

Package ‘RLumModel’

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Type Package

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Luminescence

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Description A collection of function to simulate luminescence signals in the
mineral quartz based on published models.

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License GPL-3

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RoxygenNote 5.0.1

Suggests testthat

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RLumModel-package*Modelling Ordinary Differential Equations Leading to Luminescence*

Description

A collection of function to simulate luminescence signals in the mineral quartz based on published models.

Details

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Project source code repository

<https://github.com/R-Lum/RLumModel>

Related projects

<http://www.r-luminescence.de>
<http://cran.r-project.org/package=Luminescence>
<http://shiny.r-luminescence.de>
<http://cran.r-project.org/package=RLumShiny>

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ExampleData.ModelOutput

Example data (TL curve) simulated from Bailey (2001 ,fig. 1)

Description

Example data (TL curve) simulated from Bailey (2001 ,fig. 1)

Format

A RLum.Analysis object containing one TL curve as RLum.Data.Curve.

Note

This example has only one record (TL). The used sequence was `sequence <- list(IRR = c(temp = 20, dose = 10, DoseRate = 1), TL = c(temp_begin = 20, temp_end = 400, heating_rate = 5))`

Source

`model_LuminescenceSignals()`

References

Bailey, R.M., 2001. Towards a general kinetic model for optically and thermally stimulated luminescence of quartz. Radiation Measurements 33, 17-45.

Examples

```
data(ExampleData.ModelOutput,envir = environment())
plot_RLum.Analysis(model.output)
```

model_LuminescenceSignals

Model Luminescence Signals

Description

This function models luminescence signals for quartz based on published physical models. It is possible to simulate TL, (CW-) OSL, RF measurements in a arbitrary sequence. This sequence is defined as a `list` of certain abrivations. Furthermore it is possible to load a sequence direct from the Riso Sequence Editor. The output is an `RLum.Analysis` object and so the plots are done by the `plot_RLum.Analysis` function. If a SAR sequence is simulated the plot output can be disabled and SAR analyse functions can be used.

Usage

```
model_LuminescenceSignals(model, sequence, lab.dose_rate = 1,
  simulate_sample_history = FALSE, plot = TRUE, verbose = TRUE,
  show.structure = FALSE, ...)
```

Arguments

model **character (required)**: set model to be used. Available models are: "Bailey2001", "Bailey2002", "Bailey2004", "Pagonis2007", "Pagonis2008"

sequence **list (required)**: set sequence to model as **list** or as *.seq file from the Riso sequence editor. To simulate SAR measurements there is an extra option to set the sequence list (cf. details).

lab.dose_rate **numeric** (with default): laboratory dose rate in XXX Gy/s for calculating seconds into Gray in the *.seq file.

simulate_sample_history **logical** (with default): FALSE (with default): simulation begins at laboratory conditions, TRUE: simulations begins at crystallization (all levels 0) process

plot **logical** (with default): Enables or disables plot output

verbose **logical** (with default): Verbose mode on/off

show.structure **logical** (with default): Shows the structure of the result. Recommended to show record.id to analyse with **plot_concentrations**.

... further arguments and graphical parameters passed to **plot.default**. See details for further information.

Details

Defining a sequence

| Arguments | Description | Sub-arguments |
|-----------|-----------------------------------|--|
| TL | thermally stimulated luminescence | 'temp begin', 'temp end', 'heating rate' |
| OSL | optically stimulated luminescence | 'temp', 'duration', 'optical_power' |
| ILL | illumination | 'temp', 'duration', 'optical_power' |
| LM_OSL | linear modulated OSL | 'temp', 'duration', optional: 'start_power', 'end_power' |
| RL/RF | radioluminescence | 'temp', 'dose', 'dose_rate' |
| IRR | irradiation | 'temp', 'dose', 'dose_rate' |
| CH | cutheat | 'temp', optional: 'duration', 'heating_rate' |
| PH | preheat | 'temp', 'duration' optional: 'heating_rate' |
| PAUSE | pause | 'temp', 'duration' |

Defining a SAR-sequence

| Abrivation | Description | examples |
|--------------|--|--------------------------------|
| RegDose | Dose points of the regenerative cycles | c(0, 80, 140, 260, 320, 0, 80) |
| TestDose | Test dose for the SAR cycles | 50 |
| PH | Temperature of the preheat | 240 |
| CH | Temperature of the cutheat | 200 |
| OSL_temp | Temperature of OSL read out | 125 |
| OSL_duration | Duration of OSL read out | default: 40 |

| | | |
|---------------|--|-------------|
| Irr_temp | Temperature of irradiation | default: 20 |
| PH_duration | Duration of the preheat | default: 10 |
| dose_rate | Dose rate of the laboratory irradiation source | default: 1 |
| optical_power | Percentage of the full illumination power | default: 90 |
| Irr_2recover | Dose to be recovered in a dose-recovery-test | 20 |

Value

This function returns an [RLum.Analysis](#) object with all TL, (LM-) OSL and RF/RL steps in the sequence. Every entry is an [RLum.Data.Curve](#) object and can be plotted, analysed etc. with further RLum-functions.

Function version

0.1.0

Author(s)

Johannes Friedrich, University of Bayreuth (Germany), Sebastian Kreutzer, IRAMAT-CRP2A, Université Bordeaux Montaigne (France)

References

- Bailey, R.M., 2001. Towards a general kinetic model for optically and thermally stimulated luminescence of quartz. *Radiation Measurements* 33, 17-45.
- Bailey, R.M., 2002. Simulations of variability in the luminescence characteristics of natural quartz and its implications for estimates of absorbed dose. *Radiation Protection Dosimetry* 100, 33-38.
- Bailey, R.M., 2004. Paper I-simulation of dose absorption in quartz over geological timescales and its implications for the precision and accuracy of optical dating. *Radiation Measurements* 38, 299-310.
- Pagonis, V., Chen, R., Wintle, A.G., 2007: Modelling thermal transfer in optically stimulated luminescence of quartz. *Journal of Physics D: Applied Physics* 40, 998-1006.
- Pagonis, V., Wintle, A.G., Chen, R., Wang, X.L., 2008. A theoretical model for a new dating protocol for quartz based on thermally transferred OSL (TT-OSL). *Radiation Measurements* 43, 704-708.
- Soetaert, K., Cash, J., Mazzia, F., 2012. Solving differential equations in R. Springer Science & Business Media.

See Also

[plot](#), [RLum](#), [plot_concentrations](#), [read_SEQ2R](#)

Examples

```
##=====##
## Example 1: Simulate sample history of Bailey2001
## (cf. Bailey, 2001, Fig. 1)
##=====##

##set sequence with the following steps
## (1) Irradiation at 20 deg. C with a dose of 10 Gy and a dose rate of 1 Gy/s
```

```

## (2) TL from 20-400 deg. C with a rate of 5 K/s

sequence <-
  list(
    IRR = c(20, 10, 1),
    TL = c(20, 400, 5)
  )

##model sequence
model.output <- model_LuminescenceSignals(
  sequence = sequence,
  model = "Bailey2001",
  show.structure = TRUE
)

## Not run:
##=====##
## Example 2: Simulate sequence at labour without sample history
##=====##

##set sequence with the following steps
## (1) Irraditation at 20 deg. C with a dose of 100 Gy and a dose rate of 1 Gy/s
## (2) Preheat to 200 deg. C and hold for 10 s
## (3) LM-OSL at 125 deg. C. for 100 s
## (4) Cutheat at 200 dec. C.
## (5) Irraditation at 20 deg. C with a dose of 10 Gy and a dose rate of 1 Gy/s
## (6) Pause at 200 de. C. for 100 s
## (7) OSL at 125 deg. C for 100 s with 90 % optical power
## (8) Pause at 200 deg. C for 100 s
## (9) TL from 20-400 deg. C with a heat rate of 5 K/s
## (10) Radioluminescence at 20 deg. C with a dose of 20 Gy and a dose rate of 0.01 Gy/s

sequence <-
  list(
    IRR = c(20, 100, 1),
    PH = c(200, 10),
    LM_OSL = c(125, 100),
    CH = c(200),
    IRR = c(20, 10, 1),
    PAUSE = c(200, 100),
    OSL = c(125, 100, 90),
    PAUSE = c(200, 100),
    TL = c(20, 400, 5),
    RF = c(20, 200, 0.01)
  )

# call function "model_LuminescenceSignals", set sequence = sequence,
# model = "Pagonis2008" (palaeodose = 200 Gy) and simulate_sample_history = FALSE (default),
# because the sample history is not part of the sequence

model.output <- model_LuminescenceSignals(
  sequence = sequence,
  model = "Pagonis2008"
)

```

```

#####
## Example 3: Simulate SAR sequence
#####

##set SAR sequence with the following steps
## (1) RegDose: set regenerative dose [Gy] as vector
## (2) TestDose: set test dose [Gy]
## (3) PH: set preheat temperature in deg. C
## (4) CH: Set cutheat temperature in deg. C
## (5) OSL_temp: set OSL reading temperature in deg. C
## (6) OSL_duration: set OSL reading duration in s

sequence <- list(
  RegDose = c(0,10,20,50,90,0,10),
  TestDose = 5,
  PH = 240,
  CH = 200,
  OSL_temp = 125,
  OSL_duration = 70)

# call function "model_LuminescenceSignals", set sequence = sequence,
# model = "Pagonis2007" (palaeodose = 20 Gy) and simulate_sample_history = FALSE (default),
# because the sample history is not part of the sequence

model.output <- model_LuminescenceSignals(

  sequence = sequence,
  model = "Pagonis2007",
  plot = FALSE
)

# in environment is a new object "model.output" with the results of
# every step of the given sequence.
# Plots are done at OSL and TL steps and the growth curve

# call "analyse_SAR.CWOSL" from RLum package
results <- analyse_SAR.CWOSL(model.output,
                             signal.integral.min = 1,
                             signal.integral.max = 15,
                             background.integral.min = 601,
                             background.integral.max = 701,
                             fit.method = "EXP",
                             dose.points = c(0,10,20,50,90,0,10))

#####
## Example 4: generate sequence from *.seq file and run SAR simulation
#####

# call function "model_LuminescenceSignals", load *.seq file for sequence,
# set model = "Bailey2002" (palaeodose = 10 Gy)
# and simulate_sample_history = FALSE (default),
# because the sample history is not part of the sequence

model.output <- model_LuminescenceSignals(
  sequence = "inst/extdata/sample_SAR_cycle.SEQ",
  model = "Bailey2002",

```

```

    plot = FALSE
  )

## call RLum package function "analyse_SAR.CWOSL" to analyse the simulated SAR cycle

results <- analyse_SAR.CWOSL(model.output,
                             signal.integral.min = 1,
                             signal.integral.max = 10,
                             background.integral.min = 301,
                             background.integral.max = 401,
                             dose.points = c(0,5,10,20,50,5,0),
                             fit.method = "EXP")

print(get_RLum(results))

##=====##
## Example 5: compare different optical powers of stimulation light
##=====##

# call function "model_LuminescenceSignals", model = "Bailey2004"
# and simulate_sample_history = FALSE (default),
# because the sample history is not part of the sequence
# the optical_power of the LED is varied and then compared.

optical_power <- seq(from = 0,to = 100,by = 20)

model.output <- lapply(1:length(optical_power), function(x){

sequence <- list(IRR = c(20, 50, 1),
                 PH = c(220, 10, 5),
                 OSL = c(125, 50, optical_power[x])
                 )

return(model_LuminescenceSignals(
  sequence = sequence,
  model = "Bailey2004",
  plot = FALSE
))

})

##combine output curves
model.output.merged <- merge_RLum(model.output)

##plot
plot_RLum(
  object = model.output.merged,
  xlab = "Illumination time [s]",
  ylab = "OSL signal [a.u.]",
  main = "OSL signal dependency on optical power of stimulation light",
  legend.text = paste("Optical power density", 20*optical_power/100, "mW/cm^2"),
  combine = TRUE)

## End(Not run)

```

plot_concentrations *Plot electron/hole concentrations of a specific record.id*

Description

The functions provides a plot of all changes in time of the electron respectively hole concentration in electron traps, hole centres, in the conduction and valence band.

Usage

```
plot_concentrations(object, record.id, ...)
```

Arguments

| | |
|-----------|--|
| object | RLum.Analysis (required): S4 object of class RLum.Analysis , e.g. the values of model_LuminescenceSignals . |
| record.id | numeric (required): id of the simulated record, which is to plot. To see all record.ids use structure_RLum , see examples. |
| ... | further arguments and graphical parameters passed to plot.default and plot_RLum.Analysis . |

Details

The function produces a multiple plot output and uses in main parts the Luminescence function [plot_RLum.Analysis](#). A file output is recommended (e.g., [pdf](#)).

Value

Returns multiple plots.

Function version

0.1.0

Author(s)

Johannes Friedrich, University of Bayreuth (Germany),

References

- Bailey, R.M., 2001. Towards a general kinetic model for optically and thermally stimulated luminescence of quartz. *Radiation Measurements* 33, 17-45.
- Bailey, R.M., 2002. Simulations of variability in the luminescence characteristics of natural quartz and its implications for estimates of absorbed dose. *Radiation Protection Dosimetry* 100, 33-38.
- Bailey, R.M., 2004. Paper I-simulation of dose absorption in quartz over geological timescales and its implications for the precision and accuracy of optical dating. *Radiation Measurements* 38, 299-310.
- Pagonis, V., Chen, R., Wintle, A.G., 2007: Modelling thermal transfer in optically stimulated luminescence of quartz. *Journal of Physics D: Applied Physics* 40, 998-1006.
- Pagonis, V., Wintle, A.G., Chen, R., Wang, X.L., 2008. A theoretical model for a new dating protocol for quartz based on thermally transferred OSL (TT-OSL). *Radiation Measurements* 43, 704-708.

See Also

[plot](#), [plot_RLum.Analysis](#), [model_LuminescenceSignals](#)

Examples

```
##load data
data(ExampleData.ModelOutput, envir = environment())

##show structure
Luminescence::structure_RLum(model.output)

##plot all concentrations
plot_concentrations(object = model.output,
                    record.id = 1)

##plot only specific energy-band-level (e.g. 110 degree celsius trap, "concentration level 1")
plot_concentrations(object = model.output,
                    record.id = 1,
                    subset = list(recordType = "concentration level 1"))

##plot every level on a single plot
plot_concentrations(object = model.output,
                    record.id = 1,
                    plot.single = TRUE)
```

read_SEQ2R

Parse a Risoe SEQ-file to a sequence necessary for simulating quartz luminescence

Description

A SEQ-file created by the Risoe Sequence Editor can be imported to simulate the sequence written in the sequence editor.

Usage

```
read_SEQ2R(file, lab.dose_rate = 1, txtProgressBar = TRUE)
```

Arguments

| | |
|----------------|---|
| file | character (required) : a *.seq file created by the Risoe Sequence Editor |
| lab.dose_rate | character (with default): set the dose rate of the radiation source in the laboratory [Gy/s]. Default: 1 Gy/s |
| txtProgressBar | logical (with default): enables or disables the txtProgressBar for a visuall control of the progress. Default: txtProgressBar = TRUE |

Details**Supported versions**

Supppored and tested: version 4.36.

Value

This function returns a `list` with the parsed *.seq file and the required steps for `model_LuminescenceSignals`.

Function version

0.1.0

Author(s)

Johannes Friedrich, University of Bayreuth (Germany),

References

Riso: Sequence Editor User Manual. Available at: http://www.nutech.dtu.dk/english/-/media/Andre_Universitetsenheder/Nutech/Produkter%20og%20services/Dosimetri/radiation_measurement_instruments/tl_osl_reader/Manuals/SequenceEditor.ashx?la=da

See Also

`model_LuminescenceSignals`, `readLines`

Examples

```
##search "example_SAR_cycle.SEQ" in "extdata" in package "RLumModel"  
path <- system.file("extdata", "example_SAR_cycle.SEQ", package="RLumModel")
```

```
sequence <- read_SEQ2R(file = path)
```

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