

Package ‘photobiologyInOut’

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Type Package

Title Read Spectral and Logged Data from Foreign Files

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Description Functions for reading, and in some cases writing, foreign files containing spectral data from spectrometers and their associated software, output from daylight simulation models in common use, and some spectral data repositories. As well as functions for exchange of spectral data with other R packages. Part of the 'r4photobiology' suite, Aphalo P. J. (2015) <doi:10.19232/uv4pb.2015.1.14>.

License GPL (>= 2)

VignetteBuilder knitr

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LazyLoad yes

LazyData yes

ByteCompile true

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BugReports <https://bitbucket.org/aphalo/photobiologyinout/issues/>

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photobiologyInOut-package

photobiologyInOut: Read Spectral and Logged Data from Foreign Files

Description

Functions for reading, and in some cases writing, foreign files containing spectral data from spectrometers and their associated software, output from daylight simulation models in common use, and some spectral data repositories. As well as functions for exchange of spectral data with other R packages. Part of the 'r4photobiology' suite, Aphalo P. J. (2015) <doi:10.19232/uv4pb.2015.1.14>.

Warning!

Most of the file formats supported are not standardized, and are a moving target because of changes in instrument firmware and support software. In addition the output format, especially with models, can depend on settings that users can alter. So do check that import is working as expected, and if not, please please raise an issue and upload one example of an incorrectly decoded file.

Note

From version 0.4.4 the time zone (tz) used for decoding dates and times in files imported defaults to "UTC". In most cases you will need to pass the tz (or the locale) where the file was created as an argument to the functions!

Author(s)

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References

Aphalo, Pedro J. (2015) The r4photobiology suite. UV4Plants Bulletin, 2015:1, 21-29. <https://doi.org/10.19232/uv4pb.2015.1.14>.

See Also

Useful links:

- <https://docs.r4photobiology.info/photobiologyInOut/>
- Report bugs at <https://bitbucket.org/aphalo/photobiologyinout/issues/>

as.colorSpec

Convert into 'colorSpec::colorSpec' objects

Description

Convert spectral objects (xxxx_spct, xxxx_mspect) as defined in package 'photobiology' into colorSpec objects preserving as much information as possible.

Usage

```
## S3 method for class 'generic_mspct'
as.colorSpec(x, spct.data.var = NULL, multiplier = 1, ...)

## S3 method for class 'generic_spct'
as.colorSpec(x, spct.data.var = NULL, multiplier = 1, ...)

## S3 method for class 'chroma_spct'
as.colorSpec(x, spct.data.var = NULL, multiplier = 1, ...)
```

Arguments

x	R object
spct.data.var	character The name of the variable to read spectral data from.
multiplier	numeric A multiplier to be applied to the 'spc' data to do unit or scale conversion.
...	currently ignored.

Methods (by class)

- generic_spct:
- chroma_spct:

Warning!

Always check the sanity of the returned data values, as guessing is needed when matching the different classes, and the functions defined here are NOT guaranteed to return valid data without help from the user through optional function arguments.

Note

Objects of class `colorSpec::colorSpec` do not contain metadata or class data from which the units of expression could be obtained. When using this function the user needs to use parameter `multiplier` to convert the data to what is expected by the object constructors defined in package 'photobiology' but should only rarely need to use parameter `spct.data.var` to select the quantity. `colorSpec::colorSpec` objects may use memory more efficiently than spectral objects of the classes for collections of spectra defined in package 'photobiology' as wavelengths are assumed to be the same for all member spectra, and stored only once while this assumption is not made for collections of spectra, allowing different wavelengths and lengths for the component spectra. Wavelengths are stored for each spectrum, but as spectral classes are derived from 'tbl_df' in many cases no redundant copies of wavelength data will be made in memory in spite of the more flexible semantics of the objects.

Examples

```
if (requireNamespace("colorSpec", quietly = TRUE)) {
}
```

as.generic_mspct	<i>Convert into generic_mspct</i>
------------------	-----------------------------------

Description

Convert into generic_mspct

Usage

```
## S3 method for class 'colorSpec'  
as.generic_mspct(x, multiplier = 1, ...)
```

Arguments

x	R object
multiplier	numeric A multiplier to be applied to the spectral quantity data to do unit or scale conversion.
...	currently ignored.

as.generic_spct	<i>Coerce into generic_spct</i>
-----------------	---------------------------------

Description

Coerce into generic_spct

Usage

```
## S3 method for class 'colorSpec'  
as.generic_spct(x, multiplier = 1, ...)
```

Arguments

x	R object
multiplier	numeric A multiplier to be applied to the spectral quantity data to do unit or scale conversion.
...	currently ignored.

colorSpec2mspct *Convert 'colorSpec::colorSpec' objects*

Description

Convert 'colorSpec::colorSpec' objects into spectral objects (xxxx_spct, xxxx_mspct) as defined in package 'photobiology' and vice versa preserving as much information as possible.

Usage

```
colorSpec2mspct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.source_spct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.source_mspct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.response_spct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.response_mspct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.filter_spct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.filter_mspct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.reflector_spct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.reflector_mspct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.chroma_mspct(x, multiplier = 1, ...)  
  
colorSpec2spct(x, multiplier = 1, ...)  
  
colorSpec2chroma_spct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'  
as.chroma_spct(x, multiplier = 1, ...)  
  
## S3 method for class 'colorSpec'
```

```
as.chroma_mspct(x, multiplier = 1, ...)  
mspct2colorSpec(x, spct.data.var = NULL, multiplier = 1, ...)  
spct2colorSpec(x, spct.data.var = NULL, multiplier = 1, ...)  
chroma_spct2colorSpec(x, spct.data.var = NULL, multiplier = 1, ...)
```

Arguments

x	colorSpec object
multiplier	numeric A multiplier to be applied to the 'spc' data to do unit or scale conversion.
...	currently ignored.
spct.data.var	character The name of the variable to read spectral data from.

Warning!

Always check the sanity of the imported or exported data values, as guessing is needed when matching the different classes, and the functions defined here are NOT guaranteed to return valid data without help from the user through optional function arguments.

Note

Objects of class `colorSpec::colorSpec` do not contain metadata or class data from which the units of expression could be obtained. When using this function the user needs to use parameter `multiplier` to convert the data to what is expected by the object constructors defined in package 'photobiology' but should only rarely need to use parameter `spct.data.var` to select the quantity. `colorSpec::colorSpec` objects may use memory more efficiently than spectral objects of the classes for collections of spectra defined in package 'photobiology' as wavelengths are assumed to be the same for all member spectra, and stored only once while this assumption is not made for collections of spectra, allowing different wavelengths and lengths for the component spectra. Wavelengths are stored for each spectrum, but as spectral classes are derived from 'tbl_df' in many cases no redundant copies of wavelength data will be made in memory in spite of the more flexible semantics of the objects.

Examples

```
# example run only if 'colorSpec' is available  
if (requireNamespace("colorSpec", quietly = TRUE)) {  
  library(colorSpec)  
  colorSpec2mspct(Fs.5nm)  
  colorSpec2spct(Fs.5nm)  
  colorSpec2mspct(C.5nm)  
  colorSpec2spct(C.5nm)  
}
```

hyperSpec2mspct *Convert 'hyperSpec::hyperSpec' objects*

Description

Convert `hyperSpec::hyperSpec` objects containing VIS and UV radiation data into spectral objects (`xxxx_spct`, `xxxx_mspct`) as defined in package 'photobiology' and vice versa, preserving as much information as possible. As `hyperSpec` can contain other kinds of spectral data, it does make sense to use these functions only with objects containing data that can be handled by both packages.

Usage

```
hyperSpec2mspct(x, member.class, spct.data.var, multiplier = 1, ...)
```

```
hyperSpec2spct(x, multiplier = 1, ...)
```

```
mspct2hyperSpec(x, spct.data.var, multiplier = 1, ...)
```

```
spct2hyperSpec(x, spct.data.var = NULL, multiplier = 1, ...)
```

Arguments

<code>x</code>	hyperSpec object
<code>member.class</code>	character One of the spectrum classes defined in package 'photobiology'.
<code>spct.data.var</code>	character The name to be used for the 'spc' data when constructing the spectral objects.
<code>multiplier</code>	numeric A multiplier to be applied to the 'spc' data to do unit or scale conversion. For example "a.u." units in some examples in package 'hyperSpec' seem to have scale factors applied.
<code>...</code>	currently ignored.

Warning!

Always check the sanity of the imported or exported data values, as guessing is needed when matching the different classes, and the functions defined here are NOT guaranteed to return valid data without help from the user through optional function arguments.

Note

Objects of class `hyperSpec::hyperSpec` contain metadata or class data from which the quantity measured and the units of expression can be obtained. However, units as included in the objects are not well documented making automatic conversion difficult. When using this function the user may need to use parameter `multiplier` to scale the data to what is expected by the object constructors defined in package 'photobiology' and use parameter `spct.data.var` to select the quantity.

`hyperSpec::hyperSpec` objects may use memory more efficiently than spectral objects of the classes for collections of spectra defined in package 'photobiology' as wavelengths are assumed

to be the same for all member spectra, and stored only once while this assumption is not made for collections of spectra, allowing different wavelengths and lengths for the component spectra. Wavelengths are stored for each spectrum, but as spectral classes are derived from 'tbl_df' in many cases no redundant copies of wavelength data will be made in memory in spite of the more flexible semantics of the objects.

Examples

```
# example run only if 'hyperSpec' is available
if (requireNamespace("hyperSpec", quietly = TRUE)) {
  library(hyperSpec)
  data(laser)
  wl(laser) <-
    list(wl = 1e7 / (1/405e-7 - wl (laser)),
         label = expression (lambda / nm))
  laser.mspct <- hyperSpec2mspct(laser, "source_spct", "s.e.irrad")
  class(laser.mspct)
}
```

read_ASTER_txt

Read File downloaded from ASTER data base.

Description

Reads and parses the header of a test file as available through the ASTER reflectance database. The Name field is retrieved and copied to attribute "what.measured". The header of the file is preserved as a comment.

Usage

```
read_ASTER_txt(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  npixels = 2048
)
```

Arguments

file	character string
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".

label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Ignored.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
npixels	integer Number of pixels in spectral data.

Value

A raw_spect object.

Note

The header in these files has very little information, so the user needs to supply the number of pixels in the array as well as the date-time. The file contains a date in milliseconds but as the Raspberry Pi board contains no real-time clock, it seems to default to number of milliseconds since the Pi was switched on.

References

<https://www.r4photobiology.info> <https://speclib.jpl.nasa.gov/> Baldrige, A.; Hook, S.; Grove, C. & Rivera, G. (2009) The ASTER spectral library version 2.0. Remote Sensing of Environment. 113, 711-715

read_avaspec_csv *Read '.csv' File Saved by Avantes' Software for AvaSpec.*

Description

Reads and parses the header of a processed data file as output by the program Avaspec and then imports wavelength and spectral irradiance values. The file header has little useful metadata information.

Usage

```
read_avaspec_csv(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)

read_avaspec_xls(
```

```

    path,
    date = NULL,
    geocode = NULL,
    label = NULL,
    tz = NULL,
    locale = readr::default_locale()
  )

```

Arguments

file	character string
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone used for interpreting times saved in the file header.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
path	Path to the xls/xlsx file

Value

A source_spct object.

References

<https://www.r4photobiology.info> <https://www.avantes.com/>

read_csi_dat	<i>Read '.DAT' file(s) saved by modern Campbell Scientific loggers.</i>
--------------	---

Description

Reads and parses the header of a processed data file as output by the PC400 or PC200W programmes extracting variable names, units and quantities from the header. Uses the comment attribute to store the metadata.

Usage

```
read_csi_dat(
  file,
  geocode = NULL,
  label = NULL,
  data_skip = 0,
  n_max = Inf,
  locale = readr::default_locale()
)
```

Arguments

file	Path to file as a character string.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
data_skip	integer Number of records (rows) to skip from the actual data block.
n_max	integer Maximum number of records to read.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.

Value

read_csi_dat() returns a `tibble::tibble` object.

Note

This function is not useful for .DAT and .PRN files from old CSI loggers and software. Those were simple files, lacking metadata, which was stored in separate .FLD files.

References

<https://www.r4photobiology.info> <https://www.campbellsci.eu/>

read_fmi2mspct

Read multiple solar spectra from a data file.

Description

Read spectral irradiance file as output by Anders Lindors' model based on libRadTrans for hourly simulation, or measured data from FMI's Brewer spectrometer.

Usage

```
read_fmi2mspct(
  file,
  scale.factor = 0.001,
  geocode = NULL,
  what.measured = NULL,
  how.measured = NULL,
  date.field = 2L,
  time.field = 3L,
  date.format = "ymd",
  time.format = "hms",
  tz = NULL,
  time.shift.min = 0,
  locale = readr::default_locale(),
  .skip = 0,
  .n_max = -1
)
```

Arguments

<code>file</code>	Either a path to a file, a connection, or literal data (either a single string or a raw vector).
<code>scale.factor</code>	numeric A multiplier to be applied to the spectral irradiance values.
<code>geocode</code>	A data frame with columns <code>lon</code> and <code>lat</code> used to set attribute <code>"where.measured"</code> .
<code>what.measured</code>	character string, but if <code>NULL</code> the value of <code>file</code> is used, and if <code>NA</code> the <code>"what.measured"</code> attribute is not set.
<code>how.measured</code>	character string, but if <code>NULL</code> or <code>NA</code> the <code>"how.measured"</code> attribute is not set.
<code>date.field, time.field</code>	integer. Word positions in the header line.
<code>date.format</code>	character string. One of <code>"ymd"</code> , <code>"ydm"</code> , <code>"dmy"</code> , or <code>"mdy"</code> .
<code>time.format</code>	character string. One of <code>"hms"</code> , <code>"hm"</code> .
<code>tz</code>	character Time zone used for interpreting times saved in the file header.
<code>time.shift.min,</code>	numeric. Time shift with respect to TZ in minutes.
<code>locale</code>	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
<code>.skip</code>	Number of lines to skip before reading data.
<code>.n_max</code>	Maximum number of records to read.

Value

`read_fmi2mspct()` returns a `source_mspct` object containing `source_spct` objects as members, `time.unit` attribute set to `"second"` and `when.measured` attribute set to the date-time values extracted from the file body.

Note

See [read_table](#) for details of acceptable values for `file`. Individual spectra are names based on time and date in ISO format, at the time zone given by `tz` but the time shift subtracted. Say for times expressed in headers at UTC + 120 min, we use `tz = UTC` and `time.shift.min = 120` to convert times to UTC. This is different from using `tz = EET`, which is not invariant through the course of the year because of daylight saving time. Local time zones is not necessarily consistent across years because of changes in legislation. In contrast UTC is more consistent, making it preferable for time series.

read_fmi_cum	<i>Read daily cummulated solar spectrum data file(s).</i>
--------------	---

Description

Read one or more cumulated daily spectral irradiance file as output by Anders Lindors' model based on libRadTrans. Dates are read from the file header and parsed with the function supplied as `date.f`.

Usage

```
read_fmi_cum(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = "UTC",
  locale = readr::default_locale(),
  .skip = 3,
  .n_max = -1,
  .date.f = lubridate::ymd
)
```

```
read_m_fmi_cum(
  files,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = "UTC",
  .skip = 3,
  .n_max = -1,
  .date.f = lubridate::ymd
)
```

Arguments

<code>file</code>	Either a path to a file, a connection, or literal data (either a single string or a raw vector).
-------------------	--

date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone used for interpreting times saved in the file header.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
.skip	Number of lines to skip before reading data—i.e. the number of rows in the header.
.n_max	Maximum number of records to read.
.date.f	A function for extracting a date-time from the file header passed as character string to its first argument and which returns a POSIXct object.
files	list or vector of paths each one with the same requirements as described for argument file.

Value

read_fmi_cum() returns a source_spct object with time.unit attribute set to "day" and when.measured attribute set to the date-time extracted from the header at the top of the read file.

read_m_fmi_cum returns a source_mspect containing one source_spct object for each one of the multiple files read.

Note

See [read_table](#) for details of acceptable values for file.

Examples

```
file.name <- system.file("extdata", "2014-08-21_cum.hel",
                        package = "photobiologyInOut", mustWork = TRUE)
fmi.spct <- read_fmi_cum(file = file.name)
```

read_FReD_csv

Read '.CSV' FReD database.

Description

Reads a CSV data file downloaded from the FReD (Floral Reflectance Database) and then imports wavelengths and spectral reflectance values and flower ID.

Usage

```
read_FReD_csv(
  file,
  date = NA,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

Arguments

file	character string
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone used for interpreting times saved in the file header.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.

Value

A reflectance_spct object.

References

<http://www.reflectance.co.uk> Arnold SEJ, Faruq S, Savolainen V, McOwan PW, Chittka L, 2010 FReD: The Floral Reflectance Database - A Web Portal for Analyses of Flower Colour. PLoS ONE 5(12): e14287. doi:10.1371/journal.pone.0014287

read_licor_prn

Read '.PRN' File(s) Saved by LI-COR's PC1800 Program.

Description

Reads and parses the header of a processed data file as output by the PC1800 program to extract the whole header remark field and also decode whether data is in photon or energy based units. The time field is ignored as it does not contain year information. This instrument is no longer being manufactured.

Usage

```

read_licor_prn(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  s.qty = NULL
)

read_m_licor_prn(
  files,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = Sys.timezone(),
  locale = readr::default_locale(),
  s.qty = NULL
)

```

Arguments

file	Path to file as a character string.
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone used for interpreting times saved in the file header.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
s.qty	character The name of the spectral quantity to be read. One of "s.irrad", "Tfr", or "Rfr".
files	A list or vector of character strings.

Details

Function read_m_licor_prn() calls read_licor_file() for each file in files. See [read_table](#) for a description of valid arguments for files.

Value

read_licor_prn() returns a source_spct object with time.unit attribute set to "second" and when.measured attribute set to the date-time extracted from the file name, or supplied.

Function `read_m_licor_prn()` returns a `source_mspct` object containing one spectrum per file read.

Note

The LI-1800 spectroradiometer does not store the year as part of the data, only month, day, and time of day. Because of this, in the current version, if `NULL` is the argument to `date`, year is set to 0000.

References

<https://www.r4photobiology.info> <https://www.licor.com>

read_macam_dta	<i>Read '.DTA' File Saved by Macam's Software.</i>
----------------	--

Description

Reads and parses the header of a processed data file as output by the PC program to extract the time and date fields and a user label if present, and then imports wavelengths and spectral energy irradiance values.

Usage

```
read_macam_dta(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

Arguments

<code>file</code>	character string
<code>date</code>	a <code>POSIXct</code> object to use to set the "when.measured" attribute. If <code>NULL</code> , the default, the date is extracted from the file header.
<code>geocode</code>	A data frame with columns <code>lon</code> and <code>lat</code> used to set attribute "where.measured".
<code>label</code>	character string, but if <code>NULL</code> the value of <code>file</code> is used, and if <code>NA</code> the "what.measured" attribute is not set.
<code>tz</code>	character Time zone used for interpreting times saved in the file header.
<code>locale</code>	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use <code>locale</code> to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.

Value

A source_spct object.

References

<https://www.r4photobiology.info> <http://www.irradian.co.uk/>

read_oo_jazirrad *Read Files Saved by Ocean Optics' Jaz spectrometer.*

Description

Reads and parses the header of processed data text files output by Jaz instruments extracting the spectral data from the body of the file and the metadata, including time and date of measurement from the header. Jaz modular spectrometers are manufactured by Ocean Optics, Dunedin, Florida, USA.

Usage

```
read_oo_jazirrad(  
  file,  
  date = NULL,  
  geocode = NULL,  
  label = NULL,  
  tz = NULL,  
  locale = readr::default_locale()  
)  
  
read_oo_jazpc(  
  file,  
  qty.in = "Tpc",  
  Tfr.type = c("total", "internal"),  
  Rfr.type = c("total", "specular"),  
  date = NULL,  
  geocode = NULL,  
  label = NULL,  
  tz = NULL,  
  locale = readr::default_locale()  
)  
  
read_oo_jazdata(  
  file,  
  date = NULL,  
  geocode = NULL,  
  label = NULL,  
  tz = NULL,  
  locale = readr::default_locale()  
)
```

Arguments

file	character string.
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone used for interpreting times saved in the file header.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
qty.in	character string, one of "Tpc" (spectral transmittance, %), "A" (spectral absorbance), or "Rpc" (spectral reflectance, %).
Tfr.type	character string, either "total" or "internal".
Rfr.type	character string, either "total" or "specular".

Details

Function read_oo_jazirrad can read processed irradiance output files. Function read_oo_jazpc can read processed transmittance and reflectance output files (expressed as %s). Function read_oo_jazdata can read raw-counts data.

Value

A source_spct object, a filter_spct object, a reflector_spct object or a raw_spct object.

Note

Although the parameter is called date a date time is accepted and expected. Time resolution is 1 s.

References

<https://www.r4photobiology.info> <https://oceanoptics.com/>

read_oo_pidata

Read File Saved by Ocean Optics' Raspberry Pi software.

Description

Reads and parses the header of a raw data file as output by the server running on a Raspberry Pi board to extract the whole header remark field. The time field is retrieved and decoded.

Usage

```
read_oo_pidata(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  npixels = 2048,
  spectrometer.sn = "FLMS00673"
)
```

Arguments

<code>file</code>	character string
<code>date</code>	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is set to the file modification date.
<code>geocode</code>	A data frame with columns lon and lat used to set attribute "where.measured".
<code>label</code>	character string, but if NULL the value of <code>file</code> is used, and if NA the "what.measured" attribute is not set.
<code>tz</code>	character Time zone is not saved to the file.
<code>locale</code>	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
<code>npixels</code>	integer Number of pixels in spectral data.
<code>spectrometer.sn</code>	character The serial number of the spectrometer needs to be supplied by the user as it is not included in the file header.

Value

A `raw_spect` object.

Note

The header in these files has very little information, so the user needs to supply the number of pixels in the array as well as the date-time. The file contains a time in milliseconds but as the Raspberry Pi board contains no real-time clock, it seems to default to number of milliseconds since the Pi was switched on. If no argument is passed to `date` this attribute is set to the file modification date obtained with `file.mtime()`. This date-time gives an upper limit to the real time of measurement as in some operating systems it is reset when the file is copied or even without any good apparent reason.

References

<https://www.r4photobiology.info> <https://oceanoptics.com/> <https://www.raspberrypi.org/>

read_oo_ssirrad *Read File Saved by Ocean Optics' SpectraSuite.*

Description

Reads and parses the header of a processed data file as output by SpectraSuite to extract the whole header remark field. The time field is retrieved and decoded.

Usage

```
read_oo_ssirrad(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

```
read_oo_ssdata(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

Arguments

file	character string
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone is by default read from the file.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.

Value

A source_spect object.

A raw_spect object.

References

<https://www.r4photobiology.info> <https://oceanoptics.com/>

read_qtuv_txt	<i>Read Quick TUV output file.</i>
---------------	------------------------------------

Description

Reads and parses the header of a text file output by the Quick TUV on-line web front-end at http://cprm.acom.ucar.edu/Models/TUV/Interactive_TUV/ to extract the header and spectral data. The time field is converted to a date.

Usage

```
read_qtuv_txt(
  file,
  ozone.du = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

Arguments

file	character string with the name of a text file.
ozone.du	numeric Ozone column in Dobson units.
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone is by default read from the file.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use <code>locale</code> to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.

Value

a `source_spect` object obtained by finding the center of wavelength intervals in the Quick TUV output file, and adding variables `zenith.angle` and `date`.

Note

The ozone column value used in the simulation cannot be retrieved from the file. Tested only with Quick TUV version 5.2 on 2018-07-30. This function can be expected to be robust to variations in the position of lines in the imported file and resistant to the presence of extraneous text or even summaries.

References

<https://www.r4photobiology.info> http://cprm.acom.ucar.edu/Models/TUV/Interactive_TUV/

read_tuv_usrout	<i>Read TUV output file.</i>
-----------------	------------------------------

Description

Reads and parses the header of a text file output by the TUV program to extract the header and spectral data. The time field is converted to a date.

Usage

```
read_tuv_usrout(
  file,
  ozone.du = NULL,
  date = lubridate::today(),
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

```
read_tuv_usrout2mspct(
  file,
  ozone.du = NULL,
  date = lubridate::today(),
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

Arguments

file	character string
ozone.du	numeric Ozone column in Dobson units.
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone is by default read from the file.

locale The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use `locale` to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.

Value

a `source_spect` object obtained by 'melting' the TUV file, and adding a factor `spect.idx`, and variables `zenith.angle` and `date`.

Note

The ozone column value used in the simulation cannot be retrieved from the file. Tested only with TUV version 5.0.

References

<https://www.r4photobiology.info> <https://www2.acom.ucar.edu/modeling/tuv-download>

read_uvspec_disort *Read libRadtran's uvspec output file.*

Description

Read and parse a text file output by libRadtran's `uvspec` routine for a solar spectrum simulation. The output of `uvspec` depends among other things on the solver used. We define a family of functions, each function for a different solver.

Usage

```
read_uvspec_disort(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  multiplier = 0.001,
  qty = "irradiance"
)
```

Arguments

`file` character string

`date` a `POSIXct` object to use to set the "when.measured" attribute. If `NULL`, the default, the date is extracted from the file header.

`geocode` A data frame with columns `lon` and `lat` used to set attribute "where.measured".

label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone is by default read from the file.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
multiplier	numeric A multiplier for conversion into $W\ m^{-2}\ nm^{-1}$, as the units of expression of the output from "uvspec" depend on the units in which the extraterrestrial solar spectrum data is expressed.
qty	character "uvspec" returns both irradiance and intensity with solver "disort".

Value

A source_spect object.

Note

Currently only "irradiance" is supported as qty argument as intensity is not supported by classes and methods in package 'photobiology'.

Tested only with libRadtran version 2.0

References

<https://www.r4photobiology.info> <http://www.libradtran.org>

read_uvspec_disort_vesa

Read libRadtran's uvspec output file from batch job.

Description

Reads and parses the header and body of a text file output by a script used to run libRadtran's uvspec in a batch job for a set of solar spectrum simulations. The header and time and date fields are converted into a datetime object.

Usage

```
read_uvspec_disort_vesa(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  multiplier = 1e-06,
  simplify = TRUE
)
```

Arguments

file	character string
date	a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is extracted from the file header.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
tz	character Time zone is by default read from the file.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.
multiplier	numeric A multiplier for conversion into $W m^{-2} nm^{-1}$, as the units of expression of the output from "uvspec" depend on the units in which the extraterrestrial solar spectrum data is expressed.
simplify	logical Remove redundant columns from returned object.

Value

a source_spect object, possibly containing several spectra in long form and a datetime column.

References

<https://www.r4photobiology.info> <http://www.libradtran.org>

read_yoctopuce_csv *Read '.CSV' file(s) downloaded from YoctoPuce modules.*

Description

Reads and parses the header of processed data CSV files as output by the virtual- or hardware-hubs and modules from Yoctopuce. Uses the comment attribute to store the metadata.

Usage

```
read_yoctopuce_csv(
  file,
  geocode = NULL,
  label = NULL,
  data_skip = 0,
  n_max = Inf,
  locale = readr::default_locale()
)
```

Arguments

file	Path to file as a character string.
geocode	A data frame with columns lon and lat used to set attribute "where.measured".
label	character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set.
data_skip	integer Number of records (rows) to skip from the actual data block.
n_max	integer Maximum number of records to read.
locale	The locale controls defaults that vary from place to place. The default locale is US-centric (like R), but you can use locale to create your own locale that controls things like the default time zone, encoding, decimal mark, big mark, and day/month names.

Value

read_yoctopuce_csv() returns a `tibble::tibble` object.

Note

This function should be able to read data log files from any YoctoPuce USB interface module with data logging capabilities as the format is consistent among them.

References

<https://www.r4photobiology.info> <https://www.yoctopuce.com/>

rspec2mspct	<i>Convert "pavo::rspec" objects</i>
-------------	--------------------------------------

Description

Convert between 'pavo::rspec' objects containing spectral reflectance data into spectral objects (xxxx_spct, xxxx_mspct) as defined in package 'photobiology'.

Usage

```
rspec2mspct(
  x,
  member.class = "reflector_spct",
  spct.data.var = "Rpc",
  multiplier = 1,
  ...
)

rspec2spct(x, multiplier = 1, ...)
```

Arguments

<code>x</code>	rspec object
<code>member.class</code>	character One of the spectrum classes defined in package 'photobiology'.
<code>spct.data.var</code>	character The name to be used for the 'spc' data when constructing the spectral objects.
<code>multiplier</code>	numeric A multiplier to be applied to the 'rspc' data to do unit or scale conversion.
<code>...</code>	currently ignored.

Warning!

Always check the sanity of the imported or exported data values, as guessing is needed when matching the different classes, and the functions defined here are NOT guaranteed to return valid data without help from the user through optional function arguments.

Note

Objects of class `pavo::rspec` do not contain metadata or class data from which the quantity measured and the units of expression could be obtained. When using this function the user needs to use parameter `multiplier` to convert the data to what is expected by the object constructors defined in package 'photobiology' and use parameter `spct.data.var` to select the quantity.

`pavo::rspec` objects may use memory more efficiently than spectral objects of the classes for collections of spectra defined in package 'photobiology' as wavelengths are assumed to be the same for all member spectra, and stored only once while this assumption is not made for collections of spectra, allowing different wavelengths and lengths for the component spectra. Wavelengths are stored for each spectrum, but as spectral classes are derived from 'tbl_df' in many cases no redundant copies of wavelength data will be made in memory in spite of the more flexible semantics of the objects.

Examples

```
# example run only if 'pavo' is available
if (requireNamespace("pavo", quietly = TRUE)) {
  library(pavo)
  data(sicalis, package = "pavo")
  sicalis.mspct <- rspec2mspct(sicalis)
  class(sicalis.mspct)

  data(teal, package = "pavo")
  teal.spct <- rspec2spct(teal)
  class(teal.spct)
  levels(teal.spct[["spct.idx"]])
  angles <- seq(from = 15, to = 75, by = 5) # from teal's documentation
  teal.spct[["angle"]] <- angles[as.numeric(teal.spct[["spct.idx"]])]
  teal.spct
}
}
```

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