# Testing Plan for baselineARPLss package (DATA501)

## Installation

The following steps should work either on your own computer or on a free project / instance in the R Cloud (https://posit.cloud/)

The steps to download the package and install it are the following:

```
install.packages('remotes')
library(remotes)
remotes::install_github("econdatatech/baselineARPLss")
```

Please try these steps and report the result. The expected result is an installation without errors.

# Usage

# Usage with supplied sample data

#### Baseline estimation

Try the following instructions on in your R environment and report the result. The execution might take a couple of minutes and should finish without an error.

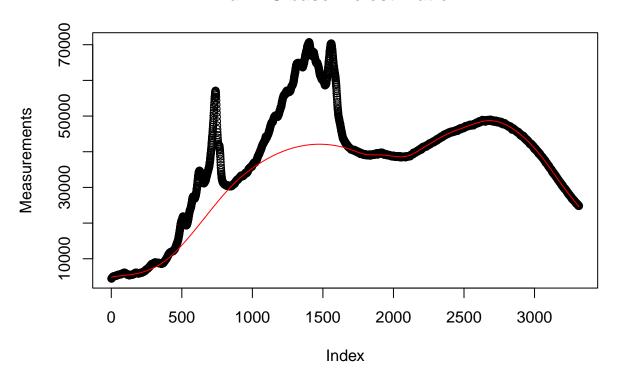
```
library(baselineARPLss)
data("Abelsonite")
res<-baseline_estimation(Abelsonite$measurement)</pre>
```

#### Result plot

Issue the below command and compare the result with the picture below.

```
plot(res)
```





Issue the below command and compare the result with the output below.

#### summary(res)

```
## [1] "The lambda parmeter value used was: 1e+06"
```

- ## [1] "The ratio parameter value used was: 1e-06"
- ## [1] "The max\_iter parameter value used was: 50"
- ## [1] "The alogrithm stopped after the following number or iterations: 50"
- ## [1] "The last weight ratio value was: 0.000153623596796175"
- ## [1] "It appears that the algorithm stopped because the maximum number of iterations was reached"

Try to obtain the help page for the baseline\_estimation function.

#### ?baseline\_estimation

#### ## starting httpd help server ... done

Give your opinion. E.g. is anything unclear or anything missing?

## Usage with new data

Select a substance or mineral of your choice from this following URL https://rruff.info/\*/display=default/ (The site might take a bit to load) Click on the magnifying glass on the right hand side of the screen for the substance of your choice. On the next page look for the following section "BROAD SCAN WITH SPECTRAL ARTIFACTS" and download the "Raman Data (RAW)" file. If such a file is not available, choose a different substance in the previous screen and try to obtain a "Raman Data (RAW)" for this substance.

Read the downloaded file into a dataframe (with a command similar to the one below)

```
df<-read.table('C:\\Users\\corvini\\Downloads\\Actinolite_R040063_Broad_Scan_532_0_unoriented_Ram
```

Try to estimate the baseline for the spectrum of this substance with the help of the documentation available (e.g. help files and vignette and manual) Report your findings.

#### Input validation

The baseline\_estimation function has a lot of parameters. Most (apart from the input data in form of the raw spectrum) have default parameters.

baseline\_estimation <- function(y, lambda = 1e6, ratio = 1e-6, max\_iter = 50, verbose=FALSE, algo="banded"

#### 3. Input Validation

Objective: Confirm that functions handle different types of input data appropriately.

Test handling of valid inputs such as numeric vectors representing Raman spectra.

Provide invalid inputs (e.g., missing values, non-numeric data, inappropriate dimensions) and check for Test behavior with edge cases like empty spectra or spectra with only a few data points.

4. Performance and Efficiency

Objective: Ensure the package performs efficiently for large datasets.

Tests:

Time the execution of baseline estimation for both small and large datasets.

Test memory usage and confirm the package does not excessively increase memory consumption for large Ra Evaluate multi-threading or parallel processing capabilities if implemented.