Package 'MsBackendSql'

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```
Title SQL-based Mass Spectrometry Data Backend
Version 1.7.1
Description SQL-based mass spectrometry (MS) data backend supporting also
     storange and handling of very large data sets. Objects from
     this package are supposed to be used with the Spectra Bioconductor
     package. Through the MsBackendSql with its minimal memory
     footprint, this package thus provides an alternative MS data
     representation for very large or remote MS data sets.
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Contents

	MsBackendOfflineSql	2
	MsBackendSql	3
Index		12

MsBackendOfflineSql

SQL-based MS backend without active database connection

Description

The MsBackendOfflineSql backend extends the MsBackendSql() backend directly and inherits thus all of its functions as well as properties. The only difference between the two backend is that MsBackendSql keeps an active connection to the SQL database inside the object while the MsBackendOfflineSql backends reconnects to the SQL database for each query. While the performance of the latter is slightly lower (due to the need to connect/disconnect to the database for each function call) it can also be used in a parallel processing environment.

Usage

```
MsBackendOfflineSql()
## S4 method for signature 'MsBackendOfflineSql'
backendInitialize(
  object,
  drv = NULL,
  dbname = character(),
  user = character(),
  password = character(),
  host = character(),
  port = NA_integer_,
  data,
  ...
)
```

Arguments

A MsBackendOfflineSql object. object drv A DBI database driver object (such as SQLite() from the RSQLite package or MariaDB() from the RMariaDB package). See dbConnect() for more informacharacter(1) with the name of the database. Passed directly to dbConnect(). dbname user character(1) with the user name for the database. Passed directly to dbConnect(). character(1) with the password for the database. Note that this password is password stored (unencrypted) within the object. Passed directly to dbConnect(). character(1) with the host running the database. Passed directly to dbConnect(). host integer(1) with the port number (optional). Passed directly to dbConnect(). port For backendInitialize(): optional DataFrame with the full spectra data that data should be inserted into a (new) MsBackendSql database. If provided, it is assumed that the provided database connection information if for a (writeable)

.. ignored.

Creation of backend objects

An empty instance of an MsBackendOfflineSql class can be created using the MsBackendOfflineSql() function. An existing *MsBackendSql* SQL database can be loaded with the backendInitialize() function. This function takes parameters drv, dbname, user, password, host and port, all parameters that are passed to the dbConnect() function to connect to the (existing) SQL database.

empty database into which data should be inserted. data could be the output of

See MsBackendSql() for information on how to create a MsBackend SQL database.

spectraData from another backend.

Author(s)

Johannes Rainer

MsBackendSq1	Spectra MS backend storing data in a SQL database
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Description

The MsBackendSql is an implementation for the MsBackend() class for Spectra() objects which stores and retrieves MS data from a SQL database. New databases can be created from raw MS data files using createMsBackendSqlDatabase().

Usage

```
MsBackendSql()
createMsBackendSqlDatabase(
  dbcon,
  x = character(),
  backend = MsBackendMzR(),
  chunksize = 10L,
  blob = TRUE,
  partitionBy = c("none", "spectrum", "chunk"),
  partitionNumber = 10L
)
## S4 method for signature 'MsBackendSql'
show(object)
## S4 method for signature 'MsBackendSql'
backendInitialize(object, dbcon, data, ...)
## S4 method for signature 'MsBackendSql'
dataStorage(object)
## S4 method for signature 'MsBackendSql'
x[i, j, ..., drop = FALSE]
## S4 method for signature 'MsBackendSql,ANY'
extractByIndex(object, i)
## S4 method for signature 'MsBackendSql'
peaksData(object, columns = c("mz", "intensity"))
## S4 method for signature 'MsBackendSql'
peaksVariables(object)
## S4 replacement method for signature 'MsBackendSql'
intensity(object) <- value</pre>
## S4 replacement method for signature 'MsBackendSql'
mz(object) <- value</pre>
## S4 replacement method for signature 'MsBackendSql'
x$name <- value
## S4 method for signature 'MsBackendSql'
spectraData(object, columns = spectraVariables(object))
## S4 method for signature 'MsBackendSql'
reset(object)
```

```
## S4 method for signature 'MsBackendSql'
spectraNames(object)
## S4 replacement method for signature 'MsBackendSql'
spectraNames(object) <- value</pre>
## S4 method for signature 'MsBackendSql'
filterMsLevel(object, msLevel = uniqueMsLevels(object))
## S4 method for signature 'MsBackendSql'
filterRt(object, rt = numeric(), msLevel. = integer())
## S4 method for signature 'MsBackendSql'
filterDataOrigin(object, dataOrigin = character())
## S4 method for signature 'MsBackendSql'
filterPrecursorMzRange(object, mz = numeric())
## S4 method for signature 'MsBackendSql'
filterPrecursorMzValues(object, mz = numeric(), ppm = 20, tolerance = 0)
## S4 method for signature 'MsBackendSql'
uniqueMsLevels(object, ...)
## S4 method for signature 'MsBackendSql'
backendMerge(object, ...)
## S4 method for signature 'MsBackendSql'
precScanNum(object)
## S4 method for signature 'MsBackendSql'
centroided(object)
## S4 method for signature 'MsBackendSql'
smoothed(object)
## S4 method for signature 'MsBackendSql'
tic(object, initial = TRUE)
## S4 method for signature 'MsBackendSql'
supportsSetBackend(object, ...)
## S4 method for signature 'MsBackendSql'
backendBpparam(object, BPPARAM = bpparam())
## S4 method for signature 'MsBackendSql'
dbconn(x)
```

Arguments

dbcon Connection to a database.

x For createMsBackendSqlDatabase(): character with the names of the raw

data files from which the data should be imported. For other methods an MsqlBackend

instance.

backend For createMsBackendSqlDatabase(): MS backend that can be used to import

MS data from the raw files specified with parameter x.

chunksize For createMsBackendSqlDatabase(): integer(1) defining the number of

input that should be processed per iteration. With chunksize = 1 each file specified with x will be imported and its data inserted to the database. With chunksize = 5 data from 5 files will be imported (in parallel) and inserted to the database. Thus, higher values might result in faster database creation, but

require also more memory.

blob For createMsBackendSqlDatabase(): logical(1) whether individual m/z

and intensity values should be stored separately (blob = FALSE) or if the m/z and intensity values for each spectrum should be stored as a single BLOB SQL

data type (blob = TRUE, the default).

partitionBy For createMsBackendSqlDatabase(): character(1) defining if and how the

peak data table should be partitioned. "none" (default): no partitioning, "spectrum": peaks are assigned to the partition based on the spectrum ID (number), i.e. spectra are evenly (consecutively) assigned across partitions. For partitionNumber = 3, the first spectrum is assigned to the first partition, the second to the second, the third to the third and the fourth spectrum again to the first partition. "chunk": spectra processed as part of the same *chunk* are placed into the same partition. All spectra from the next processed chunk are assigned to the next partition. Note that this is only available for MySQL/MariaDB databases, i.e.,

if con is a MariaDBConnection. See details for more information.

partitionNumber

For createMsBackendSqlDatabase(): integer(1) defining the number of

partitions the database table will be partitioned into (only supported for MySQL/MariaDB

databases).

object A MsBackendSql instance.

data For backendInitialize(): optional DataFrame with the full spectra data that

should be inserted into a (new) MsBackendSql database. If provided, it is assumed that dbcon is a (writeable) connection to an empty database into which data should be inserted. data could be the output of spectraData from another

backend.

.. For [: ignored. For backendInitialize, if parameter data is used: additional

parameters to be passed to the function creating the database such as blob.

i For [: integer or logical to subset the object.

j For [: ignored.

drop For [: logical(1), ignored.

columns For spectraData(): character() optionally defining a subset of spectra vari-

ables that should be returned. Defaults to columns = spectraVariables(object)

hence all variables are returned. For peaksData accessor: optional character with requested columns in the individual matrix of the returned list. Defaults to columns = c("mz", "intensity") but all columns listed by peaksVariables

would be supported.

value For all setter methods: replacement value.

name For <-: character(1) with the name of the spectra variable to replace.

msLevel For filterMsLevel(): integer specifying the MS levels to filter the data.

rt For filterRt(): numeric(2) with the lower and upper retention time. Spectra

with a retention time \geq = rt[1] and \leq = rt[2] are returned.

msLevel. For filterRt(): integerwith the MS level(s) on which the retention time filter should be a

dataOrigin For filterDataOrigin(): character with data origin values to which the

data should be subsetted.

mz For filterPrecursorMzRange(): numeric(2) with the desired lower and up-

per limit of the precursor m/z range. For filterPrecursorMzValues(): numeric

with the m/z value(s) to filter the object.

ppm For filterPrecursorMzValues(): numeric with the m/z-relative maximal ac-

ceptable difference for a m/z value to be considered matching. Can be of length

1 or equal to length(mz).

tolerance For filterPrecursorMzValues(): numeric with the absolute difference for

m/z values to be considered matching. Can be of length 1 or equal to length (mz).

initial For tic(): logical(1) whether the original total ion count should be returned

(initial = TRUE, the default) or whether it should be calculated on the spectras'

intensities (initial = FALSE).

BPPARAM for backendBpparam(): BiocParallel parallel processing setup. See bpparam()

for more information.

Details

The MsBackendSql class is principally a *read-only* backend but by extending the MsBackendCached() backend from the Spectra package it allows changing and adding (**temporarily**) spectra variables **without** changing the original data in the SQL database.

Value

See documentation of respective function.

Creation of backend objects

New backend objects can be created with the MsBackendSql() function. SQL databases can be created and filled with MS data from raw data files using the createMsBackendSqlDatabase() function or using backendInitialize() and providing all data with parameter data. In addition it is possible to create a database from a Spectra object changing its backend to a MsBackendSql or MsBackendOfflineSql using the setBackend() function. Existing SQL databases (created previously with createMsBackendSqlDatabase() or backendInitialize() with the data parameter) can be loaded using the *conventional* way to create/initialize MsBackend classes, i.e. using backendInitialize().

• createMsBackendSqlDatabase(): create a database and fill it with MS data. Parameter dbcon is expected to be a database connection, parameter x a character vector with the file names from which to import the data. Parameter backend is used for the actual data import and defaults to backend = MsBackendMzR() hence allowing to import data from mzML, mzXML or netCDF files. Parameter chunksize allows to define the number of files (x) from which the data should be imported in one iteration. With the default chunksize = 10L data is imported from 10 files in x at the same time (if backend supports it even in parallel) and this data is then inserted into the database. Larger chunk sizes will require more memory and also larger disk space (as data import is performed through temporary files) but might eventually be faster. Parameter blob allows to define whether m/z and intensity values from a spectrum should be stored as a BLOB SQL data type in the database (blob = TRUE, the default) or if individual m/z and intensity values for each peak should be stored separately (blob = FALSE). The latter case results in a much larger database and slower performance of the peaksData function, but would allow to define custom (manual) SQL queries on individual peak values. While data can be stored in any SQL database, at present it is suggested to use MySQL/MariaDB databases. For dbcon being a connection to a MySQL/MariaDB database, the tables will use the ARIA engine providing faster data access and will use table partitioning: tables are splitted into multiple partitions which can improve data insertion and index generation. Partitioning can be defined with the parameters partitionBy and partitionNumber. By default partitionBy = "none" no partitioning is performed. For blob = TRUE partitioning is usually not required. Only for blob = FALSE and very large datasets it is suggested to enable table partitioning by selecting either partitionBy = "spectrum" or partitionBy = "chunk". The first option assignes consecutive spectra to different partitions while the latter puts spectra from files part of the same *chunk* into the same partition. Both options have about the same performance but partitionBy = "spectrum" requires less disk space. Note that, while inserting the data takes a considerable amount of time, also the subsequent creation of database indices can take very long (even longer than data insertion for blob = FALSE).

- backendInitialize(): get access and initialize a MsBackendSql object. Parameter object is supposed to be a MsBackendSql instance, created e.g. with MsBackendSql(). Parameter dbcon is expected to be a connection to an existing MsBackendSql SQL database (created e.g. with createMsBackendSqlDatabase()). backendInitialize() can alternatively also be used to create a new MsBackendSql database using the optional data parameter. In this case, dbcon is expected to be a writeable connection to an empty database and data a DataFrame with the full spectra data to be inserted into this database. The format of data should match the format of the DataFrame returned by the spectraData() function and requires columns "mz" and "intensity" with the m/z and intensity values of each spectrum. The backendInitialize() call will then create all necessary tables in the database, will fill these tables with the provided data and will return an MsBackendSql for this database. Thus, the MsBackendSql supports the setBackend method from Spectra to change from (any) backend to a MsBackendSql. Note however that chunk-wise (or parallel) processing needs to be disabled in this case by passing eventually f = factor() to the setBackend() call.
- supportsSetBackend(): whether MsBackendSql supports the setBackend() method to change the MsBackend of a Spectra object to a MsBackendSql. Returns TRUE, thus, changing the backend to a MsBackendSql is supported if a writeable database connection is provided in addition with parameter dbcon (i.e. setBackend(sps, MsBackendSql(), dbcon = con) with con being a connection to an empty database would store the full spectra data from the Spectra object sps into the specified database and would return a Spectra object that uses a MsBackendSql).

backendBpparam(): whether a MsBackendSql supports parallel processing. Takes a MsBackendSql and a parallel processing setup (see bpparam() for details) as input and always returns a SerialParam() since MsBackendSql does not support parallel processing.

• dbconn(): returns the connection to the database.

Subsetting, merging and filtering data

MsBackendSql objects can be subsetted using the [or extractByIndex() functions. Internally, this will simply subset the integer vector of the primary keys and eventually cached data. The original data in the database **is not** affected by any subsetting operation. Any subsetting operation can be *undone* by resetting the object with the reset() function. Subsetting in arbitrary order as well as index replication is supported.

Multiple MsBackendSql objects can also be merged (combined) with the backendMerge() function. Note that this requires that all MsBackendSql objects are connected to the **same** database. This function is thus mostly used for combining MsBackendSql objects that were previously splitted using e.g. split().

In addition, MsBackendSql supports all other filtering methods available through MsBackendCached(). Implementation of filter functions optimized for MsBackendSql objects are:

- filterDataOrigin(): filter the object retaining spectra with dataOrigin spectra variable values matching the provided ones with parameter dataOrigin. The function returns the results in the order of the values provided with parameter dataOrigin.
- filterMsLevel(): filter the object based on the MS levels specified with parameter msLevel. The function does the filtering using SQL queries. If "msLevel" is a *local* variable stored within the object (and hence in memory) the default implementation in MsBackendCached is used instead.
- filterPrecursorMzRange(): filters the data keeping only spectra with a precursorMz within the m/z value range provided with parameter mz (i.e. all spectra with a precursor m/z >= mz[1L] and <= mz[2L]).
- filterPrecursorMzValues(): filters the data keeping only spectra with precursor m/z values matching the to use different values for ppm and tolerance for each provided m/z value.
- filterRt(): filter the object keeping only spectra with retention times within the specified retention time range (parameter rt). Optional parameter msLevel. allows to restrict the retention time filter only on the provided MS level(s) returning all spectra from other MS levels.

Accessing and modifying data

The functions listed here are specifically implemented for MsBackendSq1. In addition, MsBackendSq1 inherits and supports all data accessor, filtering functions and data manipulation functions from MsBackendCached().

- \$, \$<-: access or set (add) spectra variables in object. Spectra variables added or modified using the \$<- are *cached* locally within the object (data in the database is never changed). To restore an object (i.e. drop all cached values) the reset function can be used.
- dataStorage(): returns a character vector same length as there are spectra in object with the name of the database containing the data.
- intensity<-: not supported.

- mz<-: not supported.
- peaksData(): returns a list with the spectras' peak data. The length of the list is equal to the number of spectra in object. Each element of the list is a matrix with columns according to parameter columns. For an empty spectrum, a matrix with 0 rows is returned. Use peaksVariables(object) to list supported values for parameter columns.
- peaksVariables(): returns a character with the available peak variables, i.e. columns that could be queried with peaksData().
- reset(): restores an MsBackendSql by re-initializing it with the data from the database. Any subsetting or cached spectra variables will be lost.
- spectraData(): gets general spectrum metadata. spectraData() returns a DataFrame with the same number of rows as there are spectra in object. Parameter columns allows to select specific spectra variables.
- spectraNames(), spectraNames<-: returns a character of length equal to the number of spectra in object with the primary keys of the spectra from the database (converted to character). Replacing spectra names with spectraNames<- is not supported.
- uniqueMsLevels(): returns the unique MS levels of all spectra in object.
- tic(): returns the originally reported total ion count (for initial = TRUE) or calculates the total ion count from the intensities of each spectrum (for initial = FALSE).

Implementation notes

Internally, the MsBackendSql class contains only the primary keys for all spectra stored in the SQL database. Keeping only these integer in memory guarantees a minimal memory footpring of the object. Still, depending of the number of spectra in the database, this integer vector might become very large. Any data access will involve SQL calls to retrieve the data from the database. By extending the MsBackendCached() object from the Spectra package, the MsBackendSql supports to (temporarily, i.e. for the duration of the R session) add or modify spectra variables. These are however stored in a data. frame within the object thus increasing the memory demand of the object.

Note

The MsBackendSql backend keeps an (open) connection to the SQL database with the data and hence does not support saving/loading of a backend to disk (e.g. using save or saveRDS). Also, for the same reason, the MsBackendSql does not support parallel processing. The backendBpparam() method for MsBackendSql will thus always return a SerialParam() object.

The MsBackendOfflineSql() could be used as an alternative as it supports saving/loading the data to/from disk and supports also parallel processing.

Author(s)

Johannes Rainer

Examples

```
####
```

Create a new MsBackendSql database

```
## Define a file from which to import the data
data_file <- system.file("microtofq", "MM8.mzML", package = "msdata")</pre>
## Create a database/connection to a database
library(RSQLite)
db_file <- tempfile()</pre>
dbc <- dbConnect(SQLite(), db_file)</pre>
## Import the data from the file into the database
createMsBackendSqlDatabase(dbc, data_file)
dbDisconnect(dbc)
## Initialize a MsBackendSql
dbc <- dbConnect(SQLite(), db_file)</pre>
be <- backendInitialize(MsBackendSql(), dbc)</pre>
be
## Original data source
head(be$dataOrigin)
## Data storage
head(dataStorage(be))
## Access all spectra data
spd <- spectraData(be)</pre>
## Available variables
spectraVariables(be)
## Access mz values
mz(be)
## Subset the object to spectra in arbitrary order
be_sub \leftarrow be[c(5, 1, 1, 2, 4, 100)]
be_sub
## The internal spectrum IDs (primary keys from the database)
be_sub$spectrum_id_
## Add additional spectra variables
be_sub$new_variable <- "B"
## This variable is *cached* locally within the object (not inserted into
## the database)
be_sub$new_variable
```

Index

[,MsBackendSql-method(MsBackendSql),3	filterMsLevel,MsBackendSql-method
<pre>\$<-,MsBackendSql-method (MsBackendSql),</pre>	(MsBackendSq1), 3
3	<pre>filterPrecursorMzRange,MsBackendOfflineSql-method</pre>
<pre>backendBpparam,MsBackendSql-method</pre>	<pre>filterPrecursorMzRange,MsBackendSql-method (MsBackendSql), 3</pre>
• **	filterPrecursorMzValues,MsBackendOfflineSql-method (MsBackendSql), 3
<pre>backendInitialize,MsBackendSql-method (MsBackendSql), 3</pre>	filterPrecursorMzValues,MsBackendSql-method (MsBackendSql), 3
backendMerge,MsBackendOfflineSql-method (MsBackendSql),3	<pre>filterRt,MsBackendOfflineSql-method (MsBackendSql), 3</pre>
backendMerge,MsBackendSql-method (MsBackendSql),3	filterRt, MsBackendSql-method (MsBackendSql), 3
bpparam(), $7,9$	(Habackeriaaq1), 3
	<pre>intensity<-,MsBackendSql-method</pre>
<pre>centroided,MsBackendSql-method (MsBackendSql), 3</pre>	(MsBackendSql), 3
createMsBackendSqlDatabase	MsBackend(), 3
(MsBackendSql), 3	MsBackendCached(), 7, 9, 10
• • • • • • • • • • • • • • • • • • • •	MsBackendOfflineSql, 2
dataStorage,MsBackendOfflineSql-method	MsBackendOfflineSql(), 10
(MsBackendSql), 3	MsBackendOfflineSql-class
dataStorage,MsBackendSql-method	(MsBackendOfflineSql), 2
(MsBackendSql), 3	MsBackendSql, 3
dbconn,MsBackendOfflineSql-method	MsBackendSql(), 2, 3
(MsBackendSql), 3	MsBackendSql-class (MsBackendSql), 3
dbconn,MsBackendSql-method	<pre>mz<-,MsBackendSql-method</pre>
(MsBackendSql), 3	(MsBackendSql), 3
dbConnect(), 3	
	peaksData,MsBackendOfflineSql-method
extractByIndex,MsBackendSql,ANY-method	(MsBackendSql), 3
(MsBackendSql), 3	peaksData,MsBackendSql-method
	(MsBackendSql), 3
filterDataOrigin,MsBackendOfflineSql-method (MsBackendSql), 3	<pre>peaksVariables,MsBackendOfflineSql-method (MsBackendSql), 3</pre>
filterDataOrigin,MsBackendSql-method	peaksVariables,MsBackendSql-method
(MsBackendSq1), 3	(MsBackendSq1), 3
filterMsLevel,MsBackendOfflineSql-method	precScanNum, MsBackendSql-method
(MsBackendSq1), 3	(MsBackendSq1), 3

INDEX 13

```
reset, MsBackendOfflineSql-method
         (MsBackendSql), 3
reset, MsBackendSql-method
         (MsBackendSq1), 3
SerialParam(), 9, 10
setBackend(), 7
{\tt show}, {\tt MsBackendOfflineSql-method}
        (MsBackendSql), 3
show, MsBackendSql-method
         (MsBackendSql), 3
smoothed, MsBackendSql-method
         (MsBackendSql), 3
Spectra(), 3
spectraData, MsBackendOfflineSql-method
        (MsBackendSql), 3
spectraData, MsBackendSql-method
         (MsBackendSql), 3
spectraNames, MsBackendSql-method
         (MsBackendSql), 3
spectraNames<-,MsBackendSql-method</pre>
         (MsBackendSql), 3
supportsSetBackend, MsBackendOfflineSql-method
         (MsBackendSql), 3
{\tt supportsSetBackend}, {\tt MsBackendSql-method}
         (MsBackendSql), 3
\verb+tic,MsBackendOfflineSql-method+
         (MsBackendSql), 3
tic,MsBackendSql-method (MsBackendSql),
unique {\tt MsLevels}, {\tt MsBackendOfflineSql-method}
         (MsBackendSq1), 3
unique {\tt MsLevels}, {\tt MsBackendSql-method}
         (MsBackendSql), 3
```