

# Package ‘pvldcurve’

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**Type** Package

**Title** Simplifies the Analysis of Pressure Volume and Leaf Drying Curves

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**VignetteBuilder** knitr

**Description** Simplifies the manufacturing, analysis and display of pressure volume and leaf drying curves. From the progression of the curves turgor loss point, osmotic potential, apoplastic fraction as well as minimum conductance and stomatal closure can be derived. Methods adapted from Bartlett, Scoffoni, Sack (2012) <[doi:10.1111/j.1461-0248.2012.01751.x](https://doi.org/10.1111/j.1461-0248.2012.01751.x)> and Sack, Scoffoni, PrometheusWikiContributors (2011) <<http://prometheuswiki.org/tiki-index.php?page=Minimum+epidermal+conductance+%28gmin%2C+a.k.a.+cuticular+conductance%29>>.

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ApplyCombMod	<i>Apply a combined exponential and linear model</i>
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## Description

a non linear model combining an exponential and a linear fit is applied to the data using the Gauss-Newton algorithm of nls. starting values are calculated based on the data. Weights are applied to the model based on the estimated insecurity of the data quality.

## Usage

```
ApplyCombMod(data, y = "y", x = "x")
```

## Arguments

data	data frame containig x and y data to which the model is ought to be applied to
y	name of column in data containing y data
x	name of column in data containing x data

**Value**

model parameters

---

ApplyCombMod2	<i>Apply a combined exponential and linear model</i>
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**Description**

a non linear model combining an exponential and a linear fit is applied to the data using nls. starting values are calculated based on the data. Constraints are applied to the model based on the known constraints of the aimed model.

**Usage**

```
ApplyCombMod2(data, y = "y", x = "x")
```

**Arguments**

data	data frame containg x and y data to which combined exponentail and linear model is ought to be applied to
y	name of column in data containing y data
x	name of column in data containing x data

**Value**

model parameters

---

Conductance	<i>Leaf Conductance</i>
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**Description**

Calculates mole-based or concentration-based conductance (stomatal or minimal conductance) ( $\text{mmol s}^{-1} \text{ m}^{-2}$  or  $\text{mm s}^{-1}$ ) of the double-sided leaf area by experimental weight loss data and weather data

**Usage**

```
Conductance(data, sample = "sample", fresh.weight = "fitted.fw",
  date.and.time = "date.and.time", leaf.area = "leaf.area",
  humidity = "humidity", temperature = "temperature",
  atmospheric.pressure = 101.35, driving.force = "mole")
```

**Arguments**

data	data frame, with columns of equal length containing at least columns with time (and date) of the fresh weight measurements, the measured fresh weights (g) and the single-sided leaf area (cm <sup>2</sup> ) of the sample as well as the average relative humidity (%) and temperature (degree Celsius) during the measurement intervals. The data is to be ordered chronologically by sample. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional name of the column in data containing the sample ID; default: "sample"
fresh.weight	optional name of the column in data containing the numeric fresh weight values (g); default: "fitted.fw" (fresh weight corrected by noises as outputted for leaf drying curves by the function FittedFW)
date.and.time	optional name of the column in data containing the time of the fresh weight measurements as class POSIXct; default: "date.and.time"
leaf.area	optional name of the column in data containing the numeric single-sided leaf area values (cm <sup>2</sup> ); default: "leaf.area"
humidity	optional name of the column in data containing the numeric humidity values (%); default: "humidity"
temperature	optional name of the column in data containing the numeric temperature values (degree Celsius); default: "temperature"
atmospheric.pressure	optional, giving the numeric atmospheric pressure in kPa, default = 101.325 (atmospheric pressure at sea level)
driving.force	optional; possible values: mole or conc; defines whether conductance is expressed on the basis of a mole fraction-based (default) or a concentration-based driving force

**Details**

Calculates mole-based conductance (mmol s<sup>-1</sup> m<sup>-2</sup>) as:

$$g = T/VPD$$

whereas T = transpiration (mmol s<sup>-1</sup> m<sup>-2</sup>) is calculated as:

$$T = \Delta FM * 1000 * (\Delta t * 60 * LA * 2 / 10000 * 18.01528)^{-1}$$

whereas  $\Delta FM$  = fresh matter reduction (g),  $\Delta t$  = time interval (min), LA = single-sided leaf area (cm<sup>2</sup>) and VPD = vapor pressure deficit (mol \* mol<sup>-1</sup>) is calculated as:

$$VPD = (1 - RH/100) * (VP_{sat}/AP)$$

whereas RH = relative humidity (%), VP<sub>sat</sub> = saturation vapor pressure (kPa), AP = atmospheric pressure (kPa), whereas:

$$VP_{sat} = 0.61121 * \exp((18.678 - T/234.5)(T * 257.14 + T))$$

where  $T$  = air temperature (degree Celsius)

Concentration based conductance ( $\text{mm s}^{-1}$ ) is derived from mole-based conductance  $g(\text{mol})$  as:

$$g(\text{conc}) = g(\text{mol}) * R * (T + 273.15) / AP / 1000$$

whereas:  $R$  = gas constant ( $8.3144598 \text{ J (mol} * \text{K)}^{-1}$ ) and  $T$  = absolute temperature (degree Celsius)

### Value

The original data frame extended by a numeric column with the mole-based or the concentration-based conductance ( $\text{mmol s}^{-1} \text{ m}^{-2}$ ,  $\text{mm s}^{-1}$ ) of the double-sided leaf area (conductance). The first value of each sample is NA since conductance values are computed from row  $i$  and  $i-1$

### Examples

```
# get example data
weight_loss_data <- leaf_drying_data
weather_data <- weather_data
df <- WeatherAllocation(weight_loss_data, weather_data) # allocate weather to weight loss data

# extend the data frame by mole-based conductance values
df_with_conductance <- Conductance(df, fresh.weight = "fresh.weight")

# extend the data frame by concentration-based conductance values
df_with_conductance <- Conductance(df, fresh.weight = "fresh.weight", driving.force = "conc")

# calculate with atmospheric pressure of 99.8 kPa
df_with_conductance <- Conductance(df, fresh.weight = "fresh.weight", atmospheric.pressure = 99.8)
```

---

ExtractFitParam

*Extracts the fitting parameters from results list*

---

### Description

Extracts the coefficients and confidence intervals from the fitting results of the functions analysing the pressure volume curve (TurgorLossPoint, OsmoticPot and ModElasticity) or the functions analysing the leaf drying curve (StomatalClosure and Gmin)

### Usage

```
ExtractFitParam(result_list)
```

### Arguments

**result\_list**      output list from the functions TurgorLossPoint, OsmoticPot, ModElasticity, StomatalClosure or Gmin

**Value**

data frame containing the coefficients and the 0.95 confidence interval of the coefficients from the fit

---

ExtractParam

*Extracts parameters from result list*

---

**Description**

Extracts the curve parameters from the result lists of the functions analysing the pressure volume curve (TurgorLossPoint, OsmoticPot and ModElasticity) or the functions analysing the leaf drying curve (StomatalClosure and Gmin)

**Usage**

```
ExtractParam(result_list)
```

**Arguments**

`result_list`      output list from the functions TurgorLossPoint, OsmoticPot, ModElasticity, StomatalClosure or Gmin

**Value**

data frame containing the results from the curve analysis only, depending on the function used, relative water deficit at turgor loss point (rwd.tlp), water potential at turgor loss point (water.pot.tlp), apoplastic fraction (apo.fract), osmotic potential at full saturation (osmotic.pot.full.sat), modulus of elasticity (modulus.elasticity), time since start of desiccation of stomatal closure (stom.clos.time), relative water deficit at stomatal closure (stom.clos.rwd), minimum conductance determined by the intercept of the leaf drying curve fit with the point of stomatal closure (stom.clos.gmin), by taking the mean from all minimum conductance values (mean.gmin) or minimum conductance determined by extrapolation of minimum conductance curve linearly to full saturation (gmin.full.sat)

**Examples**

```
# use pressure volume data provided by package
pv_data <- pressure_volume_data

# do pressure volume curve analysis
pv_data <- RelativeWaterDeficit(pv_data)
results <- OsmoticPot(pv_data, graph = FALSE)

# extract curve values
ExtractParam(results)
```

---

FitLeafArea	<i>Leaf area fitting</i>
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---

## Description

Fits randomly measured leaf area values linearly to fresh weight values. Useful if the leaf area changes during a measurement series but is only randomly measured.

## Usage

```
FitLeafArea(data, sample = "sample", fresh.weight = "fresh.weight",  
            leaf.area = "leaf.area")
```

## Arguments

data	data frame, with columns of equal length, containing at least columns with the fresh.weight (g) and the leaf.area (cm <sup>2</sup> ) values, ordered by sample by descending fresh weight. A column containing the sample IDs is optionally required if several samples were measured. At least 3 leaf area values are required.
sample	string, optional name of the column in data containing the sample ID, default: "sample"
fresh.weight	optional name of the column in data containing the numeric fresh weight values (g); default: "fresh.weight"
leaf.area	optional name of the column in data containing the numeric single-sided leaf area values (cm <sup>2</sup> ); default: "leaf.area"

## Details

fits given leaf area values linearly to the respective fresh weight values and calculates leaf area values for the fresh weight values based on the fit

## Value

the original data frame extended by a numeric column containing the fitted leaf area values (leaf.area.fitted)

## Examples

```
# get example data  
df <- data.frame(  
  sample = c(as.integer(rep(1, times = 6))),  
  fresh.weight = c(1.23, 1.19, 1.15, 1.12, 1.09, 1.0),  
  leaf.area = c(10.5, NA, NA, 9.8, NA, 8.4))  
# fit leaf area  
df_new <- FitLeafArea(df)
```

FittedFW

*Correct continuous fresh weight measurements***Description**

Corrects fresh weight, measured continuously on a desiccating leaf for determination of minimum conductance, by fitting the fresh weight values to a combined exponential and linear model

**Usage**

```
FittedFW(data, sample = "sample", fresh.weight = "fresh.weight",
  time.since.start = "time.since.start", graph = TRUE,
  show.legend = TRUE)
```

**Arguments**

data	data frame, at least with a column containing numeric fresh weight (g) and time since start (min) values, ordered by sample by descending fresh.weight. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional name of the column in data containing the sample IDs, default: "sample"
fresh.weight	optional name of the column in data containing the fresh weight values (g), default: "fresh weight"
time.since.start	optional name of the column in data containing the time since start (min) values, default: "time.since.start"
graph	set FALSE if no plots are to be returned
show.legend	set FALSE if no legend is to be shown in the plots

**Details**

Determination of minimum conductance via a leaf drying curve requires fresh weight to be measured continuously on a desiccating leaf. The fresh weight values are then used to calculate leaf conductance. If several leaves are measured on one scale, the measurements are prone to noises, which influence conductance values largely. Here, the fresh weight data is corrected for noises by fitting it to a combined linear and exponential model using the port algorithm of nls().

Before using this function, check the raw data for an initial plateau. If the exponential decline does not onset directly, fitting might not succeed.

**Value**

the original data frame (data) extended by a numeric column containing the fitted fresh weight values ("fitted.fw")



## Examples

```
# get example data
df <- TimeSinceStart(leaf_drying_data)
# remove plateauing data
df <- df[df$fw.plateau != "yes",]
# extend the data frame by saturated fresh weight
df <- FittedFW(df)
```

---

FWSaturated	<i>Saturated fresh weight estimation</i>
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---

## Description

Calculates saturated fresh weight by fitting fresh weight values above the turgor loss point linearly to water potential values.

## Usage

```
FWSaturated(data, sample = "sample",
  water.potential = "water.potential", fresh.weight = "fresh.weight",
  dry.weight = "dry.weight")
```

## Arguments

data	data frame, at least with a column containing numeric water potential (bar), fresh.weight (g) and dry.weight (g) values, ordered by sample by descending water potential. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional name of the column in data containing the sample IDs, default: "sample"
water.potential	optional name of the column in data containing the water potential values (bar), default = "water.potential"
fresh.weight	optional name of the column in data containing the fresh weight values (g), default: "fresh weight"
dry.weight	optional name of the column containing the dry weight values (g), default: "dry.weight"

## Details

Above the turgor loss point, a linear relationship between water content and water potential exists. Based on this premise, saturated water content is found where water potential is zero. First, turgor loss point is calculated based on the relative leaf water loss (fresh weight minus dry weight relativized by the maximum leaf water content value). Then, data above the turgor loss point is extracted and fresh weight is fitted linearly to water potential. The point where water potential of the linear regression line is zero is the saturated water content.

**Value**

the original data frame (data) extended by a numeric column containing the saturated fresh weight values ("fresh.weight.saturated")

**Examples**

```
# get example data
df <- pressure_volume_data
# extend the data frame by saturated fresh weight
df <- FWSaturated(df)
```

---

Gmin

---

*Minimum leaf conductance*


---

**Description**

Determines mean minimum leaf conductance and minimum leaf conductance at full turgidity and stomatal closure in experimentally obtained water loss curves.

**Usage**

```
Gmin(data, sample = "sample", time.since.start = "time.since.start",
      conductance = "conductance", RWD = "RWD.interval",
      stom.clos.threshold = FALSE, graph = TRUE, show.legend = TRUE)
```

**Arguments**

data	data frame containing columns of equal lengths giving the coordinates of the curve: time since start (minutes), conductance ( $\text{mmol m}^{-2} \text{s}^{-1}$ ) and RWD (%), ordered by sample by ascending time since start. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional column name in data containing the sample ID, default: sample
time.since.start	optional column name in data containing the numerical values for time since start of the experiment (min), default: time.since.start
conductance	optional column name in data containing the numerical conductance values ( $\text{mmol m}^{-2} \text{s}^{-1}$ ), default: "conductance"
RWD	optional column name in data containing the numerical relative water deficit values (%), default: "RWD.interval" (RWD average of an interval, as outputted by RWDInterval)
stom.clos.threshold	threshold value for stomatal closure. Automatic determination by default.
graph	set FALSE if no plots are to be returned
show.legend	set FALSE if no legend is to be shown in the plots

## Details

The coordinates of stomatal closure are determined via the function `StomatalClosure()`. Conductance data including and following stomatal closure are then extracted and the average is taken (`mean.gmin`). A linear regression is applied to the data and the y axis intercept (`gmin.full.sat`) and the coordinate at the RWD point of stomatal closure (`lin.gmin`) are calculated from the function.

Before using this function, check the raw data for an initial plateau. If the exponential decline does not onset directly, fitting might not succeed.

## Value

List splitted by sample consisting of

<code>gmin</code>	mean minimum conductance ( <code>mean.gmin</code> ) ( $\text{mmol m}^{-2} \text{s}^{-1}$ ) after stomatal closure of the measurement interval, minimum conductance at full saturation ( <code>gmin.full.sat</code> ) ( $\text{mmol m}^{-2} \text{s}^{-1}$ ) and minimum conductance at stomatal closure based on the linear fit ( <code>lin.gmin</code> ) ( $\text{mmol m}^{-2} \text{s}^{-1}$ )
<code>formula</code>	formula of the linear regression of <code>gmin</code> vs. RWD
<code>coef</code>	coefficients of linear fit
<code>conf_int</code>	upper (97.5 %) and lower (2.5 %) border of 95 % confidence interval of model parameters

If `graph = TRUE`, the plotted original data is displayed with the x-axis intercept of the point of stomatal closure and the linear regression line of `gmin` showing the point of y-intercept (`gminfullsat`).

## Examples

```
# get example data
df <- WeatherAllocation(leaf_drying_data, weather_data) # allocate weather to weight loss data
df <- TimeSinceStart(df) # calculate time since start
df <- df[df$fw.plateau != "yes",] # remove plateauing data
df <- FittedFW(df, graph = FALSE) # correct noises in fresh weight
df <- RWDInterval(df, fresh.weight = "fitted.fw") # calculate RWD based in the intervals
df <- Conductance(df, fresh.weight = "fitted.fw") # calculate conductance

# calculate gmin and plot graphs
gmin <- Gmin(df)
```

---

leaf\_drying\_data

*Experimentally determined leaf drying data*


---

## Description

A dataset containing repeatedly measured fresh weights of transpiring leaves subjected to different soil moisture conditions during their growth ( $n = 6$ ) and their saturated fresh weight and dry weight. together with their saturated fresh weight, their fresh weight before saturation, their dry weight and leaf area.

**Usage**

```
leaf_drying_data
```

**Format**

A data frame with 172 rows and 9 variables:

**Details**

- treatment: Soil moisture conditions during the last 6 days of Kohlrabi growth (10-30)
- sample: Sample ID (1 - 12)
- dry.weight: Dry weight of the sample in gramm (0.3922 - 0.6692)
- fresh.weight.harvest: Fresh weight directly after cutting before over night saturation of the sample in gramm (2.8062 - 7.1009)
- fresh.weight.saturated: Saturated fresh weight of the sample in gramm (4.2186 - 7.3179)
- leaf.area: leaf area of the sample in cm<sup>2</sup> (91 - 148)
- date.and.time: Time (and date) of the fresh weight measurement of the transpiring sample (2019-03-26 09:48:00 - 2019-03-27 09:51:00)
- fresh.weight: Fresh weight of the transpiring sample in gramm (3.5112 - 7.3179)
- fw.plateau: Indicates plateauing fresh.weight values before the onset of the exponential decline. Remove values for analysis where fw.plateau = yes

---

MergeDf

---

*Merges data to data frames*


---

**Description**

merges data frames containing all necessary informations for plotting with PlotOutput()

**Usage**

```
MergeDf(x, y, y2 = FALSE, y3 = FALSE, legend, legend.y2 = FALSE,
        legend.y3 = FALSE)
```

**Arguments**

x	vector containing the x values
y	vector containing the y values
y2	optional vector containing the values for the second y coordinates
y3	optional vector containing the values for the third y coordinates
legend	name of the y values in the legend
legend.y2	optional name of the second y values in the legend
legend.y3	optional name of the third y values in the legend

**Value**

data frame with columns containing all above information in equalized length as requested by ggplot

---

ModElasticity	<i>Modulus of elasticity</i>
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---

**Description**

Determines pressure potential and the modulus of elasticity

**Usage**

```
ModElasticity(data, sample = "sample",
  water.potential = "water.potential", RWD = "RWD", graph = TRUE,
  show.legend = TRUE)
```

**Arguments**

data	data frame containing columns of equal lengths giving the numerical coordinates of the curve: water potential (bar) and RWD (%), ordered by sample by descending water potential. A column containing the sample IDs is optionally required if several samples were measured
sample	optional column name in data containing the sample ID, default: "sample"
water.potential	optional column name in data containing the water potential values of the leaf (bar), default: "water.potential"
RWD	optional column name in data containing the relative water deficit values (%), default: "RWD"
graph	set FALSE if no plots are to be returned
show.legend	set FALSE if no legend is to be shown in the plots

**Details**

Relative water deficit at turgor loss point is determined via the function `TurgorLossPoint()` and osmotic potential is calculated via the function `OsmoticPot()`.

Pressure potential is derived by subtracting osmotic potential from water potential. The part of the pressure potential prior the turgor loss point is then fitted linearly and the modulus of elasticity (`M.Elasticity`) equals the slope of the fitted line.

Before using this function, check the raw data for an initial plateau. If the exponential decline does not onset directly, fitting might not succeed.

**Value**

List splitted by sample consisting of

<code>modulus.elasticity</code>	modulus of elasticity (bar)
<code>formula</code>	formula of the transformed linear osmotic potential fit (1/-bar) and the pressure potential (bar) fit
<code>coef</code>	coefficients of the osmotic (1/-bar) and pressure potential (bar) fit
<code>conf_int</code>	upper (97.5 %) and lower (2.5 %) border of 95 % confidence interval of model parameters

If `graph = TRUE`, the original data is displayed with the x- and y-axis intercepts of the turgor loss point, the osmotic potential fit and the linear regression line of the pressure potential.

**Examples**

```
#get example data, calculate Relative Water Deficit
data <- RelativeWaterDeficit(pressure_volume_data)[pressure_volume_data$sample == 1, ]

# determine modulus of elasticity and the fitting parameters. Do not plot results.
m_elasticity <- ModElasticity(data, graph = FALSE)
```

---

OrderCheck

*Order Check*


---

**Description**

Checks for the correct ordering of the data: increasing for `date.and.time` and `time.since.start`, decreasing for `fresh.weight` and `water.potential`. Done separatly for each sample. An individualized warning is printed if not ordered correctly.

**Usage**

```
OrderCheck(data, sample = FALSE, date.and.time = FALSE,
  fresh.weight = FALSE, water.potential = FALSE,
  time.since.start = FALSE)
```

**Arguments**

<code>data</code>	data frame containing the data to be checked
<code>sample</code>	name of the column containing the sample IDs, if present in data
<code>date.and.time</code>	name of the column containing the date and time (POSIXct) of the measurements, if present in data
<code>fresh.weight</code>	name of the column containing the numeric fresh weight values, if present in data

<code>water.potential</code>	name of the column containing the numeric water potential values, if present in data
<code>time.since.start</code>	name of the column containing the numeric time since start values, if present in data

**Value**

no return value

---

<i>OsmoticPot</i>	<i>Pressure Volume Curve Analysis</i>
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---

**Description**

Determines the coordinates of the turgor loss point, osmotic potential at full hydration and apoplastic fraction

**Usage**

```
OsmoticPot(data, sample = "sample",  
  water.potential = "water.potential", RWD = "RWD", graph = TRUE,  
  show.legend = TRUE)
```

**Arguments**

<code>data</code>	data frame containing columns of equal lengths giving the numerical coordinates of the curve: water potential (bar) and RWD (%), ordered by sample by descending water potential. A column containing the sample IDs is optionally required if several samples were measured.
<code>sample</code>	optional column name in data containing the sample ID, default: "sample"
<code>water.potential</code>	optional column name in data containing the numeric water potential values (bar), default: "water.potential"
<code>RWD</code>	optional column name in data containing the relative water deficit values (%), default: "RWD"
<code>graph</code>	set FALSE if no plots are to be returned
<code>show.legend</code>	set FALSE if no legend is to be shown in the plots

Details

RWD at turgor loss point is derived by the function TurgorLossPoint().

The pressure-volume curve data is converted to -1/bar. The osmotic potential is then derived by fitting a linear regression line with the Gauss-Newton algorithm of nls() to the water potential data following the turgor loss point. The y- and x-axis intercept of the regression line gives the osmotic potential at full hydration (op.full.sat) and the apoplastic fraction (apo.fract), respectively. Water potential at turgor loss point equals the value of the osmotic potential fit at the relative water deficit at turgor loss point.

Value

List splitted by sample consisting of

turgor.loss.point	x and y coordinates of the turgor loss point (RWD (%) and water.potential (bar), respectively)
osmotic.potential	x and y intercepts of the osmotic potential fit (apoplastic fraction (apo.fract) (%) and op.full.sat (bar), respectively)
formula	formula of the linear osmotic potential fit
coef	coefficients of the linear model
conf_int	upper (97.5 %) and lower (2.5 %) border of 95 % confidence interval of model parameters

If graph = TRUE, the plotted tranformed data is displayed with the x- and y-axis intercepts of the turgor loss point and the linear regression line of the osmotic potential showing the point of y-intercept (op.full.sat) and x-intercept (apo.fract).  
Before using this function, check the raw data for an initial plateau. If the exponential decline does not onset directly, fitting might not succeed.

Examples

```
# get example data, calculate Relative Water Deficit
data <- RelativeWaterDeficit(pressure_volume_data)

# calculate pressure volume curve characteristics and plot graphs
pv_analysis <- OsmoticPot(data)
```

---

PlotOutput	<i>Plot Output</i>
------------	--------------------

---

Description

plots the data as specified



**Usage**

```
PlotOutput(sub.sample, x, y, y2 = FALSE, y3 = FALSE, legend.y,
  legend.y2 = FALSE, legend.y3 = FALSE, x.axis, y.axis,
  x.intercept = FALSE, y.intercept = FALSE,
  legend.x.intercept = FALSE, line.x, line.y, line.y2 = FALSE,
  line.y3 = FALSE, legend.line.y, legend.line.y2 = FALSE,
  legend.line.y3 = FALSE, show.legend = show.legend)
```

**Arguments**

sub.sample	sample ID
x	vector containing the x coordinates of the data to be plotted as points
y	vector containing the y coordinates of the data to be plotted as points
y2	optional vector containing the second y coordinates of the data to be plotted as points
y3	optional vector containing the third y coordinates of the data to be plotted as points
legend.y	string, name of data points to be printed in the legend
legend.y2	string, optional name of second set of data points to be printed in the legend
legend.y3	string, optional name of third set of data points to be printed in the legend
x.axis	string, label of x axis
y.axis	string, label of y axis
x.intercept	vector containing the x coordinate of the intercept
y.intercept	optional vector containing the y coordinate of the intercept
legend.x.intercept	string, name of x.intercept to be printed in the legend
line.x	vector containing the x coordinate for the lines
line.y	vector containing the y coordinates for the line
line.y2	vector containing the y coordinates for the second line
line.y3	vector containing the y coordinates for the second line
legend.line.y	string, name of line to be printed in the legend
legend.line.y2	string, name of second line to be printed in the legend
legend.line.y3	string, name of third line to be printed in the legend
show.legend	boolean, specifies whether a legend is to be printed

**Value**

graphic

---

pressure\_volume\_data    *Pressure volume curve data*

---

### Description

A dataset containing water potential and fresh weight measurements of repeatedly measured drying kohlrabi leaves subjected to different soil moisture conditions during their growth (n = 6) and their saturated fresh weight and dry weight.

### Usage

```
pressure_volume_data
```

### Format

A data frame with 160 rows and 8 variables

### Details

- date: Date of measurement
- treatment: Soil moisture conditions during the last 6 days of Kohlrabi growth (10-30)
- sample: Sample ID (1 - 12)
- fresh.weight.harvest: Fresh weight measured at harvest (12 h prior measurement of fresh.weight.saturated) (2.9813 - 7.1557)
- fresh.weight.saturated: Saturated fresh weight of the leaf in grams (4.1276 - 7.0867)
- fresh.weight: Fresh weight of the leaf in grams (2.7215 - 6.8246)
- dry.weight: Dry weight of the leaf in grams (0.2937 - 0.7267)
- water.potential: Water potential of the leaf in bar (-16.2 - -2.4)

---

RelativeWaterContent    *Relative Water Content (RWC)*

---

### Description

Calculates relative water content (RWC, %)

### Usage

```
RelativeWaterContent(data, fresh.weight = "fresh.weight",
  dry.weight = "dry.weight",
  fresh.weight.saturated = "fresh.weight.saturated")
```

**Arguments**

<code>data</code>	data frame with columns of equal length containing at least columns with the fresh weight (g), the dry weight (g) and the saturated fresh weight (g)
<code>fresh.weight</code>	optional name of the column in data containing the numeric fresh weight values (g); default: "fresh.weight"
<code>dry.weight</code>	optional name of the column in data containing the numeric dry weight values (g); default: "dry.weight"
<code>fresh.weight.saturated</code>	optional name of the column in data containing the numeric saturated fresh weight values (g); default: "fresh.weight.saturated"

**Details**

Relative water content (%) is calculated as:

$$RWC = 100 * ((FW - DW) / (FW_s - DW))^{-1}$$

whereas FW = fresh weight, DW = dry weight and FW<sub>s</sub> = fresh weight at water saturation.

**Value**

the original data frame extended by a numeric column with the relative water content (RWC) (%).

**Examples**

```
# get example data
df <- leaf_drying_data

# extend df by RWC
df_with_RWC <- RelativeWaterContent(df)
```

---

RelativeWaterDeficit    *Relative Water Deficit (RWD)*

---

**Description**

Calculates relative water deficit (%)

**Usage**

```
RelativeWaterDeficit(data, fresh.weight = "fresh.weight",
  dry.weight = "dry.weight",
  fresh.weight.saturated = "fresh.weight.saturated")
```

**Arguments**

<code>data</code>	data frame with columns of equal length containing at least columns with the fresh weight (g), the dry weight (g) and the saturated fresh weight (g)
<code>fresh.weight</code>	optional name of the column in data containing the numeric fresh weight values (g); default: <code>fresh.weight</code>
<code>dry.weight</code>	optional name of the column in data containing the numeric dry weight values (g); default: <code>dry.weight</code>
<code>fresh.weight.saturated</code>	optional name of the column in data containing the numeric saturated fresh weight values (g); default: <code>fresh.weight.saturated</code>

**Details**

Relative water deficit (%) is calculated as:

$$RWD = 100 - 100 * ((FW - DW)(FW_s - DW)^{-1})$$

whereas FW = fresh weight, DW = dry weight and FW<sub>s</sub> = fresh weight at water saturation.

**Value**

the original data frame extended by a numeric column with the relative water deficit (RWD) (%).

**Examples**

```
# get example data
df <- leaf_drying_data

# extend df by RWD
df_with_RWD <- RelativeWaterDeficit(df)
```

---

RWDInterval

*Mean relative water deficit (RWD) of an interval*


---

**Description**

Calculates relative water deficit (%) as mean value of a measurement interval

**Usage**

```
RWDInterval(data, sample = "sample", fresh.weight = "fitted.fw",
  dry.weight = "dry.weight",
  fresh.weight.saturated = "fresh.weight.saturated")
```

**Arguments**

data	data frame with columns of equal length containing at least columns with the fresh weight (g), the dry weight (g) and the saturated fresh weight (g), ordered by sample by descending by fresh weight. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional name of the column in data containing the sample IDs, default: "sample"
fresh.weight	optional name of the column in data containing the numeric fresh weight values (g); default: "fitted.fw"
dry.weight	optional name of the column in data containing the numeric dry weight values (g); default: "dry.weight"
fresh.weight.saturated	optional name of the column in data containing the numeric saturated fresh weight values (g); default: "fresh.weight.saturated"

**Details**

First, the mean fresh weight is calculated for each measurement interval. Relative water deficit (%) is then calculated as:

$$RWD = 100 - 100 * ((mFW - DW)(FWs - DW)^{-1})$$

whereas mFW = mean fresh weight, DW = dry weight and FWs = fresh weight at water saturation.

**Value**

the original data frame extended by a numeric column with the mean relative water deficit for the measurement interval (RWD.interval) (%).

**Examples**

```
# get example data
df <- leaf_drying_data

# extend df by RWD
df_with_RWD <- RWDInterval(df, fresh.weight = "fresh.weight")
```

---

SaturationVaporPressure

*Saturation vapor pressure (VPsat)*


---

**Description**

Calculates saturation vapor pressure (kPa) using the Arden Buck equation

Usage

```
SaturationVaporPressure(data, temperature = "temperature")
```

Arguments

data	data frame with at least a numeric column containing temperature (degree Celsius)
temperature	optional name of the column in data containing the temperature values; default: "temperature"

Details

Calculates saturation vapor pressure (kPa) over liquid by temperature using the Arden Buck equation:

$$VP_{sat} = 0.61121 \exp((18.678 - T/234.5)(T/257.14 + T))$$

where T = air temperature (degree Celsius)

Value

the original data frame extended by a numeric column with the saturation vapor pressure (kPa).

Examples

```
# generate example data frame
df <- data.frame(temperature = c(23.1, 23.2))

# extend df by saturation vapor pressure
df_with_VPsat <- SaturationVaporPressure(df)
```

---

StomatalClosure	<i>Point of stomatal closure</i>
-----------------	----------------------------------

---

Description

Determines the point of stomatal closure in a set of experimentally obtained leaf drying curves. Stomatal closure happens when the curve irreversibly settles to linear water loss.

Usage

```
StomatalClosure(data, sample = "sample",
  time.since.start = "time.since.start", conductance = "conductance",
  RWD = "RWD.interval", threshold = FALSE, graph = TRUE,
  show.legend = TRUE)
```

**Arguments**

<code>data</code>	data frame containing columns of equal length giving the numerical coordinates of the curve: time since start (minutes), conductance ( $\text{mmol m}^{-2} \text{ s}^{-1}$ ) and RWD (%), ordered by sample by ascending time since start. A column containing the sample IDs is optionally required if several samples were measured.
<code>sample</code>	optional name of the column in data containing the sample ID (if available), default: "sample"
<code>time.since.start</code>	optional name of the column in data containing numeric time since start values (min), default: "time.since.start"
<code>conductance</code>	optional name of the column in data containing numeric leaf conductance values ( $\text{mmol m}^{-2} \text{ s}^{-1}$ ), default: "conductance"
<code>RWD</code>	optional name of the column in data containing numeric relative water deficit (%) values, default: "RWD.interval" (RWD average of an interval as outputted by RWDInterval)
<code>threshold</code>	sensitivity for the determination of stomatal closure. 60 by default.
<code>graph</code>	set FALSE if no plots are to be returned
<code>show.legend</code>	set FALSE if no legend is to be shown in the plots

**Details**

Before using this function, check the raw data for an initial plateau. If the exponential decline does not onset directly, fitting might not succeed.

The conductances by time since start curves are fitted using the Gauss-Newton algorithm of `nls()` to a combined exponential and linear model. The exponential and linear parts are extracted and time since start at stomatal closure is localized at the point where the slope of the exponential part of the fit is higher than a threshold value. The threshold value is calculated by the use of the parameter `b` of the exponential part of the fit ( $a * \exp(b * x)$ ):  $-(b^2 * \text{sens})$ . The sensitivity constant (`sens`) is 60 by default and can be specified individually by the argument 'threshold'.

Minimum conductance (`gmin`) at stomatal closure is the conductance value of the overall fit at stomatal closure. RWD at stomatal closure is then calculated by linear regression of RWD and time since start.

**Value**

List splitted by sample consisting of

<code>stomatal.closure</code>	coordinates of the point of stomatal closure (time.since.start, RWD, conductance)
<code>formula</code>	formula of the exponential and linear part of the combined fits
<code>coef</code>	coefficients of combined model
<code>conf_int</code>	upper (97.5 %) and lower (2.5 %) border of 95 % confidence interval of model parameters

If `graph = TRUE`, the plotted original data is displayed with the exponential and linear fit of the combined model as well as the x-coordinate (time.since.start) of the point of stomatal closure.

## Examples

```
# get example data
df <- WeatherAllocation(leaf_drying_data, weather_data) # allocate weather to weight loss data
df <- TimeSinceStart(df) # calculate time since start
df <- df[df$fw.plateau != "yes",] # remove plateauing data
df <- FittedFW(df, graph = FALSE) # correct noises in fresh weight
df <- RWDInterval(df, fresh.weight = "fitted.fw") # calculate RWD based in the intervals
df <- Conductance(df, fresh.weight = "fitted.fw") # calculate conductance

# identify stomatal closure in curve and get graphs
sc <- StomatalClosure(df)
```

---

TimeInterval	<i>Time Interval</i>
--------------	----------------------

---

## Description

Calculates time intervals (min) between temporally repeated measurements

## Usage

```
TimeInterval(data, sample = "sample", date.and.time = "date.and.time")
```

## Arguments

data	data frame containing at least a column giving the time (and date) (class POSIXct) of the measurements ordered by sample and chronologically. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional name of the column in data containing the sample ID, default: "sample"
date.and.time	optional name of the column in data containing the time (and date) as class POSIXct, default: "date.and.time"

## Value

the original data frame extended by a numerical vector containing the time intervals (min) between the measurements of a sample. The first values of each sample is NA since time intervals are computed from row  $i$  and  $i-1$ .

## Examples

```
# get example data
df <- leaf_drying_data

# extend df by time interval
df_with_ti <- TimeInterval(df)
```



---

<code>TimeSinceStart</code>	<i>Time Since Start</i>
-----------------------------	-------------------------

---

**Description**

Calculates time since start (min) of measurement for temporally repeated measurements

**Usage**

```
TimeSinceStart(data, sample = "sample",  
  date.and.time = "date.and.time")
```

**Arguments**

<code>data</code>	data frame containing at least a column giving the the time (and date) of the measurements ordered by sample and chronologically. A column containing the sample IDs is optionally required if several samples were measured.
<code>sample</code>	optional name of the column in data containing the sample IDs, default: "sample".
<code>date.and.time</code>	optional name of the column in data containing the time (and date) as class POSIXct, default: "date.and.time"

**Value**

The original data frame extended by a numerical vector containing time since start (min) of the measurements.

**Examples**

```
# get example data frame  
df <- leaf_drying_data  
  
# extend df by time since start  
df_with_tss <- TimeSinceStart(df)
```

---

<code>Transpiration</code>	<i>Leaf transpiration</i>
----------------------------	---------------------------

---

**Description**

Calculates transpiration of a plant part from experimentally determined weight loss per time unit and double-sided leaf area.

**Usage**

```
Transpiration(data, sample = "sample", date.and.time = "date.and.time",
  fresh.weight = "fitted.fw", leaf.area = "leaf.area",
  output.unit = "mmol")
```

**Arguments**

<code>data</code>	data frame with columns of equal length containing at least columns with the time (and date) of the fresh weight measurements as well as columns with the measured fresh weights (g) and the single-sided leaf area (cm <sup>2</sup> ) of the sample. The data is to be ordered chronologically by sample. A column containing the sample IDs is optionally required if several samples were measured.
<code>sample</code>	optional name of the column in data containing the sample ID; default: "sample"
<code>date.and.time</code>	optional name of the column in data containing the time of the fresh weight measurements as class POSIXct; default: "date.and.time"
<code>fresh.weight</code>	optional name of the column in data containing the numeric fresh weight values (g); default: "fitted.fw"
<code>leaf.area</code>	optional name of the column in data containing the numeric single-sided leaf area values (cm <sup>2</sup> ); default: "leaf.area"
<code>output.unit</code>	optional; possible values: "mg" or "mmol"; defines whether transpiration is given in mmol m <sup>-2</sup> s <sup>-1</sup> (Default) or in mg m <sup>-2</sup> s <sup>-1</sup>

**Details**

Transpiration (mmol s<sup>-1</sup> m<sup>-2</sup>) is calculated as:

$$T = \Delta FM * 1000 * (\Delta t * 60 * LA * 2 / 10000 * 18.01528)^{-1}$$

whereas T = transpiration,  $\Delta FW$  = reduction of fresh weight (g),  $\Delta t$  = time unit (min), LA = single-sided leaf area (cm<sup>2</sup>)

**Value**

The original data frame extended by a numeric column with the transpiration (mg s<sup>-1</sup> m<sup>-2</sup> or mmol s<sup>-1</sup> m<sup>-2</sup>) of the double-sided leaf area. The first value of each sample is NA, since transpiration values are computed from row *i* and *i*-1.

**Examples**

```
# get example data and allocate
df <- WeatherAllocation(leaf_drying_data, weather_data)

# extend df by transpiration in mmol s^-1 m^-2
df_with_transpiration <- Transpiration(df, fresh.weight = "fresh.weight")

# extend df by transpiration in mg s^-1 m^-2
df_with_transpiration <- Transpiration(df, fresh.weight = "fresh.weight", output.unit = "mg")
```

---

TurgorLossPoint	<i>Turgor Loss Point</i>
-----------------	--------------------------

---

### Description

Determines the x coordinate (RWD) of the turgor loss point in a set of experimentally obtained pressure volume curves.

### Usage

```
TurgorLossPoint(data, sample = "sample",
  water.potential = "water.potential", RWD = "RWD", graph = TRUE,
  show.legend = TRUE)
```

### Arguments

data	data frame containing columns of equal lengths giving at least the numerical coordinates of the curve: water potential (bar) and RWD (%), ordered by sample by descending water potential. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional name of the column in data containing the sample ID, default: "sample"
water.potential	optional name of the column in data containing the numeric water potential values (bar), default: "water.potential"
RWD	optional name of the column in data containing numeric relative water deficit values (%), default: "RWD"
graph	set FALSE if no plots are to be returned
show.legend	set FALSE if no legend is to be shown in the plots

### Details

Before using this function, check the raw data for an initial plateau. If the exponential decline does not onset directly, fitting might not succeed.

The data is fitted using the Gauss-Newton algorithm of `nls()` to a combined exponential and linear model. The exponential and linear parts are extracted and RWD at turgor loss point is localized at their point of minimum distance.

### Value

List splitted by sample consisting of

turgor.loss.point	coordinates of the turgor loss point (RWD)
formula	formula of the exponential and linear part of the combined fits
coef	coefficients of combined model

conf\_int            upper (97.5 %) and lower (2.5 %) border of 95 % confidence interval of model parameters

If graph = TRUE, the plotted original data is displayed with the exponential and linear fit of the combined model as well as the x-coordinate (RWD) of the turgor loss point.

## Examples

```
# get sample data
data <- RelativeWaterDeficit(pressure_volume_data)[pressure_volume_data$sample == 1, ]

# identify turgor loss point in curve
turgor_loss_point <- TurgorLossPoint(data)
```

---

ValidityCheck	<i>Ensures the validity of the input data</i>
---------------	---

---

## Description

Ensures the validity of the input data

## Usage

```
ValidityCheck(data, sample = FALSE, leaf.area = FALSE,
  date.and.time = FALSE, dry.weight = FALSE,
  fresh.weight.saturated = FALSE, fresh.weight = FALSE,
  water.potential = FALSE, RWD = FALSE, conductance = FALSE,
  time.since.start = FALSE, temperature = FALSE, humidity = FALSE)
```

## Arguments

data	data frame containing the data to be checked
sample	name of column containing the sample ID (default: sample)
leaf.area	name of column containing the leaf area (cm <sup>2</sup> ) (default: leaf.area)
date.and.time	name of column containing the date and time (default: date and time)
dry.weight	name of column containing the dry weight (g) (default: dry weight)
fresh.weight.saturated	name of column containing the saturated fresh weight (g) (default: fresh.weight.saturated)
fresh.weight	name of column containing the fresh weight (g) (default: fresh.weight)
water.potential	name of column containing the water potential (bar) (default: water.potential)
RWD	name of column containing the relative water deficit (default: RWD)
conductance	name of column containing the conductance values (default: conductance)
time.since.start	name of column containing the time since start values (default: time.since.start)
temperature	name of column containing the temperature values (default: temperature)
humidity	name of column containing the humidity values (default: humidity)

**Value**

no return value

---

ValidityCheckDetail	<i>Checks if column exists in data, is numeric and has the same lengths as the others existence</i>
---------------------	---

---

**Description**

Checks if column exists in data, is numeric and has the same lengths as the others existence

**Usage**

```
ValidityCheckDetail(data_in, value)
```

**Arguments**

data_in	data frame to be checked
value	column in data

**Value**

no return value

---

VaporPressureDeficit	<i>Vapor Pressure Deficit (VPD)</i>
----------------------	-------------------------------------

---

**Description**

Calculates mole fraction vapor pressure deficit ( $\text{mol} * \text{mol}^{-1}$ ) of the air.

**Usage**

```
VaporPressureDeficit(data, humidity = "humidity",  
  temperature = "temperature", atmospheric.pressure = 101.325)
```

**Arguments**

data	data frame at least with two numeric columns of equal length containing humidity (%) and temperature (degree Celsius)
humidity	optional name of the column in data containing the humidity values; default: "humidity"
temperature	optional name of the column in data containing the temperature values; default: "temperature"
atmospheric.pressure	optional; default = 101.325 (atmospheric pressure at sea level)

**Details**

Mole fraction vapor pressure deficit is calculated as:

$$(1 - RH/100) * (VP_{sat}/AP)$$

whereas RH = relative humidity (%), VP<sub>sat</sub> = saturation vapor pressure (kPA), AP = atmospheric pressure (kPA), whereas:

$$VP_{sat} = 0.61121 \exp((18.678 - T/234.5)(T/257.14 + T))$$

where T = air temperature (degree Celsius)

**Value**

The original data frame extended by a numeric column with the vapor pressure deficit (mol \* mol<sup>-1</sup>).

**Examples**

```
# get example data
df <- weather_data

# calculate vapor pressure deficit from weather data measured at sea level
df_with_VPD <- VaporPressureDeficit(df)

# calculate vapor pressure deficit from weather data measured at 2000 m altitude
df_with_VPD <- VaporPressureDeficit(df, atmospheric.pressure = 79.495)
```

---

WeatherAllocation

*Weather Allocation*


---

**Description**

Calculates average weather (humidity, temperature) values within a measurement period.

**Usage**

```
WeatherAllocation(weight_loss_data, weather_data)
```

**Arguments**

weight\_loss\_data

data frame containing at least a column named "date.and.time" of the class POSIXct with the time (and date) of the measuring events. A column named "sample" containing the sample IDs is optionally required if several samples were measured.

weather\_data

data frame containing at least a column named "date.and.time" of the class POSIXct with the time (and date) of the weather measurements and two columns named "humidity" and "temperature" containing the numerical weather data

**Details**

Averages within a measurement period are determined by approximate integration and normalization of the weather as a function of time.

**Value**

The original weight loss data frame extended by the approximatively integrated and normalized weather data for each period between two weight measurements. The first value of each sample is NA since weather values are averaged from row  $i$  to  $i-1$ .

**Examples**

```
# get example data
weight_loss_data <- leaf_drying_data
weather_data <- weather_data

# allocate averaged weather data to weight loss data
weight_loss_data_with_weather <- WeatherAllocation(weight_loss_data, weather_data)
```

---

weather\_data

*Weather data*

---

**Description**

A dataset containing weather data obtained by a data logger at the site of the measurement of the leaf drying data (leaf\_drying\_data)

**Usage**

```
weather_data
```

**Format**

A data frame with 253 rows and 3 variables

**Details**

- date.and.time: Time (and date) of the weather logging event (2019-03-25 00:00:00 - 2019-03-27 15:00:00)
- temperature: Air temperature in degree Celsius (18.05 - 26.81)
- humidity: Relative air humidity in percentage (35.1 - 65.07)

---

WeightDifference	<i>Weight Difference</i>
------------------	--------------------------

---

**Description**

Calculates weight changes between temporally repeated measurements

**Usage**

```
WeightDifference(data, sample = "sample",  
  fresh.weight = "fresh.weight")
```

**Arguments**

data	data frame containing at least a numeric column containing the measured weights (g), ordered chronologically by sample. A column containing the sample IDs is optionally required if several samples were measured.
sample	optional name of the column in data containing the sample ID, default: "sample"
fresh.weight	optional name of the column in data containing the numeric fresh weight (g) values, default: "fresh.weight"

**Value**

the original data frame extended by a numeric column containing the absolute differences in fresh weight (g) between the measurements of a sample. The first value of each sample is NA since weight differences are computed from row *i* and *i*-1.

**Examples**

```
# get example data  
df <- leaf_drying_data  
  
# extend df by weight difference  
df_with_WD <- WeightDifference(df)
```



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