

# Package: bdXdata (via r-universe)

July 14, 2024

**Type** Package

**Title** Shared Data in Use at the CRP2A Laboratory

**Version** 0.1.0

**Description** A collection of shared data in use at the Archéosciences Bordeaux laboratory (UMR 6034 - CNRS, Université Bordeaux Montaigne, Université de Bordeaux). This package contains data specific to the laboratory instruments (calibration, monitoring, quality control, etc.). These data sets are openly distributed in order to ensure transparency and reproducibility of the results published by the laboratory team.

**License** GPL (>= 3)

**URL** <https://crp2a.github.io/bdXdata/>, <https://github.com/crp2a/bdXdata>

**BugReports** <https://github.com/crp2a/bdXdata/issues>

**Depends** R (>= 2.10)

**Encoding** UTF-8

**LazyData** true

**Roxygen** list(markdown = TRUE)

**RoxygenNote** 7.2.1

**Repository** <https://crp2a.r-universe.dev>

**RemoteUrl** <https://github.com/crp2a/bdXdata>

**RemoteRef** HEAD

**RemoteSha** 9fe9e6067eeb6b63c57e243a923f4740f36cd7e7

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**BEGe\_12cc***BEGe 12cc Calibration Data***Description**

Background and standard intensity values for 12cc BEGe.

**Usage**

```
BEGe_12cc
```

**Format**

A [list](#) of `data.frame` with 8 variables:

- `isotope` Isotope.
- `series` Radioactive decay series.
- `energy` Theoretical energy of the peak (keV).
- `bkg` Background intensity value (c/ks).
- `bkg_error` Background intensity error (c/ks).
- `std` Standard intensity value (c/ks).
- `std_error` Standard intensity error (c/ks).
- `std_name` Name of the standard.
- `mac` Mass absorption coefficient (cm<sup>2</sup>/g).

**See Also**

Other BEGe: [BEGe\\_60cc](#)

**BEGe\_60cc***BEGe 60cc Calibration Data***Description**

Background and standard intensity values for 60cc BEGe.

**Usage**

```
BEGe_60cc
```

**Format**

A [list](#) of `data.frame` with 8 variables:

`isotope` Measured isotope.  
`series` Radioactive decay series.  
`energy` Theoretical energy of the peak.  
`bkg` Background intensity value (c/ks).  
`bkg_error` Background intensity error (c/ks).  
`std` Standard intensity value (c/ks).  
`std_error` Standard intensity error (c/ks).  
`std_name` Name of the standard.  
`mac` Mass absorption coefficient (cm<sup>2</sup>/g).

**See Also**

Other BEGe: [BEGe\\_12cc](#)

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clermont

*Clermont Reference Data*

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**Description**

Clermont Reference Data

**Usage**

`clermont`

**Format**

An object of class `data.frame` with 9 rows and 15 columns.

**Source**

Guérin, G., Mercier, N. & Adamiec, G. (2011). Dose-Rate Conversion Factors: Update. *Ancient TL*, 29(1): 5-8.

Miallier, D., Guérin, G., Mercier, N., Pilleyre, T. & Sanzelle, S. (2009). The Clermont Radiometric Reference Rocks: A Convenient Tool for Dosimetric Purposes. *Ancient TL*, 27(2): 37-44.

**See Also**

Other standards: [std\\_activity](#)

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interferences                  *Interferences*

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**Description**

Interferences

**Usage**

interferences

**Format**

A [data.frame](#) with 8 observations of 6 variables:

isotope Isotope.  
energy Theoretical energy of the peak (keV).  
target  
target\_energy (keV).  
proba  
proba\_error

**See Also**

Other references: [isotopes](#), [ref\\_mac](#)

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isotopes                  *Isotopic Data*

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**Description**

Isotopic Data

**Usage**

isotopes

**Format**

A [data.frame](#) with 18 observations of 4 variables:

isotope Isotope.  
series Series  
energy Theoretical energy of the peak (keV).  
mac Mass absorption coefficient (cm<sup>2</sup>/g).  
proba\_emission

**See Also**

Other references: [interferences](#), [ref\\_mac](#)

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`ref_mac`

*Mass Absorption Coefficient*

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**Description**

Mass Absorption Coefficient

**Usage**

`ref_mac`

**Format**

A `data.frame` with 22 observations of 18 variables:

energy Energy (keV).  
hydrogen Mass absorption coefficient of hydrogen (cm<sup>2</sup>/g).  
carbon Mass absorption coefficient of carbon (cm<sup>2</sup>/g).  
nitrogen Mass absorption coefficient of nitrogen (cm<sup>2</sup>/g).  
oxygen Mass absorption coefficient of oxygen (cm<sup>2</sup>/g).  
fluorine Mass absorption coefficient of fluorine (cm<sup>2</sup>/g).  
sodium Mass absorption coefficient of sodium (cm<sup>2</sup>/g).  
magnesium Mass absorption coefficient of magnesium (cm<sup>2</sup>/g).  
aluminium Mass absorption coefficient of aluminium (cm<sup>2</sup>/g).  
silicium Mass absorption coefficient of silicium (cm<sup>2</sup>/g).  
phosphorus Mass absorption coefficient of phosphorus (cm<sup>2</sup>/g).  
sulfur Mass absorption coefficient of sulfur (cm<sup>2</sup>/g).  
chlorine Mass absorption coefficient of chlorine (cm<sup>2</sup>/g).  
potassium Mass absorption coefficient of potassium (cm<sup>2</sup>/g).  
calcium Mass absorption coefficient of calcium (cm<sup>2</sup>/g).  
titanum Mass absorption coefficient of titanum (cm<sup>2</sup>/g).  
manganese Mass absorption coefficient of manganese (cm<sup>2</sup>/g).  
iron Mass absorption coefficient of iron (cm<sup>2</sup>/g).

**See Also**

Other references: [interferences](#), [isotopes](#)

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**std\_activity**      *Characteristics of Standard Materials*

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### Description

Characteristics of Standard Materials

### Usage

`std_activity`

### Format

A `data.frame` with 2 rows and 14 variables:

`standard` Name of the standard.  
`thickness` Thickness (cm).  
`volume` Volume (cm<sup>3</sup>).  
`mass` Mass (g).  
`activity_U` Uranium specific activity (Bq/kg).  
`activity_U_error` Uranium specific activity error (Bq/kg).  
`activity_Th` Thorium specific activity (Bq/kg).  
`activity_Th_error` Thorium specific activity error (Bq/kg).  
`activity_K` Potassium specific activity (Bq/kg).  
`activity_K_error` Potassium specific activity error (Bq/kg).  
`content_U` Uranium content (ppm).  
`content_Th` Thorium content (ppm).  
`content_K` Potassium content (%).

### See Also

Other standards: [clermont](#)

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