Package 'RLumModel'

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Type Package Title Solving Ordinary Differential Equations to Understand Luminescence Version 0.2.2 Date 2017-XX-XX **Author** Johannes Friedrich [aut, trl, cre], Sebastian Kreutzer [aut, ths], Christoph Schmidt [aut, ths] Maintainer Johannes Friedrich < johannes.friedrich@uni-bayreuth.de> **Description** A collection of functions to simulate luminescence signals in the mineral quartz based on published models. Contact Package Developer Team <developer@model.r-luminescence.de> License GPL-3 **Depends** R (>= 3.3.0), utils, Luminescence (>= 0.7.0) **Imports** deSolve (>= 1.12), methods, Rcpp URL https://CRAN.R-project.org/package=RLumModel Collate RLumModel-package.R RcppExports.R calc_signal.R calc_concentrations.R create_DRT.sequence.R create_SAR.sequence.R extract_pars.R model_LuminescenceSignals.R read_SEQ2R.R set_pars.R simulate_CW_OSL.R simulate_heating.R simulate_illumination.R simulate_irradiation.R simulate_LM_OSL.R simulate_pause.R simulate_RF.R simulate_RF_and_heating.R simulate_TL.R translate_sequence.R

RoxygenNote 6.0.1

Suggests R.rsp, testthat

VignetteBuilder R.rsp

LinkingTo Rcpp, RcppArmadillo

NeedsCompilation yes

R topics documented:

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Description

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

Details

Package: RLumModel Type: Package Version: 0.2.1

Date: 2017-XX-XX License: GPL-3

Author(s)

Authors

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Project source code repository

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

Related package projects

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

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ExampleData.ModelOutput

Example data (TL curve) simulated with parameter set from Pagonis 2007

Description

Example data (TL curve) simulated with parameter set from Pagonis 2007

Format

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

Function version

0.1.1

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))
```

Author(s)

Johannes Friedrich, University of Bayreuth (Germany)

Source

model_LuminescenceSignals()

References

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.

Examples

```
data("ExampleData.ModelOutput", envir = environment())

TL_curve <- get_RLum(model.output, recordType = "TL$", drop = FALSE)

##plot TL curve
plot_RLum(TL_curve)

TL_concentrations <- get_RLum(model.output, recordType = "(TL)", drop = FALSE)
plot_RLum(TL_concentrations)</pre>
```

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 definded 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.Analysisobject 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, own_parameters = NULL,
    own_state_parameters = NULL, own_start_temperature = NULL, ...)
```

Arguments

character (required): set model to be used. Available models are: "Bailey2001", "Bailey2002", "Bailey2004", "Pagonis2007", "Pagonis2008" and "Friedrich2017".

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 concentrations.

own_parameters list (with default): This argument allows the user to submit own parameter sets. The list has to contain the following items:

- N: Concentration of electron- and hole traps [cm^(-3)]
- E: Electron/Hole trap depth [eV
- s: Frequency factor [s^(-1)]
- A: Conduction band to electron trap and valence band to hole trap transition probability [s^(-1) * cm^(3)]. CAUTION: Not every publication uses the same definition of parameter A and B! See vignette "RLumModel Usage with own parameter sets" for further details
- B: Conduction band to hole centre transition probability [s^(-1) * cm^(3)].
- Th: Photo-eviction constant or photoionisation cross section, respectively
- E_th: Thermal assistence energy [eV]
- k B: Boltzman constant 8.617e-05 [eV/K]
- W: activation energy 0.64 [eV] (for UV)
- K: 2.8e7 (dimensionless constant)
- model: "customized"
- R (optional): Ionisation rate (pair production rate) equivalent to 1 Gy/s [s^(-1) * cm^(-3)]

For further details see Bailey 2001, Wintle 1975, vignette "RLumModel - Using own parameter sets" and example 3.

own_state_parameters

numeric (with default): Some publications (e.g. Pagonis 2009) offer state parameters. With this argument the user can submit this state parameters. **Note:** You have to submit the state parameters for the conduction band and the valence band, too. For further details see vignette ""RLumModel - Using own parameter sets" and example 3.

own_start_temperature

numeric (with default): Parameter to control the start temperature (in deg. C) of a simulation. This parameter takes effect only when 'model = "customized"' is choosen.

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' [°C], 'temp end' [°C], 'heating rate' [°C/s]
OSL	optically stimulated luminescence	'temp' [°C], 'duration' [s], 'optical_power' [%]
ILL	illumination	'temp' [°C], 'duration' [s], 'optical_power' [%]

LM_OSL linear modulated OSL 'temp' [°C], 'duration' [s], optional: 'start_power' [%], 'end_power' [%] 'temp' [°C], 'dose' [Gy], 'dose_rate' [Gy/s] RL/RF radioluminescence 'temp begin' [°C], 'temp end' [°C], 'heating rate' [°C/s], 'dose_rate' [Gy/ RF_heating RF during heating/cooling 'temp' [°C], 'dose' [Gy], 'dose_rate' [Gy/s] irradiation **IRR** 'temp' [°C], optional: 'duration' [s], 'heating_rate' [°C/s] CH cutheat 'temp' [°C], 'duration' [s], optional: 'heating_rate' [°C/s] PH preheat **PAUSE** pause 'temp' [°C], 'duration' [s]

Note: 100 % illumination power equates to 20 mW/cm^2

Defining a SAR-sequence

Abrivation	Description	examples
RegDose	Dose points of the regenerative cycles [Gy]	c(0, 80, 140, 260, 320, 0, 80)
TestDose	Test dose for the SAR cycles [Gy]	50
PH	Temperature of the preheat [°C]	240
CH	Temperature of the cutheat [°C]	200
OSL_temp	Temperature of OSL read out [°C]	125
OSL_duration	Duration of OSL read out [s]	default: 40
Irr_temp	Temperature of irradiation [°C]	default: 20
PH_duration	Duration of the preheat [s]	default: 10
dose_rate	Dose rate of the laboratory irradiation source [Gy/s]	default: 1
optical_power	Percentage of the full illumination power [%]	default: 90
Irr_2recover	Dose to be recovered in a dose-recovery-test [Gy]	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

0.1.3 (2017-04-14 16:57:26)

Author(s)

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

References

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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.

Pagonis, V., Lawless, J., Chen, R., Anderson, C., 2009. Radioluminescence in Al2O3:C - analytical and numerical simulation results. Journal of Physics D: Applied Physics 42, 175107 (9pp).

Soetaert, K., Cash, J., Mazzia, F., 2012. Solving differential equations in R. Springer Science & Business Media.

Wintle, A., 1975. Thermal Quenching of Thermoluminescence in Quartz. Geophysical Journal International 41, 107-113.

See Also

```
plot, RLum, 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(</pre>
 sequence = sequence,
 model = "Bailey2001"
##get all TL concentrations
```

```
TL_conc <- get_RLum(model.output, recordType = "(TL)", drop = FALSE)</pre>
plot_RLum(TL_conc)
##plot 110 deg. C trap concentration
TL_110 <- get_RLum(TL_conc, recordType = "conc. level 1")</pre>
plot_RLum(TL_110)
##============================##
## Example 2: 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(optical_power, function(x){</pre>
sequence <- list(IRR = c(20, 50, 1),
               PH = c(220, 10, 5),
               OSL = c(125, 50, x)
               )
data <- model_LuminescenceSignals(</pre>
         sequence = sequence,
         model = "Bailey2004",
         plot = FALSE,
         verbose = FALSE
return(get_RLum(data, recordType = "OSL$", drop = FALSE))
})
##combine output curves
model.output.merged <- merge_RLum(model.output)</pre>
##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)
## Example 3: Usage of own parameter sets (Pagonis 2009)
```

```
own_parameters <- list(</pre>
 N = c(2e15, 2e15, 1e17, 2.4e16),
 E = c(0, 0, 0, 0),
 s = c(0, 0, 0, 0),
 A = c(2e-8, 2e-9, 4e-9, 1e-8),
 B = c(0, 0, 5e-11, 4e-8),
 Th = c(0, 0),
 E_{th} = c(0, 0),
 k_B = 8.617e-5,
 W = 0.64,
 K = 2.8e7,
 model = "customized",
 R = 1.7e15
## Note: In Pagonis 2009 is B the valence band to hole centre probability,
## but in Bailey 2001 this is A_j. So the values of B (in Pagonis 2009)
\#\# are A in the notation above. Also notice that the first two entries in N, A and
## B belong to the electron traps and the last two entries to the hole centres.
own_state_parameters <- c(0, 0, 0, 9.4e15)
## calculate Fig. 3 in Pagonis 2009. Note: The labels for the dose rate in the original
## publication are not correct.
## For a dose rate of 0.1 Gy/s belongs a RF signal to \sim 1.5e14 (see Fig. 6).
sequence <- list(RF = c(20, 0.1, 0.1))
model_LuminescenceSignals(
  model = "customized",
  sequence = sequence,
  own_parameters = own_parameters,
  own_state_parameters = own_state_parameters)
## Not run:
## Example 4: Simulate Thermal-Activation-Characteristics (TAC)
##set temperature
act.temp <- seq(from = 80, to = 600, by = 20)
##loop over temperature
model.output <- vapply(X = act.temp, FUN = function(x) {</pre>
##set sequence, note: sequence includes sample history
  sequence <- list(</pre>
    IRR = c(20, 1, 1e-11),
    IRR = c(20, 10, 1),
    PH = c(x, 1),
    IRR = c(20, 0.1, 1),
```

```
TL = c(20, 150, 5)
)
##run simulation
  temp <- model_LuminescenceSignals(</pre>
    sequence = sequence,
    model = "Pagonis2007",
    simulate_sample_history = TRUE,
    plot = FALSE,
    verbose = FALSE
  )
    ## "TL$" for exact matching TL and not (TL)
  TL_curve <- get_RLum(temp, recordType = "TL$")</pre>
  ##return max value in TL curve
  return(max(get_RLum(TL_curve)[,2]))
}, FUN. VALUE = 1)
##plot resutls
plot(
  act.temp[-(1:3)],
  model.output[-(1:3)],
  type = "b",
  xlab = "Temperature [\u00B0C]",
  ylab = "TL [a.u.]"
## Example 5: 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(</pre>
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(</pre>
  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,</pre>
                         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 6: generate sequence from *.seq file and run SAR simulation
##-----##
# load example *.SEQ file and construct a sequence.
# call function "model_LuminescenceSignals", load created sequence for sequence,
# set model = "Bailey2002" (palaeodose = 10 Gy)
# and simulate_sample_history = FALSE (default),
# because the sample history is not part of the sequence
path <- system.file("extdata", "example_SAR_cycle.SEQ", package="RLumModel")</pre>
sequence <- read_SEQ2R(file = path)</pre>
model.output <- model_LuminescenceSignals(</pre>
 sequence = sequence,
 model = "Bailey2001",
 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(0,8,14,26,32,0,8),
                          fit.method = "EXP")
print(get_RLum(results))
##============================##
## Example 7: Simulate sequence at laboratory without sample history
##============================##
```

12 read_SEQ2R

```
##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) Radiofluorescence at 20 deg. C with a dose of 200 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(</pre>
  sequence = sequence,
  model = "Pagonis2008"
## End(Not run)
```

read_SEQ2R

Parse a Risoe SEQ-file to a sequence neccessary 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)
```

read_SEQ2R

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 visuell con-

trol of the progress. Default: txtProgressBar = TRUE

Details

Supported versions

Suppored 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 (2017-04-14 16:57:26)
```

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)</pre>
```

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