

# Package ‘dragon’

November 5, 2020

**Title** Deep Time Redox Analysis of the Geobiology Ontology Network

**Version** 1.0.5

**Description** Create, visualize, manipulate, and analyze bipartite mineral-chemistry networks for set of focal element(s) across deep-time on Earth. The method is described in Spielman and Moore (2020) <doi:10.3389/feart.2020.585087>.

**License** GPL-3

**Imports** config, golem (>= 0.2.1), shiny, DT (>= 0.14), ggplot2 (>= 3.3.2), readr, openxlsx, dplyr (>= 1.0.0), RColorBrewer, stringr, tidyr (>= 1.0.0), purrr, tibble, broom (>= 0.5.6), cowplot (>= 1.0.0), ggforce, magrittr, shinydashboard, shinyWidgets, colourpicker (>= 1.0), colorspace (>= 1.4), visNetwork (>= 2.0.9), igraph (>= 0.4), rlang, htmltools, stats, promises, future, lubridate, xml2, rvest, curl, tidyselect

**Encoding** UTF-8

**LazyData** true

**RoxygenNote** 7.1.1

**Suggests** testthat (>= 2.1.0), processx, knitr, rmarkdown

**VignetteBuilder** knitr

**URL** <https://github.com/sjspielman/dragon>

**BugReports** <https://github.com/sjspielman/dragon/issues>

**Depends** R (>= 3.5.0)

**Collate** 'utils\_definitions.R' 'utils\_globals.R' 'utils\_names.R'  
'utils\_style-network.R' 'utils\_ui-choices.R'  
'fct\_prepare-med-data.R' 'fct\_build-legend.R'  
'fct\_build-linear-models.R' 'fct\_build-network.R'  
'fct\_build-shiny-tables.R' 'fct\_calculate-network-info.R'  
'fct\_export-network.R' 'fct\_style-network.R' 'fct\_timeline.R'  
'mod\_choose-color-palette.R'  
'mod\_choose-custom-element-colors.R' 'app\_config.R'  
'app\_server.R' 'app\_ui.R' 'run\_app.R'

**NeedsCompilation** no

**Author** Stephanie J. Spielman [aut, cre]  
(<https://orcid.org/0000-0002-9090-4788>)

**Maintainer** Stephanie J. Spielman <spielman@rowan.edu>

**Repository** CRAN

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initialize_network	<i>Initialize a mineral-chemistry network as stand-alone network rather than for embedding into the Shiny App.</i>
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### Description

Initialize a mineral-chemistry network as stand-alone network rather than for embedding into the Shiny App.

### Usage

```
initialize_network(  
  elements_of_interest,  
  force_all_elements = FALSE,  
  elements_by_redox = FALSE,  
  age_range = c(0, 5),  
  max_age_type = "Maximum",  
  cluster_algorithm = "Louvain",  
  use_data_cache = TRUE  
)
```

### Arguments

**elements\_of\_interest**  
A array of specified elements whose minerals should be included in the network. For all elements, specify "all".

**force\_all\_elements**  
A logical. If FALSE (default), minerals containing any of 'elements\_of\_interest' will be included in network. If TRUE, only minerals with full intersection of all specified elements will be included in network.

elements_by_redox	A logical. If FALSE (default), element nodes will be constructed regardless of redox state. If TRUE, creates separate node for each element's redox state, e.g. Fe <sup>2+</sup> and Fe <sup>3+</sup> would be separate nodes.
age_range	A array of two numbers giving inclusive range of mineral ages in Ga to include in network.
max_age_type	A string indicating how mineral ages should be assessed. If "Maximum" (default), filters minerals using maximum known ages at localities. If "Minimum", filters minerals using minimum known ages at localities.
cluster_algorithm	A string giving community clustering algorithm, one of "Louvain" (default) or "Leading eigenvector".
use_data_cache	A logical. If TRUE (default) cached Mineral Evolution Database will be used to build the network. If FALSE, data will be fetched from MED here. CAUTION: Requires internet connection and will take several minutes to update.

### Value

Named list containing an igraph-formatted network ('network'), an igraph::communities object giving node cluster memberships ('clustering'), a tibble of nodes associated metadata ('nodes'), and a tibble of edges and associated metadata ('edges'), and a tibble of mineral locality information ('locality\_info')

### Examples

```
## Not run:
# Include all Iron minerals whose maximum known age is between 1-2 Gya, and apply
# Louvain community clustering
initialize_network("Fe", age_range = c(1,2))

# Include all minerals containing either Iron and Oxygen whose maximum known age
# is between 1-2 Gya
initialize_network(c("Fe", "O"), age_range = c(1,2))

# Include all minerals containing both Iron and Oxygen whose maximum known age is
# between 1-2 Gya
initialize_network(c("Fe", "O"), force_all_elements = TRUE, age_range = c(1,2))

# Build the full mineral network
initialize_network("all")

## End(Not run)
```

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run\_app

*Run the "dragon" Shiny Application*

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### Description

Run the "dragon" Shiny Application

**Usage**

```
run_app()
```

**Examples**

```
## Not run:  
library(dragon)  
dragon::run_app()  
  
## End(Not run)
```

---

```
run_dragon
```

*Run the "dragon" Shiny Application. Wrapper for dragon::run\_app().*

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**Description**

Run the "dragon" Shiny Application. Wrapper for dragon::run\_app().

**Usage**

```
run_dragon()
```

**Examples**

```
## Not run:  
library(dragon)  
dragon::run_dragon()  
  
## End(Not run)
```

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