Package 'fluxfinder'

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Type Package

Title Parsing, Computation, and Diagnostics for Greenhouse Gas Measurements

Version 1.2.2

Description Parse static-chamber greenhouse gas measurement files generated by a variety of instruments; compute flux rates using multi-observation metadata; and generate diagnostic metrics and plots. Designed to be easy to integrate into reproducible scientific workflows.

URL https://github.com/COMPASS-DOE/fluxfinder

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Encoding UTF-8

Imports broom (>= 1.0), jsonlite (>= 1.8), lubridate (>= 1.0), MASS (>= 7.0)

Suggests gasfluxes, ggplot2, knitr, rmarkdown, testthat (>= 3.0.0), withr (>= 2.0)

Config/testthat/edition 3

RoxygenNote 7.3.2

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BugReports https://github.com/COMPASS-DOE/fluxfinder/issues

NeedsCompilation no

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Contents

`ffi_compute_fluxes` *Compute fluxes for multiple groups (measurements)*

Description

Compute fluxes for multiple groups (measurements)

Usage

```
ffi_compute_fluxes(
  data,
  group_column,
  time_column,
  gas_column,
  dead_band = 0,
  normalize_time = TRUE,
  fit_function = ffi_fit_models,
  ...
)
```

Arguments

`data`	A `data.frame` (or tibble or data.table)
`group_column`	Name of the grouping column in `data`, character; pass NULL to run with no grouping
`time_column`	Name of the time column in `data`, character
`gas_column`	Name of the gas (concentration or quantity) column in `data`, character

dead_band	Length of the dead band in seconds (numeric), the equilibration period whose data is dropped. This can be either a single number OR the name of the dead band column in data
normalize_time	Normalize the values so that first is zero? Logical
fit_function	Optional flux-fit function; default is ffi_fit_models
...	Other parameters passed to fit_function

Value

A data.frame with one row per group_column value. It will always include the mean, minimum, and maximum values of time_column for that group, but other columns depend on what is returned by the fit_function.

See Also

ffi_fit_models

Examples

```
# No grouping
ffi_compute_fluxes(cars, group_column = NULL, "speed", "dist")
# With grouping
cars$Plot <- c("A", "B")
ffi_compute_fluxes(cars, "Plot", "speed", "dist")
# See the introductory vignette for a fully-worked example with real data
```

---

ffi_fit_models *Fit various models to gas concentration data*

Description

Fit various models to gas concentration data

Usage

```
ffi_fit_models(time, conc, area, volume)
```

Arguments

time	Relative time of observation (typically seconds), numeric
conc	Greenhouse gas concentration (typically ppm or ppb), numeric
area	Area covered by the measurement chamber (typically cm2), numeric
volume	Volume of the system (chamber + tubing + analyzer, typically cm3), numeric

Details

If a linear model cannot be fit, NULL is returned. If the robust linear and/or polynomial models cannot be fit, then NA is returned for their particular statistics. The HM1981 approach is only valid for saturating (exponential) data and NA is returned otherwise.

Value

A wide-form data.frame with fit statistics for linear ("lin", lm), robust linear ("rob", rlm), polynomial ("poly"), and H&M1981 ("HM81", ffi_hm1981) models. The latter is based on an exponential model drawn from one-dimensional diffusion theory; see Hutchinson and Mosier (1981) and Nakano et al. (2004).

For each model type, the following columns are returned:

- Model statistics AIC, r.squared, RMSE, and p.value;
- Flux (slope) statistics flux.estimate and flux.std.error;
- Intercept statistics int.estimate and int.std.error;
- For the robust linear regression model only, a logical value converged.

Note

Normally this is not called directly by users, but instead via ffi_compute_fluxes.

References

Nakano, T., Sawamoto, T., Morishita, T., Inoue, G., and Hatano, R.: A comparison of regression methods for estimating soil–atmosphere diffusion gas fluxes by a closed-chamber technique, Soil Biol. Biochem., 36, 107–113, 2004. doi:10.1016/j.soilbio.2003.07.005

Hutchinson, G. L. and Mosier, A. R.: Improved soil cover method for field measurement of nitrous oxide fluxes, Soil Sci. Soc. Am. J., 45, 311-316, 1981. doi:10.2136/sssaj1981.03615995004500020017x

Examples

```
# Toy data - linear
ffi_fit_models(cars$speed, cars$dist)

# Toy data - nonlinear
ffi_fit_models(Puromycin$conc, Puromycin$rate)

# Real data
f <- system.file("extdata/TG10-01087.data", package = "fluxfinder")
dat <- ffi_read_LI7810(f)[1:75,] # isolate first observation
dat$SECONDS <- dat$SECONDS - min(dat$SECONDS) # normalize time to start at 0
plot(dat$SECONDS, dat$CO2)
ffi_fit_models(dat$SECONDS, dat$CO2)
```

ffi_hm1981 *Compute flux using nonlinear Hutchinson and Mosier (1981) model*

Description

Compute flux using nonlinear Hutchinson and Mosier (1981) model

Usage

```
ffi_hm1981(time, conc, h = 1)
```

Arguments

time	Time values, numeric
conc	Gas concentration values, numeric
h	Effective chamber height

Value

Flux estimate; see references for more information.

References

Hutchinson, G. L. and Mosier, A. R.: Improved soil cover method for field measurement of nitrous oxide fluxes, Soil Sci. Soc. Am. J., 45, 311-316, 1981. doi:10.2136/sssaj1981.03615995004500020017x

Examples

```
# If data are approximately linear, then NA is returned
ffi_hm1981(cars$speed, cars$dist)
# If data are nonlinear (saturating) then flux based on gas diffusion theory
ffi_hm1981(Puromycin$conc, Puromycin$rate)
```

ffi_metadata_match *Match metadata info with a vector of data timestamps*

Description

Match metadata info with a vector of data timestamps

Usage

```
ffi_metadata_match(data_timestamps, start_dates, start_times, obs_lengths)
```

Arguments

data_timestamps	Data timestamps, either character (YYYY-MM-DD HH:MM:SS) or POSIXct
start_dates	Metadata measurement dates, either character (YYYY-MM-DD) or POSIXct
start_times	Metadata measurement start time entries, either character (HH:MM:SS) or period
obs_lengths	Observation lengths in seconds, numeric; must be same length as start_dates. This should include both the intended measurement period as well as any dead band time at the beginning

Value

A numeric vector equal in length to data_timestamps, with each entry indicating the metadata entry that should be used for that observation. NA is returned if a timestamp has no match in the metadata (i.e., does not fall within any window defined by the start_dates, start_times, and observation length parameters).

Note

If data_timestamps or start_dates cannot be parsed as YYYY-MM-DD, the preferred format, then MM/DD/YYYY (used by U.S. versions of Microsoft Excel when saving CSV files, for example) will be tried.

Examples

```
# Data timestamps
d_t <- c("2024-01-01 13:00:05", "2024-01-01 13:00:10",
"2024-01-01 13:05:05", "2024-01-01 13:10:00")
# Metadata start dates and times: two measurements, starting 5 minutes apart
s_d <- c("2024-01-01", "2024-01-01")
s_t <- c("13:00:00", "13:05:00")
ol <- c(60, 60) # Observation lengths
ffi_metadata_match(d_t, s_d, s_t, ol)
# Returns {1, 1, 2, NA} indicating that the first and second data timestamps
# correspond to metadata entry 1, the third to entry 2, and the fourth
# has no match

# This generates an error because of overlapping timestamps:
try({
s_t <- c("13:00:00", "13:01:00")
ffi_metadata_match(d_t, s_d, s_t, ol)
})
```

---

ffi_qaqc *Generate a QA/QC document*

Description

Generate a QA/QC document

Usage

```
ffi_qaqc(
  flux_data,
  group_column,
  output_file = "qaqc.html",
  output_dir = getwd(),
  open_output = TRUE
)
```

Arguments

flux_data	A data frame from ffi_compute_fluxes or similar
group_column	Name of the grouping label column in flux_data, character; pass NULL to run with no grouping
output_file	Name of the output file
output_dir	Name of the output directory; default is current working directory
open_output	Automatically open the output HTML file?

Value

The path of the output file.

Examples

```
# We don't run this example because of tempfile errors on CRAN,
# but it works fine otherwise!
## Not run:
# Toy data
cars$Plot <- c("A", "B")
fd <- ffi_compute_fluxes(cars, "Plot", "speed", "dist")
x <- ffi_qaqc(fd, group_column = "Plot")
file.remove(x) # clean up

## End(Not run)
# See the introductory vignette for a fully-worked example with real data
```

---

ffi_read_EGM4 *Read an EGM-4 data file*

Description

Read an EGM-4 data file

Usage

```
ffi_read_EGM4(file, year, tz = "UTC")
```

Arguments

file	Filename to read, character
year	Four-digit year of the data (EGM-4 output files have month, day, hour, and minute, but not year), numeric or character
tz	Time zone of the file's time data, character (optional)

Value

A data.frame with the parsed data.

Examples

```
f <- system.file("extdata/EGM4-data.dat", package = "fluxfinder")
dat <- ffi_read_EGM4(f, 2023)
dat <- ffi_read_EGM4(f, 2023, tz = "EST") # specify time zone
```

---

ffi_read_LGR915 *Read a LGR 915-0011 data file*

Description

Read a LGR 915-0011 data file

Usage

```
ffi_read_LGR915(file, date_format = "DMY", tz = "UTC")
```

Arguments

file	Filename to read, character
date_format	Date format, character: "MDY" (month-day-year) "DMY" (day-month-year), or "YMD" (year-month-day)
tz	Time zone of the file's time data, character (optional)

Details

The LGR 915-0011 was an Ultra-Portable Greenhouse Gas Analyzer made by Los Gatos Research. The date in its output files can appear in different formats, which is why the date_format parameter is needed.

Value

A data.frame with the parsed data.

Note

Some LGR 915 files can have a PGP block at the end; this is ignored.

Examples

```
f <- system.file("extdata/LGR-data.csv", package = "fluxfinder")
dat <- ffi_read_LGR915(f, date_format = "MDY")
dat <- ffi_read_LGR915(f, date_format = "MDY", tz = "EST") # specify time zone
```

`ffi_read_LI7810` *Read a LI-7810 data file*

Description

Read a LI-7810 data file

Usage

```
ffi_read_LI7810(file)
```

Arguments

`file`	Filename to read, character

Details

Currently LI-7810 and LI-7820 files are handled identically.

Value

A `data.frame` with the parsed data.

Examples

```
f <- system.file("extdata/TG10-01087.data", package = "fluxfinder")
dat <- ffi_read_LI7810(f)
```

`ffi_read_LI7820` *Read a LI-7820 data file*

Description

Read a LI-7820 data file

Usage

```
ffi_read_LI7820(file)
```

Arguments

`file`	Filename to read, character

Details

Currently LI-7810 and LI-7820 files are handled identically.

Value

A data.frame with the parsed data.

Examples

```
f <- system.file("extdata/TG20-01182.data", package = "fluxfinder")
dat <- ffi_read_LI7820(f)
```

---

ffi_read_LI850 *Read a LI-850 data file*

Description

Read a LI-850 data file

Usage

```
ffi_read_LI850(file, tz = "UTC")
```

Arguments

file	Filename to read, character
tz	Time zone of the file's time data, character (optional)

Value

A data.frame with the parsed data, including a TIMESTAMP column.

Examples

```
f <- system.file("extdata/LI850.txt", package = "fluxfinder")
dat <- ffi_read_LI850(f)
dat <- ffi_read_LI850(f, tz = "EST") # specify time zone
```

`ffi_read_LIsmartchamber` *Read a LI-8200-01S (smart chamber) data file*

Description

Read a LI-8200-01S (smart chamber) data file

Usage

```
ffi_read_LIsmartchamber(file, concentrations = TRUE)
```

Arguments

`file`	Filename to read, character
`concentrations`	Return concentration data (the default), or just summary information? Logical

Value

A `data.frame` with the parsed data.

Note

These files are in JSON format. See also `https://www.licor.com/env/products/soil-flux/smart-chamber`.

Author(s)

Ben Bond-Lamberty

Examples

```
f <- system.file("extdata/LI8200-01S.json", package = "fluxfinder")
dat <- ffi_read_LIsmartchamber(f) # returns 240 rows
ffi_read_LIsmartchamber(f, concentrations = FALSE) # only 4 rows
```

ffi_read_PicarroG2301 *Read a Picarro G2301 data file*

Description

Read a Picarro G2301 data file

Usage

```
ffi_read_PicarroG2301(file, tz = "UTC")
```

Arguments

file	Filename to read, character
tz	Time zone of the file's time data, character (optional)

Value

A data.frame with the parsed data.

References

https://www.picarro.com/environmental/products/g2301_gas_concentration_analyzer

Examples

```
f <- system.file("extdata/PicarroG2301-data.dat", package = "fluxfinder")
dat <- ffi_read_PicarroG2301(f)
dat <- ffi_read_PicarroG2301(f, tz = "EST") # specify time zone
```

ideal-gas-law *Convert gas concentration to quantity using the Ideal Gas Law*

Description

Convert gas concentration to quantity using the Ideal Gas Law

Usage

```
ffi_ppm_to_umol(ppm, volume, temp, atm)

ffi_ppb_to_nmol(ppb, volume, temp, atm)
```

Arguments

`ppm`	Gas concentration (ppmv), numeric
`volume`	System volume (chamber + tubing + analyzer, m3), numeric
`temp`	Optional chamber temperature (degrees C), numeric
`atm`	Optional atmospheric pressure (Pa), numeric
`ppb`	Gas concentration (ppbv), numeric

Value

The quantity value, in micromoles (for `ffi_ppm_to_umol`) or nanomoles (for `ffi_ppb_to_nmol`).

Note

If `temp` and/or `atm` are not provided, the defaults are NIST normal temperature and pressure.

References

Steduto et al.: Automated closed-system canopy-chamber for continuous field-crop monitoring of CO2 and H2O fluxes, Agric. For. Meteorol., 111:171-186, 2002. doi:10.1016/S01681923(02)00023-0

Examples

```
ffi_ppm_to_umol(400, 0.1)
ffi_ppb_to_nmol(400, 0.1)
```

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