

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**Title** Statistical Toolbox for Sedimentary Provenance Analysis

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**Description** Bundles a number of established statistical methods to facilitate the visual interpretation of large datasets in sedimentary geology. Includes functionality for adaptive kernel density estimation, principal component analysis, correspondence analysis, multidimensional scaling, generalised procrustes analysis and individual differences scaling using a variety of dissimilarity measures. Univariate provenance proxies, such as single-grain ages or (isotopic) compositions are compared with the Kolmogorov-Smirnov, Kuiper or Sircombe-Hazelton L2 distances. Categorical provenance proxies such as chemical compositions are compared with the Aitchison and Bray-Curtis distances, and point-counting data with the chi-square distance. Also included are tools to plot compositional and point-counting data on ternary diagrams and point-counting data on radial plots, to calculate the sample size required for specified levels of statistical precision, and to assess the effects of hydraulic sorting on detrital compositions. Includes an intuitive query-based user interface for users who are not proficient in R.

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<i>ALR</i>	<i>Additive logratio transformation</i>
------------	---

---

**Description**

Calculates Aitchison’s additive logratio transformation for a dataset of class `compositional` or a compositional data matrix.

**Usage**

```
ALR(x, ...)  
  
## Default S3 method:  
ALR(x, inverse = FALSE, ...)  
  
## S3 method for class 'compositional'  
ALR(x, ...)
```

**Arguments**

- `x` an object of class `compositional` OR a matrix of numerical values
- `...` optional arguments
- `inverse` perform the inverse inverse logratio transformation?

**Value**

a matrix of ALR coordinates OR an object of class `compositional` (if `inverse=TRUE`).

**Examples**

```
# logratio plot of trace element concentrations:  
data(Namib)  
alr <- ALR(Namib$Trace)  
pairs(alr[,1:5])  
title('log(X/Pb)')
```

---

amalgamate

*Group components of a composition*


---

## Description

Adds several components of a composition together into a single component

## Usage

```
amalgamate(X, ...)

## Default S3 method:
amalgamate(X, ...)

## S3 method for class 'compositional'
amalgamate(X, ...)

## S3 method for class 'counts'
amalgamate(X, ...)

## S3 method for class 'SRDcorrected'
amalgamate(X, ...)
```

## Arguments

<code>X</code>	a compositional dataset
<code>...</code>	a series of new labels assigned to strings or vectors of strings denoting the components that need amalgamating

## Value

an object of the same class as `X` with fewer components

## Examples

```
data(Namib)
HMcomponents <- c("Zr", "tm", "rt", "TiOx", "sph", "ap", "ep",
                  "gt", "st", "amp", "cpx", "opx")
am <- amalgamate(Namib$PTHM, feldspars=c("KF", "P"),
                 lithics=c("Lm", "Lv", "Ls"), heavies=HMcomponents)
plot(ternary(am))
```

---

as.acomp	<i>create an acomp object</i>
----------	-------------------------------

---

**Description**

Convert an object of class `compositional` to an object of class `acomp` for use in the `compositions` package

**Usage**

```
as.acomp(x)
```

**Arguments**

x	an object of class <code>compositional</code>
---	---

**Value**

a `data.frame`

**Examples**

```
data(Namib)
qfl <- ternary(Namib$PT,c('Q'),c('KF','P'),c('Lm','Lv','Ls'))
plot(qfl,type="QFL.dickinson")
qfl.acomp <- as.acomp(qfl)
## uncomment the next two lines to plot an error
## ellipse using the compositions package:
# library(compositions)
# ellipses(mean(qfl.acomp),var(qfl.acomp),r=2)
```

---

as.compositional	<i>create a compositional object</i>
------------------	--------------------------------------

---

**Description**

Convert an object of class `matrix`, `data.frame` or `acomp` to an object of class `compositional`

**Usage**

```
as.compositional(x, method = NULL, colmap = "rainbow")
```

**Arguments**

x	an object of class <code>matrix</code> , <code>data.frame</code> or <code>acomp</code>
method	dissimilarity measure, either <code>'aitchison'</code> for Aitchison's CLR-distance or <code>'bray'</code> for the Bray-Curtis distance.
colmap	the colour map to be used in pie charts.

**Value**

an object of class `compositional`

**Examples**

```
data(Namib)
PT.acomp <- as.acomp(Namib$PT)
PT.compositional <- as.compositional(PT.acomp)
print(Namib$PT$x - PT.compositional$x)
## uncomment the following lines for an illustration of using this
## function to integrate the provenance package with compositions
# library(compositions)
# data(Glacial)
# a.glac <- acomp(Glacial)
# c.glac <- as.compositional(a.glac)
# summaryplot(c.glac, ncol=8)
```

---

as.counts	<i>create a counts object</i>
-----------	-------------------------------

---

**Description**

Convert an object of class `matrix` or `data.frame` to an object of class `counts`

**Usage**

```
as.counts(x, method = "chisq", colmap = "rainbow")
```

**Arguments**

x	an object of class <code>matrix</code> or <code>data.frame</code>
method	either "chisq" (for the chi-square distance) or "bray" (for the Bray-Curtis distance)
colmap	the colour map to be used in pie charts.

**Value**

an object of class `counts`

**Examples**

```
X <- matrix(c(0,100,0,30,11,2,94,36,0), nrow=3, ncol=3)
rownames(X) <- 1:3
colnames(X) <- c('a', 'b', 'c')
comp <- as.counts(X)
d <- diss(comp)
```

---

as.data.frame	<i>create a data.frame object</i>
---------------	-----------------------------------

---

**Description**

Convert an object of class `compositional` to a `data.frame` for use in the `robCompositions` package

**Usage**

```
## S3 method for class 'compositional'
as.data.frame(x, ...)

## S3 method for class 'counts'
as.data.frame(x, ...)
```

**Arguments**

<code>x</code>	an object of class <code>compositional</code>
<code>...</code>	optional arguments to be passed on to the generic function

**Value**

a `data.frame`

**Examples**

```
data(Namib)
Major.frame <- as.data.frame(Namib$Major)
## uncomment the next two lines to plot an error
## ellipse using the robCompositions package:
# library(robCompositions)
# plot(pcaCoDa(Major.frame))
```

---

botev	<i>Compute the optimal kernel bandwidth</i>
-------	---

---

**Description**

Uses the diffusion algorithm of Zdravko Botev (2011) to calculate the bandwidth for kernel density estimation

**Usage**

```
botev(x)
```

**Arguments**

x                      a vector of ordinal data

**Value**

a scalar value with the optimal bandwidth

**Author(s)**

Dzdravko Botev

**References**

Botev, Z. I., J. F. Grotowski, and D. P. Kroese. "Kernel density estimation via diffusion." *The Annals of Statistics* 38.5 (2010): 2916-2957.

**Examples**

```
fname <- system.file("DZ.csv", package="provenance")
bw <- botev(read.distributional(fname)$x$N1)
print(bw)
```

---

bray.diss	<i>Bray-Curtis dissimilarity</i>
-----------	----------------------------------

---

**Description**

Calculates the Bray-Curtis dissimilarity between two samples

**Usage**

```
bray.diss(x, y)
```

**Arguments**

x                      a vector containing the first compositional sample  
y                      a vector of length(x) containing the second compositional sample

**Value**

a scalar value

**Examples**

```
data(Namib)
print(bray.diss(Namib$HM$x["N1",], Namib$HM$x["N2",]))
```



---

CA	<i>Correspondence Analysis</i>
----	--------------------------------

---

**Description**

Performs Correspondence Analysis of point-counting data

**Usage**

```
CA(x, nf = 2, ...)
```

**Arguments**

x	an object of class counts
nf	number of correspondence factors (dimensions)
...	optional arguments to the corresp function of the MASS package

**Value**

an object of classes CA, which is synonymous to the MASS packages' correspondence class.

**Examples**

```
data(Namib)
plot(CA(Namib$PT))
```

---

central	<i>Calculate central compositions</i>
---------	---------------------------------------

---

**Description**

Computes the geometric mean composition of a continuous mixture of point-counting data.

**Usage**

```
central(x, ...)
```

**Arguments**

x	an object of class counts
...	optional arguments

**Details**

The central composition assumes that the observed point-counting distribution is the combination of two sources of scatter: counting uncertainty and true geological dispersion.

**Value**

an [5 x n] matrix with n being the number of categories and the rows containing:

**theta** the ‘central’ composition.

**err** the standard error for the central composition.

**sigma** the overdispersion parameter, i.e. the coefficient of variation of the underlying logistic normal distribution. `central` computes a continuous mixture model for each component (column) separately. Covariance terms are not reported.

**LL** the lower limit of a ‘1 sigma’ region for `codetheta`.

**UL** the upper limit of a ‘1 sigma’ region for `codetheta`.

**mswd** the mean square of the weighted deviates, a.k.a. reduced chi-square statistic.

**p.value** the p-value for age homogeneity

---

CLR

*Centred logratio transformation*


---

**Description**

Calculates Aitchison’s centered logratio transformation for a dataset of class `compositional` or a compositional data matrix.

**Usage**

```
CLR(x, ...)
```

```
## Default S3 method:
```

```
CLR(x, inverse = FALSE, ...)
```

```
## S3 method for class 'compositional'
```

```
CLR(x, ...)
```

**Arguments**

`x` an object of class `compositional` OR a matrix of numerical values

`...` optional arguments

`inverse` perform the inverse inverse logratio transformation?

**Value**

a matrix of CLR coordinates OR an object of class `compositional` (if `inverse=TRUE`)

**Examples**

```
# The following code shows that applying provenance's PCA function
# to compositional data is equivalent to applying R's built-in
# princomp function to the CLR transformed data.
data(Namib)
plot(PCA(Namib$Major))
dev.new()
clrdat <- CLR(Namib$Major)
biplot(princomp(clrdat))
```

---

combine

---

*Combine samples of distributional data*


---

**Description**

Lumps all single grain analyses of several samples together under a new name

**Usage**

```
combine(X, ...)
```

**Arguments**

X	a distributional dataset
...	a series of new labels assigned to strings or vectors of strings denoting the samples that need amalgamating

**Value**

a distributional data object with fewer samples than X

**Examples**

```
data(Namib)
combined <- combine(Namib$DZ,
  east=c('N3','N4','N5','N6','N7','N8','N9','N10'),
  west=c('N1','N2','N11','N12','T8','T13'))
summaryplot(KDEs(combined))
```

densities

*A list of rock and mineral densities***Description**

List of rock and mineral densities using the following abbreviations: Q (quartz), KF (K-feldspar), P (plagioclase), F (feldspar), Lvf (felsic/porphyritic volcanic rock fragments), Lvm (microlithic / porphyritic / trachitic volcanic rock fragments), Lcc (calcite), Lcd (dolomite), Lp (marl), Lch (chert), Lms (argillaceous / micaceous rock fragments), Lmv (metavolcanics), Lmf (metasediments), Lmb (metabasites), Lv (volcanic rock fragments), Lc (carbonates), Ls (sedimentary rock fragments), Lm (metamorphic rock fragments), Lu (serpentinite), mica, opaques, FeOx (Fe-oxides), turbids, zr (zircon), tm (tourmaline), rt (rutile), TiOx (Ti-oxides), sph (titanite), ap (apatite), mon (monazite), oth (other minerals), ep (epidote), othLgM (prehnite + pumpellyite + lawsonite + carpholite), gt (garnet), ctd (chloritoid), st (staurolite), and (andalusite), ky (kyanite), sil (sillimanite), amp (amphibole), px (pyroxene), cpx (clinopyroxene), opx (orthopyroxene), ol (olivine), spinel and othHM (other heavy minerals).

**Author(s)**

Alberto Resentini and Pieter Vermeesch

**References**

Resentini, A, Malusa M G and Garzanti, E. "MinSORTING: An Excel worksheet for modelling mineral grain-size distribution in sediments, with application to detrital geochronology and provenance studies." *Computers & Geosciences* 59 (2013): 90-97.

Garzanti, E, Ando, S and Vezzoli, G. "Settling equivalence of detrital minerals and grain-size dependence of sediment composition." *Earth and Planetary Science Letters* 273.1 (2008): 138-151.

**See Also**

restore, minsorting

**Examples**

```
data(Namib,densities)
N8 <- subset(Namib$HM,select="N8")
distribution <- minsorting(N8,densities,phi=2,sigmaphi=1,medium="air",by=0.05)
plot(distribution)
```

---

diss	<i>Calculate the dissimilarity matrix between two distributional or compositional datasets</i>
------	--

---

### Description

Calculate the dissimilarity matrix between two datasets of class distributional or compositional using the Kolmogorov-Smirnov, Sircombe-Hazelton, Aitchison or Bray Curtis distance

### Usage

```
diss(x, method)

## S3 method for class 'distributional'
diss(x, method = NULL)

## S3 method for class 'compositional'
diss(x, method = NULL)

## S3 method for class 'counts'
diss(x, method = NULL)
```

### Arguments

x	an object of class distributional, compositional or counts
method	(optional) either "KS", "Kuiper", "SH", "aitchison", "bray" or "chisq"

### Value

an object of class diss

### Examples

```
data(Namib)
print(round(100*diss(Namib$DZ)))
```

---

endmembers	<i>Petrographic end-member compositions</i>
------------	---

---

### Description

A compositional dataset comprising the mineralogical compositions of the following end-members: undissected\_magmatic\_arc, dissected\_magmatic\_arc, ophiolite, recycled\_clastic, undissected\_continental\_block, transitional\_continental\_block, dissected\_continental\_block, subcreted\_axial\_belt and subducted\_axial\_belt

**Author(s)**

Alberto Resentini and Pieter Vermeesch

**References**

Resentini, A, Malusa M G and Garzanti, E. "MinSORTING: An Excel worksheet for modelling mineral grain-size distribution in sediments, with application to detrital geochronology and provenance studies." Computers & Geosciences 59 (2013): 90-97.

Garzanti, E, Ando, S and Vezzoli, G. "Settling equivalence of detrital minerals and grain-size dependence of sediment composition." Earth and Planetary Science Letters 273.1 (2008): 138-151.

**See Also**

minsorting

**Examples**

```
data(endmembers,densities)
ophiolite <- subset(endmembers,select="ophiolite")
plot(minsorting(ophiolite,densities,by=0.05))
```

---

get.f

---

*Calculate the largest fraction that is likely to be missed*


---

**Description**

For a given sample size, returns the largest fraction which has been sampled with  $(1-p) \times 100\%$  likelihood.

**Usage**

```
get.f(n, p = 0.05)
```

**Arguments**

n	the number of grains in the detrital sample
p	the required level of confidence

**Value**

the largest fraction that is sampled with at least  $(1-p) \times 100\%$  certainty

**References**

Vermeesch, Pieter. "How many grains are needed for a provenance study?." Earth and Planetary Science Letters 224.3 (2004): 441-451.

### Examples

```
print(get.f(60))
print(get.f(117))
```

---

get.n

*Calculate the number of grains required to achieve a desired level of sampling resolution*

---

### Description

Returns the number of grains that need to be analysed to decrease the likelihood of missing any fraction greater than a given size below a given level.

### Usage

```
get.n(p = 0.05, f = 0.05)
```

### Arguments

p	the probability that all n grains in the sample have missed at least one fraction of size f
f	the size of the smallest resolvable fraction ( $0 < f < 1$ )
n,	the number of grains in the sample

### Value

the number of grains needed to reduce the chance of missing at least one fraction f of the total population to less than p

### References

Vermeesch, Pieter. "How many grains are needed for a provenance study?." Earth and Planetary Science Letters 224.3 (2004): 441-451.

### Examples

```
# number of grains required to be 99% that no fraction greater than 5% was missed:
print(get.n(0.01))
# number of grains required to be 90% that no fraction greater than 10% was missed:
print(get.n(p=0.1, f=0.1))
```

---

<code>get.p</code>	<i>Calculate the probability of missing a given population fraction</i>
--------------------	---

---

**Description**

For a given sample size, returns the likelihood of missing any fraction greater than a given size

**Usage**

```
get.p(n, f = 0.05)
```

**Arguments**

<code>n</code>	the number of grains in the detrital sample
<code>f</code>	the size of the smallest resolvable fraction ( $0 < f < 1$ )

**Value**

the probability that all `n` grains in the sample have missed at least one fraction of size `f`

**References**

Vermeesch, Pieter. "How many grains are needed for a provenance study?." *Earth and Planetary Science Letters* 224.3 (2004): 441-451.

**Examples**

```
print(get.p(60))
print(get.p(117))
```

---

GPA	<i>Generalised Procrustes Analysis of configurations</i>
-----	--

---

**Description**

Given a number of (2D) configurations, this function uses a combination of transformations (reflections, rotations, translations and scaling) to find a 'consensus' configuration which best matches all the component configurations in a least-squares sense.

**Usage**

```
GPA(X, scale = TRUE)
```

**Arguments**

<code>X</code>	a list of dissimilarity matrices
<code>scale</code>	boolean flag indicating if the transformation should include the scaling operation



**Value**

a two column vector with the coordinates of the group configuration

**See Also**

procrustes

---

indscal

*Individual Differences Scaling of provenance data*


---

**Description**

Performs 3-way Multidimensional Scaling analysis using Carroll and Chang (1970)'s Individual Differences SCALing method as implemented using De Leeuw and Mair (2011)'s stress majorization algorithm.

**Usage**

```
indscal(..., type = "ordinal")
```

**Arguments**

... a sequence of datasets of class `distributional` or `compositional`  
type is either "ratio" or "ordinal"

**Value**

an object of class `INDSCAL`, i.e. a list containing the following items:

delta: Observed dissimilarities

obsdiss: List of observed dissimilarities, normalized

confdiss: List of configuration dissimilarities

conf: List of matrices of final configurations

gspace: Joint configurations aka group stimulus space

cweights: Configuration weights

stress: Stress-1 value

spp: Stress per point

sps: Stress per subject (matrix)

ndim: Number of dimensions

model: Type of smacof model

niter: Number of iterations

nobj: Number of objects

**Author(s)**

Jan de Leeuw and Patrick Mair

**References**

de Leeuw, J., & Mair, P. (2009). Multidimensional scaling using majorization: The R package smacof. Journal of Statistical Software, 31(3), 1-30, < <http://www.jstatsoft.org/v31/i03/>>

**Examples**

```
data(Namib)
plot(indscal(Namib$DZ,Namib$HM))
```

---

KDE

---

*Create a kernel density estimate*


---

**Description**

Turns a vector of numbers into an object of class KDE using a combination of the Botev (2010) bandwidth selector and the Abramson (1982) adaptive kernel bandwidth modifier.

**Usage**

```
KDE(x, from = NA, to = NA, bw = NA, adaptive = TRUE, log = FALSE,
    n = 512, ...)
```

**Arguments**

x	a vector of numbers
from	minimum age of the time axis. If NULL, this is set automatically
to	maximum age of the time axis. If NULL, this is set automatically
bw	the bandwidth of the KDE. If NULL, bw will be calculated automatically using <code>botev()</code>
adaptive	boolean flag controlling if the adaptive KDE modifier of Abramson (1982) is used
log	transform the ages to a log scale if TRUE
n	horizontal resolution of the density estimate
...	optional arguments to be passed on to density

**Value**

an object of class KDE, i.e. a list containing the following items:

x: horizontal plot coordinates

y: vertical plot coordinates

bw: the base bandwidth of the density estimate

ages: the data values from the input to the KDE function

**See Also**

KDEs

**Examples**

```
data(Namib)
samp <- Namib$DZ$x[['N1']]
dens <- KDE(samp,0,3000,kernel="epanechnikov")
plot(dens)
```

KDEs

*Generate an object of class KDEs***Description**

Convert a dataset of class distributional into an object of class KDEs for further processing by the summaryplot function.

**Usage**

```
KDEs(x, from = NA, to = NA, bw = NA, samebandwidth = TRUE,
      adaptive = TRUE, normalise = FALSE, log = FALSE, n = 512, ...)
```

**Arguments**

x	an object of class distributional
from	minimum limit of the x-axis.
to	maximum limit of the x-axis.
bw	the bandwidth of the kernel density estimates. If bw = NA, the bandwidth will be set automatically using botev()
samebandwidth	boolean flag indicating whether the same bandwidth should be used for all samples. If samebandwidth = TRUE and bw = NULL, then the function will use the median bandwidth of all the samples.
adaptive	boolean flag switching on the adaptive bandwidth modifier of Abramson (1982)
normalise	boolean flag indicating whether or not the KDEs should all integrate to the same value.
log	boolean flag indicating whether the data should be plotted on a logarithmic scale.
n	horizontal resolution of the density estimates
...	optional parameters to be passed on to density

**Value**

an object of class KDEs, i.e. a list containing the following items:

kdes: a named list with objects of class KDE

from: the beginning of the common time scale

to: the end of the common time scale

themax: the maximum probability density of all the KDEs

pch: the plot symbol to be used by plot.KDEs

xlabel: the x-axis label to be used by plot.KDEs

**See Also**

KDE

**Examples**

```
data(Namib)
KDEs <- KDEs(Namib$DZ, 0, 3000, pch=NA)
summaryplot(KDEs, ncol=3)
```

---

KS.diss

*Kolmogorov-Smirnov dissimilarity*


---

**Description**

Returns the Kolmogorov-Smirnov dissimilarity between two samples

**Usage**

```
KS.diss(x, y)
```

**Arguments**

x                    the first sample as a vector

y                    the second sample as a vector

**Value**

a scalar value representing the maximum vertical distance between the two cumulative distributions

**Examples**

```
data(Namib)
print(KS.diss(Namib$DZ$x[['N1']], Namib$DZ$x[['T8']]))
```

---

Kuiper.diss	<i>Kuiper dissimilarity</i>
-------------	-----------------------------

---

**Description**

Returns the Kuiper dissimilarity between two samples

**Usage**

```
Kuiper.diss(x, y)
```

**Arguments**

x	the first sample as a vector
y	the second sample as a vector

**Value**

a scalar value representing the sum of the maximum vertical distances above and below the cumulative distributions of x and y

**Examples**

```
data(Namib)
print(Kuiper.diss(Namib$DZ$x[['N1']], Namib$DZ$x[['T8']]))
```

---

lines.ternary	<i>Ternary line plotting</i>
---------------	------------------------------

---

**Description**

Add lines to an existing ternary diagram

**Usage**

```
## S3 method for class 'ternary'
lines(x, ...)
```

**Arguments**

x	an object of class ternary, or a three-column data frame or matrix
...	optional arguments to the generic lines function

### Examples

```
tern <- ternary(Namib$PT, 'Q', c('KF', 'P'), c('Lm', 'Lv', 'Ls'))
plot(tern, pch=21, bg='red', labels=NULL)
middle <- matrix(c(0.01, 0.49, 0.01, 0.49, 0.98, 0.02), 2, 3)
lines(ternary(middle))
```

---

MDS

---

*Multidimensional Scaling*


---

### Description

Performs classical or nonmetric Multidimensional Scaling analysis of provenance data

### Usage

```
MDS(x, ...)

## Default S3 method:
MDS(x, fn, classical = FALSE, k = 2,
     bootstrap = FALSE, nb = 10, ...)

## S3 method for class 'compositional'
MDS(x, classical = FALSE, k = 2,
     bootstrap = FALSE, nb = 10, ...)

## S3 method for class 'counts'
MDS(x, classical = FALSE, k = 2, bootstrap = FALSE,
     nb = 10, ...)

## S3 method for class 'distributional'
MDS(x, classical = FALSE, k = 2,
     bootstrap = FALSE, nb = 10, ...)

## S3 method for class 'diss'
MDS(x, classical = FALSE, k = 2, ...)
```

### Arguments

x	an object of class <code>distributional</code> , <code>compositional</code> , <code>counts</code> or <code>diss</code>
...	optional arguments to be passed onto <code>diss</code> (if x is of class <code>compositional</code> or <code>distributional</code> ) or onto <code>cmdscale</code> or <code>isoMDS</code> (if x is of class <code>dist</code> ).
fn	a dissimilarity function. One of either <code>diss.compositional</code> , <code>diss.counts</code> , or <code>diss.distributional</code>
classical	boolean flag indicating whether classical (TRUE) or nonmetric (FALSE) MDS should be used
k	the desired dimensionality of the solution

bootstrap	resample the data to calculate confidence polygons for the MDS configuration
nb	number of bootstrap resamples

**Value**

an object of class MDS, i.e. a list containing the following items:

- points: a two column vector of the fitted configuration
- classical: a boolean flag indicating whether the MDS configuration was obtained by classical (TRUE) or nonmetric (FALSE) MDS.
- diss: the dissimilarity matrix used for the MDS analysis
- stress: (only if classical=TRUE) the final stress achieved (in percent)

**Examples**

```
data(Namib)
plot(MDS(Namib$Major,classical=TRUE))
```

minsorting

*Assess settling equivalence of detrital components***Description**

Models grain size distribution of minerals and rock fragments of different densities

**Usage**

```
minsorting(X, dens, sname = NULL, phi = 2, sigmaphi = 1,
  medium = "freshwater", from = -2.25, to = 5.5, by = 0.25)
```

**Arguments**

X	an object of class compositional
dens	a vector of mineral and rock densities
sname	sample name if unspecified, the first sample of the dataset will be used
phi	the mean grain size of the sample in Krumbein's phi units
sigmaphi	the standard deviation of the grain size distribution, in phi units
medium	the transport medium, one of either "air", "freshwater" or "seawater"
from	the minimum grain size to be evaluated, in phi units
to	the maximum grain size to be evaluated, in phi units
by	the grain size interval of the output table, in phi units

**Value**

an object of class `minsorting`, i.e. a list with two tables:

`mfract`: the grain size distribution of each mineral (sum of the columns = 1)

`mcomp`: the composition of each mineral (sum of the rows = 1)

**Author(s)**

Alberto Resentini and Pieter Vermeesch

**References**

Resentini, A, Malusa, M G and Garzanti, E. "MinSORTING: An Excel worksheet for modelling mineral grain-size distribution in sediments, with application to detrital geochronology and provenance studies." *Computers & Geosciences* 59 (2013): 90-97.

Garzanti, E, Ando, S and Vezzoli, G. "Settling equivalence of detrital minerals and grain-size dependence of sediment composition." *Earth and Planetary Science Letters* 273.1 (2008): 138-151.

**See Also**

`restore`

**Examples**

```
data(endmembers,densities)
distribution <- minsorting(endmembers,densities,sname='ophiolite',phi=2,
                           sigmaphi=1,medium="seawater",by=0.05)
plot(distribution,cumulative=FALSE)
```

---

Namib

*An example dataset*

---

**Description**

A large dataset of provenance data from Namibia comprised of 14 sand samples from the Namib Sand Sea and 2 samples from the Orange River.

**Details**

Namib is a list containing the following 6 items:

DZ: a distributional dataset containing the zircon U-Pb ages for ca. 100 grains from each sample, as well as their (1-sigma) analytical uncertainties.

PT: a compositional dataset with the bulk petrography of the samples, i.e. the quartz ('Q'), K-feldspar ('KF'), plagioclase ('P'), and lithic fragments of metamorphic ('Lm'), volcanic ('Lv') and sedimentary ('Ls') origin.

HM: a compositional dataset containing the heavy mineral composition of the samples, comprised of zircon ('zr'), tourmaline ('tm'), rutile ('rt'), Ti-oxides ('TiOx'), titanite ('sph'), apatite ('ap'),



epidote ('ep'), garnet ('gt'), staurolite ('st'), andalusite ('and'), kyanite ('ky'), sillimanite ('sil'), amphibole ('amp'), clinopyroxene ('cpx') and orthopyroxene ('opx').

PTHM: a compositional dataset combining the variables contained in PT and HM plus 'mica', 'opaques', 'turbids' and 'other' transparent heavy minerals ('LgM'), normalised to 100.

Major: a compositional dataset listing the concentrations (in wt TiO<sub>2</sub>, P<sub>2</sub>O<sub>5</sub> and MnO).

Trace: a compositional data listing the concentrations (in ppm) of Rb, Sr, Ba, Sc, Y, La, Ce, Pr, Nd, Sm, Gd, Dy, Er, Yb, Th, U, Zr, Hf, V, Nb, Cr, Co, Ni, Cu, Zn, Ga and Pb.

### Author(s)

Pieter Vermeesch and Eduardo Garzanti

### References

Vermeesch, P. and Garzanti, E., Making geological sense of 'Big Data' in sedimentary provenance analysis, *Chemical Geology* 409 (2015) 20-27

### Examples

```
data(Namib)
samp <- Namib$DZ$x[['N1']]
dens <- KDE(samp,0,3000)
plot(dens)
```

---

PCA

*Principal Component Analysis*

---

### Description

Performs PCA of compositional data using a centred logratio distance

### Usage

```
PCA(x, ...)
```

### Arguments

x                    an object of class `compositional`  
 ...                  optional arguments to R's `princomp` function

### Value

an object of classes `PCA`, which is synonymous to the stats packages' `prcomp` class.

Examples

```
data(Namib)
plot(MDS(Namib$Major,classical=TRUE))
dev.new()
plot(PCA(Namib$Major),asp=1)
print("This example demonstrates the equivalence of classical MDS and PCA")
```

---

plot.CA	<i>Point-counting biplot</i>
---------	------------------------------

---

Description

Plot the results of a correspondence analysis as a biplot

Usage

```
## S3 method for class 'CA'
plot(x, levels = NULL, labelcol = c("black", "blue"),
      vectorcol = "red", ...)
```

Arguments

x	an object of class CA
levels	a vector of length nrow(dat\$x) with values that are to be used for the colour scale
labelcol	two element vector with the colours that are to be assigned to the labels of the samples corresponding to the minimum and maximum value of levels. If levels==NULL, then the first item of the vector is used for all sample labels.
vectorcol	colour of the vector loadings for the variables
...	optional arguments of the generic <a href="#">biplot</a> function

See Also

CA

Examples

```
data(Namib)
plot(CA(Namib$PT))
```

---

plot.compositional      *Plot a pie chart*


---

**Description**

Plots an object of class compositional as a pie chart

**Usage**

```
## S3 method for class 'compositional'
plot(x, sname, annotate = TRUE, colmap = NULL,
     ...)
```

**Arguments**

x	an object of class compositional
sname	the sample name
annotate	a boolean flag controlling if the pies of the pie-chart should be labeled
colmap	an optional string with the name of one of R's built-in colour palettes (e.g., heat.colors, terrain.colors, topo.colors, cm.colors), which are to be used for plotting the data.
...	optional parameters to be passed on to the graphics object

**Examples**

```
data(Namib)
plot(Namib$Major, 'N1', colmap='heat.colors')
```

---

plot.distributional      *Plot continuous data as histograms or cumulative age distributions*


---

**Description**

Plot one or several samples from a distributional dataset as a histogram or Cumulative Age Distributions (CAD).

**Usage**

```
## S3 method for class 'distributional'
plot(x, snames = NULL, annotate = TRUE,
     CAD = FALSE, pch = NA, verticals = TRUE, colmap = NULL, ...)
```

**Arguments**

x	an object of class <code>distributional</code>
snames	a string or a vector of string with the names of the samples that need plotting if <code>snames</code> is a vector, then the function will default to a CAD.
annotate	boolean flag indicating whether the x- and y-axis should be labeled
CAD	boolean flag indicating whether the data should be plotted as a cumulative age distribution or a histogram. For multi-sample plots, the function will override this value with TRUE.
pch	an optional symbol to mark the sample points along the CAD
verticals	boolean flag indicating if the horizontal lines of the CAD should be connected by vertical lines
colmap	an optional string with the name of one of R's built-in colour palettes (e.g., <code>heat.colors</code> , <code>terrain.colors</code> , <code>topo.colors</code> , <code>cm.colors</code> ), which are to be used for plotting the data.
...	optional arguments to the generic plot function

**Examples**

```
data(Namib)
plot(Namib$DZ, c('N1', 'N2'))
```

---

plot.GPA

---

*Plot a Procrustes configuration*


---

**Description**

Plots the group configuration of a Generalised Procrustes Analysis

**Usage**

```
## S3 method for class 'GPA'
plot(x, pch = NA, pos = NULL, col = "black",
      bg = "white", cex = 1, ...)
```

**Arguments**

x	an object of class <code>GPA</code>
pch	plot symbol
pos	position of the sample labels relative to the plot symbols if <code>pch != NA</code>
col	plot colour (may be a vector)
bg	background colour (may be a vector)
cex	relative size of plot symbols
...	optional arguments to the generic plot function

**See Also**

procrustes

**Examples**

```
data(Namib)
GPA <- procrustes(Namib$DZ,Namib$HM)
coast <- c('N1','N2','N3','N10','N11','N12','T8','T13')
snames <- names(Namib$DZ)
bgcol <- rep('yellow',length(snames))
bgcol[which(snames %in% coast)] <- 'red'
plot(GPA,pch=21,bg=bgcol)
```

plot.INDSCAL

*Plot an INDSCAL group configuration and source weights***Description**

Given an object of class INDSCAL, generates two plots: the group configuration and the subject weights. Together, these describe a 3-way MDS model.

**Usage**

```
## S3 method for class 'INDSCAL'
plot(x, asp = 1, pch = NA, pos = NULL,
     col = "black", bg = "white", cex = 1, xlab = "X", ylab = "Y",
     xaxt = "n", yaxt = "n", ...)
```

**Arguments**

x	an object of class INDSCAL
asp	the aspect ratio of the plot
pch	plot symbol (may be a vector)
pos	position of the sample labels relative to the plot symbols if pch != NA
col	plot colour (may be a vector)
bg	background colour (may be a vector)
cex	relative size of plot symbols
xlab	a string with the label of the x axis
ylab	a string with the label of the y axis
xaxt	if = 'y', adds ticks to the x axis
yaxt	if = 'y', adds ticks to the y axis
...	optional arguments to the generic plot function

**See Also**

indscal

**Examples**

```
data(Namib)
coast <- c('N1','N2','N3','N10','N11','N12','T8','T13')
snames <- names(Namib$DZ)
pch <- rep(21,length(snames))
pch[which(snames %in% coast)] <- 22
plot(indscal(Namib$DZ,Namib$HM),pch=pch)
```

---

plot.KDE	<i>Plot a kernel density estimate</i>
----------	---------------------------------------

---

**Description**

Plots an object of class KDE

**Usage**

```
## S3 method for class 'KDE'
plot(x, pch = "|", xlab = "age [Ma]", ylab = "", ...)
```

**Arguments**

- x                    an object of class KDE
- pch                the symbol used to show the samples. May be a vector. Set pch = NA to turn them off.
- xlab               the label of the x-axis
- ylab               the label of the y-axis
- ...                optional parameters to be passed on to the graphics object

**See Also**

KDE

**Examples**

```
data(Namib)
samp <- Namib$DZ$x[['N1']]
dens <- KDE(samp,from=0,to=3000)
plot(dens)
```

---

plot.KDEs	<i>Plot one or more kernel density estimates</i>
-----------	--

---

**Description**

Plots an object of class KDEs

**Usage**

```
## S3 method for class 'KDEs'
plot(x, sname = NA, annotate = TRUE, pch = "|", ...)
```

**Arguments**

x	an object of class KDEs
sname	optional sample name. If sname=NA, all samples are shown on a summary plot
annotate	add a time axis?
pch	symbol to be used to mark the sample points along the x-axis. Change to NA to omit.
...	optional parameters to be passed on to the summaryplot function

**See Also**

KDEs summaryplot

**Examples**

```
data(Namib)
kdes <- KDEs(Namib$DZ)
plot(kdes, ncol=2)
```

---

plot.MDS	<i>Plot an MDS configuration</i>
----------	----------------------------------

---

**Description**

Plots the coordinates of a multidimensional scaling analysis as an X-Y scatter plot or 'map' and, if x\$classical = FALSE, a Shepard plot.

**Usage**

```
## S3 method for class 'MDS'
plot(x, nnlines = FALSE, pch = NA, pos = NULL,
     cex = 1, col = "black", bg = "white", oma = rep(1, 4),
     mar = rep(2, 4), mgp = c(2, 1, 0), xpd = NA, ...)
```

**Arguments**

x	an object of class MDS
nnlines	if TRUE, draws nearest neighbour lines
pch	plot character (see ?plot for details). May be a vector.
pos	position of the sample labels relative to the plot symbols if pch != NA
cex	relative size of plot symbols (see ?par for details)
col	plot colour (may be a vector)
bg	background colour (may be a vector)
oma	A vector of the form c(bottom, left, top, right) giving the size of the outer margins in lines of text.
mar	A numerical vector of the form c(bottom, left, top, right) that gives the number of lines of margin to be specified on the four sides of the plot.
mgp	The margin line (in mex units) for the axis title, axis labels and axis line. See ?par for further details.
xpd	A logical value or NA. See ?par for further details.
...	optional arguments to the generic plot function

**See Also**

MDS

**Examples**

```
data(Namib)
mds <- MDS(Namib$DZ)
coast <- c('N1', 'N2', 'N3', 'N10', 'N11', 'N12', 'T8', 'T13')
snames <- names(Namib$DZ)
bgcol <- rep('yellow', length(snames))
bgcol[which(snames %in% coast)] <- 'red'
plot(mds, pch=21, bg=bgcol)
```

plot.minsorting

*Plot inferred grain size distributions***Description**

Plot the grain size distributions of the different minerals under consideration

**Usage**

```
## S3 method for class 'minsorting'
plot(x, cumulative = FALSE, components = NULL,
     ...)
```



**Arguments**

x	an object of class minsorting
cumulative	boolean flag indicating whether the grain size distribution should be plotted as a density or cumulative probability curve.
components	string or list of strings with the names of a subcomposition that needs plotting
...	optional parameters to be passed on to graphics::matplot (see ?par for details)

**See Also**

minsorting

**Examples**

```
data(endmembers,densities)
OPH <- subset(endmembers,select="ophiolite")
distribution <- minsorting(OPH,densities,phi=2,sigmaphi=1,
                           medium="air",by=0.05)
plot(distribution,components=c('F','px','opaques'))
```

---

plot.PCA

---

*Compositional biplot*


---

**Description**

Plot the results of a principal components analysis as a biplot

**Usage**

```
## S3 method for class 'PCA'
plot(x, levels = NULL, labelcol = c("black", "blue"),
     vectorcol = "red", choices = 1L:2L, scale = 1, pc.biplot = FALSE,
     ...)
```

**Arguments**

x	an object of class PCA
levels	a vector of length nrow(dat\$x) with values that are to be used for the colour scale
labelcol	two element vector with the colours that are to be assigned to the labels of the samples corresponding to the minimum and maximum value of levels. If levels==NULL, then the first item of the vector is used for all sample labels.
vectorcol	colour of the vector loadings for the variables
choices	see the help pages of the <a href="#">biplot</a> function.
scale	see the help pages of the <a href="#">biplot</a> function.
pc.biplot	see the help pages of the <a href="#">biplot</a> function.
...	optional arguments of the generic <a href="#">biplot</a> function

**See Also**

PCA

**Examples**

```
data(Namib)
plot(PCA(Namib$Major))
```

---

plot.ternary	<i>Plot a ternary diagram</i>
--------------	-------------------------------

---

**Description**

Plots triplets of compositional data on a ternary diagram

**Usage**

```
## S3 method for class 'ternary'
plot(x, type = "grid", pch = NA, pos = NULL,
     labels = names(x), showpath = FALSE, bg = NA,
     col = "cornflowerblue", ticks = seq(0, 1, 0.25), ticklength = 0.02,
     lty = 2, lwd = 1, ...)
```

**Arguments**

x	an object of class ternary, or a three-column data frame or matrix
type	adds annotations to the ternary diagram, one of either empty, grid, QFL.descriptive, QFL.folk or QFL.dickinson
pch	plot character, see ?par for details (may be a vector)
pos	position of the sample labels relative to the plot symbols if pch != NA
labels	vector of strings to be added to the plot symbols
showpath	if x has class SRDcorrected, and showpath==TRUE, the intermediate values of the SRD correction will be plotted on the ternary diagram as well as the final composition
bg	background colour for the plot symbols (may be a vector)
col	colour to be used for the background lines (if applicable)
ticks	vector of tick values between 0 and 1
ticklength	number between 0 and 1 to mark the length of the ticks
lty	line type for the annotations (see type)
lwd	line thickness for the annotations
...	optional arguments to the generic points function

**See Also**

ternary

**Examples**

```
data(Namib)
tern <- ternary(Namib$PT,'Q',c('KF','P'),c('Lm','Lv','Ls'))
plot(tern,type='QFL.descriptive',pch=21,bg='red',labels=NULL)
```

---

points.ternary	<i>Ternary point plotting</i>
----------------	-------------------------------

---

**Description**

Add points to an existing ternary diagram

**Usage**

```
## S3 method for class 'ternary'
points(x, ...)
```

**Arguments**

x	an object of class ternary, or a three-column data frame or matrix
...	optional arguments to the generic points function

**Examples**

```
tern <- ternary(Namib$PT,'Q',c('KF','P'),c('Lm','Lv','Ls'))
plot(tern,pch=21,bg='red',labels=NULL)
# add the geometric mean composition as a yellow square:
gmean <- ternary(exp(colMeans(log(tern$x))))
points(gmean,pch=22,bg='yellow')
```

---

procrustes	<i>Generalised Procrustes Analysis of provenance data</i>
------------	---

---

**Description**

Given a number of input datasets, this function performs an MDS analysis on each of these and the feeds the resulting configurations into the GPA() function.

**Usage**

```
procrustes(...)
```

**Arguments**

... a sequence of datasets of classes distributional, n counts and compositional

**Value**

an object of class GPA, i.e. a list containing the following items:

points: a two column vector with the coordinates of the group configuration

labels: a list with the sample names

**Author(s)**

Pieter Vermeesch

**References**

Gower, J.C. (1975). Generalized Procrustes analysis, *Psychometrika*, 40, 33-50.

**See Also**

GPA

**Examples**

```
data(Namib)
gpa <- procrustes(Namib$DZ, Namib$HM)
plot(gpa)
```

---

provenance

*Menu-based interface for provenance*

---

**Description**

For those less familiar with the syntax of the R programming language, the `provenance()` function provides a user-friendly way to access the most important functionality in the form of a menu-based query interface. Further details and examples are provided on <http://provenance.london-geochron.com>

`provenance` provides statistical tools to interpret large amounts of distributional (single grain analyses) and compositional (mineralogical and bulk chemical) data from the command line, or using a menu-based user interface.

**Usage**

```
provenance()
```

**Details**

A list of documented functions may be viewed by typing `help(package='provenance')`. Detailed instructions are provided at <http://provenance.london-geochron.com> and in the Sedimentary Geology paper by Vermeesch, Resentini and Garzanti (2016).

**Author(s)**

Pieter Vermeesch

**Maintainer:** Pieter Vermeesch <p.vermeesch@ucl.ac.uk>

**References**

Vermeesch, P., Resentini, A. and Garzanti, E., an R package for statistical provenance analysis, Sedimentary Geology, doi:10.1016/j.sedgeo.2016.01.009.

Vermeesch, P., Resentini, A. and Garzanti, E., 2016, An R package for statistical provenance analysis, Sedimentary Geology, 336, 14-25.

**See Also**

<http://provenance.london-geochron.com>

Useful links:

- <http://provenance.london-geochron.com>

---

radialplot

---

*Visualise point-counting data on a radial plot*


---

**Description**

Implementation of a graphical device developed by Rex Galbraith to display several estimates of the same quantity that have different standard errors.

**Usage**

```
radialplot(x, num = 1, den = 2, from = NA, to = NA, t0 = NA,
  sigdig = 2, show.numbers = FALSE, pch = 21, levels = NA,
  clabel = "", bg = c("white", "red"), title = TRUE, ...)
```

**Arguments**

x	an object of class counts
num	index or name of the numerator variable
den	index or name of the denominator variable
from	minimum limit of the radial scale
to	maximum limit of the radial scale

<code>t0</code>	central value
<code>sigdig</code>	the number of significant digits of the numerical values reported in the title of the graphical output.
<code>show.numbers</code>	boolean flag (TRUE to show sample numbers)
<code>pch</code>	plot character (default is a filled circle)
<code>levels</code>	a vector with additional values to be displayed as different background colours of the plot symbols.
<code>clabel</code>	label of the colour legend
<code>bg</code>	a vector of two background colours for the plot symbols. If <code>levels=NA</code> , then only the first colour is used. If <code>levels</code> is a vector of numbers, then <code>bg</code> is used to construct a colour ramp.
<code>title</code>	add a title to the plot?
<code>...</code>	additional arguments to the generic <code>points</code> function

### Details

The radial plot (Galbraith, 1988, 1990) is a graphical device that was specifically designed to display heteroscedastic data, and is constructed as follows. Consider a set of dates  $\{t_1, \dots, t_i, \dots, t_n\}$  and uncertainties  $\{s[t_1], \dots, s[t_i], \dots, s[t_n]\}$ . Define  $z_i = z[t_i]$  to be a transformation of  $t_i$  (e.g.,  $z_i = \log[t_i]$ ), and let  $s[z_i]$  be its propagated analytical uncertainty (i.e.,  $s[z_i] = s[t_i]/t_i$  in the case of a logarithmic transformation). Create a scatterplot of  $(x_i, y_i)$  values, where  $x_i = 1/s[z_i]$  and  $y_i = (z_i - z_o)/s[z_i]$ , where  $z_o$  is some reference value such as the mean. The slope of a line connecting the origin of this scatterplot with any of the  $(x_i, y_i)$ s is proportional to  $z_i$  and, hence, the date  $t_i$ . These dates can be more easily visualised by drawing a radial scale at some convenient distance from the origin and annotating it with labelled ticks at the appropriate angles. While the angular position of each data point represents the date, its horizontal distance from the origin is proportional to the precision. Imprecise measurements plot on the left hand side of the radial plot, whereas precise age determinations are found further towards the right. Thus, radial plots allow the observer to assess both the magnitude and the precision of quantitative data in one glance.

### References

- Galbraith, R.F., 1988. Graphical display of estimates having differing standard errors. *Technometrics*, 30(3), pp.271-281.
- Galbraith, R.F., 1990. The radial plot: graphical assessment of spread in ages. *International Journal of Radiation Applications and Instrumentation. Part D. Nuclear Tracks and Radiation Measurements*, 17(3), pp.207-214.
- Galbraith, R.F. and Laslett, G.M., 1993. Statistical models for mixed fission track ages. *Nuclear Tracks and Radiation Measurements*, 21(4), pp.459-470.

### Examples

```
data(Namib)
radialplot(Namib$PT, components=c('Q', 'P'))
```

---

read.compositional	<i>Read a .csv file with compositional data</i>
--------------------	---

---

## Description

Reads a data table containing compositional data (e.g. chemical concentrations)

## Usage

```
read.compositional(fname, method = NULL, colmap = "rainbow", ...)
```

## Arguments

fname	a string with the path to the .csv file
method	either "bray" (for the Bray-Curtis distance) or "aitchison" (for Aitchison's central logratio distance). If omitted, the function defaults to 'aitchison', unless there are zeros present in the data.
colmap	an optional string with the name of one of R's built-in colour palettes (e.g., heat.colors, terrain.colors, topo.colors, cm.colors), which are to be used for plotting the data.
...	optional arguments to the built-in read.csv function

## Value

an object of class `compositional`, i.e. a list with the following items:

`x`: a data frame with the samples as rows and the categories as columns

`method`: either "aitchison" (for Aitchison's centred logratio distance) or "bray" (for the Bray-Curtis distance)

`colmap`: the colour map provided by the input argument

## Examples

```
fname <- system.file("Major.csv", package="provenance")
Major <- read.compositional(fname)
plot(PCA(Major))
```

---

read.counts	<i>Read a .csv file with point-counting data</i>
-------------	--

---

**Description**

Reads a data table containing point-counting data (e.g. petrographic, heavy mineral, palaeontological or palynological data)

**Usage**

```
read.counts(fname, method = "chisq", colmap = "rainbow", ...)
```

**Arguments**

fname	a string with the path to the .csv file
method	either "chisq" (for the chi-square distance) or "bray" (for the Bray-Curtis distance)
colmap	an optional string with the name of one of R's built-in colour palettes (e.g., heat.colors, terrain.colors, topo.colors, cm.colors), which are to be used for plotting the data.
...	optional arguments to the built-in read.csv function

**Value**

an object of class counts, i.e. a list with the following items:  
 x: a data frame with the samples as rows and the categories as columns  
 colmap: the colour map provided by the input argument

**Examples**

```
fname <- system.file("HM.csv", package="provenance")
Major <- read.counts(fname)
#plot(PCA(HM))
```

---

read.densities	<i>Read a .csv file with mineral and rock densities</i>
----------------	---

---

**Description**

Reads a data table containing densities to be used for hydraulic sorting corrections (minsorting and srd functions)

**Usage**

```
read.densities(fname, ...)
```



**Arguments**

`fname` a string with the path to the .csv file  
`...` optional arguments to the built-in `read.csv` function

**Value**

a vector with mineral and rock densities

**Examples**

```
data(Namib,densities)
N8 <- subset(Namib$HM,select="N8")
distribution <- minsorting(N8,densities,phi=2,sigmaphi=1,medium="air",by=0.05)
plot(distribution)
```

---

<code>read.distributional</code>	<i>Read a .csv file with continuous (detrital zircon) data</i>
----------------------------------	--

---

**Description**

Reads a data table containing continuous data (e.g. detrital zircon ages)

**Usage**

```
read.distributional(fname, errorfile = NA, method = "KS",
  xlab = "age [Ma]", colmap = "rainbow", ...)
```

**Arguments**

`fname` the path of a .csv file with the input data, arranged in columns.

`errorfile` the (optional) path of a .csv file with the standard errors of the input data, arranged by column in the same order as `fname`. Must be specified if the data are to be compared with the Sircombe-Hazelton dissimilarity.

`method` an optional string specifying the dissimilarity measure which should be used for comparing this with other datasets. Should be one of either "KS" (for Kolmogorov-Smirnov) or "SH" (for Sircombe and Hazelton). If `method = "SH"`, then `errorfile` should be specified. If `method = "SH"` and `errorfile` is unspecified, then the program will default back to the Kolmogorov-Smirnov dissimilarity.

`xlab` an optional string specifying the nature and units of the data. This string is used to label kernel density estimates.

`colmap` an optional string with the name of one of R's built-in colour palettes (e.g., `heat.colors`, `terrain.colors`, `topo.colors`, `cm.colors`), which are to be used for plotting the data.

`...` optional arguments to the built-in `read.csv` function

**Value**

an object of class `distributional`, i.e. a list with the following items:

`x`: a named list of vectors containing the numerical data for each sample

`err`: an (optional) named list of vectors containing the standard errors of `x`

`method`: either "KS" (for Kolmogorov-Smirnov), "Kuiper" (for the Kuiper statistic) or "SH" (for Sircombe Hazelton)

`breaks`: a vector with the locations of the histogram bin edges

`xlab`: a string containing the label to be given to the x-axis on all plots

`colmap`: the colour map provided by the input argument

**Examples**

```
agefile <- system.file("DZ.csv",package="provenance")
errfile <- system.file("DZerr.csv",package="provenance")
DZ <- read.distributional(agefile,errfile)
plot(KDE(DZ$x$N1))
```

---

restore

*Undo the effect of hydraulic sorting*

---

**Description**

Restore the detrital composition back to a specified source rock density (SRD)

**Usage**

```
restore(X, dens, target = 2.71)
```

**Arguments**

<code>X</code>	an object of class <code>compositional</code>
<code>dens</code>	a vector of rock and mineral densities
<code>target</code>	the target density (in g/cm <sup>3</sup> )

**Value**

an object of class `SRDcorrected`, i.e. an object of class `compositional` which is a daughter of class `compositional` containing the restored composition, plus one additional member called `restoration`, containing the intermediate steps of the SRD correction algorithm.

**Author(s)**

Alberto Resentini and Pieter Vermeesch

## References

Garzanti E, Ando, S and Vezzoli, G. "Settling equivalence of detrital minerals and grain-size dependence of sediment composition." *Earth and Planetary Science Letters* 273.1 (2008): 138-151.

## See Also

minsorting

## Examples

```
data(Namib,densities)
rescomp <- restore(Namib$PTHM,densities,2.71)
HMcomp <- c("zr","tm","rt","sph","ap","ep","gt",
            "st","amp","cpx","opx")
amcomp <- amalgamate(rescomp,Plag="P",HM=HMcomp,Opq="opaques")
plot(ternary(amcomp),showpath=TRUE)
```

---

SH.diss

*Sircombe and Hazelton distance*


---

## Description

Calculates Sircombe and Hazelton's L2 distance between the Kernel Functional Estimates (KFEs, not to be confused with Kernel Density Estimates!) of two samples with specified analytical uncertainties

## Usage

```
SH.diss(x, i, j, c.con = 0)
```

## Arguments

x	an object of class <code>distributional</code>
i	index of the first sample
j	index of the second sample
c.con	smoothing bandwidth of the kernel functional estimate

## Value

a scalar value expressing the L2 distance between the KFEs of samples i and j

## Author(s)

Keith Sircombe and Martin Hazelton

## References

Sircombe, K. N., and M. L. Hazelton. "Comparison of detrital zircon age distributions by kernel functional estimation." *Sedimentary Geology* 171.1 (2004): 91-111.

## See Also

KS.diss

## Examples

```
datfile <- system.file("DZ.csv",package="provenance")
errfile <- system.file("DZerr.csv",package="provenance")
DZ <- read.distributional(datfile,errfile)
d <- SH.diss(DZ,1,2)
print(d)
```

---

subset	<i>Get a subset of provenance data</i>
--------	--

---

## Description

Return a subset of provenance data according to some specified indices

## Usage

```
## S3 method for class 'distributional'
subset(x, subset = NULL, select = NULL, ...)
```

```
## S3 method for class 'compositional'
subset(x, subset = NULL, components = NULL,
       select = NULL, ...)
```

```
## S3 method for class 'counts'
subset(x, subset = NULL, components = NULL,
       select = NULL, ...)
```

## Arguments

x	an object of class <code>distributional</code>
subset	logical expression indicating elements or rows to keep: missing values are taken as false.
select	a vector of sample names
...	optional arguments for the generic subset function
components	categories to keep

**Value**

an object of class `distributional`

**See Also**

`read.distributional`

**Examples**

```
data(Namib)
coast <- c("N1", "N2", "T8", "T13", "N12", "N13")
ZTRcoast <- subset(Namib$HM, select=coast, components=c('gt', 'cpx', 'ep'))
DZcoast <- subset(Namib$DZ, select=coast)
summaryplot(ZTRcoast, KDEs(DZcoast), ncol=2)
```

---

summaryplot

*Joint plot of several provenance datasets*


---

**Description**

Arranges kernel density estimates and pie charts in a grid format

**Usage**

```
summaryplot(..., ncol = 1, pch = NA)
```

**Arguments**

<code>...</code>	a sequence of datasets of class <code>compositional</code> , <code>KDEs</code> , or <code>distributional</code>
<code>ncol</code>	the number of columns
<code>pch</code>	(optional) symbol to be used to mark the sample points along the x-axis of the <code>KDEs</code> (if appropriate).

**Value**

a summary plot of all the data comprised of `KDEs` for the datasets of class `KDEs`, pie charts for those of class `compositional` and histograms for those of class `distributional`.

**See Also**

`KDEs`

**Examples**

```
data(Namib)
KDEs <- KDEs(Namib$DZ, 0, 3000)
summaryplot(KDEs, Namib$HM, Namib$PT, ncol=2)
```

---

ternary	<i>Define a ternary composition</i>
---------	-------------------------------------

---

## Description

Create an object of class ternary

## Usage

```
ternary(X, x = 1, y = 2, z = 3)
```

## Arguments

X	an object of class compositional OR a matrix or data frame with numerical data
x	string/number or a vector of strings/numbers indicating the variables/indices making up the first subcomposition of the ternary system.
y	second (set of) variables
z	third (set of) variables

## Value

an object of class ternary, i.e. a list containing:

x: a three column matrix (or vector) of ternary compositions.

and (if X is of class SRDcorrected)

restoration: a list of intermediate ternary compositions inherited from the SRD correction

## See Also

restore

## Examples

```
data(Namib)
tern <- ternary(Namib$PT,c('Q'),c('KF','P'),c('Lm','Lv','Ls'))
plot(tern,type="QFL")
```

---

ternary.ellipse	<i>Ternary confidence ellipse</i>
-----------------	-----------------------------------

---

## Description

plot a  $100(1 - \alpha)\%$  confidence region around the data or around its mean.

## Usage

```
ternary.ellipse(x, ...)

## Default S3 method:
ternary.ellipse(x, alpha = 0.05, population = TRUE,
  ...)

## S3 method for class 'compositional'
ternary.ellipse(x, alpha = 0.05,
  population = TRUE, ...)

## S3 method for class 'counts'
ternary.ellipse(x, alpha = 0.05, population = TRUE,
  ...)
```

## Arguments

x	an object of class ternary
...	optional formatting arguments
alpha	cutoff level for the confidence ellipse
population	show the standard deviation of the entire population or the standard error of the mean?

## Examples

```
data(Namib)
tern <- ternary(Namib$Major, 'CaO', 'Na2O', 'K2O')
plot(tern)
ternary.ellipse(tern)
```

---

text.ternary	<i>Ternary text plotting</i>
--------------	------------------------------

---

**Description**

Add text an existing ternary diagram

**Usage**

```
## S3 method for class 'ternary'  
text(x, labels = 1:nrow(x$x), ...)
```

**Arguments**

x	an object of class ternary, or a three-column data frame or matrix
labels	a character vector or expression specifying the text to be written
...	optional arguments to the generic text function

**Examples**

```
data(Namib)  
tern <- ternary(Namib$Major, 'CaO', 'Na2O', 'K2O')  
plot(tern, pch=21, bg='red', labels=NULL)  
# add the geometric mean composition as a text label:  
gmean <- ternary(exp(colMeans(log(tern$x))))  
text(gmean, labels='geometric mean')
```



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