

Package: CBMutils (via r-universe)

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Type Package

Title Utilities for modelling carbon based on CBM-CFS3

Description Implementation of several components of the Carbon Budget Model of the Canadian Forest Service (v3).

URL <https://CBMutils.predictiveecology.org/>,

<https://github.com/PredictiveEcology/CBMutils>

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License GPL-3

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CBMutils-package CBMutils

Description

Utilities for modelling carbon in R/SpaDES based on the Carbon Budget Model of the Canadian Forest Service v3 (CBM-CFS3).

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See Also

Useful links:

- <https://CBMutils.predictiveecology.org/>
- <https://github.com/PredictiveEcology/CBMutils>

.pooldef	<i>Disturbance matrix pool names and ids</i>
----------	--

Description

TODO: confirm these

Usage

```
.pooldef
.pooldefids
.poolnames
.poolids
```

Format

- An object of class character of length 26.
- An object of class environment of length 26.
- An object of class character of length 25.
- An object of class environment of length 25.

Details

NOTE: A reminder that indexing in R starts at 1, whereas in C++ it starts at 0. Several C++ data structures do not include the Input category, so the indices there are defined using .poolids - 1.

barPlot	barPlot
---------	---------

Description

barPlot

Usage

```
barPlot(cbmPools, masterRaster)
```

Arguments

cbmPools	TODO
masterRaster	A SpatRaster object to use as the template

Value

TODO

biomassTurnoverMatrix *Biomass turnover matrix*

Description

Biomass turnover matrix

Usage

biomassTurnoverMatrix(turnoverParam, PoolCount)

Arguments

turnoverParam	DESCRIPTION NEEDED
PoolCount	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

biomProp *Proportions of total tree biomass in stemwood, bark, branches, and foliage*

Description

Implements equations 4-7 of Boudewyn et al. (2007), used to determine the proportions of total tree biomass in stemwood, bark, branches, and foliage ($p_{stemwood}$, p_{bark} , $p_{branches}$, $p_{foliage}$, respectively), using parameters a , b from Table 6 (table6) and volume-proportion caps from Table 7 (table7).

Usage

biomProp(table6, table7, vol)

Arguments

table6	data.frame corresponding to Table 6 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table6.csv .
table7	data.frame corresponding to Table 7 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table7.csv .
vol	gross merchantable volume per hectare (m^3/ha)

Details

TODO: will eventually add species, ecozone

Value

four-column matrix will columns corresponding to $p_{stemwood}$, p_{bark} , $p_{branches}$, and $p_{foliage}$

References

Boudewyn, P., Song, X., Magnussen, S., & Gillis, M. D. (2007). Model-based, volume-to-biomass conversion for forested and vegetated land in Canada (BC-X-411). Natural Resource Canada, Pacific Forestry Centre. <https://cfs.nrcan.gc.ca/pubwarehouse/pdfs/27434.pdf>

b_m

Calculate stemwood biomass (per ha) of live merchantable trees

Description

Implements equation 1 of Boudewyn et al. (2007) to determines the total stemwood biomass of merchantable trees (in metric tonnes per hectare; T/ha), using parameters a and b from Table 3 (table3).

Usage

```
b_m(table3, vol)
```

Arguments

table3	data.frame corresponding to Table 3 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table3.csv .
vol	gross merchantable volume per hectare (m^3/ha)

Value

stemwood biomass of merchantable trees (b_m in units T/ha)

References

Boudewyn, P., Song, X., Magnussen, S., & Gillis, M. D. (2007). Model-based, volume-to-biomass conversion for forested and vegetated land in Canada (BC-X-411). Natural Resource Canada, Pacific Forestry Centre. <https://cfs.nrcan.gc.ca/pubwarehouse/pdfs/27434.pdf>

`calcC`

Sum carbon for totalCarbon or aboveGround or belowGround

Description

Sum carbon for totalCarbon or aboveGround or belowGround

Usage

```
calcC(cbmPools, poolToSum, masterRaster)
```

Arguments

cbmPools	DESCRIPTION NEEDED
poolToSum	DESCRIPTION NEEDED
masterRaster	A SpatRaster object to use as the template

Value

DESCRIPTION NEEDED

`calcTurnoverRates`

Calculate turnover rates

Description

For all spatial units in the simulation, calculate turnover rates.

Usage

```
calcTurnoverRates(turnoverRates, spatialUnitIds, spatialUnits)
```

Arguments

turnoverRates	DESCRIPTION NEEDED
spatialUnitIds	DESCRIPTION NEEDED
spatialUnits	DESCRIPTION NEEDED

Value

extracts the turnover rates for the specific SPU you are in. These are used to in the core module to calculate the specific rates, which are the used to calculate Net Primary Productivity (NPP) both in the core module and in the next function.

carbonOutPlot	carbonOutPlot
---------------	---------------

Description

carbonOutPlot

Usage

carbonOutPlot(emissionsProducts, masterRaster)

Arguments

emissionsProducts

TODO

masterRaster A SpatRaster object to use as the template

Value

invoked for side effect of creating plot

checkProp	<i>check that the proportions add up to 1</i>
-----------	---

Description

check that the proportions add up to 1

Usage

checkProp(DTofMatrices, tarCols)

Arguments

DTofMatrices DESCRIPTION NEEDED

tarCols DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

checkTransactions	<i>Verify and check transactions</i>
-------------------	--------------------------------------

Description

Post-simulations, this function takes the last two years of simulations, and verifies that the C transfers are as expected on n random pixelGroups. It randomly selects three pixelGroups, calculated the expected transactions from time end(sim) and time end(sim) - 1, and compares them to the last year of simulations. Cases of disturbed, non disturbed, sw, hw, young and old stands were tested.

Usage

```
checkTransactions(  
  cbmPools,  
  pixelKeep,  
  opMatrixCBM,  
  allProcesses,  
  pixelGroupC,  
  pooldef,  
  timeSim,  
  outputsSim,  
  n = 3  
)
```

Arguments

cbmPools	from sim\$cbmPools
pixelKeep	from sim\$pixelKeep
opMatrixCBM	from sim\$opMatrixCBM
allProcesses	from sim\$allProcesses
pixelGroupC	from sim\$pixelGroupC
pooldef	from sim\$pooldef
timeSim	from time(sim)
outputsSim	from outputs(sim)
n	number of pixelGroups (i.e., size of subset) to check

Structure

1. Read-in data
2. randomly select 3 pixelGroups
3. read-in carbon pools at timeSim and timeSim - 1 for the three pixelGroup
4. pull-out matrices applied to these three pixelGroups between t0 and t1

5. matrices into a `data.table`
6. prep pools at `t0`
7. merge three pools at `t0` with their respective transaction matrices ids
8. apply all matrices (8.1:8.9)

TODO: I am passing a whole `simList` here for simplification, but only these objects needed:

1. `$cbmPools;`
2. `$pixelKeep;`
3. `$opMatrixCBM;`
4. `$allProcesses;`
5. `$pooldef.`

TODO: there are still a bunch of checks that need to be changed to assertions (if statements?)

All arguments come from a `simList` object of the same name, except `timeSim` (which comes from `time(sim)`) and `outputsSim` (which comes from `outputs(sim)`).

`computeBioTurnoverMatrices`

Compute Turnover Matrices

Description

Compute Turnover Matrices

Usage

```
computeBioTurnoverMatrices(turnoverParameters, PoolCount)
```

Arguments

<code>turnoverParameters</code>	DESCRIPTION NEEDED
<code>PoolCount</code>	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

```
computeDomDecayMatrices
```

Compute all dom decay matrices in coordinate matrix format

Description

The first column in the specified decayRates parameter acts as the key to each matrix.

Usage

```
computeDomDecayMatrices(decayRates, decayParameters, PoolCount)
```

Arguments

decayRates	matrix of decay rates column 1 is the key for the values in columns 1:n and columns 1:n are the DOM pool specific decay rates.
decayParameters	table of CBM decay parameters
PoolCount	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

```
computeDomTurnoverMatrices
```

Compute DOM turnover matrices

Description

Compute DOM turnover matrices

Usage

```
computeDomTurnoverMatrices(turnoverParameters, PoolCount)
```

Arguments

turnoverParameters	DESCRIPTION NEEDED
PoolCount	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

ComputeGrowthAndDeclineMatrices
Compute growth and decline matrices

Description

DESCRIPTION NEEDED

Arguments

growthIncrements	Matrix of growth increments
decline	Matrix containing proportion decline for each pool

ComputeGrowthAndDeclineMatrices2
Compute growth and decline matrices

Description

DESCRIPTION NEEDED

Usage

```
ComputeGrowthAndDeclineMatrices2(
  growthIncrements,
  ages,
  gcids,
  pools,
  rootParameters,
  turnoverParams,
  biomassToCarbonRate,
  swMult = 1,
  hwMult = 1
)
```

Arguments

growthIncrements	a hash table of growth increments by gcid, by age
ages	the stand age, (the inventory age) stands will be simulated to this age in the final pass
gcids	integer vector of growth curve ids
pools	Matrix (DESCRIPTION NEEDED)

```

rootParameters  data.frame specifying root parameters
turnoverParams 'data.frame specifying turnover parameters
biomassToCarbonRate
                numeric DESCRIPTION NEEDED
swMult          softwood multiplier? (DESCRIPTION NEEDED)
hwMult          hardwood multiplier? (DESCRIPTION NEEDED)

```

ComputeGrowthCoordinateMatrices*Compute growth coordinate matrices***Description**

Transforms growth matrix into coordinate format matrix in terms of CBM pools.

Arguments

pn	PoolNames object
growthIncrements	Matrix of growth increments

ComputeGrowthIncrements*Compute growth increments***Description**

Compute the total growth (above ground and belowground) increment matrices for all stands based on age and hashtable of growth increments by growth curve id.

Usage

```

ComputeGrowthIncrements(
  growthIncrements,
  ages,
  gcids,
  pools,
  rootParameters,
  biomassToCarbonRate,
  swMult = 1,
  hwMult = 1
)

```

Arguments

<code>growthIncrements</code>	a hash table of growth increments by gcid, by age
<code>ages</code>	the stand age, (the inventory age) stands will be simulated to this age in the final pass
<code>gcids</code>	DESCRIPTION NEEDED
<code>pools</code>	DESCRIPTION NEEDED
<code>rootParameters</code>	<code>data.frame</code> (DESCRIPTION NEEDED)
<code>biomassToCarbonRate</code>	DESCRIPTION NEEDED
<code>swMult</code>	softwood multiplier? (DESCRIPTION NEEDED)
<code>hwMult</code>	hardwood multiplier? (DESCRIPTION NEEDED)

`ComputeOvermatureDecline`*Compute overmature decline***Description**

Compute flows to dom pools that occur on a growth curve decline.

Usage

```
ComputeOvermatureDecline(growthIncrements, turnoverParams)
```

Arguments

<code>growthIncrements</code>	Matrix of growth increments
<code>turnoverParams</code>	<code>data.frame</code> specifying turnover parameters

`ComputeOverMatureDeclineCoordinateMatrices`*Compute overmature decline coordinate matrices***Description**

Transforms overmature decline matrix into coordinate format matrix in terms of CBM pools.

Arguments

<code>pn</code>	PoolNames object
<code>decline</code>	Matrix containing proportion decline for each pool

```
computeSlowDecayMatrices
```

Compute slow decay matrices

Description

Compute slow decay matrices

Usage

```
computeSlowDecayMatrices(decayRates, decayParameters, PoolCount)
```

Arguments

decayRates	DESCRIPTION NEEDED
decayParameters	DESCRIPTION NEEDED
PoolCount	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

```
computeSlowMixingMatrix
```

Compute slow mixing matrix

Description

Compute slow mixing matrix

Usage

```
computeSlowMixingMatrix(slowMixingRate, PoolCount)
```

Arguments

slowMixingRate	DESCRIPTION NEEDED
PoolCount	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

convertM3biom

Calculate biomass from gross merchantable volume

Description

Implements the flowchart from figure 3 of Boudewyn et al. (2007) to determined the total above ground biomass (T/ha) from gross merchantable volume (m^3/ha).

Usage

```
convertM3biom(
  meta,
  gCvalues,
  spsMatch,
  ecozones,
  params3,
  params4,
  params5,
  params6,
  params7
)
```

Arguments

meta	DESCRIPTION NEEDED
gCvalues	DESCRIPTION NEEDED
spsMatch	DESCRIPTION NEEDED
ecozones	DESCRIPTION NEEDED
params3	data.frame corresponding to Table 3 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table3.csv .
params4	data.frame corresponding to Table 4 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table4.csv .
params5	data.frame corresponding to Table 5 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table5.csv .
params6	data.frame corresponding to Table 6 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table6.csv .
params7	data.frame corresponding to Table 7 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table7.csv .

Value

three-column matrix with columns corresponding to biomass (T/ha) for total merchantable, foliage, and other.

References

Boudewyn, P., Song, X., Magnussen, S., & Gillis, M. D. (2007). Model-based, volume-to-biomass conversion for forested and vegetated land in Canada (BC-X-411). Natural Resource Canada, Pacific Forestry Centre. <https://cfs.nrcan.gc.ca/pubwarehouse/pdfs/27434.pdf>

cumPoolsCreate *Create cumPools data.table*

Description

Create cumPools data.table

Usage

```
cumPoolsCreate(  
  fullSpecies,  
  gcMeta,  
  userGcM3,  
  stable3,  
  stable4,  
  stable5,  
  stable6,  
  stable7,  
  thisAdmin  
)
```

Arguments

fullSpecies	DESCRIPTION NEEDED
gcMeta	DESCRIPTION NEEDED
userGcM3	DESCRIPTION NEEDED
stable3	DESCRIPTION NEEDED
stable4	DESCRIPTION NEEDED
stable5	DESCRIPTION NEEDED
stable6	DESCRIPTION NEEDED
stable7	DESCRIPTION NEEDED
thisAdmin	DESCRIPTION NEEDED

Value

cumPools data.table

`cumPoolsSmooth` *Smooth the cumPools data.table*

Description

This uses the Chapman Richards equation to smooth the curves that are in `colsToUse`.

Usage

```
cumPoolsSmooth(
  cumPoolsRaw,
  colsToUse = c("totMerch", "fol", "other"),
  colsToUseNew = paste0(colsToUse, "_New")
)
```

Arguments

<code>cumPoolsRaw</code>	Data.table with a numeric column called <code>age`~`</code> and columns called <code>colsToUse`~`</code> .
<code>colsToUse</code>	A character vector of columns to smooth.
<code>colsToUseNew</code>	A character vector of column names for the new smoothed columns

Value

#' A `data.table` with original columns plus new columns named '`colsToUseNew`'.

Author(s)

Celine Boisvenue and Eliot McIntire

`dataset-class` *Dataset class*

Description

DESCRIPTION NEEDED

decayRate	<i>Calculate the decay rate based on mean annual temperature and other parameters</i>
-----------	---

Description

This is the ecological theory for decomposing being used in CBM.

Usage

```
decayRate(meanAnnualTemp, baseDecayRate, q10, tref, max)
```

Arguments

meanAnnualTemp	scalar temperature in degrees Celsius
baseDecayRate	scalar base decay rate constant
q10	the scalar q10 value
tref	the reference temperature
max	the maximum allowed decay rate

Value

the scalar rate of decay

domDecayMatrix	<i>Compute a single dom decay matrix based on the specified table of decay rates</i>
----------------	--

Description

Compute a single dom decay matrix based on the specified table of decay rates

Usage

```
domDecayMatrix(decayRates, decayParameters, PoolCount)
```

Arguments

decayRates	vector of decay rates (each element represents a dom pool)
decayParameters	table of cbm decay parameters
PoolCount	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

`domDecayMatrixItem` *Calculate a portion of the DOM decay matrix*

Description

Calculate a portion of the DOM decay matrix

Usage

```
domDecayMatrixItem(mat, decayRates, propToAtmosphere, src, dst, emission)
```

Arguments

<code>mat</code>	the datatable
<code>decayRates</code>	vector of annual decay rates by dom pool
<code>propToAtmosphere</code>	vector of the proportions of decay emitted to the atmosphere as CO2 by this process
<code>src</code>	the integer code for the dom pool being decayed
<code>dst</code>	the integer code for the dom pool receiving non-emitted decayed matter
<code>emission</code>	the integer code for the CO2 pool

Value

A modified copy of the input `mat`

`domTurnOverMatrix` *DOM turnover matrix*

Description

DOM turnover matrix

Usage

```
domTurnOverMatrix(turnoverParam, PoolCount)
```

Arguments

<code>turnoverParam</code>	DESCRIPTION NEEDED
<code>PoolCount</code>	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

gcidsCreate	<i>Standardized way to create gcids</i>
-------------	---

Description

This uses the underscore delimiter between column names.

Usage

```
gcidsCreate(...)
```

Arguments

... The data.table with ONLY the columns on which to determine unique gcids

getDecayRates	<i>Get decay rates</i>
---------------	------------------------

Description

Returns a vector of decay rates where the indices of the vector are the DOM pools of CBM.

Usage

```
getDecayRates(meanAnnualTemp, decayParameters, domPools)
```

Arguments

meanAnnualTemp scalar temperature in deg Celcius

decayParameters

table of decay parameters for calculating the temperature dependent decay rate

domPools DESCRIPTION NEEDED

Value

the vector of decay rates

```
getIdentityCoordinateMatrix  
    getIdentityCoordinateMatrix
```

Description

DESCRIPTION NEEDED

Usage

```
getIdentityCoordinateMatrix(size)
```

Arguments

size	Numeric indicating the number of rows and columns in the matrix.
------	--

```
getTable  
    getTable
```

Description

DESCRIPTION NEEDED

Usage

```
getTable(filename, dbPath, sqlDir)
```

Arguments

filename	DESCRIPTION NEEDED
dbPath	DESCRIPTION NEEDED
sqlDir	DESCRIPTION NEEDED

hash	<i>Hashing functions</i>
------	--------------------------

Description

DESCRIPTION NEEDED

Usage

hash(x)

matrixHash(x)

Arguments

x DESCRIPTION NEEDED

histDist	<i>Historical disturbances</i>
----------	--------------------------------

Description

Identifies the stand-replacing wildfire disturbance in each spatial unit.

Usage

histDist(mySpu = c(27, 28))

Arguments

mySpu Numeric spatial unit id(s).

Details

Historical disturbances in CBM-CFS3 are used for "filling-up" the soil-related carbon pools. Boudewyn et al. (2007) translate the m3/ha curves into biomass per ha in each of four pools: total biomass for stem wood, total biomass for bark, total biomass for branches and total biomass for foliage. Biomass in coarse and fine roots, in aboveground- and belowground- very-fast, -fast, -slow, in medium-soil, and in snags still needs to be estimated. In all spatial units in Canada, the historical disturbance is set to fire. A stand-replacing fire disturbance is used in a disturb-grow cycle, where stands are disturbed and regrown with turnover, overmature, decay, functioning until the dead organic matter pools biomass values stabilize (+/- 10%).

TODO: (I think but that is in the Rcpp-RCMBGrowthIncrements.cpp so can't check).:

By default the most recent is selected, but the user can change that.

```
loadDisturbanceMatrixIds  
Load disturbance matrix IDs
```

Description

Load disturbance matrix IDs

Usage

```
loadDisturbanceMatrixIds(disturbanceMatrixValues, dbPools)
```

Arguments

disturbanceMatrixValues	DESCRIPTION NEEDED
dbPools	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

```
m3ToBiomIncOnlyPlots m3ToBiomIncOnlyPlots
```

Description

```
m3ToBiomIncOnlyPlots
```

Usage

```
m3ToBiomIncOnlyPlots(inc)
```

Arguments

inc	TODO
-----	------

Value

TODO

m3ToBiomPlots*Plot all columns that are not id_col*

Description

Plot all columns that are not id_col

Usage

```
m3ToBiomPlots(
  inc = "increments",
  id_col = "gcids",
  nrow = 5,
  ncol = 5,
  filenameBase = "rawCumBiomass_",
  path = NULL,
  title = "Cumulative merch fol other by gc id",
  scales = "free_y"
)
```

Arguments

inc	DESCRIPTION NEEDED
id_col	DESCRIPTION NEEDED
nrow	DESCRIPTION NEEDED
ncol	DESCRIPTION NEEDED
filenameBase	DESCRIPTION NEEDED
path	DESCRIPTION NEEDED
title	DESCRIPTION NEEDED
scales	DESCRIPTION NEEDED

m3ToVolCheckPlots*m3ToVolCheckPlots*

Description

Check the growth increments by above ground pool. This is necessary because the Boudevyn et al parameters used to translate the m3/ha into biomass/ha do not always work. Returns a list of plots, each plot show the merch, fol, and other increments for a specific growth curve.

Usage

```
m3ToVolCheckPlots(sim)
```

Arguments

`sim` A SpaDES CBM simulation object (`simList`).

`matrixDT` *Make a data.table out of the carbon transfer matrices*

Description

Make a data.table out of the carbon transfer matrices

Usage

```
matrixDT(matricesIn, indicesIn)
```

Arguments

<code>matricesIn</code>	DESCRIPTION NEEDED
<code>indicesIn</code>	DESCRIPTION NEEDED

Value

a data.table object summarizing the carbon transfers

`nmfac` *Expansion factor for non-merchantable live tree biomass*

Description

Implements equation 2 of Boudewyn et al. (2007), used to determine the total stem wood biomass (in metric tonnes per hectare; T/ha) of non-merchantable trees (B_n), together with the stemwood biomass of live merchantable and non-merchantable trees (B_{nm}), using parameters a , b , and k from Table 4 (table4).

Usage

```
nmfac(table4, eq1, vol)
```

Arguments

<code>table4</code>	data.frame corresponding to Table 4 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table4.csv .
<code>eq1</code>	stemwood biomass of merchantable trees (T/ha) from equation 1 of Boudewyn et al. (2007) (i.e., the result of <code>b_m()</code>).
<code>vol</code>	gross merchantable volume per hectare (m^3/ha)

Value

two-column matrix with columns corresponding to b_n and b_{nm}

References

Boudewyn, P., Song, X., Magnussen, S., & Gillis, M. D. (2007). Model-based, volume-to-biomass conversion for forested and vegetated land in Canada (BC-X-411). Natural Resource Canada, Pacific Forestry Centre. <https://cfs.nrcan.gc.ca/pubwarehouse/pdfs/27434.pdf>

NPP

Calculate post-simulation net primary productivity

Description

Calculate post-simulation net primary productivity

Usage

```
NPP(pools, pixelGroupSpu, year1, year2, turnoverRates)
```

Arguments

pools	DESCRIPTION NEEDED
pixelGroupSpu	DESCRIPTION NEEDED
year1, year2	Consecutive start and end years (e.g., 1990, 1991)
turnoverRates	DESCRIPTION NEEDED

Examples

```
## Not run:
spadesCBMout <- simInitAndSpades() ## TODO: fill this in
NPP(pools = spadesCBMout$cbmPools,
    pixelGroupSpu = spadesCBMout$pixelGroupC[, .(pixelGroup, spatial_unit_id)],
    year1 = 1990, year2 = 1991,
    turnoverRates = spadesCBMout$turnoverRates)

## End(Not run)
```

NPPplot

NPPplot

Description

NPPplot

Usage`NPPplot(spatialDT, NPP, masterRaster)`**Arguments**

<code>spatialDT</code>	TODO
<code>NPP</code>	TODO
<code>masterRaster</code>	A SpatRaster object to use as the template

Value

TODO

plotCarbonRasters

*Plotting pools***Description**

Plotting pools

Usage`plotCarbonRasters(pixelkeep, cbmPools, poolsToPlot, years, masterRaster)`**Arguments**

<code>pixelkeep</code>	DESCRIPTION NEEDED
<code>cbmPools</code>	DESCRIPTION NEEDED
<code>poolsToPlot</code>	DESCRIPTION NEEDED
<code>years</code>	DESCRIPTION NEEDED
<code>masterRaster</code>	A SpatRaster object to use as the template

Examples

```
## Not run:
# include 'totalCarbon' in poolsToPlot to add plot of total carbon
plotCarbonRasters(
  cbmPools = spadesCBMout$cbmPools,
  poolsToPlot = c("totalCarbon", "BelowGroundSlowSoil"),
  masterRaster = spadesCBMout$masterRaster,
  pixelkeep = spadesCBMout$pixelKeep,
  years = c(1990, 2000, 2005)
)
## End(Not run)
```

prepInputsEcozones prepInputsEcozones

Description

`prepInputsEcozones`

Usage

```
prepInputsEcozones(url = NULL, dPath, rasterToMatch)
```

Arguments

<code>url</code>	A URL to the data
<code>dPath</code>	destination path
<code>rasterToMatch</code>	A RasterLayer with NA representing "off study area" and otherwise any arbitrary value for "in study area". Equivalent to <code>rasterToMatch</code> argument in <code>reproducible::prepInputs()</code> .

prepInputsVRI prepInputsVRI

Description

Read in the BC VRI, with growth curve information (from ws3), and creates a raster stack of the age and gcID.

Usage

```
prepInputsVRI(url, dPath, rasterToMatch)
```

Arguments

<code>url</code>	A URL to the data
<code>dPath</code>	destination path
<code>rasterToMatch</code>	A RasterLayer with NA representing "off study area" and otherwise any arbitrary value for "in study area". Equivalent to <code>rasterToMatch</code> argument in reproducible::prepInputs() .

`prepInputsVRIage`*Read in the BC VRI, with growth curve information (from ws3), and creates a raster stack of the age***Description**

Read in the BC VRI, with growth curve information (from ws3), and creates a raster stack of the age

Usage

```
prepInputsVRIage(
  VRIurl,
  dPath,
  rasterToMatch,
  targetFile,
  field = "PROJ_AGE_1"
)
```

Arguments

<code>VRIurl</code>	a URL to the data
<code>dPath</code>	destination path
<code>rasterToMatch</code>	A RasterLayer with NA representing "off study area" and otherwise any arbitrary value for "in study area". Equivalent to <code>rasterToMatch</code> argument in reproducible::prepInputs() .
<code>targetFile</code>	A gdb.zip file that is the inventory file.
<code>field</code>	The age column name in the inventory file.

*query**query*

Description

DESCRIPTION NEEDED

Usage

```
query(dbPath, sql)
```

Arguments

dbPath	Path to sqlite database file.
sql	SQL statement with to execute in the database.

*readSqlFile**readSqlFile*

Description

DESCRIPTION NEEDED

Usage

```
readSqlFile(filePath)
```

Arguments

filePath	Path to sqlite database file.
----------	-------------------------------

*retrieveSpuRaster**Produce a raster with spUnits*

Description

Produce a raster with spUnits

Usage

```
retrieveSpuRaster(  
  spatialUnitsFile = shapefile("data/spUnit_Locator.shp"),  
  UserArea,  
  rasterRes = c(250, 250)  
)
```

Arguments

spatialUnitsFile	DESCRIPTION NEEDED
UserArea	DESCRIPTION NEEDED
rasterRes	The desired raster resolution.

Examples

```
## Not run:
test1 <- raster::shapefile("data/forIan/SK_data/CBM_GIS/SpadesCBM_TestArea.shp")
out1 <- retrieveSpuRaster(UserArea = test1, rasterRes = c(250, 250))
if (interactive()) Plot(out1)

## End(Not run)
```

sapfac

*Expansion factor for sapling-sized trees***Description**

Implements equation 3 of Boudewyn et al. (2007), used to determine the total stem wood biomass (in metric tonnes per hectare; T/ha) of sapling-sized trees (B_s), using parameters a , b , and k from Table 5 (table5).

Usage

```
sapfac(table5, eq2, vol)
```

Arguments

table5	data.frame corresponding to Table 5 from Boudewyn et al. (2007), available from https://nfi.nfis.org/resources/biomass_models/appendix2_table5.csv .
eq2	two-column matrix giving stemwood biomass of non-merchantable trees (i.e., b_n given in units T/ha), and merchantable + non-merchantable trees (i.e., b_{nm} given in units T/ha), from equation 2 of Boudewyn et al. (2007) (i.e., the result of nmfac()).
vol	gross merchantable volume per hectare (m^3/ha)

Value

stemwood biomass of sapling-sized trees (b_s in units T/ha)

References

Boudewyn, P., Song, X., Magnussen, S., & Gillis, M. D. (2007). Model-based, volume-to-biomass conversion for forested and vegetated land in Canada (BC-X-411). Natural Resource Canada, Pacific Forestry Centre. <https://cfs.nrcan.gc.ca/pubwarehouse/pdfs/27434.pdf>

`seeDist`

See disturbances

Description

Get the descriptive name of the disturbance, the source pools, the sink pools, and the proportions transferred.

Usage

```
seeDist(
  distId = c(161, 230, 313, 361),
  dbPath = file.path("data", "cbm_defaults", "cbm_defaults.db")
)
```

Arguments

<code>distId</code>	Description needed
<code>dbPath</code>	Path to sqlite database file.

Value

A list of `data.frames`, one per disturbance matrix id.

`simDist`

get the descriptive name and proportions transferred for disturbances in a simulation requires a simulation list post simulations (from spades()) and returns a list of data.frames. Each data had the descriptive name of a disturbance used in the simulations, the disturbance matrix identification number from cbm_defaults, the pool from which carbon is taken (source pools) in this specific disturbance, the pools into which carbon goes, and the proportion in which the carbon-transfers are completed.

Description

get the descriptive name and proportions transferred for disturbances in a simulation requires a simulation list post simulations (from spades()) and returns a list of `data.frames`. Each data had the descriptive name of a disturbance used in the simulations, the disturbance matrix identification number from `cbm_defaults`, the pool from which carbon is taken (source pools) in this specific disturbance, the pools into which carbon goes, and the proportion in which the carbon-transfers are completed.

Usage

```
simDist(sim)
```

Arguments

sim A SpaDES CBM simulation (`simList`) object.

slowDecayMatrix *Slow decay matrix*

Description

Slow decay matrix

Usage

```
slowDecayMatrix(decayRates, decayParameters, PoolCount)
```

Arguments

decayRates	DESCRIPTION NEEDED
decayParameters	DESCRIPTION NEEDED
PoolCount	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

spatialPlot *spatialPlot*

Description

`spatialPlot`

Usage

```
spatialPlot(pixelkeep, cbmPools, poolsToPlot, years, masterRaster)
```

Arguments

pixelkeep	TODO
cbmPools	TODO
poolsToPlot	TODO
years	TODO
masterRaster	A SpatRaster object to use as the template

Value

TODO

spatialRaster *Plot*

Description

the spatialPlot function plots rasters directly, but does not return rasters, whereas it does so here.

Usage

```
spatialRaster(pixelkeep, cbmPools, poolsToPlot, years, masterRaster)
```

Arguments

pixelkeep	DESCRIPTION NEEDED
cbmPools	DESCRIPTION NEEDED
poolsToPlot	DESCRIPTION NEEDED
years	DESCRIPTION NEEDED
masterRaster	A SpatRaster object to use as the template

Value

RasterLayer but invoked for its side-effect of plotting the rasters.

See Also

spatialPlot

spatialUnitDecayRates *Spatial unit decay rates*

Description

Spatial unit decay rates

Usage

```
spatialUnitDecayRates(climate, decayparameters, domPools)
```

Arguments

climate	DESCRIPTION NEEDED
decayparameters	DESCRIPTION NEEDED
domPools	DESCRIPTION NEEDED

Value

DESCRIPTION NEEDED

Spinup	<i>Spinup</i>
--------	---------------

Description

Spinup a landscape by running rotations of stand replacing disturbances repeatedly until the pre-disturbance slow pools and the last rotation pre-disturbance slow pools are within a tolerance. Stand then grows to inventory age, and is ready for CBM simulation.

Usage

```
Spinup(
  pools,
  opMatrix,
  constantProcesses,
  growthIncrements,
  ages,
  gcids,
  historicdmids,
  lastPassdmids,
  delays,
  minRotations,
  maxRotations,
  returnIntervals,
  rootParameters,
  turnoverParams,
  biomassToCarbonRate,
  debug = FALSE
)
```

Arguments

pools	Matrix (DESCRIPTION NEEDED)
opMatrix	Matrix (DESCRIPTION NEEDED)
constantProcesses	a list of constant process C dynamics matrices

```

growthIncrements           a hash table of growth increments by gcid, by age
ages                      the stand age, (the inventory age) stands will be simulated to this age in the final
                           pass
gcids                     the growth curve ids (referenced by growthIncrements)
historicdmids             (DESCRIPTION NEEDED)
lastPassdmids              (DESCRIPTION NEEDED)
delays                     (DESCRIPTION NEEDED)
minRotations               (DESCRIPTION NEEDED)
maxRotations               (DESCRIPTION NEEDED)
returnIntervals             (DESCRIPTION NEEDED)
rootParameters  data.frame specifying root parameters
turnoverParams   data.frame specifying turnover parameters
biomassToCarbonRate        (DESCRIPTION NEEDED)
debug                     logical indicating whether to run spinup in debug mode.

```

spuDist	<i>Identify the ID number (CBM-CFS3 legacy) possible in the current spatial unit</i>
---------	--

Description

You give is spatial units you are targeting mySpu and it gives you the disturbance matrix id that are possible/default in that specific spu and a descriptive name of that disturbance matrix it creates a data.frame of length number of disturbances, with three columns: `spatial_unit_id`, `disturbance_matrix_id`, and a description of the disturbance.

Usage

```
spuDist(mySpu, dbPath)
```

Arguments

<code>mySpu</code>	Numeric spatial unit id(s).
<code>dbPath</code>	Path to sqlite database file.

Details

TODO: can we have a Canada-wide SPU map and they locate themselves on the map? this needs to be done before simulations are run so the user can provide this info (location info) for the simulations - Ian is working on this.

the function has the defaults from the SK managed forest example. These can be changed by feeding in other SPU.

Examples

```
## Not run:
## using raster
library(terra)
spuRaster <- rast(file.path("data/forIan/SK_data/CBM_GIS/spUnits_TestArea.tif"))
spatial_unit_id <- values(spuRaster) # 28 27
mySpu <- unique(spatial_unit_id)

## using growth curves
f <- file.path("spadesCBMinputs/data/SK_ReclineRuns30m/LookupTables/yieldRCBM.csv")
gcIn <- as.matrix(read.csv(f))
mySpu <- unique(gcIn[, 1])

## End(Not run)
```

StepPools

Step through pools, computing carbon transfers

Description

Computes by row (stand/pixelGroup) across all columns (pools).

Usage

```
StepPools(pools, opMatrix, flowMatrices)
```

Arguments

pools	A numeric matrix describing the amount of carbon in each pool (column) across multiple stands/pixel groups (rows).
opMatrix	A numeric matrix with columns corresponding to flows, and rows corresponding to stands/pixel groups. Values are indices into the flowMatrices.
flowMatrices	A list of lists (or environments) containing all carbon flow matrices.

Examples

```
nPixGrp <- 3

pools <- matrix(c(1.0, 10.0, 0.0, 1.0, 20.0, 0.0, 1.0, 5.0, 0.0),
                 ncol = 3, nrow = nPixGrp, byrow = TRUE)
colnames(pools) <- c("input", "pool1", "pool2")
pools

##      input pool1 pool2
## [1,]     1    10     0
## [2,]     1    20     0
## [3,]     1     5     0
```

```

op <- matrix(rep(c(1, 1), nPixGrp), ncol = 2, nrow = nPixGrp, byrow = TRUE)
colnames(op) <- c("disturbance", "growth")

cnames <- c("row", "col", "value")
dist <- matrix(c(2, 3, 1, 1, 1, 1), ncol = 3, nrow = 2, byrow = TRUE)
colnames(dist) <- cnames
grow <- matrix(c(1, 2, 0.1, 1, 3, 0.2, 2, 3, 0.3, 3, 3, 1.0), ncol = 3, nrow = 4, byrow = TRUE)
colnames(grow) <- cnames
flow <- list(Disturbance = list(dist), Growth = list(grow))

new_pools <- StepPools(pools, op, flow)
new_pools

##      input pool1 pool2
## [1,]    1   0.1  10.2
## [2,]    1   0.1  20.2
## [3,]    1   0.1   5.2

```

StepPoolsRef*Step through pools, computing carbon transfers***Description**

Computes by row (stand/pixelGroup) across all columns (pools) in place.

Usage

```
StepPoolsRef(pools, opMatrix, flowMatrices)
```

Arguments

- | | |
|--------------|--|
| pools | A numeric matrix describing the amount of carbon in each pool (column) across multiple stands/pixel groups (rows). |
| opMatrix | A numeric matrix with columns corresponding to flows, and rows corresponding to stands/pixel groups. Values are indices into the flowMatrices. |
| flowMatrices | A list of lists (or environments) containing all carbon flow matrices. |

ws3Build**ws3Build****Description**

ws3, the harvest optimizer created by Greg Paradis provided the BC disturbance rasters (harvest or harvest and fires) per TSA and not for the whole study area. This function reads-in the file structure, pulls in all the TSA-level rasters, knits them together and post-processes them to match the rasterToMatch.

Usage

```
ws3Build(masterRaster, tsaDirs, years, pathsTifs)
```

Arguments

masterRaster	A SpatRaster object to use as the template
tsaDirs	character vector of paths to the TSA directories
years	numeric or integer vector specifying the years for the data
pathsTifs	character vector of paths for the geotiff files

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