

Package: GENERead (via r-universe)

September 22, 2026

Type Package
Title Package for Reading Binary Files
Version 2.0.10
Date 2024-03-05
Maintainer Jia Ying Chua <jiayingc@activinsights.com>
Description Functions and analytics for GENE-compatible accelerometer data into R objects. See topic 'GENERead' for an introduction to the package. See <<https://activinsights.com/technology/geneactiv/>> for more details on the GENEActiv device.
License GPL-2 | GPL-3
Encoding UTF-8
Depends bitops, mmap
Suggests MASS, methods
RoxygenNote 7.3.1
NeedsCompilation yes
Author Zhou Fang [aut], Joss Langford [aut], Charles Sweetland [aut], Jia Ying Chua [aut, cre]
Repository <https://ai-jyc.r-universe.dev>
Date/Publication 2024-03-07 02:27:01 UTC
RemoteUrl <https://github.com/cran/GENERead>
RemoteRef HEAD
RemoteSha 92870d1fc3791059f70c28b64c387599f776023d

Contents

GENERead-package	2
AccData	6
convert.time	7
epoch	7

GENEActiv.calibrate	9
get.intervals	11
GRtime	13
hanning.window	14
header.info	15
parse.time	16
plot.stft	17
print.stft	20
recalibrate	20
stft	22
svm	25
uniform.window	26

Index	27
--------------	-----------

GENEAread-package	<i>A package to process binary accelerometer output files.</i>
-------------------	--

Description

This is a package to process binary output files from the GENE A accelerometer data.

The main functions are:

- read.bin
- stft
- epoch

A function to process binary accelerometer files and convert the information into R objects.

Usage

```
read.bin(
  binfile,
  outfile = NULL,
  start = NULL,
  end = NULL,
  Use.Timestamps = FALSE,
  verbose = TRUE,
  do.temp = TRUE,
  do.volt = TRUE,
  calibrate = TRUE,
  downsample = NULL,
  blocksize,
  virtual = FALSE,
  mmap.load = (.Machine$sizeof.pointer >= 8),
  pagerefs = TRUE,
  ...
)
```

Arguments

<code>binfile</code>	A filename of a file to process.
<code>outfile</code>	An optional filename specifying where to save the processed data object.
<code>start</code>	Either: A representation of when in the file to begin processing, see Details.
<code>end</code>	Either: A representation of when in the file to end processing, see Details.
<code>Use.Timestamps</code>	To use timestamps as the start and end time values this has to be set to TRUE. (Default FALSE)
<code>verbose</code>	A boolean variable indicating whether some information should be printed during processing should be printed.
<code>do.temp</code>	A boolean variable indicating whether the temperature signal should be extracted
<code>do.volt</code>	A boolean variable indicating whether the voltage signal should be extracted.
<code>calibrate</code>	A boolean variable indicating whether the raw accelerometer values and the light variable should be calibrated according to the calibration data in the headers.
<code>downsample</code>	A variable indicating the type of downsampling to apply to the data as it is loaded. Can take values: NULL: (Default) No downsampling Single numeric: Reads every downsample-th value, starting from the first. Length two numeric vector: Reads every downsample[1]-th value, starting from the downsample[2]-th. Non-integer, or non-divisor of 300 downsampling factors are allowed, but will lead to imprecise frequency calculations, leap seconds being introduced, and generally potential problems with other methods. Use with care.
<code>blocksize</code>	Integer value giving maximum number of data pages to read in each pass. Defaults to 10000 for larger data files. Sufficiently small sizes will split very large data files to read chunk by chunk, reducing memory requirements for the <code>read.bin</code> function (without affecting the final object), but conversely possibly increasing processing time. Can be set to Inf for no splitting.
<code>virtual</code>	logical. If set TRUE, do not do any actual data reading. Instead construct a <code>VirtualAccData</code> object containing header information to allow use with <code>get.intervals</code>
<code>mmap.load</code>	Default is <code>(.Machine\$sizeof.pointer >= 8)</code> . see mmap for more details
<code>pagerefs</code>	A variable that can take two forms, and is considered only for <code>mmap.load = TRUE</code> NULL or FALSE, in which case <code>pagerefs</code> are dynamically calculated for each record. (Default) A vector giving sorted byte offsets for each record for <code>mmap</code> reading of data files. TRUE, in which case a full page reference table is computed before any processing occurs. Computing <code>pagerefs</code> takes a little time and so is a little slower. However, it is safer than dynamic computations in the case of missing pages and high temperature variations. Further, once page references are calculated, future reads are much faster, so long as the previously computed references are supplied.
<code>...</code>	Any other optional arguments can be supplied that affect manual calibration and data processing. These are: • <code>mmap</code>: logical. If TRUE (Default on 64bit R), use the <code>mmap</code> package to process the <code>binfile</code>

- gain: a vector of 3 values for manual gain calibration of the raw (x,y,z) axes. If gain=NULL, the gain calibration values are taken from within the output file itself.
- offset: a vector of 3 value for manual offset calibration of the raw (x,y,z) axes. If offset=NULL, the offset calibration values are taken from within the output file itself.
- luxv: a value for manual lux calibration of the light meter. If luxv=NULL, the lux calibration value is taken from within the output file itself.
- voltv: a value for manual volts calibration of the light meter. If voltv=NULL, the volts calibration value is taken from within the output file itself.
- warn: if set to true, give a warning if input file is large, and require user confirmation.

Details

The main tasks performed by the package are listed below. The relevant topic contains documentation and examples for each.

- Extraction of file header material is accomplished by [header.info](#).
- Input and downsampling of data is accomplished by [read.bin](#).
- Selection of time intervals is accomplished via [get.intervals](#).
- Computation of epochal summaries is accomplished by [epoch](#) and other functions documented therein.
- Computation of STFT analyses is accomplished by [stft](#).

The package provides definitions and methods for the following S3 classes:

- GRtime: Provides numeric storage and streamlined plotting for times. [GRtime](#)
- AccData: Stores GENEa accelerometer data, allowing plotting, subsetting and other computation. [AccData](#)
- VirtAccData: A virtual AccData object, for just-in-time data access via [get.intervals](#).
- stft: Processed STFT outputs, for plotting via [plot.stft](#).

The read.bin package reads in binary files compatible with the GeneActiv line of Accelerometers, for further processing by the other functions in this package. Most of the default options are those required in the most common cases, though users are advised to consider setting start and end to smaller intervals and/or choosing some level of downsampling when working with data files of longer than 24 hours in length.

The function reads in the desired analysis time window specified by start and end. For convenience, a variety of time window formats are accepted:

Large integers are read as page numbers in the dataset. Page numbers larger than that which is available in the file itself are constrained to what is available. Note that the first page is page 1. Small values (between 0 and 1) are taken as proportions of the data. For example, 'start = 0.5' would specify that reading should begin at the midpoint of the data. Strings are interpreted as dates and times using `parse.time`. In particular, times specified as "HH:MM" or "HH:MM:SS" are taken as the earliest time interval containing these times in the file. Strings with an integer prepended, using a space separator, as interpreted as that time after the appropriate number of midnights have

passed - in other words, the appropriate time of day on the Nth full day. Days of the week and dates in "day/month", "day/month/year", "month-day", "year-month-day" are also handled. Note that the time is interpreted in the same time zone as the data recording itself.

Actual data reading proceeds by two methods, depending on whether `mmap` is true or false. With `mmap = FALSE`, data is read in line by line using `readLine` until blocksize is filled, and then processed. With `mmap = TRUE`, the `mmap` package is used to map the entire data file into an address file, byte locations are calculated (depending on the setting of `pagerefs`), blocksize chunks of data are loaded, and then processed as raw vectors.

There are advantages and disadvantages to both methods: the `mmap` method is usually much faster, especially when we are only loading the final parts of the data. `readLines` will have to process the entire file in such a case. On the other hand, `mmap` requires a large amount of memory address space, and so can fail in 32 bit systems. Finally, reading of compressed bin files can only be done with the `readLine` method. Generally, if `mmap` reading fails, the function will attempt to catch the failure, and reprocess the file with the `readLine` method, giving a warning. Once data is loaded, calibration is then either performed using values from the binary file, or using manually inputted values (using the `gain`, `offset`, `luxv` and `voltv` arguments).

Main tasks performed

NA

Classes implemented

NA

WARNING

Reading in an entire .bin file will take a long time if the file contains a lot of datasets. Reading in such files without downsampling can use up all available memory. See `memory.limit`. This function is specific to header structure in GENEActiv output files. By design, it should be compatible with all firmware and software versions to date (as of version of current release). If order or field names are changed in future .bin files, this function may have to be updated appropriately.

Author(s)

Zhou Fang <zhou@activinsights.co.uk>

Activinsights Ltd. <joss.langford@activinsights.co.uk>

Charles Sweetland <charles@sweetland-solutions.co.uk>

Examples

```
requireNamespace("GENEAread")
binfile = system.file("binfile/TESTfile.bin", package = "GENEAread")[1]
#Read in the entire file, calibrated
procfile <- read.bin(binfile)
# print(procfile)
# procfile$data.out[1:5,]
# Uncalibrated, mmap off
procfile2 <- read.bin(binfile, calibrate = FALSE)
```

```

# procfile2$data.out[1:5,]
#Read in again, reusing already computed mmap pagerefs
# procfile3 <- read.bin(binfile, pagerefs = procfile2$pagerefs )
#Downsample by a factor of 10
procfilelo<-read.bin(binfile, downsample = 10)
# print(procfilelo)
object.size(procfilelo) / object.size(procfile)
#Read in a 1 minute interval
procfileshort <- read.bin(binfile, start = "16:50", end = "16:51")
# print(procfileshort)
##NOT RUN: Read, and save as a R workspace
#read.bin(binfile, outfile = "tmp.Rdata")
#print(load("tmp.Rdata"))
#print(processedfile)

```

AccData

Accelerometer Data Object

Description

Accelerometer Data Output from read.bin function

Format

An AccData object

Source

Output of [read.bin](#)

See Also

[read.bin](#)

Examples

```

requireNamespace("GENEAread")
binfile = system.file("binfile/TESTfile.bin", package = "GENEAread")[1]
#Read in the entire file, calibrated
procfile<-read.bin(binfile)
print(procfile)

plot(procfile$temperature,
xlim = c(min(procfile$data.out[,1]),
          max(procfile$data.out[,1])),
ylim = c(10,40))

plot(procfile$data.out[,c(1,7)])

```

convert.time	<i>convert.time</i>
--------------	---------------------

Description

extends time display from package chron to use h:m:s for > 1 day times.

Usage

```
convert.time(x, format = NULL)
```

Arguments

- | | |
|--------|--|
| x | Object to process. For convert.time, must be numeric. |
| format | A character string indicating the form of output. See strptime for details. If NULL, will be automatically chosen. |

Details

convert.time converts numerics to GRtime objects. The format argument allows a format string to be attached specifying the default format to display in. as.GRtime is a wrapper to convert.time, that when supplied with character input, coerces the value first to numeric using parse.time.

epoch	<i>Compute epochal summary statistics.</i>
-------	--

Description

Computes epochal summary statistics for an "AccData" object, matrix, or vector, and collates into a matrix or vector.

Usage

```
apply.epoch(x, epoch.size = 10, incl.date = FALSE, FUN)
```

Arguments

- | | |
|------------|---|
| x | The object to compute statistics for. Can be an "AccData" object, a matrix, or a vector. |
| epoch.size | Numeric giving intervals to consider and aggregate. For "AccData" x taken as seconds. Otherwise, considered as rows, or as individual readings. |
| incl.date | logical. If TRUE, include a column of times or original indices with the results. |
| FUN | A function to be applied to each epoch. |

Details

These functions compute epochal summary statistics for "AccData" objects, matrices and vectors.

`apply.epoch` is the general function - according to the size of `epoch.size`, it splits up the `x` into collections of consecutive rows, each with the same size. These are then successively supplied to `FUN` as its first argument. If the result of `FUN` is a single value, then the results are concatenated into a vector output. Otherwise, an array is formed with each row corresponding to a single epochal group. For `AccData`, the sampling frequency of the dataset is used to interpret the epoch size in seconds. Otherwise, the raw record indices are used. If `incl.date` is set, the original timestamp vector of the data, or the original indices, are downsampled and included as the first column of the output.

The remaining functions are wrappers that compute various commonly useful statistics – in particular, applied to "AccData" objects and arrays, they by default compute the epochal SVM mean, standard deviation, median, median absolute deviation, and autocorrelation, and sample quantiles respectively. (Arrays are treated as each column representing the `x`, `y`, and `z` components respectively.) Applied to vector input, processing will occur without the SVM calculation. This behaviour may be overridden by the `sqrt` setting, which will force the function to use the squared (default for arrays and "AccData") or original unit (default for vectors) values in the statistical analysis.

Value

A vector or array giving the computed epochal summaries. With `incl.date = TRUE`, the result is given as a `data.frame` suitable for plotting.

Examples

```
## Not run:
dat <- read.bin(system.file("binfile/TESTfile.bin", package = "GENEAread")[1]
, calibrate = TRUE)

#look for the epochs that exceed a certain threshold 50% of the time
plot(apply.epoch(dat, epoch.size = 3 ,
                FUN = function(t) mean(abs(svm(t) -1)>0.2)> 0.5 ), type = "l")

plot(dat[,1], svm(dat), log = "y", pch = ".")
lines(mean.epoch(dat, incl.date = TRUE), lwd = 2)
lines(mean.epoch(dat, epoch.size = 30, incl.date = TRUE), col = 2, lwd = 2)
# This should give all the same results, but by a different way
lines(apply.epoch(dat, epoch.size = 30,
                FUN = function(A) mean(svm(A, FALSE))), incl.date = TRUE), col = 3)

epsize = 30
lines(apply.epoch(dat, epoch.size = epsize,
                FUN = function(t) median(t[,1])),
                apply.epoch(dat, epoch.size = epsize,
                FUN = function(A) mean(svm(A, FALSE))), col = 4)

#note this is different
lines(apply.epoch(dat, epoch.size = epsize,
                FUN = function(t) median(t[,1])),
                apply.epoch(dat, epoch.size = epsize,
                FUN = function(A) mean(svm(A, sqrt = TRUE)))^2,
                col = 5)
```

```
#plot some statistics
par(mfrow = c(5,1), mar = c(1,4.5,1,1))
plot(sd.epoch(dat), type="l")
plot(median.epoch(dat), type= "l")
plot(mad.epoch(dat), type= "l")
plot(acf.epoch(dat), type = "l")
plot(autocor.epoch(dat), type= "l")
tmp = quantile.epoch(dat, quantiles= c(0.1, 0.25, 0.5, 0.75, 0.9)); matplot(tmp, type = "l")

## End(Not run)
```

GENEActiv.calibrate *GENEActiv.calibrate*

Description

Function starts by identifying ten second windows of non-movement. Next, the average acceleration per axis per window is used to estimate calibration error (offset and scaling) per axis. The function provides recommended correction factors to address the calibration error and a summary of the callibration procedure.

Usage

```
GENEActiv.calibrate(
  binfile,
  use.temp = TRUE,
  spherecrit = 0.3,
  minloadcrit = 72,
  printsummary = TRUE,
  chunksize = c(),
  window sizes = c(5, 900, 3600)
)
```

Arguments

<code>binfile</code>	A filename of a file to process.
<code>use.temp</code>	use temperature sensor data for calibration
<code>spherecrit</code>	The minimum required acceleration value (in g) on both sides of 0 g for each axis. Used to judge whether the sphere is sufficiently populated
<code>minloadcrit</code>	The minimum number of hours the code needs to read for the autocalibration procedure to be effective (only sensitive to multitudes of 12 hrs, other values will be ceiled). After loading these hours only extra data is loaded if calibration error has not been reduced to under 0.01 g
<code>printsummary</code>	if TRUE will print a summary when done

chunksize	number between 0.2 and 1 to specify the size of chunks to be loaded as a fraction of a 12 hour period, e.g. 0.5 equals 6 hour chunks. The default is 1 (12 hrs). For machines with less than 4Gb of RAM memory a value below 1 is recommended.
window sizes	Three values to indicate the lengths of the windows as in c(window1,window2,window3): window1 is the short epoch length in seconds and by default 5 this is the time window over which acceleration and angle metrics are calculated, window2 is the long epoch length in seconds for which non-wear and signal clipping are defined, default 900. However, window3 is the window length of data used for non-wear detection and by default 3600 seconds. So, when window3 is larger than window2 we use overlapping windows, while if window2 equals window3 non-wear periods are assessed by non-overlapping windows.

Details

The outputs from the function are as follows

- scale scaling correction values, e.g. c(1,1,1)
- offset offset correction values, e.g. c(0,0,0)
- tempoffset correction values related to temperature, e.g. c(0,0,0)
- cal.error.start absolute difference between Euclidean norm during all non-movement windows and 1 g before autocalibration
- cal.error.end absolute difference between Euclidean norm during all non-movement windows and 1 g after autocalibration
- spheredata average, standard deviation, Euclidean norm and temperature (if available) for all ten second non-movement windows as used for the autocalibration procedure
- npoints number of 10 second no-movement windows used to populate the sphere
- nhoursused number of hours of measurement data scanned to find the ten second time windows with no movement
- meantempcal mean temperature corresponding to the data as used for autocalibration.

Author(s)

Vincent T van Hees <vincentvanhees@gmail.com> Zhou Fang Charles Sweetland <charles@Sweetland-solutions.co.uk>

References

van Hees VT, Fang Z, Langford J, Assah F, Mohammad A, da Silva IC, Trenell MI, White T, Wareham NJ, Brage S. Auto-calibration of accelerometer data for free-living physical activity assessment using local gravity and temperature: an evaluation on four continents. *J Appl Physiol* (1985). 2014 Aug 7

get.intervals	<i>Extract an interval of data.</i>
---------------	-------------------------------------

Description

Function for extracting sub intervals of data, and implementation of just-in-time loading.

Usage

```
get.intervals(x, start=0, end = 1, length = NULL,
time.format = c("auto", "seconds", "days", "proportion", "measurements", "time"),
incl.date = FALSE, simplify = TRUE ,read.from.file=FALSE, size=Inf, ...)
```

Arguments

x	Object to process. Can be array,
start	Start of interval.
end	End of interval.
length	Length of interval.
time.format	Method with which start and end should be understood.
incl.date	logical. Include a column denoting time?
simplify	logical. If TRUE, output an array. Otherwise output a AccData object.
read.from.file	logical. If TRUE, re-read the relevant time interval from the original bin file.
size	Desired number of samples in output.
...	Additional arguments to be passed to read.bin , if read.from.file is TRUE.

Details

The function extracts the desired analysis time window specified by start and end. If length is specified, then the end is set to a point length units after start. The times are interpreted in terms of time.format. For convenience, a variety of time window formats are accepted:

- "seconds": Seconds since start of dataset.
- "days": Days since start of dataset.
- "proportion": Proportional point within dataset, given as a numeric between 0 and 1.
- "measurements": Raw number of samples since start of dataset.
- "time": Time string, as understood via [parse.time](#).
- "auto": Default - attempt to determine time format from size and type of start.

Some capacity for using mixed types of inputs for start and length in particular is present.

The input object x is typically an "AccData" object, though arrays are also accepted. "VirtAccData" are dealt with by using the timestamp and call information recorded within them to do a new read

of the original bin file, assuming this is still available. This is useful for 'just in time' reads of data. "AccData" can be dealt with in this way by setting `read.from.file`.

Note that for `read.from.file`, only "time" and "proportion" `time.format` are presently supported.

With `simplify = FALSE`, an "AccData" S3 object with the desired records. Otherwise, an array containing either 3 or 4 columns, containing the x, y, z acceleration vectors and optionally a time vector.

See Also

[read.bin](#), [AccData](#), [get.intervals](#)

Examples

```
binfile = system.file("binfile/TESTfile.bin", package = "GENEAread")[1]

#Read in a highly downsampled version of the file
procfile<-read.bin(binfile, downsample = 100)
print(procfile)

#Overlay some segments in different colour
lines(get.intervals(procfile, start = 0.4, end = 0.5,
                    time.format = "prop", incl.date = TRUE)[,1:2],
      col=2)

lines(get.intervals(procfile, start = 0.4, end = 5,
                    time.format = "sec", incl.date = TRUE)[,1:2],
      col=3)

lines(get.intervals(procfile, start = "16:51", end = "16:52",
                    time.format = "time", incl.date = TRUE)[,1:2],
      col=4)

# Note that measurements will depend on the downsampling rate,
# not the original sampling rate of the data
lines(get.intervals(procfile, start = 100, length = 10,
                    time.format = "measurement", incl.date = TRUE)[,1:2],
      col=5)

#This is also understood
lines(get.intervals(procfile, start = "16:52:10", 30,
                    incl.date = TRUE)[,1:2],
      col=6)

#Now load in virtually
virtfile<-read.bin(binfile, virtual = TRUE)
#Notice that get.intervals with simplify = FALSE gives a genuine AccData object
realfile = get.intervals(virtfile, start = 0.5, end = 1, simplify = FALSE)
virtfile
realfile
#get.intervals calls read.bin automatically
points(get.intervals(virtfile, start = "16:52:10", "16:52:40",
                    incl.date = TRUE)[,1:2], col=4, pch = ".")
```

```
#Alternatively, re-read procfile at a different resampling rate.
lines(get.intervals(procfile, start = "16:49:00", "16:49:30",
                    incl.date = TRUE, read.from.file = TRUE, downsample = 300)[,1:2],
      col=2)
```

GRtime

Date time handling for the GENEAREad package.

Description

Stores date time data as a numeric, with facility for pretty printing and axis commands.

Usage

```
as.GRtime(x, format = NULL, ...)
```

Arguments

x	Object to process. For <code>convert.time</code> , must be numeric. For <code>as.GRtime</code> may be numeric or character. For <code>format.GRtime</code> , a GRtime object, or a numeric.
format	A character string indicating the form of output. See strptime for details. If NULL, will be automatically chosen.
...	Additional arguments to be passed to parse.time , as.numeric , format.POSIXct , axis , pretty.POSIXt .

Details

The GRtime class handles dates and times for the GENEAREad class. The class treats dates as numerics denoting seconds since the UNIX epoch, with potentially a string attached specifying the format to print in. Unlike POSIXct, we avoid some of the processing, especially with respect to time zones, and allow some more flexibility in time computation and display. A range of operators are defined.

`convert.time` converts numerics to GRtime objects. The `format` argument allows a format string to be attached specifying the default format to display in. `as.GRtime` is a wrapper to `convert.time`, that when supplied with character input, coerces the value first to numeric using `parse.time`.

`format.GRtime` formats GRtime objects for pretty printing. If `format` is provided as argument, that is used. Else, if the `format` attribute is set on `x`, that is used. Finally, if formats are not provided, and `x` is of length greater than one, the range of values of `x` is used to decide the units displayed. Numerics are also accepted - they are coerced to GRtime.

`axis.GRtime` is used to plot GRtime axis, choosing, by default, breakpoints that give 'pretty' sub intervals. Note that [plot.default](#) uses `axis.GRtime` by default if supplied with a GRtime object in one of the directions. However, [image.default](#) based functions do not use the class axis functions, so axes must be plotted manually.

`pretty.GRtime` computes 'pretty' breakpoints, using the algorithm of `pretty.POSIXt`. Attributes are preserved.

- For `convert.time`, `as.GRtime` and `pretty.GRtime`, a `GRtime` object.
- For `format.GRtime` a character string representation.
- For `axis.GRtime` a list containing positions and labels for axis markers.

See Also

[parse.time](#), [get.intervals](#), [AccData](#)

Examples

```
as.GRtime("00:01")
#format is automatically set
convert.time(1:10)
convert.time(1:10*1000)
#we add a different default format
convert.time(1:10*1000, "%H:%M:%OS3") -> t
t
str(t)
# we override format with our own
format(t, format = "%a %d/%m/%y %H:%M:%OS3")

# plot calls axis.GRtime automatically. Notice
# that the format attribute is used.
plot(t, 1:10)
#strip out the default format
t2 = convert.time(t, format = NULL)
plot(t2, 1:10)

# image plots are a bit more complex

Z = matrix(rnorm(100), 10)
image(x = t, y = t2, z = Z, axes = FALSE)
Axis(x = t, side = 1) #Axis also works
box() #complete the bounding box

# custom axes
plot(t2, 1:10, xaxt = "n")
```

hanning.window

Hanning Window

Description

A hanning window used by the STFT function

Usage

`hanning.window(n)`

Arguments

n number of points inside the window

header.info *Get header info from GENE output (.bin) file*

Description

Function to extract relevant header fields and values from a file.

Usage

```
header.info(binfile, more=TRUE)
```

Arguments

binfile The file from which to extract the header

more logical. If TRUE, extract additional data from file useful for calibration and data reading.

Details

The function extracts useful information from a .bin file, such as information about the GENE device used to produce the output, and characteristics of the subject who wore the device. The function also accepts data that has been compressed in 'gzip', 'bzip2' or 'xz' formats. See file. With more set to TRUE, additional data is extracted, mainly for internal use in read.bin.

This function is specific to header structure in GENEActiv output files. By design, it should be compatible with all firmware and software versions to date (as of version of current release). If order or field names are changed in future .bin files, this function may have to be updated appropriately. The function works by looking for appropriate section headings in the .bin files.

Value

A data.frame with extracted header information, each row a particular header field with its value. If more is TRUE, an attribute "calibration" is attached to the object, consisting of a list with measurement offsets, sampling frequency estimates, start times and time zones, data position offsets, and if mmap is detected, byte locations and increments for mmap reading.

Warning

NA

See Also

[read.bin](#)

Examples

```
fileheader <- header.info(system.file("binfile/TESTfile.bin",
                                     package = "GENEAread")[1],
                           more = TRUE)

print(fileheader)
attr(fileheader, "calibration")
```

 parse.time

Parses a character time representation to another format.

Description

Converts a character vector in a variety of forms into either the raw second, second classed as POSIXct, or days since Unix epoch.

Usage

```
parse.time(t="", format=c("seconds", "days", "POSIX"), tzzone = 0,
           start = NULL, startmidnight = NULL)
```

Arguments

t	A character string representation of a date-time expression.
format	A character string indicating which representation to output. Can be either seconds, days or POSIX.
tzzone	The time zone the time is given in, expressed as an offset from UTC in hours.
start	Earliest allowable time stamp in the data, as seconds since Unix epoch.
startmidnight	Midnight of day '0' in the data, as seconds since Unix epoch.

Details

The function processes character vectors of the form "DATE TIME" – that is to say, a maximum of two terms separated by a space per value.

"TIME" is given in 24 hour format, separated by colons, in "hh:mm", "hh:mm:ss", "hh:mm:ss:ms" or "hh:mm:ss.ms" format. If omitted, the time is taken to be 00:00:00.000.

"DATE" can be a date representation as "YYYY-MM-DD", "DD/MM/YY" or "DD/MM/YYYY" (noting the use of a colon or backslash separator to distinguish between the two). Alternatively, with start and/or startmidnight supplied, an integer "NN" or string "DOW" corresponding to a day of the week can be used instead. Then, the function will find the first timestamp matching the correct "TIME", that falls NN midnights after startmidnight and is after start, or, in the latter case, the first timestamp after the day of start that matches the appropriate day of the week. If a blank "DATE" is supplied, the function will either use the UNIX epoch, or find the first match, corresponding to the case NN = 0.

Once this is done the time is converted to the required format: POSIX is the usual R POSIXct format; days is the julian days since UNIX epoch 1970-1-1; seconds is the number of seconds (including subseconds) since 1970-1-1. Note that for formats other than POSIX, the output is in the same timezone as tzone. POSIX stores the time internally as the time in UTC, and applies a format that gives this time local to the user.

Value

A converted date-time string in the specified format. In the case of "seconds", or "days", a numeric. For POSIX, a [POSIXct](#) object.

See Also

[convert.time](#), [get.intervals](#)

Examples

```
t1 = parse.time("2012-06-21 13:04:01"); print(t1)
parse.time("21/06/12 13:04:01") #gives the same result

parse.time(c("19/07/70", "20/07/70"), format = "days")
#results here will depend on your locale
parse.time(c("19/07/70", "20/07/70"), format = "POSIX", tzone = -4)

#one is the same day, one can only find a match the next day
parse.time("13:05", start = t1) - t1
parse.time("13:00", start = t1) - t1
#asking to wait 1 midnight means both times are considered as
#times on the same, full day of data
parse.time(c("1 13:05", "1 13:00"), start = t1) - t1
#2012-06-21 is a Thursday, so this is equivalent
parse.time(c("Fri 13:05", "Fri 13:00"), start = t1) - t1
#Longer form days of the week are also understood. Note that
#the first day does not get matched.
parse.time(c("Thursday 13:05", "Thursday 13:00"), start = t1) - t1
```

plot.stft

plot stft

Description

Processes a dataset, creating an object contained processed time-frequency analyses. These can then be plotted.

Usage

```
## S3 method for class 'stft'
plot(
  x,
  mode = c("decibels", "modulus", "pval"),
  log = "",
  showmax = TRUE,
  median = FALSE,
  xaxis = TRUE,
  toptthresh,
  reassign = (!(is.null(x$LGD)) && !("mv" %in% x$type)),
  ylim,
  xlim,
  new = TRUE,
  zlim.raw,
  zlim.quantile,
  cex,
  col = gray(63:0/63),
  ...
)
```

Arguments

x	"stft" class object to be processed.
mode	What should be plotted? • "decibels": log10 of FFT modulus • "modulus": Raw FFT modulus • "pvalue": P-value of each frequency's modulus assuming that window was in fact white noise of equal standard deviation
log	For log = "y", use a log scale on the y axis.
showmax	Vector or logical. Compute and plot the principle frequency components?
median	logical. If TRUE, smooth the STFT plot in the time direction with a running median.
xaxis	logical. If TRUE, plot pretty time axes.
toptthresh	For finite values, crop plot for frequencies higher than this value, and show a summary plot up top.
reassign	logical. Plot reassigned stft, if available?
xlim, ylim	Parameters controlling axes limits of plot.
new	logical. If TRUE, make a new plot. Otherwise overlay on to existing plot.
zlim.raw	Raw values at which to threshold values for computation of heatmap colours.
zlim.quantile	Quantile values at which to threshold values for computation of heatmap colours.
cex	Size of points for reassigned STFT plotting.
col	Vector of colours to be used for plotting.
...	Additional arguments to be passed to methods.

Details

STFT objects are created by the `stft` function. These methods print some useful summary statistics about them, and produce plots. `mode` determines the type of plot. "decibel" and "modulus" work with the raw values, while "pvalue" conducts some degree of normalisation in each time window and so is perhaps more useful for data showing a large variation in sd across different points in time. If the `null.calc` was set in the original `stft` argument, that is used - otherwise, an Exponential distribution is fit to each window, and the `pvalues` computed from that.

By default, the function uses some empirical quantile based colour thresholds designed to give somewhat reasonable and informative plots. This can be overridden, however, by setting different `zlim.raw` or `zlim.quantile` results. This can be useful for comparing two different datasets.

Reassigned stft plots are constructed, by default, when they are available, and when the original was not a "mv" stft. Unlike the heatmap used in the usual stft plot, a 2d scatterplot is used instead. This means that if there are few data points, it can be advantageous to set a higher `cex` value for larger points and better display.

With Accelerometer data, often the frequencies of interest are concentrated at the lower frequencies. `Topthresh` crops the frequency display to show only those frequencies. A summary plot is shown on the top, to compensate. Choosing a grid of frequencies, this plot draws one line to represent the energies present in the signal at that particular frequency, and higher. Black lines are drawn for frequencies less than $2/3$ the `topthresh`, red lines for $2/3 - 1$ times `topthresh`, and blue lines for frequencies higher than `topthresh`. Alternative, set `log = "y"` to put frequencies on a log scale.

Value

These functions are run for their side effects.

See Also

`stft`, `image.default`

Examples

```
## Not run: # Real data
binfile = system.file("binfile/TESTfile.bin", package = "GENEAread")[1]

#Read in the entire file, calibrated
procfile<-read.bin(binfile)
#Create stft object
obj = stft(procfile, type = "svm", quiet = TRUE)
#Look at it
print(obj)

plot(obj, cex = 5)
plot(obj, showmax = FALSE, cex = 5) #suppress principals

#pval plot
plot(obj, mode = "pval", cex = 5)
#disable reassigned stft
plot(obj, mode = "pval", reassign = FALSE)
#median smoothing
```

```

plot(obj, mode = "pval", reassign = FALSE, median = TRUE)
#log scale frequency, no top bar
dev.new(); plot(obj, mode = "pval", reassign = FALSE, toptthresh = Inf, log = "y")

## End(Not run)

```

print.stft	<i>print.stft</i>
------------	-------------------

Description

print the stft object

Usage

```

## S3 method for class 'stft'
print(x, ...)

```

Arguments

x	stft object
...	ignored

recalibrate	<i>recalibrate</i>
-------------	--------------------

Description

Taking a GENEActiv binfile and using the recalibration script to create a new calibrated binfile

Usage

```

recalibrate(
  datadir,
  outputdir,
  use.temp = TRUE,
  spherecrit = 0.3,
  minloadcrit = 72,
  printsummary = TRUE,
  chunksize = c(0.5),
  windowsizes = c(60, 900, 3600)
)

```

Arguments

<code>datadir</code>	The location of the directory/file containing GENEActiv binfile.
<code>outputdir</code>	The location of the directory/file for the calibrated files to be saved.
<code>use.temp</code>	Use temperature sensor data if available (Geneactive only)
<code>spherecrit</code>	the minimum required acceleration value (in g) on both sides of 0 g for each axis. Used to judge whether the sphere is sufficiently populated
<code>minloadcrit</code>	the minimum number of hours the code needs to read for the autocalibration procedure to be effective (only sensitive to multitudes of 12 hrs, other values will be ceiled). After loading these hours only extra data is loaded if calibration error has not been reduced to under 0.01 g.
<code>printsummary</code>	if TRUE will print a summary when done chunksize number between 0.2 and 1 to specify the size of chunks to be loaded as a fraction of a 12 hour period, e.g. 0.5 equals 6 hour chunks. The default is 1 (12 hrs). For machines with less than 4Gb of RAM memory a value below 1 is recommended.
<code>chunksize</code>	number between 0.2 and 1 to specify the size of chunks to be loaded as a fraction of a 12 hour period, e.g. 0.5 equals 6 hour chunks. The default is 1 (12 hrs). For machines with less than 4Gb of RAM memory a value below 1 is recommended.
<code>window sizes</code>	Three values to indicate the lengths of the windows as in <code>c(window1,window2,window3)</code> : window1 is the short epoch length in seconds and by default 5 this is the time window over which acceleration and angle metrics are calculated, window2 is the long epoch length in seconds for which non-wear and signal clipping are defined, default 900. However, window3 is the window length of data used for non-wear detection and by default 3600 seconds. So, when window3 is larger than window2 we use overlapping windows, while if window2 equals window3 non-wear periods are assessed by non-overlapping windows.

Details

Takes each binfile found in the data directory, calibrates according to the routine by Vincent T. van Hees and saves the calibrated file to the specified output directory

Value

Saves a calibrated binfile to an output folder

Examples

```
## Not run:
DataDirectory = "C:/Users/DataDirectory"
ReCalibrate(DataDirectory)

## End(Not run)
```

stft

Computes Short Time Fourier Transforms

Description

Processes a dataset, creating an object contained processed time-frequency analyses. These can then be plotted.

Usage

```
stft(X, start=0, end=1, length=NULL, time.format = c("auto"),
     type = c("mv", "svm", "sum"), mv.indices,
     date.col, reassign = TRUE, plot.it = FALSE,...)
```

Arguments

X	The dataset to be processed.
start, end, length, time.format	A specification for the segment to process, as in <code>get.intervals</code> .
type	The type of STFT to compute.
mv.indices	For type = "mv" or type = "sum", the indices to process and the order to process them in.
date.col	logical. Whether the first column should be ignored and treated as a timestamp. If unset, is automatically chosen.
reassign	logical. If TRUE, compute the time-reassigned STFT. For type <code>c("mv", "sum")</code> , this is done with the first coordinate in <code>mv.indices</code> .
plot.it	logical. Whether to plot the STFT immediately when processing is complete, using the default <code>plot.stft</code> options.
...	Additional optional arguments to control the STFT computation. These are: • win: Window size in seconds for STFT computation. Increased window size mean better frequency resolution, but poorer time resolution. Defaults to 10 seconds. • inc: Increment between successive time steps for processing. Defaults to <code>win/2</code>. • coef: Number of fourier frequencies to compute. Small values will remove the higher frequencies from the processed object. Defaults to the maximum, <code>win/2</code>. • wtype: String giving the name of a window function, providing coefficients for filtering before processing. "hanning.window" is the default, with "uniform.window" also available. • freq: Sampling frequency of data set. If not given, is taken from X itself, or assumed to be 1 if unavailable.

- **centre:** If TRUE (Default), centre the data in each window before processing is done. Useful for avoiding excessively large DC offset coefficients in results.
- **calc.null:** If TRUE (Defaults to FALSE), compute a 'null' STFT by resampling the data completely, then doing a STFT.
- **pvalues:** If TRUE (Defaults to FALSE) Compute bootstrapped pvalues for each position by resampling within each window and applying a wilcox test.
- **time:** Allows the user to set an overriding timestamp vector to be used for processing.
- **quiet:** If TRUE, suppress output.

Details

This function accepts input in a variety of forms and computes short time fourier transforms to extract frequency structure from the data. *X* may be an array, a vector, or an *AccData* object. If *date.col* is TRUE, the first column of an array *X* would be used to determine timestamps. Otherwise indices would be used. If *date.col* is not set, the function will attempt to determine whether the first column is timestamp-like. The timestamp column is removed from *X* (and so not included in consideration of *mv.indices*, for instance). With vectors, the basic method is to compute the STFT by creating windows of size *win* seconds every *inc* seconds, and computing the fourier transform. With multi-dimensional data and *AccData*, processing is done on the dimensions that are in *mv.indices*, or the first three non-date columns if that is unavailable. Three methods are possible:

- 1. *type* = "mv": The one dimensional method is first applied to each of the chosen column indices. These are then collated by choosing, for each time-frequency combination, the maximum such value across each of the indices.
- 2. *type* = "svm": The SVM is computed first for each time step by computing the square rooted sum of squares. This is then dealt with using the one dimensional method.
- 3. *type* = "sum": As in "mv", the 1d method is applied. The square of the modulus of the result is then summed and square rooted.

If *reassign* is set, the time reassigned *stft* is also computed for the first element of *mv.indices* or the *svm* as appropriate, by using finite differencing. This gives potentially better resolution results for data with a clear signal component.

Value

A "stft" class object - a list with the following components:

- **call:** The function call.
- **type:** Type of STFT computed.
- **values:** Mod of FFT computed, with each row corresponding to a specific time increment.
- **increment, windowsize, centre, sampling.frequency:** Various control parameters used in the computation.
- **null.logmean, null.logsd:** Log of the square rooted mean and standard deviation of the Mod FFT squared for the randomised data, if *calc.null* = TRUE.

- p.values: Wilcoxon pvalues, if pvalues = TRUE.
- principals: Principal frequencies.
- frequency: Frequencies at which FFT is computed.
- time: Timestamps for FFT windows.
- LGD: Local group delay matrix for reassigned STFT.
- CIF: Channelized instantaneous frequency matrix for reassigned STFT.

References

Fulop, S.A. & Fitz, K. (2006). Algorithms for computing the time-corrected instantaneous frequency (reassigned) spectrogram, with applications J Acoustical Society of America 119(1), 360–371. Nelson, D.J. (2001). Cross-spectral methods for processing speech J Acoustical Society of America 110(1), 2575-2592.

Examples

```
## Not run:
#Some artificial data
time = 1:5000
#sum of two sine curves at 0.3 Hz and 0.05 Hz
f1 = 0.3; f2 = 0.05
sin1 = sin(time * f1 * 2*pi)
sin2 = sin(time * f2 * 2*pi)
#add a bit of noise
signal = sin1 + sin2 + 1*rnorm(5000)
#non-reassigned
stft(signal, plot = TRUE, reassign = FALSE, win = 100)
#reassigned
stft(signal, plot = TRUE, reassign = TRUE, win = 100)

#add a third component: varying frequency.
stft(signal + sin(cumsum(seq(f2, f1, length = 5000))*2*pi),
      plot = TRUE, reassign = TRUE, win = 100)

# Real data
binfile = system.file("binfile/TESTfile.bin", package = "GENEAread")[1]

# Read in the entire file, calibrated
procfile<-read.bin(binfile)
# Default is mv
stft(procfile, plot.it = TRUE)
# Try sum?
stft(procfile, plot.it = TRUE, type = "sum", reassign = FALSE)

# Just look at the last 50% of the data
stft(procfile, start = 0.5, plot.it = TRUE)

# not reassigned, svm
stft(procfile, type = "svm", reassign = FALSE, plot.it = TRUE)
# a narrower 5 second window means better time resolution
```

```

stft(procfile, type = "svm", reassign = FALSE, plot.it = TRUE, win = 5)
# choose increments so as not to overlap
stft(procfile, type = "svm", reassign = FALSE, plot.it = TRUE, win = 5, inc = 5)
# uniform windows
stft(procfile, type = "svm", reassign = FALSE, plot.it = TRUE, wtype = "uniform.window")
# Svm, reassigned, quietly
obj = stft(procfile, type = "svm", quiet = TRUE)
plot(obj, cex = 3, showmax = FALSE, mode = "pval")

#example code
plot(stft(subs(mag, 0.94,0.96), win = 1024, plot = F, coef = 512), zlog = T, log="y")
plot(stft(subs(mag, 0.7,8), win = 1024, plot = F, coef = 512), zlog = T, log="y")
plot(stft(subs(mag, 0.0001,0.005), win = 1024, plot = F, coef = 512), zlog = T)
plot(stft(subs(mag, 0.7,0.8), win = 1024, plot = F), zlog = T, log = "y")

plot(stft(rep(1, 1000) +
  c(sin(1:500/ 10 * 2*pi), rep(0, 500)) +
  c(rep(0, 300),sin(1:500/ 20 * 2*pi), rep(0, 200)),
  freq = 1, plot.it = F), log="x")

stft(sin(1:1000 / (1 +sqrt(1000:1)) * 2 * pi), freq = 1)
stft(rep(1, 1000) + sin(1:1000/ 10 * 2*pi), freq = 1)

## End(Not run)

```

svm

sum vector magnitude

Description

svm acts identically to 'mean.epoch', with the epoch set to the sampling period. In other words, it computes the instantaneous sum of vector magnitudes of the acceleration at each record point. The function takes "AccData", array and vector input. Note that if provided with an array with 4 or more columns, columns 2 to 4 are used – the first column is regard as a timestamp and hence ignored.

Usage

```
svm(obj, sqrt)
```

Arguments

obj	AccData object
sqrt	Function to use to calculate SVM

Examples

```

dat <- read.bin(system.file("binfile/TESTfile.bin", package = "GENEaread")[1], calibrate = TRUE)
svm(dat)

```

uniform.window	<i>Uniform Window</i>
----------------	-----------------------

Description

A uniform window used by the STFT function

Usage

```
uniform.window(n)
```

Arguments

n	number of points inside the window
---	------------------------------------

Index

* Internal

hanning.window, [14](#)
uniform.window, [26](#)

* datasets

AccData, [6](#)

AccData, [4](#), [6](#), [12](#), [14](#)

acf.epoch (epoch), [7](#)

apply.epoch (epoch), [7](#)

as.GRtime (GRtime), [13](#)

as.numeric, [13](#)

autocor.epoch (epoch), [7](#)

axis, [13](#)

Axis.GRtime (GRtime), [13](#)

axis.GRtime (GRtime), [13](#)

c.GRtime (GRtime), [13](#)

convert.time, [7](#), [17](#)

epoch, [4](#), [7](#)

format.GRtime (GRtime), [13](#)

format.POSIXct, [13](#)

GENEActiv.calibrate, [9](#)

GENEAread (GENEAread-package), [2](#)

GENEAread-package, [2](#)

get.intervals, [4](#), [11](#), [12](#), [14](#), [17](#)

GRtime, [4](#), [13](#)

hanning.window, [14](#)

header.info, [4](#), [15](#)

image.default, [13](#), [19](#)

mad.epoch (epoch), [7](#)

mean.epoch (epoch), [7](#)

median.epoch (epoch), [7](#)

mmap, [3](#), [5](#)

Ops.GRtime (GRtime), [13](#)

parse.time, [11](#), [13](#), [14](#), [16](#)

plot.default, [13](#)

plot.stft, [4](#), [17](#)

POSIXct, [17](#)

pretty.GRtime (GRtime), [13](#)

pretty.POSIXt, [13](#)

print.GRtime (GRtime), [13](#)

print.stft, [20](#)

print.VirtAccData (get.intervals), [11](#)

quantile.epoch (epoch), [7](#)

read.bin, [4](#), [6](#), [11](#), [12](#), [15](#)

read.bin (GENEAread-package), [2](#)

readLines, [5](#)

recalibrate, [20](#)

sd.epoch (epoch), [7](#)

stft, [4](#), [19](#), [22](#)

strptime, [7](#), [13](#)

svm, [25](#)

uniform.window, [26](#)

VirtAccData (get.intervals), [11](#)