

# Package ‘PBSawatea’

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**Title** Tools for Running Awatea and Visualizing the Results

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**Depends** R (>= 2.12.0), methods, PBSmodelling, scape, scapeMCMC, lattice, xtable, gdata, gplots

**Suggests** PBStools, Hmisc, coda, grid

**Description** Provides tools for running population models  
using Awatea, which is a variant of Coleraine modified  
by Allan Hicks and Paul Starr for marine fish stocks.

**License** GPL (>= 2)

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allEqual	<i>Are All Values Equal to the First?</i>
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**Description**

A short-cut function for `all(x==x[1])`, which asks are all values in *x* equal to the first value in *x*.

**Usage**

`allEqual(x)`

**Arguments**

`x`                      vector of values.

**Value**

TRUE or FALSE

**See Also**

[all](#), [clearAll](#), [clipVector](#)

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AWATEAdata-class	<i>S4: AWATEA Data Class</i>
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**Description**

The set of functions described here provide an easy method of dealing with a complicated scheme for manipulating data when using the catch-at-age model called **Awatea**, a variant of **Coleraine**. Awatea is maintained by Allan Hicks (University of Washington) and Paul Starr (Canadian Ground-fish Research and Conservation Society).

To use **PBSawatea**'s management functions, the function `runADMB` calls `readAD` which creates an `AWATEAdata` object. The `AWATEAdata` object contains a distinct R environment where data are stored and accessed by various package functions.

**Details**

An Awatea data file can be loaded into an `AWATEAdata` object using `readAD`, and if a complementary results file is available, the user can run `reweight` for a single re-weighting of the abundance data and the composition data.

Alternatively, starting with a single input data file, a user can run the function `runADMB` to perform multiple (iterative) re-weightings automatically.

## Objects from the Class

This S4 object can be created by calls of the form:

```
dat = new("AWATEAdata", txtnam=character(), input=character(), vlst=list(),
dnam=character(), nvars=numeric(), vdesc=character(), vars=list(),
gcomm=character(), vcomm=character(), resdat=list() , likdat=list(),
pardat=list(), stddat=list(), cordat=list(), evadat=list(), reweight=list())
```

The function `readAD` populates the first nine slots of the S4 object by reading in an Awatea data file from the user's current working directory. If Awatea output files exist (with the same prefix as the input file and suffixes = { `.res`, `.lik`, `.par`, `.std`, `.cor`, `.eva`}), then slots `resdat`, `likdat`, `pardat`, `stddat`, `cordat`, and `evadat`, respectively, will also be populated.

## Slots

`txtnam` the file name (including extension) of the Awatea input data file.  
`input` a string vector comprising the line-by-line information from the data file.  
`vlst` a list named by line number (e.g. "L001: . . .") of comments and data.  
`dnam` a vector of data descriptors referenced by numeric line number.  
`nvars` a numeric scalar indicating the number of input variables.  
`vdesc` as for `dnam` but indexed by variable identifier (e.g. "v001").  
`vars` variable values indexed by variable identifier.  
`gcomm` commented lines indexed by numeric line number.  
`vcomm` value descriptors indexed by numeric line number.  
`resdat` list of Awatea results from the file `results.dat` after calling `importRes`.  
`likdat` list of MPD likelihoods from the file `likelihood.dat` after calling `importLik`.  
`pardat` list of parameter values from the file `Awatea.par` after calling `importPar`.  
`stddat` list of estimated parameter values from the file `Awatea.std` after calling `importStd`.  
`cordat` list of correlation objects from the file `Awatea.cor` after calling `importCor`.  
`evadat` eigenvalues of the Hessian from the file `Awatea.eva` after calling `importEva`.  
`reweight` list of re-weighted abundance and composition data (see [reweight](#) for details).

## Methods

`fix`, signature(`x`="AWATEAdata") : replace data elements of an input list  
`reweight`, signature(`x`="AWATEAdata") : reweight abundance and composition data  
`view`, signature(`x`="AWATEAdata") : view the basic input list  
`write`, signature(`x`="AWATEAdata") : write a new input data file

## Note

Some of the output values in `results.dat`, `likelihood.dat`, `Awatea.par`, `Awatea.std`, and `Awatea.cor` contain redundant information.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[readAD](#) for loading the ADMB data file  
[importRes](#) for reading in the file results.dat  
[importLik](#) for reading in the file likelihood.dat  
[importPar](#) for reading in the file Awatea.par  
[importStd](#) for reading in the file Awatea.std  
[importCor](#) for reading in the file Awatea.cor  
[importEva](#) for reading in the file Awatea.eva  
[reweight](#) for re-weighting abundance and composition data  
[runADMB](#) for running the Awatea model through multiple iterations.

---

calc.projExpect

---

*Calculate Expectations and Probabilities*


---

**Description**

Calculate the expectation of projection to reference, and probability of being greater than reference.

**Usage**

```
calc.projExpect ( obj, projObj, refYrs )
calc.projExpect2( obj, projObj, refList )
calc.projProbs  ( obj, projObj, refYrs )
calc.projProbs2 ( obj, projObj, refList )
calc.refProbs   ( projObj=currentProj$B, refPlist=refPointsList )
```

**Arguments**

obj	matrix of biomass MCMCs.
projObj	matrix of biomass projections.
refYrs	numeric vector of reference years
refList	list of reference years (numeric vectors).
refPlist	list of reference points.

**Details**

calc.projExpect... Calculate the expectation of projection to reference.  
 ..... Compare reference years to projection years.  
 calc.projExpect2... Calculate expectation (projection biomass / reference biomass).  
 calc.projProbs... Calculate the probability of being greater than refYrs.  
 ..... Compare reference years to projection years.  
 calc.projProbs2... Calculate the probability of being greater than refYrs.  
 ..... Compare reference years to projection years.  
 calc.refProbs... Calculate the reference probabilities (based on calc.projProbs2).

**Value**

Decision tables

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[calc.refVal](#)

---

calc.refVal

*Calculate Reference Value for Performance Measure*

---

**Description**

Calculate the reference value for performance measures.

**Usage**

```
calc.refVal(obj, refYrs, fun=mean)
```

**Arguments**

obj	scape Biomass matrix with $n$ rows and $m$ columns, where $n$ = number of MCMC samples, and $m$ = number of years.
refYrs	numeric years in reference period.
fun	the function to apply to reference period $i$ .

**Value**

Returns a vector of length `nrow(obj)` reference values.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

See Also

[calc.projExpect](#), [findTarget](#)

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closeAllWin	<i>Close All Open Devices</i>
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Description

Close all open devices.

Usage

```
closeAllWin()
```

See Also

[closeWin](#)

---

compB0	<i>Compare Reference Criteria and Points Relative to B0</i>
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---

Description

Compare COSWEIC reference criteria and DFO reference points relative to  $B_0$ . The figure concept comes from Chris Woods (PBS).

Usage

```
compB0(B, Mnames=NULL, ratios=c(0.4,0.8),
  include=list(A1=TRUE, A2=TRUE, SSPM=TRUE, Bmsy=TRUE, Bt=TRUE),
  t.yr=2011, boxwidth=0.6, figgy=FALSE, width=12, height=9, ...)
```

Arguments

B	list of list of MCMC samples (see <b>Details</b> ); the first level of the list is the model run, while the second level contains MCMC samples (one of which should be $B_0$ which acts as the divisor to the other MCMCs).
Mnames	optional model names for the boxplot.
ratios	reference levels of $B_{MSY}$ (usually 0.4 and 0.8).
include	list of logicals specifying whether to plot COSEWIC's reference criteria A1 and A2, the Schaefer surplus production model SSPM, the biomass at maximum sustainable yield Bmsy, and the biomass of a given time period Bt, where t.yr specifies the year. All values are cast in terms of $B_0$ .
t.yr	numeric year that represents the time period for $B_t$ .

boxwidth	width of the box in x-units.
figgy	logical: if TRUE, send figure to four output files (.eps, .pdf, .png, and .wmf).
width	width of the output files in inches.
height	height of the output files in inches.
...	additional values for <code>plotBox</code> 's argument pars.

### Details

An example of the input list B:

```
List of 2
..$ 29.01:List of 3
....$ B0.MCMC : num [1:1000]
....$ Bt.MCMC : num [1:1000]
....$ Bmsy.MCMC: num [1:1000]
..$ 30.01:List of 3
....$ B0.MCMC : num [1:1000]
....$ Bt.MCMC : num [1:1000]
....$ Bmsy.MCMC: num [1:1000]
```

The function creates a figure comparing COSEWIC criteria and reference points for the Schaefer surplus production model and specified ratios of  $B_{MSY}$  from catch-at-age models. The coordinate space (y-axis) is relative to  $B_0$ .

### Value

Invisibly returns a list object of `xBox` and `BarBox` used to create the boxplot.

### Note

Uses a modified version of boxplot called `plotBox`.

### Author(s)

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

### See Also

`plotBox`, `importMCMC`, `msyCalc`  
`plotmath`, `boxplot`, `bxp`



compBmsy

*Compare Biomass Posteriors Relative to Bmsy***Description**

Compare posterior distributions of biomass from MCMCs for species and model runs.

**Usage**

```
compBmsy(Bspp, spp="POP", Mnams=c("Est M", "Fix M"),
         ratios=c(0.4, 0.8), t.yr=2011, figgy=FALSE, width=12, height=9, ...)
```

**Arguments**

Bspp	list of species MCMC results, which are lists of model runs, each listing data frames of $B_t$ and $B_{MSY}$ .
spp	species code(s) (e.g., 3-letter), which are the first-level names of Bspp.
Mnams	optional model names for the boxplots.
ratios	reference levels of $B_{MSY}$ (usually 0.4 and 0.8).
t.yr	numeric year that represents the time period for $B_t$ .
figgy	logical: if TRUE, send figure to four output files (.eps, .pdf, .png, and .wmf).
width	width of the output files in inches.
height	height of the output files in inches.
...	additional values for <a href="#">plotBox</a> 's argument pars.

**Details**

An example of the input list Bspp:

```
List of 2
..$ POP:List of 2
....$ run23:List of 3
.....$ B0.MCMC : num [1:1000]
.....$ Bt.MCMC : num [1:1000]
.....$ Bmsy.MCMC: num [1:1000]
....$ run16:List of 3
.....$ B0.MCMC : num [1:1000]
.....$ Bt.MCMC : num [1:1000]
.....$ Bmsy.MCMC: num [1:1000]
....- attr(*, "spp")= chr "POP"
..$ YMR:List of 2
....$ 29.01:List of 3
.....$ B0.MCMC : num [1:1000]
.....$ Bt.MCMC : num [1:1000]
.....$ Bmsy.MCMC: num [1:1000]
....$ 30.01:List of 3
```

```
.....$ B0.MCMC : num [1:1000]
.....$ Bt.MCMC  : num [1:1000]
.....$ Bmsy.MCMC: num [1:1000]
....- attr(*, "spp")= chr "YMR"
```

The function creates a set of horizontal boxes delimited by the quantiles (0.025, 0.25, 0.5, 0.75, 0.975) that illustrate the posterior MCMC samples of biomass relative to  $B_{MSY}$ . The default reference points ( $0.4B_{MSY}$  and  $0.8B_{MSY}$ ) are shown by vertical dashed lines.

**Value**

Invisibly returns the boxplot list object Bmsy.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC.

**See Also**

[compB0](#), [plotBox](#)

---

cquantile	<i>Running Quantile</i>
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---

**Description**

Creates a set of running quantiles from MCMC traces.  
(Uses subfunction found in **coda**'s function cumuplot.)

**Usage**

```
cquantile(z, probs)
cquantile.vec(z, prob)
```

**Arguments**

- z                    an MCMC object.
- probs                vector of quantiles.
- prob                single quantile.

**Value**

cquantile.....running quantile matrix  
cquantile.vec...running quantile vector

**Note**

Arni Magnusson describes a running quantile as:  
*“the evolution of the sample quantiles as a function of the number of iterations”*

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[cumuplot](#), [plotTracePOP](#)

---

findTarget

*Find Time to Achieve a Target Reference Point*


---

**Description**

Find the time (years) to achieve a recovery target (including a moving target) with a given confidence. Produce decision tables showing the probability of exceeding the reference point.

**Usage**

```
findTarget(Vmat, yrU=as.numeric(dimnames(Vmat)[[2]]), yrG=90,
           ratio=0.5, target=B0.MCMC, conf=0.95, plotit=FALSE, retVal="N")
```

**Arguments**

Vmat	matrix of projected biomass values $B_{Nt}$ , where $N$ = number of MCMCs and $t$ = projection year.
yrU	user-specified projection years.
yrG	number of years $G$ for a moving target window ( <i>e.g.</i> , 3 YMR generations = 90y); might not work for all possibilities.
ratio	recovery target ratio $R$ .
target	recovery target values $T_N$ = B0.MCMC for ratios of $B_0$ ; = Bmsy.MCMC for ratios of $B_{MSY}$ ; = Bt.MCMC for moving window of $B_{N,t-G}$ .
conf	confidence level $C$ required.
plotit	logical: if TRUE, plot the probability $p_t$ of exceeding target reference point.
retVal	character name of object to return: retVal="N" : creates global object "Ttab" (see below); retVal="p.hi" : creates global object "Ptab" (see below).

### Details

As this function uses Bayesian output, there are  $N$  (e.g., 1000) values of some target  $T_N$ , which can remain fixed ( $B_0$ ,  $B_{MSY}$ ) or move forward in time  $G$  years before the projection year  $t$  (that is  $T_{N,t-G}$ ). For simplification, we'll just call all targets  $T_N$ .

The probability of exceeding a target ratio  $R$  is:

$$p_t = \frac{1}{N} \sum^N \left[ \frac{B_{Nt}}{T_N} \geq R \right],$$

where  $R$  = target ratio of the reference point (e.g.,  $0.4B_{MSY}$  ( $R=0.4$ ),  $0.2B_0$  ( $R=0.2$ ),  $0.5B_{t-G}$  ( $R=0.5$ )).

At a glance, we can see for any given projection year  $t$  whether the probability of achieving a target ratio is greater than the confidence required:

$$p_t \geq C,$$

where  $C$  is the confidence level acceptable.

### Value

If `retVal="N"` then the function returns a data frame object called "Ttab" in the user's global environment. This table reports the number of years to achieve the target reference point at various catch levels with a specified confidence.

If `retVal="p.hi"` then the function returns a list object called "Ptab" in the user's global environment. This list contains data frames (tables) that report the probability of achieving various reference points at specified catch levels.

Any other `retVal` will return a list of the specified object, if it exists in the function.

### Author(s)

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

### See Also

[runSweaveMCMC](#)

---

get.resFile

*Get Awatea Results Files for Menu*

---

### Description

A function that retrieves the names of Awatea results files (`.res$`) for use in the `mainMenu` command. When choice is made, the function loads the results file and assigns it to the global environment as `currentRes`.

**Usage**

```
get.resFile(resFile=NULL)
```

**Arguments**

resFile	supposedly the name of a results file, but the code suggests that argument is ignored.
---------	--

**Value**

A results file chosen from a menu.

**Note**

AME: made changes so that options are compatible with those in `load.allResFiles`. Previously, trouble occurred when overwriting.

**See Also**

[mainMenu](#), [importCol2](#)

---

getYrIdx

*Select a Subset of Years for Plotting*

---

**Description**

Select a subset of years for which many years are available. The default is to select 5-year increments.

**Usage**

```
getYrIdx(yrNames, mod=5)
```

**Arguments**

yrNames	vector (character or numeric) of years.
mod	select the years modulo mod.

**Value**

Subset of input years that are modulo mod.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[findPat](#), [pad0](#)

---

gfcode

---

*Code: Groundfish Species Codes and Names*


---

### Description

Data frame of groundfish species described primarily by Hart codes. The information resides in a data table called C\_Species in the relational database PacHarvest.

### Usage

```
data(gfcode)
```

### Format

Data frame comprising 549 rows (species) and 5 columns:

code	Hart code
latin	scientific (latin) name
name	common species name
code2	2-letter code if available
code3	3-letter code if available

### Details

The data frame comprises 5,49 rows (species) and 5 columns. Hart codes for fish species essentially reference page numbers in Hart (1973).

Quick reference to species name: `species["424",]$name` yields Quillback rockfish.

Quick reference to latin name: `species["424",]$latin` yields *Sebastes maliger*.

### Source

Norm Olsen, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

### References

Hart, J.L. (1973) Pacific Fishes of Canada. *Bulletin of the Fisheries Research Board of Canada* **180**, 740 pp.

---

graphics

---

*Open a Windows Device*


---

### Description

Open a windows device in portrait or landscape mode.

**Usage**

```
graphics(view = "portrait")
```

**Arguments**

view            if "portrait", set width = 8.5 in and height = 11 in  
                  if "landscape", set width = 11 in and height = 8.5 in

**See Also**

[resetGraph](#)

---

importCol2

---

*Import Coleraine Model Results (AME version)*


---

**Description**

Import Coleraine model results from .res file, and rearrange into a standard format suitable for plotting.

**Usage**

```
importCol2(res.file, info="", Dev=FALSE, CPUE=FALSE, Survey=FALSE,
           CAc=FALSE, CAS=FALSE, CLc=FALSE, CLs=FALSE, LA=FALSE,
           quiet=TRUE, extra=TRUE)
```

**Arguments**

res.file	name of Coleraine model results file to import.
info	optional string containing information to store with model results.
Dev	logical: whether recruitment deviates were estimated in model.
CPUE	logical: whether model was fitted to catch-per-unit-effort data.
Survey	logical: whether model was fitted to survey abundance index data.
CAC	logical: whether model was fitted to commercial catch-at-age data.
CAS	logical: whether model was fitted to survey catch-at-age data.
CLc	logical: whether model was fitted to commercial catch-at-length data.
CLs	logical: whether model was fitted to survey catch-at-length data.
LA	logical: whether model was fitted to length-at-age data.
quiet	logical: whether to report progress while parsing file.
extra	logical: if TRUE, import likelihoods, parameters, priors, and recruitment residuals.

## Details

This function was modified from the original `importCol` function in the **scape** package to grab extra data.

## Value

A list of class `scape` containing at least `N`, `B`, and `Se1`. The other elements may or may not be included in the list, depending on how `importRes` was called:

<code>N</code>	predicted numbers at age
<code>B</code>	predicted biomass, recruitment, and observed landings (year things)
<code>Se1</code>	predicted selectivity and observed maturity (age things)
<code>Dev</code>	predicted recruitment deviates from the stock-recruitment curve
<code>CPUE</code> , <code>Survey</code>	commercial and survey abundance index and fit
<code>CAC</code> , <code>CAS</code>	commercial and survey C@A (catch at age) and fit
<code>CLC</code> , <code>CLS</code>	commercial and survey C@L (catch at length) and fit
<code>LA</code>	observed L@A and fit

## Note

This `import` function is implemented for the Coleraine statistical catch-at-age software, and can serve as a template for **scape** users who would like to implement `import` functions for specific stock assessment software.

The functions `ll` (package **gdata**) and `head` are recommended for browsing model results, e.g. `ll(x.cod)`; `ll(x.cod$N)`; `head(x.cod$N)`.

## References

Hilborn, R., Maunder, M., Parma, A., Ernst, B., Payne, J., and Starr, P. (2003) Coleraine: A generalized age-structured stock assessment model. User's manual version 2.0. *University of Washington Report SAFS-UW-0116*. Available at:  
<http://fish.washington.edu/research/coleraine/pdf/coleraine.pdf>.

## See Also

`importRes`, `read.table`, `readLines`, and `scan` to import any data.  
[scape-package](#) gives an overview of the package **scape**.



---

`importCor`*Import Awatea Correlation File*

---

**Description**

Import an Awatea correlation file that results from an MPD minimisation.

**Usage**

```
importCor(cor.file)
```

**Arguments**

`cor.file`            correlation file (e.g., `Awatea.cor`).

**Details**

Extracts the data from character vectors and makes various data objects.

**Value**

The output is a list object with the following components:

<code>cfile</code>	character vector representing the file line-by-line
<code>cor</code>	data.frame of the correlation file
<code>cor.mat</code>	matrix representing the correlation matrix only
<code>index</code>	character vector used in the row and column names of <code>cor</code> and <code>cor.mat</code>
<code>cor.name</code>	Awatea parameter names in the correlation file
<code>cor.value</code>	Awatea parameter values in the correlation file
<code>cor.std.dev</code>	Awatea parameter standard deviations in the correlation file
<code>hessian_log_determinant</code>	log of the determinant of the hessian reported in the header

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[importPar](#), [importRes](#), [importStd](#)

---

importEva	<i>Import Awatea Hessian Eigenvalues</i>
-----------	--

---

**Description**

Import a vector of eigenvalues of the Hessian from an MPD analysis file (Awatea.eva).

**Usage**

```
importEva(eva.file)
```

**Arguments**

eva.file            vector of hessian eigenvalues; first element should be a positive definite number.

**Details**

The eigenvalues of the Hessian represent the 2nd derivatives of the negative log-likelihood function. If they are all positive it indicates a minimum. The first element of this vector should be positive definite before an MCMC is run.

**Value**

The output is a list object with only one component:

eva                a numeric vector representing the eigenvalues of the Hessian.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[importCor](#), [importLik](#), [importPar](#), [importRes](#), [importStd](#)

---

importLik	<i>Import Awatea Likelihoods</i>
-----------	----------------------------------

---

**Description**

Import an Awatea file of the likelihoods from an MPD estimation.

**Usage**

```
importLik(lik.file)
```

**Arguments**

lik.file                likelihoods file (usually likelihood.dat)

**Details**

Extracts the data from character vectors and makes various data objects.

**Value**

The output is a list object with the following components:

lik	character vector representing the file line-by-line
Total_likelihood	values of the final likelihood (objective function value)
Worst_gradient	worst gradient?
CPUE	commercial CPUE likelihood(s)
Survey_Index	survey index likelihood(s)
CA_Commercial	commercial catch-at-age likelihood(s)
CL_Commercial	commercial catch-at-length likelihood(s)
CA_survey	survey catch-at-age likelihood(s)
CL_no_sex_data_survey	likelihood(s) for catch-at-length from surveys with no sex data?
CL_data_survey	likelihood(s) for catch-at-length from surveys with sex data?
Von_B_Likelihood	von Bertalanffy likelihood
Priors	priors?
Penalty_for_U	penalty for U?

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[importCor](#), [importPar](#), [importRes](#), [importStd](#)

---

`importMCMC.ddiff`*Import Functions for PJS Delay Difference Model*

---

### Description

Make a **scapeMCMC** object identical in format to the result of `importMCMC` (or `importProj`) from PJS delay difference model output.

The difference is that  $B$  is biomass defined by a delay difference model.

### Usage

```
importMCMC.ddiff()  
importProj.ddiff(yrVal="2006")
```

### Arguments

`yrVal` character year for delay-difference model (?).

### Value

`importMCMC.ddiff` returns a list object containing:

- L...likelihood MCMCs,
- P...parameter MCMCs,
- B...spawning biomass MCMCs,
- R...recruitment MCMCs.

`importProj.ddiff` returns a list object containing:

- B...projected biomass,
- Y...projected yields.

### Note

Get the biomass projection – PJS does 1 year ahead projection. The column "X" appears as the last column because trailing ", " exist in the `mcmcprojbiom.csv` file.

Note also that "cat=" in .csv file becomes "cat." in `read.table`.

### Author(s)

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

### See Also

[importMCMC](#), [importProj](#), [read.table](#)

---

importPar	<i>Import Awatea Parameter File</i>
-----------	-------------------------------------

---

**Description**

Import an Awatea file of all parameters resulting from an MPD minimisation.

**Usage**

```
importPar(par.file)
```

**Arguments**

par.file	parameter file (e.g., Awatea.par).
----------	------------------------------------

**Details**

Extracts the data from character vectors and makes various data objects.

**Value**

The output is a list object with the following components:

par	character vector representing the file line-by-line
npars	number of parameters reported in header
fval	objective function value reported in header
maxgrad	maximum gradient component reported in header
parameters	numerous reported parameters (e.g., R0) that will depend on the model

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[importCor](#), [importRes](#), [importStd](#)

---

importProjRec	<i>Import Projected Recruitment</i>
---------------	-------------------------------------

---

### Description

Import the projected recruitments (actually, the values are random normals  $N(0, 1)$ ).

### Usage

```
importProjRec(dir, info="", coda=FALSE, quiet=TRUE)
```

### Arguments

dir	directory where MCMC projections reside.
info	user-supplied information, if desired.
coda	logical: if TRUE, use the function <code>mcmc</code> in the package <b>coda</b> to generate MCMCs.
quiet	logical: if TRUE, print progress messages to the R console.

### Details

The values saved by the Awatea code are random normals  $N(0, 1)$ , which for a particular MCMC sample are the same for all the catch strategies.

### Value

A list object comprising:

B	data frame of spawning biomass (dim = MCMC samples by projected years)
Y	data frame of yield (dim = MCMC samples by projected years)
eps	data frame of $\epsilon_t$ (dim = MCMC samples by projected years)

### Note

The function `importProj` does not import recruitment residuals.

This function grabs the `tempdev` values from Awatea, which are just  $N(0, 1)$  values, then multiplies them by  $\sigma_R$  to yield  $\epsilon_t \sim N(0, \sigma_R^2)$ .

The parameter value for  $\sigma_R$  can be found in `currentRes$extra$residuals$p_log_RecDev[6]`.

### Author(s)

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

### See Also

[importProj](#)  
**coda:** [mcmc](#)

importRes

*Import Coleraine Model Results (RH version)***Description**

Import Coleraine model results from .res file, and rearrange into a standard format suitable for plotting.

**Usage**

```
importRes(res.file, info="", Dev=FALSE, CPUE=FALSE, Survey=FALSE,
          CAc=FALSE, CAs=FALSE, CLc=FALSE, CLs=FALSE, LA=FALSE, quiet=TRUE,
          extra=TRUE, sep=" ")
```

**Arguments**

res.file	name of Coleraine model results file to import.
info	optional string containing information to store with model results.
Dev	logical: whether recruitment deviates were estimated in model.
CPUE	logical: whether model was fitted to catch-per-unit-effort data.
Survey	logical: whether model was fitted to survey abundance index data.
CAc	logical: whether model was fitted to commercial catch-at-age data.
CAs	logical: whether model was fitted to survey catch-at-age data.
CLc	logical: whether model was fitted to commercial catch-at-length data.
CLs	logical: whether model was fitted to survey catch-at-length data.
LA	logical: whether model was fitted to length-at-age data.
quiet	logical: whether to report progress while parsing file.
extra	logical: if TRUE, import likelihoods, parameters, priors, and recruitment residuals.
sep	the field separator character (usually " " or "\t").

**Details**

This function was modified from the original importCol function in the **scape** package to grab extra data and to deal with anomalous characters in Coleraine results files.

**Value**

A list of class `list` containing at least `N`, `B`, and `Se1`. The other elements may or may not be included in the list, depending on how `importRes` was called:

<code>N</code>	predicted numbers at age
<code>B</code>	predicted biomass, recruitment, and observed landings (year things)

Se1	predicted selectivity and observed maturity (age things)
Dev	predicted recruitment deviates from the stock-recruitment curve
CPUE, Survey	commercial and survey abundance index and fit
CAC, CAs	commercial and survey C@A (catch at age) and fit
CLc, CLs	commercial and survey C@L (catch at length) and fit
LA	observed L@A and fit

### Note

This import function is implemented for the Coleraine statistical catch-at-age software, and can serve as a template for **scape** users who would like to implement import functions for specific stock assessment software.

The functions `ll` (package **gdata**) and `head` are recommended for browsing model results, e.g. `ll(x.cod)`; `ll(x.cod$N)`; `head(x.cod$N)`.

### References

Hilborn, R., Maunder, M., Parma, A., Ernst, B., Payne, J., and Starr, P. (2003) Coleraine: A generalized age-structured stock assessment model. User's manual version 2.0. *University of Washington Report SAFS-UW-0116*. Available at:  
<http://fish.washington.edu/research/coleraine/pdf/coleraine.pdf>.

### See Also

`importCor`, `importPar`, `importStd`  
`runADMB`, `readAD`, `reweight`, `importCol2`  
`read.table`, `readLines`, and `scan` to import any data.  
**scape-package** gives an overview of the package **scape**.

---

importStd

---

*Import Awatea Estimated Parameter File*


---

### Description

Import an Awatea file of estimated parameters resulting from an MPD minimisation.

### Usage

```
importStd(std.file, vnam="name")
```

### Arguments

std.file	parameter file (e.g., Awatea.std).
vnam	field name listing parameter names in data file.



**Details**

Extracts the data from character vectors and makes various data objects.

**Value**

The output is a list object with the following components:

std	data.frame representation of the .std file
parameters	data.frame for each of the estimated parameters in std

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[importCor](#), [importPar](#), [importRes](#)

---

load.allResFiles	<i>Load All Awatea .res Files</i>
------------------	-----------------------------------

---

**Description**

Load all Awatea .res files in the working directory into a list object.

**Usage**

```
load.allResFiles(resList = NULL)
```

**Arguments**

resList	AME: sets directory to path above current and sets the pattern to "results.dat\$"; probably deprecated.
---------	---

**Value**

List of multiple calls to [importCol2](#).

**Note**

If deprecated, remove from package **PBSawatea**.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[importCol2](#)

MAfun2

*Calculate Mean Age by Year***Description**

Calculate mean ages from proportions-at-age (modified from a subfunction in runADMB).

**Usage**

```
MAfun2(padata, brks=NULL)
```

**Arguments**

padata	proportion-at-age data CAC or CAs from a call to importCol2.
brks	breaks specified as numeric years to split the commercial data up into regimes that may account for index discontinuities ( <b>not used</b> ).

**Details**

Mean age function supplied by Chris Francis (2011).

padata has fields:

Series...series identifier

Year.....numeric year

Age.....age bin

Obs.....observed proportions

Fit.....predicted (fitted) proportions

SS.....sample size (effective  $N$ )

**Value**

List object of observed and expected mean ages, variance of expected ages, and a few bits and bobs.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC.

**References**

Francis, R.I.C.C. (2011, in press) Data weighting in statistical fisheries stock assessment models. *Canadian Journal of Fisheries and Aquatic Sciences*.

**See Also**

[runADMB](#), [importCol2](#)

---

`mainMenu`*Create a Menu of Options and Actions*

---

### Description

From a main menu, the user can choose various options and actions.

We tend to do everything from the command line so the menu functionality fosters the warning: *CAVEAT EMPTOR*.

### Usage

```
mainMenu()  
loadMenu()  
mpdMenu()  
mcmcMenu()  
utilMenu()
```

### Details

#### Main menu items:

- Import files
- MPD plots
  - Plot all MPD graphs
  - Save all MPD plots to PNG
- MCMC plots
  - Plot all MCMC plots
  - Save all MCMC plots to PNG
- Close all graphics windows
- Help & Utilities

#### Load menu items:

- Get Awatea res file
- Get Awatea MCMC file
- Get Awatea projection file
- Load all res files in working directory
- Get PJS Delay Difference MCMC+Projection

#### MPD menu items:

- Plot biomass, recruitment, catch
- Plot numbers at age
- Plot selectivity and maturity
- Plot commercial catch-at-age results
- Plot survey catch-at-age results
- Plot survey catch-at-length results
- Plot abundance index
- All residual plots
- Plot multi-panel biomass, recruitment, catch
- Plot multi-panel exploitation rate
- Plot alternative numbers at age

**MCMC menu items:**

Plot biomass and projections by policy  
 Probability of projection biomass > reference  
 Expectation of projection biomass / reference  
 Plot biomass posterior densities (plotDens)  
 Plot recruitment posterior densities (plotDens)  
 Plot parameter posterior densities (plotDens)  
 Plot cumulative quantiles (plotCumu)  
 Plot traces (plotTrace)  
 Plot PJS traces (plt.allTraces)

**Utils menu items:**

scape Help  
 scapeMCMC Help  
 Portrait graphsheet  
 Landscape graphsheet

**See Also**

[get.resFile](#), [importCol2](#)

---

makeErrMat

*Make Ageing Error Matrix for Awatea*

---

**Description**

Make a simple ageing error matrix for Awatea.

**Usage**

```
makeErrMat(N=60, ondiag=0.8, offdiag=0.1, corner=0.9)
```

**Arguments**

N	numeric scalar indicating number of age classes, which determines the dimension of the matrix.
ondiag	numeric value to appear along the matrix diagonal.
offdiag	numeric value to appear one cell to the left and right of the matrix diagonal.
corner	numeric value to appear in the top left and bottom right corners of the matrix.

**Value**

Simple symmetric ageing error matrix.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**[plotProp](#), [weightBio](#)

---

`msyCalc`*Calculate the Maximum Sustainable Yield*

---

**Description**

Load in `MSY.out` and calculate the MSY (maximum sustainable yield).

**Usage**

```
msyCalc(dir = getwd(), error.rep = 1)
```

**Arguments**

<code>dir</code>	working directory.
<code>error.rep</code>	numeric: if 1, report errors (reaching bounds), if 0 do not.

**Value**

Returns a list object containing:

- `yield...` maximum sustainable yield,
- `u.....` exploitation rate at MSY,
- `VB.....` vulnerable biomass at MSY,
- `B.....` spawning biomass at MSY,
- `nProj...` number of projections needed to reach MSY.

**Note**

See `msyTestCreating.r` for full details when figuring this out.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**[findTarget](#)

---

`out.pmTables`*Write Decision Tables to Comma-Delimited Files*

---

**Description**

Write decision tables to comma-delimited text files (.csv).

**Usage**

```
out.pmTables(obj, fileName="pm", dec=3)
```

**Arguments**

<code>obj</code>	list object containing tables (matrices or data frames).
<code>fileName</code>	prefix for output file names.
<code>dec</code>	number of decimal places to retain.

**Value**

Comma-delimited text files (.csv).

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[write.table](#), [writeList](#)

---

`panLab`*Write Text to Figure in Relative (0:1) Coordinates*

---

**Description**

Write text to a figure by first setting the coordinate space to lie between 0 and 1:  
`par(usr=c(0,1,0,1))`.

**Usage**

```
panLab(x, y, txt, ...)
```

**Arguments**

x	relative x-coordinate.
y	relative y-coordinate.
txt	text to add to figure.
...	additional arguments sent to function text.

**Note**

Currently, this function does not reset the coordinate space to the original.  
Use [addLabel](#) instead.

**See Also**

[addLabel](#), [addLegend](#)

---

panLegend

*Place a Legend in a Figure using Relative (0:1) Coordinates*

---

**Description**

Place a legend in a figure by first setting the coordinate space to lie between 0 and 1:  
`par(usr=c(0,1,0,1))`.

**Usage**

```
panLegend(x, y, legTxt, ...)
```

**Arguments**

x	relative x-coordinate.
y	relative y-coordinate.
legTxt	legend text to add to figure.
...	additional arguments sent to function legend.

**Note**

Currently, this function does not reset the coordinate space to the original.  
Use [addLegend](#) instead.

**See Also**

[addLabel](#), [addLegend](#)

## Description

**PBSawatea** contains the code used for the modelling of populations of Pacific Ocean Perach (*Sebastes alutus*) and Yellowmouth Rockfish (*S. reedi*) along the British Columbia (BC) coast.

Implementation is done using a modified version of the Coleraine statistical catch-at-age software (Hilborn *et al.* 2003) called Awatea (Alan Hicks, NOAA, pers. comm.). Awatea is a platform for implementing the AD (Automatic Differentiation) Model Builder software (Otter Research 1999), which provides (a) maximum posterior density estimates using a function minimiser and automatic differentiation, and (b) an approximation of the posterior distribution of the parameters using the Markov Chain Monte Carlo (MCMC) method, specifically using the Hastings-Metropolis algorithm (Gelman *et al.* 2004).

Running of Awatea is streamlined using code written in R (R Development Core Team 2009), rather than the original Microsoft Excel implementation. Figures and tables of output are automatically produced through R using code adapted from the R packages **scape** (Magnusson 2009) and **scapeMCMC** (Magnusson and Stewart 2007). We use the R function Sweave (Leisch 2008) in the package **utils** to automatically collate, via LATEX, the large amount of figures and tables into a single portable document file (.pdf) for each model run.

We provide master Sweave files used in folder `../library/PBSawatea/snw` to build the .pdf document. The user must copy these to a local working directory if they are not already there.

## References

- Gelman, A., Carlin, J.B., Stern, H.S. and Rubin, D.B. (2004) Bayesian data analysis, 2nd edition. Chapman and Hall/CRC, New York, 668 p.
- Hilborn, R., Maunder, M., Parma, A., Ernst, B. Payne, J., and Starr, P. (2003) Coleraine: a generalized age-structured stock assessment model. *School of Aquatic and Fishery Sciences*, University of Washington, 54 p.
- Leisch, F. (2008) Sweave, R package.
- Magnusson, A. (2009) Scape – statistical catch-at-age plotting environment, R package.
- Magnusson, A. and Stewart, I. (2007) MCMCscape – MCMC diagnostic plots. R package.
- Otter Research Ltd. (1999) An introduction to AD Model Builder for use nonlinear modeling and statistics. Otter Research Ltd., British Columbia. 194 p.
- R Development Core Team (2011) R: A language and environment for statistical computing. R Foundation for Statistical Computing, Vienna, Austria.  
ISBN 3-900051-07-0  
<http://www.R-project.org>



plotB2

*Plot Biomass, Recruitment, and Landings (AME Version)***Description**

Plot scape model predicted biomass, stock recruitment, and landings.

AME: This is an alteration of Arni Magnussons plotB function to accommodate PJS's request not to show biomass prior to fishery and survey indices period.

**Usage**

```
plotB2(model, what="d", series=NULL, years=NULL, axes=TRUE, div=1,
       legend="bottom", main="", xlab="", ylab="", cex.main=1.2,
       cex.legend=1, cex.lab=1, cex.axis=0.8, las=1,
       tck=c(1,what=="d")/2, tick.number=5, lty.grid=3, col.grid="white",
       pch=16, cex.points=0.8, col.points="black", lty.lines=1:3,
       lwd.lines=2, col.lines="black", ratio.bars=3, col.bars="grey",
       plot=TRUE, ...)
```

**Arguments**

model	fitted scape model.
what	what to plot: "d"[efault], "s"[tock recruitment], or "l"[andings].
series	vector of strings indicating which column names in model\$B data frame to plot (all by default).
years	vector of numbers indicating which years to include (all by default).
axes	whether to plot axis values.
div	denominator to shorten values on the y axis, or a vector with two elements referring to x and y axis.
legend	legend location: "bottom", "left", "top", "right", or "" to suppress legend.
main	main title.
xlab	x-axis label.
ylab	y-axis label.
cex.main	size of main title.
cex.legend	size of legend text.
cex.lab	size of axis labels.
cex.axis	size of tick labels.
las	orientation of tick labels: 0=parallel, 1=horizontal, 2=perpendicular, 3=vertical.
tck	tick mark length.
tick.number	number of tick marks.
lty.grid	line type of gridlines.

<code>col.grid</code>	color of gridlines.
<code>pch</code>	symbol for points.
<code>cex.points</code>	size of points.
<code>col.points</code>	color of points.
<code>lty.lines</code>	line type of main lines.
<code>lwd.lines</code>	line width of main lines.
<code>col.lines</code>	color of main lines.
<code>ratio.bars</code>	width of bars.
<code>col.bars</code>	color of bars.
<code>plot</code>	whether to draw plot.
<code>...</code>	passed to <code>xyplot</code> and <code>panel.superpose</code> .

**Details**

The "d"[efault] plot shows spawning biomass and vulnerable biomass as lines, and landings as bars, on the same scale.

**Value**

When `plot=TRUE`, a trellis plot is drawn and a data frame is returned, containing the data used for plotting. When `plot=FALSE`, a trellis object is returned.

**Note**

The `Args` function from the **gdata** package is recommended for reviewing the arguments, instead of `args`.

**See Also**

[xyplot](#), [panel.barchart](#), and [panel.superpose](#) are the underlying drawing functions.  
[plotCA](#), [plotCL](#), [plotIndex](#), [plotIndex2](#) and [plotLA](#) plot model fit and data.  
[plotB](#), [plotN](#), and [plotSel](#) plot derived quantities.  
[scape-package](#) gives an overview of the **scape** package.

---

<code>plotBars</code>	<i>Barplots of Annual Age Proportions</i>
-----------------------	---

---

**Description**

Plot barplots of specific-year age proportions.

**Usage**

```
plotBars(res, type="N", prop=TRUE, year=min(res[[type]][["Year"]]),
  sex=c(2,1), age=NULL, fill=c("orange","cyan"),
  eps=FALSE, pix=FALSE, win=TRUE, ...)
```

**Arguments**

res	Awatea results file from a call to importRes: importRes("POPrun05/pop-3CD.05.01.res", Dev=T, CPUE=T, Survey=T, CAc=T, CAs=T)
type	type of annual value (e.g., "N" = numbers, "B" = biomass).
prop	logical: if TRUE, convert values from type to proportions.
year	numeric scalar or vector of years.
sex	sex code where 1 = males and 2 = females; note that Awatea uses females before males.
age	specify age vector if subset of available is desired.
fill	bar colour by sex for barplots (one colour per sex).
eps	logical: if TRUE, send figure to a postscript (.eps) file.
pix	logical: if TRUE, send figure to a portable network graphics (.png) file.
win	logical: if TRUE, send figure to the R windows device.
...	extra parameters (not currently used).

**Details**

The plot is used primarily to see how an age composition for a year compares with an equilibrium age structure (represented by an exponential decay from 1 using  $e^{-M}$ ).

**Value**

Aside from the figure plots, the function invisibly returns a list of:

dat	data file from importRes for the type and year specified.
mat	three-dimensional array (age, sex, year) for the year(s) specified.
xpos	x-position generated by barplot for the ages specified.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotBox](#), [compB0](#)

---

plotBmcmcPOP

*Plot Spawning and Vulnerable Biomass*


---

## Description

Plot spawning and vulnerable biomass from posterior as boxplots, and add catch bars on same graph.

## Usage

```
plotBmcmcPOP( obj, currentRes1=currentRes,
               p=c(0.025,0.25,0.5,0.75,0.975),
               xyType="quantBox",
               lineType=c(3,2,1,2,3),
               refLines=NULL, xlim=NULL, ylim=NULL,
               userPrompt=FALSE, save=TRUE, xLab=c(1939,1939,1939),
               yLab=c(10000,70000,170000),
               textLab=c("catch","spawning","vulnerable"),
               yaxis.by=10000, tcl.val=-0.2, ...)
```

## Arguments

obj	MCMC data frame of $B$ (currentMCMC\$B).
currentRes1	list/scape object created by <a href="#">importCol2</a> .
p	quantiles to use in quantBox.
xyType	type of plot (currently only uses quantBox).
lineType	line types to use in quantBox.
refLines	reference lines to add to plot.
xlim	limits of the x-axis.
ylim	limits of the y-axis.
userPrompt	<b>not used</b>
save	<b>not used</b>
xLab	x-coordinates for labels.
yLab	y-coordinates for labels.
textLab	text labels to display on plot.
yaxis.by	increments along the y-axis to place tick marks.
tcl.val	tick length.
...	additional arguments passed to the function rect.

**Note**

Combines ideas from `plt.quantBio` and `plotB2`. Don't need lattice, just one figure, no panels. Vulnerable biomass has no posterior saved, which must be why it's not been done before. Hmmm.... still worth seeing spawning though?

Taking what is needed from `plt.quantBio`, this basically works:

```
plt.quantBio(currentMCMC$B, xyType=rpType),
```

though it creates 2x3 plots. The object should be the specific MCMC posterior by year (so just a data frame), *e.g.*, `currentMCMC$B`.  
`currentRes1` is local `currentRes`.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotVBcatch](#), [plotB2](#)

---

plotBox

*Plot Boxes using Quantiles*

---

**Description**

Produce box-and-whisker plot(s) of the given (grouped) values. This function is simply a modified version of [boxplot](#) that sets the whiskers to specified quantiles rather than 1.5 IRQ.

**Usage**

```
plotBox(x, ..., range=1.5, width=NULL, varwidth=FALSE,
        notch=FALSE, outline=TRUE, names, plot=TRUE,
        border=par("fg"), col=NULL, log="",
        pars=list(boxwex=0.8, staplewex=0.5, outwex=0.5, whisklty=1),
        horizontal=FALSE, add=FALSE, at=NULL,
        quants=c(0.025,0.25,0.5,0.75,0.975), outliers=FALSE)
```

**Arguments**

**x** for specifying data from which the boxplots are to be produced. Either a numeric vector, or a single list containing such vectors. Additional unnamed arguments specify further data as separate vectors (each corresponding to a component boxplot). *NAs* are allowed in the data.  
 Arguments `formula`, `data`, `codesubset`, and `na.action` are itemized below in **Notes**, but have not been tested for `plotBox`.

...	For the formula method, named arguments to be passed to the default method. For the default method, unnamed arguments are additional data vectors (unless <code>x</code> is a list when they are ignored), and named arguments are arguments and <a href="#">graphical parameters</a> to be passed to <code>bxp</code> in addition to the ones given by argument <code>pars</code> (and override those in <code>pars</code> ). Note that <code>bxp</code> may or may not make use of graphical parameters it is passed: see its documentation.
range	this determines how far the plot whiskers extend out from the box. If <code>range</code> is positive, the whiskers extend to the most extreme data point which is no more than <code>range</code> times the interquartile range from the box. A value of zero causes the whiskers to extend to the data extremes.
width	a vector giving the relative widths of the boxes making up the plot.
varwidth	if <code>varwidth</code> is TRUE, the boxes are drawn with widths proportional to the square-roots of the number of observations in the groups.
notch	if <code>notch</code> is TRUE, a notch is drawn in each side of the boxes. If the notches of two plots do not overlap this is 'strong evidence' that the two medians differ (Chambers <i>et al.</i> , 1983, p. 62). See <a href="#">boxplot.stats</a> for the calculations used.
outline	if <code>outline</code> is not true, the outliers are not drawn (as points whereas S+ uses lines).
names	group labels which will be printed under each boxplot. Can be a character vector or an <a href="#">expression</a> (see <a href="#">plotmath</a> ).
plot	if TRUE (the default) then a boxplot is produced. If not, the summaries which the boxplots are based on are returned.
border	an optional vector of colors for the outlines of the boxplots. The values in <code>border</code> are recycled if the length of <code>border</code> is less than the number of plots.
col	if <code>col</code> is non-null it is assumed to contain colors to be used to colour the bodies of the box plots. By default they are in the background colour.
log	character indicating if <code>x</code> or <code>y</code> or both coordinates should be plotted in log scale.
pars	a list of (potentially many) more graphical parameters, e.g., <code>boxwex</code> or <code>outpch</code> ; these are passed to <code>bxp</code> (if <code>plot</code> is true); for details, see there. Some explicit settings: <code>boxwex</code> – a scale factor to be applied to all boxes. When there are only a few groups, the appearance of the plot can be improved by making the boxes narrower. <code>staplewex</code> – staple line width expansion, proportional to box width. <code>outwex</code> – outlier line width expansion, proportional to box width. <code>whisklty</code> – whisker line type.
horizontal	logical indicating if the boxplots should be horizontal; default FALSE means vertical boxes.
add	logical, if true <i>add</i> boxplot to current plot.
at	numeric vector giving the locations where the boxplots should be drawn, particularly when <code>add</code> = TRUE; defaults to <code>1:n</code> where <code>n</code> is the number of boxes.
quants	numeric vector of 5 quantiles to specify (i) the extent of the lowest whisker, (ii) the lower boundary of the box, (iii) the middle line of the box, (iv) the upper boundary of the box, and (v) the extent of the upper whisker.

outliers            logical: if TRUE show the outliers (but used primarily to suppress outliers when FALSE).

## Details

The generic function `boxplot` currently has a default method (`boxplot.default`) and a formula interface (`boxplot.formula`).

If multiple groups are supplied either as multiple arguments or via a formula, parallel boxplots will be plotted, in the order of the arguments or the order of the levels of the factor (see [factor](#)).

Missing values are ignored when forming boxplots.

## Value

List with the following components:

stats	a matrix, each column contains the extreme of the lower whisker, the lower hinge, the median, the upper hinge and the extreme of the upper whisker for one group/plot. If all the inputs have the same class attribute, so will this component.
n	a vector with the number of observations in each group.
conf	a matrix where each column contains the lower and upper extremes of the notch.
out	the values of any data points which lie beyond the extremes of the whiskers.
group	a vector of the same length as <code>out</code> whose elements indicate to which group the outlier belongs.
names	a vector of names for the groups.

## Additional arguments

Additional arguments used by `boxplot` but not tested in `plotBox`:

formula.....	A formula, such as <code>y ~ grp</code> , where <code>y</code> is a numeric vector of data values to be split into groups according to the grouping variable <code>grp</code> (usually a factor).
data.....	A data.frame (or list) from which the variables in <code>formula</code> should be taken.
subset.....	An optional vector specifying a subset of observations to be used for plotting.
na.action...	A function which indicates what should happen when the data contain NAs. The default is to ignore missing values in either the response or the group.

## References

Becker, R. A., Chambers, J. M. and Wilks, A. R. (1988) *The New S Language*. Wadsworth & Brooks/Cole.

Chambers, J. M., Cleveland, W. S., Kleiner, B. and Tukey, P. A. (1983) *Graphical Methods for Data Analysis*. Wadsworth & Brooks/Cole.

Murrell, P. (2005) *R Graphics*. Chapman & Hall/CRC Press.

See also [boxplot.stats](#).

**See Also**

See [compB0](#) to show MCMC distributions of  $B$  relative to  $B_0$ .

[boxplot.stats](#) which does the computation, [bxp](#) for the plotting and more examples; and [stripchart](#) for an alternative (with small data sets).

---

plotBVBnorm

*Plot Spawning and Vulnerable Biomass Relative to Virgin*


---

**Description**

Plot spawning and vulnerable biomass boxplots relative to virgin levels  $B_0$  and  $V_0$ , respectively.

**Usage**

```
plotBVBnorm(mcmcObj,
             p=c(0.025,0.25,0.5,0.75,0.975),
             xyType="quantBox",
             lineType=c(3,2,1,2,3),
             refLines=NULL, xLim=NULL, yLim=NULL,
             userPrompt=FALSE, save=TRUE, xLeg=0.7, yLeg=0.9,
             yaxis.by=0.02, tcl.val=-0.2,
             B.col="black", VB.col="black", ...)
```

**Arguments**

mcmcObj	MCMC list object (currentMCMC).
p	quantiles to use in quantBox.
xyType	type of plot (currently only uses quantBox).
lineType	line types to use in quantBox.
refLines	reference lines to add to plot.
xLim	limits of the x-axis.
yLim	limits of the y-axis.
userPrompt	<b>not used</b>
save	<b>not used</b>
xLeg	x-coordinate for legend.
yLeg	y-coordinate for legend.
yaxis.by	increments along the y-axis to place tick marks.
tcl.val	tick length.
B.col	colour for spawning biomass.
VB.col	colour for vulnerable biomass.
...	<b>not used</b>



**Note**

AME: tried in separate file, but then changed that to lattice and wouldn't be good format for Arni's boxplots.

Based on plotVBcatch (tweaking some).  
currentRes1 is local currentRes.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotVBcatch](#), [plotBmcmcPOP](#)

---

plotChains

*Plot Cumulative Frequency of MCMC Chains*

---

**Description**

Plot cumulative frequency of  $n$  chains by partitioning one trace.  
(Modified from the function plotTracePOP.)

**Usage**

```
plotChains(mcmc, nchains=3, pdisc=0.1, axes=FALSE, same.limits=FALSE,
           between=list(x=axes,y=axes), div=1, span=1/4, log=FALSE,
           base=10, main=NULL, xlab=NULL, ylab=NULL, cex.main=1.2,
           cex.lab=1, cex.strip=0.8, cex.axis=0.8,
           las=0, tck=0.5, tick.number=5, lty.trace=1, lwd.trace=1,
           col.trace="grey", lty.median=1, lwd.median=1,
           col.median="black", lty.quant=2, lwd.quant=1,
           col.quant="black", plot=TRUE, probs=c(0.025, 0.5, 0.975), ...)
```

**Arguments**

mcmc	MCMC chain(s) as a vector, data frame or mcmc object.
ncchains	number of chains to create from one trace.
pdisc	proportion of the initial trace to discard before creating chains.
axes	whether axis values should be plotted.
same.limits	whether panels should have same x-axis limits.
between	list with x and y indicating panel spacing.
div	denominator to shorten values on the y axis.
span	smoothness parameter ( <b>not used</b> ).
log	whether values should be log-transformed.

<code>base</code>	logarithm base.
<code>main</code>	main title.
<code>xlab</code>	x-axis title.
<code>ylab</code>	y-axis title.
<code>cex.main</code>	size of main title.
<code>cex.lab</code>	size of axis labels.
<code>cex.strip</code>	size of strip labels.
<code>cex.axis</code>	size of tick labels.
<code>las</code>	orientation of tick labels: 0=parallel, 1=horizontal, 2=perpendicular, 3=vertical.
<code>tck</code>	tick mark length.
<code>tick.number</code>	number of tick marks.
<code>lty.trace</code>	line type of trace.
<code>lwd.trace</code>	line width of trace.
<code>col.trace</code>	colour of trace.
<code>lty.median</code>	line type of median.
<code>lwd.median</code>	line width of median.
<code>col.median</code>	colour of median.
<code>lty.quant</code>	line type of quantile trace.
<code>lwd.quant</code>	line width of quantile trace.
<code>col.quant</code>	colour of quantile trace.
<code>plot</code>	whether to draw plot.
<code>probs</code>	quantile values for quantile trace.
<code>...</code>	passed to <code>panel.trace</code> ( <b>not used</b> ).

**Value**

When `plot=TRUE`, a trellis plot is drawn and a data frame is returned, containing the data used for plotting. When `plot=FALSE`, a trellis object is returned.

**Note**

This idea stemmed from a discussion with PJS.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotTracePOP](#), [plotDensPOP](#)

---

plotCPUE*Plot CPUE and Add Error Bars*

---

**Description**

Plot CPUE and fit with error bars.

**Usage**

```
plotCPUE(obj, main="", save=NULL, bar=1.96, yLim=NULL, ...)
```

**Arguments**

obj	data frame of CPUE indices from Awatea's results file ( <i>e.g.</i> , <code>currentRes\$CPUE</code> ).
main	title for figure
save	<b>not used</b>
bar	standard deviation of the normal distribution (1.96 is the approximate value of the 97.5 percentile point).
yLim	limits of the y-axis.
...	<b>not used</b>

**Value**

A postscript file:  
`CPUEser.eps`...CPUE indices with error bars.

**Note**

Copied code from `plotIndexNotLattice`.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotIndexNotLattice](#)

---

plotDensPOP

*Plot MCMC Density (AME Version)*


---

## Description

Plot Markov-chain Monte Carlo density. This is an approximation of the posterior probability density function.

## Usage

```
plotDensPOP(mcmc, probs=c(0.025,0.975), points=FALSE, axes=TRUE,
  same.limits=FALSE, between=list(x=axes,y=axes), div=1,
  log=FALSE, base=10, main=NULL, xlab=NULL, ylab=NULL,
  cex.main=1.2, cex.lab=1, cex.strip=0.8, cex.axis=0.7,
  las=0, tck=0.5, tick.number=5,
  lty.density=1, lwd.density=3, col.density="black",
  lty.median=2, lwd.median=1, col.median="darkgrey",
  lty.outer=3, lwd.outer=1, col.outer="darkgrey", pch="|",
  cex.points=1, col.points="black", plot=TRUE, MPD.height=0.04,
  mpd = mcmc[1, ], ...)
```

```
plotDensPOPpars(mcmc, probs=c(0.025,0.975), points=FALSE,
  axes=TRUE, same.limits=FALSE, between=list(x=axes,y=axes),
  div=1, log=FALSE, base=10, main=NULL, xlab=NULL, ylab=NULL,
  cex.main=1.2, cex.lab=1, cex.strip=0.8, cex.axis=0.7,
  las=0, tck=0.5, tick.number=5,
  lty.density=1, lwd.density=3, col.density="black",
  lty.median=2, lwd.median=1, col.median="darkgrey",
  lty.outer=3, lwd.outer=1, col.outer="darkgrey", pch="|",
  cex.points=1, col.points="black", plot=TRUE, MPD.height=0.04,
  mpd = mcmc[1, ], ...)
```

```
plotDensPOPparsPrior(mcmc, probs=c(0.025,0.975), points=FALSE,
  axes=TRUE, same.limits=FALSE, between=list(x=axes,y=axes),
  div=1, log=FALSE, base=10, main=NULL, xlab=NULL, ylab=NULL,
  cex.main=1.2, cex.lab=1, cex.strip=0.8, cex.axis=0.7,
  las=0, tck=0.5, tick.number=5,
  lty.density=1, lwd.density=3, col.density="black",
  lty.median=2, lwd.median=1, col.median="darkgrey",
  lty.outer=3, lwd.outer=1, col.outer="darkgrey", pch="|",
  cex.points=1, col.points="black", plot=TRUE, MPD.height=0.04,
  mpd = mcmc[1, ], ...)
```

## Arguments

mcmc	MCMC chain(s) as a vector, data frame or mcmc object.
probs	vector of outer quantiles to draw, besides the median.

points	whether data points should be plotted along the x axis.
axes	whether axis values should be plotted.
same.limits	whether panels should have same x-axis limits.
between	list with x and y indicating panel spacing.
div	denominator to shorten values on the x axis.
log	whether values should be log-transformed.
base	logarithm base.
main	main title.
xlab	x-axis label.
ylab	y-axis label.
cex.main	size of main title.
cex.lab	size of axis labels.
cex.strip	size of strip labels.
cex.axis	size of tick labels.
las	orientation of tick labels: 0=parallel, 1=horizontal, 2=perpendicular, 3=vertical.
tck	tick mark length.
tick.number	number of tick marks.
lty.density	line type of density curve.
lwd.density	line width of density curve.
col.density	colour of density curve.
lty.median	line type of median.
lwd.median	line width of median.
col.median	colour of median.
lty.outer	line type of outer quantiles.
lwd.outer	line width of outer quantiles.
col.outer	colour of outer quantiles.
pch	symbol for data points.
cex.points	size of data points.
col.points	colour of data points.
plot	whether to draw plot.
MPD.height	how far up to put MPD.
mpd	vector of MPD values (from which MCMC search starts).
...	passed to densityplot and panel.densityplot.

### Details

The function `plotDensPOPpars` differs from `plotDensPOP` only by a few tweaks to the internal list object `myscales`.

The function `plotDensPOPparsPrior` adds the priors automatically.

Value

When `plot=TRUE`, a trellis plot is drawn and a data frame is returned, containing the data used for plotting. When `plot=FALSE`, a trellis object is returned.

Note

The `Args` function from the **gdata** package is recommended for reviewing the arguments, instead of `args`.

AME: edited `plotDens` function to have less whitesapce, not repeat x-axis labels, and make y-axes the same scales. Cannot just do it through the options.

For Recruits and Biomass, use `plotDensPOppars` for parameters. Tried y-axes the same scales, but 1973–1975 are so narrow that they make all the others really small: `same.limits=TRUE, ylim=c(0, 0.0005)`.

See Also

`xyplot` and `panel.densityplot` are the underlying drawing functions, and `densplot` is a similar non-trellis plot.

`plotTrace`, `plotAuto`, `plotCumu`, and `plotSplom` are diagnostic plots.

`plotDens` and `plotQuant` are posterior plots.

`scapeMCMC-package` gives an overview of the package.

---

plotIndex2	<i>Plot Abundance Index (AME Version)</i>
------------	---

---

Description

Plot scape model fit to abundance index data.  
Revised version of Arni's function to confine plotting to data region.

Usage

```
plotIndex2(model, what="c", series=NULL, axes=TRUE, same.limits=FALSE,
  between=list(x=axes,y=axes), ylim=NULL, q=1, bar=1, log=FALSE,
  base=10, main="", xlab="", ylab="", cex.main=1.2, cex.lab=1,
  cex.strip=0.8, cex.axis=0.8, las=1, tck=c(1,0)/2,
  tick.number=5, lty.grid=3, col.grid="white", pch=16,
  cex.points=1.2, col.points="black", lty.lines=1, lwd.lines=4,
  col.lines="dimgrey", lty.bar=1, plot=TRUE, ...)
```

Arguments

model	fitted scape model containing element CPUE and/or Survey.
what	what to plot: "c"[ommercial] or "s"[urvey] abundance index.
series	vector of strings indicating which gears or surveys to plot (all by default).

<code>axes</code>	whether to plot axis values.
<code>same.limits</code>	whether panels should have same y-axis limits.
<code>between</code>	list with x and y indicating panel spacing.
<code>ylim</code>	vector with lower and upper y-axis limits.
<code>q</code>	denominator to scale the y axis, e.g. to vulnerable biomass. Similar to the <code>div</code> argument in <code>plotN</code> and <code>plotB</code> .
<code>bar</code>	extent of error bars relative to standard error.
<code>log</code>	whether to log-transform values.
<code>base</code>	logarithm base.
<code>main</code>	main title.
<code>xlab</code>	x-axis label.
<code>ylab</code>	y-axis label.
<code>cex.main</code>	size of main title.
<code>cex.lab</code>	size of axis labels.
<code>cex.strip</code>	size of strip labels.
<code>cex.axis</code>	size of tick labels.
<code>las</code>	orientation of tick labels: 0=parallel, 1=horizontal, 2=perpendicular, 3=vertical.
<code>tck</code>	tick mark length.
<code>tick.number</code>	number of tick marks.
<code>lty.grid</code>	line type of gridlines.
<code>col.grid</code>	color of gridlines.
<code>pch</code>	symbol for points.
<code>cex.points</code>	size of points.
<code>col.points</code>	color of points and error bars.
<code>lty.lines</code>	line type of main lines.
<code>lwd.lines</code>	line width of main lines.
<code>col.lines</code>	color of main lines.
<code>lty.bar</code>	line type of error bars.
<code>plot</code>	whether to draw plot.
<code>...</code>	passed to <code>xyplot</code> , <code>panel.xyplot</code> , and <code>panel.xYplot</code> .

### Value

When `plot=TRUE`, a trellis plot is drawn and a data frame is returned, containing the data used for plotting. When `plot=FALSE`, a trellis object is returned.

### Note

The `Args` function from the **gdata** package is recommended for reviewing the arguments, instead of `args`.

**See Also**

[xyplot](#), [panel.xyplot](#), and [panel.xYplot](#) are the underlying drawing functions.  
[plotCA](#), [plotCL](#), [plotIndex](#), and [plotLA](#) plot model fit and data.  
[plotB](#), [plotB2](#), [plotN](#), and [plotSel](#) plot derived quantities.  
[scape-package](#) gives an overview of the package.

---

plotIndexNotLattice      *Plot Survey Indices*

---

**Description**

Plot index series with error bars. Create postscript files automatically.

**Usage**

```
plotIndexNotLattice(obj, objCPUE, main="", save=NULL, bar=1.96,
  ssnames=paste("Ser",1:9, sep=""), ...)
```

**Arguments**

obj	data frame of survey indices from Awatea's results file ( <i>e.g.</i> , <code>currentRes\$Survey</code> ).
objCPUE	data frame of CPUE indices from Awatea's results file ( <i>e.g.</i> , <code>currentRes\$CPUE</code> ).
main	title for figure
save	<b>not used</b>
bar	standard deviation of the normal distribution (1.96 is the approximate value of the 97.5 percentile point).
ssnames	survey series names for figure labelling and image creation.
...	<b>not used</b>

**Value**

Four postscript files:  
`survIndSer.eps`...each survey panel focuses on the years of the survey;  
`survIndSer2.eps`...each panel uses a fixed set of years that span all surveys;  
`survIndSer3.eps`...one panel showing all series normalised to their means;  
`survIndSer4.eps`...compares normalised GIG series with CPUE series.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plt.idx](#)



---

plotRmcmcPOP	<i>Plot Recruitment Posterior Quantiles</i>
--------------	---

---

**Description**

Plot recruitment posteriors quantiles as one graph over time.

**Usage**

```
plotRmcmcPOP( obj,
               p=c(0.025,0.25,0.5,0.75,0.975),
               xyType="quantBox",
               lineType=c(3,2,1,2,3),
               refLines=NULL, xLim=NULL, yLim=NULL,
               userPrompt=FALSE, save=TRUE, tcl.val=-0.2,
               yaxis.by=10000, yLab="Recruitment", ...)
```

**Arguments**

obj	MCMC data frame of $R$ (currentMCMC\$R).
p	quantiles to use in quantBox.
xyType	type of plot (currently only uses quantBox).
lineType	line types to use in quantBox.
refLines	reference lines to add to plot.
xLim	limits of the x-axis.
yLim	limits of the y-axis.
userPrompt	<b>not used</b>
save	<b>not used</b>
tcl.val	tick length.
yaxis.by	increments along the y-axis to place tick marks.
yLab	label for the y-axis.
...	additional arguments passed to the function rect.

**Note**

AME: Plot recruitment posteriors quantiles as one graph over time.

Already have the full posterior densities done.

Using plotBmcmcPOP as template, but will be simpler as no extra stuff. Probably not simplifying down as much as could due to time constraints.

Adding yLab and then using for exploitation plot also.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

See Also

[plotBmcmcPOP](#), [plotB2](#)

---

plotSnail	<i>Plot Snail Trails of Exploitation vs. Biomass</i>
-----------	--

---

Description

Plot the historical progression of the ratio  $u_t/u_{MSY}$  against  $B_t/B_{MSY}$ .

Usage

```
plotSnail(BoverBmsy, UoverUmsy, p=c(0.1,0.9),
          xLim=NULL, yLim=NULL, Lwd=2)
```

Arguments

- |           |   |
|-----------|---|
| BoverBmsy | numeric matrix of $B_t$ over $B_{MSY}$ .        |
| UoverUmsy | numeric matrix of $u_t$ over $u_{MSY}$ .        |
| p         | quantiles to show the bulk of the distribution. |
| xLim      | limits of the x-axis.                           |
| yLim      | limits of the y-axis.                           |
| Lwd       | line width of the snail trail.                  |

Details

The graph attempts to show the time history of the exploitation rate compared to the spawning biomass using a precautionary framework recast in Bayesian terms.

Note

The term *snail trail* comes from PJS.

Author(s)

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

See Also

[plotBVBnorm](#)

plotTracePOP

*Plot MCMC Traces (AME Version)***Description**

Plot Markov-chain Monte Carlo traces. This is a diagnostic plot for deciding whether a chain shows unwanted trends by showing the trace of moving quantiles. (Modified from the **scapeMCMC** function plotTrace.)

**Usage**

```
plotTracePOP(mcmc, axes=FALSE, same.limits=FALSE,
             between=list(x=axes,y=axes), div=1, span=1/4, log=FALSE,
             base=10, main=NULL, xlab=NULL, ylab=NULL, cex.main=1.2,
             cex.lab=1, cex.strip=0.8, cex.axis=0.8,
             las=0, tck=0.5, tick.number=5, lty.trace=1, lwd.trace=1,
             col.trace="grey", lty.median=1, lwd.median=1,
             col.median="black", lty.quant=2, lwd.quant=1,
             col.quant="black", plot=TRUE, probs=c(0.025, 0.5, 0.975),
             mpd=mcmc[1, ], ...)
```

**Arguments**

mcmc	MCMC chain(s) as a vector, data frame or mcmc object.
axes	whether axis values should be plotted.
same.limits	whether panels should have same x-axis limits.
between	list with x and y indicating panel spacing.
div	denominator to shorten values on the y axis.
span	smoothness parameter ( <b>not used</b> ).
log	whether values should be log-transformed.
base	logarithm base.
main	main title.
xlab	x-axis title.
ylab	y-axis title.
cex.main	size of main title.
cex.lab	size of axis labels.
cex.strip	size of strip labels.
cex.axis	size of tick labels.
las	orientation of tick labels: 0=parallel, 1=horizontal, 2=perpendicular, 3=vertical.
tck	tick mark length.
tick.number	number of tick marks.

lty.trace	line type of trace.
lwd.trace	line width of trace.
col.trace	colour of trace.
lty.median	line type of median.
lwd.median	line width of median.
col.median	colour of median.
lty.quant	line type of quantile trace.
lwd.quant	line width of quantile trace.
col.quant	colour of quantile trace.
plot	whether to draw plot.
probs	quantile values for quantile trace.
mpd	vector of MPD values (from which MCMC search starts).
...	passed to panel.trace ( <b>not used</b> ).

### Value

When plot=TRUE, a trellis plot is drawn and a data frame is returned, containing the data used for plotting. When plot=FALSE, a trellis object is returned.

### Note

The Args function from the **gdata** package is recommended for reviewing the arguments, instead of args.

### See Also

[xyplot](#) and [panel.loess](#) are the underlying drawing functions, and [traceplot](#) is a similar non-trellis plot.

[plotTrace](#), [plotAuto](#), [plotCumu](#), and [plotSplom](#) are diagnostic plots.

[plotDensPOP](#), [plotDens](#), [plotQuant](#), and [plotChains](#) are posterior plots.

[scapeMCMC-package](#) gives an overview of the **scapeMCMC** package.

---

plotVBcatch

*Plot Vulnerable Biomass*

---

### Description

Plot vulnerable biomass from posterior as boxplots, and add catch bars on same graph.

**Usage**

```
plotVBcatch( obj, currentRes1=currentRes,
             p=c(0.025,0.25,0.5,0.75,0.975),
             xyType="quantBox",
             lineType=c(3,2,1,2,3),
             refLines=NULL, xLim=NULL, yLim=NULL,
             userPrompt=FALSE, save=TRUE, xLab = c(1939,1939),
             yLab=c(10000,220000),
             textLab=c("catch","vulnerable"),
             yaxis.by=10000, tcl.val=-0.2, ...)
```

**Arguments**

obj	MCMC data frame of $B$ (currentMCMC\$B).
currentRes1	list/scape object created by <a href="#">importCol2</a> .
p	quantiles to use in quantBox.
xyType	type of plot (currently only uses quantBox).
lineType	line types to use in quantBox.
refLines	reference lines to add to plot.
xLim	limits of the x-axis.
yLim	limits of the y-axis.
userPrompt	<b>not used</b>
save	<b>not used</b>
xLab	x-coordinates for labels.
yLab	y-coordinates for labels.
textLab	text labels to display on plot.
yaxis.by	increments along the y-axis to place tick marks.
tcl.val	tick length.
...	additional arguments passed to the function rect.

**Note**

AME: This function is essentially a tweak of plotBmcmcPOP.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotBmcmcPOP](#), [plotB2](#)

---

plt.ageResidsPOP      *Plot Model Residuals*


---

**Description**

Plot model residuals as boxplots or qq-plots.

**Usage**

```
plt.ageResidsPOP (obj, ages=c(2,60), pct=c(5,25,50,75,95), main=NULL)
plt.ageResidsqqPOP(obj, ages=c(2,60), pct=c(5,25,50,75,95), main=NULL)
plt.yearResidsPOP (obj, ages=c(2,60), pct=c(5,25,50,75,95),
                  main=NULL, fill.in=TRUE, ... )
plt.cohortResids (obj, ages=c(2,59), pct=c(5,25,50,75,95), main=NULL)
```

**Arguments**

obj	output from <a href="#">stdRes.CA</a> .
ages	age classes to plot.
pct	quantiles to show in boxplot or qq-plot.
main	title for plot if desired.
fill.in	logical: if TRUE, add missing years to boxplot.
...	additional arguments for boxplot.

**Details**

```
plt.ageResidsPOP....plot age class residuals as boxplots.
plt.ageResidsqqPOP...plot age class residuals as qq-plot.
plt.yearResidsPOP....plot age residuals by year as boxplots.
plt.cohortResids....plot age residuals by cohort as boxplots.
```

**Note**

Some trouble noted adding text and legend.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plt.allTraces](#), [plt.expRate](#), [plt.idx](#), [plt.mcmcGraphs](#), [plotIndexNotLattice](#), [plotChains](#), [plotCPUE](#)

---

plt.allTraces	<i>Plot MCMC Traces</i>
---------------	-------------------------

---

**Description**

Plot traces from MCMC samples.

**Usage**

```
plt.allTraces(obj, bioYrList=NULL, recYrList=NULL, save=TRUE)
```

**Arguments**

obj	vector of MCMC samples.
bioYrList	years to plot spawning biomass traces.
recYrList	years to plot recruitment traces.
save	logical: if TRUE, save figure to a raster file (.jpg).

**Note**

Appears to be some figure requested by PJS.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotChains](#)

---

plt.expRate	<i>Plot Exploitation Rate</i>
-------------	-------------------------------

---

**Description**

Plot exploitation rate against year.

**Usage**

```
plt.expRate(obj, yLim=c(0,0.5), xLim=c(1954,2005))
```

**Arguments**

obj	an object from load.allResFiles.
yLim	limits of the y-axis.
xLim	limits of the x-axis.

**Details**

Simple points and lines plot.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plt.idx](#)

---

plt.idx	<i>Plot Survey Index Residuals</i>
---------	------------------------------------

---

**Description**

Plot the survey index residuals as a quantile-quantile plot (see [qqnorm](#)).

**Usage**

```
plt.idx(obj, main="Residuals", save=NULL, ssnames=paste("Ser",1:9,sep=""), ...)
```

**Arguments**

- |         |  |
|---------|--|
| obj     | a data frame with columns Year, stdRes, and Fit.             |
| main    | title for the plot.  |
| save    | logical: if TRUE, save the figure to a raster file (.png).   |
| ssnames | survey series names for figure labelling and image creation. |
| ...     | <b>not used</b>  |

**Details**

QQ-plots show sample quantiles vs. theoretical quantiles.

**Note**

The save option has been disabled for some reason.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plt.stdResids](#), [plotIndexNotLattice](#)



plt.mpdGraphs

*Plot a Set of Figures for MPD and MCMC***Description**

Plot a set of postscript figures .eps for the MPD (mode of the posterior distribution) and MCMC (Monte Carlo Markoff Chain) results.

**Usage**

```
plt.mpdGraphs(obj, save=FALSE, ssnames=paste("Ser",1:9,sep=""))
plt.mcmcGraphs(mcmcObj, projObj, save=FALSE,
               ylim.recruitsMCMC = NULL, ylim.exploitMCMC = NULL,
               ylim.VBcatch = NULL, ylim.BVBnorm = NULL,
               xlim.snail = NULL, ylim.snail = NULL,
               plotPolicies = names(projObj$Y[1:6]),
               onePolicy=names(projObj$Y[2]), mpd = list())
```

**Arguments**

obj	an Awatea results object (e.g., currentRes)
mcmcObj	an Awatea MCMC object (e.g., currentMCMC)
projObj	an Awatea projected biomass object (e.g., currentProj)
save	<b>not used</b>
ssnames	survey series names for figure labelling and image creation
ylim.recruitsMCMC	y-limits for the recruitsMCMC plot
ylim.exploitMCMC	y-limits for the recruitsMCMC plot
ylim.VBcatch	y-limits for the VBcatch plot
ylim.BVBnorm	y-limits for the BVBnorm plot
xlim.snail	x-limits for the snail-trail plot
ylim.snail	y-limits for the snail-trail plot
plotPolicies	six-policies projections to plot
onePolicy	the one policy from the above six to use for some figures
mpd	list of MPD values for parameters, biomass, and recruitment.

**Details**

Creates a whole heap o' postscript files.

**Value**

**plt.mpdGraphs** creates the following postscript files:

exploit.eps.....annual exploitation rate,  
 recruits.eps.....annual recruitment at age 1,  
 selectivity.eps.....selectivity curves for commercial gear(s) and survey(s),  
 ageComm.eps.....fits to annual commercial age composition (panel plots),  
 ageSurv.eps.....fits to annual survey age composition (panel plots),  
 survIndSer.eps.....four figures of survey indices (calls [plotIndexNotLattice](#)),  
 CPUEser.eps.....CPUE indices with error bars (calls [plotCPUE](#)),  
 commAgeResids.eps.....standardised residuals for commercial gear,  
 survAgeResidsSer.eps.....standardised residuals for surveys,  
 meanAge.eps.....mean age for catch and surveys,  
 stockRecruit.eps.....stock recruitment function.

**plt.mcmcGraphs** creates the following postscript files:

recruitsMCMC.eps.....boxplots of annual MCMC recruitment,  
 exploitMCMC.eps.....boxplots of annual MCMC exploitation rate,  
 pdfBiomass.eps.....density panel plots of annual female spawning biomass,  
 pdfRecruitment.eps.....density panel plots of annual recruitment,  
 selectivityMCMC.eps....**not currently implemented**,  
 traceRecruits.eps.....panel plots of annual recruitment traces with running quantiles,  
 traceBiomass.eps.....panel plots of annual spawning biomass with running quantiles,  
 traceParams.eps.....panel plots of parameter traces with running quantiles,  
 splitChain.eps.....panel plots of cumulative parameter estimate chains,  
 VBcatch.eps.....boxplots of annual vulnerable biomass and barplots of catch,  
 BVBnorm.eps.....spawning and vulnerable biomass relative to their virgin levels,  
 Bproj.eps.....boxplots of spawning biomass – MCMCs and projections,  
 snail.eps.....time series of  $u_t/u_{MSY}$  vs.  $B_t/B_{MSY}$ ,  
 pairs[1,2,3].eps.....pairs plot of parameter MCMC samples.

**Note**

The function `plt.mpdGraphs` was recently cleaned up by AME. Be wary not to use `currentRes` (globally available when the function should be using `obj`, which is the results object passed to the function).

The same is true for `plt.mcmcGraphs` where the global objects `currentMCMC` and `currentProj` are use within the function rather than using the local objects `mcmcObj` and `projObj`.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plt.mcmcGraphs](#), [plotIndexNotLattice](#), [plotCPUE](#), [plotChains](#), [plotSnail](#)

---

plt.numR	<i>Plot Numbers at Age at Equilibrium</i>
----------	---

---

**Description**

Plot numbers at age at equilibrium. Plot recruitment (age 1).

**Usage**

```
plt.numR(obj, minYr = NULL)
```

**Arguments**

obj	an Awatea results object from load.allResFiles.
minYr	minimum year to display in plot.

**Note**

Not sure what this plot is used for (RH).

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plotRmcmcPOP](#)

---

plt.quantBio	<i>Plot Quantile Boxes of MCMC and Projected Biomass</i>
--------------	--

---

**Description**

Plots MCMC and projected biomass as quantile boxes, the former in black, the latter in red.

**Usage**

```
plt.quantBio(obj, projObj=NULL, policy=NULL,
  p=c(0.025,0.25,0.5,0.75,0.975), xyType="lines",
  lineType=c(3,2,1,2,3), refLines=NULL,
  xLim=NULL, yLim=NULL, userPrompt=FALSE,
  save=TRUE, yaxis.lab="Spawning biomass" )

plt.quantBioBB0(obj, projObj=NULL, policy=NULL,
  p=c(0.025,0.25,0.5,0.75,0.975), xyType="lines",
  lineType=c(3,2,1,2,3), refLines=NULL,
```

```
xLim=NULL, yLim=NULL, userPrompt=FALSE, save=TRUE,
main="", cex.main="", tcl.val=-0.2,
xaxis.by=1, yaxis.by=10000, xaxis.lab="Year",
yaxis.lab="Spawning biomass" )
```

### Arguments

obj	an Awatea MCMC object ( <i>e.g.</i> , currentMCMC).
projObj	an Awatea projected biomass object ( <i>e.g.</i> , currentProj).
policy	numeric vector specifying catch policy.
p	quantiles to use from the biomass samples.
xyType	string specifying type of plot.
lineType	line types for the quantiles if xyType="lines".
refLines	reference points.
xLim	limits of the x-axis.
yLim	limits of the y-axis.
userPrompt	logical: if TRUE prompts user before figure is drawn.
save	logical: if TRUE save figure as a raster file .png.
main	character string specifying a title for the plot.
cex.main	font size for figure title.
tcl.val	tick length.
xaxis.by	tick mark intervals for x-axis.
yaxis.by	tick mark intervals for y-axis.
xaxis.lab	label for x-axis.
yaxis.lab	label for y-axis.

### Value

List of the reconstructed (MCMC) and projected results.

### Note

plt.quantBioBB0 performs similarly as for plt.quantBio but uses  $B_t/B_0$  instead of  $B_t$ .

### Author(s)

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

### See Also

[plotBmcmcPOP](#), [plotVBcatch](#), [plotBVBnorm](#), [plotRmcmcPOP](#)

---

plt.ssbVbCatch	<i>Plot Annual Spawning and Vulnerable Biomass</i>
----------------	--

---

**Description**

Plot MPD values of spawning biomass (SB) and vulnerable biomass(VB), as well as catch, against year.

**Usage**

```
plt.ssbVbCatch(obj, x1=1966, xLim=c(1954,2005), yLim=c(0,25000))
```

**Arguments**

obj	an Awatea results object from <code>load.allResFiles</code> .
x1	year to start plotting SB and VB lines.
xLim	limits of the x-axis.
yLim	limits of the y-axis.

**Note**

This analysis uses the MPD (mode of the posterior distribution) values for  $B$  and  $V$ .

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[load.allResFiles](#), [get.resFile](#), [plt.expRate](#)

---

plt.stdResids	<i>Plot Diagnostics for Standardised Residuals</i>
---------------	--

---

**Description**

Plot standardised residuals against year, fitted value, and theoretical residuals.

**Usage**

```
plt.stdResids(obj, pct=c(5,25,50,75,95),
  main=NULL, yLim=NULL, xLim=xLim)
```

**Arguments**

obj	a data frame with columns Year, stdRes, and Fit.
pct	percentiles to display as horizontal lines on the quantile-quantile plot.
main	title for the figure.
yLim	limits of the y-axis.
xLim	limits of the x-axis.

**Details**

Figure provides three panels of standardised residuals *vs.*  
 (i) years, (ii) fitted or predicted values, and (iii) theoretical quantiles.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plt.ssbVbCatch](#)

---

readAD

---

*Read ADMB Input and Create AWATEA Class Object*


---

**Description**

Read the ADMB input file and create an AWATEA class object.

**Usage**

```
readAD(txt)
```

**Arguments**

txt	string name of an Awatea input file.
-----	--------------------------------------

**Details**

The Awatea input file contains headers (lines prefixed with "#") and data that are read sequentially into the model by the binary executable `Awatea.exe`.

**Value**

An AWATEA class cobject with the slots:

- txtnam.....character: name of the input file,
- input.....character: vector of strings that are the lines of the input file,
- vlst.....list: each line of the input file with a label specifying line number and  
.....indicating whether the line is a Comment or Data,
- dnam.....character: vector of strings specifying data contents labelled by line number,
- nvars.....numeric: number of data variables,
- vdesc.....character: vector of strings specifying data contents labelled by  
.....variable number (*e.g.*, v001),
- vars.....list: numeric values of data labelled by variable number,
- gcomm.....character: vector of comments labelled by line number,
- vcomm.....character: vector of variable names labelled by line number,
- output.....list: Awatea results file imported by function [importRes](#),
- reweight...list: empty (later populated by function [reweight](#)).

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC.

**References**

Hilborn, R., Maunder, M., Parma, A., Ernst, B. Payne, J., and Starr, P. (2003) Coleraine: a generalized age-structured stock assessment model. School of Aquatic and Fishery Sciences, University of Washington, 54 p.

**See Also**

[runADMB](#), [reweight](#)

---

refPoints	<i>Calculate Reference Points</i>
-----------	-----------------------------------

---

**Description**

Calculate reference points relative to either  $B_{MSY}$  or  $B_0$ .

**Usage**

```
refPoints(mcmcObj=currentMCMC, projObj=currentProj,
          msyObj=currentMSY, refLevels=c(0.4,0.8,1))

refPointsB0(mcmcObj=currentMCMC, projObj=currentProj,
            B0Obj=B0.MCMC, refLevels=B0refLevels, refNames=B0refNames)
```

**Arguments**

mcmcObj	MCMC list object ( <i>e.g.</i> , currentMCMC).
projObj	projected biomass list object ( <i>e.g.</i> , currentProj).
msyObj	MSY list object ( <i>e.g.</i> , currentMSY).
refLevels	reference levels relative to $B_{MSY}$ (or $B_0$ ).
B0Obj	vector of $B_0$ MCMC values ( <i>e.g.</i> , B0.MCMC).
refNames	names of the $B_0$ reference levels refLevels.

**Value**

List of reference points relative to either  $B_{MSY}$  or  $B_0$ .

**Note**

Call from Sweave as `refPoints()` or, in full:  
`refPoints(currentMCMC, currentProj, currentMSY, refLevels=c(0.4,0.8,1))`

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[calc.refVal](#)

---

reweight	<i>Reweight Abundance and Composition Data</i>
----------	--

---

**Description**

Weight the abundance data by adjusting survey and CPUE CVs, and weight the composition data by adjusting the effective sample size  $N$ .

**Usage**

```
reweight(obj, cvpro=FALSE, mean.age=TRUE, ...)
```

**Arguments**

obj	an AWATEA class object created initially by readAD.
cvpro	CV process error added to CV observation error: $c_t = \sqrt{c_o^2 + c_p^2}$ ; if FALSE index CVs are reweighted using the standard deviation of normalized residuals.



mean.age      logical: if TRUE, use mean-age residuals to reweight the effective  $N$  for the age composition data (see Francis 2011);  
                  if FALSE, reweight  $N$  using  $\Sigma(P(1 - P))/\Sigma(O - P)^2$ , where  $O$  = observed proportions-at age and  $P$  = predicted/fitted proportions-at-age.  
 ...            additional arguments to reweight.

### Details

For the reweight to work, a corresponding Awatea results file (.res) with the same prefix as the input file must be available in the working directory before calling readAD. This will populate the output slot with fitted data that the reweight needs.

### Value

An AWATEA class cobject with the slots outlined in [readAD](#) with the following slot populated by this function:

```
reweight...list of reweight results:
..nrwt....the number of the current reweighting,
..survey...survey indices with CV values (observed, fitted, normalised residuals, reweighted),
..cpue....if used in the model, CPUE indices with CV values (as above),
..wNcpa....reweighted effective  $N$  for commercial compositions (proportions-at-age),
..wNspa....reweighted effective  $N$  for survey compositions,
..SDNR....standard deviation of normalised residuals for abundance and composition data,
..wj.....weights for composition data from a mean-age weighted calculation (Francis 2011).
```

### Author(s)

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC.

### References

Francis, R.I.C.C. (2011, in press) Data weighting in statistical fisheries stock assessment models. *Canadian Journal of Fisheries and Aquatic Sciences*.

### See Also

[runADMB](#), [readAD](#), [stdRes.CA](#), [stdRes.index](#)

---

runADMB

*Run AD Model Builder Code for Awatea*

---

### Description

Run compiled AD Model Builder code called Awatea to reconstruct a population trajectory for a marine fish stock.

**Usage**

```
runADMB(filename.ext, wd=getwd(), strSpp="XYZ", runNo=1, rwtNo=0,
  doMPD=FALSE, N.reweight=0, cvpro=FALSE, mean.age=TRUE,
  doMCMC=FALSE, mcmc=1e6, mcsave=1e3, ADargs=NULL, verbose=FALSE,
  doMSY=FALSE, msyMaxIter=15000., msyTolConv=0.01, endStrat=0.301, stepStrat=0.001,
  delim="-", awateaPath="E:/Projects/ADMB/Coleraine", clean=FALSE, ...)
```

**Arguments**

filename.ext	character file name including its extension.
wd	character string specifying the working directory, if different from the current working directory.
strSpp	string 3-letter code of the species.
runNo	the model run number.
rwtNo	the reweight number.
doMPD	logical: if TRUE, perform an MPD analysis.
N.reweight	the number of reweights to perform in the MPD analysis.
cvpro	CV process error added to CV observation error: $c_t = \sqrt{c_o^2 + c_p^2}$ ; if FALSE index CVs are reweighted using the standard deviation of normalized residuals.
mean.age	logical: if TRUE, use mean-age residuals to reweight the effective $N$ for the age composition data (see Francis 2011); if FALSE, reweight $N$ using $\Sigma(P(1 - P))/\Sigma(O - P)^2$ , where $O$ = observed proportions-at age and $P$ = predicted/fitted proportions-at-age.
doMCMC	logical: if TRUE, perform an MCMC analysis.
mcmc	number of MCMC iterations to perform.
mcsave	frequency of MCMC iterations to save.
ADargs	additional arguments for a call to Awatea.
verbose	logical: if TRUE, spew Awatea messages to the R console.
doMSY	logical: if TRUE, perform an MSY analysis.
msyMaxIter	maximum iterations for the MSY calculations.
msyTolConv	tolerance for convergence in the MSY calculations.
endStrat	maximum fishing mortality for the MSY analysis.
stepStrat	fishing mortality step size for the MSY analysis.
delim	character that delimits the components of the filename prefix.
awateaPath	system path on which the executable file Awatea.exe exists.
clean	logical: if TRUE, clean all Awatea files before running an MPD analysis.
...	additional arguments (not currently used for any purpose).

## Details

This function is primarily used to automate MPD reweightings and to perform the MSY calculations. MCMCs are better run from a command line console on a super computer as they generally require > 12 h to complete.

Once an MCMC has been created, the user can run various projections separately using  
`awatea -ind filename.ext -mceval`  
 on the command line.

## Text Output Files Produced

`Results.dat` contains an exhaustive listing of the maximum likelihood estimates for the model and derived parameters, a re-listing of some of the data and fixed parameters, and a listing of most of the prediction made. This includes numbers at age, fecundity, vulnerable biomass, survey trajectories, and so on. It is always placed in the same directory as the text input file. If there is an existing `Results.dat` file in that directory, it will be overwritten.

Other output files generated during the parameter estimation process include `Awatea.par` (contains the maximum likelihood estimates of the free parameters), `Awatea.cor` (shows standard deviation and correlations between the estimated parameters), and `Awatea.std` (standard deviation of the estimated parameters).

Be aware that using argument `ADargs=list("-nohess")` will shorten the time to perform an MPD (useful if there will be multiple re-weightings), but the output files `Awatea.cor` and `Awatea.std` will **not** be produced.

## Note

Be careful when setting the `clean` argument to `TRUE`. Some of the patterns for matching to `Awatea` files might inadvertently clean non-related files. The patterns are currently set to:  
`"^Awatea", "^admodel", "\\p.st$", "\\out$", "\\rpt$", "\\tmp$",`  
`"^variance$", "^results.dat$", "^likelihood.dat$"`

## Author(s)

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC.

## References

- Edwards, A.M., Starr, P.J., and Haigh, R. (2010, in revision). Stock assessment for Pacific Ocean Perch (*Sebastes alutus*) in Queen Charlotte Sound, British Columbia. *Canadian Science Advisory Secretariat, Research Document*.
- Edwards, A.M., Haigh, R., and Starr, P.J. (2011, in revision). Stock assessment and recovery potential assessment for Yellowmouth Rockfish (*Sebastes reedi*) along the Pacific coast of Canada. *Canadian Science Advisory Secretariat, Research Document*.
- Francis, R.I.C.C. (2011, in press) Data weighting in statistical fisheries stock assessment models. *Canadian Journal of Fisheries and Aquatic Sciences*.
- Hilborn, R., Maunder, M., Parma, A., Ernst, B. Payne, J., and Starr, P. (2003) Coleraine: a generalized age-structured stock assessment model. School of Aquatic and Fishery Sciences, University of Washington, 54 p.

**See Also**

[importRes](#), [readAD](#), [reweight](#), [runSweave](#)

---

runMCMC

*Wrapper to Function <runSweaveMCMC>*


---

**Description**

This small utility function simply provides a wrapper to the function `runSweaveMCMC` so that multiple documents can be built at once.

**Usage**

```
runMCMC(strSpp="XYZ", prefix=c("spp","area"),
        runs=7, rewt=0:6, cpue=FALSE, estM=TRUE,
        delim="-", mcs=1:1000)
```

**Arguments**

<code>strSpp</code>	three-letter code that identifies the species.
<code>prefix</code>	character vector of filename prefix components delimited by <code>delim</code> .
<code>runs</code>	the run number(s).
<code>rewt</code>	the reweight number(s).
<code>cpue</code>	logical: if TRUE, retain the CPUE figures in the Sweave file.
<code>estM</code>	logical: if TRUE the routine will assume natural mortality $M$ was estimated in the run, and the Sweave code uses the parameters "M_1" and "M_2"; if FALSE, the function removes these parameters from the Sweave file.
<code>delim</code>	character that delimits the components of the filename prefix.
<code>mcs</code>	vector of subsamples to select from the MCMC sample chain.

**Details**

Essentially loops through `runSweaveMCMC` using `(i in runs)` and `(j in rewt)`.

**Value**

Produces multiple Sweave documents and PDF files for MCMCs.

**Note**

Sweave files can be found in the library directory:

```
.../R/.../library/PBSawatea/snw/ymrrun-masterMCMC.Snw
.../R/.../library/PBSawatea/snw/ymrrun-masterMCMC20yrProjs.Snw
```

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[runSweaveMCMC](#), [runSweave](#), [runMPD](#)

---

runMPD

*Wrapper to Function <runSweave>*


---

**Description**

This small utility function simply provides a wrapper to the function [runSweave](#) so that multiple documents can be built at once.

**Usage**

```
runMPD(prefix=c("spp","area"), runs=1, rwts=0, ...)
```

**Arguments**

prefix	character vector of filename prefix components delimited by delim.
runs	the run number(s).
rwts	the reweight number(s).
...	arguments passed on to the function <a href="#">runSweave</a> .

**Details**

Essentially loops through [runSweave](#) using (i in runs) and (j in rwts).

**Value**

Produces multiple Sweave documents and PDF files.

**Note**

A recent Seave file called `ymrrun-master.Snw` can be found in the library directory:  
`.../R/.../library/PBSawatea/snw`

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[runSweave](#), [runSweaveMCMC](#), [runMCMC](#)

runSweave

*Run Customised Sweave Files for Awatea MPD Runs***Description**

Create and run customised Sweave files for Awatea MPD runs.

**Usage**

```
runSweave(wd=getwd(), cpue=FALSE, strSpp="XYZ",
  filename="spp-area-00.txt", runNo=1, rwtNo=0, running.awatea=0,
  Nsurvey=3, Snames=paste("Ser",1:Nsurvey,sep=""),
  SApos=rep(TRUE,Nsurvey), delim="-", debug=FALSE)
```

**Arguments**

wd	working directory in which Awatea input and master Sweave files occur.
cpue	logical: if TRUE the routine will leave calls to CPUE figures in the Sweave file, otherwise, they are removed.
strSpp	three-letter code that identifies the species.
filename	name of Awatea input file.
runNo	the run number that identifies a unique combination of input values.
rwtNo	the reweight number of the MPD run to build and collate figures.
running.awatea	numeric: if 0 load a previous .rep file, if 1 re-run Awatea
Nsurvey	number of surveys specified in the input file; controls image creation and place-holders.
Snames	survey names for figure labelling and image creation.
SApos	logical: surveys that have age composition data in the model.
delim	character that delimits the components of the filename prefix.
debug	logical: if TRUE, function only builds the target .Snw file but does not execute it.

**Details**

The values specified by the arguments (or derived variables) are directly substituted into the Sweave file wherever similarly named variables preceded by the @ symbol occur.

The cpue switch signals the removal of pieces of Sweave code if cpue=TRUE.

The Nsurvey argument essentially copies one Sweave line into Nsurvey lines for a set of lines identified by unique Sweave snippets.

**Value**

A customised Sweave file for runNo and rwtNo is created in a subdirectory (e.g., ./YMRrun/MPD.29.01/) from where it is run.

**Note**

A fairly recent Seave file called ymrrun-master.Snw can be found in the library directory:  
 .../R/.../library/PBSawatea/snw

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[runMPD](#), [runSweaveMCMC](#), [runADMB](#)

---

runSweaveMCMC

---

*Run Customised Sweave Files for Awatea MCMC Runs*


---

**Description**

Create and run customised Sweave files for Awatea MCMC runs.

**Usage**

```
runSweaveMCMC(wd=getwd(), cpue=FALSE, estM=TRUE, strSpp="XYZ",
  filename="spp-area-00.txt", runNo=1, rwtNo=0, running.awatea=0,
  delim="-", mcsb=1:1000)
```

**Arguments**

wd	working directory in which Awatea input and master Sweave files occur.
cpue	logical: if TRUE the routine will leave calls to CPUE figures in the Sweave file, otherwise, they are removed.
estM	logical: if TRUE the routine will assume natural mortality $M$ was estimated in the run, and the Sweave code uses the parameters "M_1" and "M_2"; if FALSE, the function removes these parameters from the Sweave file.
strSpp	three-letter code that identifies the species.
filename	name of Awatea input file.
runNo	the run number that identifies a unique combination of input values.
rwtNo	the reweight number of the MPD run to build and collate figures.
running.awatea	numeric: if 0 load a previous .rep file, if 1 re-run Awatea
delim	character that delimits the components of the filename prefix.
mcsb	vector of subsamples to select from the MCMC sample chain.

### Details

The values specified by the arguments (or derived variables) are directly substituted into the Sweave file wherever similarly named variables preceded by the @ symbol occur.

The cpue switch signals the removal of pieces of Sweave code if cpue=TRUE.

The estM switch signals the removal of "M\_1" and "M\_2" from the Sweave code if estM=FALSE.

### Value

A customised Sweave file for runNo and rwtNo is created in a subdirectory (e.g., ./YMRrun/MCMC.29.01/) from where it is run.

### Note

Seave files can be found in the library directory:

.../R/.../library/PBSawatea/snw/ymrrun-masterMCMC.Snw

.../R/.../library/PBSawatea/snw/ymrrun-masterMCMC20yrProjs.Snw

### Author(s)

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

### See Also

[runMCMC](#), [runSweave](#), [runADMB](#)

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 srFun

---

*Stock Recruitment Function*


---

### Description

Take a vector of spawners in year  $t-1$  and calculate recruits in year  $t$ .

### Usage

```
srFun(spawners, h = h.mpd, R0 = R0.mpd, B0 = B0.mpd)
```

### Arguments

spawners	a vector of spawners where either : each element corresponds to spawners in year $t-1$ or the vector calculates a single year but multiple MCMCs.
h	steepness parameter value.
R0	recruitment at $t = 0$ (virgin conditions).
B0	spawning biomass at virgin conditons.



**Details**

(AME wording) To input a vector of spawners in year  $t-1$  and calculate recruits in year  $t$ . Output for recruits is vector, each element corresponds to spawners the the year before, so will usually want to shift the output by 1 so that recruits in year  $t$  are based on spawners in year  $t-1$ .

Can also have each input as a vector (used when calculating a single year but multiple MCMCs, as in first year of projections is based on penultimate year of MCMC calculations).

**Value**

A vector of recruitments in year  $t$ .

**Note**

This function was originally a subfunction in `plt.mpdGraphs`.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[plt.mpdGraphs](#)

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stdRes.CA

---

*Calculate Standardised Residuals for Robust Normal Likelihood*


---

**Description**

Calculate the standardised residuals for Awatea's implementation of the Fournier *robustified* normal likelihood for proportions-at-length.

Based on PJS's summary of the CASAL document and ACH's change to length.

**Usage**

```
stdRes.CA(obj, trunc=3, myLab="Age Residuals", prt=TRUE)
```

**Arguments**

obj	scape/list object of Awatea's results file (.res).
trunc	maximum standardised residual; values greater than this are set to trunc.
myLab	general label for the output.
prt	logical: if TRUE, print the results.

**Value**

List object of standardised residuals.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[stdRes.index](#), [importCol2](#), [reweight](#)

---

stdRes.index

*Calculate Standardised Residuals for Abundance Indices*

---

**Description**

Calculate the standardised residuals for commercial and survey indices.

**Usage**

```
stdRes.index(obj, label=NULL, prt=TRUE)
```

**Arguments**

obj	data frame of observed and fitted index values from Awatea's results file (.res).
label	general label for the output.
prt	logical: if TRUE, print the results.

**Value**

Input data frame with additional column of standardised residuals.

**Author(s)**

Andrew Edwards, Pacific Biological Station, Fisheries and Oceans Canada, Nanaimo BC

**See Also**

[stdRes.CA](#), [importCol2](#), [reweight](#)

---

tabSAR	<i>Make Files of Reference Point Tables</i>
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---

**Description**

Generate comma-delimited, two-dimensional output tables (.csv) from objects reporting probabilities and times for DFO reference points and COSEWIC criteria.

**Usage**

```
tabSAR(models=paste("input-ymr", pad0(c(29, 30), 2), pad0(1, 2), sep="."),
        pnam="refProbs3Gen90", tnam=c("Ttab0.5", "Ttab0.8", "Ttab0.95"),
        cats=seq(0, 2500, 500), digits=2)
```

**Arguments**

models	names of binary system files that store the decision tables.
pnam	name of list object containing matrices of reference probabilities.
tnam	names of matrices reporting times to reach reference points/criteria.
cats	catch strategies (subset) to report in output tables.
digits	number of digits to retain after the decimal.

**Details**

At the moment, the reference point objects are a list of matrices for the probabilities and individual matrices for the times to reach the references. The latter should also be a list of matrices to parallel the probability object.

**Value**

No value is returned in R; however, comma-delimited text files (.csv) are created in the user's working directory. The probability tables have the suffix \_prob.csv while the time-to-target tables have the suffix \_targ.csv.

**Note**

The subset of the master decision tables in the stock assessment are primarily used in the SAR (Science Advisory Report) for managers.

**Author(s)**

Rowan Haigh, Pacific Biological Station, Fisheries and Oceans Canada, Nanimo BC

**See Also**

[refPoints](#), [findTarget](#),  
[cat](#), [round](#), [pad0](#), [show0](#)

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