Package 'PhenoSpectra'

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Title Multispectral Data Analysis and Visualization

Version 0.1.0
Description Provides tools for processing, analyzing, and visualizing spectral data collected from 3D laser-based scanning systems. Supports applications in agriculture, forestry, environmental monitoring, industrial quality control, and biomedical research. Enables evaluation of plant growth, productivity, resource efficiency, disease management, and pest monitoring. Includes statistical methods for extracting insights from multispectral and hyperspectral data and generating publication-ready visualizations. See Zieschank & Junker (2023) <doi:10.3389 fpls.2023.1141554=""> and Saric et al. (2022) <doi:10.1016 j.tplants.2021.12.003=""> for related work.</doi:10.1016></doi:10.3389>
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feature_selection

Feature Selection for Spectral Data

Description

This function filters healthy vs diseased samples, selects the most discriminative spectral variables, applies FDR correction, and exports the results.

Usage

```
feature_selection(
  file_path,
  output_path = "selected_features.xlsx",
  fdr_threshold = 0.01
)
```

Arguments

```
file_path Path to the cleaned dataset (output of qaqcs function).

output_path Path to save the selected features table.

fdr_threshold Threshold for filtering significant features (default: 0.01).
```

Value

A data.table containing selected spectral variables.

```
# Create mock spectral data
library(openxlsx)
mock_data <- data.frame(
    treatment = sample(0:1, 100, replace = TRUE),
    var1 = rnorm(100),
    var2 = rnorm(100),
    var3 = rnorm(100)
)
temp_file <- tempfile(fileext = ".xlsx")
write.xlsx(mock_data, temp_file)

# Perform feature selection
output_path <- tempfile(fileext = ".xlsx")
selected_features <- feature_selection(temp_file, output_path, fdr_threshold = 0.01)
head(selected_features)</pre>
```

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predict_SDS

Predict Spectral Disease Severity (SDS)

Description

This function predicts Spectral Disease Severity (SDS) using a standard linear regression model (lm()). It automatically handles column names with special characters by using backticks and constrains predictions to the range [0, 100].

Usage

```
predict_SDS(cleaned_data, sf_test, fixed_effects = NULL)
```

Arguments

cleaned_data A dataframe containing spectral measurements and treatment labels.

sf_test A dataframe containing selected important features (from statistical tests).

fixed_effects A character vector of fixed effects to include (default: NULL). Example: c("Scan.date").

Value

A dataframe with predicted SDS values for all treatments, constrained between 0 and 100.

```
# Create mock spectral data
library(openxlsx)
cleaned_data <- data.frame(
    treatment = sample(0:1, 100, replace = TRUE),
    var1 = rnorm(100),
    var2 = rnorm(100),
    var3 = rnorm(100),
    Scan.date = sample(
        seq.Date(
            from = as.Date('2023-01-01'),
            to = as.Date('2023-12-31'),
            by = 'day'
        ),
        100
      ),
      Scan.time = format(Sys.time(), "%H:%M:%S")
)</pre>
```

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qaqcs

Perform QA/QC on spectral data while preserving original column names

Description

Perform QA/QC on spectral data while preserving original column names

Usage

```
qaqcs(
  file_path,
  output_path,
  handle_missing = "NA",
  handle_outliers = "NA",
  group_by_col = "treatment"
)
```

Arguments

```
file_path Path to the input file
output_path Path to save the cleaned data
handle_missing Method to handle missing values ('impute', 'remove', or 'NA')
handle_outliers

Method to handle outliers ('impute', 'remove', or 'NA')
group_by_col Column name for grouping
```

Value

A list with cleaned data and a summary table

```
library(openxlsx)
# Create mock raw data
raw_data <- data.frame(
    treatment = sample(0:1, 100, replace = TRUE),
    var1 = rnorm(100),
    var2 = rnorm(100),
    var3 = rnorm(100)
)

# Save mock data to a temporary file
raw_data_file <- tempfile(fileext = ".xlsx")
output_file <- tempfile(fileext = ".xlsx")
write.xlsx(raw_data, raw_data_file)</pre>
```

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```
# Run QA/QC with missing values imputed and outliers removed
cleaned_result <- qaqcs(
   file_path = raw_data_file,
   output_path = output_file,
   handle_missing = "impute",
   handle_outliers = "remove",
   group_by_col = "treatment"
)
head(cleaned_result$cleaned_data)</pre>
```

reads

Read and merge spectral data using data.table exclusively

Description

Read and merge spectral data using data.table exclusively

Usage

```
reads(directory, pattern, output_path)
```

Arguments

pattern Path to the directory containing files

File pattern to search for (e.g., 'input')

output_path Path to save the processed output

Value

A merged data.table

```
library(data.table)
# Create mock data files with all required columns
mock_data1 <- data.frame(
    treatment = sample(0:1, 50, replace = TRUE),
    var1 = rnorm(50),
    var2 = rnorm(50),
    Scan.date = sample(
        seq.Date(
            from = as.Date('2023-01-01'),
            to = as.Date('2023-12-31'),
            by = 'day'
        ),
        50,
        replace = TRUE
    ),
    Scan.time = format(Sys.time(), "%H:%M:%S"),</pre>
```

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```
timestamp = Sys.time() # Add timestamp column
mock_data2 <- data.frame(</pre>
  treatment = sample(0:1, 50, replace = TRUE),
  var1 = rnorm(50),
  var2 = rnorm(50),
  Scan.date = sample(
    seq.Date(
      from = as.Date('2023-01-01'),
      to = as.Date('2023-12-31'),
      by = 'day'
    ),
    50,
    replace = TRUE
  ),
  Scan.time = format(Sys.time(), "%H:%M:%S"),
  timestamp = Sys.time() # Add timestamp column
)
# Save mock data to temporary CSV files
temp_dir <- tempdir()</pre>
file1 <- file.path(temp_dir, "input_file1.csv")</pre>
file2 <- file.path(temp_dir, "input_file2.csv")</pre>
fwrite(mock_data1, file1)
fwrite(mock_data2, file2)
# Run the reads() function on mock CSV data
merged_data <- reads(</pre>
  directory = temp_dir,
  pattern = "input",
  output_path = tempfile(fileext = ".csv")
)
head(merged_data)
```

setup_environment

Setup Environment for PhenoSpectra

Description

Automatically installs and loads all required packages.

This function installs missing packages required by the PhenoSpectra package.

Usage

```
setup_environment()
setup_environment()
```

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Value

NULL. This function installs required packages if missing and loads them.

```
# Example usage:
setup_environment() # Automatically installs and loads required packages
# Run to manually install missing packages
setup_environment()
```

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