

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](https://doi.org/10.19232/uv4pb.2015.1.14)>.

License GPL (>= 2)

VignetteBuilder knitr

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BugReports <https://github.com/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](https://doi.org/10.19232/uv4pb.2015.1.14).

Data acquisition

The support for Ocean Insight, formerly Ocean Optics, spectrometers in package 'photobiologyInOut' is limited to the import of data acquired with Ocean Optics' software as is. In contrast, package 'ooacquire', part of these same suite, makes it possible to control, modify settings and acquire spectral data from Ocean Optics spectrometers directly from within R. 'ooacquire' also supports the conversion of raw-counts data into physical quantities.

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. [doi:10.19232/uv4pb.2015.1.14](https://doi.org/10.19232/uv4pb.2015.1.14).

See Also

Useful links:

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

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, ...)

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

## S3 method for class 'colorSpec'
as.generic_mspct(x, 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. |

Details

Objects of class `colorSpec::colorSpec` do not contain metadata or class data from which the units of expression could be obtained. When using function `colorSpec2mspct` 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. When using `as.colorSpec` methods to convert collections of spectra into `colorSpec` objects, if the wavelengths of the individual spectra differ, only the shared range of wavelengths is retained and within this range, wavelength values are made consistent by interpolation.

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

In `generic_mspct` objects, wavelengths are stored for each spectrum, individual `generic_spct` objects. However, 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)
}
```

mspct2colorSpec

Convert into 'colorSpec::colorSpec' objects

Description

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

Usage

```
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, ...)
```

```
## 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

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

Methods (by class)

- `as.colorSpec(generic_spct):`
- `as.colorSpec(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)) {
  library(colorSpec)
  as.colorSpec(polyester.spct)
  as.colorSpec(sun.spct)
  as.colorSpec(filter_mspct(list(polyester.spct, yellow_gel.spct)))
}
```

mspct2fdata

Convert spectra into 'fda.usc::fdata' objects

Description

Convert spectral objects (xxxx_spct, xxxx_mspct) as defined in package 'photobiology' into `fda.usc::fdata` objects, preserving as much information as possible. As `fdata` objects can contain other kinds of data, the reverse conversion is supported (experimentally) and mainly for 'fdata' objects returned by the functional data analysis methods from package `fda.usc` to spectral data previously exported in the opposite direction.

Usage

```
mspct2fdata(x, spct.data.var = NULL, multiplier = 1, ...)
```

```
spct2fdata(x, spct.data.var = NULL, multiplier = 1, ...)
```

```
fdata2spct(x, multiplier = 1, member.class = NULL, drop.idx = FALSE, ...)
```

```
fdata2mspct(x, multiplier = 1, member.class = NULL, drop.idx = FALSE, ...)
```

Arguments

| | |
|----------------------------|--|
| <code>x</code> | generic_mspct or generic_spct object or an object belonging to a derived class, or an object of class 'fdata' depending on the function. |
| <code>spct.data.var</code> | character The name of the column containing data to export. If NULL the first spectral data column found is used. |
| <code>multiplier</code> | numeric A multiplier to be applied to the 'spc' data to do unit or scale conversion. |
| <code>...</code> | possibly additional named arguments passed to object constructors. |
| <code>member.class</code> | character Name of the class of the spectrum or of the members of the collection of spectra. |
| <code>drop.idx</code> | logical Flag indicating whether to drop or keep <code>idx.var</code> in the collection members. |

Warning!

When converting multiple spectra, all the spectra to be included in the `fdata` object must share the same wavelength values. Spectra that do not fulfil this condition will be skipped. The data variable needs also to be present in all individual spectra as no conversions are applied automatically by this function. If a different name, indicating a different quantity or a different base of expression is encountered, the affected spectrum is skipped with a warning.

Examples

```

if (requireNamespace("fda.usc", quietly = TRUE)) {
# from spectra to fdata
  sun.fdata <- spct2fdata(sun.spct)
  str(sun.fdata)
  polyester.fdata <- spct2fdata(polyester.spct)
  str(polyester.fdata)
# from fdata to spectra
  fdata2spct(sun.fdata)
  fdata2spct(sun.fdata, drop.idx = TRUE)
  fdata2spct(polyester.fdata, drop.idx = TRUE)
}

```

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 = Inf
)

```

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_spct object.

Note

The header in these files has metadata information, but mostly on the origin of the data. For a date and/or geocode are to be added to the return object it must be supplied by the user. as well as the date-time. Some metadata is extracted and added as attributes, while the whole header is copied to the comment attribute.

References

<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

Examples

```
file.name <-
  system.file("extdata", "drygrass-spectrum.txt",
             package = "photobiologyInOut", mustWork = TRUE)

fred.spct <- read_ASTER_txt(file = file.name, npixels = Inf)

fred.spct
getWhatMeasured(fred.spct)
cat(comment(fred.spct))
```

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_spect object.

References

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

Examples

```

file.name <-
  system.file("extdata", "spectrum-avaspec.csv",

```

```

package = "photobiologyInOut", mustWork = TRUE)

avaspec.spct <- read_avaspec_csv(file = file.name)

avaspec.spct
getWhatMeasured(avaspec.spct)
cat(comment(avaspec.spct))

```

```
read_cid_spectravue_csv
```

Read File Saved by CID's SpectraVue.

Description

Read wavelength and spectral data from Measurements.CSV files exported from CID Bio-Sciences' **SpectraVue CI-710s** (not the older CI-710!) leaf spectrometer, importing them into R. Available metadata is also extracted from the file. `read_cid_spectravue_csv()` only accepts "row oriented" CSV files. These may contain multiple spectra, one per row.

Usage

```

read_cid_spectravue_csv(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  range = c(380, 1100),
  simplify = TRUE,
  absorbance.to = "object",
  strict.range = NA,
  ...
)

```

Arguments

| | |
|----------------------|--|
| <code>file</code> | character |
| <code>date</code> | a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date and time are extracted from the file. |
| <code>geocode</code> | A data frame with columns lon and lat used to set attribute "where.measured". If NULL, the default, the geocode is extracted from the file, if present, and if NA the "where.measured" attribute is not set. |
| <code>label</code> | character string. If NULL, the default, the value of the "tag" present in the file is used, and if NA the "what.measured" attribute is not set. |
| <code>tz</code> | character Time zone is by default that of the machine's locale. |

| | |
|---------------|--|
| 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. |
| range | numeric A vector of length two, or any other object for which function <code>coderange()</code> will return range of wavelengths expressed in nanometres. |
| simplify | logical If TRUE, single spectra are returned as individual spectra instead of collections of length one. |
| absorbance.to | character Affects only absorbance measurements: "object", "all" or "A". |
| strict.range | logical Flag indicating whether off-range values result in an error (TRUE) instead of a warning (FALSE), or the test is disabled (NA). |
| ... | additional arguments passed to the constructor of the 'filter_spct' object. |

Details

SpectraVue's row-wise spectral Measurements.CSV files contain columns with metadata on the right edge, followed by columns with data for each of the 2048 pixels or wavelengths. The value in column "Mode" indicates the quantity measured, decoded into variables Tpc, Rpc or A. The data the rows in the CSV file are read and stored in `filter_spct`, `reflector_spct` or `object_spct` objects. These objects are collected into a single `filter_mspct`, `reflector_mspct`, `object_mspct` or `generic_spct` object and returned.

Spectral data outside the range 400 nm to 1000 nm are very noisy and thus outside the valid range for the measurements. Out-of-range spectral data can also be cause by calibration drift. Consequently, reading of data is done always with the range check disabled, while whether a check is used before returning the collection of spectra depends on the argument passed to `strict.range` which by default is set to disable checks. This is done, because in most cases measurements from this instrument tend to require further processing before they comply with theoretical expectations of $Tfr + Rfr + Afr = 1$.

Value

An object of class `filter_spct`, `relector_spct`, `object_spct` or `generic_mspct`.

Internal vs. total transmittance and absorbance

Spectravue returns transmittance values labelled with Transmittance as Mode. Transmittance values are not *total* as most of the scattered light transmitted is not detected. Absorbance (Abs) values returned labelled with Absorbance as Mode are for absorbance computed from the Transmittance. This estimate of absorbance overestimates real absorbance in the case of scattering materials like plant leaves. It is best to save spectral data acquired as absorbance into objects of class 'object_spct' containing reflectance and transmittance, the default, as this preserves all the acquired data.

Specular vs. total reflectance

This function assumes that SpectraVue returns close to *total* reflectance readings. Given the optics of the instrument this is likely only an approximation.

Warning!!

CID's support has answered on 2022-05-19 that the extremely biased (plainly wrong!) values of transmittance measured by this instrument are due to a design flaw and that they are working on a solution for the problem. In practice, reflectance seems biased but usable as an instrument-specific quantity with arbitrary units. Transmittance and absorbance seem useless as values are wrong by about an order of magnitude.

Note

SpectraVue creates three or four .CSV files for each series of measurements saved. Of these files, this function reads the one with name ending in Measurements.CSV. The first part of the file name gives the time of the session, but as the files can contain multiple spectra measured at different times, the time metadata is extracted separately for each spectrum. We provide a default argument for range that discards data for short and long wavelengths because values outside this range are according to the instrument's manual outside the usable range and in practice extremely noisy.

References

<https://cid-inc.com/>

Examples

```
# read file containing a single reflectance spectrum

file.name <-
  system.file("extdata", "cid-spectravue-Rpc-Measurements.csv",
             package = "photobiologyInOut", mustWork = TRUE)

cid_filter.spct <-
  read_cid_spectravue_csv(file = file.name)
summary(cid_filter.spct)

cid_filter.spct <-
  read_cid_spectravue_csv(file = file.name, simplify = FALSE)
summary(cid_filter.spct)

# read data measured as absorbance (A, Rpc and Tpc)

file.name <-
  system.file("extdata", "cid-spectravue-Abs-Measurements.csv",
             package = "photobiologyInOut", mustWork = TRUE)
cid.object_spct <-
  read_cid_spectravue_csv(file = file.name)
summary(cid.object_spct)

cid_A.filter_spct <-
  read_cid_spectravue_csv(file = file.name, absorbance.to = "A")
summary(cid_A.filter_spct)
```

| | |
|--------------|---|
| 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(),
  na = c("", "NA", "NAN"),
  ...
)
```

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. |
| na | character Vector of strings to interpret as missing values. Set this option to character() to indicate no missing values. |
| ... | Further named arguments currently passed to read_csv(). |

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.campbellsci.eu/>

| | |
|----------------|--|
| read_fmi2mspct | <i>Read multiple solar spectra from a data file.</i> |
|----------------|--|

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

| | |
|------------------------|--|
| file | Either a path to a file, a connection, or literal data (either a single string or a raw vector). |
| scale.factor | numeric A multiplier to be applied to the spectral irradiance values. |
| geocode | A data frame with columns lon and lat used to set attribute "where.measured". |
| what.measured | character string, but if NULL the value of file is used, and if NA the "what.measured" attribute is not set. |
| how.measured | character string, but if NULL or NA the "how.measured" attribute is not set. |
| date.field, time.field | integer. Word positions in the header line. |
| date.format | character string. One of "ymd", "ydm", "dmy", or "mdy". |
| time.format | character string. One of "hms", "hm". |
| tz | character Time zone used for interpreting times saved in the file header. |

| | |
|-----------------|---|
| time.shift.min, | numeric. Time shift with respect to TZ in minutes. |
| 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. |
| .n_max | 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

| | |
|---------|---|
| file | 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_mspct 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_foreign2mspct *Read multiple foreign files with spectral data*

Description

Read spectra from a homogeneous list of files based on a path and a list of filenames or a path and a search pattern for files. The imported spectra are returned as a single object of one of the collection of spectra classes from package 'photobiology'.

Usage

```
read_foreign2mspct(path = ".", list = NULL, pattern = NULL, .fun, ...)
```

Arguments

| | |
|---------|--|
| path | character A path point to the location of the files. |
| list | character A vector or list of character strings pointing to files relative to path, |
| pattern | character A search pattern to select files within path. See list.files which is used internally. Argument ignored is list is non-null. |
| .fun | function One of the functions exported by this package for reading spectral data. |
| ... | Named arguments passed ot the call to .fun. |

Details

This function iterates over a list of file names reading them with the function passed as argument to '.fun' and combines the spectra as a collection of spectra of a class suitable for the spectral objects returned by the argument to '.fun'. This function can either return for each file read either a single spectrum as an object of class 'generic_spct' or a class derived from it, or a collection of spectra of class 'generic_mspct' or a class derived from it. The class of the returned object depends on the class of the member spectra.

Value

An object of class 'generic_mspct' or a class derived from it, containing a collection of member spectra of class 'generic_spct' or of one of the classes derived from it.

| | |
|---------------|-----------------------------------|
| read_FReD_csv | <i>Read '.CSV' FReD database.</i> |
|---------------|-----------------------------------|

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. Those relevant should match the format of the CSV file being read. |

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

Examples

```

file.name <-
  system.file("extdata", "FReDflowerID_157.csv",
             package = "photobiologyInOut", mustWork = TRUE)

fred.spct <- read_FReD_csv(file = file.name)

fred.spct
getWhatMeasured(fred.spct)
cat(comment(fred.spct))

```

read_li180_txt

Read '.TXT' File(s) Saved by LI-COR's LI-180 spectroradiometer.

Description

Reads and parses the header of a data file as output by the LI-180 spectrometer (not to be confused with the LI-1800 spectrometer released in the 1980's by LI-COR) to extract the whole header remark field and also decode whether data is in photon or energy based units. This is a new instrument released in year 2020.

Usage

```

read_li180_txt(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  s.qty = "s.e.irrad"
)

read_m_li180_txt(
  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 <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. |
| s.qty | character The name of the spectral quantity to be read. One of "s.e.irrad" or "s.q.irrad". |
| files | A list or vector of character strings. |

Details

Function `read_m_licor_espd()` calls `red_licor_espd()` for each file in `files`. See [read.table](#) for a description of valid arguments for files.

Value

`read_licor_espd()` 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 header, or supplied by the user. Spectrometer model, serial number and integration time are stored in attributes. The whole file header is saved as a comment while the footer is discarded.

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

Note

The LI-180 spectroradiometer stores little information of the instrument and settings, possibly because they cannot be altered by the user or configured. The length of the file header does not seem to be fixed, so the start of the spectral data is detected by searching for "380nm".

References

LI-COR Biosciences, Environmental. <https://www.licor.com/env/>

Examples

```
file.name <-  
  system.file("extdata", "LI-180-irradiance.txt",  
             package = "photobiologyInOut", mustWork = TRUE)
```

```

licor180.spct <- read_li180_txt(file = file.name)

licor180.spct
getWhenMeasured(licor180.spct)
getWhatMeasured(licor180.spct)
cat(comment(licor180.spct))

```

| | |
|----------------|--|
| read_licor_prn | <i>Read '.PRN' File(s) Saved by LI-COR's PC1800 Program.</i> |
|----------------|--|

Description

Read and parse 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 irradiance in photon or energy based units, transmittance, reflectance or absorbance and then extract the wavelength and spectral data. PC1800 is an MS-DOS program provided for use with the LI-1800 spectrometer. This instrument was released in the 1980's and was sold until the early 2000's. It was very popular and several of them remain in use. (It should not be confused with the LI-180, a new spectrometer released in 2020.)

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 `red_licor_prn()` 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

LI-COR Biosciences, Environmental. <https://www.licor.com/env/>

Examples

```
file.name <-
  system.file("extdata", "spectrum.PRN",
             package = "photobiologyInOut", mustWork = TRUE)

licor.spct <- read_licor_prn(file = file.name)

licor.spct
getWhenMeasured(licor.spct)
getWhatMeasured(licor.spct)
cat(comment(licor.spct))
```

read_macam_dta *Read '.DTA' File Saved by Macam's Software.*

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

| | |
|---------|---|
| 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 source_spct object.

References

<https://www.irradian.co.uk/>

Examples

```
file.name <-
  system.file("extdata", "spectrum.DTA",
             package = "photobiologyInOut", mustWork = TRUE)

macam.spct <- read_macam_dta(file = file.name)

macam.spct
getWhenMeasured(macam.spct)
getWhatMeasured(macam.spct)
cat(comment(macam.spct))
```

| | |
|------------------|--|
| read_oo_jazirrad | <i>Read Files Saved by Ocean Optics' Jaz spectrometer.</i> |
|------------------|--|

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 were manufactured by Ocean Optics. The company formerly named Ocean Optics is now called Ocean Insight.

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 if seconds are entered with a decimal fraction, such as "2021-10-05 10:10:10.1234".

References

<https://www.oceaninsight.com/>

Examples

```

file.name <-
  system.file("extdata", "spectrum.jaz",
             package = "photobiologyInOut", mustWork = TRUE)

jaz.spct <- read_oo_jazpc(file = file.name)

jaz.spct
getWhenMeasured(jaz.spct)
getWhatMeasured(jaz.spct)
cat(comment(jaz.spct))

```

| | |
|----------------|--|
| read_oo_pidata | <i>Read File Saved by Ocean Optics' Raspberry Pi software.</i> |
|----------------|--|

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. The company formerly named Ocean Optics is now called Ocean Insight.

Usage

```

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

```

Arguments

| | |
|---------|--|
| file | character string |
| date | a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date is set to the file modification date. |
| 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 not saved to 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. |
| npixels | integer Number of pixels in spectral data. |
| spectrometer.sn | 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. 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. The user may wish to supply the date-time as an argument, but 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. The user may need to supply the number of pixels in the array although the default of `npixels = Inf` usually works and triggers no warnings.

References

<https://www.oceaninsight.com/> <https://www.raspberrypi.org/>

Examples

```
file.name <-
  system.file("extdata", "spectrum.pi",
             package = "photobiologyInOut", mustWork = TRUE)

oopi.spct <- read_oo_pidata(file = file.name)

oopi.spct
getWhenMeasured(oopi.spct)
getWhatMeasured(oopi.spct)
cat(comment(oopi.spct))
```

| | |
|-----------------|---|
| read_oo_ssirrad | <i>Read File Saved by Ocean Optics' SpectraSuite.</i> |
|-----------------|---|

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. SpectraSuite was a program, now replaced by OceanView. The company formerly named Ocean Optics is now called Ocean Insight.

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.oceaninsight.com/>

Examples

```
file.name <-
  system.file("extdata", "spectrum.SSIrrad",
             package = "photobiologyInOut", mustWork = TRUE)

ooss.spct <- read_oo_ssirrad(file = file.name)

ooss.spct
getWhenMeasured(ooss.spct)
getWhatMeasured(ooss.spct)
cat(comment(ooss.spct))
```

| | |
|---------------|------------------------------------|
| 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 UCAR 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 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 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.acom.ucar.edu/Models/TUV/Interactive_TUV/

read_spectrapen_csv *Read '.CSV' File Saved by PSI's Software.*

Description

Reads and parses the header of a processed .CSV file as output by the by the PSI (Photon Systems Instruments, Czech Republic) SpectraPen miniature spectrometer.

Usage

```
read_spectrapen_csv(
  file,
  start.row = 1,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale()
)
```

Arguments

| | |
|-----------|---|
| file | character string. |
| start.row | integer The first line to read, counting from the top of the file. |
| 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". If NULL, the default, the location is extracted from the file header. |
| label | character string, but if NULL the label from file header is used, if the label is missing, the index 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 `source_mspct` object.

References

<https://psi.cz/>

Examples

```
# fetch path to example file to read
file.name <-
  system.file("extdata", "spectrum-psi-spectrapen-SP.csv",
             package = "photobiologyInOut", mustWork = TRUE)

spectrapen.mspct <- read_spectrapen_csv(file = file.name)

spectrapen.mspct
getWhenMeasured(spectrapen.mspct)
getWhatMeasured(spectrapen.mspct)
cat(comment(spectrapen.mspct))
```

| | |
|-----------------|------------------------------|
| 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_spct object obtained by 'melting' the TUV file, and adding a factor spct.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://www2.acom.ucar.edu/modeling/tropospheric-ultraviolet-and-visible-tuv-radiation-model>

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

References

<http://www.libradtran.org>

| | |
|------------------|--|
| read_wasatch_csv | <i>Read File Saved by Wasatch's Enlighten.</i> |
|------------------|--|

Description

Read wavelength and spectral data from the data section of a file as output by Enlighten importing them into R. Parse the header of a file to extract the acquisition time, instrument name and serial number, as well additional metadata related to the instrument and its settings. Function read_wasatch_csv() only accepts "column oriented" CSV files.

Usage

```
read_wasatch_csv(
  file,
  date = NULL,
  geocode = NULL,
  label = NULL,
  tz = NULL,
  locale = readr::default_locale(),
  s.qty = NULL,
  extra.cols = "keep",
  scale.factor = 1,
  simplify = TRUE,
  ...
)
```

Arguments

| | |
|---------|---|
| file | character |
| date | a POSIXct object to use to set the "when.measured" attribute. If NULL, the default, the date and time are 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. |

| | |
|---------------------------|--|
| <code>tz</code> | character Time zone is by default that of the machine's locale. |
| <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. |
| <code>s.qty</code> | character, possibly named. The name of the quantity using the conventions accepted used in package 'photobiology' that is to be imported from column "Processed" from the file. |
| <code>extra.cols</code> | character What to do non-processed data columns if present in file. One of "keep", "drop.pixel", "drop" or "split". |
| <code>scale.factor</code> | numeric vector of length 1, or length equal to the number of rows (= detector pixels). Numeric multiplier applied to returned spectral values. |
| <code>simplify</code> | logical If TRUE, single spectra are returned as individual spectra instead of collections of length one. |
| <code>...</code> | additional arguments passed to the constructor of the spectrum object. |

Details

Enlighten's column-wise CSV files contain at least two columns, Wavelength and Processed. In the header the Technique used is recorded. Additional data columns can be present. Column Pixel contains the pixel index in the array as integers. Columns Raw, Dark and Reference contain detector counts data. Technique is used to guess the type of spectrum stored in the column named Processed, which can be detector counts or derived values. By default the data are read into a single spectrum object and all columns retained, but only the data in Processed are interpreted as spectral data corresponding to the class of the object. If passed `extra.cols = "drop"`, only Wavelength and Processed are copied to the returned object, while if passed `extra.cols = "drop.pixel"` only the contents of column Pixel are discarded. If passed `extra.cols = "split"` all columns containing spectral data are each read into a separate spectrum, these are collected and a "generic_mspect" object containing them returned. `extra.cols` can be a named vector of mappings, of length at least one but possibly longer. If longer a "generic_mspect" is returned, otherwise a spectrum object as inferred from the name each column is mapped to.

Value

An object of a class derived from `generic_spct` such as `raw_spct` or `filter_spct`. `generic_spct` is derived from tibble and data frame.

Acknowledgements

We thank Ruud Niesen from Photon Mission (<https://photonmission.com/>) for organizing the loan of the spectrometer used to produce the various files needed for the development of this function.

Note

Enlighten, the free software from Wasatch Photonics can save spectra in a variety of additional formats: different types of CSV files, plain text and JSON. Plain text files contain no metadata or even column headers and if the need arises can be read with R function `read.table()`. JSON files contain the most detailed metadata.

References

<https://wasatchphotonics.com/> <https://wasatchphotonics.com/product-category/software/>

Examples

```
file.name <-
  system.file("extdata", "enlighten-wasatch-scope.csv",
             package = "photobiologyInOut", mustWork = TRUE)

wasatch.raw.spct <-
  read_wasatch_csv(file = file.name)
summary(wasatch.raw.spct)

wasatch.raw.spct <-
  read_wasatch_csv(file = file.name, s.qty = "counts")
summary(wasatch.raw.spct)

wasatch.raw.spct <-
  read_wasatch_csv(file = file.name, s.qty = c(Processed = "counts"))
summary(wasatch.raw.spct)

wasatch.raw.spct <-
  read_wasatch_csv(file = file.name, extra.cols = "drop")
summary(wasatch.raw.spct)
```

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. |

Details

Yoctopuce modules are small USB connected and USB powered, but isolated, very high quality miniature data acquisition and interface modules. All modules capable of data acquisition can log measured data autonomously and these data can be locally or remotely downloaded as a CSV file. (It is also possible and very easy to access these modules from R using package 'reticulate' and the Python library provided by Yoctopuce, or to send commands and retrieve data through the built-in HTML server of the modules or dedicated hubs.)

Value

read_yoctopuce_csv() returns a `tibble::tibble` object, with the number of columns dependent on the CSV file read.

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.yoctopuce.com/>

Examples

```
# We read a CSV file previously downloaded from a YoctoMeteo module.

file.name <-
  system.file("extdata", "yoctopuce-data.csv",
             package = "photobiologyInOut", mustWork = TRUE)

yoctopc.tb <- read_yoctopuce_csv(file = file.name)

yoctopc.tb
cat(comment(yoctopc.tb))
```

rspec2mspct *Convert "pavo::rspec" objects*

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

| | |
|---------------|--|
| x | rspec object |
| member.class | character One of the spectrum classes defined in package 'photobiology'. |
| spct.data.var | character The name to be used for the 'spc' data when constructing the spectral objects. |
| multiplier | numeric A multiplier to be applied to the 'rspc' data to do unit or scale conversion. |
| ... | 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.mspect <- rspec2mspect(sicalis)
  class(sicalis.mspect)

  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
}
```

spct_CCT

Correlated color temperature

Description

Wrapper on function `computeCCT` from package 'colorSpec' that accepts `source_spct` objects.

Usage

```
spct_CCT(
  spct,
  isotherms = "robertson",
  locus = "robertson",
  strict = FALSE,
  named = FALSE
)
```

Arguments

| | |
|------------------------|---|
| <code>spct</code> | source_spct A single light source spectrum. |
| <code>isotherms</code> | character A vector whose elements match one of the available isotherm families: 'robertson', 'mccamy', and 'native'. Matching is partial and case-insensitive. When more than one family is given, a matrix is returned, see Value. When <code>isotherms = 'native'</code> the isotherms are defined implicitly as lines perpendicular to the locus, see Details in <code>CCTfromXYZ</code> . The character NA (<code>NA_character_</code>) is taken as a synonym for 'native'. |

| | |
|--------|---|
| locus | character Valid values are 'robertson' and 'precision', see above. Matching is partial and case-insensitive. |
| strict | logical The CIE considers the CCT of a chromaticity uv to be meaningful only if the distance from uv to the Planckian locus is less than or equal to 0.05 (in CIE UCS 1960). If strict=FALSE, then this condition is ignored. Otherwise, the distance is computed along the corresponding isotherm, and if it exceeds 0.05 the returned CCT is set to NA. |
| named | logical Whether to set the name attribute of the returned value to the name of the spectrum passed as argument if possible. |

Details

Please see [computeCCT](#) for the details of the computations and references.

Value

A numeric value for "color temperature " in degrees Kelvin.

Examples

```
spct_CCT(white_led.source_spct)
spct_CCT(sun.spct)
```

spct_CRI

Color reproduction index

Description

Wrapper on function [computeCRI](#) from package 'colorSpec' that accepts [source_spct](#) objects.

Usage

```
spct_CRI(spct, adapt = TRUE, attach = FALSE, tol = 0.0054, named = FALSE)
```

Arguments

| | |
|--------|--|
| spct | source_spct A single light source spectrum. |
| adapt | logical If TRUE, then a special chromatic adaption is performed, see Details in computeCRI . |
| attach | logical If TRUE, then a large list of intermediate calculations is attached to the returned number, as attribute data. This attached list includes data for all special 14 color samples, although the last 6 do not affect the returned CRI |

| | |
|-------|--|
| tol | numeric For the CRI to be meaningful the chromaticities of the test and reference illuminants must be sufficiently close in the CIE 1960 uniform chromaticity space. If the tolerance is exceeded, the function returns NA. The default $\text{tol}=5.4\text{e-}3$ is the one recommended by the CIE, but the argument allows the user to override it. |
| named | logical Whether to set the name attribute of the returned value to the name of the spectrum passed as argument if possible. |

Details

Please see [computeCRI](#) for the details of the computations and references.

Value

A numeric value between zero and 100, or NA if the light is not white enough.

Examples

```
spct_CRI(white_led.source_spct)
spct_CRI(sun.spct)
```

| | |
|----------|--|
| spct_SSI | <i>Spectral (color) similarity index</i> |
|----------|--|

Description

Wrapper on function [computeSSI](#) from package 'colorSpec' that accepts [source_spct](#) objects.

Usage

```
spct_SSI(
  spct,
  reference.spct = NULL,
  digits = 0,
  isotherms = "mccamy",
  locus = "robertson",
  named = FALSE
)
```

Arguments

| | |
|----------------------|---|
| spct, reference.spct | source_spct Single light source spectra. |
| digits | integer The number of digits after the decimal point in the returned vector. According to Holm the output should be rounded to the nearest integer, which corresponds to $\text{digits} = 0$. To return full precision, set $\text{digits} = \text{Inf}$. |

| | |
|-----------|--|
| isotherms | character This is only used when reference=NULL. It is passed to computeCCT in order to compute the CCT of each test spectrum. |
| locus | character This is only used when reference=NULL. It is passed to computeCCT in order to compute the CCT of each test spectrum. |
| named | logical Whether to set the name attribute of the returned value to the name of the spectrum passed as argument if possible. |

Details

Please see [computeSSI](#) for the details of the computations and references.

Value

A numeric value between zero and 100.

Examples

```
spct_SSI(white_led.source_spct, sun.spct)
```

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