

# toxvaldbmain

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**Type** Package

**Title** Builds the ToxValDB v9.6.2 Database

**Version** 9.6.2

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## Description

ToxValDB is a database containing quantitative records from in vivo toxicology studies from many sources. The database has 2 main parts - toxval\_source containing source data in separate tables, and the main toxval schema which combines data from multiple sources into a single format. This package moves data from toxval\_source to toxval. Data is read from files or other databases into toxval\_source and then pulled into toxval where terms are converted to standard values. This version is setup to build ToxValDB v9.6.2.

**Imports** DBI,  
RMySQL,  
openxlsx,  
dplyr,  
tidyr,  
stringr,  
tibble,  
janitor,  
XML,  
miniUI,  
RCurl,  
gsubfn,  
textclean,  
magrittr

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**Encoding** UTF-8

**LazyData** true

**RoxygenNote** 7.3.2

**Suggests** knitr,  
rmarkdown

**VignetteBuilder** knitr

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---

|                     |                     |
|---------------------|---------------------|
| <i>cas_checkSum</i> | <i>cas_checkSum</i> |
|---------------------|---------------------|

---

**Description**

Check CAS RN validity via checksum method  
For a suspected CAS RN, determine validity by calculating final digit checksum

**Usage**

`cas_checkSum(x)`

**Arguments**

`x` chr. Input vector of values to check. Standard CAS notation using hyphens is fine, as #' all non-digit characters are stripped for checksum calculation. Each element of `x` should contain #' only one suspected CAS RN to check.

**Details**

This function performs a very specific type of check for CAS validity, namely whether the final digit checksum follows the CAS standard. By default, it also ensures that the digit length is compatible with CAS standards.

**Value**

A logical vector of length `x` denoting whether each `x` is a valid CAS by the checksum method. NA input values will be set 0 (FAIL).

**See Also**

[str\\_detect](#), [str\\_pad](#)

**Examples**

```
cas_good <- c("71-43-2", "18323-44-9", "7732-18-5") # benzene, clindamycin, water
cas_bad  <- c("61-43-2", "18323-40-9", "7732-18-4") # single digit change from good
cas_checkSum(c(cas_good, cas_bad))
```

---

```
check.toxval_type.route.units
```

*Output distinct combinations of toxval\_type, exposure\_route, and toxval\_units to check*

---

## Description

Output distinct combinations of toxval\_type, exposure\_route, and toxval\_units to check

## Usage

```
check.toxval_type.route.units(
    toxval.db,
    source = NULL,
    subsource = NULL,
    load.dict = NULL
)
```

## Arguments

|           |   |
|-----------|---|
| toxval.db | The version of toxvaldb to use.   |
| source    | Source to check   |
| subsource | Subsource to check (NULL default)   |
| load.dict | Name of dictionary containing expected combinations, or NULL if dictionary should not be read |

---

```
chem.check.v2
```

*Check the chemicals from a file Names with special characters are cleaned and trimmed CASRN are fixed (dashes put in, trimmed) and check sums are calculated The output is sent to a file called chemcheck.xlsx in the source data file One option for using this is to edit the source file until no errors are found*

---

## Description

Check the chemicals from a file Names with special characters are cleaned and trimmed CASRN are fixed (dashes put in, trimmed) and check sums are calculated The output is sent to a file called chemcheck.xlsx in the source data file One option for using this is to edit the source file until no errors are found

## Usage

```
chem.check.v2(res0, in_source = NULL, verbose = FALSE)
```

**Arguments**

|           |  |
|-----------|--|
| res0      | The data frame in which chemicals names and CASRN will be replaced |
| in_source | The source to be processed. If source=NULL, process all sources    |
| verbose   | If TRUE, print diagnostic messages                                 |

**Value**

Return a list with fixed CASRN and name and flags indicating if fixes were made: res0=res0,name.OK=name.OK,casrn.OK=c

---

|                      |                             |
|----------------------|-----------------------------|
| clean.last.character | <i>clean.last.character</i> |
|----------------------|-----------------------------|

---

**Description**

Clean unneeded characters from the end of a string

**Usage**

```
clean.last.character(x)
```

**Arguments**

|   |                      |
|---|----------------------|
| x | String to be cleaned |
|---|----------------------|

**Details**

DETAILS

**Value**

The cleaned string

**See Also**

[str\\_trim](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

`clean.toxval.by.source`*Delete a portion of the contents of the toxval database*

---

**Description**

Delete a portion of the contents of the toxval database

**Usage**

```
clean.toxval.by.source(toxval.db, source)
```

**Arguments**

|                        |   |
|------------------------|---|
| <code>toxval.db</code> | The version of toxval from which the data is deleted. |
| <code>source</code>    | The data source name                                  |

**Value**

The database will be altered

---

`compare.versions`*Compare versions of toxval*

---

**Description**

Compare versions of toxval

**Usage**

```
compare.versions(db1, db2)
```

**Arguments**

|                  |                                 |
|------------------|---------------------------------|
| <code>db1</code> | The old version of the database |
| <code>db2</code> | The new version of the database |



---

|                  |                            |
|------------------|----------------------------|
| duplicate.hunter | <i>Diagnose duplicates</i> |
|------------------|----------------------------|

---

**Description**

Diagnose duplicates

**Usage**

```
duplicate.hunter(toxval.db, source = NULL, source_table = NULL)
```

**Arguments**

|           |                          |
|-----------|--------------------------|
| toxval.db | Database version         |
| source    | The source to be updated |

**Value**

Write a file with the results

---

|                              |                                     |
|------------------------------|-------------------------------------|
| ecotox.select.study.duration | <i>ECOTOX Select study_duration</i> |
|------------------------------|-------------------------------------|

---

**Description**

Function to select the appropriate study\_duration value from conc1\_\* fields

**Usage**

```
ecotox.select.study.duration(in_data, dur_col = NULL)
```

**Arguments**

|         |                        |
|---------|------------------------|
| in_data | Input ECOTOX dataframe |
|---------|------------------------|

**Value**

Processed dataframe with new study\_duration\_values, units, and qualitiifer fields

```
ecotox.select.toxval.numeric
```

*ECOTOX Select toxval\_numeric*

---

**Description**

Function to select the appropriate toxval\_numeric value from conc1\_\* fields

**Usage**

```
ecotox.select.toxval.numeric(in_data)
```

**Arguments**

|         |                        |
|---------|------------------------|
| in_data | Input ECOTOX dataframe |
|---------|------------------------|

**Value**

Processed dataframe with new toxval\_numeric, units, and qualitiifer fields

---

```
ecotox.species.dictionary
```

*Extract the ECOTOX species dictionary from the ECOTOX data*

---

**Description**

Extract the ECOTOX species dictionary from the ECOTOX data

**Usage**

```
ecotox.species.dictionary(toxval.db, do.load = F, sys.date = "2023-05-03")
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | The version of toxval into which the tables are loaded.                   |
| do.load   | If TRUE, load the data from the input file and put into a global variable |
| source.db | The version of toxval source - used to manage chemicals                   |
| verbose   | Whether the loaded rows should be printed to the console.                 |
| log       | If TRUE, send output to a log file  |

---

`export.all.by.source`    *Build a data frame of the data from toxval and export by source as a series of xlsx files*

---

### Description

Build a data frame of the data from toxval and export by source as a series of xlsx files

### Usage

```
export.all.by.source(
  toxval.db,
  source = NULL,
  subsources = NULL,
  include.qc.status = TRUE
)
```

### Arguments

|                                |   |
|--------------------------------|---|
| <code>toxval.db</code>         | Database version  |
| <code>source</code>            | The source to be updated  |
| <code>subsources</code>        | The subsources to be updated  |
| <code>include.qc.status</code> | Boolean whether to include the <code>qc_status</code> field, or filter out "fail" records. Default is TRUE. #' @return for each source writes an Excel file with the name <code>../export/export_by_source_data/toxval_all_toxval.db_source.xlsx</code> |

---

`export.delete.qc_status.fail.by.source`  
*export.delete.qc\_status.fail.by.source*

---

### Description

Function to export log of `qc_status` "fail" records and delete fail records from ToxValDB.

### Usage

```
export.delete.qc_status.fail.by.source(toxval.db, source, subsources)
```

### Arguments

|                         |  |
|-------------------------|--|
| <code>toxval.db</code>  | The database version to use                                    |
| <code>source</code>     | The source name  |
| <code>subsources</code> | The specific subsources to process, if desired (Default: NULL) |

**Value**

None. XLSX log generated and SQL UPDATE statements are pushed to the database.

---

```
export.for.missing.species
```

*Export the records with missing species and point of departure values*

---

**Description**

Export the records with missing species and point of departure values

**Usage**

```
export.for.missing.species(toxval.db, source = NULL)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | Database version  |
| source    | The source to be updated #' @return for each source writes an Excel file with the name ../export/export_by_source_data/toxval_all_toxval.db_source.xlsx |

---

```
export.for.oppt
```

*Build a data frame of the PODs and exports as xlsx*

---

**Description**

Build a data frame of the PODs and exports as xlsx

**Usage**

```
export.for.oppt(toxval.db, file.name = "TSCA PICS")
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | Database version  |
| file.name | If not NA, this is a file containing chemicals, and only those chemicals will be exported |

**Value**

writes an Excel file with the name ../export/toxval\_pod\_summary\_[human\_eco]\_Sys.Date().xlsx

---

export.for.toxval.qc    *Build a data frame of the PODs and exports as xlsx*

---

**Description**

Build a data frame of the PODs and exports as xlsx

**Usage**

```
export.for.toxval.qc(toxval.db, source = NULL)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | Database version  |
| human_eco | Either 'human health' or 'eco'  |
| file.name | If not NA, this is a file containing chemicals, and only those chemicals will be exported |

**Value**

writes an Excel file with the name ../export/toxval\_pod\_summary\_[human\_eco]\_Sys.Date().xlsx

---

export.for.toxvaldb.manuscript  
                          *Build a data frame of the data for the toxval manuscript*

---

**Description**

Build a data frame of the data for the toxval manuscript

**Usage**

```
export.for.toxvaldb.manuscript(toxval.db)
```

**Arguments**

|           |                          |
|-----------|--------------------------|
| toxval.db | Database version         |
| source    | The source to be updated |

**Value**

Write a file with the results

---

```
export.input.dict.file.list
```

*export.input.dict.file.list*

---

**Description**

Pull list of directories, files, and dictionaries used throughout the ToxValDB workflow.

**Usage**

```
export.input.dict.file.list()
```

**Value**

Dataframe with "folder\_name" field of files and folders in the toxval.config()\$datapath directory.

**See Also**

[read\\_excel](#) [write\\_xlsx](#) [filter](#), [pull](#), [bind\\_rows](#), [distinct](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

```
export.missing.dictionary.entries
```

*Find "original" values that have not been included in the dictionaries and export them*

---

**Description**

Find "original" values that have not been included in the dictionaries and export them

**Usage**

```
export.missing.dictionary.entries(
  toxval.db,
  source = NULL,
  subsources = NULL,
  report.only = FALSE
)
```

**Arguments**

|             |  |
|-------------|--|
| toxval.db   | The version of toxval in which the data is altered.                          |
| source      | The source to be fixed. If source=NULL, fix all sources                      |
| subsource   | The subsource to be fixed (NULL default)                                     |
| report.only | Whether to report or write/export data. Default is FALSE (write/export data) |

**Value**

An excel file in dictionaries with the missing entries (if report.only=TRUE, return tibble) "missing dictionary entries Sys.Date.xlsx"

---

|   |
|---|
| export.missing.strains                      |
| <i>Set the strain information in toxval</i> |

---

**Description**

Set the strain information in toxval

**Usage**

```
export.missing.strains(toxval.db, date_string = "2024-02-27")
```

**Arguments**

|             |   |
|-------------|---|
| toxval.db   | The version of the database to use        |
| date_string | The date of the latest dictionary version |

---

|   |
|---|
| export.missing.toxval_type  |
| <i>Export any toxval_types that are not in the toxval_type dictionary</i> |

---

**Description**

Export any toxval\_types that are not in the toxval\_type dictionary

**Usage**

```
export.missing.toxval_type(toxval.db, report.only = FALSE)
```

**Arguments**

|             |  |
|-------------|--|
| toxval.db   | The version of toxval in which the data is altered.                          |
| report.only | Whether to report or write/export data. Default is FALSE (write/export data) |

**Value**

An excel file in dictionaries with the missing entries (if report.only=TRUE, return tibble) "dictionary/missing/missing\_toxval\_type Sys.Date.xlsx"

---

```
export.toxvaldb.subset
```

*Build a data frame of the data from toxval for a subset of chemicals*

---

**Description**

Build a data frame of the data from toxval for a subset of chemicals

**Usage**

```
export.toxvaldb.subset(toxval.db, filename)
```

**Arguments**

|           |  |
|-----------|--|
| toxval.db | Database version   |
| filename  | The name of the file the be imported - should be a short name that will be used in the output filename. This is an xlsx file and needs a column labeled dtxsid |

**Value**

Write a file with the results

---

```
export.update.source.info
```

*export.update.source.info*

---

**Description**

Pull updated field information for source info dictionary

**Usage**

```
export.update.source.info(toxval.db, source.db, dict_date = "2024-08-28")
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | Database version                            |
| source.db | The source to be updated                    |
| dict_date | Date stamp of the dictionary file to update |

**Value**

Overwrite existing source info dictionary with updated fields



---

`fill.chemical.by.source`*Fill the chemical table*

---

**Description**

Fill the chemical table

**Usage**

```
fill.chemical.by.source(toxval.db, source, verbose = TRUE)
```

**Arguments**

|                        |  |
|------------------------|--|
| <code>toxval.db</code> | The version of toxvaldb to use.              |
| <code>source</code>    | The source to be used                        |
| <code>verbose</code>   | If TRUE, print out extra diagnostic messages |

---

`fill.toxval.defaults` *Set Toxval Defaults*

---

**Description**

Set Toxval Defaults

**Usage**

```
fill.toxval.defaults(toxval.db, mat)
```

**Arguments**

|                        |   |
|------------------------|---|
| <code>toxval.db</code> | The version of toxval from which to set defaults. |
| <code>mat</code>       | An input matrix of data                           |

**Value**

The data matrix after fixing

---

```
fill.toxval.defaults.global.by.source
```

*Set Toxval Defaults globally, replacing blanks with -*

---

### Description

Set Toxval Defaults globally, replacing blanks with -

### Usage

```
fill.toxval.defaults.global.by.source(
    toxval.db,
    source = NULL,
    subsources = NULL
)
```

### Arguments

|            |   |
|------------|---|
| toxval.db  | The version of toxval from which to set defaults. |
| source     | The source to be fixed                            |
| subsources | The subsources to be fixed (NULL default)         |

---

```
fix.all.param.by.source
```

*Alter the contents of toxval according to an excel dictionary file with fields - exposure\_method, exposure\_route, sex, strain, study\_duration\_class, study\_duration\_units, study\_type, toxval\_type, exposure\_form, toxval\_subtype*

---

### Description

Alter the contents of toxval according to an excel dictionary file with fields - exposure\_method, exposure\_route, sex, strain, study\_duration\_class, study\_duration\_units, study\_type, toxval\_type, exposure\_form, toxval\_subtype

### Usage

```
fix.all.param.by.source(
    toxval.db,
    source = NULL,
    subsources = NULL,
    fill.toxval_fix = TRUE
)
```

**Arguments**

|                 |   |
|-----------------|---|
| toxval.db       | The version of toxval in which the data is altered.               |
| source          | The source to be fixed. If source=NULL, fix all sources           |
| subsource       | The subsource to be fixed (NULL default)                          |
| fill.toxval_fix | If TRUE (default) read the dictionaries into the toxval_fix table |

**Value**

The database will be altered

---

fix.associated.pod.critical\_effect

*Appends associated POD information to critical\_effect for derived toxval\_types*

---

**Description**

Appends associated POD information to critical\_effect for derived toxval\_types

**Usage**

```
fix.associated.pod.critical_effect(res, map_fields)
```

**Arguments**

|            |   |
|------------|---|
| res        | The data to be altered                            |
| map_fields | The fields used to map entries to associated PODs |

---

fix.casrn

*fix.casrn*

---

**Description**

Fix a CASRN that has one of several problems

**Usage**

```
fix.casrn(casrn, cname = "", verbose = FALSE)
```

**Arguments**

|         |                                 |
|---------|---------------------------------|
| casrn   | Input CASRN to be fixed         |
| cname   | An optional chemical name       |
| verbose | if TRUE, print the input values |

**Details**

DETAILS

**Value**

the fixed CASRN

**See Also**[reexports](#)**Examples**

```
## Not run:
if(interactive()){
  fix.casrn("107028")
  # Expected output "107-02-8"
}

## End(Not run)
```

---

```
fix.critical_effect.icf.by.source
```

*standardize critical\_effect in toxval table based on icf dictionary and  
toxval critical effects dictionary*

---

**Description**

standardize critical\_effect in toxval table based on icf dictionary and toxval critical effects dictionary

**Usage**

```
fix.critical_effect.icf.by.source(toxval.db, source, subsource = NULL)
```

**Arguments**

|           |  |
|-----------|--|
| toxval.db | The version of toxvaldb to use.          |
| source    | The source to be fixed                   |
| subsource | The subsource to be fixed (NULL default) |

---

```
fix.dedup.hierarchy.by.source
```

*Set qc\_status as "fail" for lesser priority duplicates*

---

## Description

Set qc\_status as "fail" for lesser priority duplicates

## Usage

```
fix.dedup.hierarchy.by.source(
  toxval.db,
  source = NULL,
  subsources = NULL,
  priority_list = NULL,
  subsources_priority_list = NULL,
  criteria = c("dtxsid"),
  report.only = FALSE
)
```

## Arguments

|                          |  |
|--------------------------|--|
| toxval.db                | The version of toxvaldb to use.  |
| source                   | Source to be fixed   |
| subsources               | Subsources to be fixed (NULL default)  |
| priority_list            | Named list describing source priority, with low priority index and high priority value |
| subsources_priority_list | Named list of subsources to deprecate per source, source index/subsource value         |
| criteria                 | List of parameters used to make deduping decisions                                     |
| report.only              | Whether to report or write/export data. Default is FALSE (write/export data)           |

---

```
fix.derived.toxval_type.by.source
```

*Set select normalized toxval fields to '-' if the record is a select toxval\_type*

---

## Description

Set select normalized toxval fields to '-' if the record is a select toxval\_type

**Usage**

```
fix.derived.toxval_type.by.source(  
  toxval.db,  
  source = NULL,  
  subsources = NULL,  
  report.only = FALSE  
)
```

**Arguments**

|             |   |
|-------------|---|
| toxval.db   | The version of the database to use                                    |
| source      | The source to be fixed  |
| subsources  | The subsources to be fixed (NULL default)                             |
| report.only | Whether to update database or only report the results. Default FALSE. |

---

fix.empty.by.source    *Set all empty cells in toxval to '-'*

---

**Description**

Set all empty cells in toxval to '-'

**Usage**

```
fix.empty.by.source(toxval.db, source = NULL, subsources = NULL)
```

**Arguments**

|            |   |
|------------|---|
| toxval.db  | The version of toxval in which the data is altered. |
| source     | The source to be fixed                              |
| subsources | The subsources to be fixed (NULL default)           |

**Value**

The database will be altered

---

```
fix.empty.record_source.by.source
```

*Set all empty cells in record\_source to '-'*

---

### Description

Set all empty cells in record\_source to '-'

### Usage

```
fix.empty.record_source.by.source(toxval.db, source = NULL)
```

### Arguments

|           |   |
|-----------|---|
| toxval.db | The version of toxval in which the data is altered. |
| source    | The source to be fixed                              |

### Value

The database will be altered

---

```
fix.exposure.params
```

*Fix the exposure fields: exposure\_method, exposure\_route, exposure\_form based on a 3 column dictionary ~/dictionary/exposure\_route\_method\_form.xlsx*

---

### Description

Fix the exposure fields: exposure\_method, exposure\_route, exposure\_form based on a 3 column dictionary ~/dictionary/exposure\_route\_method\_form.xlsx

### Usage

```
fix.exposure.params(
  toxval.db,
  source = NULL,
  subsource = NULL,
  report.only = FALSE
)
```

**Arguments**

|                              |  |
|------------------------------|--|
| <code>toxval.db</code>       | The version of toxval in which the data is altered.                            |
| <code>source</code>          | The source to be fixed. If <code>source=NULL</code> , fix all sources          |
| <code>subsource</code>       | The subsource to be fixed (NULL default)                                       |
| <code>report.only</code>     | Whether to report or write/export data. Default is FALSE (write/export data)   |
| <code>fill.toxval_fix</code> | If TRUE (default) read the dictionaries into the <code>toxval_fix</code> table |

**Value**

The database will be altered (if `report.only=TRUE`, return missing entries)

---

```
fix.exposure_route.not_specified.by.source
      fix.exposure_route.not_specified.by.source
```

---

**Description**

Function to use a dictionary file to update cases where `exposure_route` reported as "not specified".

**Usage**

```
fix.exposure_route.not_specified.by.source(toxval.db, source, subsource)
```

**Arguments**

|                        |   |
|------------------------|---|
| <code>toxval.db</code> | The database version to use                                   |
| <code>source</code>    | The source name   |
| <code>subsource</code> | The specific subsource to process, if desired (Default: NULL) |

**Details**

DETAILS

**Value**

None. SQL UPDATE statements are pushed to the database.

**See Also**

[pull](#), [mutate-joins](#), [select](#), [rename](#), [mutate](#), [filter](#), [distinct](#) [read\\_excel](#) [write\\_xlsx](#)



**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

```
fix.generation.by.source
```

*Alter the contents of toxval according to an excel dictionary file with field generation*

---

**Description**

Alter the contents of toxval according to an excel dictionary file with field generation

**Usage**

```
fix.generation.by.source(toxval.db, source, subsource = NULL)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | The version of toxval in which the data is altered. |
| source    | The source to be processed                          |
| subsource | The subsource to be processed (NULL default)        |

**Value**

The database will be altered

---

```
fix.human_eco.by.source
```

*Fix the human\_eco flag*

---

**Description**

Fix the human\_eco flag

**Usage**

```
fix.human_eco.by.source(toxval.db, source = NULL, subsource = NULL)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | The version of toxval in which the data is altered. |
| source    | The source to be fixed. If NULL, fix all sources    |
| subsource | The subsource to be fixed (NULL default)            |

**Value**

The database will be altered

---

|                               |                                      |
|-------------------------------|--------------------------------------|
| <code>fix.non_ascii.v2</code> | <code><i>fix.non_ascii.v2</i></code> |
|-------------------------------|--------------------------------------|

---

**Description**

Flag and fix non-ascii characters in the database

**Usage**

```
fix.non_ascii.v2(df, source)
```

**Arguments**

|        |                               |
|--------|-------------------------------|
| df     | The dataframe to be processed |
| source | Current ToxVal source         |
| The    | source to be fixed            |

**Details**

DETAILS

**Value**

The dataframe with non ascii characters replaced with cleaned versions

**See Also**

[read.xlsx](#), [write.xlsx](#) [str\\_trim](#) [stri\\_escape\\_unicode](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

```
fix.priority_id.by.source
```

*Fix the priority\_id in the toxval table based on source*

---

### Description

Fix the priority\_id in the toxval table based on source

### Usage

```
fix.priority_id.by.source(toxval.db, source = NULL, subsource = NULL)
```

### Arguments

|           |  |
|-----------|--|
| toxval.db | The version of toxvaldb to use.                      |
| source    | The source to be fixed, If NULL, set for all sources |
| subsource | The subsource to be fixed (NULL default)             |

---

```
fix.qc_status.by.source
```

*Fix the qa\_status flag*

---

### Description

Fix the qa\_status flag

### Usage

```
fix.qc_status.by.source(
  toxval.db,
  source.db,
  source = NULL,
  subsource = NULL,
  reset = FALSE
)
```

### Arguments

|           |   |
|-----------|---|
| toxval.db | The version of toxval in which the data is altered. |
| source    | The source to be fixed                              |
| subsource | The subsource to be fixed (NULL default)            |
| reset     | If TRUE, reset all values to 'pass' before setting  |
| sourcedb  | The source database name                            |

**Value**

The database will be altered

---

|                     |                            |
|---------------------|----------------------------|
| fix.replace.unicode | <i>fix.replace.unicode</i> |
|---------------------|----------------------------|

---

**Description**

A function to check all character fields and handle unicode symbols, either by removing them or replacing them with alphabetic equivalents.

**Usage**

```
fix.replace.unicode(df)
```

**Arguments**

|    |  |
|----|--|
| df | Character vector to check/replace unicode symbols. |
|----|--|

**Details**

DETAILS

**Value**

Returns a modified version of the input vector with unicode replacements.

**See Also**

[stri\_escape\_unicode][stringi::stri\_escape\_unicode] [str\_extract][stringr::str\_extract]

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

```
fix.risk_assessment_class.by.source
```

*Set the risk assessment class of toxval according to an excel dictionary.  
Values may beset multiple times, so the excel sheet should be ordered  
so that the last ones to be set are last*

---

### Description

Set the risk assessment class of toxval according to an excel dictionary. Values may beset multiple times, so the excel sheet should be ordered so that the last ones to be set are last

### Usage

```
fix.risk_assessment_class.by.source(  
    toxval.db,  
    source = NULL,  
    subsource = NULL,  
    restart = TRUE,  
    report.only = FALSE  
)
```

### Arguments

|             |  |
|-------------|--|
| toxval.db   | The version of toxval in which the data is altered.                          |
| source      | The source to be updated   |
| subsource   | The subsource to be updated (NULL default)                                   |
| restart     | If TRUE, delete all values and start from scratch                            |
| report.only | Whether to report or write/export data. Default is FALSE (write/export data) |

---

```
fix.single.param.by.source
```

*Alter the contents of toxval according to an excel dictionary*

---

### Description

Alter the contents of toxval according to an excel dictionary

### Usage

```
fix.single.param.by.source(  
    toxval.db,  
    param,  
    source,  
    subsource = NULL,
```

```
ignore = FALSE,  
report.only = FALSE,  
units.data = NULL  
)
```

**Arguments**

|             |  |
|-------------|--|
| toxval.db   | The version of toxval in which the data is altered.                                  |
| param       | The parameter value to be fixed  |
| source      | The source to be fixed   |
| subsource   | The subsource to be fixed (NULL default)   |
| ignore      | If TRUE allow missing values to be ignored   |
| report.only | Whether to report or write/export data. Default is FALSE (write/export data)         |
| units.data  | A dataframe containing current units data if units are to be reported (NULL default) |

**Value**

The database will be altered

---

|   |
|---|
| fix.species.common_name                     |
| <i>Fix issues with species common names</i> |

---

**Description**

Fix issues with species common names

**Usage**

```
fix.species.common_name(toxval.db)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | The version of toxval into which the tables are loaded. |
|-----------|---|

---

`fix.species.duplicates`*Check for species duplicates - same common name but multiple species\_ids*

---

**Description**

Check for species duplicates - same common name but multiple species\_ids

**Usage**

```
fix.species.duplicates(toxval.db, source = NULL, subsource = NULL)
```

**Arguments**

|                        |   |
|------------------------|---|
| <code>toxval.db</code> | The version of the database to use                          |
| <code>source</code>    | The source to be checked. If source=NULL, check all sources |
| <code>subsource</code> | The subsource to be checked (NULL default)                  |

---

`fix.species.v2`      *Set the species\_id column in toxval*

---

**Description**

This function replaces fix.species This function precedes toxvaldb.load.species

**Usage**

```
fix.species.v2(  
  toxval.db,  
  source = NULL,  
  subsource = NULL,  
  date_string = "2023-05-18"  
)
```

**Arguments**

|                          |  |
|--------------------------|--|
| <code>toxval.db</code>   | The version of the database to use       |
| <code>source</code>      | The source to be fixed                   |
| <code>subsource</code>   | The subsource to be fixed (NULL default) |
| <code>date_string</code> | The date version of the dictionary       |

---

|               |                                      |
|---------------|--------------------------------------|
| fix.strain.v2 | Set the strain information in toxval |
|---------------|--------------------------------------|

---

**Description**

Set the strain information in toxval

**Usage**

```
fix.strain.v2(  
  toxval.db,  
  source = NULL,  
  subsources = NULL,  
  date_string = "2024-04-08",  
  reset = FALSE  
)
```

**Arguments**

- |             |  |
|-------------|--|
| toxval.db   | The version of the database to use                   |
| source      | The source to be fixed. If NULL, fix for all sources |
| subsource   | The subsource to be fixed (NULL default)             |
| date_string | The date of the latest dictionary version            |

---

|                           |  |
|---------------------------|--|
| fix.study_duration.params | <i>Fix the study duration fields: study_duration_value, study_duration_units, study_duration_class based on a 3 column dictionary ~/dictionary/exposure_route_method_form.xlsx</i> |
|---------------------------|--|

---

**Description**

Fix the study duration fields: study\_duration\_value, study\_duration\_units, study\_duration\_class based on a 3 column dictionary ~/dictionary/exposure\_route\_method\_form.xlsx

**Usage**

```
fix.study_duration.params(  
  toxval.db,  
  source = NULL,  
  subsources = NULL,  
  report.only = FALSE  
)
```



**Arguments**

|                 |  |
|-----------------|--|
| toxval.db       | The version of toxval in which the data is altered.                          |
| source          | The source to be fixed. If source=NULL, fix all sources                      |
| subsource       | The subsource to be fixed (NULL default)                                     |
| report.only     | Whether to report or write/export data. Default is FALSE (write/export data) |
| fill.toxval_fix | If TRUE (default) read the dictionaries into the toxval_fix table            |

**Value**

The database will be altered

---

|                 |                                  |
|-----------------|----------------------------------|
| fix.study_group | <i>Set the study_group field</i> |
|-----------------|----------------------------------|

---

**Description**

Set the study\_group field

**Usage**

```
fix.study_group(
  toxval.db,
  source = NULL,
  subsource = NULL,
  report.only = FALSE
)
```

**Arguments**

|             |  |
|-------------|--|
| toxval.db   | Database version   |
| source      | The source to be updated   |
| subsource   | The subsource to be updated (NULL default)                                   |
| report.only | Whether to apply study_group fix or just report what fixes would be applied. |

**Value**

for each source writes an Excel file with the name ../export/export\_by\_source\_data/toxval\_all\_toxval.db\_source.xlsx

---

fix.study\_type.by.source

*Fix the study\_type using manual curation on a source-by-source basis*


---

## Description

This function replaces the original `export.for_study_type` and `fix.study_type.manual`, with the intention of making it easier to fix the study types on a source-by-source basis. All of the work will happen in the directory `~/Repo/dictionary/study_type_by_source`. Each source will have its own file and will not have a date attached to make maintenance easier. To start the process, run this with `mode="export"`. This will write a source-specific file to the `export_temp` directory. Open either the `xlsx` or `csv` (if the `xlsx` is corrupted) and place this file into the main directory (`study_type_by_source`) and edit it there as documented in the main documentation. Next run this function with `mode="import"`. This will load your changes into the database. It is suggested that before working on a new source that the old version in the `study_type_by_source` get pushed to the `old_versions` directory.

## Usage

```
fix.study_type.by.source(
    toxval.db,
    mode = "export",
    source = NULL,
    subsources = NULL,
    custom.query.filter = NULL,
    report.only = FALSE
)
```

## Arguments

|                                  |   |
|----------------------------------|---|
| <code>toxval.db</code>           | The version of <code>toxval</code> in which the data is altered.  |
| <code>mode</code>                | Either <code>export</code> or <code>import</code>   |
| <code>source</code>              | The source you want to work on. If <code>NULL</code> , this will run all sources  |
| <code>subsources</code>          | The subsources to be fixed  |
| <code>custom.query.filter</code> | Additional filters for the query. Example: <code>custom.query.filter = paste0(" and b.human_eco='human health' and ", "e.toxval_type_supercategory in ('Point of Departure','Lethality Effect Level','Toxicity Value'))"</code> |
| <code>report.only</code>         | Whether to report or write/export data. Default is <code>FALSE</code> (write/export data)   |

## Value

The database will be altered

---

|                 |  |
|-----------------|--|
| fix.trim_spaces | <i>Trim leading and trailing blanks from all character columns</i> |
|-----------------|--|

---

**Description**

Trim leading and trailing blanks from all character columns

**Usage**

```
fix.trim_spaces(res)
```

**Arguments**

|                 |   |
|-----------------|---|
| toxval.db       | The version of toxval in which the data is altered.               |
| source          | The source to be fixed. If source=NULL, fix all sources           |
| fill.toxval_fix | If TRUE (default) read the dictionaries into the toxval_fix table |

**Value**

The database will be altered

---

|                     |                                     |
|---------------------|-------------------------------------|
| fix.units.by.source | <i>Do all of the fixes to units</i> |
|---------------------|-------------------------------------|

---

**Description**

1. All of these steps operate on the toxval\_units column.
2. Replace variant unit names with standard ones, running fix.single.param.new.by.source.R This fixes issues like variant names for mg/kg-day and uses the dictionary file dictionary/toxval\_units\_5.xlsx
3. Fix special characters in toxval\_units
4. Fix issues with units containing extra characters for some ECOTOX records
5. Convert units that are multiples of standard ones (e.g. ppb to ppm). This uses the dictionary file dictionary/toxval\_units conversions 2018-09-12.xlsx
6. Run conversions from molar to mg units, using MW. This uses the dictionary file dictionary/MW conversions.xlsx
7. Convert ppm to mg/m3 for inhalation studies. This uses the conversion Concentration (mg/m3) = 0.0409 x concentration (ppm) x molecular weight. See <https://cfpub.epa.gov/ncer/abstracts/index.cfm/fuseaction/dispatch> This function requires that the DSSTox external chemical\_id be set
8. Convert ppm to mg/kg-day in toxval according to a species-specific conversion factor for oral exposures. This uses the dictionary file dictionary/ppm to mgkgday by animal.xlsx See: [www10.plala.or.jp/biostatistics/1-3.doc](http://www10.plala.or.jp/biostatistics/1-3.doc) This probably assumes feed rather than water
9. Make sure that eco studies are in mg/L and human health in mg/m3

**Usage**

```
fix.units.by.source(
  toxval.db,
  source = NULL,
  subsources = NULL,
  do.convert.units = FALSE,
  report.only = FALSE,
  report.extra = FALSE
)
```

**Arguments**

|                               |   |
|-------------------------------|---|
| <code>toxval.db</code>        | The version of toxvaldb to use.   |
| <code>source</code>           | Source to be fixed  |
| <code>subsources</code>       | Subsources to be fixed (NULL default)   |
| <code>do.convert.units</code> | If TRUE, so unit conversions, as opposed to just cleaning   |
| <code>report.only</code>      | Whether to report or write/export data. Default is FALSE (write/export data)  |
| <code>report.extra</code>     | If reporting, then choose whether to record extra conversion information (e.g. <code>toxval_type</code> , <code>mw</code> , <code>species_id</code> , etc.) |

---

|                                 |   |
|---------------------------------|---|
| <code>generate.originals</code> | <i>Duplicate any columns with '_original' Set Toxval Defaults</i> |
|---------------------------------|---|

---

**Description**

Duplicate any columns with '\_original' Set Toxval Defaults

**Usage**

```
generate.originals(toxval.db, mat)
```

**Arguments**

|                        |   |
|------------------------|---|
| <code>toxval.db</code> | The version of toxval from which to set defaults. |
| <code>mat</code>       | The matrix of data to be altered                  |

**Value**

The altered input matrix

---

`getDBConn`*getDBConn*

---

**Description**

Get the names the database server, user, and pass or returns error message

**Usage**

```
getDBConn()
```

**Details**

DETAILS

**Value**

print the database connection information

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

`import.dictionary`*import the toxval and toxval\_type dictionaries*

---

**Description**

import the toxval and toxval\_type dictionaries

**Usage**

```
import.dictionary(toxval.db)
```

**Arguments**

`toxval.db`      The name of the database

---

|   |   |
|---|---|
| <code>import.source.info.by.source</code> | <i>Load Source Info for each source into toxval The information is in the file ~/dictionary/source_info 2023-11-30.xlsx</i> |
|---|---|

---

**Description**

Load Source Info for each source into toxval The information is in the file ~/dictionary/source\_info 2023-11-30.xlsx

**Usage**

```
import.source.info.by.source(  
    toxval.db,  
    source = NULL,  
    dict_date = "2025-05-16"  
)
```

**Arguments**

|                        |   |
|------------------------|---|
| <code>toxval.db</code> | The version of toxval into which the source info is loaded.     |
| <code>source</code>    | The specific source to be loaded, If NULL, load for all sources |
| <code>dict_date</code> | Date stamp of the dictionary file to update                     |

---

|                          |   |
|--------------------------|---|
| <code>load.dsstox</code> | <i>Load DSSTox if needed from a file into a global variables (DSSTOX)</i> |
|--------------------------|---|

---

**Description**

Load DSSTox if needed from a file into a global variables (DSSTOX)

**Usage**

```
load.dsstox(reprocess = FALSE)
```

---

|             |  |
|-------------|--|
| log_message | Function to combine output log with output message |
|-------------|--|

---

**Description**

Function to combine output log with output message  
Function to combine output log with output message

**Usage**

log\_message(log\_df, message\_df\_col)  
  
log\_message(log\_df, message\_df\_col)

**Arguments**

log\_df                 Dataframe to which the log information will be appended  
message\_df\_col       New message to add

---

|                |   |
|----------------|---|
| mv_orchestrate | Orchestrate Materialized View(s) Creation |
|----------------|---|

---

**Description**

Function to create a new materialized view database table by querying an input database version based on an input field\_dictionary.xlsx file.

**Usage**

mv\_orchestrate(toxval.db, include.qc.status = FALSE)

**Arguments**

toxval.db             Version of the database to store the new view.  
include.qc.status     Boolean whether to include the qc\_status field, or filter out "fail" records. Default is FALSE.

**Details**

DETAILS

**Value**

None. SQL statements are executed to generate a database table.

**See Also**

[read\\_excel](#), [unite](#), [rename](#), [mutate](#), [case\\_when](#), [select](#), [mutate-joins](#), [filter](#), [bind\\_rows](#), [character\(0\)](#), [MySQLDriver-class](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

|                                   |                             |
|-----------------------------------|-----------------------------|
| <code>printCurrentFunction</code> | <i>printCurrentFunction</i> |
|-----------------------------------|-----------------------------|

---

**Description**

Print the name of the current function

**Usage**

```
printCurrentFunction(comment.string = NA)
```

**Arguments**

`comment.string` An optional string to be printed

**Details**

DETAILS

**Value**

None

**See Also**

[flush.console](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```



---

|                |                       |
|----------------|-----------------------|
| pull_jira_info | <i>pull_jira_info</i> |
|----------------|-----------------------|

---

## Description

Script to process CSV export of Jira into a status log

## Usage

```
pull_jira_info(
  jira_project = "TOXVAL",
  in_file = NULL,
  auth_token = NULL,
  ticket_filter_list = NULL
)
```

## Arguments

|                    |                                       |
|--------------------|---------------------------------------|
| jira_project       | Jira project code (e.g. CVTDB)        |
| in_file            | File path to Jira ticket summary CSV. |
| auth_token         | Authorization token for Jira          |
| ticket_filter_list | List of Jira tickets to filter to     |

## Details

DETAILS

## Value

Summary DataFrame of Jira tickets by Epic, Label, and Status

## See Also

[download.file][utils::download.file], [unzip][utils::unzip] [read\_csv][readr::read\_csv], [cols][readr::cols] [select][dplyr::select], [contains][dplyr::contains], [mutate][dplyr::mutate], [everything][dplyr::everything], [filter][dplyr::filter], [distinct][dplyr::distinct], [left\_join][dplyr::left\_join], [group\_by][dplyr::group\_by], [summarise][dplyr::summarise], [n][dplyr::n] [unite][tidyr::unite] [str\_squish][stringr::str\_squish]

## Examples

```
## Not run:
if(interactive()){
  out = pull_jira_info(jira_project="project_name")
}

## End(Not run)
```

---

|                                 |
|---------------------------------|
| qa_toxval_url_validation        |
| <i>qa_toxval_url_validation</i> |

---

**Description**

Function to pull URLs and log their HTTP statuses

**Usage**

```
qa_toxval_url_validation(  
  tbl_list = c("toxval", "record_source", "source_info"),  
  db,  
  log_suffix  
)
```

**Arguments**

|            |   |
|------------|---|
| tbl_list   | List of ToxVal tables to pull URLs from, Default: c("toxval", "record_source", "source_info")   |
| db         | ToxVal database name to pull URLs from  |
| log_suffix | SUffix to add to end of log file to uniquely identify file, Default: Date stamp from Sys.Date() |

**Details**

DETAILS

**Value**

None. Log file is generated

**See Also**

[read\\_excel filter](#), [mutate](#), [reexports](#), [bind\\_pivot\\_longer](#) [keep GET](#), [status\\_code](#) [write\\_xlsx](#)

**Examples**

```
## Not run:  
if(interactive()){  
  qa_toxval_url_validation(tbl_list = c("toxval", "record_source", "source_info"), db="res_toxval_v94", log_suffix="")  
}  
  
## End(Not run)
```

---

runInsert

*runInsert*


---

### Description

Insert a record into a database. if auto.increment=TRUE, return the auto incremented primary key of the record. otherwise, return -1

### Usage

```
runInsert(query, db, do.halt = F, verbose = F, auto.increment.id = F)
```

### Arguments

|                   |   |
|-------------------|---|
| query             | a properly formatted SQL query as a string                                |
| db                | the name of the database  |
| do.halt           | if TRUE, halt on errors or warnings                                       |
| verbose           | if TRUE, print diagnostic information                                     |
| auto.increment.id | PARAM_DESCRIPTION, Default: F   |
| auto.increment    | if TRUE, add the auto increment primary key even if not part of the query |

### Details

DETAILS

### Value

Returns the database table auto incremented primary key ID

### See Also

[character\(0\)](#), [MySQLDriver-class](#)

### Examples

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

|                |                       |
|----------------|-----------------------|
| runInsertTable | <i>runInsertTable</i> |
|----------------|-----------------------|

---

## Description

Inserts multiple rows into a database table

## Usage

```
runInsertTable(mat, table, db, do.halt = TRUE, verbose = FALSE, get.id = TRUE)
```

## Arguments

|         |   |
|---------|---|
| mat     | data frame containing the data, with the column names corresponding |
| table   | name of the database table to which data will be inserted           |
| db      | the name of the database  |
| do.halt | if TRUE, halt on errors or warnings                                 |
| verbose | if TRUE, print diagnostic information                               |
| get.id  | Whether to return ID or not, Default: T                             |

## Details

DETAILS

## Value

ID or None

## See Also

[character\(0\), MySQLDriver-class](#)

## Examples

```
## Not run:  
if(interactive()){  
  #EXAMPLE1  
}  
  
## End(Not run)
```

---

`runQuery`*runQuery*

---

## Description

Runs a database query and returns a result set

## Usage

```
runQuery(query = NULL, db, do.halt = TRUE, verbose = FALSE)
```

## Arguments

|                      |  |
|----------------------|--|
| <code>query</code>   | a properly formatted SQL query as a string |
| <code>db</code>      | the name of the database                   |
| <code>do.halt</code> | if TRUE, halt on errors or warnings        |
| <code>verbose</code> | if TRUE, print diagnostic information      |

## Details

DETAILS

## Value

Query results

## See Also

[character\(0\), MySQLDriver-class flush.console](#)

## Examples

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

runStatement

*runStatement*

---

**Description**

Run a SQL statement, such as an ALTER or UPDATE

**Usage**

```
runStatement(query, db, do.halt = FALSE, verbose = FALSE)
```

**Arguments**

|         |  |
|---------|--|
| query   | a properly formatted SQL query as a string |
| db      | the name of the database                   |
| do.halt | if TRUE, halt on errors or warnings        |
| verbose | if TRUE, print diagnostic information      |

**Details**

DETAILS

**Value**

None. SQL statement is run.

**See Also**

[character\(0\)](#), [MySQLDriver-class](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

|           |                  |
|-----------|------------------|
| runUpdate | <i>runUpdate</i> |
|-----------|------------------|

---

**Description**

Runs a database query and returns a result set

**Usage**

```
runUpdate(  
    table,  
    updateQuery = NULL,  
    updated_df = NULL,  
    db,  
    do.halt = TRUE,  
    verbose = FALSE,  
    trigger_check = TRUE  
)
```

**Arguments**

|               |  |
|---------------|--|
| table         | table to update  |
| updateQuery   | a properly formatted SQL query as a string in the form of an UPDATE INNER JOIN |
| updated_df    | a dataframe of updated data to temporarily write to database for INNER JOIN    |
| db            | the name of the database   |
| do.halt       | if TRUE, halt on errors or warnings  |
| verbose       | if TRUE, print diagnostic information  |
| trigger_check | if FALSE, audit triggers are ignored/bypassed                                  |

**Details**

DETAILS

**Value**

None

**See Also**

[character\(0\), MySQLDriver-class dbSendStatement](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

```
set.critical_effect_category
```

*Sets the final category for each term/study\_type pair in the critical\_effect\_terms table*

---

**Description**

Sets the final category for each term/study\_type pair in the critical\_effect\_terms table

**Usage**

```
set.critical_effect_category(toxval.db)
```

**Arguments**

toxval.db      The version of toxval into which the tables are loaded.

**Value**

None. SQL update statement is performed

---

```
set.experimental_record.by.source
```

*Sets experimental\_record flag by source for records in toxval*

---

**Description**

Sets experimental\_record flag by source for records in toxval

**Usage**

```
set.experimental_record.by.source(toxval.db, source = NULL)
```

**Arguments**

toxval.db      The version of toxval into which the tables are loaded.  
 source        Name of source to set. Default NULL means set experimental record for all sources



**Value**

None. SQL update statement is performed

---

set.initial.qc\_status    *Set toxval qc\_status to source table qc\_status*

---

**Description**

Set toxval qc\_status to source table qc\_status

**Usage**

```
set.initial.qc_status(toxval.db, source.db, source, subsource = NULL)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | The database version to use                         |
| source.db | The source database name                            |
| source    | The source name                                     |
| subsource | The subsource to update, if desired (Default: NULL) |

---

set.qc.category.by.source  
*Function for assigning QC Categories to sources in toxval via the qc\_category field*

---

**Description**

Function for assigning QC Categories to sources in toxval via the qc\_category field

**Usage**

```
set.qc.category.by.source(  
  toxval.db,  
  source.db,  
  source = NULL,  
  confluence_url = "https://confluence.epa.gov/x/VuCkFg",  
  confluence_access_token,  
  jira_access_token,  
  jira_rdata = "qc_category"  
)
```

**Arguments**

|                         |  |
|-------------------------|--|
| toxval.db               | The version of toxval into which the tables are loaded.      |
| source.db               | The source database to use.                                  |
| source                  | The source to set a qc_category for                          |
| confluence_url          | URL to QC tracking Confluence page                           |
| confluence_access_token | A personal access token for authentication in Confluence     |
| jira_access_token       | A personal access token for authentication in Jira           |
| jira_rdata              | Local RData file of Jira data to use for jira_tickets object |

---

```
set.study_type.by.study_group
```

*Normalize study\_type within study\_group*

---

**Description**

Normalize study\_type within study\_group

**Usage**

```
set.study_type.by.study_group(
  toxval.db,
  source = NULL,
  subsource = NULL,
  report.only = TRUE,
  filter.matching = FALSE
)
```

**Arguments**

|                 |  |
|-----------------|--|
| toxval.db       | The version of toxvaldb to use.  |
| source          | Source to check  |
| subsource       | Subsource to check (NULL default)  |
| report.only     | Whether to alter ToxVal (FALSE) or simply record suggestions (TRUE), default TRUE    |
| filter.matching | Whether to ignore entries whose current study_type matches suggestion, default FALSE |

---

|   |
|---|
| set.supersource.export.names  |
| <i>Populate export_source_name and supersource fields in ToxVal</i> |

---

**Description**

Populate export\_source\_name and supersource fields in ToxVal

**Usage**

```
set.supersource.export.names(  
    toxval.db,  
    source = NULL,  
    version_date = "2025-05-16"  
)
```

**Arguments**

- |              |  |
|--------------|--|
| toxval.db    | The version of toxval from which to set defaults.      |
| source       | The source to be fixed (if NULL then edit all sources) |
| version_date | The date of the source_info dictionary to be used      |

---

|  |
|--|
| set_extraction_doc_clowder_id  |
| <i>Inserts clowder document information into record_source table</i> |

---

**Description**

Inserts clowder document information into record\_source table

**Usage**

```
set_extraction_doc_clowder_id(toxval.db, source.db, source = NULL)
```

**Arguments**

- |           |  |
|-----------|--|
| toxval.db | The version of toxval into which the tables are loaded.                  |
| source.db | The source database to use.  |
| source    | Name of source to set. Default NULL means set Clowder ID for all sources |

**Value**

None. SQL insert statement is performed

set\_toxval\_relationship\_by\_toxval\_type

*Generic function for setting record relationships based on standardized rules*

---

### Description

Generic function for setting record relationships based on standardized rules

### Usage

```
set_toxval_relationship_by_toxval_type(res, toxval.db)
```

### Arguments

|           |   |
|-----------|---|
| res       | The data that has relationships to be represented       |
| toxval.db | The version of toxval into which the tables are loaded. |

---

source\_chemical.chemidplus

*special process to deal with source chemicals for ChemIDPlus*

---

### Description

special process to deal with source chemicals for ChemIDPlus

### Usage

```
source_chemical.chemidplus(  
  toxval.db,  
  source.db,  
  res,  
  source,  
  chem.check.halt = FALSE,  
  casrn.col = "casrn",  
  name.col = "name",  
  verbose = F  
)
```

**Arguments**

|                 |   |
|-----------------|---|
| toxval.db       | The version of toxval into which the source info is loaded. |
| source.db       | The source database version                                 |
| source          | The xource to be processed (ECOTOX)                         |
| chem.check.halt | If TRUE, halt if there are errors in the chemical checking  |
| casrn.col       | Name of the column containing the CASRN                     |
| name.col        | Name of the column containing chemical names                |
| verbose         | If TRUE, output extra diagnostics information               |

---

source\_chemical.ecotox

*special process to deal with source chemicals for ECOTOX*


---

**Description**

special process to deal with source chemicals for ECOTOX

**Usage**

```
source_chemical.ecotox(
  toxval.db,
  source.db,
  res,
  source,
  chem.check.halt = FALSE,
  casrn.col = "casrn",
  name.col = "name",
  verbose = F
)
```

**Arguments**

|                 |   |
|-----------------|---|
| toxval.db       | The version of toxval into which the source info is loaded. |
| source.db       | The source database version                                 |
| source          | The xource to be processed (ECOTOX)                         |
| chem.check.halt | If TRUE, halt if there are errors in the chemical checking  |
| casrn.col       | Name of the column containing the CASRN                     |
| name.col        | Name of the column containing chemical names                |
| verbose         | If TRUE, output extra diagnostics information               |

---

source\_chemical.extra *special process to deal with source chemicals for extra source (cancer, genetox, skin\_eye, etc)*

---

### Description

special process to deal with source chemicals for extra source (cancer, genetox, skin\_eye, etc)

### Usage

```
source_chemical.extra(
  toxval.db,
  source.db,
  res,
  source,
  chem.check.halt = FALSE,
  casrn.col = "casrn",
  name.col = "name",
  verbose = FALSE
)
```

### Arguments

|           |   |
|-----------|---|
| toxval.db | The version of toxval into which the source info is loaded.   |
| source.db | The source database version   |
| source    | The source to be processed #' @param chem.check.halt If TRUE, halt if there are errors in the chemical checking |
| casrn.col | Name of the column containing the CASRN   |
| name.col  | Name of the column containing chemical names  |
| verbose   | If TRUE, output extra diagnostics information   |

---

source\_chemical.toxrefdb  
*Special process to deal with source chemicals for ToxRefDB. This will put the chemicals into the source database source\_chemical table*

---

### Description

Special process to deal with source chemicals for ToxRefDB. This will put the chemicals into the source database source\_chemical table

**Usage**

```
source_chemical.toxrefdb(
  toxval.db,
  source.db,
  res,
  source = "ToxRefDB",
  chem.check.halt = FALSE,
  casrn.col = "casrn",
  name.col = "name",
  verbose = FALSE
)
```

**Arguments**

|                 |   |
|-----------------|---|
| toxval.db       | The version of toxval into which the source info is loaded.   |
| source.db       | The source database version                                   |
| res             | The dataframe to which the chemical_id will be added          |
| source          | The name of the source  |
| chem.check.halt | If TRUE, stop if there are problems with the chemical mapping |
| casrn.col       | The name of the column containing the CASRN                   |
| name.col        | The name of the column containing the chemical name           |
| verbose         | If TRUE, write out diagnostic messages                        |

**Value**

Returns the input dataframe with the chemical\_id added

---

source\_hash\_vectorized

*source\_hash\_vectorized*

---

**Description**

Generate the hash key for a source table based on hashing columns

**Usage**

```
source_hash_vectorized(res, hashing_cols)
```

**Arguments**

|              |  |
|--------------|--|
| res          | The data frame to be processed                             |
| hashing_cols | Optional list of columns to use for generating source_hash |

Details

DETAILS

Value

Input dataframe with new source\_hash field

See Also

[digest distinct](#)

Examples

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

|            |  |
|------------|--|
| species.qc | Run some checks on the species information |
|------------|--|

---

Description

Run some checks on the species information

Usage

```
species.qc(toxval.db)
```

Arguments

toxval.db      The version of toxval into which the tables are loaded.

---

|                         |                                    |
|-------------------------|------------------------------------|
| species.strain.mismatch | Find species and strain mismatches |
|-------------------------|------------------------------------|

---

Description

Find species and strain mismatches

Usage

```
species.strain.mismatch(toxval.db)
```



**Arguments**

toxval.db      Database version

**Value**

Write a file with the results

---

|               |                      |
|---------------|----------------------|
| toxval.config | <i>toxval.config</i> |
|---------------|----------------------|

---

**Description**

Define a set of global variables. These include the source path (datapath) and the source databases (e.g. dev\_toxval\_version and dev\_toxval\_source\_version).

**Usage**

toxval.config()

**Details**

DETAILS

**Value**

Returns a set of parameters to be used throughout the package

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

|                |  |
|----------------|--|
| toxval.init.db | <i>Initialize the database. THis sill load the species, info and dictionary tables</i> |
|----------------|--|

---

### Description

Initialize the database. THis sill load the species, info and dictionary tables

### Usage

```
toxval.init.db(toxval.db, version_info, reset = F, date_string = "2022-05-25")
```

### Arguments

|              |   |
|--------------|---|
| toxval.db    | The version of toxval into which the tables are loaded. |
| version_info | Info string describing the version.                     |
| reset        | If TRUE, delete all content from the database           |
| date_string  | The date of the dictionary versions                     |

---

|                        |  |
|------------------------|--|
| toxval.load.alaska_dec | <i>Load the alaska_dec (old ACToR - flex)data from toxval sourcedb to toxval</i> |
|------------------------|--|

---

### Description

Load the alaska\_dec (old ACToR - flex)data from toxval sourcedb to toxval

### Usage

```
toxval.load.alaska_dec(toxval.db, source.db, log = F)
```

### Arguments

|           |  |
|-----------|--|
| toxval.db | The database version to use                |
| source.db | The source database                        |
| log       | If TRUE, output log inoformation to a file |

---

|                              |  |
|------------------------------|--|
| <code>toxval.load.all</code> | <i>Load and process all information into ToxValDB. The entire process can be run with one command: <code>toxval.load.all(toxval.db=...,source.db=..., do.all=TRUE)</code> It can also be run in stages, but needs to be run in the order of the do.X parameters listed here. If any earlier step is run, all of the subsequent steps need to be rerun.</i> |
|------------------------------|--|

---

## Description

Load and process all information into ToxValDB. The entire process can be run with one command: `toxval.load.all(toxval.db=...,source.db=..., do.all=TRUE)` It can also be run in stages, but needs to be run in the order of the do.X parameters listed here. If any earlier step is run, all of the subsequent steps need to be rerun.

## Usage

```
toxval.load.all(
  toxval.db,
  source.db,
  confluence_access_token,
  jira_access_token,
  log = FALSE,
  do.init = FALSE,
  do.reset = FALSE,
  do.load = FALSE,
  do.post = FALSE,
  do.extra = FALSE
)
```

## Arguments

|                                      |   |
|--------------------------------------|---|
| <code>toxval.db</code>               | The version of toxval into which the tables are loaded.                           |
| <code>source.db</code>               | The version of toxval_source database from which information is pulled.           |
| <code>confluence_access_token</code> | A personal access token for authentication in Confluence                          |
| <code>jira_access_token</code>       | A personal access token for authentication in Jira                                |
| <code>log</code>                     | If TRUE write the output from each load script to a log file                      |
| <code>do.init</code>                 | If True, clean out all of the database tables                                     |
| <code>do.reset</code>                | If TRUE, empty the database to restart  |
| <code>do.load</code>                 | If TRUE, load all of the source   |
| <code>do.post</code>                 | If TRUE, do th post-processingwork of fixing study type and risk_assessment_class |
| <code>do.extra</code>                | If TRUE, load the non-toxval data (genetox, bcfbaf, skin/eye)                     |

**Value**

Nothing is returned

---

```
toxval.load.atsdr.pfas.2021
```

*Load data ATSDR PFAS 2021 data from toxval\_source to toxval*

---

**Description**

Load data ATSDR PFAS 2021 data from toxval\_source to toxval

**Usage**

```
toxval.load.atsdr.pfas.2021(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

```
toxval.load.atsdr_mrls
```

*Load the ATSDR MRLs data from toxval\_source to toxval*

---

**Description**

Load the ATSDR MRLs data from toxval\_source to toxval

**Usage**

```
toxval.load.atsdr_mrls(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The version of toxval into which the tables are loaded.     |
| source.db          | The source database to use.                                 |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                       |  |
|-----------------------|--|
| toxval.load.caloeehha | <i>Load new_caloeehha from toxval_source to toxval</i> |
|-----------------------|--|

---

**Description**

Load new\_caloeehha from toxval\_source to toxval

**Usage**

```
toxval.load.caloeehha(
  toxval.db,
  source.db,
  log = FALSE,
  remove_null_dtxsid = TRUE
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The version of toxval into which the tables are loaded.     |
| source.db          | The source database to use.                                 |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                    |  |
|--------------------|--|
| toxval.load.cancer | <i>prepare the cancer call data. The data comes form a series of files<br/>       ../NIOSH/NIOSH_CARC_2018.xlsx ../IRIS/iris_cancer_call_2018-10-03.xlsx ../PPRTV_ORNL/PPRTV_ORNL cancer calls 2018-10-25.xlsx ../cancer_summary/cancer/NTP/NTP cancer clean.xlsx ../cancer_summary/cancer/IARC/IARC cancer 2018-10-29.xlsx ../cancer_summary/cancer/HealthCanada/HealthCanada_TRVs_2010_AppendixA v2.xlsx ../cancer_summary/cancer/EPA_OPP_CARC/EPA_CARC.xlsx ../cancer_summary/cancer/CalEPA/calepa_p65_cancer_only.xlsx</i> |
|--------------------|--|

---

**Description**

extract all of the chemicals with cancer slope factor or unit risk with appropriate units

**Usage**

```
toxval.load.cancer(toxval.db, source.db)
```

**Arguments**

|           |                                    |
|-----------|------------------------------------|
| toxval.db | The version of the database to use |
|-----------|------------------------------------|

---

```
toxval.load.chemidplus
```

*Load ChemID Plus Acute data to toxval*

---

**Description**

Load ChemID Plus Acute data to toxval

**Usage**

```
toxval.load.chemidplus(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE,  
  do.init = FALSE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                   |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                            |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value   |
| do.init            | if TRUE, read the data in from the file and set up the matrix |

---

|                    |  |
|--------------------|--|
| toxval.load.copper | <i>Load Copper Manufacturers data to toxval from toxval_source</i> |
|--------------------|--|

---

**Description**

Load Copper Manufacturers data to toxval from toxval\_source

**Usage**

```
toxval.load.copper(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

- |                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                   |                          |
|-------------------|--------------------------|
| toxval.load.dedup | <i>toxval.load.dedup</i> |
|-------------------|--------------------------|

---

**Description**

Perform deduping on data before it is loaded to toxval

**Usage**

```
toxval.load.dedup(  
  res,  
  dedup_fields = NULL,  
  hashing_cols = toxval.config()$hashing_cols,  
  delim = " |::| "  
)
```

**Arguments**

- |              |  |
|--------------|--|
| res          | dataframe containing the source data to dedup  |
| dedup_fields | vector containing field names to dedup, Default: NULL (all fields but hashing cols)      |
| hashing_cols | vector containing field names of hashing columns, Default: toxval.config()\$hashing_cols |
| delim        | string used to separate collapsed values, Default: '  ::  '                              |

**Details**

DETAILS

**Value**

dataframe containing deduped source data

**See Also**

[select](#), [group\\_by](#), [summarise](#), [context](#), [filter](#), [mutate](#), [across](#), [reexports](#), [na\\_if](#), [distinct](#)

**Examples**

```
## Not run:
if(interactive()){
  #EXAMPLE1
}

## End(Not run)
```

---

|                 |   |
|-----------------|---|
| toxval.load.dod | <i>Load the DOD data from toxval_source to toxval</i> |
|-----------------|---|

---

**Description**

Load the DOD data from toxval\_source to toxval

**Usage**

```
toxval.load.dod(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

- |           |   |
|-----------|---|
| toxval.db | The version of toxval into which the tables are loaded. |
| source.db | The source database to use.                             |
| log       | If TRUE, send output to a log file                      |



---

`toxval.load.doe.benchmarks`*Load DOE Wildlife Benchmarks data from toxval\_source to toxval*

---

**Description**

Load DOE Wildlife Benchmarks data from toxval\_source to toxval

**Usage**

```
toxval.load.doe.benchmarks(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                                 |   |
|---------------------------------|---|
| <code>toxval.db</code>          | The database version to use                                 |
| <code>source.db</code>          | The source database   |
| <code>log</code>                | If TRUE, send output to a log file                          |
| <code>remove_null_dtxsid</code> | If TRUE, delete source records without curated DTXSID value |

---

`toxval.load.doe.pac`     *Load DOE Protective Action Criteria data from toxval\_source to toxval*

---

**Description**

Load DOE Protective Action Criteria data from toxval\_source to toxval

**Usage**

```
toxval.load.doe.pac(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The version of toxval into which the tables are loaded.     |
| source.db          | The source database to use.                                 |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

toxval.load.echa\_iuclid

*Loading the ECHA IUCLID data to toxval from toxval\_source This method is different from most because there are multiple tables (one per study type) for this source*

---

**Description**

Loading the ECHA IUCLID data to toxval from toxval\_source This method is different from most because there are multiple tables (one per study type) for this source

**Usage**

```
toxval.load.echa_iuclid(
  toxval.db,
  source.db,
  log = FALSE,
  remove_null_dtxsid = TRUE
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                    |   |
|--------------------|---|
| toxval.load.ecotox | <i>Load ECOTOX from the datahub to toxval</i> |
|--------------------|---|

---

**Description**

Load ECOTOX from the datahub to toxval

**Usage**

```
toxval.load.ecotox(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE,  
  sys.date = "2024-09-19"  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |
| sys.date           | The version of the data to be used                          |

---

|                  |  |
|------------------|--|
| toxval.load.efsa | <i>Load EFSA data from toxval_source to toxval</i> |
|------------------|--|

---

**Description**

Load EFSA data from toxval\_source to toxval

**Usage**

```
toxval.load.efsa(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | The version of toxval into which the tables are loaded. |
| source.db | The source database to use.                             |
| log       | If TRUE, send output to a log file                      |

---

`toxval.load.epa_aegl`    *Load the EPA AEGL data from toxval\_source to toxval*

---

### Description

Load the EPA AEGL data from toxval\_source to toxval

### Usage

```
toxval.load.epa_aegl(  
    toxval.db,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```

### Arguments

|                                 |   |
|---------------------------------|---|
| <code>toxval.db</code>          | The database version to use                                 |
| <code>source.db</code>          | The source database   |
| <code>log</code>                | If TRUE, send output to a log file                          |
| <code>remove_null_dtxsid</code> | If TRUE, delete source records without curated DTXSID value |

---

`toxval.load.epa_dcap`    *Load EPA DCAP source from toxval\_source to toxval*

---

### Description

Load EPA DCAP source from toxval\_source to toxval

### Usage

```
toxval.load.epa_dcap(  
    toxval.db,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtssid | If TRUE, delete source records without curated DTSSID value |

---

toxval.load.epa\_etap    *Load EPA ETAP source from toxval\_source to toxval*

---

**Description**

Load EPA ETAP source from toxval\_source to toxval

**Usage**

```
toxval.load.epa_etap(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtssid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtssid | If TRUE, delete source records without curated DTSSID value |

---

toxval.load.epa\_hhtv    *Load EPA HHTV data to toxval from toxval\_source*

---

**Description**

Load EPA HHTV data to toxval from toxval\_source

**Usage**

```
toxval.load.epa_hhtv(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtssid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

```
toxval.load.epa_ow_npdwr
```

*Loading EPA OW NPDWR to toxval from toxval\_source*

---

**Description**

Loading EPA OW NPDWR to toxval from toxval\_source

**Usage**

```
toxval.load.epa_ow_npdwr(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

```
toxval.load.epa_ow_nrwqc_hhc
```

*Loading EPA OW NRWQC-HHC to toxval from toxval\_source*

---

**Description**

Loading EPA OW NRWQC-HHC to toxval from toxval\_source

**Usage**

```
toxval.load.epa_ow_nrwqc_hhc(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtgsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtgsid | If TRUE, delete source records without curated DTGSID value |

---

|                     |   |
|---------------------|---|
| toxval.load.generic | <i>Generic structure for loading to toxval from toxval_source</i> |
|---------------------|---|

---

**Description**

Generic structure for loading to toxval from toxval\_source

**Usage**

```
toxval.load.generic(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtgsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtgsid | If TRUE, delete source records without curated DTGSID value |

```
toxval.load.genetox.all
```

*Load the Genetox data from Grace*

---

### Description

Load the Genetox data from Grace

### Usage

```
toxval.load.genetox.all(  
  toxval.db,  
  source.db,  
  sys.date = "2021-09-10",  
  verbose = FALSE  
)
```

### Arguments

|           |                                  |
|-----------|----------------------------------|
| toxval.db | The database to use.             |
| source.db | The source database to use.      |
| sys.date  | The input file version           |
| verbose   | If TRUE output debug information |

---

```
toxval.load.gestis.dnel
```

*Load GESTIS DNEL data from toxval\_source to toxval*

---

### Description

Load GESTIS DNEL data from toxval\_source to toxval

### Usage

```
toxval.load.gestis.dnel(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```



**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                  |   |
|------------------|---|
| toxval.load.hawc | <i>Load HAWC from toxval_source to toxval</i> |
|------------------|---|

---

**Description**

Load HAWC from toxval\_source to toxval

**Usage**

```
toxval.load.hawc(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                           |  |
|---------------------------|--|
| toxval.load.hawc_pfas_150 | <i>Load HAWC PFAS 150 from toxval_source to toxval</i> |
|---------------------------|--|

---

**Description**

Load HAWC PFAS 150 from toxval\_source to toxval

**Usage**

```
toxval.load.hawc_pfas_150(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |  |
|--------------------|--|
| toxval.db          | The version of toxval into which the tables are loaded.        |
| source.db          | The version of toxval_source from which the tables are loaded. |
| log                | If TRUE, send output to a log file                             |
| remove_null_dtgsid | If TRUE, delete source records without curated DTGSID value    |

---

```
toxval.load.hawc_pfas_430
```

*Load HAWC PFAS 430 from toxval\_source to toxval*

---

**Description**

Load HAWC PFAS 430 from toxval\_source to toxval

**Usage**

```
toxval.load.hawc_pfas_430(
  toxval.db,
  source.db,
  log = FALSE,
  remove_null_dtgsid = TRUE
)
```

**Arguments**

|                    |  |
|--------------------|--|
| toxval.db          | The version of toxval into which the tables are loaded.        |
| source.db          | The version of toxval_source from which the tables are loaded. |
| log                | If TRUE, send output to a log file                             |
| remove_null_dtgsid | If TRUE, delete source records without curated DTGSID value    |

---

```
toxval.load.healthcanada
```

*Load Health Canada data from toxval\_source to toxval*

---

**Description**

Generic structure for loading to toxval from toxval\_source

**Usage**

```
toxval.load.healthcanada(  
    toxval.db,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                   |   |
|-------------------|---|
| toxval.load.heast | <i>Load the HEAST data from toxval_source to toxval</i> |
|-------------------|---|

---

**Description**

Load the HEAST data from toxval\_source to toxval

**Usage**

```
toxval.load.heast(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                  |                                 |
|------------------|---------------------------------|
| toxval.load.hess | <i>Load HESS data to ToxVal</i> |
|------------------|---------------------------------|

---

**Description**

Load HESS data to ToxVal

**Usage**

```
toxval.load.hess(toxval.db, source.db, log = FALSE, remove_null_dtgsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtgsid | If TRUE, delete source records without curated DTGSID value |

---

|                   |  |
|-------------------|--|
| toxval.load.hpvis | <i>Load HPVIS from toxval_source to toxval</i> |
|-------------------|--|

---

**Description**

Load HPVIS from toxval\_source to toxval

**Usage**

```
toxval.load.hpvis(toxval.db, source.db, log = FALSE, remove_null_dtgsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtgsid | If TRUE, delete source records without curated DTGSID value |

---

|                     |   |
|---------------------|---|
| toxval.load.initial | <i>Delete the contents of the toxval database</i> |
|---------------------|---|

---

**Description**

Delete the contents of the toxval database

**Usage**

```
toxval.load.initial(toxval.db)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | The version of toxval from which the data is deleted. |
|-----------|---|

---

|                  |  |
|------------------|--|
| toxval.load.iris | <i>Load IRIS source from toxval_source to toxval</i> |
|------------------|--|

---

**Description**

Load IRIS source from toxval\_source to toxval

**Usage**

```
toxval.load.iris(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The version of toxval into which the tables are loaded.     |
| source.db          | The source database to use.                                 |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                       |  |
|-----------------------|--|
| toxval.load.mass_mmcl | <i>Load the Mass. Drinking Water Standards data from toxval_source to toxval</i> |
|-----------------------|--|

---

**Description**

Load the Mass. Drinking Water Standards data from toxval\_source to toxval

**Usage**

```
toxval.load.mass_mmcl(  
    toxval.db,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                   |  |
|-------------------|--|
| toxval.load.niosh | <i>Load NIOSH from toxval_source to toxval</i> |
|-------------------|--|

---

**Description**

Load NIOSH from toxval\_source to toxval

**Usage**

```
toxval.load.niosh(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                      |   |
|----------------------|---|
| toxval.load.ntp.pfas | <i>Load NTP PFAS data from toxval_source to toxval Generic structure for loading to toxval from toxval_source</i> |
|----------------------|---|

---

**Description**

Load NTP PFAS data from toxval\_source to toxval Generic structure for loading to toxval from toxval\_source

**Usage**

```
toxval.load.ntp.pfas(  
    toxval.db,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                 |   |
|-----------------|---|
| toxval.load.opp | <i>Load EPA OPP data to toxval from toxval_source</i> |
|-----------------|---|

---

**Description**

Load EPA OPP data to toxval from toxval\_source

**Usage**

```
toxval.load.opp(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

```
toxval.load.osha_air_limits
```

*Load the osha\_air\_limits (old ACToR - flex) data from toxval sourcedb to toxval*

---

### Description

Load the osha\_air\_limits (old ACToR - flex) data from toxval sourcedb to toxval

### Usage

```
toxval.load.osha_air_limits(toxval.db, source.db, log = F)
```

### Arguments

|           |                                    |
|-----------|------------------------------------|
| toxval.db | The database version to use        |
| source.db | The source database                |
| log       | If TRUE, send output to a log file |

---

```
toxval.load.ow_dwsha
```

*Load the ow\_dwsha data from toxval\_source to toxval*

---

### Description

Load the ow\_dwsha data from toxval\_source to toxval

### Usage

```
toxval.load.ow_dwsha(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

### Arguments

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |



---

|                  |  |
|------------------|--|
| toxval.load.penn | <i>Load Penn data from toxval_source to toxval</i> |
|------------------|--|

---

**Description**

Load Penn data from toxval\_source to toxval

**Usage**

```
toxval.load.penn(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The version of toxval into which the tables are loaded.     |
| source.db          | The source database to use.                                 |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

---

|                           |   |
|---------------------------|---|
| toxval.load.penn_dep_mscs | <i>Load the penn_dep_mscs data from toxval_source to toxval</i> |
|---------------------------|---|

---

**Description**

Load the penn\_dep\_mscs data from toxval\_source to toxval

**Usage**

```
toxval.load.penn_dep_mscs(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

```
toxval.load.pfas_150_sem_v2
```

*Load pfas\_150\_sem from toxval\_source to toxval*

---

### Description

Load pfas\_150\_sem from toxval\_source to toxval

### Usage

```
toxval.load.pfas_150_sem_v2(  
    toxval.db,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```

### Arguments

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

```
toxval.load.postprocess
```

*Do all of the post-processing steps for a source*

---

### Description

Do all of the post-processing steps for a source

### Usage

```
toxval.load.postprocess(  
    toxval.db,  
    source.db,  
    source,  
    do.convert.units = FALSE,  
    chem_source = NULL,  
    subsource = NULL,  
    remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use   |
| source.db          | The source database name  |
| source             | The source name   |
| do.convert.units   | If TRUE, convert units, mainly from ppm to mg/kg-day. This code is not debugged |
| chem_source        | Used only for source=ECHA IUCLID  |
| subsource          | The specific subsource to process, if desired (Default: NULL)                   |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value                     |

---

`toxval.load.pprtv.cphea`*Load PPRTV CPHEA from toxval\_source to toxval*

---

**Description**

Load PPRTV CPHEA from toxval\_source to toxval

**Usage**

```
toxval.load.pprtv.cphea(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                 |  |
|-----------------|--|
| toxval.load.rsl | <i>Load the RSL data from source db to toxval - the source database needs to be updated periodically</i> |
|-----------------|--|

---

**Description**

Load the RSL data from source db to toxval - the source database needs to be updated periodically

**Usage**

```
toxval.load.rsl(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                      |                               |
|----------------------|-------------------------------|
| toxval.load.skin.eye | <i>Load the Skin eye data</i> |
|----------------------|-------------------------------|

---

**Description**

Load the Skin eye data

**Usage**

```
toxval.load.skin.eye(toxval.db, source.db, verbose = FALSE)
```

**Arguments**

|           |  |
|-----------|--|
| toxval.db | Database version                                 |
| verbose   | if TRUE, print diagnostic messages along the way |

---

```
toxval.load.source_chemical
```

*Perform the DSSTox mapping*

---

**Description**

Perform the DSSTox mapping

**Usage**

```
toxval.load.source_chemical(  
    toxval.db,  
    source.db,  
    source = NULL,  
    verbose = TRUE,  
    remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The version of toxvaldb to use.                             |
| source.db          | The source database version                                 |
| source             | The source to update for                                    |
| verbose            | If TRUE, print out extra diagnostic messages                |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

```
toxval.load.species
```

*Load the species table*

---

**Description**

Load the species table

**Usage**

```
toxval.load.species(toxval.db, date_string = "2023-05-18")
```

**Arguments**

|             |   |
|-------------|---|
| toxval.db   | The version of toxval into which the tables are loaded. |
| date_string | The date string for the dictionary files                |

---

|                  |  |
|------------------|--|
| toxval.load.test | <i>Load TEST data from toxval_source to toxval</i> |
|------------------|--|

---

**Description**

Load TEST data from toxval\_source to toxval

**Usage**

```
toxval.load.test(toxval.db, source.db, log = FALSE, remove_null_dtxsid = TRUE)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                         |                                     |
|-------------------------|-------------------------------------|
| toxval.load.toxrefdb2.1 | <i>Load ToxRefdb data to toxval</i> |
|-------------------------|-------------------------------------|

---

**Description**

Load ToxRefdb data to toxval

**Usage**

```
toxval.load.toxrefdb2.1(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

|                       |   |
|-----------------------|---|
| toxval.load.usgs_hbsl | <i>Load the usgs_hbsl (old ACToR - flex) data from toxval source db to toxval</i> |
|-----------------------|---|

---

### Description

Load the usgs\_hbsl (old ACToR - flex) data from toxval source db to toxval

### Usage

```
toxval.load.usgs_hbsl(toxval.db, source.db, log = F)
```

### Arguments

|           |                                    |
|-----------|------------------------------------|
| toxval.db | The database version to use        |
| source.db | The source database                |
| log       | If TRUE, send output to a log file |

---

|                   |   |
|-------------------|---|
| toxval.load.ut_hb | <i>Load the uterotrophic and Hershberger data</i> |
|-------------------|---|

---

### Description

Load the uterotrophic and Hershberger data

### Usage

```
toxval.load.ut_hb(toxval.db, source.db, log = FALSE, remove_null_dtgsid = TRUE)
```

### Arguments

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtgsid | If TRUE, delete source records without curated DTGSID value |

---

toxval.load.who\_ipcs     *Load the who\_ipcs data from toxval source db to toxval*

---

### Description

Load the who\_ipcs data from toxval source db to toxval

### Usage

```
toxval.load.who_ipcs(  
    toxval.db,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```

### Arguments

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

toxval.load.who\_jecfa\_adi  
                          *Load WHO JECFA ADI from toxval\_source to toxval*

---

### Description

Load WHO JECFA ADI from toxval\_source to toxval

### Usage

```
toxval.load.who_jecfa_adi(  
    toxvaldb,  
    source.db,  
    log = FALSE,  
    remove_null_dtxsid = TRUE  
)
```



**Arguments**

|                    |   |
|--------------------|---|
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |
| toxval.db          | The database version to use                                 |

---

`toxval.load.who_jecfa_tox_studies`*Load WHO JECFA Tox Studies from toxval\_source to toxval*

---

**Description**

Load WHO JECFA Tox Studies from toxval\_source to toxval

**Usage**

```
toxval.load.who_jecfa_tox_studies(  
  toxval.db,  
  source.db,  
  log = FALSE,  
  remove_null_dtxsid = TRUE  
)
```

**Arguments**

|                    |   |
|--------------------|---|
| toxval.db          | The database version to use                                 |
| source.db          | The source database   |
| log                | If TRUE, send output to a log file                          |
| remove_null_dtxsid | If TRUE, delete source records without curated DTXSID value |

---

`toxval.set.mw`*Set the molecular weight in the toxval table, for use in unit conversions*

---

**Description**

Set the molecular weight in the toxval table, for use in unit conversions

**Usage**

```
toxval.set.mw(toxval.db, source = NULL, subsource = NULL)
```

**Arguments**

|           |                             |
|-----------|-----------------------------|
| toxval.db | The database version to use |
| source    | The source                  |
| subsource | The subsource               |

---

|                      |   |
|----------------------|---|
| toxval.summary.stats | <i>Generate summary statistics on the toxval database</i> |
|----------------------|---|

---

**Description**

Generate summary statistics on the toxval database

**Usage**

```
toxval.summary.stats(toxval.db, export = FALSE, summ_level = "source")
```

**Arguments**

|            |  |
|------------|--|
| toxval.db  | The version of toxval into which the tables are loaded.      |
| export     | Boolean whether to export a file. Default FALSE.             |
| summ_level | String of "source" or "supersource" how to group the summary |

**Value**

DataFrame of record qc\_status summary by source.

---

|               |  |
|---------------|--|
| toxvaldb.view | <i>Produce a view of the ToxValDB Data</i> |
|---------------|--|

---

**Description**

‘toxvaldb.view’ Produces a view for ToxValDB by merging specified tables

**Usage**

```
toxvaldb.view(toxval.db, user, password, count = 10)
```

**Arguments**

|           |   |
|-----------|---|
| toxval.db | Database version  |
| user      | The username for the MySQL database. The database instance is hard-coded in the function setDBConn(). |
| password  | The user's MySQL database password.   |
| count     | If count>0, only select this number of records from each source, used for debugging                   |

**Value**

Write a file with the results: data/view/ToxValDB View toxval.db Sys.Date().xlsx

---

```
toxvaldb_release_comparison_stats  
  toxvaldb_release_comparison_stats.R
```

---

**Description**

Pull comparison stats between toxvaldb versions.

**Usage**

```
toxvaldb_release_comparison_stats(repoDir)
```

**Arguments**

|         |                      |
|---------|----------------------|
| repoDir | Path to Repo/ folder |
|---------|----------------------|

**Value**

None. RData file is saved.

---

```
toxval_type.species.mismatch
```

*Find species and toxval\_type mismaps, e.g. species other than human  
fro RfD, RfC, cancer sloper, MSL, etc.*

---

**Description**

Find species and toxval\_type mismaps, e.g. species other than human fro RfD, RfC, cancer sloper, MSL, etc.

**Usage**

```
toxval_type.species.mismatch(toxval.db)
```

**Arguments**

|           |                          |
|-----------|--------------------------|
| toxval.db | Database version         |
| source    | The source to be updated |

**Value**

Write a file with the results

---

`%>%`*Pipe operator*

---

**Description**

See `magrittr::%>%` for details.

**Usage**

```
lhs %>% rhs
```

**Arguments**

|                  |  |
|------------------|--|
| <code>lhs</code> | A value or the <code>magrittr</code> placeholder.          |
| <code>rhs</code> | A function call using the <code>magrittr</code> semantics. |

**Value**

The result of calling `'rhs(lhs)'`.

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