# Package 'dbs'

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| <b>Description</b> The package provides the extended facilities for DBSolveOptimum users. It has two main purposes: (1) the creation and analysis of DBsolveOptimum inputs and outputs for multiple simulations and statistical analysis, (2) the import/export DBSolveOptimum model and data files from/to different formats. |
| License MIT  |
| <pre>URL https://github.com/insysbio/dbs-package,</pre>  |
| http://sourceforge.net/projects/dbsolve/   |
| <pre>BugReports https://github.com/insysbio/dbs-package/issues</pre>   |
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dbs-package

Accessory functions to support DBSolveOptimum

### **Description**

The package provides the extended facilities for DBSolveOptimum users. It has two main purposes: (1) the creation and analysis of DBsolveOptimum inputs and outputs for multiple simulations and statistical analysis, (2) the import/export DBSolveOptimum model and data files from/to different formats.

### **Details**

Package: dbs
Type: Package
Version: 0.9.1
Date: 2017-11-20
License: MIT

DBSolveOptimum is a stand-alone software tool for construction and analysis of mathematical models of biological systems. DBSolveOptimum is implemented with new tools for extended data analysis and multiple simulations, which are important for simulation of virtual clinical trials and application of modern modeling techniques, like quantitative systems pharmacology, to problems arising in drug research and development.

DBSolveOptimum is free for academic and industrial use. The latest version of DBSolveOptimum can be downloaded from: http://sourceforge.net/projects/dbsolve/

### Copyright

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insysbio.com

### Author(s)

Evgeny Metelkin, Aleksey Aleksev, Aleksey Kalinkin

### References

N.Gizzatkulov, I.Goryanin, E.Metelkin, E.Mogilevskaya, K.Peskov and O.Demin. DBSolve Optimum: a software package for kinetic modeling which allows dynamic visualization of simulation

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```
results. BMC Systems Biology, 2010, 4 (109): 1-11. PubMed
E.Metelkin, A.Alekseev, G.Lebedeva, O.Demin. DBSolve Optimum r.33: Practical Guide pdf
```

### See Also

```
calccb
calcop
parsetgen
import.slv
rct.from.slv
import.dat
C.from.slv
```

C.from.slv

Create .C code from ruSlv object

### **Description**

The function creates **deSolve**-compatible C code from .SLV image and save it to a file.

### Usage

### **Arguments**

slv object of ruSlv class, model image. file filename to save .C code.

output character vector with names of additional output values.

dbs.compatibility

logical value to use DBSolve-like method of parameters updates.

### See Also

```
import.slv
deSolve
ode
```

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```
func = "derivs",
         parms=example1_ruSlv$ode.parameters.external,
         dllname = "model",
         initfunc = "initmod",
         nout=length("D"),
         outnames = "D"
 )
dyn.unload(paste0("model", .Platform$dynlib.ext))
plot(res)
## End(Not run)
### plot simulations from SLV model
## Not run:
  filePath<-system.file(package = "dbs", "extdata/example1.slv")</pre>
  raw<-read.slv(filePath) # read from example</pre>
  compatible.slv(raw) # TRUE
  example1_ruSlv<-import.slv(raw)
  C.from.slv(example1_ruSlv, output="D")
  system("R CMD SHLIB model.c") # compilation for .DLL
  library(deSolve)
  dyn.load(paste0("model", .Platform$dynlib.ext))
  res<-ode(y=example1_ruSlv$ode.initials,
           times=seq(0,example1_ruSlv$solver.time.limit,0.1),
           func = "derivs",
           parms=example1_ruSlv$ode.parameters.external,
           dllname = "model",
           initfunc = "initmod",
           nout=length("D"),
           outnames = ^{"}D"
  )
  dyn.unload(paste0("model", .Platform$dynlib.ext))
  plot(res)
## End(Not run)
```

calccb

Calculation of confidence bands

### **Description**

The function calculates pointwice confidence bands based on Monte-Carlo simulations in DB-SolveOptimum. The lower and upper confidence band calculated as lower and upper quantile for interpolated particular x point.

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### **Arguments**

| input       | data.frame passed from DBSolveOptimum output  |
|-------------|---|
| x.col       | number or name of column in input corresponded to free variable (i.e. time).  |
| x.seq       | numerical vector of points to interpolate values in x.col.  |
| y.col       | vector of column numbers or names in input corresponded to simulated variables (model output).  |
| factor.col  | vector of column numbers or names in input corresponded to condition parameters (model input).  |
| q.seq       | sequence of probabilities for calculation of lower and upper quantile. The default vector $c(0.025,\ 0.5,\ 0.975)$ corresponds to calculation of median value and 0.95 confidence band. |
| nos.col     | number or name of column in input corresponded to enumeration of random parameter set.  |
| par_calc    | logical value to use parallel calculation for acceleration. It requires parallel, foreach, iterators packages.  |
| cpu.cores   | the number of CPU cores to use if par_calc=TRUE.  |
| silent      | logical value to suppress the messages during calculations.   |
| include.nos | vector of number of samples to analyze the approximation in the chosed number of sample.  |
|             | other arguments passed to quantile  |

#### Value

The returned value is data.frame class object. The columns describe:

names(x.col) free variable values passed from argument x.col.
var\_id names of simulated variables as passed from y.col.
quant\_ columns represent calculated quntiles for interpolated points.
names(factor.col)

condition variable values passed from argument factor.col.

group unique identifier for combination of factor.col.

The value has the additional attributes:

col.def definition of columns, type of data in columns.

col.title titles for columns. May be usefull for visualization.

var.title titles for simulated variables. May be usefull for visualization.

group.title titles for condition groups. May be usefull for visualization.

approx\_nos\_ column(or columns) represents interpolated points for chosed number of sample (presented only if include.nos has values in).

•

### See Also

foreach
quantile
approxfun

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#### **Examples**

```
### calculation of confidence bands based on example4.slv
## Not run: example4_parset_bs.cond_res<-read.delim("example4_parset_bs.cond_res.txt") # read from DBSolve ou
example4_cb<-calccb(input=example4_parset_bs.cond_res,</pre>
                    x.col="t",
                    x.seq=seq(0,96,by=0.5),
                    y.col=c("C0","C1"),
                    factor.col = c("Dose","T"))
## Not run: write.delim(example4_cb, "example4_cb.txt") # save results
### plot all results with lattice
library(lattice)
xyplot(quant_0.025+quant_0.5+quant_0.975~t|var_id+group,
       data=example4_cb,
       type="1",
       1ty=c(2,1,2),
       xlab="Time, h",
       ylab="Concentration of drug, ng/ml",
       main="All CB simulations")
###You can also plot all results using ggplot2:
library(ggplot2)
ggplot(example4_cb,aes(t,quant_0.025))+
        geom_line(linetype="dashed", color="blue")+
        geom_line(aes(t,quant_0.5),color="black",linetype="solid")+
        geom_line(aes(t,quant_0.975),color="blue",linetype="dashed")+
        facet_wrap(~var_id+group)+
        ggtitle("All CB simulations")+
        scale_x_continuous(name="Time,h")+
        scale_y_continuous(name="Concentration of drug, ng/ml")
### plot dbsolve output results
example4_cb<-calccb(input=example4_parset.cond_res,</pre>
                    x.col="t",
                    x.seq=seq(0,96,by=0.5),
                    y.col=c("C0","C1"),
                    factor.col = c("Dose","T"),
                    par_calc = FALSE)
```

calcop

Calculate interpolation for optimal values

### **Description**

The function interpolates simulations in DBSolveOptimum for the series of conditions and create mod.frame object.

```
calcop(input, x.col, x.seq, y.col, factor.col = c())
```

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#### **Arguments**

| input      | data.frame passed from DBSolveOptimum output   |
|------------|--|
| x.col      | number or name of column in input corresponded to free variable (i.e. time).                   |
| x.seq      | numerical vector of points to interpolate values in x.col.                                     |
| y.col      | vector of column numbers or names in input corresponded to simulated variables (model output). |
| factor.col | vector of column numbers or names in input corresponded to condition parameters (model input). |

#### Value

The returned value is mod.frame class object which is extension of data.frame class with the additional attributed. The columns describe:

```
names(x.col) free variable values passed from argument x.col.

var_id names of simulated variables as passed from y.col.

simulation column represents simulation value for interpolated points.

names(factor.col)

condition variable values passed from argument factor.col.

group unique identifier for combination of factor.col.
```

The value has the additional attributes:

```
col.def definition of columns, type of data in columns.

col.title titles for columns. May be usefull for visualization.

var.title group.title titles for condition groups. May be usefull for visualization.
```

#### See Also

```
calccb
approxfun
```

```
### calculation based on example4.slv
## Not run: example4_output_op<-read.delim("dbs_output_op.txt") # read from output</pre>
example4_op<-calcop(input=example4_output_op,</pre>
                    x.col="t",
                    x.seq=seq(0,96,by=0.5),
                    v.col=c("C0","C1"),
                    factor.col = c("Dose", "T"))
## Not run: write.delim(example4_op, "example4_op.txt") # save results
### plot all results with lattice
library(lattice)
xyplot(simulation~t|var_id+group,
       data=example4_op,
       type="1",
       lty=1,
       xlab="Time, h",
       ylab="Concentration of drug, ng/ml",
       main="All CB simulations")
```

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| clean.comments | Clean C-style text |
|----------------|--------------------|
|----------------|--------------------|

### **Description**

Function takes character vector of C-style lines (as passed from readLines), delete comments, spaces, empty lines, multiple semicolomns and line breaks.

### Usage

```
clean.comments(input)
```

### **Arguments**

input

character vector of C-style text lines, vector can be passed from readLines func-

### Value

Character vector of cleaned C-style text. It can be saved to file by cat(..., sep = "\n") function.

### **Examples**

```
## Not run: cTextExample
(cln<-clean.comments(cTextExample))
cat(cln, file="cln.txt", sep = "\n") # save text to file</pre>
```

hessian2stat

Functions for calculating statistical characteristics from hessian or parameter set

### **Description**

The functions calculate statistical characteristics based on normal distribution or transformed normal distribution.

### Usage

```
hessian2stat(hessian, optim, transform="", level=0.95, ...)
parset2stat(parset, transform="", level=0.95, norm.test=shapiro.test, ...)
```

### **Arguments**

hessian Hessian matrix of second derivatives (-2logL vs parameters). It must be positive-

definite.

optim Optimal values of parameters (best-fit values).

transform String vector for what distribution is used on parameter. For default all parame-

ters are set to normal distribution.

nan.plot

level confidence level for calculation of confidence intervals.
... other arguments passed to solve.

parset data.frame with parameter set (mandatory for parset2stat)
norm.test function to perform normality test

### Value

The output for this function is covariance matrix calculated based on assymptotic approach, and it takes into account the transformation of parameter space.

### **Examples**

```
### calculate statistical characteristics based on calculated hessian, see 'example4.slv' from DBSolve manual
## Not run: example4_parset_bs<-read.delim("example4_hessian.txt") # read hessian from file
optimal<-c(kcat=7.130016e-01, Vd=5.205980e+00, Km=5.240306e+00, kabs=2.014304e+00)
hessian2stat(hessian=as.matrix(example4_hessian), optim=optimal, transform="log")

### calculate statistical characteristics based on parameter set, see 'example4.slv' from DBSolve manual
## Not run: example4_parset_bs<-read.delim("example4_parset_bs.txt") # read hessian from file
parset2stat(parset=example4_parset_bs[,1:4])</pre>
```

nan.plot

Plot all points including infinite

### Description

The functions to plot all points of the dataset even they are out of the xlim and ylim range.

### Usage

### Arguments

| Х        | numeric vector to plot on x-axis.   |    |
|----------|---|----|
| У        | numeric vector to plot on y-axis.   |    |
| xlim     | the x limits (x1, x2) for main region.  |    |
| ylim     | the y limits (y1, y2) for main region.  |    |
| log      | a character string which contains "x" if the x axis is to be logarithmic, "y" if the y axis is to be logarithmic and "xy" or "yx" if both axes are to be logarithmic. |    |
|          | other graphical parameters passed to plot.default or points   |    |
| force.bo | logical value to forcefully create the extended region. F means the extender region is created if some of points is out of the main region.                           | ed |
| delta    | the relative size of region based on main region size.  |    |

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#### **Details**

If someone use the default plot and points method the points out of xlim and ylim bacame invisible. In some cases it is not OK. The presented functions creates the exteded region and plot all points including infinete values there.

The function work only for numerical vectors but not for data.frame, matrix, etc.

### Note

These are experimental functions so the troubles are possible.

#### See Also

```
plot.default
points
```

### **Examples**

```
### comparison of exponential plots
x<-seq(0,20, 0.1)
y<-5*exp(x)
plot(x,y) # default plot without limits
plot(x,y, ylim=c(0,100)) # default plot with y limits
nan.plot(x,y, ylim=c(0,100))</pre>
```

parsetgen

Functions for DBSolveOptimum to generate parameters and conditions using multivariate normal distribution

### Description

The function parsetgen generates random dataset.

The function parsetgen.cond add to the dataset from parsetgen conditions columns and nos column

### Usage

```
parsetgen(cov, mu, transform="", samples=1024)
parsetgen.cond(parset, cond=data.frame(), max.samples=nrow(parset), uniq.nos=FALSE)
```

### **Arguments**

| COV       | covariance matrix with dimension nxn and whose element in i,j positions is the covariance between the i and j elements.                     |
|-----------|---|
| mu        | named vector length of n with named parameters and their expectations   |
| transform | character vector containing the distribution for each parameter. It can be "" for normal distribution and "log" for log-normal distribution |
| samples   | number of the samples that we want to get in the ouput dataset (1024 as default)  |
| parset    | output data.frame from parsetgen function or users data.frame   |

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cond conditional data.frame. If empty, the result of parsetgen.cond is input data.frame with nos column

max.samples number of resulted samples for each condition. If nrow(parset)<max.samples, that provide an error. If nrow(parset)>max.samples, that cuts the data frame by number of samples

uniq.nos number of unique parameter set. If FALSE, then nos will be with repeating, and if TRUE - nos will have unique number.

### Value

The output is data.frame with the columns named as names in your expect vector, and for parsetgen.cond it will be data.frame with the "nos" column, and names(parset) columns and names(cond) column

#### See Also

rmvnorm

```
### create parameter set for Monte-Carlo simulation using covariance matrix (example4)
example4_parset<-parsetgen(example4_stat$muCov, example4_stat$mu, example4_stat$transform)
example4_parset.cond<-parsetgen.cond(</pre>
  example4_parset,
  cond=data.frame(Dose=c(1,5,10,1,5,10), T=c(12,12,12,24,24,24))
write.delim(example4_parset.cond, "example4_parset.cond.txt")
### create parameter set for Monte-Carlo simulation using parameter set based on bootstrapping(example4)
example4_parset_bs.cond<-parsetgen.cond(
  example4_parset_bs[,c(1,2,3,4,6)],
  cond=data.frame(Dose=c(1,5,10,1,5,10), T=c(12,12,12,24,24,24))
write.delim(example4_parset_bs.cond, "example4_parset_bs.cond.txt")
### Making some parameter set with three expectation and cov as vector
  expect<-c(kcat=0.5, Vd=23.4, Km=12.4) #making expect vector
  cov<-c(23.5, 37.9, 23.5) #making vector for coviance matrix
  transform<-c("log", "", "log") #the transform vector with distributions
  output<-parsetgen(cov, expect, transform)</pre>
  #Making some dataset, with three expectation and cov as vector
  expect<-c(kcat=0.5, Vd=23.4, Km=12.4) #making expect vector
  cov<-matrix(c(23.5,0,1, 0,37.9,0, 1,0,23.5), ncol=3) #coviance matrix
  \label{transform} {\sf transform} {\sf <-c("log","","log")} \  \, {\sf \#the transform vector with distributions}
  output<-parsetgen(cov, expect, transform)</pre>
  #Using the parsetgen.cond function. Suppose we have output from parsetgen function
  cond < -data.frame(cond1 = c(1, 2.5, 1), cond2 = c(0, 0, 1), cond3 = c(0, 0, 0))
  output1<-parsetgen.cond(output, cond=cond, max.samples=1024,uniq.nos=FALSE)
```

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rct.from.slv

Create .RCT from ruSlv

### **Description**

The function analyzes ruSlv stoicheometry matrix and creates reaction list in .RCT format.

### Usage

```
rct.from.slv(y)
```

### **Arguments**

У

the object of class ruSlv, model image.

### **Details**

The function uses the stoicheometry matrix, names of rates and metabolites only. The true structure of differential equation is not taken into account.

#### Value

character vector representing list of reactions which can be saved using cat(..., sep="\n").

### See Also

```
import.slv
cat
```

### **Examples**

```
### create .RCT file from 'example4.slv'
rct<-rct.from.slv(example4_ruSlv)
cat(rct, file="example.rct", sep="\n")</pre>
```

read.dat

Import .DAT files

### Description

The set of functions to import experimental dataset from DBSolveOptimum format .DAT file. read.dat reads .DAT file and perform initial parsing. import.dat creates ruData object from read.dat output.

```
read.dat(file)
import.dat(dat)
```

read.list

### **Arguments**

file filename of .DAT file.

dat list object of format ruData.raw which is output of read.dat function.

### Value

The returned value of import.dat is an object of class ruData which mode is list and structure corresponds to ruList. Second level is lists each of which has the following components:

data\_id character identifier of the dataset

data experimental data of data. frame class

conditions data.frame describing conditions

solver character identifier of solver type: ode, explicit, implicit

error.type description of error model, currently possible values are: "additive T", "additive

F"

#### See Also

```
write.list
```

### **Examples**

```
### read and save data from 'example4.dat'
filePath<-system.file(package = "dbs", "extdata/example4.dat")
dat_raw<-read.dat(filePath)
example4_ruData<-import.dat(dat_raw)
write.list(example4_ruData, "example4_ruData.txt")</pre>
```

read.list

Read, write and check object of ruList format

### **Description**

The common functions for manipulating objects of structure ruList: reading from file, writing to file of format ruList.txt and checking ruList object for appropriateness.

### Usage

```
read.list(file)
write.list(x, file="")
check.list(x)
```

### **Arguments**

x the object of format ruList file a character string naming a file 14 read.slv

#### **Details**

ruList and ruList.txt is the internal dbs-package format for representing complex objects.

The structure of ruList format can be described as three-level nested object: (1) list with any number of elements of 2-d level, (2) list with any number of elements of 3-d level, (3) objects of classes: data.frame, matrix, numeric, character, integer, logical, mod.frame. Any level may have attributes of classes: data.frame, matrix, numeric, character, integer, logical.

The ruList.txt is a human readable representation of ruList object saved to .TXT file.

#### Value

```
read.list returns the object of ruList format check.list returns logical TRUE if the object x has appropriate structure
```

### Note

The current version of check.list does not check attributes of 3-d level.

The current version of write.list does not check x argument for consistency. Be carefull.

#### See Also

list

### **Examples**

```
### write, read and check ruList
models<-list(example4_ruSlv)
write.list(models, "models.txt")
models1<-read.list("models.txt")
check.list(models1) # output: TRUE
all.equal(models1, models) # output: 'names for target but not for current'</pre>
```

read.slv

Import .SLV files to R

### **Description**

A set of functions to import model files of DBSolveOptimum (.SLV) to R-environment as the object of class ruSlv.

```
read.slv function reads .SLV file and creates list of format ruSlv.raw
```

compatible.slv function checks compatibility of ruSlv.raw passed from read.slv for compatibility with current version of import.slv.

import.slv function analyzes ruSlv.raw passed from read.slv and creates the object of class ruSlv.

```
read.slv(file)
compatible.slv(x)
import.slv(x)
```

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#### **Arguments**

 $\label{eq:saved_from_def} \textbf{file} \qquad \qquad \textbf{valid} \; . \\ \textbf{SLV} \; \textbf{file} \; \textbf{saved} \; \textbf{from} \; \textbf{DBSolveOptimum}.$ 

x object of format ruSlv. raw to use for model import.

#### **Details**

In many cases DBSolveOptimum features are not enough to manipulate the model structure and to simulate specific conditions. This set of functions transforms all the structure of .SLV file to the R-environment for easy manipulation.

read.slv function reads model code from file and perform initial parsing based on SLV version described in file 'slv25tab.csv'.

import.slv analyzes different parts of read.slv output and create an object of class ruSlv which is an image of initial .SLV file. compatible.slv is developed to check the structure of ruSlv.raw object for further parsing using import.slv. It is part of import.slv function so it is not necessary to use it separately.

#### Value

read.slv returns list of the format ruSlv.raw compatible.slv

returns TRUE if x output is compatible with current version of import.slv and

FALSE otherwise

import.slv returns the object of mode list and ruSlv class attribute

#### Note

The current version of dbs-package officially supports only .SLV version 25. Try compatible.slv to check.

Known restrictions:

- 1. Cannot read 'events'
- 2. Cannot read 'fit conditions'

### See Also

```
rct.from.slv
C.from.slv
```

```
### import 'example4.slv'
filePath<-system.file(package = "dbs", "extdata/example4.slv")
raw<-read.slv(filePath) # read from example
compatible.slv(raw) # TRUE
example4_ruSlv<-import.slv(raw)

### import 'example1.slv'
filePath<-system.file(package = "dbs", "extdata/example1.slv")
raw<-read.slv(filePath) # read from example
compatible.slv(raw) # TRUE
example1_ruSlv<-import.slv(raw)</pre>
```

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replacePow

Replacing ^ to Pow

### Description

Function replacePow(e) used in import.slv to replace  $^{\wedge}$  to pow, for convert to C model file without errors.

### Usage

```
replacePow(e)
```

### **Arguments**

е

expression with ^ and sqrt as expression type.

### **Details**

For using function replacePow(e), You need to parse your text and after using function deparse to obtain the text.

### See Also

```
C.from.slv
read.slv
```

### **Examples**

```
#Example to using the replacePow function  text = "Cr1 = fdr^x + x^y + sqrt(25^x(x+3)) - sqrt(sqrt(13-x) + (4*x-5)^2(2+y))" \\ e=parse(text = text) #Its neccesary to use parse and deparse functions print(deparse(replacePow(e)[[1]])) #printing text with pow
```

replacePowlist

Function for replacing \alpha's to pow inside the list

### **Description**

The function replacing ^'s to pow based on parsing the string vector

### Usage

```
replacePowlist(1,x)
```

### **Arguments**

- list with vectors containing expressions to replace
- x names of the vector inside list

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### Value

The output is the list with replaced ^'s.

### See Also

```
read.slv
import.slv
```

### **Examples**

```
##Example of using the function

temp1<-list(one=c("eee^3","2^rrr"))
replacePowlist(temp1,1)</pre>
```

signup

Up- and down- rounding

### Description

signup rounds the values in its first argument to the specified number of significant digits upwards. signdown rounds the values in its first argument to the specified number of significant digits downwards.

### Usage

```
signup(x, digits = 6)
signdown(x, digits = 6)
```

### **Arguments**

```
x a numeric vector.digits integer indicating the number of significant digits to be used.
```

### See Also

```
signif
```

```
signup(c(1.111, 1.2345e5, 9.8765e-5), 3)
signdown(c(1.111, 1.2345e5, 9.8765e-5), 3)
```

18 write.delim

write.delim

Data output with tab delimiters

### **Description**

prints its required argument x (after converting it to a data frame if it is not one nor a matrix) to a file with tab delimiter without quotes and row names

### Usage

```
write.delim(x, file = "", ...)
```

### Arguments

```
    the object to be written, preferably a matrix or data frame. If not, it is attempted to coerce x to a data frame.
    a character string naming a file
    other arguments passed to write.table
```

### **Details**

```
The function is equivalent to write.table(x = x, file = file, quote = F, sep = "\t", row.names = F, ...)
```

### See Also

```
write.table
```

```
### create and write data.frame
df<-data.frame(number=1:5, words=c("one", "two", "three", "four", "five"))
write.delim(df, "df.txt")</pre>
```

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