

# Package ‘CRHMr’

March 19, 2019

**Type** Package

**Title** Pre- and post- processing for CRHM

**Version** 2.8.3

**Date** 2019-03-19

**Author** Kevin Shook, Centre for Hydrology, University of Saskatchewan

**Maintainer** Kevin Shook <kevin.shook@usask.ca>

**Description** The package is pronounced 'Krimmer' in honour of my wife's family.  
The CRHMr functions are used to create time series of forcing meteorological data for the Cold Regions Hydrological Modelling (CRHM) platform. Functions are also provided to read and aggregate CRHM output.

**Depends** R (>= 3.1)

**Imports** ggplot2(>= 2.0.0), lubridate(>= 1.3), plyr, reshape2, scales,  
signal, stringr(>= 1.0), zoo

**Suggests** tidyhydat, RSQLite, EcoHydRology,  
knitr,  
rmarkdown

**License** GPL-3

**LazyData** true

**Encoding** UTF-8

**URL** <https://github.com/CentreForHydrology/CRHMr>

**RoxygenNote** 6.1.1

**NeedsCompilation** no

**VignetteBuilder** knitr

## R topics documented:

CRHMr-package	4
aggDataframe	5
appendObs	7
assembleObs	8

automatePrj . . . . .	9
BadLake7376 . . . . .	10
basinWaterBalancePlot . . . . .	10
changeEatoRH . . . . .	11
changeRHtoEa . . . . .	12
convertDPtoRH . . . . .	13
createObsDataframe . . . . .	14
CRHM_summary . . . . .	15
CRHM_vars . . . . .	16
cumsumDataframe . . . . .	17
cumulativeDischargePlot . . . . .	18
cumulDailyWater . . . . .	19
dailyByHourlyP . . . . .	20
datetimeToDate . . . . .	21
dateToDatetime . . . . .	22
deDupe . . . . .	23
deleteSpikes . . . . .	24
deltaStorage . . . . .	25
distributeInst . . . . .	26
distributeMean . . . . .	27
distributeP . . . . .	28
distributeQli . . . . .	29
distributeQsi . . . . .	30
doubleMass . . . . .	32
emissivity . . . . .	33
extendObs . . . . .	33
findDupes . . . . .	34
findGaps . . . . .	35
findSpikes . . . . .	36
GMToffset . . . . .	37
hruGroupWaterSummary . . . . .	38
hydrograph . . . . .	39
hydroYear . . . . .	41
impute . . . . .	42
insertMissing . . . . .	43
interpolate . . . . .	44
keepGood . . . . .	45
list2df . . . . .	46
logAction . . . . .	48
longwave . . . . .	48
makeRegular . . . . .	49
maxObs . . . . .	50
minObs . . . . .	51
monthlyPrecipTotals . . . . .	53
monthlyQQplot . . . . .	54
nancumsum . . . . .	55
parseNums . . . . .	55
parseText . . . . .	56

PcpFiltPosTh . . . . .	57
phaseCorrect . . . . .	58
plotObs . . . . .	59
plotPrecipsBy Year . . . . .	60
plotTempsBy Year . . . . .	61
qair2ea . . . . .	62
qair2rh . . . . .	63
qqplotValues . . . . .	64
readCampbell . . . . .	65
readCLASSfile . . . . .	66
readDebugFile . . . . .	67
readExportFile . . . . .	68
readObsFile . . . . .	69
readOutputFile . . . . .	70
readOutputUnits . . . . .	71
readPrjDimensions . . . . .	72
readPrjHRUnames . . . . .	72
readPrjModuleParams . . . . .	73
readPrjNumVals . . . . .	74
readPrjOutputVariables . . . . .	75
readPrjParameters . . . . .	76
readPrjTextVals . . . . .	77
regress . . . . .	77
rh2vp . . . . .	79
ribbonPlotAirTemps . . . . .	80
runCRHM . . . . .	81
saturatedVP . . . . .	82
setPrjBasinName . . . . .	82
setPrjDates . . . . .	83
setPrjHRUnames . . . . .	84
setPrjObs . . . . .	85
setPrjOutputVariables . . . . .	86
setPrjParameters . . . . .	87
setPrjRunID . . . . .	88
simpleDailyWater . . . . .	89
simpleMaxSolar . . . . .	90
simpleRibbonPlot . . . . .	91
summariseObsFiles . . . . .	93
tMinMaxToHourly . . . . .	94
trimObs . . . . .	95
user . . . . .	96
vectorsToVelocity . . . . .	97
vp2rh . . . . .	98
weighingGauge-methods . . . . .	98
weighingGauge1 . . . . .	100
weighingGauge2 . . . . .	101
weighingGauge3 . . . . .	102
weighingGauge4 . . . . .	103

weighingGauge5 . . . . .	104
weighingGaugeInterval . . . . .	106
weighingGaugePlot . . . . .	107
wellHydrograph . . . . .	108
wetland-methods . . . . .	109
wetland_h_s . . . . .	109
wetland_pond_area . . . . .	110
wetland_pond_depth . . . . .	111
wetland_pond_size . . . . .	112
wg . . . . .	113
win.eol . . . . .	114
writeChangeLogFile . . . . .	114
writeCLASSfile . . . . .	116
writeObsFile . . . . .	116
yearlyPeaks . . . . .	117
zeroMissingPrecip . . . . .	118

<b>Index</b>	<b>120</b>
--------------	------------

---

CRHMr-package	<i>Contains functions to perform pre- and post- processing on data used with the Cold Regions Hydrological Modelling (CRHM) platform</i>
---------------	--

---

## Description

All data in **CRHMr**, whether model input or output must be stored in the standard type of data frame. The file input functions ([readObsFile](#) [readExportFile](#) [readOutputFile](#)) will automatically create standard **CRHMr** data frames. If you are reading data from another type of file using standard R functions, then you will have to force it to be in the standard format.

The first column of the data frame is labelled `datetime`, and as is the date and time stored as a `POSIXct` value, with the correctly-specified time zone. The only exception is for daily CRHM data or for aggregated values which use a daily or longer time step. In these cases the first column will have the appropriate name.

Because CRHM allows there to be many values of a variable, each variable will have a trailing number such as `.1`. This is added automatically when importing CRHM data. Most of the functions allow you to select a variable by its column number. In all cases, the column number does *NOT* include the datetime column.

To make your research more reproducible, each function writes information about what it did, including the date and time it was run, to a log file (`CRHMr.log`) in the default directory. It is suggested that you also run the function [user](#) when you first start to use **CRHMr** as it writes information about your computer to the log file. This may be helpful when trying to figure out bugs. Please send the output of your log file (including the user output) when reporting bugs.

The package contains functions to do the following:

1. Read in CRHM data into a **CRHMr** data frame. This includes data from `.obs` files (observation data used by CRHM), and model run outputs, either output automatically, or manually exported.

2. Manipulate obs data. Includes functions to plot the values, to find missing values, or datetimes, and to infill gaps by interpolation (linea or spline) and by imputation from other datasets.
3. Convert Ea values to RH and vice-versa. Interpolation and imputation require Ea values, as RH values depend on the air temperature. For safety, **CRHMr** functions do not permit both Ea and RH values in a data frame, as it would be impossible to know which was correct.
4. Write a data frame data to a CRHM obs file.
5. Execute CRHM. This allows CRHM to be run automatically, which is very useful for doing sensitivity analyses. There are functions to prepare a CRHM model to be executed, including setting the run start and end dates, and to run CRHM from inside R.
6. Examine output from CRHM runs. Includes functions to read in model output and to aggregate, summarize and plot the values.

## References

To cite **CRHMr** in publications, use the command `citation('CRHMr')` to get the current version of the citation.

The CRHM program is described in:

Pomeroy, John W, D M Gray, T Brown, N Hedstrom, W L Quinton, R J Granger, and S K Carey. 2007. "The Cold Regions Hydrological Model : A Platform for Basing Process Representation and Model Structure on Physical Evidence". *Hydrological Processes* 21 (19): 2650-2567.

The CRHM model may be downloaded from <http://www.usask.ca/hydrology/CRHM.php>.

---

aggDataframe

*Aggregates a data frame to a shorter time period*


---

## Description

CRHM data (observations and outputs) are generally produced at hourly time steps. This function aggregates CRHM data to daily, monthly or yearly values. The data can be aggregated by their maxima, minima, means, sums and any combination of these statistics.

## Usage

```
aggDataframe(CRHMrdataframe, columns = 1, period = "annual",
  funs = c("mean"), AggFilename = "", startMonth = 10,
  useSecondYear = TRUE, omitMissing = FALSE, logfile = "")
```

## Arguments

CRHMrdataframe	Required. A valid <b>CRHMr</b> data frame.
columns	The columns to be aggregated, not including the datetime. The default is the first column. This can be a vector, i.e. <code>c(1,2,3)</code> .
period	The period of aggregation. Must be one of 'hourly', 'daily', 'monthly', 'yearly' (or 'annual') or 'hydro'. Default is 'yearly'. Multiple-hour aggregation is not yet supported.

funcs	A character vector containing the function(s) for aggregation. The default is 'mean', but can also include 'min', 'max', 'sum', 'length', delta, and first. The function(s) will be applied to all of the specified columns
AggFilename	Optional. File name for the aggregated data.
startMonth	Optional. Starting month, to be used when aggregating by hydrological year.
useSecondYear	Optional. Logical. Should the hydrological year be based on the first or second calendar year. In other words would January 1, 2015 be the hydrological year 2014 or 2015? The default is TRUE (i.e., the hydrological year would be 2015). Note that the Campbell Scientific program SPLIT uses the first calendar year (i.e., the hydrological year would be 2014). To emulate this program, set useSecondYear to be FALSE.
omitMissing	Optional. If FALSE missing (i.e. NA_real_) values will be included in the aggregation, causing the aggregated values to also be NA_real_. If TRUE, then the missing values will be omitted from the calculations. This option is only used by the functions max, min, mean and sum.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

Returns a data frame with the aggregated values.

**Note**

The period of aggregation must be greater than the time step of the CRHM data. This function does NOT remove NA values before differencing.

**Author(s)**

Kevin Shook

**See Also**

[interpolate](#) [impute](#) [yearlyPeaks](#) [hydroYear](#)

**Examples**

```
Badlake.t.monthly <- aggDataframe(BadLake7376, period='monthly',
  columns=1, funcs=c('mean'))
Badlake.radiation.daily <- aggDataframe(BadLake7376, period='daily',
  columns=c(6,7,8), funcs=c('mean'))
```

---

appendObs*Appends two CRHM obs data frames*

---

### Description

This function joins two data frames of CRHM obs. The data frames must have the same number of columns, and the variables must be in the same order. The usual reason is for joining two data frames of differing periods. The primary and secondary data frames may be of any time periods. Where there are two values for a given datetime, the primary values are used. Rows of missing values at the beginning and end of the time series are deleted.

### Usage

```
appendObs(primaryObs, secondaryObs, trim = TRUE, quiet = TRUE,  
          logfile = "")
```

### Arguments

primaryObs	Required. The primary <b>CRHMr</b> data frame of obs values.
secondaryObs	Required. The secondary obs data frame. Note that both data frames must have the same time intervals.
trim	Optional. If set to TRUE (the default) then rows missing all values at the beginnings and/or ends of the obs data frames will be omitted.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns a data frame of values combined from the primaryObs and secondaryObs data frames. If unsuccessful, returns the value FALSE.

### Note

In addition to the usual notation in the logfile, this function also writes a separate file which summarises the new data frame. The summaries are also printed to the screen, if quiet=FALSE, the logfile also contains a complete listing of the source of each value in the infilled data frame. Each value is listed as being 'primary' (from the primaryObs data frame), 'secondary' (derived from the secondaryObs data frame) or 'NA' (missing).

### Author(s)

Kevin Shook

## Examples

```
## Not run:
broadview <- appendObs(broad2855, broad2856)
## End(Not run)
```

---

assembleObs

*Assembles two CRHM obs data frames*


---

## Description

This function joins two data frames of CRHM obs. The data frames can have different columns (it is assumed that they will), but should have some dates in common. Both data frames must have the same time step.

## Usage

```
assembleObs(obs1, obs2, quiet = TRUE, logfile = "")
```

## Arguments

obs1	Required. The first <b>CRHM</b> data frame of obs values.
obs2	Required. The second obs data frame.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If successful, returns the merges values of both data frames. Note that where the datetimes in the data frames are not the same, the merged values will be set to NA\_real\_. If unsuccessful, returns the value FALSE.

## Author(s)

Kevin Shook

## Examples

```
## Not run:
MSC.trh <- assembleObs(MSC.t, MSC.rh)
## End(Not run)
```



---

automatePrj	<i>Automates a CRHM .prj file</i>
-------------	-----------------------------------

---

## Description

Converts a CRHM .prj file to be able to be run automatically. It adds the settings Auto\_Run, Auto\_Exit and Log\_All to the end of the .prj file, if they are required.

## Usage

```
automatePrj(inputPrjFile = "", outputPrjFile = "", quiet = TRUE,  
            logfile = "")
```

## Arguments

inputPrjFile	Required. Name of the the CRHM .prj to be converted.
outputPrjFile	Optional. If omitted, the input .prj file will be overwritten.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If successful, returns TRUE. If unsuccessful, returns FALSE.

## Author(s)

Kevin Shook

## See Also

[runCRHM](#)

## Examples

```
## Not run:  
result <- automatePrj('Bad Lake 1974-1975.prj', 'Bad Lake 1974-1975 auto.prj')  
## End(Not run)
```

---

BadLake7376

*Observed meteorological data for Bad Lake*


---

### Description

A data frame containing meteorological data from Bad Lake, Saskatchewan, from January 1, 1973 to January 1, 1976. This data was collected by the Division of Hydrology at the University of Saskatchewan.

### Usage

```
BadLake7376
```

### Format

A data frame with 26280 rows and 9 variables (including the datetime):

**datetime** date and time as a POSIXct object

**t.1** air temperature (C)

**rh.1** relative humidity (percent)

**u.1** wind speed (m/s)

**SunAct.1** Actual daily sunshine hours, repeated for each hour of the day

**ppt.1** interval (hourly) precipitation (mm)

**Qsi.1** Incoming shortwave radiation ( $\text{W/m}^2$ )

**Qso.1** Outgoing shortwave radiation ( $\text{W/m}^2$ )

**Qn.1** Net radiation ( $\text{W/m}^2$ )

### Source

This data is also distributed with the example files of the CRHM program. CRHM can be obtained from <http://www.usask.ca/hydrology/CRHM.php>.

---

basinWaterBalancePlot

*Stacked plot of yearly water balance variables*


---

### Description

Creates a stacked bar plot of all water balance components computed for a basin. Basin inputs are plotted as positive values, outputs are plotted as negative.

### Usage

```
basinWaterBalancePlot(yearlyWater, negCols = "")
```

**Arguments**

yearlyWater	Required. A data frame of water balance components. The first column must be the year. Note that this function does not allow you to select columns - all columns will be plotted.
negCols	Optional. Columns to be plotted as negative values. If not specified (the default) the columns will be guessed from their names.

**Value**

If successful returns a **ggplot2** object showing stacked bars of the water balance components for each year. If unsuccessful returns FALSE.

**Examples**

```
## Not run:
# get daily water balance from CRHM output
daily <- simpleDailyWater(Bol84, prjFile = 'Bologna1984_02.prj',
  basinMean = TRUE, summarize = TRUE)

# get yearly values
daily.datetime <- dateToDatetime(daily, timezone='Etc/GMT+7')
yearly <- aggDataframe(daily.datetime, period='hydro', funs='sum')

# plot
p <- basinWaterBalancePlot(yearly)

## End(Not run)
```

---

changeEatoRH

*Changes Ea values to RH in a CRHM data frame*


---

**Description**

This function converts values of Ea to RH. Note that the specified obs data frame must contain both Ea and air temperatures. For safety, **CRHMr** does not permit values of both RH and Ea in a data frame. The names of the variables containing air temperature and ea values must be of the form t.x and ea.x, respectively, where x is an number, even if the column numbers are specified.

**Usage**

```
changeEatoRH(obs, t.cols = 1, ea.cols = 1, quiet = TRUE,
  logfile = "")
```

**Arguments**

obs	Required. A <b>CRHM</b> r data frame of observations.
t.cols	Optional. A vector containing the column numbers (not including the datetime) holding the air temperatures. If no columns are specified then the locations of the temperatures are guessed from the column names. The air temperatures must be in °C.
ea.cols	Optional. A vector containing the column numbers (not including the datetime) holding the vapour pressures. If no columns are specified then the locations of the vapour pressures are guessed from the column name. The ea values must be in kPa.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, the function returns the original data frame with the ea columns converted to RH values in percent. If unsuccessful, it returns the value FALSE.

**Author(s)**

Kevin Shook

**See Also**

[changeRHtoEa](#)

**Examples**

```
# First, convert RH values to Ea
BadLake7376.ea <- changeRHtoEa(BadLake7376)
# now convert Ea values back to RH
BadLake7376.rh <- changeEatoRH(BadLake7376.ea)
```

---

changeRHtoEa

---

*Changes RH values to Ea in a CRHM dataframe*


---

**Description**

This function converts values of RH to Ea. **CRHM**r does not allow interpolation or imputation of RH values, so you must convert RH values to Ea before infilling. Note that the specified obs dataframe must contain both RH and air temperatures. For safety, **CRHM**r does not permit values of both RH and Ea in a data frame. The names of the variables containing air temperature and RH values must be of the form t.x and rh.x, respectively, where x is a number, even if the column numbers are specified.

**Usage**

```
changeRHtoEa(obs, t.cols = 1, rh.cols = 1, quiet = TRUE,
  logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> data frame of observations.
t.cols	Optional. A vector containing the column numbers (not including the datetime) holding the air temperatures. If no columns are specified then the locations of the temperatures are guessed from the column names. The air temperatures must be in °C.
rh.cols	Optional. A vector containing the column numbers (not including the datetime) holding the relative humidities. If no columns are specified then the locations of the RH values are guessed from the column names.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, the function returns the original data frame with the RH columns converted to ea values in kPa. If unsuccessful, it returns the value FALSE.

**Author(s)**

Kevin Shook

**See Also**

[changeEatoRH](#)

**Examples**

```
BadLake7376.ea <- changeRHtoEa(BadLake7376)
```

---

 convertDPtoRH

*Converts dew point temperature and ambient temperature to RH*

---

**Description**

This is an internal **CRHMr** function and should normally not need to be called directly. Note that it does *NO* data checking.

**Usage**

```
convertDPtoRH(t, td)
```

**Arguments**

t	Required. Air temperature in °C.
td	Required. Dew point temperature in °C.

**Value**

Returns the RH as a percentage.

**Author(s)**

Kevin Shook

**References**

Lawrence, M. 2005. The relationship between relative humidity and the dewpoint temperature in moist air. American Meteorological Society: 225-233. <http://andrew.rsmas.miami.edu/bmcnoldy/Humidity.htm>

**Examples**

```
convertDPtoRH(10, 9)
```

---

createObsDataframe	<i>Creates an empty data fame for CRHM obs</i>
--------------------	--

---

**Description**

Creates a data fame to hold observations. All values are initially set to NA\_real\_. Once the data fame has been created, it may be filled from other data fames by imputation.

**Usage**

```
createObsDataframe(start.date, end.date, timestep = 1,
  variables = c("t", "rh", "u", "p"), reps = 1, timezone = "",
  logfile = "")
```

**Arguments**

start.date	Required. Starting date, in the format 'Y-m-d', 'd/m/Y', 'd-m-Y', 'd B Y', 'd b Y', 'd-B-Y', 'd-b-Y', 'B d, Y', or 'b d, Y', where 'Y' = year, 'm' = month number, 'B' = capitalized month name, 'b' = lower case month name, 'd' = day number.
end.date	Required. Ending date,, in the format 'd/m/Y', 'd-m-Y', 'd B Y', 'd b Y', 'd-B-Y', 'd-b-Y', 'B d, Y', or 'b d, Y', where 'Y' = year, 'm' = month number, 'B' = capitalized month name, 'b' = lower case month name, 'd' = day number.
timestep	Optional. Time interval between values in hours. The default value is 1 hour. The interval can be less than 1 hour.

variables	Optional. vector containing the names of variables to be created. The default values are 't', 'rh', 'u', and 'p'.
reps	Optional. The number of repetitions for each value. The default is 1.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Details

The names for the variables in the data fame are derived from the specified variables, and the number of reps. For example if the default variable names are used, with 2 reps, then the variables created will be t.1, t.2, rh.1, rh.2, u.1, u.2, p.1, and p.2.

### Value

If successful, the function returns a CRHM data fame with the specified number of each variable, with values set to be NA. If unsuccessful, returns the value FALSE.

### Author(s)

Kevin Shook

### See Also

[impute](#)

### Examples

```
vars <- c('t','rh','u')
stoon <- createObsDataframe('1/1/1960', '31/12/1960', variables=vars, reps=2, timezone='etc/GMT+6')
```

---

CRHM\_summary

*Summarizes a **CRHM**r data frame*

---

### Description

Summarizes the values in a **CRHM**r data frame. This is an internal **CRHM**r function and should *never* need to be called directly.

**Usage**

```
CRHM_summary(CRHMdataframe)
```

**Arguments**

CRHMdataframe    Required. Name of the **CRHM** data frame to be summarized.

**Value**

If successful, returns a data frame containing the number of rows, complete rows, starting and ending datetimes, and the variables in the specified data frame. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**Examples**

```
## Not run:
summary <- CRHMsummary(BadLake7376)
## End(Not run)
```

---

CRHM\_vars

*CRHM variables*


---

**Description**

A data frame containing information about all of the variables in CRHM that are likely to be output from a model run.

**Usage**

```
CRHM_vars
```

**Format**

A data frame with 370 rows and 5 variables:

**name** name of the variable

**type** type of variable (its dimensions)

**description** description of the variable

**units** variable units

**end\_of\_day** TRUE if the value is only output at the end of the day (00:00), otherwise FALSE

**Source**

The variable information was abstracted from the CRHM c++ source code.



---

cumsumDataframe	<i>Cumulative sums for a data frame</i>
-----------------	---

---

## Description

Finds the cumulative sum of all columns in a data frame. Note that all columns must be numeric - date and datetime variables cannot be summed.

## Usage

```
cumsumDataframe(df)
```

## Arguments

df	Required. Data frame to be summed.
----	------------------------------------

## Value

Returns a dataframe containing the cumulative sums of all columns in the original data frame.

## Note

This function is used by other **CRHMr** functions, so it does *NO* parameter testing.

## Author(s)

Kevin Shook

## See Also

[cumulDailyWater](#)

## Examples

```
## Not run:  
cumul <- cumsumDataframe(modelOutput)  
## End(Not run)
```

---

cumulativeDischargePlot

*Plots cumulative CRHM and/or WSC flows*


---

## Description

Creates a **ggplot** object of annual cumulative flows from a data frame of CRHM output and/or WSC daily flows. The WSC flows are obtained from the **tidyhydat** package. The WSC data is truncated so that it only includes the range of the CRHM data. If more than a single year of data is specified, then the plot will be faceted by year.

## Usage

```
cumulativeDischargePlot(CRHMflows = NULL, CRHMflowsLabel = "",
  CRHMflowCol = 1, WSCdailyFlowsID = NULL, WSCdailyFlowsLabel = "",
  facetCols = 3, quiet = TRUE, hydat_path = NULL)
```

## Arguments

CRHMflows	Optional. A data frame of CRHM modelled flows. The flows must be in m <sup>3</sup> /s.
CRHMflowsLabel	Optional. Labels for the CRHM data. If not specified, and CRHM data are plotted, then the name of the CRHM variable will be used.
CRHMflowCol	Optional. Column containing the flowrates, not including the datetime. Default is 1.
WSCdailyFlowsID	Optional. If NULL (the default) the WSC daily flows will not be plotted. If a WSC station ID is specified, e.g. WSCdailyFlowsID = "05CC001", then the daily flows will be obtained from <b>tidyhydat</b> and plotted.
WSCdailyFlowsLabel	Optional. Labels for the daily flows. If not specified, then the WSC station number will be used, followed by 'daily'.
facetCols	Optional. Number of columns to wrap the facets (if they are used). Setting a value less than 1 will cause an error.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
hydat_path	Optional. Path to the HYDAT database. This can usually be left blank.

## Value

If successful, returns a **ggplot2** object containing the cumulative discharge plot. If unsuccessful, the value FALSE will be returned.

**Author(s)**

Kevin Shook

**See Also**[hydrograph](#)**Examples**

```
## Not run:
p <- cumulativeDischargePlot(CRHMflows = crhm_output, WSCdailyFlowsID = "05FA025",
  facetCols=2, quiet=FALSE)
## End(Not run)
```

---

cumulDailyWater	<i>Cumulative water fluxes</i>
-----------------	--------------------------------

---

**Description**

Accumulates water fluxes by year or water year. The values are cumulative for each day.

**Usage**

```
cumulDailyWater(dailyWaterVals, waterYear = FALSE, logfile = "")
```

**Arguments**

dailyWaterVals	Required. CRHM daily water fluxes as calculated by simpleDailyWater.
waterYear	Optional. If FALSE the fluxes are accumulated by calendar year. If TRUE the fluxes are accumulated by the water year, based on the first month of the model output.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a data frame containing the date and the cumulative values of each variable. If unsuccessful, returns FALSE.

**Note**

The output from this function can be plotted using simpleRibbonPlot.

**Author(s)**

Kevin Shook

**See Also**[simpleDailyWater](#) [simpleRibbonPlot](#)

**Examples**

```
## Not run:
waterYearCumul <- cumulDailyWater(CRHM_daily, waterYear=TRUE)

## End(Not run)
```

---

dailyByHourlyP	<i>Distributes daily precipitation according to hourly values</i>
----------------	---

---

**Description**

Used to distribute daily precipitation totals, according to hourly values. Typically, this is used when the hourly values are of lower quality than the daily values, such as when they are imputed from another location.

**Usage**

```
dailyByHourlyP(dailyObs, dailyPcol = 0, hourlyObs, hourlyPcol = 0,
  missingRatio = 1, quiet = TRUE, logfile = "")
```

**Arguments**

dailyObs	obs Required. A <b>CRHM</b> r data frame containing daily precipitation values.
dailyPcol	Optional. The number of the column containing the daily precipitation. If not specified, the column will be guessed from the variable name. Note that the variable name <b>MUST</b> include the letter 'p', even if the number is specified.
hourlyObs	obs Required. A <b>CRHM</b> r data frame containing hourly precipitation values.
hourlyPcol	Optional. The number of the column containing the hourly precipitation. If not specified, the column will be guessed from the variable name. Note that the variable name <b>MUST</b> include the letter 'p', even if the number is specified.
missingRatio	Optional. The value of the ratio used to multiply the hourly precipitation when daily precipitation data are not available. The default is 1. If you want to omit these values, use NA_real_.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a modified version of the obs data frame, consisting of the original hourly precipitation data, and the adjusted\_precip.

**Author(s)**

Kevin Shook

**Examples**

```
## Not run: hourly <- dailyByHourlyP(dailyObs = daily, hourlyObs = hourly)
```

---

datetimeToDate	<i>Converts an R datetime to a date</i>
----------------	---

---

**Description**

This is used when a **CRHMr** date is stored as a datetime (POSIXct) variable.

**Usage**

```
datetimeToDate(dataframe, logfile = "")
```

**Arguments**

dataframe	Required. A data frame. The first name of the column must be a POSIXct variable, which can be converted to a standard R date.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a data frame where the first column is a standard **CRHMr** date. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**See Also**

[dateToDatetime](#)

**Examples**

```
Badlake.radiation.daily <- aggDataframe(BadLake7376, period='daily',
columns=c(6,7,8), funs=c('mean'))
BadLake.datetime <- dateToDatetime(Badlake.radiation.daily, timezone='CST')
BadLake.date <- datetimeToDate(BadLake.datetime)
```

---

dateToDatetime	<i>Converts an R date to a datetime</i>
----------------	---

---

### Description

The `aggDataframe` function produces non-standard data frames, which have dates rather than datetimes as their first column. This function will convert a data frame with dates in the first column to have datetimes with the time being set to a constant value.

### Usage

```
dateToDatetime(dataframe, hour = 0, timezone = "", logfile = "")
```

### Arguments

<code>dataframe</code>	Required. A data frame. The first name of the column must be 'date', and it must be standard R date.
<code>hour</code>	Optional. The hour to be used for the datetime values. Default is 0.
<code>timezone</code>	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
<code>logfile</code>	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns a data frame where the first column is a standard **CRHMr** datetime. If unsuccessful, returns the value FALSE.

### Author(s)

Kevin Shook

### See Also

[aggDataframe datetimeToDate](#)

### Examples

```
Badlake.radiation.daily <- aggDataframe(BadLake7376, period='daily',
columns=c(6,7,8), funs=c('mean'))
BadLake.datetime <- dateToDatetime(Badlake.radiation.daily, timezone='CST')
```

---

deDupe*Removes duplicated datetimes in obs data frame*

---

### Description

Removes duplicated datetime values. Many time series, especially from Environment Canada, may contain duplicated datetimes. This function replaces the duplicated values. It is important to use this function before interpolating or imputing values, and especially before writing the data frame to an obs file.

### Usage

```
deDupe(obs, action = "mean", quiet = TRUE, logfile = "")
```

### Arguments

obs	Required. A <b>CRHMr</b> data frame containing the obs values.
action	Optional. The action used to replace the duplicate values. Must be one of 'min', 'max', 'mean', 'skip', 'delete', 'split' or 'second'. Default is 'mean'.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If there are no duplicates, returns 'No duplicates'. If duplicates exist, and are successfully removed, the de-duplicated data frame is returned. If the de-duplication is unsuccessful, then the value FALSE is returned.

### Note

If action='min', 'max', or 'mean', the action function is applied to all of the values for each duplicate datetime. If action='skip' or 'delete', then the values of the duplicate datetimes are deleted. If action='split' then the original values are kept, and the duplicate values are written to an obs file. The name of the obs file is the name of the obs variable followed by '\_dupes.obs'. If action='second', then the second duplicate values are used. This can be useful when dealing with duplicates caused by daylight savings time.

### Author(s)

Kevin Shook

### See Also

[findDupes](#)

**Examples**

```
BadLake.deduped <- deDupe(BadLake7376, action='mean')
```

---

deleteSpikes	<i>Deletes spikes in an obs data frame</i>
--------------	--

---

**Description**

Finds spikes, and sets their values to be NA\_real\_.

**Usage**

```
deleteSpikes(obs, colnum = 1, threshold = 0, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> obs data frame.
colnum	Optional. The number of the column to test for spikes, not including the date-time.
threshold	Required. The threshold for the <i>change</i> in the observed values. The threshold is actually a rate, i.e. the change per unit time. So if you are looking at air temperature, and the threshold is set to 5, then any change in temperature of +/- 5 degrees in one time interval will be considered to be a spike.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a data frame consisting of the datetime and the original obs values, where all of the spike values have been set to be NA\_real\_. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**See Also**

[findSpikes](#)

**Examples**

```
BadLake <- BadLake7376
# finds all windspeeds which change by more than 10 m/s per interval,
# and sets their values to NA_real_
BadLake$u.nospikes <- deleteSpikes(BadLake7376, 3, 10)
```



---

deltaStorage	<i>Calculates changes in storage over specified intervals</i>
--------------	---

---

## Description

Calculates changes in storage over specified intervals

## Usage

```
deltaStorage(CRHMdataframe = NULL, columns = 1, period = "yearly",
             startMonth = 10, useSecondYear = TRUE, logfile = "")
```

## Arguments

CRHMdataframe	Required. A standard CRHM data frame.
columns	Required. The columns to be differenced, not including the datetime. The default is the first column. The columns can be a single number or a vector, i.e. <code>c(1,2,3)</code> .
period	Required. The period of aggregation. Must be one of 'hourly', 'daily', 'monthly', 'yearly' (or 'annual') or 'hydro'. Default is 'yearly'.
startMonth	Optional. Starting month, to be used when differencing by hydrological year. Default is 10 (October).
useSecondYear	Optional. Logical. Should the hydrological year be based on the first or second calendar year. In other words would January 1, 2015 be the hydrological year 2014 or 2015? The default is TRUE (i.e., the hydrological year would be 2015).
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Note

The period of differencing must be greater than the time step of the CRHM data. This function does NOT remove NA values before aggregation. Use `na.omit` or one of the infilling functions ([interpolate](#) or [impute](#)) if you want to remove missing values.

## Author(s)

Kevin Shook

## See Also

[simpleDailyWater](#)

## Examples

```
## Not run: delta <- deltaStorage(crhmm, c(23,24,25))
```

---

distributeInst	<i>Distribute instantaneous values</i>
----------------	--

---

### Description

Distributes instantaneous values to a shorter time interval. The missing datetimes are inserted and then the values are interpolated. This function is typically used to downscale obs values such as t, ea, and u.

### Usage

```
distributeInst(obs, obsCols = "all", timeStep = 0,
  interpolationMethod = "linear", maxLength = 5, quiet = TRUE,
  logfile = "")
```

### Arguments

obs	Required. The <b>CRHMr</b> data frame of obs values.
obsCols	Optional. A vector containing the columns to be imputed in the obs data frame, not including the datetime. The default 'all' specifies all columns.
timeStep	Required. The time step (in hours) for the interpolated values. This value must be smaller than the time step in the original time series.
interpolationMethod	Optional. A vector containing the methods to be used for interpolation for each of the variables. Currently supported methods are 'linear' and 'spline'. The default is to use linear interpolation. If fewer methods than columns are specified, the methods are recycled.
maxLength	Optional. The maximum gap length to be interpolated. Defaults to 5 time steps.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not use

### Value

If successful, returns a dataframe of the selected columns interpolated to the specified time step. If unsuccessful, returns FALSE.

### Author(s)

Kevin Shook

### See Also

[distributeMean](#)

**Examples**

```
## Not run:
hourly_vals <- distributeInst(vegreville, c(1,2,3), timeStep = 1)
## End(Not run)
```

---

distributeMean	<i>Distribute mean values</i>
----------------	-------------------------------

---

**Description**

Distributes mean values to a shorter time interval. The missing datetimes are inserted and then the values are repeated. This function is typically used to downscale obs values such as t, ea, and u.

**Usage**

```
distributeMean(obs, obsCols = "all", timeStep = 0, maxLength = 5,
  quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. The <b>CRHMr</b> data frame of obs values.
obsCols	Optional. A vector containing the columns to be imputed in the obs data frame, not including the datetime. The default 'all' specifies all columns.
timeStep	Required. The time step (in hours) for the interpolated values. This value must be smaller than the time step in the original time series.
maxLength	Optional. The maximum gap length to be interpolated. Defaults to 5 time steps.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not use

**Value**

If successful, returns a dataframe of the selected columns interpolated to the specified time step. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[distributeInst](#)

## Examples

```
## Not run:
hourly_vals <- distributeMean(vegreville, c(1,2,3), timeStep = 1)
## End(Not run)
```

---

distributeP

*Distributes total precipitation over intervals*


---

## Description

Distributes total precipitation (often daily) evenly over all of the applicable intervals. The function has 2 uses. The primary use is to distribute total precipitation when the precipitation time step is greater than the time step of the obs dataframe. For example, there may be 3-hour total precipitation within an obs dataframe holding hourly values for all other variables. The secondary use is to disaggregate precipitation to a shorter time step. Normally this can be done inside CRHM by putting the data in a separate obs file. However, if you want to do this in this function, you must specify the new time step.

## Usage

```
distributeP(obs, p.cols = 0, timestep = 0, quiet = TRUE,
  logfile = "")
```

## Arguments

obs	Required. The <b>CRHM</b> r data frame of obs values.
p.cols	Optional. The number of the column(s) to be used. If not specified, the column will be guessed from the variable name. Note that the variable name <b>MUST</b> include the letter 'p', even if the number is specified.
timestep	Optional. The new time step in hours. If zero (the default) the time step is not used. The time step is only used when disaggregating precipitation to a shorter time step. Note that the specified time step must be shorter than that used in the original data frame.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If successful, returns a modified version of the obs data frame. If no time step is specified, the missing precipitation values in the specified data frame are replaced by average precipitation values. If a time step is specified, a new data frame containing **ONLY** the datetime and the distributed precipitation will be returned. If unsuccessful, returns the value FALSE.

**Note**

This function is potentially destructive. If you have a missing precipitation total, then the next precipitation total will be distributed over a longer time period. For example, you may have 3-hour precipitation data, normally at 03:00, 06:00, 09:00 etc., to be distributed hourly. If the 06:00 data are missing, then the 09:00 precipitation will be distributed over the interval from 04:00 to 09:00.

**Author(s)**

Kevin Shook

**See Also**

[impute](#) [interpolate](#) [distributeInst](#) [distributeQli](#) [distributeQsi](#)

**Examples**

```
## Not run:
distributed <- distributeP(obs, 3)
## End(Not run)
```

---

distributeQli

*Distributes incoming longwave radiation to shorter time interval*


---

**Description**

Downscales incoming longwave radiation to shorter time intervals, based on the air temperatures, using the Stefan-Boltzmann law  $Q_{li} = \epsilon \sigma T^4$ . The atmospheric emissivity ( $\epsilon$ ) is calculated from the longwave radiation and the temperature at the original time intervals. The air temperatures are interpolated to the shorter intervals (if required), and are then used to calculate the incoming longwave, using the longer-interval emissivities. Finally, the downscaled longwave radiation is adjusted so that its mean values over the original time intervals are the same as the original values.

**Usage**

```
distributeQli(QliObs, QliColnum = 1, tObs, tColnum = 1, timeStep = 1,
  interpolationMethod = "linear", maxlength = 0, quiet = TRUE,
  logfile = "")
```

**Arguments**

QliObs	Required. A <b>CRHMr</b> obs data frame containing incoming longwave radiation in in W/m <sup>2</sup> .
QliColnum	Optional. The number of the columns (not including the datetime) containing the incoming longwave values. Default is 1.
tObs	Required. A <b>CRHMr</b> obs data frame containing air temperatures. The values may be in K or °C (the function will detect and adjust).

tColnum	Optional. The number of the columns (not including the datetime) containing the air temperature values. Default is 1.
timeStep	Optional. The number of hours to be used as the downscaled time step. Default is 1.
interpolationMethod	Optional. The method to be used for interpolating the air temperatures (if required). Currently supported methods are 'linear' and 'spline'. The default is to use linear interpolation.
maxlength	Optional. The maximum gap length to be interpolated. If not specified (the default), then the maximum gap length permitted is twice the time interval of the air temperatures.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used

**Value**

If successful, returns an obs data frame containing the downscaled incoming longwave radiation (Qli) in  $\text{W/m}^2$ . If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**See Also**

[emissivity longwave interpolate](#)

**Examples**

```
## Not run:
hourlyLongwave <- distributeQli(LWobs, 1, tobs, 1)
## End(Not run)
```

---

distributeQsi

*Distributes incoming shortwave radiation to shorter time interval*


---

**Description**

Downscales incoming shortwave radiation to shorter time intervals, based on the extra-terrestrial radiation. The mean atmospheric transmittance is computed from the ratio of the incoming shortwave to the extra-terrestrial radiation over the original time intervals. The extra-terrestrial radiation over the shorter time intervals is then multiplied by the mean transmittance values to calculate the shorter interval incoming shortwave radiation.

**Usage**

```
distributeQsi(QsiObs, QsiColnum = 1, latitude = "",
  sunTimeOffset = 2, timeStep = 1, solarMethod = "simpleMaxSolar",
  details = FALSE, quiet = TRUE, logfile = "")
```

**Arguments**

QsiObs	Required. A <b>CRHMR</b> obs data frame containing incoming shortwave radiation in in $W/m^2$ .
QsiColnum	Optional. The number of the columns (not including the datetime) containing the incoming shortwave values. Default is 1.
latitude	Required. The latitude for which values are to be calculated.
sunTimeOffset	Optional. The offset (in hours) is added to the solar time to convert it to local time. The default value '2' shifts the daily peak to occur at 2pm.
timeStep	Optional. The number of hours to be used as the downscaled time step. Default is 1.
solarMethod	The method to be used for calculating the extra-terrestrial radiation. The default method is 'simpleMaxSolar'. Note that this method is only valid for latitudes between 49 and 55°N. The other supported method is 'PotSolarInst', which requires the package <b>EcoHydRology</b> to be installed.
details	Optional. If TRUE then the calculated atmospheric transmittance and extra-terrestrial radiation (So_h_W) will be returned along with the downscaled Qsi. If FALSE (the default) then only the Qsi values and the datetime are returned.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used

**Value**

If successful, returns an obs data frame containing the downscaled incoming shortwave radiation (Qsi) in  $W/m^2$  and, optionally, the transmittance and extra-terrestrial solar radiation (So\_h\_W). If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**See Also**

[simpleMaxSolar](#) [distributeQli](#)

**Examples**

```
## Not run:
hourlyShortwave <- distributeQsi(QsiObs=SWobs, QsiColnum=1, latitude=51.1, timeStep=1)
## End(Not run)
```

---

doubleMass	<i>Calculates double mass values</i>
------------	--------------------------------------

---

### Description

Calculates cumulative sums for two variables and optionally returns a plot. Only values coinciding in time are used for the cumulative sums.

### Usage

```
doubleMass(primaryObs, primaryCol = 1, secondaryObs, secondaryCol = 1,
  plot = TRUE, logfile = "")
```

### Arguments

primaryObs	Required. The primary <b>CRHMr</b> data frame of obs values.
primaryCol	Optional. The column number to be used, not including the datetime. If not specified, defaults to the first column.
secondaryObs	Required. The secondary <b>CRHMr</b> data frame of obs values.
secondaryCol	Optional. The column number to be used, not including the datetime. If not specified, defaults to the first column.
plot	Optional. Should a plot be created? Default is FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If `plot = TRUE`, then a **ggplot2** object will be returned, with the `primaryObs` values plotted on the x-axis and the `secondaryObs` values plotted on the y-axis. Otherwise, a dataframe is returned with the cumulative values of each of the variables.

### Author(s)

Kevin Shook

### Examples

```
## Not run:
p <- doubleMass(primaryObs=CamrosePPT, secondaryObs=WetaskiwinPPT, plot=TRUE)
## End(Not run)
```



---

emissivity	<i>Calculates atmospheric emissivity</i>
------------	--

---

### Description

Calculates atmospheric emissivity from specified longwave radiation and air temperatures. This is an internal **CRHMr** function and should *never* need to be called directly, as it does *NO* data checking.

### Usage

```
emissivity(longwave, tK)
```

### Arguments

longwave	Required. Longwave radiation in W/m <sup>2</sup> .
tK	Required. Air temperatures in K.

### Value

Returns the atmospheric emissivity.

### Author(s)

Kevin Shook

### See Also

[distributeQli longwave](#)

### Examples

```
emissivity(100, 293)
```

---

extendObs	<i>Extends an obs data frame</i>
-----------	----------------------------------

---

### Description

This function extends a **CRHMr** obs data frame to include the specified dates. All added values are set to be NA\_real\_. The purpose of this function is to create locations for data that can be infilled or imputed from another data set.

### Usage

```
extendObs(obs, startDate = "", endDate = "", timezone = "",
  quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. The <b>CRHMr</b> obs data frame to be extended.
startDate	Optional. The starting date for the obs data frame. Can either be a year as a number (e.g. 1995) or a date in 'Y-m-d' format, i.e. '1995-06-01'. Either the startDate or the endDate, or both, must be specified.
endDate	Optional. The ending date for the obs data frame. Can either be a year as a number (e.g. 1995) or a date in 'Y-m-d' format, i.e. '1995-07-01'.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a modified version of the obs data frame. The missing values in the specified data frame are replaced by NA\_real\_ values. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**Examples**

```
BadLake7378 <- extendObs(BadLake7376, endDate='1978-01-01', timezone='CST')
```

---

findDuples

*Finds duplicated datetimes in obs dataframe*


---

**Description**

Finds duplicate datetime values. All of the duplicate values are written to a .csv file. Many time series, especially from Environment Canada, may contain duplicate datetimes.

**Usage**

```
findDuples(obs, dupefile = "", quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHM</b> r data frame containing obs values.
dupefile	Optional. The name of the output file. If omitted the dupe file will be the name of the obs data frame, followed by ‘_dupedatetimes.csv’.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If there are duplicate datetimes, returns TRUE. If there are no duplicates, returns FALSE.

**Note**

If quiet=FALSE, the function gives a list of the duplicate datetimes. Use this function before removing duplicates and/or interpolation and imputation. Most importantly, use this function before writing values to an obs file. Duplicate datetime values in an obs file will prevent CRHM from executing.

**Author(s)**

Kevin Shook

**See Also**

[findGaps](#) [interpolate](#) [impute](#) [writeObsFile](#)

**Examples**

```
findDupes(BadLake7376, quiet=FALSE)
```

---

findGaps	<i>Finds gaps in obs variables</i>
----------	------------------------------------

---

**Description**

Finds successive missing values in an obs data frame. This function is useful to show where data are missing, before you use the [interpolate](#) or [impute](#) functions. All of the gaps in each variable are written to a .csv file.

**Usage**

```
findGaps(obs, gapfile = "", minlength = 1, quiet = TRUE,
         logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> data frame containing obs values.
gapfile	Optional. The name of the output file. If omitted the output file will be the name of the obs data frame, followed by ‘_gaps.csv’.
minlength	The minimum gap length included in the analysis (in time steps). Default is 1 time step.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, and there are gaps, returns TRUE. If successful, and there are no gaps, returns ‘No gaps’.  
If unsuccessful, returns FALSE.

**Note**

If quiet=FALSE, the functions gives a summary of the gaps (number of gaps, their total length) for each year.

**Author(s)**

Kevin Shook

**See Also**

[insertMissing](#) [interpolate](#) [impute](#)

**Examples**

```
findGaps(BadLake7376, quiet=FALSE)
```

---

findSpikes

---

*Finds spikes in an obs data frame*


---

**Description**

Finds spikes (short bursts of very large or small values) in the specified time variable.

**Usage**

```
findSpikes(obs, colnum = 1, threshold = 0, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> obs data frame.
colnum	Optional. The number of the column to test for spikes, not including the <code>datetime</code> .
threshold	Required. The threshold for the <i>change</i> in the observed values. The threshold is actually a rate, i.e. the change per unit time. So if you are looking at air temperature, and the threshold is set to 5, then any change in temperature of +/- 5 degrees in one time interval will be considered to be a spike.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful and there are no spikes, returns `0`. If there are spikes, returns their `datetime` values. If unsuccessful returns `FALSE`.

**Note**

If `quiet=FALSE`, the function displays a list of the datetimes of the spikes.

**Author(s)**

Kevin Shook

**See Also**

[deleteSpikes](#) [findGaps](#) [findDupes](#)

**Examples**

```
# check for spikes in wind speed - look for changes greater than 10 m/s per interval
findSpikes(BadLake7376,3, 10)
```

---

GMToffset

*Finds offset in hours from UTC for given time zone.*


---

**Description**

Finds offset in hours from UTC for given time zone.

**Usage**

```
GMToffset(timezone)
```

**Arguments**

timezone	Required. A string containing the time zone. Works with designated time zones such as 'CST' and 'MST' for Central Standard or Mountain Standard time or offset time zones such as 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. Does NOT work with location time zones like 'America/Regina'.
----------	---

**Value**

Returns the offset in hours as a numeric value.

**Note**

This code was based on the function `GMTOffsetFromTz` in the package **oce**, but works for more time zones and returns a negative value in the western hemisphere.

**Author(s)**

Kevin Shook

**Examples**

```
GMTOffset('MST')
GMTOffset('etc/GMT+6')
```

---

`hruGroupWaterSummary`    *Summarize daily CRHM water outputs by HRU group*

---

**Description**

Calculates daily totals of water storages and fluxes for groups of HRUs, for models *without* sub-basins.

**Usage**

```
hruGroupWaterSummary(dailyWater, vars = "all", prjFile, HRUgroups,
  basinMean, quiet = TRUE, logfile = "")
```

**Arguments**

<code>dailyWater</code>	Required. The output from the <code>simpleDailyWater</code> command.
<code>vars</code>	Optional. Variable column numbers to be used (not including the datetime). The default value 'all' selects all columns.
<code>prjFile</code>	Required. The CRHM model .prj file.
<code>HRUgroups</code>	Required. This can be either a vector or list. If a vector, it contains strings used to group the selected HRU names in the .prj file. If a list, it contains vectors of the HRU numbers to be grouped. See the examples for both usages.
<code>basinMean</code>	Required. This refers to how the daily water values were calculated by <code>simpleDailyWater</code> . If TRUE then the aggregated water depths/volumes are computed over the whole basin. If FALSE then the the aggregated water depths/volumes are computed over area of the HRU groups.
<code>quiet</code>	Optional. Suppresses display of messages, except for errors. The default is TRUE. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet = FALSE</code> .
<code>logfile</code>	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful returns a data frame containing each variable summarized by each HRU group. The names of the variables are in the form: `hruGroup.variable.group_total` or `hruGroup.variable.basin_total`. If unsuccessful, returns the value FALSE.

**Examples**

```
## Not run:
# First, calculate daily water values. In these examples, the water is
# in mm per HRU

# b <- readOutputFile('Bologna1984_30y_output.txt', timezone='MST')
# get daily values without aggregation; values calculated w.r.t the HRU area
dailyHRU <- simpleDailyWater(b, prjFile='Bologna1984_02.prj', basinMean=FALSE)

# summarize by groups of HRUs with similar names
daily_group <- hruGroupSummary(daily, prjFile = 'Bologna1984_02.prj',
HRUgroups=c('firn','ice','bl'), basinMean=FALSE)

# summarize by named groups of HRUs with specified numbers
groups <- list(group1=c(1,2,3), group2=c(4,5,6))
daily_group <- hruGroupSummary(daily, prjFile = 'Bologna1984_02.prj',
HRUgroups=groups, basinMean=FALSE)

## End(Not run)
```

hydrograph

*Plots hydrograph of CRHM output and/or WSC daily and/or peak flows*

**Description**

Creates a **ggplot** hydrograph from any of CRHM flows, WSC daily flows and/or WSC peak flows.

**Usage**

```
hydrograph(CRHMflows = NULL, CRHMflowsLabels = "", CRHMcols = NULL,
CRHMDaily = FALSE, WSCdailyFlowsID = NULL, WSCdailyFlowsLabel = "",
WSCpeakFlowsID = NULL, WSCpeakFlowsLabel = "",
forceMissingPeakTimes = FALSE, commonTime = FALSE,
fakeDates = FALSE, quiet = TRUE, hydat_path = NULL)
```

**Arguments**

CRHMflows	Optional. A data frame of CRHM modelled flows. The flows must be in m <sup>3</sup> /s. Remember to divide interval flows by their length in seconds, i.e. 3600 if hourly.
CRHMflowsLabels	Optional. Labels for the CRHM data. If not specified, and CRHM data are plotted, then the name(s) of the CRHM variables will be used.

CRHMcols	Required. Column(s) containing the flowrates. As always, the numbers do not include the datetime.
CRHMDaily	Optional. Should CRHM flows be plotted as daily values? Default is FALSE.
WSCdailyFlowsID	Optional. If NULL (the default) the WSC daily flows will not be plotted. If a WSC station ID is specified, e.g. WSCdailyFlowsID = "05CC001", then the daily flows will be obtained from <b>tidyhydat</b> and plotted.
WSCdailyFlowsLabel	Optional. Labels for the daily flows. If not specified, then the WSC station number will be used, followed by 'daily'.
WSCpeakFlowsID	Optional. If NULL (the default) the WSC peak flows will not be plotted. If a WSC station number is specified, e.g. WSCpeakFlows = "05CC001", then the daily flows will be obtained from <b>tidyhydat</b> and plotted.
WSCpeakFlowsLabel	Optional. Labels for the annual peak flows. If not specified, then the WSC station ID will be used, followed by 'annual peak'.
forceMissingPeakTimes	Optional. Some peaks may be missing their time of day and/or time zone. If TRUE, the missing peak times will be set to noon on the date of record. The missing time zone will be set to that of the other values or the user's time zone if there are none specified. If FALSE (the default value) peaks missing times and/or timezones will be deleted from the plot.
commonTime	Optional. If set to TRUE then the hydrographs will only plotted over their common time range. Default is FALSE.
fakeDates	Optional. If set to TRUE then all dates have their year replaced with 2000, and the actual year is added as a variable in the plotted data. This allows the plot to be faceted by year, as shown in the examples.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
hydat_path	Optional. Path to the HYDAT database. This can usually be left blank.

**Value**

If successful, returns a **ggplot2** object. If unsuccessful, returns FALSE.

**Note**

The CRHM flows are plotted as lines, the daily flows are plotted as steps, and the annual peaks are plotted as points.

**Author(s)**

Kevin Shook



## Examples

```
## Not run:
p <- hydrograph(CRHMflows = "BadLakeModelOutput", CRHMflowsLabels = "CRHM Bad Lake model",
CRHMcols = 1, WSCpeakFlowsID = "05HF014")
# once the ggplot graph has been returned, it can easily be modified:
mintime <- as.POSIXct(as.Date('1975-04-01', format='%Y-%m-%d'), tz='etc/GMT+6')
maxtime <- as.POSIXct(as.Date('1975-04-30', format='%Y-%m-%d'), tz='etc/GMT+6')
# load in all of ggplot2 to modify plots
# library(ggplot2)
p <- p + xlim(mintime, maxtime) + ylim(0, 4)

# re-plot with fake dates
p2 <- hydrograph(CRHMflows = "BadLakeModelOutput", CRHMflowsLabels = "CRHM Bad Lake model",
CRHMcols = 1, WSCdailyFlowsID = "05HF014", WSCpeakFlowsID = "05HF014",
commonTime=TRUE, fakeDates=TRUE)
p2 <- p2 + facet_wrap(~year, scales='free_y', ncol=8)
# set axis limits to be the specified date range, and set labels to be the month names
mintime <- as.POSIXct(as.Date('2000-03-01', format='%Y-%m-%d'), tz='etc/GMT+6')
maxtime <- as.POSIXct(as.Date('2000-06-30', format='%Y-%m-%d'), tz='etc/GMT+6')
p2 <- p2 + scale_x_datetime(date_breaks = "1 month",
limits=c(mintime, maxtime), date_labels = "%b")

## End(Not run)
```

---

hydroYear

*Returns a vector of the hydrological year for the input data frame*


---

## Description

Calculates the hydrological year, based on the starting month.

## Usage

```
hydroYear(CRHMdata, startMonth = 10, useSecondYear = TRUE,
logfile = "")
```

## Arguments

CRHMdata	Required. A <b>CRHM</b> r data frame.
startMonth	Optional. The starting month for the hydrological year. Default is 10 (October).
useSecondYear	Optional. Logical. Determines if the hydrological year is based on the first or second calendar year. In other words would the hydrological year for January 1, 2015 be 2014 or 2015? The default is TRUE (i.e., the hydrological year would be 2015). Note that the Campbell Scientific program SPLIT uses the first calendar year (i.e., the hydrological year would be 2014). To emulate this program, set useSecondYear to be FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns the hydrological year as a vector. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook. Inspired by a function by Chris Marsh.

**See Also**

[aggDataframe](#)

**Examples**

```
# create a new data frame, then add the hydro year
BadLake <- BadLake7376
BadLake$hydroyear <- hydroYear(BadLake, startMonth=9)
```

---

impute

---

*Imputes CRHM obs values*


---

**Description**

This function fills in missing (NA) values in a **CRHM** obs data frame by imputation. The primary values are the data that will be infilled. The secondary values are those used to fill in the gaps. The secondary values are adjusted using the specified multipliers (regression slopes) and offsets (regression intercepts).

**Usage**

```
impute(primaryObs, primaryCols = 1, secondaryObs, secondaryCols = 1,
        multipliers = 1, offsets = 0, quiet = TRUE, logfile = "")
```

**Arguments**

primaryObs	Required. The primary <b>CRHM</b> data frame of obs values.
primaryCols	Optional. A vector containing the columns to be imputed in the primary data frame, not including the datetime. If not specified, defaults to the first column.
secondaryObs	Required. The secondary CRHM obs data frame.
secondaryCols	Optional. A vector containing the columns to be imputed in the secondary data frame, not including the datetime. If not specified, defaults to the first column.
multipliers	Optional. A vector of multipliers applied to secondary observations. These may be obtained using the regress function. The default value is 1.
offsets	Optional. A vector of offsets added to secondary observations. These may be obtained using the regress function. The default value is 0.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful returns a modified version of the primaryObs data frame. The missing values in the primary data frame are replaced by corresponding values from the secondary data frame, after adjustment using the specified multipliers and offsets. If unsuccessful, returns the value FALSE.

**Note**

In addition to the usual notation in the logfile, this function also writes a separate logfile which summarizes the primaryObs data frame, and the new infilled data frame. The summaries are also printed to the screen, if quiet=FALSE. The logfile also contains a complete listing of the source of each value in the infilled data frame. Each value is listed as being 'original' (unmodified from the primaryObs data frame), 'imputed' (derived from the adjusted values of the secondaryObs data frame) or 'NA' (missing).

**Author(s)**

Kevin Shook

**See Also**

[interpolate regress distributeP](#)

**Examples**

```
## Not run:
v.filled <- impute(veg, c(1,2,3), st,
c(2,3,1), c(0.895,0.708,1.1209), c(-0.8128, 0.0607, 9.005))
## End(Not run)
```

---

insertMissing	<i>Inserts missing datetimes</i>
---------------	----------------------------------

---

**Description**

Many time series have missing rows. This function finds missing rows, and adds rows with the appropriate datetime values and NA\_real\_ values for all variables. Once the missing values have been added, you can use [interpolate](#) or [impute](#) to infill them.

**Usage**

```
insertMissing(obs, quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> data frame containing obs values.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful returns a modified version of the obs data frame with missing datetime values inserted. All of the missing variables in the inserted rows are replaced by `NA_real_` values. If unsuccessful, returns the value `FALSE`.

**Note**

This function should be used before doing interpolation or imputation of variables. Note that this function is also called by the function [findGaps](#).

**Author(s)**

Kevin Shook

**See Also**

[findGaps](#) [interpolate](#) [impute](#)

**Examples**

```
BadLake.withmissing <- insertMissing(BadLake7376, quiet=FALSE)
```

---

interpolate

*Fills missing obs values by interpolation*


---

**Description**

Missing (NA) values in a **CRHMr** obs data frame are filled by linear or spline interpolation. The user can set the maximum gap length (in time steps) allowed to be filled. Note that this function will *NOT* interpolate RH or precipitation values.

**Usage**

```
interpolate(obs, varcols = 1, methods = "linear", maxlength = 5,
  quiet = TRUE, logfile = "")
```

**Arguments**

<code>obs</code>	Required. The <b>CRHMr</b> data frame of obs values.
<code>varcols</code>	Optional. A vector containing the columns to be imputed in the obs data frame, not including the datetime. If not specified, defaults to the first column.
<code>methods</code>	Optional. A vector containing the methods to be used for interpolation for each of the variables. Currently supported methods are 'linear' and 'spline'. The default is to use linear interpolation. If fewer methods than columns are specified, the methods are recycled.
<code>maxlength</code>	Optional. The maximum gap length to be interpolated. Defaults to 5 time steps.

quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a modified version of the obs data frame. The missing values in the specified data frame are replaced by interpolated values. If unsuccessful, returns the value FALSE.

**Note**

In addition to the usual notation in the logfile, this function also writes a separate logfile which summarises the original obs data frame, and the new infilled data frame. The summaries are also printed to the screen, if quiet=FALSE. The logfile also contains a complete listing of the source of each value in the infilled data frame. Each value is listed as being 'original' (unmodified from the primary obsframe), 'linear interpolation' (infilled by linear interpolation), 'spline interpolation' (infilled by spline interpolation) or 'NA' (missing).

**Author(s)**

Kevin Shook

**See Also**

[impute.distributeP](#)

**Examples**

```
## Not run:
v.filled <- interpolate(v, c(1,3))
## End(Not run)
```

---

keepGood

*Keeps rows where there are some obs values*

---

**Description**

Removes rows where all values are missing, keeps the remainder.

**Usage**

```
keepGood(obs, quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> obs dataframe.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns the dataframe containing all rows with any non-missing values. If unsuccessful (including if there are only missing values in the original data), returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[trimObs](#)

**Examples**

```
BadLake.good <- keepGood(BadLake7376)
```

---

list2df	<i>Convert a list of vectors to a data frame.</i>
---------	---

---

**Description**

This function will convert a list of vectors to a data frame. This function will handle three different types of lists of vectors. First, if all the elements in the list are named vectors, the resulting data frame will have a number of columns equal to the number of unique names across all vectors. In cases where some vectors do not have names in other vectors, those values will be filled with NA.

**Usage**

```
list2df(x, row.names = NULL, optional = FALSE, ...)
```

**Arguments**

x	a list to convert to a data frame.
row.names	a vector equal to length(x) corresponding to the row names. If NULL, the row names will be set to names(x).
optional	not used.
...	other parameters passed to <a href="#">data.frame</a> .

## Details

The second case is when all the vectors are of the same length. In this case, the resulting data frame is equivalent to applying `rbind` across all elements.

The third case handled is when there are varying vector lengths and not all the vectors are named. This condition should be avoided. However, the function will attempt to convert this list to a data frame. The resulting data frame will have a number of columns equal to the length of the longest vector. For vectors with length less than this will fill the row with NAs. Note that this function will print a warning if this condition occurs.

## Value

a data frame.

## Author(s)

Jason Bryer «[jason@bryer.org](mailto:jason@bryer.org)»

## References

<http://stackoverflow.com/questions/4227223/r-list-to-data-frame>

## Examples

```
## Not run:
test1 <- list( c(a='a',b='b',c='c'), c(a='d',b='e',c='f'))
as.data.frame(test1)

test2 <- list( c('a','b','c'), c(a='d',b='e',c='f'))
as.data.frame(test2)

test3 <- list('Row1'=c(a='a',b='b',c='c'), 'Row2'=c(var1='d',var2='e',var3='f'))
as.data.frame(test3)

test4 <- list('Row1'=letters[1:5], 'Row2'=letters[1:7], 'Row3'=letters[8:14])
as.data.frame(test4)

test5 <- list(letters[1:10], letters[11:20])
as.data.frame(test5)

test6 <- list(list(letters), letters)
as.data.frame(test6)
## End(Not run)
```

---

logAction	<i>Writes a comment to the <b>CRHMr</b> log file</i>
-----------	--

---

### Description

This is an internal **CRHMr** function and should normally not need to be called directly. This function is called by almost all **CRHMr** functions, and writes the comment, and the date and time to the logfile.

### Usage

```
logAction(comment = "", logfile = "")
```

### Arguments

comment	Required. Comment string to be written. This normally includes the function name, as well as the obs file being processed.
logfile	Optional. Name of the file to be used for logging the action. Normally not used. If not specified, then 'CRHMr.log' in the current working directory will be used.

### Value

If successful, returns TRUE. If unsuccessful, returns FALSE.

### Author(s)

Kevin Shook

### Examples

```
## Not run:
comment <- paste('tMinMaxToHourly dataframe:', obsName, sep='')
result <- logAction(comment)
## End(Not run)
```

---

longwave	<i>Calculates incoming longwave radiation</i>
----------	---

---

### Description

Calculates the incoming longwave radiation from the specified emissivity and air temperatures. This is an internal **CRHMr** function and should *never* need to be called directly, as it does *NO* data checking.

### Usage

```
longwave(em, tK)
```



**Arguments**

em	Required. Emissivity (0-1).
tK	Required. Air temperatures in K.

**Value**

Returns the longwave radiation in  $\text{W/m}^2$ .

**Author(s)**

Kevin Shook

**See Also**

[distributeQli emissivity](#)

**Examples**

```
longwave(0.2, 293)
```

---

makeRegular

---

*Makes the datetimes of an obs data frame fit the time step exactly*


---

**Description**

Makes the datetimes of an obs data frame fit the time step exactly

**Usage**

```
makeRegular(obs, timezone = "", quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> obs data frame.
timezone	Required. Timezone to be assigned to data.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a modified version of the obs data frame containing the adjusted datetime and all of the variables. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**Examples**

```
regular <- makeRegular(BadLake7376, timezone='CST')
```

---

maxObs	<i>Sets maximum values for an obs data frame</i>
--------	--

---

**Description**

Tests values in a **CRHMr** obs data frame to see if they exceed maximum thresholds. Values exceeding the thresholds can be set to either the maximum allowable value or to `NA_real_`, which is useful for infilling or imputing values.

**Usage**

```
maxObs(obs, varcols = "", maxvals = "", actions = "max",
       quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. The <b>CRHMr</b> obs data frame.
varcols	Optional. A vector containing the columns to be imputed in the obs data frame, not including the datetime. If not specified, defaults to all columns.
maxvals	Optional. A vector containing the maximum permissible values for each of the specified columns. If omitted, the default values are used.
actions	Optional. A vector containing the methods to be used for replacing values that exceed the maximum threshold values. Currently supported actions are 'max' (using the threshold) and 'NA' (inserting <code>NA_real_</code> values). The default is 'max'. If fewer actions are specified than variables, then the actions are recycled.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet=FALSE</code> .
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a modified version of the obs data frame. The values exceeding the specified max values are replaced by either `NA_real_` or the threshold max values. If unsuccessful, returns the value `FALSE`.

**Note**

The default threshold values are

t	40 C
ea	3.5 kPa
rh	100 percent
ppt	100 mm
p	20 mm
u	10 m/s
SunAct	18 hr
qsi	1262 W/m <sup>2</sup>
qso	1000 W/m <sup>2</sup>
qn	250 W/m <sup>2</sup>

Author(s)

Kevin Shook

See Also

[minObs](#)

Examples

```
# use all of the default values
bad.max <- maxObs(BadLake7376)
summary(bad.max)
# use specified columns with default max values, replace with 'NA' values
bad.max2 <- maxObs(BadLake7376, varcols=c(1,2), actions='NA')
# use specfied columns with specified max values and actions
bad.max3 <- maxObs(BadLake7376, maxvals=c(35, 10), varcols=c(1,3), actions=c('NA', 'max'))
summary(bad.max3)
```

---

minObs	<i>Sets minimum values for an obs data frame</i>
--------	--

---

Description

Tests values in a **CRHMr** obs data frame to see if they exceed minimum thresholds. Values exceeding the thresholds can be set to either the minimum allowable value or to NA\_real, which is useful for infilling or imputing values.

Usage

```
minObs(obs, varcols = "", minvals = "", actions = "min",
       quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. The <b>CRHMr</b> obs data frame.
varcols	Optional. A vector containing the columns to be imputed in the obs data frame, not including the datetime. If not specified, defaults to all columns.
minvals	Optional. A vector containing the minimum permissible values for each of the specified columns. If omitted, the default values are used.
actions	Optional. A vector containing the methods to be used for replacing values that exceed the minimum threshold values. Currently supported actions are 'min' (using the threshold) and 'NA' (using NA_real_ values). The default is 'min'. If fewer actions are specified than variables, then the actions are recycled.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a modified version of the obs data frame. The values smaller than the specified min values are replaced by either NA\_real\_ or the threshold min values. If unsuccessful, returns the value FALSE.

**Note**

The default threshold values are

t	-40 C
ea	0.01 kPa
rh	0.5 percent
ppt	0 mm
p	0 mm
u	0 m/s
SunAct	0 hr
qsi	0 W/m <sup>2</sup>
qso	0 W/m <sup>2</sup>
qn	-60 W/m <sup>2</sup>

**Author(s)**

Kevin Shook

**See Also**

[maxObs](#)

**Examples**

```
# use all of the default values
```

```
bad.min <- minObs(BadLake7376)
# use specified columns with default min values, replace with 'NA' values
bad.min2 <- minObs(BadLake7376, varcols=c(1,2), actions='NA')
# use specified columns with specified min values and actions
bad.min3 <- minObs(BadLake7376, minvals=c(-30, 22), varcols=c(1,2), actions=c('min', 'NA'))
```

---

monthlyPrecipTotals	<i>Returns mean monthly precipitation totals</i>
---------------------	--

---

## Description

Calculates mean total precipitation for each calendar month in the observation data frame. Optionally plots the values. Can work with p and/or ppt values.

## Usage

```
monthlyPrecipTotals(obs, precipCols = 0, plot = TRUE, logfile = "")
```

## Arguments

obs	Required. A <b>CRHMr</b> data frame containing p and/or ppt values..
precipCols	Optional. A vector containing the columns to be analyzed, not including the datetime column. If not specified, all precipitation columns are used.
plot	Optional. If TRUE, which is the default, a <b>ggplot2</b> object of the monthly mean precipitations is returned. If FALSE, the values are returned as a data frame.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If successful, returns either a data frame or a **ggplot2** object of the monthly mean precipitations. If unsuccessful, returns FALSE.

## Author(s)

Kevin Shook

## See Also

[aggDataFrame monthlyQQplot](#)

## Examples

```
## Not run:
marmot_monthly_plot <- monthlyPrecipTotals(marmot)
marmot_monthly_means <- monthlyPrecipTotals(marmot, plot=FALSE)
## End(Not run)
```

---

monthlyQQplot	<i>Quantile-quantile plot by month</i>
---------------	--

---

### Description

Does QQ plots of monthly values of a single variable type in two data frames.

### Usage

```
monthlyQQplot(primaryCRHM, primaryCol = 1, secondaryCRHM,
               secondaryCol = 1, samePeriod = TRUE, logfile = "")
```

### Arguments

primaryCRHM	Required. The primary <b>CRHM</b> r data frame. Quantiles of this data will be plotted on the X axis.
primaryCol	Optional. The column in the primary data frame, not including the datetime. If not specified, defaults to the first column.
secondaryCRHM	Required. The secondary <b>CRHM</b> r data frame. Quantiles of this data will be plotted on the Y axis.
secondaryCol	Optional. The column in the secondary data frame, not including the datetime. If not specified, defaults to the first column.
samePeriod	Optional. Logical. Should the same period of time be used for both variables? Default is TRUE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns a **ggplot2** object of faceted monthly QQ plots (3 rows x 4 columns) of the specified variables. If unsuccessful, returns FALSE.

### Author(s)

Kevin Shook

### See Also

[monthlyPrecipTotals](#)

### Examples

```
BadLake <- BadLake7376
BadLake$year <- as.numeric(format(BadLake7376$datetime, format='%Y'))
badlake73 <- subset(BadLake, year==1973)
badlake75 <- subset(BadLake, year==1975)
p <- monthlyQQplot(badlake73, 1, badlake75, 1, samePeriod=FALSE)
print(p)
```

---

nancumsum	<i>Finds cumulative sum of non-missing values in an array</i>
-----------	---

---

**Description**

Finds cumulative sum of non-missing values in an array

**Usage**

```
nancumsum(x)
```

**Arguments**

x	Required. An array or vector of numeric values
---	--

**Value**

Returns an array or vector with cumulative sums of each column.

**Author(s)**

Kevin Shook. The code is copied from a MATLAB program nancumsum.m written by Alan Barr.

**See Also**

[PcpFiltPosTh](#)

**Examples**

```
a <- nancumsum(BadLake7376[,2])
```

---

parseNums	<i>Parses a string containing numbers</i>
-----------	---

---

**Description**

This is an internal **CRHM** function and should normally not need to be called directly. It is used for reading data from messy files such as CRHM projects.

**Usage**

```
parseNums(numString)
```

**Arguments**

numString	Required. A character string containing numbers separated by any number of spaces.
-----------	--

**Value**

Returns a numeric vector.

**Author(s)**

Kevin Shook

**Examples**

```
parseNums(' 1 2 3 4 5 ')
```

---

parseText

*Parses a string containing several sub-strings*

---

**Description**

This is an internal **CRHMr** function and should normally not need to be called directly. It is used for reading data from messy files such as CRHM projects.

**Usage**

```
parseText(textString)
```

**Arguments**

textString	Required. A character string containing strings separated by any number of spaces.
------------	--

**Value**

Returns a character vector.

**Author(s)**

Kevin Shook

**Examples**

```
parseText(' red green blue black')
```



---

PcpFiltPosTh

*Cleans up a cumulative precipitation time series*


---

## Description

Cleans up a cumulative precipitation time series by transferring changes below a specified threshold to neighbouring periods, and eliminating large negative changes associated with gauge servicing (bucket emptying). The filtering is done using a brute-force algorithm that identifies small or negative changes (below dPcpTh) then transfers them to neighbouring positive changes thus aggregating all changes to values above dPcpTh. The transfers are made in ascending order, starting with the lowest (most negative). The cumulative total remains unchanged. This description (and most of the comments in the code) were written by Alan Barr.

## Usage

```
PcpFiltPosTh(Pcp, dPcpTh = 0.1, dpServicingTh = -100, quiet = TRUE)
```

## Arguments

Pcp	Required. Measured cumulative precipitation (mm).
dPcpTh	Optional. Minimum interval precipitation (mm). Default is 0.1 mm.
dpServicingTh	Optional. Threshold for change in storage due to servicing (mm). Default is -100 mm.
quiet	Optional. Suppresses display of messages, except for errors. If set quiet=FALSE, the output will pause for 5 seconds ever 100 iterations, which may be useful for debugging your parameters. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default), which causes fewer parameters to be output, without pausing, every 1000 iterations.

## Value

If unsuccessful, returns the value FALSE. If successful, returns a time series of cleaned values. Writes parameter values to screen.

## Author(s)

Kevin Shook. The code is copied from a MATLAB program PcpFiltPosTh.m written by Alan Barr, 28 Aug 2012.

## Examples

```
## Not run:
testp <- PcpFiltPosTh(d12$precip, dPcpTh = 0.05, dpServicingTh = -50)
## End(Not run)
```

---

phaseCorrect

---

*Corrects precipitation for undercatch and calculates phase*


---

## Description

Applies an iterative solution to calculate the psychrometric hydrometeor temperature. The hydrometeor temperature is then used to partition the raw precipitation into rainfall and snowfall. The snowfall is adjusted for gauge undercatch, if selected.

## Usage

```
phaseCorrect(obs, Tcol = 1, RHcol = 2, Ucol = 3, Pcol = 4,
             RH_type = 1, shield = 1, quiet = TRUE, logfile = "")
```

## Arguments

obs	Required. A <b>CRHMr</b> data frame of observations.
Tcol	Required. Column number containing the air temperature (not including the datetime). Default is column 1. The values must be in °C.
RHcol	Required. Column number containing the RH (not including the datetime). Default is column 2. The values must be as percentages.
Ucol	Required. Column number containing the wind speeds (not including the datetime). Default is column 3. The values must be in m/s.
Pcol	Required. Column number containing the precipitaion (not including the datetime). Default is column 4. The values must be in mm.
RH_type	Optional. Set to 1 if RH is relative to water, any othe value if relative to ice. Default is 1.
shield	Optional. Set to 1 to use Alter shield undercatch correction from MacDonald and Pomeroy (2007). Set to 2 to use Alter shield undercatch correction from Smith. Set to any other value for unadjusted precipitation. Default is 1.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If unsuccessful returns FALSE. If successful, returns the original dataframe with additional columns for the rain, snow, total precipitation and rain ratio.

## Author(s)

Phillip Harder

## References

Harder, P., and J. Pomeroy (2013), “Estimating precipitation phase using a psychrometric energy balance method”, *Hydrol. Process.*, 27(13), 1901-1914, doi:10.1002/hyp.9799.

Macdonald, J., and J. Pomeroy (2007), “Gauge Undercatch of Two Common Snowfall Gauges in a Prairie Environment”, *Proc. 64th East. Snow Conf. St. John’s, Canada.*, 119-126.

## Examples

```
corrected <- phaseCorrect(BadLake7376, Tcol=1, RHcol=2, Ucol=3, Pcol=5)
```

---

plotObs	<i>Creates a <b>ggplot</b> object from an obs data frame</i>
---------	--

---

## Description

This function creates a **ggplot2** object from an obs data frame. If the specified variables are all of single type (e.g. temperature), it returns a single-pane plot. If the specified variables are of more than one type (e.g. temperature, precipitation), then the plot is faceted. The function returns an object that you can modify, or save with the standard **ggplot2** commands.

## Usage

```
plotObs(obs, varcols = c(0), startDate = "", endDate = "",
        plotType = "lines")
```

## Arguments

obs	Required. The <b>CRHMr</b> obs data frame to be plotted.
varcols	Optional. A vector containing the numbers of the columns (not including the datetime) to be plotted. If not specified, all variables are plotted.
startDate	Optional. The starting date for the plot. Can either be a year as a number (e.g. 1995) or a date in ‘Y-m-d’ format, i.e. ‘1995-06-01’. If not specified, the first datetime value in the obs data is used.
endDate	Optional. The ending date for the plot. Can either be a year as a number (e.g. 1995) or a date in ‘Y-m-d’ format, i.e. ‘1995-12-31’. If not specified, the first datetime value in the obs data is used.
plotType	Optional. Defines the type of the plot. Either ‘lines’ or ‘points’. Default is ‘lines’. If you want to show the extent of data, use ‘points’ as lines will NOT be drawn when points are separated by NA_real_ values.

## Details

Units are automatically specified for obs variables of the following types: t, rh, u, p, ea, and solar radiation.

**Value**

If successful returns a **ggplot2** object. If unsuccessful, returns the value FALSE.

**Note**

Note that missing (NA\_real\_) values in the obs data frame will result in a warning message.

**Author(s)**

Kevin Shook

**See Also**

[summariseObsFiles](#)

**Examples**

```
p <- plotObs(BadLake7376)
print(p)
```

---

plotPrecipsByYear	<i>Creates a <b>ggplot2</b> object of cumulative precipis by year</i>
-------------------	---

---

**Description**

Calculates cumulative precipitations and creates a **ggplot2** faceted by year. Note that this function would work for any other cumulative variable, if the Y-axis label is changed.

**Usage**

```
plotPrecipsByYear(obs = "", obsNames = "", precipCols = 1,
  commonDates = TRUE)
```

**Arguments**

obs	Required. Either a <b>CRHMr</b> obs data frame or a list of data frames.
obsNames	Optional. A vector containing names for the data frames. If there are more data frames than names, the data will be named obs_1, obs_2 etc.
precipCols	Optional. Column(s) containing precipitation. Can be a single value or a vector. Values will be recycled if necessary. Default is 1.
commonDates	Optional. If TRUE (the default) then the range of dates will be restricted to the common first and last dates of the datasets.

**Value**

If successful, returns a faceted **ggplot2** object of cumulative precipitation vs time. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook.

**See Also**

[plotTempsByYear](#)

**Examples**

```
## Not run: l <- list(meas, WRF)
names <- c("Measured", "WRF")
p <- plotPrecipsByYear(l, names)
library(ggplot2)
p <- p + scale_colour_manual(values=c("red", "blue"))
## End(Not run)
```

---

plotTempsByYear	<i>Creates a <b>ggplot2</b> object of daily mean air temps by year</i>
-----------------	--

---

**Description**

Aggregates air temperatures to mean daily values and creates a **ggplot2** faceted by year. Note that this function would work for any other non-cumulative variable, if the Y-axis label is changed.

**Usage**

```
plotTempsByYear(obs = "", obsNames = "", tempCols = 1,
  commonYears = TRUE)
```

**Arguments**

obs	Required. Either a <b>CRHMr</b> obs data frame or a list of data frames.
obsNames	Optional. A vector containing names for the data frames. If there are more data frames than names, the data will be named obs 1, obs 2 etc.
tempCols	Optional. Column(s) containing air temperatures. Can be a single value or a vector. Values will be recycled if necessary. Default is 1.
commonYears	Optional. If TRUE (the default) then the range of years will be restricted to the common first and last years.

**Value**

If successful, returns a faceted **ggplot2** object of mean daily air temperatures vs time. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook.

**See Also**

[plotPrecipsByYear](#)

**Examples**

```
## Not run: l <- list(meas, WRF)
names <- c("Measured", "WRF")
p <- plotTempsByYear(l, names)
library(ggplot2)
p <- p + scale_colour_manual(values=c("red", "blue"))
## End(Not run)
```

---

qair2ea

*Converts specific humidity to vapour pressure.*


---

**Description**

Converts specific humidity to vapour pressure.

**Usage**

```
qair2ea(qair, press = 101325)
```

**Arguments**

qair	Required. Specific humidity (dimensionless).
press	Optional. Air pressure in Pa. Default is 101325, the standard atmospheric air pressure.

**Value**

If successful, returns the vapour pressure in Pa. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**References**

R code for conversion of air pressure and absolute humidity was taken from project PEcAn The Predictive Ecosystem Analyzer <http://pecanproject.github.io>. The source code is available at <https://github.com/PecanProject/pecan/blob/master/modules/data.atmosphere/R/metutils.R>.

**See Also**

[qair2rh](#)

**Examples**

```
ea <- qair2ea(0.0001, 101325)
```

---

qair2rh	<i>qair2rh</i>
---------	----------------

---

**Description**

From Bolton 1980 The computation of Equivalent Potential Temperature [http://www.eol.ucar.edu/projects/ceop/dm/documents/refdata\\_report/eqns.html](http://www.eol.ucar.edu/projects/ceop/dm/documents/refdata_report/eqns.html).

**Usage**

```
qair2rh(qair, temp, press = 101325, allowStupid = FALSE)
```

**Arguments**

qair	Required. Specific humidity, dimensionless (e.g. kg/kg) ratio of water mass / total air mass.
temp	Required. Air temp in degrees C.
press	Optional. Air pressure in Pa. Default is 101325, the standard atmospheric air pressure.
allowStupid	Optional. If TRUE then impossible values (smaller than zero or greater than 1) will be allowed. The default value, FALSE, restricts the returned value to be between 0 and 1.

**Details**

Converts specific humidity to relative humidity

**Value**

If successful, returns rh (relative humidity), the ratio of actual water mixing ratio to saturation mixing ratio. Note that this value is a fraction (i.e. between 0 and 1) rather than a percentage. If unsuccessful, returns the value FALSE.

**Author(s)**

David LeBauer. Modified by Kevin Shook for air temps < 0 °C and to work for vectors or scalars.

**References**

<https://github.com/PecanProject/pecan/blob/master/modules/data.atmosphere/R/metutils.R>.

**See Also**[qair2ea](#)**Examples**

```
qair2rh(0.0001, 10)
```

---

**qqplotValues***Calculates quantile values for a Quantile-Quantile plot*

---

**Description**

The built-in qqplot function does not work with **ggplot2**. This just calls the qqplot function to calculate the quantiles without plotting them.

**Usage**

```
qqplotValues(x, y)
```

**Arguments**

x	Required. A numeric vector.
y	Required. A numeric vector.

**Value**

Returns a dataframe with the quantiles of x and y.

**Author(s)**

Kevin Shook

**Examples**

```
quantiles <- qqplotValues(runif(20), runif(50))
```



---

readCampbell	<i>Reads in Campbell datalogger data</i>
--------------	--

---

### Description

Reads a datafile produced by Campbell Scientific dataloggers into a **CRHMr** data frame. Both old (multiple table, no header) and new (single table, with header) data files can be read in.

### Usage

```
readCampbell(campbellFile, missingValue = -6999, timezone = "",
             quiet = TRUE, logfile = "")
```

### Arguments

campbellFile	Required. The name of the file containing the Campbell Scientific datalogger output.
missingValue	Optional. Default is -6999. All values less than or equal to this will be set to be NA_real_.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If unsuccessful, returns FALSE. If successful, returns a **CRHMr** data frame. If there are two tables in the campbellFile, then they are returned as a list.

### Note

If the campbellFile does not contain headers, then the variable names will be assigned by R. In any case, you will probably want to rename the variables using the command names().

### Author(s)

Kevin Shook

See Also

[readObsFile](#) [readExportFile](#) [readOutputFile](#)

Examples

```
## Not run:
# read in an old format Campbell file
oldformatdata <- readCampbell('OldDataLogger.dat', timezone='MST')
# now extract the data frames from the returned list
obs1 <- oldformatdata[[1]]
obs2 <- oldformatdata[[2]]
# read in a new format Campbell file
obs3 <- readCampbell('BowHut_Fifteen.dat', timezone='MST')
## End(Not run)
```

---

readCLASSfile	<i>Reads forcing data from CLASS formatted file</i>
---------------	---

---

Description

Reads data from files created for CLASS into a stanard **CRHMr** obs data frame.

Usage

```
readCLASSfile(CLASSfile = "", timezone = "", convertNames = TRUE,
  quiet = TRUE, logfile = "")
```

Arguments

CLASSfile	Required. File of meteorological data in CLASS format.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use ‘CST’ and ‘MST’ for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use ‘etc/GMT+6’ or ‘etc/GMT+7’ for Central Standard and Mountain Standard time. DO NOT use ‘America/Rgina’ as the time zone, as it includes historical changes between standard and daylight savings time.
convertNames	Optional. If TRUE, then the CLASS variables will be converted to Qsi.1, Qli.1, p.1, t.1, ea.1, u.1, and AirPressure.1. Otherwise, the CLASS names SWin, LWin, Prec, T, qa, U, and P will be used.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns a **CRHM** data frame. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook, with assistance from Andrew Ireson.

**Examples**

```
## Not run:
class <- readCLASSfile('driv.met', timezone='Etc/GMT+7')
## End(Not run)
```

---

readDebugFile	<i>Reads in CRHM debug file</i>
---------------	---------------------------------

---

**Description**

Reads in CRHM debug file

**Usage**

```
readDebugFile(debugFile = "", returnData = "all")
```

**Arguments**

debugFile	Required. CRHM debug file name.
returnData	Optional. The data to be returned. If 'all' (the default) then all sections will be returned as a list of data frames. Other options are 'header', 'initialconditions', 'finalconditions', 'basinsummary', and 'groupsummary'. Note that if 'groupsummary' is specified, and there are no groups, then an error will result.

**Value**

If successful, returns a list of data frames, or a single specified data frame. If unsuccessful, returns FALSE.

**Examples**

```
## Not run:
allDebug <- readDebug("07-08_debug.txt")
# show header
View(allDebug$header)
# show group summary
View(allDebug$groupSummary)
## End(Not run)
```

---

readExportFile	<i>Reads exported CRHM output into a data frame</i>
----------------	---

---

### Description

Reads output from a CRHM model, that has been manually exported, into a **CRHMr** data frame. CRHM model output can be exported in several formats. This function is intended to read in all of them. Note that because of the way that R reads in files, the variable names will have appended periods, e.g. t.1.

### Usage

```
readExportFile(exportFile, timezone = "", quiet = TRUE, logfile = "")
```

### Arguments

exportFile	Required. The name of the CRHM export file to be read in.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns a **CRHMr** data frame. If unsuccessful, returns the value FALSE.

### Author(s)

Kevin Shook

### See Also

[readOutputFile](#) [readObsFile](#)

### Examples

```
## Not run:
brandon <- readExportFile('Brandon_export.txt', 'etc/GMT+6')
## End(Not run)
```

---

readObsFile	<i>Reads a CRHM obs file into a dataframe</i>
-------------	---

---

### Description

Reads a file of CRHM observation data into a **CRHM** data frame. Note that because of the way that R reads in files, the variable names will have appended periods, e.g. t.1.

### Usage

```
readObsFile(obsFile = "", timezone = "", quiet = TRUE,
            logfile = "")
```

### Arguments

obsFile	Required. Name of the CRHM obs file to be read in.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns a **CRHM** data frame. If unsuccessful, returns the value FALSE.

### Author(s)

Kevin Shook

### See Also

[readExportFile](#) [readOutputFile](#)

### Examples

```
# output example obs data to a file
BadLake7376.obs <- writeObsFile(BadLake7376, 'BadLake7376.obs')
# now read data back in
BadLake <- readObsFile('BadLake7376.obs', 'etc/GMT+6')
```

---

readOutputFile	<i>Reads CRHM model output into a data frame</i>
----------------	--

---

### Description

Reads a file containing output from a CRHM model into a **CRHM***r* data frame. Note that because of the way that R reads in files, the variable names will have appended periods, e.g. `t.1`. This version reads in both old CRHM output (without units) and new CRHM output which contains the variable units in the second line.

### Usage

```
readOutputFile(outputFile, timezone = "", quiet = TRUE, logfile = "")
```

### Arguments

<code>outputFile</code>	Required. The name of the CRHM output file to be read in.
<code>timezone</code>	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
<code>quiet</code>	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet=FALSE</code> .
<code>logfile</code>	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns a **CRHM***r* data frame. If unsuccessful, returns the value `FALSE`.

### Author(s)

Kevin Shook

### See Also

[readExportFile](#) [readObsFile](#) [readOutputUnits](#)

### Examples

```
## Not run:
stoon <- readOutputFile('CRHM_output_1.txt', 'etc/GMT+6')
## End(Not run)
```

---

readOutputUnits	<i>Reads variable units</i>
-----------------	-----------------------------

---

### Description

Gets units for all variables in a specified CRHM output file. If the file is the new CRHM format, then the units are read directly from the second line of the file. If the file is in the older CRHM format, then units are found by matching the variables to those in the internal data frame CRHM\_vars.

### Usage

```
readOutputUnits(outputFile, quiet = TRUE)
```

### Arguments

outputFile	Required. A CRHM output file.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet = FALSE.

### Value

If successful, returns a **CRHM** data frame containing the name of each variable and its units. If unsuccessful, returns the value FALSE.

### Author(s)

Kevin Shook

### See Also

[readOutputFile](#)

### Examples

```
## Not run:  
units <- readOutputFile('CRHM_output_1.txt')  
## End(Not run)
```

---

readPrjDimensions	<i>Reads the dimensions from a .prj file</i>
-------------------	--

---

**Description**

Reads the dimensions from a .prj file

**Usage**

```
readPrjDimensions(prjFile = "", logfile = "")
```

**Arguments**

prjFile	Required. Name of .prj file.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns the dimension values as a data frame with the columns dimension and value. The dimensions are nhru (the number of HRUs), nlay (the number of layers) and nobs (the number of obs files). If unsuccessful, the value FALSE is returned.

**Author(s)**

Kevin Shook

**See Also**

[readPrjParameters](#)

**Examples**

```
## Not run: model_dims <- readPrjDimensions("Bad74_Frozen.prj")
```

---

readPrjHRUnames	<i>Reads names of HRUs from .prj file</i>
-----------------	---

---

**Description**

This function reads names of HRUs from a .prj file. The function readPrjParameters will not work, as it only returns numeric values.

**Usage**

```
readPrjHRUnames(prjFile = "", logfile = "")
```



**Arguments**

prjFile	Required. Name of .prj file.
logfile	logfile Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns the parameter values as a vector. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[setPrjHRUNames](#)

**Examples**

```
## Not run:
# read in as a vector
HRUNames <- readPrjHRUNames("Bad_Lake_1974-1975.prj")
## End(Not run)
```

---

readPrjModuleParams	<i>Reads all parameters for all HRUs</i>
---------------------	--

---

**Description**

Reads all parameters for all HRUs

**Usage**

```
readPrjModuleParams(moduleName, prjFile)
```

**Arguments**

moduleName	Required. A text pattern of the module name. For the Albedo Richards module it would be 'albedo_Richard'.
prjFile	Required. Name of .prj file.

**Value**

A data frame with the names of the parameters, and the values of the parameters for each HRU.

**Author(s)**

Ezequiel Tourn

## Examples

```
## Not run:
fileName <- "Vuriloche.prj"
shared_parameters <- readPrjModuleParams(moduleName = 'Shared',
prjFile = fileName)
## End(Not run)
```

---

readPrjNumVals

*Reads numeric values from a CRHM project text object*


---

## Description

Reads specified number of numeric values from a project file variable. This is an internal function and is not able to be called by a function outside of **CRHMr**, as it is not exported. This documentation is for maintenance purposes.

## Usage

```
readPrjNumVals(prj, searchString, value_count)
```

## Arguments

prj	Required. A CRHM file read into memory as a sequence of lines, as returned from readPrj.
searchString	Required. The identifying string.
value_count	Required. The number of values to be read in.

## Value

If successful, returns a vector with the specified number of values. If unsuccessful, returns the value FALSE.

## Examples

```
## Not run:
hru_area <- readPrjNumVals(Bologna, 'hru_area', 19)
## End(Not run)
```

---

`readPrjOutputVariables`*Reads output variables from a CRHM model .prj file*

---

**Description**

Reads output variables from a CRHM model .prj file

**Usage**

```
readPrjOutputVariables(prjFile, asDataframe = FALSE, logfile = "")
```

**Arguments**

<code>prjFile</code>	Required. Name of CRHM .prj file to read.
<code>asDataframe</code>	Optional. If FALSE (the default), returns the set of output variables as a vector. If TRUE, returns a data frame containing the module, variable and HRUs as a string. The vector form is useful if you want to delete some variables. The data frame is more useful if you want to change the variable or its HRUs.
<code>logfile</code>	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns the output variables. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[setPrjOutputVariables](#)

**Examples**

```
## Not run:
# read in as a vector
variables <- readPrjOutputVariables('Bad_Lake_1974-1975.prj')

# read in as a data frame
variables <- readPrjOutputVariables('Bad_Lake_1974-1975.prj', asDataframe=TRUE)
# change value
variables$HRUs[1] <- '1 2'
## End(Not run)
```

---

readPrjParameters	<i>Reads values of specified parameter from .prj file</i>
-------------------	---

---

### Description

This function reads numeric parameter values only. Attempting to read in character values, such as HRU names, will return NULL.

### Usage

```
readPrjParameters(prjFile = "", paramName = "", logfile = "")
```

### Arguments

prjFile	Required. Name of .prj file.
paramName	Required. Name of parameter to be read.
logfile	logfile Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns the parameter values as a vector. If unsuccessful, returns FALSE.

### Author(s)

Kevin Shook

### See Also

[setPrjOutputVariables](#)

### Examples

```
## Not run:  
# read in as a vector  
variables <- readPrjParameters("Bad_Lake_1974-1975.prj", "Soil soil_moist_init")  
## End(Not run)
```

---

readPrjTextVals	<i>Reads text values from a CRHM project text object</i>
-----------------	--

---

### Description

Reads text values from a project file variable. The number of values to be read in does not need to be specified. This is an internal function and is not able to be called by a function outside of **CRHMr**, as it is not exported. This documentation is for maintenance purposes.

### Usage

```
readPrjTextVals(prj, searchString, valueCount)
```

### Arguments

prj	Required. A CRHM file read into memory as a sequence of lines, as returned from readPrj.
searchString	Required. The identifying string.
valueCount	Required. The number of values to be read in.

### Value

If successful, returns a vector with the numeric values. If unsuccessful, returns the value FALSE.

### Examples

```
## Not run:
hru_names <- readPrjTextVals(Bologna, 'hru_names', 19)
## End(Not run)
```

---

regress	<i>Regresses one set of CRHM observations against another</i>
---------	---

---

### Description

The purpose of this function is to generate the multiplier and an offset values (if required) used by the impute function. Therefore, it performs a linear (least-squares) regression of the primary dataset (the values to be produced by imputation) against the secondary dataset. The function removes missing values and merges the data sets together before doing the regression to ensure that the variables are aligned by datetime.

### Usage

```
regress(primaryCRHM, primary.columns = 1, secondaryCRHM,
        secondary.columns = 1, forceOrigin = FALSE, plot = FALSE,
        quiet = TRUE, logfile = "")
```

## Arguments

<code>primaryCRHM</code>	Required. The primary <b>CRHM</b> r data frame.
<code>primary.columns</code>	Optional. Optional. A vector containing the columns to be imputed in the primary data frame, not including the datetime. If not specified, defaults to the first column.
<code>secondaryCRHM</code>	Required. The secondary <b>CRHM</b> r data frame.
<code>secondary.columns</code>	Optional. A vector containing the columns to be imputed in the secondary data frame, not including the datetime. If not specified, defaults to the first column.
<code>forceOrigin</code>	Optional. If TRUE then the regressions will be forced through the origin. Note that in this case the values of $r^2$ and the intercept will <i>not</i> be calculated as they have no meaning.
<code>plot</code>	Optional. Set <code>plot=TRUE</code> if you want a plot for each regression. Default is FALSE. The function has to guess the axis units from the variable names and it may be mistaken; you can always change the labels.
<code>quiet</code>	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet=FALSE</code> .
<code>logfile</code>	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If successful, and `plot=TRUE`, returns a plot of all of the regressions. If successful, and `plot=FALSE`, returns a data frame containing the regression parameters. If unsuccessful, returns FALSE.

## Note

If successful, exports a data frame with the columns `xvals` (secondary data frame), `yvals` (primary data frame), `first date` (in common data frame), `last date` (in common data frame), and `slope` for each variable to the regression file. If `forceOrigin=FALSE` then the parameters `intercept`, and `r2` will also be returned. This function will not permit regression on RH values, so convert these to Ea values first.

## Author(s)

Kevin Shook

## See Also

[impute](#)

## Examples

```
# this would be used to get the regression coefficients used to impute data from Saskatoon
# to fill gaps in the Regina data
## Not run:
```

```
reg <- regress(regina, c(1,2,3), saskatoon, c(1,2,3))  
## End(Not run)
```

---

**rh2vp***Converts air temperature and relative humidity to vapour pressure*

---

### Description

This is an internal **CRHMr** function and should *never* need to be called directly, as it does *NO* data checking. Use the function `changeRHtoEa` instead.

### Usage

```
rh2vp(airtemp, rh)
```

### Arguments

<code>airtemp</code>	Required. Air temperature in °C.
<code>rh</code>	Required. Relative humidity in percent.

### Value

Returns vapour pressure in kPa. If the air temperatures are scalar, returns a scalar value. If the air temperatures are a vector, returns a vector.

### Author(s)

Kevin Shook

### See Also

[changeRHtoEa](#)

### Examples

```
rh2vp(-5, 50)
```

---

ribbonPlotAirTemps	<i>Plots min, max and mean air temperatures for each day of the year</i>
--------------------	--

---

## Description

Creates **ggplot2** ribbon plots of the daily minimum, maximum and mean values of air temperatures over a year. The shaded area indicates the range of values in the data frame for each day.

## Usage

```
ribbonPlotAirTemps(obs = "", obsNames = "", tempCols = 1,
  facet = TRUE, facetColour = "", facetCols = 2)
```

## Arguments

obs	Required. Either a <b>CRHMr</b> obs data frame or a list of data frames.
obsNames	Optional. A vector containing names for the data frames. If there are more data frames than names, the data will be named obs_1, obs_2 etc.
tempCols	Optional. Column(s) containing temperatures. Can be a single value or a vector. Values will be recycled if necessary. Default is 1.
facet	Optional. If FALSE (the default), then all variables will be on a single plot. If TRUE, then each variable will be on a separate facet.
facetColour	Optional. The colour to be used for shading the ribbon plot in each facet. If not specified (the default), then each plot will be coloured separately.
facetCols	Optional. The number of columns to be used for the facets, if selected.

## Value

If unsuccessful, returns FALSE. If successful, returns a **ggplot2** object.

## Author(s)

Kevin Shook

## See Also

[plotTempsByYear](#)

## Examples

```
p <- ribbonPlotAirTemps(BadLake7376, tempCols = 1, facet=FALSE)
```



---

`runCRHM`*Runs a CRHM model*

---

### Description

Runs CRHM with a specified .prj file and (optionally) .obs file(s) and .par file(s). Under OSX and Linux, CRHM requires either 1) the program Wine to run CRHM.exe or 2) a native-compiled version. This function will use Wine if it is required. Note that none of the file names or their paths can contain spaces.

### Usage

```
runCRHM(CRHMfile = "", prjFile = "", obsFiles = "", parFiles = "",
        outFile = "", logfile = "")
```

### Arguments

CRHMfile	Required. CRHM executable file (usually CRHM.exe).
prjFile	Required. Name of .prj file. If the file does not contain any displayed observations or variables, or if any of the settings Auto_Run, Auto_Exit or Log_All are missing, then the function will terminate with an error message.
obsFiles	Optional. Name(s) of obs file(s).
parFiles	Optional. Name(s) of parameter file(s).
outFile	Optional. Name(s) of output file(s). If not specified, the original CRHM output file name (e.g. 'CRHM_output_1.txt') is kept.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Details

If Wine is required, the function will copy the specified .obs, and .par files to the directory of the .prj file, before execution of CRHM. After the program has finished, the copies will be deleted. Running a CRHM model requires that the .prj file has been setup to run automatically, as shown in the function `automatePrj`.

### Value

If successful, returns TRUE. If unsuccessful, returns FALSE.

### Author(s)

Kevin Shook

### See Also

[automatePrj](#)

**Examples**

```
## Not run:
result <- runCRHM('c:/CRHM/CRHM.exe', 'c:/BadLake/BadLake1975.prj',
outfile='c:/BadLake/BadLake1975Output.txt')
## End(Not run)
```

saturatedVP

*Converts calculates saturated vapour pressure from air temperature***Description**

This is an internal **CRHM** function and should normally not need to be called directly.

**Usage**

```
saturatedVP(airTemp)
```

**Arguments**

airTemp            Required. Air temperature in °C.

**Value**

If successful, returns the saturated vapour pressure in kPa. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**Examples**

```
saturatedVP(-5)
```

setPrjBasinName

*Sets the basin name in a CRHM model .prj file***Description**

Sets the basin name in a CRHM model .prj file

**Usage**

```
setPrjBasinName(inputPrjFile = "", basinName = "",
outputPrjFile = "", logfile = "")
```

**Arguments**

inputPrjFile	Required. Name of the .prj file.
basinName	Required. New name for basin.
outputPrjFile	Optional. If omitted, the input .prj file will be overwritten.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[runCRHM](#) [setPrjHRUnames](#) [setPrjDates](#) [setPrjParameters](#)

**Examples**

```
## Not run:
result <- setPrjBasinName('c:/CRHM/Bad Lake 1974-1975.prj', 'NewBadLakemodel')
## End(Not run)
```

---

setPrjDates	<i>Sets the model run start and end dates in a CRHM model .prj file</i>
-------------	---

---

**Description**

Sets the model run start and end dates in a CRHM model .prj file

**Usage**

```
setPrjDates(inputPrjFile = "", startDate = "", endDate = "",
  outputPrjFile = "", initialStateFile = "", finalStateFile = "",
  logfile = "")
```

**Arguments**

inputPrjFile	Required. Name of the .prj file.
startDate	Required. Model run starting date in format 'YYYY MM DD'.
endDate	Required. Model run ending date in format 'YYYY MM DD'.
outputPrjFile	Optional. If omitted, the input .prj file will be overwritten.
initialStateFile	Optional. If specified, the specified file will be set as the initial state file, i.e. the state variables to be read in at the beginning of the model run.

**finalStateFile** Optional. If specified, the specified file will be set as the final state file, i.e. the state variables to be exported at the end of the model run.

**logfile** Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns TRUE. If unsuccessful, returns FALSE.

### Note

If you are using Windows, the paths for the state files must use '\\' instead of '\' for the directory delimiter. The reason is that R, like many programming languages, interprets '\' as an escape code. This is different from the file path set in 'inputPrjFile', which must be an R path using the '/' symbol. See the example below.

### Author(s)

Kevin Shook

### See Also

[runCRHM](#) [setPrjObs](#) [setPrjParameters](#)

### Examples

```
## Not run:
result <- setPrjDates('c:/CRHM/Bad Lake 1974-1975.prj',
  '1974 10 1', '1975 10 1', initialStateFile='c:\\CRHM\\Model_Initial_State.int',
  finalStateFile='c:\\CRHM\\Model_Final_State.int')
## End(Not run)
```

---

setPrjHRUnames	<i>Sets the names of HRUs in a CRHM model .prj file</i>
----------------	---

---

### Description

Replaces the existing names of all HRUs, in a .prj file.

### Usage

```
setPrjHRUnames(inputPrjFile = "", HRUnames = "", outputPrjFile = "",
  logfile = "")
```

### Arguments

**inputPrjFile** Required. Name of the .prj file.

**HRUnames** Required. Vector containing new HRU names. There must be at least as many values in this vector as HRU names in the original .prj file.

**outputPrjFile** Optional. If omitted, the input .prj file will be overwritten.

**logfile** Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Note**

Currently, this function has only been tested on simple .prj files, i.e. without REW Grp parameters.

**Author(s)**

Kevin Shook

**See Also**

[runCRHM](#) [setPrjParameters](#) [setPrjDates](#)

**Examples**

```
## Not run:
result <- setPrjHRNames('c:/CRHM/Bad74_Frozen.prj', c('fallow', 'stubble', 'grass'))
## End(Not run)
```

---

setPrjObs

*Sets the names of the obs files in a CRHM model .prj file*


---

**Description**

Sets the names of the obs files in a CRHM model .prj file

**Usage**

```
setPrjObs(inputPrjFile = "", obsFiles = "", outputPrjFile = "",
  logfile = "")
```

**Arguments**

inputPrjFile	Required. Name of the .prj file.
obsFiles	Optional. A string or vector of strings containing the names of the obs file(s). If none are supplied, then the existing obs files will be deleted.
outputPrjFile	Optional. If omitted, the input .prj file will be overwritten.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[runCRHM](#) [setPrjDates](#) [setPrjParameters](#)

**Examples**

```
## Not run:
obsFiles <- c("BadLakeHourly.obs", "BadLakeppt.obs")
result <- setPrjObs("c:/CRHM/Bad Lake 1974-1975.prj", obsFiles)
## End(Not run)
```

---

setPrjOutputVariables *Sets the output variables in a CRHM model .prj file*

---

**Description**

Sets the output variables in a CRHM model .prj file

**Usage**

```
setPrjOutputVariables(inputPrjFile = "", variables = "",
  outputPrjFile = "", logfile = "")
```

**Arguments**

inputPrjFile	Required. Name of the .prj file.
variables	Required. Variables to be written to output file. These may be either a character vector combining the variable module, variable name and HRUs or a dataframe with the same information in 3 separate columns. Either format can be returned by setOutputVariables. If the 'null' is specified for variables, then all variables are removed.
outputPrjFile	Optional. If omitted, the input .prj file will be overwritten.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[readPrjOutputVariables](#) [setPrjDates](#)

## Examples

```
## Not run:
# read in existing variables as a vector
variables <- readPrjOutputVariables('Bad_Lake_1974-1975.prj', asDataframe=FALSE)

# delete the first set of variables and write to a new file
variables <- variables[-1]
setPrjOutputVariables('Bad_Lake_1974-1975.prj', variables, 'Bad_Lake_1974-1975_revised.prj')

# read in existing variables as a dataframe and write to new file
variables <- readPrjOutputVariables('Bad_Lake_1974-1975.prj', asDataframe=TRUE)
newVariables <- c('evap', 'evapGrangerD', '1 2 3 4 5')
variables <- rbind(variables, newVariables)
setPrjOutputVariables('Bad_Lake_1974-1975.prj', variables, 'Bad_Lake_1974-1975_revised2.prj')

# deletes all output variables and overwrites the file
result <- setPrjOutputVariables('Bad_Lake_1974-1975.prj', 'null')
## End(Not run)
```

---

setPrjParameters	<i>Sets the parameter values in a CRHM model .prj file</i>
------------------	--

---

## Description

Replaces the existing values of a single parameter, for all HRUs, in a .prj file.

## Usage

```
setPrjParameters(inputPrjFile = "", paramName = "", paramVals = "",
  outputPrjFile = "", quiet = TRUE, logfile = "")
```

## Arguments

inputPrjFile	Required. Name of the .prj file.
paramName	Required. Name of parameter to set.
paramVals	Required. Vector containing new parameter values. There must be at least as many values in this vector as parameter values in the original .prj file.
outputPrjFile	Optional. If omitted, the input .prj file will be overwritten.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Note**

Currently, this function has only been tested on simple .prj files, i.e. without REW Grp parameters. It *might* work for complex files if the 'REW Grp' specification is included in the parameter name.

**Author(s)**

Kevin Shook

**See Also**

[runCRHM setPrjDates](#)

**Examples**

```
## Not run:
result <- setPrjParameters('c:/CRHM/Bad74_Frozen.prj',
'fetch', c(1500, 1500, 1500))
## End(Not run)
```

---

setPrjRunID

*Sets the Run ID in a CRHM model .prj file*


---

**Description**

Sets the Run ID in a CRHM model .prj file

**Usage**

```
setPrjRunID(inputPrjFile = "", RunID = 0, outputPrjFile = "",
logfile = "")
```

**Arguments**

inputPrjFile	Required. Name of the .prj file.
RunID	Required. New RunID for basin. Must be an integer.
outputPrjFile	Optional. If omitted, the input .prj file will be overwritten.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook



**See Also**

[runCRHM](#) [setPrjHRUnames](#) [setPrjDates](#) [setPrjParameters](#)

**Examples**

```
## Not run:
result <- setPrjRunID('c:/CRHM/Bad Lake 1974-1975.prj', '5')
## End(Not run)
```

---

simpleDailyWater	<i>Aggregates simple CRHM model output flows to daily values</i>
------------------	--

---

**Description**

Calculates daily values of the CRHM outputs of water storages and fluxes, for models *without* sub-basins, for each HRU by its area, and calculates the net fluxes for the basin and each sub-basin. Note that this function requires a built-in dataframe `CRHM_vars` which contains information about CRHM variables. If your model uses a variable which is not in the dataframe, then you will get an error message.

**Usage**

```
simpleDailyWater(CRHMoutput, vars = "all", prjFile = "",
  basinMean = TRUE, summarize = TRUE, units = "mm", quiet = TRUE,
  logfile = "")
```

**Arguments**

CRHMoutput	Required. The CRHM model output as a standard <b>CRHM</b> r dataframe (obs, export or output data)
vars	Optional. Variable column numbers to be used (not including the datetime). The default 'all' selects all columns.
prjFile	Required. The CRHM .prj file.
basinMean	Optional. If TRUE (the default) then the fluxes and storages will be reported as depths over the entire basin. If FALSE, then fluxes and storages are reported as depths over each HRU's area.
summarize	Optional. If TRUE (the default), then a single basin-wide value of each flux and storage is returned. If FALSE then the values of the fluxes and storages are returned for each HRU.
units	Optional. The units for output. The default is 'mm', but 'm3/s' can also be specified.
quiet	Optional. Suppresses display of messages, except for errors. The default is TRUE. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful returns a data frame containing the daily values of all of the water fluxes and storages.

**Author(s)**

Kevin Shook

**See Also**

[cumulDailyWater](#)

**Examples**

```
## Not run:
b <- readOutputFile('Bologna1984_30y_output.txt', timezone='MST')
# get daily values without aggregation; values calculated w.r.t the HRU area
dailyHRU <- simpleDailyWater(b, prjFile='Bologna1984_02.prj', basinMean=FALSE)

# get daily values without aggregation; values calculated w.r.t the basin area
dailyHRU_per_basin <- simpleDailyWater(b, prjFile='Bologna1984_02.prj', basinMean=TRUE)

get daily basin values
daily_basin <- simpleDailyWater(b, prjFile='Bologna1984_02.prj', basinMean=TRUE, summarize=TRUE)

## End(Not run)
```

---

simpleMaxSolar

*Estimates extra-terrestrial shortwave radiation*

---

**Description**

Simple estimation of extra-terrestrial radiation on a horizontal plane. Daily and sub-daily values in MJ/m<sup>2</sup> and W/m<sup>2</sup> are provided,

**Usage**

```
simpleMaxSolar(datetime, latitude, hoursOffset = 2, quiet = TRUE,
  logfile = "")
```

**Arguments**

datetime	Required. Can either be a vector of <b>CRHMr</b> datetimes or any dataframe with a datetime in column 1.
latitude	Required. The latitude for which values are to calculated.
hoursOffset	Optional. The offset (in hours) is added to the solar time to convert it to local time. The default value '2' shifts the daily peak to occur at 2pm.

quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns a dataframe containing these variables:

1. datetime
2. So\_MJ, the daily extra-terrestrial radiation in MJ/m<sup>2</sup>
3. So\_W, the daily extra-terrestrial radiation in W/m<sup>2</sup>
4. R\_h\_d, the conversion from daily to hourly fluxes
5. So\_h\_W, the sub-daily extra-terrestrial radiation in W/m<sup>2</sup>
6. So\_h\_MJ, the sub-daily extra-terrestrial radiation in MJ/m<sup>2</sup>

If unsuccessful, the value FALSE will be returned.

### Note

This version is valid for latitudes between 49 °N and 55°N. It is not known how well it will perform outside this range.

### Author(s)

Kevin Shook

### References

This code is based on Shook, K., and J. Pomeroy (2011), “Synthesis of incoming shortwave radiation for hydrological simulation”, Hydrol. Res., 42(6), 433, doi:10.2166/nh.2011.074.  
Please cite this paper if you use this function in a publication.

### Examples

```
maxSolar <- simpleMaxSolar(BadLake7376, 51.366)
```

---

simpleRibbonPlot

*Simple plots of CRHM output max, min and mean values*

---

### Description

Creates **ggplot2** ribbon plots of the daily minimum, maximum and mean values for all selected variables over a year, for models *without* sub-basins. The shaded area indicates the range of values in the data frame for each day.

**Usage**

```
simpleRibbonPlot(CRHMoutput, columns = "all", waterYear = FALSE,
  ylabel = "Basin total (mm)", skipLeapDay = TRUE, facet = FALSE,
  facetColour = "blue", facetCols = 3)
```

**Arguments**

CRHMoutput	Required. A data frame containing either instantaneous or cumulative values. The data may be either daily or hourly.
columns	The columns to be aggregated, not including the datetime. The default is the 'all', but can be a vector, i.e. c(1,2,3). Note that all variables must have the same units.
waterYear	Optional. If True, the statistics are calculated and plotted by water year, if FALSE (the default) the statistics are calculated and plotted by calendar year. If you are plotting cumulative values using <code>cumulDailyWater</code> , this parameter must be set to be the same as that function.
ylabel	Optional. Label to be used for the y-axis.
skipLeapDay	Optional. If TRUE (the default) then values for Feb. 29 will be omitted from the analyses and plots. If FALSE, then these values will be included, which may result in strange-looking plots.
facet	Optional. If FALSE (the default), then all variables will be on a single plot. If TRUE, then each variable will be on a separate facet.
facetColour	Optional. The colour to be used for shading the ribbon plot in each facet, if selected.
facetCols	Optional. The number of columns to be used for the facets, if selected.

**Value**

If unsuccessful, returns FALSE. If successful, returns a **ggplot2** object.

**Author(s)**

Kevin Shook

**See Also**

[cumulDailyWater](#)

**Examples**

```
## Not run:
p <- ribbonPlot(waterYearCumul, waterYear=TRUE, units='mm', facet=TRUE, facetCols=2)
## End(Not run)
```

---

summariseObsFiles	<i>Summarises all obs files in a directory</i>
-------------------	--

---

## Description

This function summarises all of the obs files in specified directory, optionally plotting the values in each obs file in a separate graph.

## Usage

```
summariseObsFiles(file.dir = "", timezone = "",  
  summaryFile = "obsSummary.csv", plot = FALSE)
```

## Arguments

file.dir	Optional Directory containing .obs files. If not specified, defaults to current directory. Note that this is an R path, which uses the '/' symbol on ALL operating systems.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time zone, as it includes historical changes between standard and daylight savings time.
summaryFile	Optional. File to contain summary. Defaults to 'obsSummary.csv' in the same directory as the obs files.
plot	Optional. Logical. Do you want to plot graphs of each obs file? Defaults to FALSE.

## Value

If successful returns no value. If unsuccessful, returns FALSE.

## Author(s)

Kevin Shook

## See Also

[plotObs](#)

## Examples

```
## Not run:
summariseObsFiles('c:/data/BadLake/MS', timezone='etc/GMT+6', plot=TRUE)
## End(Not run)
```

---

tMinMaxToHourly	<i>Converts daily to hourly temperatures</i>
-----------------	--

---

## Description

Converts daily values of tmin, tmax and (if they exist) tmean to hourly values, by spline interpolation. Only single columns of the daily temperatures can be used.

## Usage

```
tMinMaxToHourly(obs, tmin.col = 0, tmax.col = 0, tmean.col = 0,
  tmin.time = "07:00", tmax.time = "15:00", tmean.time = "12:00",
  timezone = "", quiet = TRUE, logfile = "")
```

## Arguments

obs	Required. A <b>CRHMr</b> obs dataframe.
tmin.col	Optional. The number of the tmin column (omitting the datetime) in the obs dataframe. If omitted, the column named 'tmin' will be used. There must be a tmin column in the dataset
tmax.col	Optional. The number of the tmax column (omitting the datetime) in the obs dataframe. If omitted, the column named 'tmax' will be used. There must be a tmax column in the dataset.
tmean.col	Optional. The number of the tmax column (omitting the datetime) in the obs dataframe. If omitted, the column named 'tmean' will be searched for. This column need not be present in the dataset. If the value is set to -1, then the mean temperature will not be used in the interpolation.
tmin.time	Optional. The time of day at which the minimum air temperature occurs, specified in 'HH:MM' format. If not specified, defaults to '07:00'.
tmax.time	Optional. The time of day at which the maximum air temperature occurs, specified in 'HH:MM' format. If not specified, defaults to '15:00'.
tmean.time	Optional. The time of day at which the maximum air temperature occurs, specified in 'HH:MM' format. If not specified, defaults to '12:00'.
timezone	Required. The name of the timezone of the data as a character string. This should be the timezone of your data, but omitting daylight savings time. Note that the timezone code is specific to your OS. To avoid problems, you should use a timezone without daylight savings time. Under Linux, you can use 'CST' and 'MST' for Central Standard or Mountain Standard time, respectively. Under Windows or OSX, you can use 'etc/GMT+6' or 'etc/GMT+7' for Central Standard and Mountain Standard time. DO NOT use 'America/Regina' as the time

	zone, as it includes historical changes between standard and daylight savings time.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns an obs dataframe with the hourly air temperatures. If not successful, returns FALSE.

**Note**

Interpolation over intervals longer than 1 day is not allowed.

**Author(s)**

Kevin Shook

**See Also**

[interpolate](#)

**Examples**

```
## Not run:
MSC.hourly <- tMinMaxToHourly(MSC.daily, timezone='etc/GMT+7')
## End(Not run)
```

---

trimObs

*Trims a dataframe of missing values at the beginning and end*

---

**Description**

Removes values before 01:00 on the first complete day of the data, and after 00:00 on the last complete rows. Because it is destructive, this command should *only* be used immediately before exporting observations to a CRHM .obs file. For daily obs, removes values before the first complete row and after the last complete row.

**Usage**

```
trimObs(obs, quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> obs dataframe.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns the trimmed dataframe. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**See Also**

[writeObsFile](#)

**Examples**

```
BadLake.trimmed <- trimObs(BadLake7376)
```

---

user

*Logs information about the user*


---

**Description**

Writes information about the user and the user's computer to the logfile.

**Usage**

```
user(logfile = "")
```

**Arguments**

logfile	Optional. Name of the file to be used for logging the action. Normally not used.
---------	--

**Value**

Returns nothing.

**Note**

This function records information about your computer (the user, the operating system, the version of R, and all of the packages in use). It will make it easier to debug your problems and may help you to figure out why R code that used to work is no longer giving the correct answer.



**Author(s)**

Kevin Shook

**Examples**

```
user()
```

---

vectorsToVelocity	<i>Convert wind vectors to velocity</i>
-------------------	---

---

**Description**

Converts wind u and v vectors to a single wind velocity. The vectors may be in any units, i.e. km/h or m/s.

**Usage**

```
vectorsToVelocity(uObs, uColnum = 1, vObs, vColnum = 1, quiet = TRUE,
  logfile = "")
```

**Arguments**

uObs	Required. A <b>CRHMr</b> obs data frame containing the wind u vectors.
uColnum	Optional. The column number containing the u values, not including the date-time. Default is column 1.
vObs	Required. A <b>CRHMr</b> obs data frame containing the wind v vectors. Note that u and v must have the same units!
vColnum	Optional. The column number containing the u values, not including the date-time. Default is column 1.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used

**Value**

If successful, returns an obs data frame containing the wind speed (in the original vector units) and the wind direction in degrees from North. If unsuccessful, returns the value FALSE.

**Author(s)**

Kevin Shook

**Examples**

```
## Not run:
windspeed <- vectorsToVelocity(windObs, uColnum=1, uObs, vColnum=2)
## End(Not run)
```

vp2rh

*Converts air temperature and vapour pressure to relative humidity***Description**

This is an internal **CRHMr** function and should *never* need to be called directly, as it does *NO* data checking. Use the function `changeEatoRH` instead.

**Usage**

```
vp2rh(airtemp, vapourPressure)
```

**Arguments**

`airtemp` Required. Air temperature in °C.  
`vapourPressure` Required. Vapour pressure in kPa.

**Value**

Returns RH in percent.

**Author(s)**

Kevin Shook

**See Also**

[changeEatoRH](#)

**Examples**

```
vp2rh(-5, 0.2007355)
```

---

weighingGauge-methods *Processes weighing gauge precipitation*

---

**Description**

These functions clean accumulated precipitation data from a weighing gauge. The functions are wrapper for other R and **CRHMr** functions as well as functions adapted from MATLAB code by Alan Barr. The typical sequence of operation of the functions is indicated by their numbers, however it is NOT always necessary to execute any given function.

**weighingGaugePlot** Plots cumulative and interval precipitation

**weighingGauge1** Gap fills the data

**weighingGauge2** Removes spikes

**weighingGauge3** Applies the Savitzky-Golay polynomial filter

**weighingGauge4** Calls PcpFiltPosTh

**weighingGauge5** Simplified reset and jitter removal

**weighingGaugeInterval** Deaccumulates the precipitation

### Author(s)

Kevin Shook

### See Also

[weighingGaugePlot](#) [weighingGauge1](#) [weighingGauge2](#) [weighingGauge3](#) [weighingGauge4](#) [weighingGauge5](#)  
[weighingGaugeInterval](#)

### Examples

```
## Not run:
# check data - shows jitter, missing values and a large reset
weighingGaugePlot(wg)
# see jitter and missing values only
weighingGaugePlot(wg, endDate = '2010-10-01')

# fill gaps, and check to see if all missing values have been interpolated
wg1 <- weighingGauge1(wg, maxGapLength=10)
weighingGaugePlot(wg1)

# remove the spikes, and check
wg2 <- weighingGauge2(wg1, spikeThreshold=300, maxSpikeGap=3)
weighingGaugePlot(wg2)

# apply the smoothing filter - only for use with high-frequency data
wg3 <- weighingGauge3(wg2, filterLength=5)
weighingGaugePlot(wg3)

# apply Alan Barr's filter
wg4 <- weighingGauge4(wg3, smallDropThreshold = 0.05, serviceThreshold = -50)
weighingGaugePlot(wg4)

# deaccumulate precipitation
wg5 <- weighingGaugeInterval(wg4)

# use simplified method and plot results
# note that the infilled values are used
wg6 <- weighingGauge5(wg1)
weighingGaugePlot(wg6)
wg7 <- weighingGaugeInterval(wg6)
## End(Not run)
```

weighingGauge1

*Fills gaps in weighing gauge precipitation data***Description**

This method fills missing values in accumulated precipitation data from a weighing gauge using linear interpolation. This function is a wrapper for functions and was suggested by Craig Smith. This function need not be called if your data does not contain any gaps. It is a good idea to use `weighingGaugePlot` to look at your data before calling this function.

**Usage**

```
weighingGauge1(obs, precipCol = 1, maxGapLength = 5, quiet = TRUE,
  logfile = "")
```

**Arguments**

<code>obs</code>	Required. A standard <b>CRHMr</b> obs file.
<code>precipCol</code>	Optional. The number of the column containing the weighing gauge cumulative precipitation data, not including the datetime. Default is column 1. These values are cumulative precipitation.
<code>maxGapLength</code>	Optional. The maximum gap length included in the analysis (in time intervals). Default is 5 time intervals. Used by <code>interpolate</code> .
<code>quiet</code>	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet=FALSE</code> .
<code>logfile</code>	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If unsuccessful, returns `FALSE`. If successful, returns a modified version of the obs dataframe containing only the datetime and the gap-filled total precipitation values.

**Note**

This version does not explicitly incorporate the effects of evaporation.

**Author(s)**

Kevin Shook

**See Also**

[weighingGauge2](#) [weighingGauge3](#) [weighingGauge4](#) [weighingGaugePlot](#) [weighingGaugeInterval](#)

## Examples

```
## Not run:
test1 <- weighingGauge1(wg, maxGapLength=6)
## End(Not run)
```

---

weighingGauge2

*Removes spikes from weighing gauge cumulative precipitation*


---

## Description

Removes positive and negative spikes from deaccumulated weighing gauge precipitation. The gaps left from spike removal are infilled by linear interpolation. This function is a wrapper for other functions and was suggested by Craig Smith. This function need not be called if your data does not contain any spikes. It is a good idea to use weighingGaugePlot to look at your data before calling this function. Note that this function will also remove gauge resets, if they are within the range of the spikeThreshold.

## Usage

```
weighingGauge2(obs, precipCol = 1, spikeThreshold = 1000,
  maxSpikeGap = 3, quiet = TRUE, logfile = "")
```

## Arguments

obs	Required. A standard <b>CRHMr</b> obs file.
precipCol	Optional. The number of the column containing the weighing gauge cumulative precipitation data, not including the datetime. Default is column 1. This function can only work on 1 column of precipitation.
spikeThreshold	Optional. Threshold for single-interval precipitation (mm). Any spikes (changes in cumulative precipitation) whose absolute value is greater than the threshold will be deleted. The default value is 1000 mm.
maxSpikeGap	Optional. Maximum length of spikes (in time steps) to be filled. Default is 3.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

## Value

If unsuccessful, returns FALSE. If successful, returns a modified version of the obs dataframe containing only the datetime and the adjusted precipitation values.

## Note

If you don't specify a large enough value for maxSpikeGap, then all of the values after the spike will be set to NA.

**Author(s)**

Kevin Shook

**See Also**[weighingGauge1](#) [weighingGauge3](#) [weighingGauge4](#) [weighingGaugePlot](#) [weighingGaugeInterval](#)**Examples**

```
## Not run:
test2 <- weighingGauge2(wg, spikeThreshold = 300)
## End(Not run)
```

weighingGauge3

*Filters weighing gauge cumulative precipitation***Description**

This function is a wrapper for the Savitzky-Golay polynomial filter function `sgolayfilt` from the package **signal**. It passes the specified cumulative precipitation data, the filter length, the filter order (set to 1) and the derivative order (set to 0) to the filter. It is a good idea to use `weighingGaugePlot` to look at your data before and after calling this function. Note that this function is only intended to be used with very high frequency (i.e. 1 minute) gauged data, and should not be used with longer-period accumulations.

**Usage**

```
weighingGauge3(obs, precipCol = 1, filterLength = 0, quiet = TRUE,
  logfile = "")
```

**Arguments**

<code>obs</code>	Required. A standard <b>CRHMr</b> obs file.
<code>precipCol</code>	Optional. The number of the column containing the weighing gauge cumulative precipitation data, not including the datetime. Default is column 1. This function can only work on 1 column of precipitation.
<code>filterLength</code>	Required. The number of time intervals used by the Savitzky-Golay polynomial filter. Note that this must be an ODD number. The appropriate value will depend on the type of your data, and the noise present.
<code>quiet</code>	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet=FALSE</code> .
<code>logfile</code>	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If unsuccessful, returns FALSE. If successful, returns a modified version of the obs dataframe containing the adjusted precipitation values. Note that the name of the precipitation variable is modified: '\_sg\_filtered' is appended.

**Note**

This function is potentially destructive, as it smooths the accumulated precipitation. You may need to set the filter length by trial-and error. You can check the effects using the function weighingGaugePlot.

**Author(s)**

Kevin Shook

**See Also**

[weighingGauge1](#) [weighingGauge2](#) [weighingGauge4](#) [weighingGaugePlot](#) [weighingGaugeInterval](#)

**Examples**

```
## Not run:
test3 <- weighingGauge3(wg, filterLength=5)
## End(Not run)
```

---

weighingGauge4

*Runs Alan Barr's filter on weighing gauge cumulative precipitation*

---

**Description**

This function is a wrapper for Alan Barr's function PcpFiltPosTh which removes jitter from precipitation data. The called function is iterative and may be slow to execute.

**Usage**

```
weighingGauge4(obs, precipCol = 1, smallDropThreshold = 0.1,
  serviceThreshold = -100, serviceGapLength = 3, quiet = TRUE,
  logfile = "")
```

**Arguments**

obs	Required. A standard <b>CRHMr</b> obs file.
precipCol	Optional. The number of the column containing the weighing gauge cumulative percipitaion data, not including the datetime. Default is column 1. This function can only work on 1 column of precipitation.
smallDropThreshold	Optional. Minimum interval precipitation (mm). Default is 0.1.

serviceThreshold	Optional. Threshold for removing change in storage due to servicing (mm). Note that this must be a NEGATIVE value. Default is -100.
serviceGapLength	Optional. Maximum gap length (in time intervals) to be filled by linear interpolation after removing servicing (i.e. gauge reset) data. The default value is 3, but you may have to adjust it for very high-frequency data.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If unsuccessful, returns FALSE. If successful, returns a modified version of the obs dataframe containing the adjusted precipitation values. Note that the name of the precipitation variable is modified: '\_PcpFiltPost' is appended.

**Note**

This function is potentially destructive, as it smooths the accumulated precipitation. You may need to set the filter length by trial-and error. You can check the effects using the function `weighingGaugePlot`. This version does not explicitly incorporate the effects of evaporation.

**Author(s)**

Kevin Shook

**See Also**

[PcpFiltPosTh](#) [weighingGauge1](#) [weighingGauge2](#) [weighingGauge3](#) [weighingGaugePlot](#) [weighingGaugeInterval](#)

**Examples**

```
## Not run:
test4 <- weighingGauge4(wg, smallDropThreshold = 0.05, serviceThreshold = -50)
## End(Not run)
```

---

weighingGauge5

*Simplified weighing gauge processing*

---

**Description**

This function removes gauge resets (large negative changes usually due to servicing) and uses a simple method (cumulative maxima) to remove jitter. It is a good idea to use `weighingGauge1` to infill missing values before calling this function and `weighingGaugePlot` to look at your data before and after calling this function.



**Usage**

```
weighingGauge5(obs, precipCol = 1, resetThreshold = 50, quiet = TRUE,  
  logfile = "")
```

**Arguments**

obs	Required. A standard <b>CRHMr</b> obs file.
precipCol	Optional. The number of the column containing the weighing gauge cumulative precipitation data, not including the datetime. Default is column 1. These values are cumulative precipitation. This function can only work on 1 column of precipitation. The precipitation cannot contain missing values - infill them using <code>weighingGauge1</code> first.
resetThreshold	Optional. The minimum (absolute value) of negative values considered to be gauge resets. Default is 50 mm.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet=FALSE</code> .
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If unsuccessful, returns `FALSE`. If successful, returns a modified version of the obs dataframe containing only the datetime and the modified cumulative precipitation values.

**Note**

This version does not explicitly incorporate the effects of evaporation.

**Author(s)**

Kevin Shook, based on an algorithm developed by Thai Nguyen of Alberta Environment.

**See Also**

[weighingGauge1](#) [weighingGaugePlot](#) [weighingGaugeInterval](#)

**Examples**

```
## Not run:  
wg1 <- weighingGauge1(wg, maxGapLength=500)  
wg5 <- weighingGauge5(wg1, resetThreshold=70)  
## End(Not run)
```

---

weighingGaugeInterval *Deaccumulates weighing gauge precipitation*

---

### Description

Converts weighing gauged cumulative precipitation to interval values. The interval precipitation is checked to make sure that it does not contain any negative values. Negative precipitation will result in the function displaying an error message and returning the value FALSE.

### Usage

```
weighingGaugeInterval(obs, precipCol = 1, quiet = TRUE, logfile = "")
```

### Arguments

obs	Required. A standard <b>CRHMr</b> obs file.
precipCol	Optional. The number of the column containing the weighing gauge cumulative precipitation data, not including the <code>datetime</code> . Default is column 1.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave <code>quiet=TRUE</code> (i.e. the default). If you are working interactively, you will probably want to set <code>quiet=FALSE</code> .
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If unsuccessful, returns FALSE. If successful, returns a modified version of the obs dataframe containing the `datetime`, the cumulative and interval precipitation.

### Author(s)

Kevin Shook

### See Also

[weighingGauge1](#) [weighingGauge2](#) [weighingGauge3](#) [weighingGauge4](#) [weighingGauge5](#) [weighingGaugePlot](#)

### Examples

```
## Not run:
test5 <- weighingGaugeStep5(wg)
## End(Not run)
```

---

weighingGaugePlot	<i>Plots weighing gauge precipitation</i>
-------------------	---

---

## Description

Plots cumulative precipitation data from a weighing gauge. The plot contains facets with the cumulative and interval precipitations plotted against time.

## Usage

```
weighingGaugePlot(obs, precipCol = 1, startDate = "", endDate = "",  
  showMissing = TRUE)
```

## Arguments

obs	Required. A standard <b>CRHMr</b> obs dataframe.
precipCol	Optional. The column containing the cumulative precipitation data. Default is column 1.
startDate	Optional. The starting date for the plot. Can either be a year as a number (e.g. 1995) or a date in 'Y-m-d' format, i.e. '1995-06-01'. If not specified, the first datetime value in the obs data is used.
endDate	Optional. The ending date for the plot. Can either be a year as a number (e.g. 1995) or a date in 'Y-m-d' format, i.e. '1995-12-31'. If not specified, the first datetime value in the obs data is used.
showMissing	Optional. If set to TRUE (the default), then the position of missing values will be indicated by red points on the x-axis.

## Value

If successful returns a ggplot2 object. If unsuccessful, returns the value FALSE.

## Author(s)

Kevin Shook.

## See Also

[weighingGauge1](#) [weighingGauge2](#) [weighingGauge3](#) [weighingGauge4](#) [weighingGauge5](#)

## Examples

```
## Not run:  
p <- weighingGaugePlot(test1)  
## End(Not run)
```

---

wellHydrograph	<i>Plots hydrograph of daily well depths</i>
----------------	--

---

## Description

Creates a **ggplot2** object of hydrograph of well depths.

## Usage

```
wellHydrograph(wellObs, depthCol = 1, fakeDates = FALSE,
  hydroYear = TRUE, startMonth = 10, quiet = TRUE)
```

## Arguments

wellObs	Required. <b>CRHMr</b> obs data frame of daily depths.
depthCol	Optional. Column containing depths. Default is 1.
fakeDates	Optional. If set to TRUE then all dates have their year replaced with 2000, and the actual year is added as a variable in the plotted data. This allows the plot to be faceted by year, as shown in the examples.
hydroYear	Optional. If TRUE, the depths are plotted against the hydrological year.
startMonth	Optional. Sets the first month of the hydrological year (if required).
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.

## Value

If successful, returns a **ggplot2** object. If unsuccessful, returns FALSE.

## Author(s)

Kevin Shook

## Examples

```
## Not run:
p <- wellHydrograph(well, fakeDates = FALSE)
#' # re-plot with fake dates
p <- wellHydrograph(well, fakeDates = TRUE, hydroYear = TRUE)
p <- p + ggplot2::facet_wrap(~year)
p <- p + ggplot2::scale_x_date(date_breaks = "3 months",
  date_labels = "%b")

## End(Not run)
```

wetland-methods

*Processes CRHM wetland (depressional-storage) model outputs***Description**

These methods are used to process output from CRHM models which contain many HRUs functioning as depressional storage, particularly where the Pothole Cascade Model (PCM) is implemented. Note that the term *pond* refers to the water within a depression.

**wetland\_h\_s** Calculates h and s parameters for the H-v relationship

**wetland\_pond\_area** Estimates the area(s) of pond(s) from the H-v relationship

**wetland\_pond\_depth** Estimates the max depth(s) of pond(s) from the H-v relationship

**wetland\_pond\_size** Estimates the sizes(depths, areas and perimeters) of all ponds for each time step from CRHM output

wetland\_h\_s

*Calculates Hayashi-van der Kamp parameters for a Prairie depressions***Description**

Computing pond area and depth for Prairie depressions from the Hayashi-van der Kamp method requires values of the parameters h and s. Unfortunately, these values can't be solved explicitly from the pond max area and volume. This function finds the values of the parameters by iterative solution.

**Usage**

```
wetland_h_s(max_volume, max_area, p, max_iterations = 100,
            max_error = 1e-04, return_stats = FALSE)
```

**Arguments**

max_volume	Required. Maximum volume of a pond in m <sup>3</sup> .
max_area	Required. Maximum area of a pond in m <sup>2</sup> .
p	Required. Scaling exponent for pond depth-area relationship. See reference.
max_iterations	Optional. Maximum number of iterations for solution. Default is 100.
max_error	Optional. Maximum error for estimation of parameters, as a fraction of max_area.
return_stats	If TRUE, then the error of solution and the number of iterations will also be returned. Default is FALSE Default is 0.001.

**Value**

If successful, returns a vector of the parameters h (the maximum pond depth) in m and s (the pond area for a depth of 1 m) in m<sup>2</sup>). Optionally returns the error and number of iterations required for solution. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**References**

Hayashi, M., and G. van der Kamp (2000), *Simple equations to represent the volume–area–depth relations of shallow wetlands in small topographic depressions*, *J. Hydrol.*, 237(1–2), 74–85, doi:10.1016/S0022-1694(00)00300-0.

**See Also**

Other wetland\_functions: [wetland\\_pond\\_area](#), [wetland\\_pond\\_depth](#), [wetland\\_pond\\_size](#)

**Examples**

wetland\_h\_s(5000, 500, 1.72)

---

wetland_pond_area	<i>Calculates Hayashi-van der Kamp area of water in a Prairie pond</i>
-------------------	--

---

**Description**

Calculates Hayashi-van der Kamp area of water in a Prairie pond

**Usage**

wetland\_pond\_area(pond\_depth = NULL, p = NULL, s = NULL)

**Arguments**

- |            |  |
|------------|--|
| pond_depth | Required. Depth of water in pond (in m). May be a vector of values. Note that in this context the depth refers to the deepest point in the pond. |
| p          | Required. Value of p exponent for ponds. May be a vector of values.  |
| s          | Required. Value of s parameter (pond area for a depth of 1 m) in m <sup>2</sup> ) for ponds. May be a vector of values.                          |

**Value**

If successful, returns a vector of the pond areas, for the given depths. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**References**

Hayashi, M., and G. van der Kamp (2000), Simple equations to represent the volume–area–depth relations of shallow wetlands in small topographic depressions, *J. Hydrol.*, 237(1–2), 74–85, doi:10.1016/S0022-1694(00)00300-0.

**See Also**

Other wetland\_functions: [wetland\\_h\\_s](#), [wetland\\_pond\\_depth](#), [wetland\\_pond\\_size](#)

**Examples**

```
wetland_pond_area(1, 1.72, 25)
```

---

wetland_pond_depth	<i>Calculates Hayashi-van der Kamp maximum depth of water in a Prairie pond</i>
--------------------	---

---

**Description**

Calculates Hayashi-van der Kamp maximum depth of water in a Prairie pond

**Usage**

```
wetland_pond_depth(pond_volume = NULL, p = NULL, s = NULL)
```

**Arguments**

pond_volume	Required. Volume of water in ponds (in m <sup>3</sup> ). May be a vector of values.
p	Required. Value of p exponent for ponds. May be a vector of values.
s	Required. Value of s parameter (pond area for a depth of 1 m) in m <sup>2</sup> ) for pond. May be a vector of values.

**Value**

If successful, returns a vector of the pond maximum depths, for the given volumes. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

References

Hayashi, M., and G. van der Kamp (2000), Simple equations to represent the volume–area–depth relations of shallow wetlands in small topographic depressions, *J. Hydrol.*, 237(1–2), 74–85, doi:10.1016/S0022-1694(00)00300-0.

See Also

Other wetland\_functions: [wetland\\_h\\_s](#), [wetland\\_pond\\_area](#), [wetland\\_pond\\_size](#)

Examples

```
wetland_pond_depth(100, 1.72, 100)
```

---

wetland_pond_size	<i>Estimates pond area, depth and perimeter (circular)</i>
-------------------	--

---

Description

Estimates pond area, depth and perimeter (circular)

Usage

```
wetland_pond_size(CRHM_output = NULL, HRU_name = NULL,  
  HRU_SD_max = NULL, HRU_area = NULL, wetland_HRU = NULL,  
  area_p_vals = NULL, logfile = "")
```

Arguments

CRHM_output	Required. Data frame of CRHM outputs including Sd values for all wetland HRUs.
HRU_name	Required. Names of all HRUs. Can be obtained using readPrjParameters.
HRU_SD_max	Required. SDmax values for all HRUs. Can be obtained using readPrjParameters.
HRU_area	Required. Areas of all HRUs. Can be obtained using readPrjParameters.
wetland_HRU	Optional. A vector of the number of each wetland HRUs.
area_p_vals	Optional. A data frame containing two columns. The first column is the wetland area (in m <sup>2</sup> ), the second column is the corresponding value of p to use. If this parameter is not specified, then the default values will be used. The default is to use p = 1.72 for areas greater than 0 and smaller than 10,000 m <sup>2</sup> , and p = 3.3 for areas greater than or equal to 10,000 m <sup>2</sup> . These values were estimated for Smith Creek basin.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

Value

If successful, returns a data frame consisting of the datetime (or date), and the pond area (m<sup>2</sup>), depth (m), and perimeter (m) (assuming the pond is circular) for each time step in CRHM\_output.



**Note**

Because this function may be slow and generates a very large quantity of output, it is a good idea to restrict the range of CRHM\_output *before* calling the function.

**Author(s)**

Kevin Shook

**See Also**

Other wetland\_functions: [wetland\\_h\\_s](#), [wetland\\_pond\\_area](#), [wetland\\_pond\\_depth](#)

**Examples**

```
## Not run:
CRHM_output <- readOutputFile("CRHM_output_1.txt", "Etc/GMT+6")
HRU_name <- readPrjParameters("SmithCreek.prj", "hru_names")
HRU_SD_max <- readPrjParameters("SmithCreek.prj", "Sdmax")
HRU_area <- readPrjParameters("SmithCreek.prj", "hru_area")
area <- c(0, 5000, 10000)
p <- c(1.4, 1.7, 3.3)
area_p_vals <- data.frame(area, p)
pond_sizes <- wetland_pond_size(CRHM_output, HRU_name, HRU_SD_max, HRU_area, area_p_vals)
## End(Not run)
```

---

wg

*Synthetic weighing precipitation gauge data*


---

**Description**

A dataframe containing data for use with the weighingGauge functions.

**Usage**

wg

**Format**

A dataframe with 17460 rows and 2 variables (including the datetime):

**datetime** date and time as a POSIXct object

**p** cumulative precipitation (mm)

**Source**

This data was derived from real data provided by Craig Smith. It has been adjusted by incorporating missing values and a simulated gauge reset.

---

win.eol

*Gets the Windows end of line characters*


---

### Description

Finds the end of line (eol) characters required for writing Windows files, such as CRHM obs files. No parameters are required. This is an internal **CRHMr** function and should *never* need to be called directly.

### Usage

```
win.eol()
```

### Value

Returns the Windows end of line characters (cr and lf).

### Note

This function is used to make the creation of Windows-specific files work on all platforms. CRHM requires its obs and project files to use the Windows end of line characters, which are expressed differently on UNIX-based operating systems such as Linux and OSX.

### Author(s)

Kevin Shook

### See Also

[automatePrj](#) [setPrjDates](#) [writeObsFile](#) [runCRHM](#)

### Examples

```
windowsEndOfLine <- win.eol()
```

---

writeChangeLogFile

*Writes information about changes to a **CRHMr** dataframe to a file*


---

### Description

This is an internal **CRHMr** function and should normally not need to be called directly. This function writes a file containing the a summary of the files as well as the changed data values and a description of their source.

**Usage**

```
writeChangeLogFile(action, original.data.info, changed.data.info,  
  changed.data.type, comment1 = "", comment2 = "", comment3 = "",  
  comment4 = "", quiet)
```

**Arguments**

action	Required. A character string indicating the action taken - ususally this is the name of the function making the changes.
original.data.info	Required. Information about the original dataframe from CRHM_summary.
changed.data.info	Required. Information about the changed dataframe from CRHM_summary.
changed.data.type	Required. A vector indicating the type/source of data in the changed dataframe.
comment1	Optional. A comment about the data and/or changes.
comment2	Optional. A comment about the data and/or changes.
comment3	Optional. A comment about the data and/or changes.
comment4	Optional. A comment about the data and/or changes.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.

**Value**

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Author(s)**

Kevin Shook

**Examples**

```
## Not run:  
writeChangeLogFile('appendObs', obs1.info, obs2.info, output.type,  
  'primary data: obs1', 'secondary data: obs2')  
## End(Not run)
```

---

writeCLASSfile	<i>Writes forcing data to a CLASS file</i>
----------------	--

---

### Description

Writes a standard **CRHMr** data frame to a CLASS forcing data file.

### Usage

```
writeCLASSfile(CLASS, CLASSfile = "", logfile = "")
```

### Arguments

CLASS	Required. A <b>CRHMr</b> data frame containing the fields SWin, LWin, Prec, T, qa, U, and P, in addition to the datetime.
CLASSfile	Required. A the file to hold the CLASS data.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

### Value

If successful, returns the value TRUE. If unsuccessful, returns the value FALSE.

### Examples

```
## Not run:
result <- writeCLASSfile(class, 'driv.met')
## End(Not run)
```

---

writeObsFile	<i>Writes a dataframe to a .obs file</i>
--------------	--

---

### Description

Writes a CRHM dataframe to an obs file. It will not allow ea values - they must be converted to RH values first. The CRHM filter command to convert RH values to ea is added automatically to the header of the .obs file.

### Usage

```
writeObsFile(obs, obsfile = "", comment = "", quiet = TRUE,
  logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> obs data frame.
obsfile	Required. Name for file to hold observations.
comment	Optional. A comment to be added to the header of the .obs file. If omitted, the header will consist of the name of the obs data frame followed by 'created by CRHMr function writeObsFile'.
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If successful, returns TRUE. If unsuccessful, returns FALSE.

**Note**

If there are missing values in the obs dataframe, the last line of the header will contain the line '\$\$ Missing'.

**Author(s)**

Kevin Shook

**See Also**

[trimObs](#) [changeEatoRH](#)

**Examples**

```
# write sample obs dataframe to .obs file
result <- writeObsFile(BadLake7376, 'BadLake7376.obs')
```

---

yearlyPeaks

*Finds datetimes of annual max values*


---

**Description**

Finds the annual maxima (and, optionally, minima) of the specified columns in a CRHM dataframe and their corresponding datetimes.

**Usage**

```
yearlyPeaks(CRHMrdata, columns = 0, minvals = FALSE, logfile = "")
```

**Arguments**

CRHMdata	Required. A <b>CRHM</b> r dataframe.
columns	Optional. A vector containing the columns to be analyzed, not including the datetime column. If not specified, all data columns are used.
minvals	Optional. Logical. Specifies if yearly minimum values should also be found. Default value is FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

Returns a dataframe with maximum (and minimum if selected) values, and their corresponding datetimes, for each year for each selected variable.

**Author(s)**

Kevin Shook

**See Also**

[aggDataframe](#)

**Examples**

```
## Not run:
p <- yearlyPeaks(BrandonObs, columns=c(1,2), minvals=TRUE)
## End(Not run)
```

---

zeroMissingPrecip	<i>Sets missing precipitation values to zero</i>
-------------------	--

---

**Description**

Removes missing precipitation (NA\_real\_values (p and/or ppt) in an obs dataframe, by setting them to zero. This function is potentially dangerous, as setting the missing values to zero prevents any further imputation/interpolation of values.

**Usage**

```
zeroMissingPrecip(obs, precipCols = 0, quiet = TRUE, logfile = "")
```

**Arguments**

obs	Required. A <b>CRHMr</b> obs dataframe.
precipCols	Optional. A vector containing the numbers of the columns (excluding the date-time) to be processed. If omitted, all of the columns whose names contain 'p' or 'ppt' will be processed
quiet	Optional. Suppresses display of messages, except for errors. If you are calling this function in an R script, you will usually leave quiet=TRUE (i.e. the default). If you are working interactively, you will probably want to set quiet=FALSE.
logfile	Optional. Name of the file to be used for logging the action. Normally not used.

**Value**

If the action is successful, returns a dataframe with all missing precipitation values set to zero. If the action is unsuccessful, then the value FALSE is returned. If quiet=FALSE, then a summary of the returned obs data will be returned.

**Note**

As described, this function is potentially dangerous, and its action *CANNOT* be undone. It should only be used when there is no further hope for imputing/interpolating precipitation data. Note that its use is really only justified in arid or semi-arid climates, where precipitation is rare. Because it is so destructive, the function should only be used immediately before exporting to an obs file.

**Author(s)**

Kevin Shook

**See Also**

[writeObsFile](#)

**Examples**

```
BadLake7376.clean <- zeroMissingPrecip(BadLake7376, quiet=FALSE)
```

# Index

## \*Topic **datasets**

- BadLake7376, [10](#)
- CRHM\_vars, [16](#)
- wg, [113](#)
  
- aggDataframe, [5](#), [22](#), [42](#), [53](#), [118](#)
- appendObs, [7](#)
- assembleObs, [8](#)
- automatePrj, [9](#), [81](#), [114](#)
  
- BadLake7376, [10](#)
- basinWaterBalancePlot, [10](#)
  
- changeEatoRH, [11](#), [13](#), [98](#), [117](#)
- changeRHtoEa, [12](#), [12](#), [79](#)
- convertDPtoRH, [13](#)
- createObsDataframe, [14](#)
- CRHM\_summary, [15](#)
- CRHM\_vars, [16](#)
- CRHMr-package, [4](#)
- cumsumDataframe, [17](#)
- cumulativeDischargePlot, [18](#)
- cumulDailyWater, [17](#), [19](#), [90](#), [92](#)
  
- dailyByHourlyP, [20](#)
- data.frame, [46](#)
- datetimeToDate, [21](#), [22](#)
- dateToDatetime, [21](#), [22](#)
- deDupe, [23](#)
- deleteSpikes, [24](#), [37](#)
- deltaStorage, [25](#)
- distributeInst, [26](#), [27](#)
- distributeMean, [26](#), [27](#)
- distributeP, [28](#), [43](#), [45](#)
- distributeQli, [29](#), [31](#), [33](#), [49](#)
- distributeQsi, [30](#)
- doubleMass, [32](#)
  
- emissivity, [30](#), [33](#), [49](#)
- extendObs, [33](#)
  
- findDupes, [23](#), [34](#), [37](#)
- findGaps, [35](#), [35](#), [37](#), [44](#)
- findSpikes, [24](#), [36](#)
  
- GMToffset, [37](#)
  
- hruGroupWaterSummary, [38](#)
- hydrograph, [19](#), [39](#)
- hydroYear, [6](#), [41](#)
  
- impute, [6](#), [15](#), [25](#), [29](#), [35](#), [36](#), [42](#), [43–45](#), [78](#)
- insertMissing, [36](#), [43](#)
- interpolate, [6](#), [25](#), [29](#), [30](#), [35](#), [36](#), [43](#), [44](#), [44](#), [95](#)
  
- keepGood, [45](#)
  
- list2df, [46](#)
- logAction, [48](#)
- longwave, [30](#), [33](#), [48](#)
  
- makeRegular, [49](#)
- maxObs, [50](#), [52](#)
- minObs, [51](#), [51](#)
- monthlyPrecipTotals, [53](#), [54](#)
- monthlyQQplot, [53](#), [54](#)
  
- nancumsum, [55](#)
  
- parseNums, [55](#)
- parseText, [56](#)
- PcpFiltPosTh, [55](#), [57](#), [104](#)
- phaseCorrect, [58](#)
- plotObs, [59](#), [93](#)
- plotPrecipsByYear, [60](#), [62](#)
- plotTempsByYear, [61](#), [61](#), [80](#)
  
- qair2ea, [62](#), [64](#)
- qair2rh, [62](#), [63](#)
- qqplotValues, [64](#)
  
- readCampbell, [65](#)



readCLASSfile, 66  
readDebugFile, 67  
readExportFile, 4, 66, 68, 69, 70  
readObsFile, 4, 66, 68, 69, 70  
readOutputFile, 4, 66, 68, 69, 70, 71  
readOutputUnits, 70, 71  
readPrjDimensions, 72  
readPrjHRUnames, 72  
readPrjModuleParams, 73  
readPrjNumVals, 74  
readPrjOutputVariables, 75, 86  
readPrjParameters, 72, 76  
readPrjTextVals, 77  
regress, 43, 77  
rh2vp, 79  
ribbonPlotAirTemps, 80  
runCRHM, 9, 81, 83–86, 88, 89, 114  
  
saturatedVP, 82  
setPrjBasinName, 82  
setPrjDates, 83, 83, 85, 86, 88, 89, 114  
setPrjHRUnames, 73, 83, 84, 89  
setPrjObs, 84, 85  
setPrjOutputVariables, 75, 76, 86  
setPrjParameters, 83–86, 87, 89  
setPrjRunID, 88  
simpleDailyWater, 19, 25, 89  
simpleMaxSolar, 31, 90  
simpleRibbonPlot, 19, 91  
summariseObsFiles, 60, 93  
  
tMinMaxToHourly, 94  
trimObs, 46, 95, 117  
  
user, 4, 96  
  
vectorsToVelocity, 97  
vp2rh, 98  
  
weighingGauge-methods, 98  
weighingGauge1, 99, 100, 102–107  
weighingGauge2, 99, 100, 101, 103, 104, 106, 107  
weighingGauge3, 99, 100, 102, 102, 104, 106, 107  
weighingGauge4, 99, 100, 102, 103, 103, 106, 107  
weighingGauge5, 99, 104, 106, 107  
weighingGaugeInterval, 99, 100, 102–105, 106  
  
weighingGaugePlot, 99, 100, 102–106, 107  
wellHydrograph, 108  
wetland-methods, 109  
wetland\_h\_s, 109, 111–113  
wetland\_pond\_area, 110, 110, 112, 113  
wetland\_pond\_depth, 110, 111, 111, 113  
wetland\_pond\_size, 110–112, 112  
wg, 113  
win.eol, 114  
writeChangeLogFile, 114  
writeCLASSfile, 116  
writeObsFile, 35, 96, 114, 116, 119  
  
yearlyPeaks, 6, 117  
  
zeroMissingPrecip, 118