

# The siar Package

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**Type** Package

**Title** Stable Isotope Analysis in R

**Version** 4.0

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**Depends** hrcde, coda, MASS

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**Description** This package takes data on organism isotopes and fits a Bayesian model to their dietary habits based upon a Gaussian likelihood with a mixture dirichlet-distributed prior on the mean. Latest version (4.0) includes the feature to add in concentration dependence. See siardemo() for an example.

**License** GPL ( $\geq 2$ )

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allgroups

*The entire set of Geese isotope data*


---

## Description

A 5 column matrix containing isotopic estimates for 251 geese collected at 8 different time points. The first column indicates the time point group, the second and third are d15N (Nitrogen) and d13C (Carbon) isotopic values for the Geese plasma, the third and fourth are d15N and d13C values for the Geese cells. Note that these are raw values; they have not undergone fractionation correction.

## Usage

```
data(allgroups)
```

## Format

A data frame with 251 observations on the following 5 variables.

**Group** Group number / time point

**d15NP1** d15N plasma

**d13CP1** d13C plasma

**d15NCe** d15N cells

**d13CCe** d13C cells

## Examples

```
#see siarmenu() and option 9 for a demo using part of this data
```

---

concdepdemo*Concentration dependence values for the geese demo data*

---

**Description**

A 5 column, 4 row matrix containing the mean and standard deviation of the concentration dependence values for each of the 2 isotopes used for each different source. Note that the standard deviation is not currently implemented and is set to 0 in this example.

**Usage**

```
data(concdepdemo)
```

**Format**

A data frame with 4 observations on the following 5 variables.

**Sources** The source name.

**Meand15N** The mean concentration dependence value for 15N

**SDd15N** The standard deviation concentration dependence value for 15N

**Meand13C** The mean concentration dependence value for 13C

**SDd13C** The standard deviation concentration dependence value for 13C

**Examples**

```
#see siarmenu() and option 9 for a demo using this data
```

---

correctionsdemo*Fractionation correction values for the geese data*

---

**Description**

A 5 column, 4 row matrix containing the mean and standard deviation of the correction values for each of the 2 isotopes used for each different source

**Usage**

```
data(correctionsdemo)
```

**Format**

A data frame with 4 observations on the following 5 variables.

**Source** The source name.

**Mean15N** The mean correction value for 15N

**Sd15N** The standard deviation correction value for 15N

**Mean13C** The mean correction value for 13C

**Sd13C** The standard deviation correction value for 13C

**Examples**

```
#see siarmenu() and option 9 for a demo using this data
```

---

geese1demo

*A single group of the geese data*

---

**Description**

A 2 column, 9 row matrix containing the plasma data for the first group of geese

**Usage**

```
data(geese1demo)
```

**Format**

A data frame with 9 observations on the following 2 variables.

**d15NP1** d15N plasma

**d13CP1** d13C plasma

**Examples**

```
#see siarmenu() and option 9 for a demo using this data
```

---

`geese2demo`*A multi-group version of the geese plasma data*

---

**Description**

A 3 column, 251 row matrix which contains the isotopic plasma values of 251 geese over 2 isotopes

**Usage**

```
data(geese2demo)
```

**Format**

A data frame with 251 observations on the following 3 variables.

**Group** The group number / time point

**d15NP1** The d15N plasma value

**d13CP1** The d13C plasma value

**Examples**

```
#see siarmenu() and option 9 for a demo using similar data
```

---

`newgraphwindow`*Opens a new graphics window on a variety of platforms*

---

**Description**

Opens a new graphics window on a variety of platforms

**Usage**

```
newgraphwindow()
```

**Details**

Not intended for use outside [siarmenu](#)

**Author(s)**

Andrew Parnell

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panelcontour	Adds contours to a matrix plot
--------------	--------------------------------

---

**Description**

A simple function used by pairs to produce neat looking matrix plots. Not intended for use by those using siar.

**Usage**

```
panelcontour(x, y, ...)
```

**Arguments**

x	A numeric vector containing data with which to produce a contour plot
y	A numeric vector containing data with which to produce a contour plot
...	Other arguments

**Author(s)**

Andrew Parnell

---

---

panelcor	Adds correlations to a matrix plot
----------	------------------------------------

---

**Description**

A simple function used by pairs to produce neat looking matrix plots. Not intended for use by those using siar.

**Usage**

```
panelcor(x, y, digits = 2, prefix = "", cex.cor, ...)
```

**Arguments**

x	A numeric vector containing data with which to produce correlations
y	A numeric vector containing data with which to produce correlations
digits	Number of digits to display on plot
prefix	Text to add before the correlation
cex.cor	Multiplier for the size of the text on the plot
...	Other arguments

**Author(s)**

Unknown

---

panelhist	<i>Adds histograms to the diagonal of a matrix plot</i>
-----------	---

---

**Description**

A simple function used by pairs to produce neat looking matrix plots. Not intended for use by those using siar.

**Usage**

```
panelhist(x, ...)
```

**Arguments**

x	A numeric vector containing data over which to compute a histogram
...	Other arguments

**Author(s)**

Unknown

---

siar-package	<i>Stable Isotope Analysis in R.</i>
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---

**Description**

This package takes data on animal isotopes and fits a Bayesian model to their dietary habits based upon a Gaussian likelihood with a dirichlet prior mixture on the mean. The main function, siarmcmcdirichletv4(), allows the user to specify the data and choose the size of the MCMC run. A wrapper for the package, siarmenu() gives a walkthrough of all the functions contained in the package and produces some pretty plots. Some example data on Geese plasma is included for illustration.

**Details**

Package:	siar
Type:	Package
Version:	3.2
Date:	2008-06-12
License:	GPL (>= 2)

For a demo of how to use the package, type siarmenu() and then choose option 9,

**Author(s)**

Andrew Parnell <Andrew.Parnell@tcd.ie>

**See Also**

[siarmenu](#) [siarmcmcdirichletv4](#)

**Examples**

```
## See siarmenu()
```

---

siaraddcross	<i>Plotting tool for adding isotope bi-plot data to a figure</i>
--------------	--

---

**Description**

A sub-function for siarplotdata() and not intended for calling directly by the user.

**Author(s)**

Andrew Parnell and Andrew Jackson

---

siardemo	<i>Runs the siar model and some nice plots for the siar package</i>
----------	---

---

**Description**

A simple function which utilises the loaded in Geese plasma data to run the MCMC on dietary proportions. Can be accessed either directly or through the menu function

**Usage**

```
siardemo(siarversion = 0)
```

**Arguments**

siarversion    Not required

**Author(s)**

Andrew Parnell



---

siarelicit

*Elicit prior parameters for the Dirichlet distribution.*


---

### Description

Allows users to enter mean estimated proportions and a standard deviation term so that useful prior distributions can be entered into the [siarmcmcdirichletv4](#) function.

### Usage

```
siarelicit(siardata)
```

### Arguments

**siardata** A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the [siarmenu](#) function.

### Details

Uses standard results from the Dirichlet distribution to turn the estimated mean proportions  $M_I$  and a variance term  $V_1$  to give the estimated parameters  $a_i$  via:

$$a_i = M_i \left( \frac{M_1(1 - M_1)}{V_1} - 1 \right)$$

Note that  $V$  can be given for any of the  $k$  sources.

### Author(s)

Andrew Parnell

---

siarhdrs

*Creates hdrs and convergence diagnostics from siar output*


---

### Description

Creates highest density regions and convergence diagnostics from siar output. Accessed by the siar menu function and not really intended for use outside that environment

### Usage

```
siarhdrs(siardata)
```

**Arguments**

`siardata` A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the [siarmenu](#) function.

**Details**

Not intended for use outside [siarmenu](#)

**Author(s)**

Andrew Parnell

---

siarhistograms	<i>Produce neat siar histograms</i>
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---

**Description**

Produces neat and colourful histograms for siar output.

**Usage**

```
siarhistograms(siardata, siarversion)
```

**Arguments**

`siardata` A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the [siarmenu](#) function.

`siarversion` The siar version number as a string.

**Details**

Not intended for use outside [siarmenu](#)

**Author(s)**

Andrew Parnell

---

siarloaddata	<i>Loads in siar data</i>
--------------	---------------------------

---

**Description**

Loads in siar data via a neat menu-driven interface.

**Usage**

```
siarloaddata(siarversion)
```

**Arguments**

siarversion	
siarversion	The siar version number as a string.

**Details**

Not intended for use outside [siarmenu](#)

**Author(s)**

Andrew Parnell

---

siarmatrixplot	<i>Matrix plots of siar output</i>
----------------	------------------------------------

---

**Description**

Produces matrix plots of siar output

**Usage**

```
siarmatrixplot(siardata, siarversion=0)
```

**Arguments**

siardata	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
siarversion	The siar version number.

**Details**

Not intended for use outside [siarmenu](#)

**Author(s)**

Andrew Parnell

---

 siarmcmcdirichletv4

*MCMC for stable isotope data*


---

**Description**

Runs an MCMC on stable isotope data from certain organisms to determine their dietary habits.

**Usage**

```
siarmcmcdirichletv4(data, sources, corrections = 0, concdep = 0, iterations=200000,
```

**Arguments**

data	A matrix with each food source as a separate row and each isotope as a separate column.
sources	A matrix containing the mean and standard deviations of the fractionated correction values for each of the isotopes. Also allows corrections = 0 for pre-corrected data.
corrections	A matrix containing the mean and standard deviations of the fractional correction values for each of the isotopes. Also allows corrections = 0 for pre-corrected data.
concdep	A matrix containing the mean and standard deviations of the concentration dependence values for each of the isotopes. Also allows concdep = 0 for data with no required concentration dependence. Note that version 4.0 does not use the standard deviations.
iterations	The number of iterations to run.
burnin	The size of the burnin
howmany	How often to report the number of iterations.
thinby	The amount of thinning of the iterations.
prior	The dirichlet distribution prior parameters, the default is rep(1,numsources). New parameters can be estimated via the function <a href="#">siarelicit</a> .
siardata	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.

## Details

The model assumes that each target value comes from a Gaussian distribution with an unknown mean and standard deviation. The structure of the mean is a weighted combination of the food sources' isotopic values. The weights are made up dietary proportions (which are given a Dirichlet prior distribution) and the concentration dependencies given for the different food sources. The standard deviation is divided up between the uncertainty around the fractionation corrections (if corrections are given) and the natural variability between target individuals within a defined group (or between all individuals if no grouping structure is specified). The default iterations numbers work well for the demo data sets, but advanced users will want to adjust them to suit their analysis.

## Value

A parameter matrix consisting of (iterations-burnin)/thinby rows with numgroups\*(numsources+numiso) columns, where numsources is the number of food sources, numiso is the number of isotopes, and numgroups is the number of groups. The parameter matrix is structured so that, for each group, the first columns are those of the proportions of each food source eaten, the next columns are the standard deviations for each isotope. This format repeats across rows to each group. The parameters may then subsequently be used for plotting, convergence checks, summaries, etc, etc.

## Author(s)

Andrew Parnell

## See Also

[siarmenu](#), [siarelicit](#)

## Examples

```
# Should take around 10 seconds to run
#out <- siarmcmcdirichletv4(geseldemo, sourcesdemo, correctionsdemo, concdepdemo)
```

---

siarmenu

*A list of menu options for running the siar package*

---

## Description

Brings up a list of menu options which allow the user to run MCMC and produce some plots

## Usage

```
siarmenu()
```

**Details**

The internal workings of this function uses a list called `siardata` containing some or all of the following parts: `targets`, `sources`, `corrections`, `PATH`, `TITLE`, `numgroups`, `numdata`, `numsources`, `numiso`, `SHOULDRUN`, `GRAPHSONLY`, `EXIT`, and `output`. `Targets`, `sources` and `corrections` are the isotopic values, source values and fractionation correction values respectively. `PATH` is the path used to get to the files. `TITLE` is the title to be used on most of the graphs. `Numgroups`, `numdata`, `numsources` and `numiso` are the number of groups, number of data points, number of sources and number of isotopes respectively. `SHOULDRUN`, `GRAPHSONLY` and `EXIT` are used to determine which parts of the menu system can be accessed.

**Author(s)**

Andrew Parnell

**See Also**

[siarmcmcdirichletv4](#)

**Examples**

```
#siarmenu()
```

---

`siarmultigrouprun`    *siar MCMC for multi-group data*

---

**Description**

Runs the siar MCMC with Dirichlet mixture mean for isotopic data

**Usage**

```
siarmultigrouprun(siardata)
```

**Arguments**

<code>siardata</code>	A list containing some or all of the following parts: <code>targets</code> , <code>sources</code> , <code>corrections</code> , <code>PATH</code> , <code>TITLE</code> , <code>numgroups</code> , <code>numdata</code> , <code>numsources</code> , <code>numiso</code> , <code>SHOULDRUN</code> , <code>GRAPHSONLY</code> , <code>EXIT</code> , and <code>output</code> . For more details of these inputs see the <a href="#">siarmenu</a> function.
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**Details**

Not intended for use outside [siarmenu](#)

**Author(s)**

Andrew Parnell

---

siarplotdata	<i>Produces plots of target data and sources</i>
--------------	--

---

## Description

Produces colourful scatter plots of siar target data and sources.

## Usage

```
siarplotdata(siardata, siarversion = 0, grp=1:siardata$numgroups, panel=NULL, isos=c
```

## Arguments

siardata	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
siarversion	The siar version number as a string.
grp	A vector containing the groups of target consumer data to be rendered on the graph. Default value NULL draws all groups. Groups are identified by their own data marker.
panel	A scalar value that determines if the groups of consumer data are to be drawn on the same graph (default=NULL) or on separate panels within a single figure. Number of rows and columns of panels can be specified by a 2 element vector. Alternatively, giving a single value e.g. panel=1 will cause the program to attempt to fit a "reasonable" number of panels to each row and column.
isos	A two element vector containing the reference to each isotope combination for the x and y axis to be rendered in the figure. Note, only relevant for datasets containing >2 isotopes. By default, if there are more than two isotopes, separate figures will be created for all possible combinations of isotopes.
leg	A scalar determining how the legend is to be created. Default leg=1 prompts the user to locate the legend on each figure. leg = 2, puts the legend in a new figure automatically (useful if you want to omit the legend but still want to retain access to the information). leg = 0 omits the legend entirely.

## Details

Can be called at any time after running [siarloaddata](#) or when running [siarmenu](#)

## Author(s)

Andrew Parnell and Andrew Jackson

---

```
siarplotdatawrapper
```

*Handles repeated plotting instructions for siarplotdata()*

---

### Description

A sub-function for siarplotdata() and not intended for calling directly by the user.

### Author(s)

Andrew Parnell and Andrew Jackson

---

```
siarplottarget
```

*Plots the consumers' data in isotope space*

---

### Description

A sub-function for siarplotdata() and not intended for calling directly by the user.

### Author(s)

Andrew Parnell and Andrew Jackson

---

```
siarproportionbygroupplot
```

*siar proportion plots by group*

---

### Description

Plots boxplots or line plots representing defined credible intervals for each source (x-axis) for a given group. The representation is basically the same as siarhistograms but allows easier comparison of source contribution within a group. Similar in style to siarproportionbysourceplot().

### Usage

```
siarproportionbygroupplot(siardata, siarversion=0, probs=c(95, 75, 50), xlabels=NULL, g
```



**Arguments**

<code>siardata</code>	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
<code>siarversion</code>	The siar version number.
<code>probs</code>	Define the extent probability intervals for a given parameter.
<code>xlabels</code>	Specifies the text to associate with each group defined as ticks on the x-axis.
<code>grp</code>	Specifies which group the graph is drawn for. Default prompts the user for input from the command line.
<code>type</code>	Determines the style of graph. <code>type="boxes"</code> draws boxplot style (default), <code>type="lines"</code> draws overlain lines increasing in thickness
<code>clr</code>	Determines the set of colours to use for the boxes. Default is greyscale.
<code>scl</code>	Specifies a proportional scaling factor to increase ( $scl > 1$ ) or decrease ( $scl < 1$ ) the default width of lines or boxes. Default = 1.
<code>xspc</code>	Sets the amount of blank space either side of the first and last (on the x-axis) graphic object.
<code>prn</code>	If <code>prn=TRUE</code> the values for the defined probability densities ( <code>probs</code> ) are returned to the command window. Default is <code>prn=FALSE</code> with no such output.
<code>leg</code>	Determines whether a legend is to be drawn ( <code>leg=TRUE</code> ) or not (default <code>leg=FALSE</code> ). Note, currently only supported for <code>type="lines"</code> .

**Author(s)**

Andrew Jackson & Andrew Parnell

---

```
siarproportionbysourceplot
      siar proportion plots by source
```

---

**Description**

Useful for siar data with multiple groups where the variability of each source over time is of interest.

**Usage**

```
siarproportionbysourceplot(siardata, siarversion=0, probs=c(95, 75, 50), xlabels=NULL,
```

**Arguments**

siardata	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
siarversion	The siar version number.
probs	Define the extent probability intervals for a given parameter.
xlabels	Specifies the text to associate with each group defined as ticks on the x-axis.
grp	Specifies which source group the graph is drawn for. Default prompts the user for input from the command line.
type	Determines the style of graph. type="boxes" draws boxplot style (default), type="lines" draws overlain lines increasing in thickness
clr	Determines the set of colours to use for the boxes. Default is greyscale.
scl	Specifies a proportional scaling factor to increase (scl > 1) or decrease (scl < 1) the default width of lines or boxes. Default = 1.
xspc	Sets the amount of blank space either side of the first and last (on the x-axis) graphic object.
prn	If prn=TRUE the values for the defined probability densities (probs) are returned to the command window. Default is prn=FALSE with no such output.
leg	Determines whether a legend is to be drawn (leg=TRUE) or not (default leg=FALSE). Note, currently only supported for type="lines".

**Author(s)**

Andrew Parnell

---

siarsaveoutput	<i>Saves siar output to a file</i>
----------------	------------------------------------

---

**Description**

Saves created siar output to a file

**Usage**

```
siarsaveoutput (siardata)
```

**Arguments**

siardata	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
----------	---

**Details**

Not intended for use outside [siarmenu](#)

---

`siarsinglegrouprun` *siar MCMC for single group data*

---

**Description**

Runs the siar MCMC with Dirichlet mixture mean for isotopic data

**Usage**

```
siarsinglegrouprun(siardata)
```

**Arguments**

<code>siardata</code>	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
-----------------------	---

**Details**

Not intended for use outside [siarmenu](#)

**Author(s)**

Andrew Parnell

---

`siarsolomcmcv4` *MCMC for stable isotope data with only single target organisms*

---

**Description**

Runs an MCMC on stable isotope data from certain organisms to determine their dietary habits. This version requires only a single target organism per group

**Usage**

```
siarsolomcmcv4(data, sources, corrections = 0, concdep = 0, iterations=200000, burn
```

## Arguments

<code>data</code>	A matrix with each food source as a separate row and each isotope as a separate column.
<code>sources</code>	A matrix containing the mean and standard deviations of the fractionated correction values for each of the isotopes. Also allows corrections = 0 for pre-corrected data.
<code>corrections</code>	A matrix containing the mean and standard deviations of the fractional correction values for each of the isotopes. Also allows corrections = 0 for pre-corrected data.
<code>concdep</code>	A matrix containing the mean and standard deviations of the concentration dependence values for each of the isotopes. Also allows concdep = 0 for data with no required concentration dependence. Note that version 4.0 does not use the standard deviations.
<code>iterations</code>	The number of iterations to run.
<code>burnin</code>	The size of the burnin
<code>howmany</code>	How often to report the number of iterations.
<code>thinby</code>	The amount of thinning of the iterations.
<code>prior</code>	The dirichlet distribution prior parameters, the default is <code>rep(1, numsources)</code> . New parameters can be estimated via the function <a href="#">siarelicit</a> .
<code>siardata</code>	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.

## Details

The model assumes that each target value comes from a Gaussian distribution with an unknown mean and standard deviation. The structure of the mean is a weighted combination of the food sources' isotopic values. The weights are made up dietary proportions (which are given a Dirichlet prior distribution) and the concentration dependencies given for the different food sources. The standard deviation is divided up between the uncertainty around the fractionation corrections (if corrections are given) and the natural variability between target individuals within a defined group (or between all individuals if no grouping structure is specified). The default iterations numbers work well for the demo data sets, but advanced users will want to adjust them to suit their analysis.

Note that this version is analogous to the Moore and Semmens (2008) MixSIR model except with a Dirichlet prior distribution.

## Value

A parameter matrix consisting of  $(\text{iterations} - \text{burnin}) / \text{thinby}$  rows with  $\text{numgroups} * (\text{numsources} + \text{numiso})$  columns, where `numsources` is the number of food sources, `numiso` is the number of isotopes, and `numgroups` is the number of groups. The parameter matrix is structured so that, for each group, the first columns are those of the proportions of each food source eaten, the next columns are the standard deviations for each isotope. This format repeats across rows to each group. The parameters may then subsequently be used for plotting, convergence checks, summaries, etc, etc.

**Author(s)**

Andrew Parnell

**References**

Moore and Semmens (2008), Incorporating uncertainty and prior information into stable isotope mixing models, Ecology Letters.

**See Also**

[siarmenu](#), [siarelicit](#)

---

siarsolomultigrouprun

*siar MCMC for multi-group data with only one target organism per group*

---

**Description**

Runs the siar MCMC with Dirichlet mixture mean for isotopic data with only one target organism per group.

**Usage**

```
siarsolomultigrouprun(siardata)
```

**Arguments**

siardata	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
----------	---

**Author(s)**

Andrew Parnell

---

```
siarsolosinglegroupun
```

*siar MCMC for single group data with only one organism*

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### Description

Runs the siar MCMC with Dirichlet mixture mean for isotopic data when there is only one target organism.

### Usage

```
siarsolosinglegroupun(siardata)
```

### Arguments

siardata	A list containing some or all of the following parts: targets, sources, corrections, PATH, TITLE, numgroups, numdata, numsources, numiso, SHOULD RUN, GRAPH-ONLY, EXIT, and output. For more details of these inputs see the <a href="#">siarmenu</a> function.
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### Details

Not intended for use outside [siarmenu](#)

### Author(s)

Andrew Parnell

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sourcesdemo	<i>Source (in this case plant) isotope values</i>
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### Description

A 3 column, 4 row matrix containing 4 different plants and their measurements on 2 different isotopes

### Usage

```
data(sourcesdemo)
```

**Format**

A data frame with 5 observations on the following 3 variables.

**Sources** The plants name

**Meand15N** d15N mean

**SDd15N** d15N standard deviation

**Meand13C** d13C mean

**SDd13C** d13C standard deviation

**Examples**

```
#see siarmenu() and option 9 for a demo using this data
```

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