Package 'CalcThemAll.PRM'

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Title Calculate Pesticide Risk Metric (PRM) Values from Multiple
Pesticides...Calc Them All

Version 1.1.1

Description Contains functions which can be used to calculate Pesticide Risk Metric
values in aquatic environments from concentrations of multiple pesti-
cides with known species sensitive
distributions (SSDs). Pesticides provided by this package have all be validated
however if the user has their own pesticides with SSD values they can append them
to the pesticide_info table to include them in estimates.

Encoding UTF-8

License GPL (>= 3)

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Imports dplyr, lubridate, magrittr, MASS, stats, VGAM, plotly, zoo, DT

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R topics documented:

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add_your_own_pesticide

Add new pesticides to the pesticide_info table

Description

Add new pesticides to the pesticide_info table

Usage

```
add_your_own_pesticide(
  pesticides,
  relative_LORs,
  pesticide_types,
  distribution_types,
  shape_locations = NA,
  shape_location_2s = NA,
  scales = NA,
  scale_2s = NA,
  weights = NA,
  pesticide_info = CalcThemAll.PRM::pesticide_info
)
```

Arguments

pesticides   A vector of pesticide names

relative_LORs   A vector of relative limit of reporting replacement values

pesticide_types

              A vector of the new pesticide's types

distribution_types

              A vector of the new pesticide's species sensitivity distribution types

shape_locations

              A vector of shape/location values (if applicable, else put NA)

shape_location_2s

              A vector of secondary shape/location values (if applicable, else put NA)

scales   A vector of scale values (if applicable, else put NA)

scale_2s   A vector of secondary scale values (if applicable, else put NA)

weights   A vector of weight values (if applicable, else put NA)

pesticide_info A data set to add pesticides too

Value

A data frame

Examples

```
new <- add_your_own_pesticide(pesticides = "Poison", relative_LORs = 0.023,
pesticide_types = "Poison", distribution_types = "Log-Normal", scales = 0.09,
shape_locations = 0.014)
multiple_new <- add_your_own_pesticide(pesticides = c("Poison", "Acid", "Sludge"),
relative_LORs = c(0.03, 0.01, 0.5), pesticide_types = c("Poison", "Acid", "Sludge"),
distribution_types = c("Log-Normal", "Log-Logistic Log-Logistic", "Burr Type III"),
scales = c(0.3, 0.002, 2),
scale_2s = c(NA, 0.04, NA),  shape_locations = c(1, 0.07, 3),
shape_location_2s = c(NA, 0.14, 2.3), weights = c(NA, 0.08, NA))
```

Burr_Type_III_Formula   *Burr Type III (Inverse Burr) Equation Formula*

Description

Burr Type III (Inverse Burr) Equation Formula

Usage

```
Burr_Type_III_Formula(concentration, scale, shape_location, shape_location_2)
```

Arguments

concentration	The vector of concentration values for a selected pesticide, that has a Burr Type III shaped species sensitivity distribution, to run the equation on.
scale	The Burr Type III scale/b value for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" data frame.
shape_location	The Burr Type III c/shape value for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.
shape_location_2	
	The Burr Type III k/shape value for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
#Chlorpyrifos is used as its species sensitivity distribution fits Burr Type III
Chlorpyrifos <- c(0.000000001, 0.5, 2.7, 11)
Burr_Type_III_Formula(concentration = Chlorpyrifos,
scale = pesticide_info$scale[pesticide_info$pesticide == "Chlorpyrifos"],
shape_location = pesticide_info$shape_location[pesticide_info$pesticide == "Chlorpyrifos"],
shape_location_2 = pesticide_info$shape_location_2[pesticide_info$pesticide == "Chlorpyrifos"])
```

---

calculate_daily_average_PRM

Calculate Daily Average Pesticide Risk Metric Values For Each Pesticide Type

---

Description

Calculate Daily Average Pesticide Risk Metric Values For Each Pesticide Type

Usage

```
calculate_daily_average_PRM(
  LOR_treated_data,
  include_PAF = FALSE,
  pesticide_info = CalcThemAll.PRM::pesticide_info
)
```

Arguments

LOR_treated_data

> A data set of LOR treated pesticide concentration values in individual columns that match the pesticide names in the "pesticide_info" data frame. This data set should also include a "Date", "Sampling Year" and "Site Name" column.

include_PAF If "TRUE" Percentage Affected Fraction values are included in the output along with Daily PRM in a list format. These values can be useful for plotting relative individual pesticide contribution to overall PRM, however most will not need this so default is "FALSE".

pesticide_info The reference table which contains all relevant information for calculations. It is recommended that the "pesticide_info" data set included in this package be used and if you wish to include more pesticides you can appended them with the relevant information to this table. If you are creating your own table you must ensure that the pesticide name column is title "pesticide" and the relative LOR replacement column is "relative_LOR" for the function to run.

Value

If include_PAF is "FALSE" returns a data frame of daily average PRM values for each pesticide type for each sample. Wet season average calculations can be run on the returned data. If include_PAF is "TRUE" returns a list with daily PRM values in a data frame as the first object and a data frame of PAF values as the second object.

Examples

```
Canto_pesticides_LOR_treated <- treat_LORs_all_data(raw_data = Canto_pesticides,
pesticide_info = CalcThemAll.PRM::pesticide_info)
Canto_daily_PRM <- calculate_daily_average_PRM(LOR_treated_data = Canto_pesticides_LOR_treated)
head(Canto_daily_PRM)
```

---

calculate_wet_season_average_PRM

Calculate Wet Season Average Pesticide Risk Metric Values Using Multiple Imputation

---

Description

Calculate Wet Season Average Pesticide Risk Metric Values Using Multiple Imputation

Usage

```
calculate_wet_season_average_PRM(
  daily_PRM_data,
  PRM_group = "Total PRM",
  imputations = 1000,
  min_sampling_days = 12,
```

```
    wet_season_length = 182
)
```

Arguments

`daily_PRM_data`	A data set of calculated daily average PRM values. This data set should also include a "Date", "Sampling Year" and "Site Name" column.
`PRM_group`	This specifies the name of the column to run the calculations on. The daily average calculations gives PRM for each pesticide type and total in different columns so this selects which to run. "Total" is set as the default as it is the PRM of all pesticides.
`imputations`	This sets the number of imputations to run. The more imputations the greater the reliability, however it also increases calculation time. You can increase imputations beyond 1000 however the improvement of the confidence interval on imputed values may not be sufficient to warrant increased computing time. We recommend 1000 :)

`min_sampling_days`
> This is the minimum number of sampling days a site-year combination must have to calculate a wet season average PRM. No less than 12 is the default (1 for each month) for reliability but more is recommended.

`wet_season_length`
> The length of the wet season in days.

Value

A data frame

Examples

```
Canto_pesticides_LOR_treated <- treat_LORs_all_data(raw_data = Canto_pesticides,
pesticide_info = CalcThemAll.PRM::pesticide_info)
Canto_daily_PRM <- calculate_daily_average_PRM(LOR_treated_data = Canto_pesticides_LOR_treated)
Celestial_City_2019_2020_daily_PRM <- Canto_daily_PRM %>%
dplyr::filter(`Site Name` == "Celestial City" & `Sampling Year` == "2019-2020")
CC2019_2020_wet_season_PSII_PRM <- calculate_wet_season_average_PRM(daily_PRM_data =
Celestial_City_2019_2020_daily_PRM, PRM_group = "PSII Herbicide PRM")
CC2019_2020_wet_season_PSII_PRM
```

Canto_pesticides   *Canto Region Pesticide Concentration Values (Example Data Set)*

Description

A subset of pesticide concentration data for all pesticides in "pesticide_info" created for this package with fabricated sites

Usage

```
Canto_pesticides
```

Format

Canto_pesticides:

A data frame with 808 rows and 64 columns:

Site Name Site name

Date Sampling date

Ametryn, Atrazine, Chlorpyrifos, Diuron, Fipronil, Fluroxypyr, Haloxyfop (acid), Hexazinone, Imazapic, Imidaclop
pesticide concentration values in ug/L ...

---

find_Sampling_Year   *Find Sampling Year*

Description

Find Sampling Year

Usage

```
find_Sampling_Year(dates, wet_season_split = 7)
```

Arguments

dates   A date vector of sampling dates. Must be in yyyy-mm-dd format.

wet_season_split

The first month of the sampling year in numeric e.g. July = 7. July (7) is used
as the default as this is the first month of the Queensland wet season.

Value

A factored character vector

Examples

```
dates <- as.Date(c("2014-03-04", "2014-12-30", "2015-06-12"))
sampling_years <- find_Sampling_Year(dates) #cut of date
```

---

find_season   *Find Sample's Season (Wet or Dry)*

Description

Find Sample's Season (Wet or Dry)

Usage

```
find_season(wet_season_start_dates, sampling_dates, wet_season_length = 182)
```

Arguments

wet_season_start_dates

> A vector of dates signifying the first day of the wet season for site year combinations.

sampling_dates A date vector of sampling dates. Must be in yyyy-mm-dd format.

wet_season_length

> The length of the wet season in days.

Value

A character vector

Examples

```
dates <- as.Date(c("2014-12-04", "2014-10-30", "2015-11-12"))
wet_start_dates <- as.Date(c("2014-10-04", "2014-12-30", "2015-09-12"))
Seasons <- find_season(wet_start_dates, sampling_dates = dates)
#cut of date for the sampling year will be last day of June
```

---

find_wet_season_end   *Find Wet Season End Date*

Description

Find Wet Season End Date

Usage

```
find_wet_season_end(wet_season_start, wet_season_length = 182)
```

Arguments

wet_season_start

A vector of dates signifying the first day of the wet season for site year combinations.

wet_season_length

The length of the wet season in days.

Value

A character vector

Examples

```
wet_season_start_dates <- as.Date(c("2014-10-04", "2014-12-30", "2015-09-12"))
wet_season_end_dates <- find_wet_season_end(wet_season_start_dates)
#cut of date for the sampling year will be last day of June
```

Gamma_Formula *Gamma CDF Equation Formula*

Description

Gamma CDF Equation Formula

Usage

```
Gamma_Formula(concentration, shape_location, scale)
```

Arguments

concentration   The vector of concentration values for a selected pesticide, that has a Gamma shaped species sensitivity distribution, to run the equation on.

shape_location   The k/shape value for the Gamma equation for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

scale   The scale/theta value for the Gamma equation for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
gamma_pesticide_concentrations <- c(0.000000001, 0.5, 2.7, 11)
Gamma_Formula(concentration = gamma_pesticide_concentrations,
shape_location = 0.23, scale = 1.3)
```

---

imputation_beta   *Imputation Function - Beta Version*

Description

Imputation Function - Beta Version

Usage

```
imputation_beta(impute_variable, wet_season_length = 182)
```

Arguments

impute_variable
                    The variable you wish to impute.
wet_season_length
                    The length of the wet season in days.

Value

A data frame.

---

imputation_kernel   *Imputation Function - Kernal Version*

Description

Imputation Function - Kernal Version

Usage

```
imputation_kernel(impute_variable, wet_season_length = 182)
```

Arguments

impute_variable
                    The variable you wish to impute.
wet_season_length
                    The length of the wet season in days.

Value

A data frame.

Inverse_Weibull_Formula

Inverse Weibull Formula

Description

Inverse Weibull Formula

Usage

```
Inverse_Weibull_Formula(concentration, shape_location, scale)
```

Arguments

concentration   The vector of concentration values for a selected pesticide, that has a Inverse Weibull shaped species sensitivity distribution, to run the equation on.

shape_location   The Inverse Weibull shape/alpha value for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

scale   The Inverse Weibull scale/beta value for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
Hexazinone <- c(0.000000001, 0.5, 2.7, 11)
#Hexazinone is used as its species sensitivity distribution plots fits Inverse Weibull
Inverse_Weibull_Formula(concentration = Hexazinone,
shape_location = pesticide_info$shape_location[pesticide_info$pesticide == "Hexazinone"],
scale = pesticide_info$scale[pesticide_info$pesticide == "Hexazinone"])
```

Log_Gumbel_Formula   *Log Gumbel CDF Equation Formula*

Description

Log Gumbel CDF Equation Formula

Usage

```
Log_Gumbel_Formula(concentration, shape_location, scale)
```

Arguments

concentration	The vector of concentration values for a selected pesticide, that has a Log Gumbel shaped species sensitivity distribution, to run the equation on.
shape_location	The mu/location value for the Log Gumbel equation for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.
scale	The beta/scale value for the Log Gumbel equation for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
#This Example should produce roughly 1% and 5% PRM values
LogGumbel_pesticide_concentrations <- c(0.095957794, 0.245881898)
Log_Gumbel_Formula(concentration = LogGumbel_pesticide_concentrations,
shape_location = 0.9980581, scale = 2.188285)
```

Log_Logistic_Formula   *Log Logistic Formula*

Description

Log Logistic Formula

Usage

```
Log_Logistic_Formula(concentration, scale, shape_location)
```

Arguments

concentration	The vector of concentration values for a selected pesticide, that has a Log Logistic shaped species sensitivity distribution, to run the equation on.
scale	The Log Logistic alpha/scale value for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

shape_location The Log Logistic beta/shape value for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
Imazapic <- c(0.000000001, 0.5, 2.7, 11)
#Imazapic is used as its species sensitivity distribution plots fits Log Logistic
Log_Logistic_Formula(concentration = Imazapic,
scale = pesticide_info$scale[pesticide_info$pesticide == "Imazapic"],
shape_location = pesticide_info$shape_location[pesticide_info$pesticide == "Imazapic"])
```

Log_Logistic_Log_Logistic_Formula

Log Logistic Log Logistic (double curve) CDF Formula

Description

Log Logistic Log Logistic (double curve) CDF Formula

Usage

```
Log_Logistic_Log_Logistic_Formula(
  concentration,
  scale,
  shape_location,
  scale_2,
  shape_location_2,
  weight
)
```

Arguments

concentration The vector of concentration values for a selected pesticide, that has a Log Logistic Log Logistic shaped species sensitivity distribution, to run the equation on.

scale The alpha/scale value for the first Log Logistic equation for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

shape_location   The beta/shape value for the first Log Logistic equation for the selected pesti-
                  cide. These can be found in the "pesticide_info" data frame provided in this
                  package. If you are including other pesticides you will need to append them
                  with their respective distribution variables to the "pesticide_info" table.

scale_2   The alpha/scale value for the second Log Logistic equation for the selected pes-
                  ticide. These can be found in the "pesticide_info" data frame provided in this
                  package. If you are including other pesticides you will need to append them
                  with their respective distribution variables to the "pesticide_info" table.

shape_location_2

                  The beta/shape value for the second Log Logistic equation for the selected pes-
                  ticide. These can be found in the "pesticide_info" data frame provided in this
                  package. If you are including other pesticides you will need to append them
                  with their respective distribution variables to the "pesticide_info" table.

weight   The weight parameter for combining the two equations for the selected pesticide.
                  These can be found in the "pesticide_info" data frame provided in this package.
                  If you are including other pesticides you will need to append them with their
                  respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
#This Example should produce roughly 1% and 5% PRM values
LogL_LogL_pesticide_concentrations <- c(0.00341453, 0.009854566)
Log_Logistic_Log_Logistic_Formula(concentration = LogL_LogL_pesticide_concentrations,
scale = 0.5823392, shape_location = -3.499604, scale_2 = 1.144555,
shape_location_2 = 1.100755, weight = 0.3585467)
```

Log_Normal_Formula   *Log-Normal CDF Equation Formula*

Description

Log-Normal CDF Equation Formula

Usage

```
Log_Normal_Formula(concentration, shape_location, scale)
```

Arguments

concentration	The vector of concentration values for a selected pesticide, that has a Log Normal shaped species sensitivity distribution, to run the equation on.
shape_location	The mu/location value for the Log Normal equation for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.
scale	The sigma/scale value for the Log Normal equation for the selected pesticide. These can be found in the "pesticide_info" data frame provided in this package. If you are including other pesticides you will need to append them with their respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
LogN_pesticide_concentrations <- c(0.000000001, 0.5, 2.7, 11)
Log_Normal_Formula(concentration = LogN_pesticide_concentrations,
shape_location = 0.23, scale = 1.3)
```

Log_Normal_Log_Normal_Formula

Log-Normal Log-Normal (double curve) CDF Equation Formula

Description

Log-Normal Log-Normal (double curve) CDF Equation Formula

Usage

```
Log_Normal_Log_Normal_Formula(
  concentration,
  shape_location,
  scale,
  shape_location_2,
  scale_2,
  weight
)
```

Arguments

concentration	The vector of concentration values for a selected pesticide, that has a Log Normal Log Normal species sensitivity distribution, to run the equation on.

shape_location   The mu/shape value for the first Log Normal equation for the selected pesticide.
                  These can be found in the "pesticide_info" data frame provided in this package.
                  If you are including other pesticides you will need to append them with their
                  respective distribution variables to the "pesticide_info" table.

scale   The sigma/scale value for the first Log Normal equation for the selected pes-
                  ticide. These can be found in the "pesticide_info" data frame provided in this
                  package. If you are including other pesticides you will need to append them
                  with their respective distribution variables to the "pesticide_info" table.

shape_location_2
                  The mu/shape value for the second Log Normal equation for the selected pes-
                  ticide. These can be found in the "pesticide_info" data frame provided in this
                  package. If you are including other pesticides you will need to append them
                  with their respective distribution variables to the "pesticide_info" table.

scale_2   The sigma/scale value for the second Log Normal equation for the selected pes-
                  ticide. These can be found in the "pesticide_info" data frame provided in this
                  package. If you are including other pesticides you will need to append them
                  with their respective distribution variables to the "pesticide_info" table.

weight   The weight parameter for combining the two equations for the selected pesticide.
                  These can be found in the "pesticide_info" data frame provided in this package.
                  If you are including other pesticides you will need to append them with their
                  respective distribution variables to the "pesticide_info" table.

Value

a numeric vector

Examples

```
#This Example should produce roughly 1% and 5% PRM values
LogN_LogN_pesticide_concentrations <- c(4.79E-05, 0.000225588)
Log_Normal_Log_Normal_Formula(concentration = LogN_LogN_pesticide_concentrations,
shape_location = -5.596431, scale = 2.061943,
shape_location_2 = 0.01174954, scale_2 = 0.9134796, weight = 0.5733126)
```

---

pesticide_info   *Pesticide Information for Pesticide Risk Metric Calculations (Refer-*
                               ence Table)

Description

A reference table for PRM calculations in this package that include pesticide information such as
type, species sensitivity distribution distributions and relevant equation variables.

Usage

```
pesticide_info
```

Format

pesticide_info:
A data frame with 62 rows and 9 columns:

pesticide Pesticide name

relative_LOR The relative Limit of Reporting replacement value

pesticide_type The pesticide method of effect

distribution_type The species sensitivity distribution shape/type

shape_location, shape_location2, scale, scale2, weight Species sensitivity distribution variables
used in PRM calculations ...

Source

https://www.publications.qld.gov.au/dataset/method-development-pesticide-risk-metric-baseline-condi
resource/c65858f9-d7ba-4aef-aa4f-e148f950220f

---

plot_daily_PRM *Plot Daily Average PRM Values for a Single Site/Sampling Year*

Description

Plot Daily Average PRM Values for a Single Site/Sampling Year

Usage

```
plot_daily_PRM(
  daily_PRM_data,
  wet_season_start = NULL,
  wet_season_length = 182,
  PRM_group = "Total PRM",
  title = FALSE,
  legend = "Numerical"
)
```

Arguments

daily_PRM_data A data set of calculated daily average PRM values for a single site and sampling
year. This data set should also include a "Date", "Sampling Year" and "Site
Name" column.

wet_season_start

The date of the start of the wet season for this site sampling year. If not applica-
ble leave as NA

wet_season_length

The length of the wet season in days.

PRM_group This specifies the name of the column to plot. The daily average calculations
gives PRM for each pesticide type and a total in different columns. "Total" is set
as the default as it is the PRM value for all pesticides.

title   TRUE or FALSE value to include a title.

legend   Does the legend show "Numerical" or "Categorical" values for PRM values on
                         the plot.

Value

A plotly plot

Examples

```
Canto_pesticides_LOR_treated <- treat_LORs_all_data(raw_data = Canto_pesticides,
pesticide_info = CalcThemAll.PRM::pesticide_info)
Canto_daily_PRM <- calculate_daily_average_PRM(LOR_treated_data = Canto_pesticides_LOR_treated)
Violet_Town_2017_2018_PRM <- Canto_daily_PRM %>%
dplyr::filter(.data$`Sampling Year` == "2017-2018" & .data$`Site Name` == "Violet Town")
plot_daily_PRM(Violet_Town_2017_2018_PRM, "2017-10-02", PRM_group = "Total PRM")
```

---

plot_wet_season_window
                        Plot Wet Season Window Box on plot_daily_PRM

Description

Plot Wet Season Window Box on plot_daily_PRM

Usage

```
plot_wet_season_window(wet_season_start = 0, wet_season_length = 182)
```

Arguments

wet_season_start
                        The date of the start of the wet season for this site sampling year.

wet_season_length
                        The length of the wet season in days.

Value

A plotly shape

Examples

```
shape <- plot_wet_season_window(wet_season_start = "2017-08-01")
```

PRM_DT *PRM Data Table Function*

Description

PRM Data Table Function

Usage

```
PRM_DT(PRM_data, fill_cols = NULL, colour_cols = NULL)
```

Arguments

PRM_data   a data frame of either daily average or wet season PRM values

fill_cols   A vector of column names of pesticide groups to fill colour with corresponding PRM risk category

colour_cols   A vector of column names of pesticide groups to colour text with corresponding PRM risk category

Value

a data table colour coded to PRM risk

Examples

```
Canto_pesticides_LOR_treated <- treat_LORs_all_data(raw_data = Canto_pesticides,
pesticide_info = CalcThemAll.PRM::pesticide_info)
Canto_daily_PRM <- calculate_daily_average_PRM(LOR_treated_data = Canto_pesticides_LOR_treated)
PRM_DT(PRM_data = Canto_daily_PRM, fill_cols = "Total PRM",
colour_cols = c("PSII Herbicide PRM", "Other Herbicide PRM", "Insecticide PRM"))
```

---

treat_LORs *Treat a single observations LORs*

Description

Treat a single observations LORs

Usage

```
treat_LORs(
  sample_data,
  pesticide_info = CalcThemAll.PRM::pesticide_info,
  treatment_method = "Zero"
)
```

Arguments

sample_data	A single observation containing a concentration value for each pesticide being used in the metric. LOR values should be in "<0.05" format and no values should be empty "".
pesticide_info	The reference table which contains all relevant information for calculations. It is recommended that the "pesticide_info" dataset included in this package be used and if you wish to include more or less pesticides you can appended them with the relevant information to this table. If you are creating your own table you must ensure that the pesticide name column is title "pesticides" and the relative LOR replacement column is "relative_LOR" for the function to run.

treatment_method

Select how to treat the LOR values with either "WQI" representing the Queensland Department of Environment & Science Water Quality Monitoring & Investigations team's method for replacing LORs or "Zero" which replaces them with a negligible numeric value. Zero is the default here as this function on its own only treats a single observation and therefore the first detection in the WQI method cannot be used.

Value

returns the provided data set with the first row's LOR values treated.

Examples

```
first_sample <- Canto_pesticides[1,] #this selects only the first row (sample)
LOR_treated_first_sample <- treat_LORs(sample_data = first_sample,
pesticide_info = CalcThemAll.PRM::pesticide_info, treatment_method = "Zero")
print(LOR_treated_first_sample)
```

---

treat_LORs_all_data   *Treat a whole data set's LOR values*

Description

Treat a whole data set's LOR values

Usage

```
treat_LORs_all_data(
  raw_data,
  pesticide_info = CalcThemAll.PRM::pesticide_info,
  wet_season_split = 7,
  treatment_method = "WQI"
)
```

Arguments

raw_data   A data set of raw pesticide concentration values in individual columns that match
                  the pesticide names in the "pesticide_info" data frame. This data set should also
                  include a "Date" column and "Site Name" column. A reference data set can be
                  seen in the "Canto_pesticides" data frame provided in this package, your data
                  should mirror these column headings.

pesticide_info   The reference table which contains all relevant information for calculations. It is
                  recommended that the "pesticide_info" data set included in this package be used
                  and if you wish to include more or less pesticides you can appended them with
                  the relevant information to this table. If you are creating your own table you
                  must ensure that the pesticide name column is title "pesticides" and the relative
                  LOR replacement column is "relative_LOR" for the function to run.

wet_season_split
                  The first month of the sampling year in numeric e.g. July = 7. July (7) is used as
                  the default as this is the first month of the Queensland wet season. This is only
                  required for the LOR replacement method and if needed.

treatment_method
                  Select how to treat the LOR values with either the default "WQI" representing
                  the Queensland Department of Environment & Science Water Quality Moni-
                  toring & Investigations team's method for replacing LORs or "Zero" which re-
                  places them with a negligible numeric value.

Value

returns the raw_data frame with the LOR values replaced by their specified treatment values. PRM
calculations can now be run on the returned data.

Examples

```
Canto_pesticides_LOR_treated <- treat_LORs_all_data(raw_data = Canto_pesticides,
pesticide_info = CalcThemAll.PRM::pesticide_info)
head(Canto_pesticides_LOR_treated)
```

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