

# Package ‘RStoolbox’

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---

 classifyQA

*Classify Landsat QA bands*


---

**Description**

extracts five classes from QA band: background, cloud, cirrus, snow and water.

**Usage**

```
classifyQA(img, type = c("background", "cloud", "cirrus", "snow",
  "water"), confLayers = FALSE, sensor = "OLI",
  legacy = "collection1", ...)
```

**Arguments**

img	RasterLayer. Landsat 8 OLI QA band.
type	Character. Classes which should be returned. One or more of c("background", "cloud", "cirrus", "snow", "water").
confLayers	Logical. Return one layer per class classified by confidence levels, i.e. cloud:low, cloud:med, cloud:high.
sensor	Sensor to encode. Options: c("OLI", "TIRS", "ETM+", "TM", "MSS").
legacy	Encoding systematic Options: c("collection1", "pre_collection"). Default is "collection1" for the Landsat Collection 1 8-bit quality designations. Use "pre_collection" for imagery downloaded before the Collection 1 quality designations were introduced
...	further arguments passed to <a href="#">writeRaster</a>

**Details**

By default each class is queried for \*high\* confidence. See [encodeQA](#) for details. To return the different confidence levels per condition use confLayers=TRUE. This approach corresponds to the way LandsatLook Quality Images are produced by the USGS.

**Value**

Returns a RasterLayer with maximal five classes:

class	value
background	1L
cloud	2L
cirrus	3L
snow	4L
water	5L

Values outside of these classes are returned as NA. If `confLayers = TRUE` then a RasterStack with one layer per condition (except 'background') is returned, whereby each layer contains the confidence level of the condition.

Confidence	value
low	1L
med	2L
high	3L

### See Also

[encodeQA](#) [decodeQA](#)

### Examples

```
library(raster)
qa <- raster(ncol = 100, nrow=100, val = sample(1:2^14, 10000))

## QA classes
qacs <- classifyQA(img = qa)
## Confidence levels
qacs_conf <- classifyQA(img = qa, confLayers = TRUE)
```

---

cloudMask

*Simple Cloud Detection*

---

### Description

Developed for use with Landsat data cloudMask relies on the distinctive difference between the blue (or any other short-wave band) and thermal band for semi-automated creation of a cloud mask. Since it relies on thermal information it doesn't work well for sensors without thermal bands.

### Usage

```
cloudMask(x, threshold = 0.8, blue = "B1_sre", tir = "B6_sre",
  buffer = NULL, plot = FALSE, verbose)
```

### Arguments

<code>x</code>	RasterBrick or RasterStack with reflectance and brightness temperature OR the mask of a previous run of cloudMask with <code>returnDiffLayer=TRUE</code> .
<code>threshold</code>	Numeric. cloud detection threshold. If not provided it will be guessed. Everything <i>below</i> this threshold will be considered a cloud pixel (unless it is removed by filtering afterwards).
<code>blue</code>	Character or integer. Bandname or number for the blue band
<code>tir</code>	Character or integer. Bandname or number for the thermal band

buffer	Integer. Number of pixels to use as a buffer that will be added to the identified cloud centers.
plot	Logical. Plots of the cloud mask for all sub-steps (sanitizing etc.) Helpful to find proper parametrization.
verbose	Logical. Print messages or suppress.

**Value**

Returns a RasterStack with two layers: CMASK contains the binary cloud mask (1 = cloud, NA = not-cloud) and NDTCI contains the cloud index.

**Note**

Typically clouds are cold in the thermal region and have high reflectance in short wavelengths (blue). By calculating a normalized difference index between the two bands and thresholding a rough cloud mask can be obtained. Before calculating the spectral cloud index (let's call it Normalized Difference Thermal Cloud Index (NDTCI)) the thermal band will be matched to the same value range as the blue band. Therefore, it doesn't matter whether you provide DN, radiance or brightness temperature.

This approach to cloud masking is very simplistic. And aims only at rough removal of potentially clouded areas. Nevertheless, it is able to provide satisfactory results. More sophisticated approaches, including cloud cast shadow detection can be found elsewhere, e.g. [fmask](#).

It can make sense to find a suitable threshold on a cropped version of the scene. Also make sure you make use of the `returnDiffLayer` argument to save yourself one processing step. Buffering should be seen as final polishing, i.e. as long as the pure cloud centers are not detected properly, you might want to turn it off. since it takes some time to calculate. Once your mask detects obvious cloud pixels properly re-enable buffering for fine tuning if desired. Finally, once a suitable threshold is established re-run `cloudMask` on the whole scene with this threshold and go get a coffee.

**See Also**

[cloudShadowMask](#)

**Examples**

```
library(raster)
library(ggplot2)
## Import Landsat example subset
data(lsat)
## We have two tiny clouds in the east
ggRGB(lsat, stretch = "lin")

## Calculate cloud index
cldmsk <- cloudMask(lsat, blue = 1, tir = 6)
ggR(cldmsk, 2, geom_raster = TRUE)

## Define threshold (re-use the previously calculated index)
## Everything above the threshold is masked
```

```

## In addition we apply a region-growing around the core cloud pixels
cldmsk_final <- cloudMask(cldmsk, threshold = 0.1, buffer = 5)

## Plot cloudmask
ggRGB(lsat, stretch = "lin") +
  ggR(cldmsk_final[[1]], ggLayer = TRUE, forceCat = TRUE, geom_raster = TRUE) +
  scale_fill_manual(values = "red", na.value = NA)

#' ## Estimate cloud shadow displacement
## Interactively (click on cloud pixels and the corresponding shadow pixels)
## Not run: shadow <- cloudShadowMask(lsat, cldmsk_final, nc = 2)

## Non-interactively. Pre-defined shadow displacement estimate (shiftEstimate)
shadow <- cloudShadowMask(lsat, cldmsk_final, shiftEstimate = c(-16,-6))

## Plot
csmask <- raster::merge(cldmsk_final[[1]], shadow)
ggRGB(lsat, stretch = "lin") +
  ggR(csmask, ggLayer = TRUE, forceCat = TRUE, geom_raster = TRUE) +
  scale_fill_manual(values = c("blue", "yellow"),
    labels = c("shadow", "cloud"), na.value = NA)

```

---

cloudShadowMask

*Cloud Shadow Masking for Flat Terrain*


---

## Description

Intended for interactive use, `cloudShadowMask` will ask the user to select a few corresponding cloud/cloudShadow pixels which will be used to estimate coordinates for a linear cloudmask shift.

## Usage

```
cloudShadowMask(img, cm, nc = 5, shiftEstimate = NULL,
  preciseShift = NULL, quantile = 0.2, returnShift = FALSE)
```

## Arguments

<code>img</code>	Raster* object containing the scene
<code>cm</code>	Raster* object. Cloud mask (typically the result of <code>cloudMask</code> )
<code>nc</code>	Integer. Number of control points. A few points (default) are fine because the final shift is estimated by <code>coregisterImages</code> .
<code>shiftEstimate</code>	NULL or numeric vector of length two (x,y). Estimated displacement of shadows in map units. If NULL, the user will be asked to select control points interactively.
<code>preciseShift</code>	NULL or numeric vector of length two (x,y). Use this if cloud/cloud-shadow displacement is already known, e.g. from a previous run of <code>cloudShadowMask</code> .

quantile	Numeric (between 0 and 1). Quantile threshold used for image co-registration. By default the 20% quantile of the total intensity (sum) of the image is used as potential shadow mask.
returnShift	Logical. Return a numeric vector containing the shift parameters. Useful if you estimate parameters on a subset of the image.

### Details

This is a very simplistic approach to cloud shadow masking (simple shift of the cloud mask). It is not image based and accuracy will suffer from clouds at different altitudes. However, just as cloudMask this is a quick and easy to use tool for Landsat data if you're just working on a few scenes and don't have fMask or CDR data at hand. Although for some test scenes it does perform surprisingly well.

### Value

Returns a RasterLayer with the cloud shadow mask (0 = shadow, NA = not-shadow).

### See Also

[cloudMask](#)

### Examples

```
library(raster)
library(ggplot2)
## Import Landsat example subset
data(lsat)
## We have two tiny clouds in the east
ggRGB(lsat, stretch = "lin")

## Calculate cloud index
cldmsk <- cloudMask(lsat, blue = 1, tir = 6)
ggR(cldmsk, 2, geom_raster = TRUE)

## Define threshold (re-use the previously calculated index)
## Everything above the threshold is masked
## In addition we apply a region-growing around the core cloud pixels
cldmsk_final <- cloudMask(cldmsk, threshold = 0.1, buffer = 5)

## Plot cloudmask
ggRGB(lsat, stretch = "lin") +
  ggR(cldmsk_final[[1]], ggLayer = TRUE, forceCat = TRUE, geom_raster = TRUE) +
  scale_fill_manual(values = "red", na.value = NA)

#' ## Estimate cloud shadow displacement
## Interactively (click on cloud pixels and the corresponding shadow pixels)
## Not run: shadow <- cloudShadowMask(lsat, cldmsk_final, nc = 2)

## Non-interactively. Pre-defined shadow displacement estimate (shiftEstimate)
```

```
shadow <- cloudShadowMask(lsat, cldmsk_final, shiftEstimate = c(-16,-6))

## Plot
csmask <- raster::merge(cldmsk_final[[1]], shadow)
ggRGB(lsat, stretch = "lin") +
  ggR(csmask, ggLayer = TRUE, forceCat = TRUE, geom_raster = TRUE) +
  scale_fill_manual(values = c("blue", "yellow"),
    labels = c("shadow", "cloud"), na.value = NA)
```

---

 coregisterImages

*Image to Image Co-Registration based on Mutual Information*


---

## Description

Shifts a slave image to match the reference image (master). Match is based on maximum mutual information.

## Usage

```
coregisterImages(slave, master, shift = 3, shiftInc = 1,
  nSamples = 1e+05, reportStats = FALSE, verbose, nBins = 100, ...)
```

## Arguments

slave	Raster* object. Slave image to shift to master. Slave and master must have equal numbers of bands.
master	Raster* object. Reference image. Slave and master must have equal numbers of bands.
shift	Numeric or matrix. If numeric, then shift is the maximal absolute radius (in pixels of slave resolution) which slave is shifted (seq(-shift, shift, by=shiftInc)). If shift is a matrix it must have two columns (x shift and y shift), then only these shift values will be tested.
shiftInc	Numeric. Shift increment (in pixels, but not restricted to integer). Ignored if shift is a matrix.
nSamples	Integer. Number of samples to calculate mutual information.
reportStats	Logical. If FALSE it will return only the shifted images. Otherwise it will return the shifted image in a list containing stats such as mutual information per shift and joint histograms.
verbose	Logical. Print status messages. Overrides global RStoolbox.verbose option.
nBins	Integer. Number of bins to calculate joint histogram.
...	further arguments passed to <a href="#">writeRaster</a> .



## Details

Currently only a simple linear x - y shift is considered and tested. No higher order shifts (e.g. rotation, non-linear transformation) are performed. This means that your imagery should already be properly geometrically corrected.

**Mutual information** is a similarity metric originating from information theory. Roughly speaking, the higher the mutual information of two data-sets, the higher is their shared information content, i.e. their similarity. When two images are exactly co-registered their mutual information is maximal. By trying different image shifts, we aim to find the best overlap which maximises the mutual information.

## Value

reportStats=FALSE returns a Raster\* object (x-y shifted slave image). reportStats=TRUE returns a list containing a data.frame with mutual information per shift (\$MI), the shift of maximum MI (\$bestShift), the joint histograms per shift in a list (\$jointHist) and the shifted image (\$coregImg).

## Examples

```
library(raster)
library(ggplot2)
library(reshape2)
data(rlogo)
reference <- rlogo
## Shift reference 2 pixels to the right and 3 up
missreg <- shift(reference, x = 2, y = 3)

## Compare shift
p <- ggR(reference, sat = 1, alpha = .5)
p + ggR(missreg, sat = 1, hue = .5, alpha = 0.5, ggLayer=TRUE)

## Coregister images (and report statistics)
coreg <- coregisterImages(missreg, master = reference,
                          nSamples = 500, reportStats = TRUE)

## Plot mutual information per shift
ggplot(coreg$MI) + geom_raster(aes(x,y,fill=mi))

## Plot joint histograms per shift (x/y shift in facet labels)

df <- melt(coreg$jointHist)
df$L1 <- factor(df$L1, levels = names(coreg$jointHist))
df[df$value == 0, "value"] <- NA ## don't display p = 0
ggplot(df) + geom_raster(aes(x = Var2, y = Var1, fill=value)) + facet_wrap(~L1) +
  scale_fill_gradientn(name = "p", colours = heat.colors(10), na.value = NA)

## Compare correction
ggR(reference, sat = 1, alpha = .5) +
  ggR(coreg$coregImg, sat = 1, hue = .5, alpha = 0.5, ggLayer=TRUE)
```

---

`decodeQA`*Decode QA flags to bit-words*

---

**Description**

Intended for use with Landsat 16-bit QA bands. Decodes pixel quality flags from integer to bit-words.

**Usage**

```
decodeQA(x)
```

**Arguments**

x	Integer (16bit)
---	-----------------

**See Also**

[encodeQA](#)

**Examples**

```
decodeQA(53248)
```

---

`encodeQA`*Encode QA Conditions to Integers*

---

**Description**

Intended for use with Landsat 16-bit QA bands. Converts pixel quality flags from human readable to integer, which can then be used to subset a QA image. Please be aware of the default settings which differ for different parameters. Depending on, which sensor and legacy is selected, some quality parameters are not used, since the sequences of available bitwise quality designations differ per sensor and collection.

**Usage**

```
encodeQA(fill = "no", terrainOcclusion = "no", radSaturation = "na",  
cloudMask = "all", cloud = "all", cloudShadow = "all",  
snow = "all", cirrus = "all", droppedPixel = "no", water = "all",  
droppedFrame = "no", sensor = "OLI", legacy = "collection1")
```

**Arguments**

fill	Designated fill. Options: c("yes", "no", "all").
terrainOcclusion	Terrain induced occlusion. Options: c("yes", "no", "all").
radSaturation	Number of bands that contain radiometric saturation. Options: c("na", "low", "med", "high", "all") for no bands, 1-2 bands, 3-4 bands, 5 or more bands contain saturation.
cloudMask	Cloud mask. Options: c("yes", "no", "all").
cloud	Cloud confidence. Options: c("na", "low", "med", "high", "all").
cloudShadow	Cloud shadow confidence. Options: c("yes", "no", "all").
snow	Snow / ice confidence. Options: c("na", "low", "med", "high", "all").
cirrus	Cirrus confidence. Options: c("na", "low", "med", "high", "all").
droppedPixel	Dropped pixel. Options: c("yes", "no", "all").
water	Water confidence. Options: c("na", "low", "med", "high", "all").
droppedFrame	Dropped frame. Options: c("yes", "no", "all").
sensor	Sensor to encode. Options: c("OLI", "TIRS", "ETM+", "TM", "MSS").
legacy	Encoding systematic Options: c("collection1", "pre_collection"). Default is "collection1" for the Landsat Collection 1 8-bit quality designations. Use "pre_collection" for imagery downloaded before the Collection 1 quality designations were introduced

**Value**

Returns the Integer value for the QA values

**Note**

Only currently populated bits are available as arguments.

**References**

<https://landsat.usgs.gov/qualityband> for pre collection quality designations (legacy = "pre\_collection")  
<https://landsat.usgs.gov/collectionqualityband> for Collection 1 quality designations (legacy = "collection1")

**Examples**

```
encodeQA(snow = "low", cirrus = c("med", "high"), cloud = "high")
```

---

 estimateHaze
 

---

*Estimate Image Haze for Dark Object Subtraction (DOS)*


---

### Description

estimates the digital number (DN) pixel value of \*dark\* objects for the visible wavelength range.

### Usage

```
estimateHaze(x, hazeBands, darkProp = 0.01, maxSlope = TRUE,
  plot = FALSE, returnTables = FALSE)
```

### Arguments

x	Raster* object or a previous result from estimateHaze with returnTables = TRUE from which to estimate haze
hazeBands	Integer or Character. Band number or bandname from which to estimate atmospheric haze (optional if x contains only one layer)
darkProp	Numeric. Proportion of pixels estimated to be dark.
maxSlope	Logical. Use darkProp only as an upper boundary and search for the DN of maximum slope in the histogram below this value.
plot	Logical. Option to display histograms and haze values
returnTables	Logical. Option to return the frequency table per layer. Only takes effect if x is a Raster* object. If x is a result of estimateHaze tables will always be returned.

### Details

It is assumed that any radiation originating from \*dark\* pixels is due to atmospheric haze and not the reflectance of the surface itself (the surface is dark, i.e. it has a reflectance close to zero). Hence, the haze values are estimates of path radiance, which can be subtracted in a dark object subtraction (DOS) procedure (see [radCor](#))

Atmospheric haze affects almost exclusively the visible wavelength range. Therefore, typically, you'd only want to estimate haze in blue, green and red bands, occasionally also in the nir band.

### Value

If returnTables is FALSE (default). Then a vector of length(hazeBands) containing the estimated haze DNs will be returned. If returnTables is TRUE a list with two components will be returned. The list element 'SHV' contains the haze values, while 'table' contains another list with the sampled frequency tables. The latter can be re-used to try different darkProp thresholds without having to sample the raster again.

**Examples**

```

data(lsat)

## Estimate haze for blue, green and red band
haze <- estimateHaze(lsat, hazeBands = 1:3, plot = TRUE)
haze

## Find threshold interactively
#### Return the frequency tables for re-use
#### avoids having to sample the Raster again and again
haze <- estimateHaze(lsat, hazeBands = 1:3, returnTables = TRUE)
## Use frequency table instead of lsat and fiddle with
haze <- estimateHaze(haze, hazeBands = 1:3, darkProp = .1, plot = TRUE)
haze$SHV

```

fCover

*Fractional Cover Analysis***Description**

fCover takes a classified high resolution image, e.g. vegetation and non-vegetation based on Landsat and calculates cover fractions for pixels of a coarser resolution, e.g. MODIS.

**Usage**

```

fCover(classImage, predImage, nSamples = 1000, classes = 1,
        model = "rf", tuneLength = 3, method = "cv", maxNA = 0,
        clamp = TRUE, filename = NULL, verbose, ...)

```

**Arguments**

classImage	high resolution RasterLayer containing a landcover classification, e.g. as obtained by <a href="#">superClass</a> .
predImage	coarse resolution RasterLayer for which fractional cover will be estimated.
nSamples	Integer. Number of pixels to sample from predImage to train the regression model
classes	Integer. Classes for which fractional cover should be estimated (one or more).
model	Character. Which model to fit for image regression. See <a href="#">train</a> for options. Defaults to randomForest ('rf')
tuneLength	Integer. Number of levels for each tuning parameters that should be generated by train. See Details and <a href="#">train</a> .
method	Character. Resampling method for parameter tuning. Defaults to 10 fold cross-validation. See <a href="#">trainControl</a> for options.
maxNA	Numeric. Maximal proportion of NAs allowed in training pixels.
clamp	Logical. Enforce results to stay within [0,1] interval. Values <0 are reset to 0 and values >1 to 1.

filename	Character. Filename of the output raster file (optional).
verbose	Logical. Print progress information.
...	further arguments to be passed to <a href="#">trainControl</a> and <a href="#">writeRaster</a>

## Details

fCover gets the pixel values in a high resolution classified image that correspond to randomly selected moderate resolution pixels and then calculates the percent of the classified image pixels that represent your cover type of interest. In other words, if your high resolution image has a pixel size of 1m and your moderate resolution image has a pixel size of 30m the sampling process would take a block of 900 of the 1m resolution pixels that correspond to a single 30m pixel and calculate the percentage of the 1m pixels that are forest. For example, if there were 600 forest pixels and 300 non-forest pixels the value given for the output pixel would be 0.67 since  $600/900 = 0.67$ .

fCover relies on the `train()` function from the `caret` package which provides access to a huge number of classifiers. Please see the available options at [train](#). The default classifier (`randomForest`) we chose has been shown to provide very good results in image regression and hence it is recommended you start with this one. If you choose a different classifier, make sure it can run in regression mode.

Many models require tuning of certain parameters. Again, this is handled by [train](#) from the `caret` package. With the `tuneLength` argument you can specify how many different values of each tuning parameter should be tested. The Random Forest algorithm for example can be tuned by varying the `mtry` parameter. Hence, by specifying `tuneLength = 10`, ten different levels of `mtry` will be tested in a cross-validation scheme and the best-performing value will be chosen for the final model.

If you register a parallel backend before, model fitting will run in parallel.

If the total no-data values for a block of high resolution pixels is greater than `maxNA` then it will not be included in the training data set since there is too much missing data to provide a reliable cover percentage. If the no-data proportion is less than `maxNA` the no-data pixels are removed from the total number of pixels when calculating the percent cover.

## Value

Returns a list with two elements: `models` contains the fitted models evaluated in tenfold cross-validation (`caret` train objects); `fCover` contains a `RasterStack` with a fractional cover layer for each requested class.

## See Also

[superClass](#)

## Examples

```
library(raster)
library(caret)
## Create fake input images
data(rlogo)
lsat <- rlogo
agg.level <- 9
modis <- aggregate(lsat, agg.level)
```

```

## Perform classification
lc      <- unsuperClass(lsat, nClass=2)

## Calculate the true cover, which is of course only possible in this example,
## because the fake coarse resolution imagery is exactly res(lsat)*9
trueCover <- aggregate(lc$map, agg.level, fun = function(x, ...){sum(x == 1, ...)/sum(!is.na(x))})

## Run with randomForest and support vector machine (radial basis kernel)
## Of course the SVM is handicapped in this example due to poor tuning (tuneLength)
par(mfrow=c(2,3))
for(model in c("rf", "svmRadial")){
  fc <- fCover(
    classImage = lc$map ,
    predImage = modis,
    classes=1,
    model=model,
    nSample = 50,
    number = 5,
    tuneLength=2
  )

  ## How close is it to the truth?
  compare.rf <- trueCover - fc$map
  plot(fc$map, main = paste("Fractional Cover: Class 1\nModel:", model))
  plot(compare.rf, main = "Diffence\n true vs. predicted")
  plot(trueCover[,fc$map[,], xlim = c(0,1), ylim =c(0,1),
        xlab = "True Cover", ylab = "Predicted Cover" )
  abline(coef=c(0,1))
  rmse <- sqrt(cellStats(compare.rf^2, sum)/ncell(compare.rf))
  r2 <- cor(trueCover[, fc$map[,], "complete.obs")
  text(0.9,0.1,paste0(paste(c("RMSE:", "R2:"),
                             round(c(rmse, r2),3)),collapse="\n"), adj=1)
}

## Reset par
par(mfrow=c(1,1))

```

---

fortify.raster

*Fortify method for classes from the raster package.*


---

## Description

Fortify method for classes from the raster package.

## Usage

```

## S3 method for class 'RasterLayer'
fortify(x, maxpixels = 50000)

```

```
## S3 method for class 'RasterBrick'
fortify(...)

## S3 method for class 'RasterStack'
fortify(...)
```

### Arguments

x	Raster* object to convert into a dataframe.
maxpixels	Integer. Maximum number of pixels to sample
...	not used by this method

### Value

Returns a data.frame with coordinates (x,y) and corresponding raster values.

### Examples

```
library(ggplot2)
data(rlogo)
r_df <- fortify(rlogo)
head(r_df)
```

---

getMeta

*Extract bandwise information from ImageMetaData*

---

### Description

This is an accessor function to quickly access information stored in ImageMetaData, e.g. scale factor per band. Intended for use with imagery which was imported using stackMeta. Will return parameters using the actual band order in img.

### Usage

```
getMeta(img, metaData, what)
```

### Arguments

img	Raster* or character vector with band names.
metaData	ImageMetaData or path to meta data file.
what	Character. Parameter to extract. Either data descriptors, or conversion parameters (see Details for options)



**Details**

Possible metadata parameters (what argument):

Data descriptors

```
'FILES'
'QUANTITY'
'CATEGORY'
'NA_VALUE'
'SATURATE_VALUE'
'SCALE_FACTOR'
'DATA_TYPE'
'SPATIAL_RESOLUTION'
```

### Conversion parameters

```
'CALRAD'  Conversion parameters from DN to radiance
'CALBT'   Conversion parameters from radiance to brightness temperature
'CALREF'  Conversion parameters from DN to reflectance (Landsat 8 only)
```

### Value

If what is one of c('CALRAD', 'CALBT', 'CALREF') a data.frame is returned with bands in rows (order corresponding to img band order). Otherwise a named numeric vector with the corresponding parameter is returned (layernames as names).

### Examples

```
## Import example data
mtlFile <- system.file("external/landsat/LT52240631988227CUB02_MTL.txt", package="RStoolbox")
meta <- readMeta(mtlFile)
lsat <- stackMeta(mtlFile)

## Get integer scale factors
getMeta(lsat, metaData = meta, what = "SCALE_FACTOR")

## Conversion factors for brightness temperature
getMeta("B6_dn", metaData = meta, what = "CALBT")

## Conversion factors to top-of-atmosphere radiance
## Band order not corresponding to metaData order
getMeta(lsat[[5:1]], metaData = meta, what = "CALRAD")

## Get integer scale factors
getMeta(lsat, metaData = meta, what = "SCALE_FACTOR")

## Get file basenames
getMeta(lsat, metaData = meta, what = "FILES")
```

---

getValidation	<i>Extract validation results from superClass objects</i>
---------------	---

---

## Description

Extract validation results from superClass objects

## Usage

```
getValidation(x, from = "testset", metrics = "overall")
```

## Arguments

x	superClass object or caret::confusionMatrix
from	Character. 'testset' extracts the results from independent validation with testset. 'cv' extracts cross-validation results.
metrics	Character. Only relevant in classification mode (ignored for regression models). Select 'overall' for overall accuracy metrics, 'classwise' for classwise metrics, 'confmat' for the confusion matrix itself and 'caret' to return the whole caret::confusionMatrix object.

## Value

Returns a data.frame with validation results. If metrics = 'confmat' or 'caret' will return a table or the full caret::confusionMatrix object, respectively.

## Examples

```
library(pls)
## Fit classifier (splitting training into 70% training data, 30% validation data)
train <- readRDS(system.file("external/trainingPoints.rds", package="RStoolbox"))
SC <- superClass(rlogo, trainData = train, responseCol = "class",
                 model="pls", trainPartition = 0.7)
## Independent testset-validation
getValidation(SC)
getValidation(SC, metrics = "classwise")
## Cross-validation based
getValidation(SC, from = "cv")
```

**Description**

Plot single layer imagery in grey-scale. Can be used with any Raster\* object.

**Usage**

```
ggR(img, layer = 1, maxpixels = 5e+05, alpha = 1, hue = 1,
    sat = 0, stretch = "none", quantiles = c(0.02, 0.98),
    coord_equal = TRUE, ggLayer = FALSE, ggObj = TRUE,
    geom_raster = FALSE, forceCat = FALSE)
```

**Arguments**

img	raster
layer	Character or numeric. Layername or number. Can be more than one layer, in which case each layer is plotted in a subplot.
maxpixels	Integer. Maximal number of pixels to sample.
alpha	Numeric. Transparency (0-1).
hue	Numeric. Hue value for color calculation [0,1] (see \[grDevices]hsv). Change if you need anything else than greyscale. Only effective if sat > 0.
sat	Numeric. Saturation value for color calculation [0,1] (see \[grDevices]hsv). Change if you need anything else than greyscale.
stretch	Character. Either 'none', 'lin', 'hist', 'sqrt' or 'log' for no stretch, linear, histogram, square-root or logarithmic stretch.
quantiles	Numeric vector with two elements. Min and max quantiles to stretch to. Defaults to 2% stretch, i.e. c(0.02,0.98).
coord_equal	Logical. Force addition of coord_equal, i.e. aspect ratio of 1:1. Typically useful for remote sensing data (depending on your projection), hence it defaults to TRUE. Note however, that this does not apply if (ggLayer=FALSE).
ggLayer	Logical. Return only a ggplot layer which must be added to an existing ggplot. If FALSE a stand-alone ggplot will be returned.
ggObj	Logical. Return a stand-alone ggplot object (TRUE) or just the data.frame with values and colors
geom_raster	Logical. If FALSE uses annotation_raster (good to keep aesthetic mappings free). If TRUE uses geom_raster (and aes(fill)). See Details.
forceCat	Logical. If TRUE the raster values will be forced to be categorical (will be converted to factor if needed).

## Details

When `img` contains factor values and `annotation=TRUE`, the raster values will automatically be converted to numeric in order to proceed with the brightness calculation. <sup>ã</sup>

The raster package provides a class lookup-table for categorical rasters (e.g. what you get if you run `superClass` in classification mode). If your raster has a lookup-table `ggR` will automatically treat it as categorical (see [factor](#)). However, the factor status of Raster objects is easily lost and the values are interpreted as numeric. In such cases you should make use of the `forceCat = TRUE` argument, which makes sure that `ggplot2` uses a discrete scale, not a continuous one.

The `geom_raster` argument switches from the default use of `annotation_raster` to `geom_raster`. The difference between the two is that `geom_raster` performs a meaningful mapping from pixel values to fill colour, while `annotation_raster` is simply adding a picture to your plot. In practice this means that whenever you need a legend for your raster you should use `geom_raster = TRUE`. This also allows you to specify and modify the fill scale manually. The advantage of using `annotation_raster` (`geom_raster = TRUE`) is that you can still use the `scale_fill*` for another variable. For example you could add polygons and map a value to their fill colour. For more details on the theory behind aesthetic mapping have a look at the [ggplot2](#) manuals.

## Value

`ggObj = TRUE:` `ggplot2` plot  
`ggLayer = TRUE:` `ggplot2` layer to be combined with an existing `ggplot2`  
`ggObj = FALSE:` data.frame in long format suitable for plotting with `ggplot2`, includes the pixel values and the calculated

## See Also

[ggRGB](#), [fortify](#)

## Examples

```
library(ggplot2)
library(raster)
data(rlogo)

## Simple grey scale annotation
ggR(rlogo)

## With linear stretch contrast enhancement
ggR(rlogo, stretch = "lin", quantiles = c(0.1,0.9))

## Don't plot, just return a data.frame
df <- ggR(rlogo, ggObj = FALSE)
head(df)

## ggplot with geom_raster instead of annotation_raster
## and default scale_fill*
```

```

ggR(rlogo, geom_raster = TRUE)

## with different scale
ggR(rlogo, geom_raster = TRUE) +
  scale_fill_gradientn(name = "mojo", colours = rainbow(10)) +
  ggtitle("**Funkadelic**")

## Layermode (ggLayer=TRUE)
data <- data.frame(x = c(0, 0:100,100), y = c(0,sin(seq(0,2*pi,pi/50))*10+20, 0))
ggplot(data, aes(x, y)) + ggR(rlogo, geom_raster= FALSE, ggLayer = TRUE) +
  geom_polygon(aes(x, y), fill = "blue", alpha = 0.4) +
  coord_equal(ylim=c(0,75))

## Categorical data
## In this case you probably want to use geom_raster=TRUE
## in order to perform aesthetic mapping (i.e. a meaningful legend)
rc <- raster(rlogo)
rc[] <- cut(rlogo[[1]][[]], seq(0,300, 50))
ggR(rc, geom_raster = TRUE)

## Legend customization etc. ...
ggR(rc, geom_raster = TRUE) + scale_fill_discrete(labels=paste("Class", 1:6))

## Creating a nicely looking DEM with hillshade background
data(srtm)
terr <- terrain(srtm, c("slope", "aspect"))
hill <- hillShade(terr[["slope"]], terr[["aspect"]])
ggR(hill)

ggR(hill) +
  ggR(srtm, geom_raster = TRUE, ggLayer = TRUE, alpha = 0.3) +
  scale_fill_gradientn(colours = terrain.colors(100), name = "elevation")

```

---

ggRGB

---

*Create ggplot2 Raster Plots with RGB from 3 RasterLayers*


---

## Description

Calculates RGB color composite raster for plotting with ggplot2. Optional values for clipping and stretching can be used to enhance the imagery.

## Usage

```

ggRGB(img, r = 3, g = 2, b = 1, scale, maxpixels = 5e+05,
  stretch = "none", ext = NULL, limits = NULL,
  clipValues = "limits", quantiles = c(0.02, 0.98), ggObj = TRUE,
  ggLayer = FALSE, alpha = 1, coord_equal = TRUE,
  geom_raster = FALSE, nullValue = 0)

```

**Arguments**

<code>img</code>	RasterStack or RasterBrick
<code>r</code>	Integer or character. Red layer in <code>x</code> . Can be set to <code>NULL</code> , in which case the red channel will be set to zero.
<code>g</code>	Integer or character. Green layer in <code>x</code> . Can be set to <code>NULL</code> , in which case the green channel will be set to zero.
<code>b</code>	Integer or character. Blue layer in <code>x</code> . Can be set to <code>NULL</code> , in which case the blue channel will be set to zero.
<code>scale</code>	Numeric. Maximum possible pixel value (optional). Defaults to 255 or to the maximum value of <code>x</code> if that is larger than 255
<code>maxpixels</code>	Integer. Maximal number of pixels used for plotting.
<code>stretch</code>	Character. Either <code>'none'</code> , <code>'lin'</code> , <code>'hist'</code> , <code>'sqrt'</code> or <code>'log'</code> for no stretch, linear, histogram, square-root or logarithmic stretch.
<code>ext</code>	Extent object to crop the image
<code>limits</code>	Vector or matrix. Can be used to reduce the range of values. Either a vector of two values for all bands ( <code>c(min, max)</code> ) or a 3x2 matrix with min and max values (columns) for each layer (rows).
<code>clipValues</code>	Matrix, numeric vector, string or NA. Values to reset out of range (out of <code>limits</code> ) values to. By default ( <code>'limits'</code> ) values are reset to <code>limits</code> . A single value (e.g. NA) will be recycled to all lower/higher clippings, A vector of length two ( <code>c(min,max)</code> ) can be used to specify lower and higher replace values, applied to all bands. A two column matrix (typically with three rows) can be used to fully control lower and upper clipping values differently for each band.
<code>quantiles</code>	Numeric vector with two elements. Min and max quantiles to stretch. Defaults to 2% stretch, i.e. <code>c(0.02,0.98)</code> .
<code>ggObj</code>	Logical. If <code>TRUE</code> a ggplot2 object is returned. If <code>FALSE</code> a data.frame with coordinates and color will be returned.
<code>ggLayer</code>	Logical. If <code>TRUE</code> a ggplot2 layer is returned. This is useful if you want to add it to an existing ggplot2 object. Note that if <code>TRUE &amp; annotate = FALSE</code> you have to add a <code>scale_fill_identity()</code> manually in your call to <code>ggplot()</code> .
<code>alpha</code>	Numeric. Transparency (0-1).
<code>coord_equal</code>	Logical. Force addition of <code>coord_equal</code> , i.e. aspect ratio of 1:1. Typically useful for remote sensing data (depending on your projection), hence it defaults to <code>TRUE</code> . Note however, that this does not apply if ( <code>ggLayer=FALSE</code> ).
<code>geom_raster</code>	Logical. If <code>FALSE</code> <code>annotation_raster</code> is used, otherwise <code>geom_raster()+scale_fill_identity</code> is used. Note that you can't use <code>scale_fill*</code> in addition to the latter, because it already requires <code>scale_fill_identity()</code> .
<code>nullValue</code>	Numeric. Intensity value used for <code>NULL</code> layers in color compositing. E.g. <code>set g=NULL</code> and fix green value at 0.5 (defaults to 0).

**Details**

Functionality is based on [plotRGB](#) from the raster package.

**Value**



ggObj = TRUE: ggplot2 plot  
ggLayer = TRUE: ggplot2 layer to be combined with an existing ggplot2  
ggObj = FALSE: data.frame in long format suitable for plotting with ggplot2, includes the pixel values and the calculated c

## See Also

[ggR](#), [fortify](#)

## Examples

```
library(ggplot2)
data(rlogo)

ggRGB(rlogo, r=1, g=2, b=3)

## Define minMax ranges
ggRGB(rlogo, r=1,g=2, b=3, limits = matrix(c(100,150,10,200,50,255), ncol = 2, by = TRUE))

## Perform strong linear contrast stretch
ggRGB(rlogo, r = 1, g = 2, b = 3,stretch = "lin", quantiles = c(0.2, 0.8))

## Use only two layers for color calculation
ggRGB(rlogo, r = 1, g = 2, b = NULL)

## Return only data.frame
df <- ggRGB(rlogo, ggObj = FALSE)
head(df)

## Use in layer-mode, e.g. to add to another plot
wave <- data.frame(x = c(0, 0:100,100), y = c(0,sin(seq(0,2*pi,pi/50))*10+20, 0))
p <- ggplot(wave, aes(x, y))
p + ggRGB(rlogo, ggLayer = TRUE) +
  geom_polygon(aes(x, y), fill = "blue", alpha = 0.4) +
  coord_equal(ylim=c(0,75))
```

---

histMatch

*Image to Image Contrast Matching*

---

## Description

Performs image to image contrast adjustments based on histogram matching using empirical cumulative distribution functions from both images.

**Usage**

```
histMatch(x, ref, xmask = NULL, refmask = NULL, nSamples = 1e+05,
  intersectOnly = TRUE, paired = TRUE, forceInteger = FALSE,
  returnFunctions = FALSE, ...)
```

**Arguments**

x	Raster*. Source raster which is to be modified.
ref	Raster*. Reference raster, to which x will be matched.
xmask	RasterLayer. Mask layer for x to exclude pixels which might distort the histogram, i.e. are not present in ref. Any NA pixel in xmask will be ignored (maskvalue = NA).
refmask	RasterLayer. Mask layer for ref. Any NA pixel in refmask will be ignored (maskvalue = NA).
nSamples	Integer. Number of random samples from each image to build the histograms.
intersectOnly	Logical. If TRUE sampling will only take place in the overlap extent of the two rasters. Otherwise the full rasters will be used for sampling.
paired	Logical. If TRUE the corresponding pixels will be used in the overlap.
forceInteger	Logical. Force integer output.
returnFunctions	Logical. If TRUE the matching functions will be returned instead of applying them to x.
...	Further arguments to be passed to <a href="#">writeRaster</a> .

**Value**

A Raster\* object of x adjusted to the histogram of ref. If returnFunctions = TRUE a list of functions (one for each layer) will be returned instead.

**Note**

x and ref must have the same number of layers.

**References**

Richards and Jia: Remote Sensing Digital Image Analysis. Springer, Berlin, Heidelberg, Germany, 439pp.

**Examples**

```
library(ggplot2)
library(raster)
data(rlogo)
## Original image a (+1 to prevent log(0))
img_a <- rlogo + 1
## Degraded image b
img_b <- log(img_a)
```

```

## Cut-off half the image (just for better display)
img_b[, 1:50] <- NA

## Compare Images before histMatching
ggRGB(img_a,1,2,3)+
  ggRGB(img_b, 1,2,3, ggLayer = TRUE, stretch = "lin", q = 0:1) +
  geom_vline(aes(xintercept = 50))+
  ggtitle("Img_a vs. Img_b")

## Do histogram matching
img_b_matched <- histMatch(img_b, img_a)

## Compare Images after histMatching
ggRGB(img_a, 1, 2, 3)+
  ggRGB(img_b_matched, 1, 2, 3, ggLayer = TRUE, stretch = "lin", q = 0:1) +
  geom_vline(aes(xintercept = 50))+
  ggtitle("Img_a vs. Img_b_matched")

## Histogram comparison
opar <- par(mfrow = c(1, 3), no.readonly = TRUE)
img_a[,1:50] <- NA
redLayers <- stack(img_a, img_b, img_b_matched)[[c(1,4,7)]]
names(redLayers) <- c("img_a", "img_b", "img_b_matched")

hist(redLayers)
## Reset par
par(opar)

```

---

ImageMetaData

*ImageMetaData Class*


---

## Description

ImageMetaData Class

## Usage

```

ImageMetaData(file = NA, format = NA, sat = NA, sen = NA,
  scene = NA, proj = NA, date = NA, pdate = NA, path = NA,
  row = NA, az = NA, selv = NA, esd = NA, files = NA,
  bands = NA, quant = NA, cat = NA, na = NA, vsat = NA,
  scal = NA, dtyp = NA, calrad = NA, calref = NA, calbt = NA,
  radRes = NA, spatRes = NA)

```

## Arguments

file	Character. Metadata file
format	Character. Metadata format, e.g. xml, mtl
sat	Character. Satellite platform

sen	Character. Sensor
scene	Character. Scene_ID
proj	CRS. Projection.
date	POSIXct. Aquisition date.
pdate	POSIXct. Processing date.
path	Integer. Path.
row	Integer. Row.
az	Numeric. Sun azimuth
selv	Numeric. Sun elevation
esd	Numeric. Earth-sun distance
files	Character vector. Files containing the data, e.g. tiff files
bands	Character vector. Band names
quant	Character vector. Quantity, one of c("dn", "tra", "tre", "sre", "bt", "idx")
cat	Character vector. Category, e.g. c("image", "pan", "index", "qa")
na	Numeric vector. No-data value per band
vsat	Numeric vector. Saturation value per band
scal	Numeric vector. Scale factor per band. e.g. if data was scaled to 1000*reflectance for integer conversion.
dtype	Character vector. Data type per band. See <a href="#">dataType</a> for options.
calrad	data.frame. Calibration coefficients for dn->radiance conversion. Must have columns 'gain' and 'offset'. Rows named according to bands.
calref	data.frame. Calibration coefficients for dn->reflectance conversion. Must have columns 'gain' and 'offset'. Rows named according to bands.
calbt	data.frame. Calibration coefficients for dn->brightness temperature conversion. Must have columns 'K1' and 'K2'. Rows named according to bands.
radRes	Numeric vector. Radiometric resolution per band.
spatRes	Numeric vector. Spatial resolution per band.

---

 lsat

*Landsat 5TM Example Data*


---

### Description

Subset of Landsat 5 TM Scene: LT52240631988227CUB02 Contains all seven bands in DN format.

### Usage

```
data(lsat)
```

### Examples

```
data(lsat)
ggRGB(lsat, stretch = "lin")
```

---

mesma *Multiple Endmember Spectral Mixture Analysis (Spectral Unmixing)*

---

### Description

mesma performs a multiple endmember spectral mixture analysis on a multiband raster image.

### Usage

```
mesma(img, em, method = "NNLS", iterate = 400, tolerance = 1e-08,
      ..., verbose)
```

### Arguments

img	RasterBrick or RasterStack. Remote sensing imagery (usually hyperspectral).
em	Matrix or data.frame with spectral endmembers. Rows represent a single end-member of a class, columns represent the spectral bands (i.e. columns correspond to number of bands in img). Number of rows needs to be > 1.
method	Character. Select an unmixing method. Currently, only "NNLS" is implemented. Default is "NNLS". <ul style="list-style-type: none"> <li>• NNLS: applies a non-negative least squares (NNLS) regression which is using a sequential coordinate-wise algorithm (SCA) based on Franc et al. (2005).</li> </ul>
iterate	Integer. Set maximum iteration per pixel. Processing time could increase the more iterations are made possible. Default is 400.
tolerance	Numeric. Tolerance limit representing a nearly zero minimal number. Default is 1e-8.
...	further arguments passed to <a href="#">writeRaster</a> .
verbose	Logical. Prints progress messages during execution.

### Value

RasterBrick. The object will contain one band per endmember, with each value representing the estimated presence probability of the endmember per pixel (0 to 1), and an RMSE band.

### Note

Depending on iterate and tolerance settings, the sum of estimated presence probabilities per pixel varies around 1.

### Author(s)

Jakob Schwalb-Willmann

## References

Franc, V., Hlaváč, V., & Navara, M. (2005). Sequential coordinate-wise algorithm for the non-negative least squares problem. In: International Conference on Computer Analysis of Images and Patterns (pp. 407-414). Berlin, Heidelberg.

## Examples

```
#load packages
library(raster)
library(RStoolbox)

#load an example dataset
data(lsat)

#make up some endmember spectra: water and land
em_names <- c("water", "land")
pts <- data.frame(class=em_names, cell = c(47916,5294))
em <- lsat[pts$cell]
rownames(em) <- em_names

#unmix the image for water and land
probs <- mesma(lsat, em, method = "NNLS")

#take a look
raster::hist(probs$water)
raster::plot(probs$water, col = c("white","blue"))
raster::hist(probs$land)
raster::plot(probs$land, col = c("white","brown"))
```

---

normImage

*Normalize Raster Images: Center and Scale*

---

## Description

For each pixel subtracts the mean of the raster layer and optionally divide by its standard deviation.

## Usage

```
normImage(img, norm = TRUE, ...)
```

## Arguments

img	Raster* object. Image to transform. Transformation will be performed separately for each layer.
norm	Logical. Perform normalization (scaling) in addition to centering, i.e. divide by standard deviation.
...	further arguments passed to <a href="#">writeRaster</a> .

**Value**

Returns a Raster\* with the same number layers as input layers with each layer being centered and optionally normalized.

**Examples**

```
library(raster)
## Load example data
data(rlogo)

## Normalization: Center and Scale
rlogo_center_norm <- normImage(rlogo)
hist(rlogo_center_norm)

## Centering
rlogo_center <- normImage(rlogo, norm = FALSE)
```

---

oneHotEncode	<i>One-hot encode a raster or vector</i>
--------------	--

---

**Description**

Splits a categorical raster layer (or a vector) into a multilayer raster (or matrix).

**Usage**

```
oneHotEncode(img, classes, background = 0, foreground = 1,
             na.rm = FALSE, ...)
```

**Arguments**

img	RasterLayer or integer/numeric vector containing multiple classes
classes	integer: vector of classes which should be extracted
background	integer: background value (default = 0)
foreground	integer: foreground value (default = 1)
na.rm	logical: if TRUE, NAs will be coerced to the background value.
...	further arguments passed to <a href="#">writeRaster</a> . Ignored if img is not a RasterLayer, but a numeric/integer vector

**Value**

A RasterBrick with as many layers as there are classes. Pixels matching the class of interest are set to 1, background values by default are set to 0 (see background argument)

**Examples**

```

library(raster)

## example data
data(rlogo)
sc <- unsuperClass(rlogo, nClasses = 3)

## one-hot encode
sc_oneHot <- oneHotEncode(sc$map, classes = c(1,2,3))

## check results
sc_oneHot
plot(sc_oneHot)

```

---

panSharpen

*Pan Sharpen Imagery / Image Fusion*


---

**Description**

provides different methods for pan sharpening a coarse resolution (typically multispectral) image with a higher resolution panchromatic image. Values of the pan-chromatic and multispectral images must be of the same scale, (e.g. from 0:1, or all DNs from 0:255)

**Usage**

```
panSharpen(img, pan, r, g, b, pc = 1, method = "brovey", norm = TRUE)
```

**Arguments**

img	Raster* object. Coarse resolution multispectral image
pan	RasterLayer. High resolution image, typically panchromatic.
r	Character or Integer. Red band in img. Only relevant if method!='pca'
g	Character or Integer. Green band in img. Only relevant if method!='pca'
b	Character or Integer. Blue band in img. Only relevant if method!='pca'
pc	Integer. Only relevant if method = 'pca'. Which principal component to replace. Usually this should be the first component (default). Only if the first component is dominated by something else than brightness it might be worth a try to use the second component.
method	Character. Choose method from c("pca", "ihs", "brovey").
norm	Logical. Rescale pan image to match the 1st PC component. Only relevant if method = 'pca'. If TRUE only min and max are matched to the 1st PC. If FALSE pan will be histogram matched to the 1st PC.



## Details

Pan sharpening options:

- `method='pca'`: Performs a pca using [rasterPCA](#). The first component is then swapped for the pan band an the PCA is rotated backwards.
- `method='ihs'`: Performs a color space transform to Intensity-Hue-Saturation space, swaps intensity for the histogram matched pan and does the backwards transformation.
- `method='brovey'`: Performs Brovey reweighting. Pan and img must be at the same value scale (e.g. 0:1, or 0:255) otherwise you'll end up with psychodelic colors.

## Examples

```
library(raster)
library(ggplot2)

## Load example data
data(lsat)
## Fake panchromatic image (30m resolution covering
## the visible range (integral from blue to red))
pan      <- sum(lsat[[1:3]])
ggR(pan, stretch = "lin")

## Fake coarse resolution image (150m spatial resolution)
lowResImg <- aggregate(lsat, 5)

## Brovey pan sharpening
lowResImg_pan <- panSharpen(lowResImg, pan, r = 3, g = 2, b = 1, method = "brovey")
lowResImg_pan
## Plot
ggRGB(lowResImg, stretch = "lin") + ggtitle("Original")
ggRGB(lowResImg_pan, stretch="lin") + ggtitle("Pansharpened (Brovey)")
```

---

pifMatch

*Pseudo-Invariant Features based Image Matching*

---

## Description

Match one scene to another based on linear regression of pseudo-invariant features (PIF).

## Usage

```
pifMatch(img, ref, method = "cor", quantile = 0.95,
  returnPifMap = TRUE, returnSimMap = TRUE, returnModels = FALSE)
```

**Arguments**

<code>img</code>	RasterStack or RasterBrick. Image to be adjusted.
<code>ref</code>	RasterStack or RasterBrick. Reference image.
<code>method</code>	Method to calculate pixel similarity. Options: euclidean distance ('ed'), spectral angle ('sam') or pearson correlation coefficient ('cor').
<code>quantile</code>	Numeric. Threshold quantile used to identify PIFs
<code>returnPifMap</code>	Logical. Return a binary raster map of pixels which were identified as pseudo-invariant features.
<code>returnSimMap</code>	Logical. Return the similarity map as well
<code>returnModels</code>	Logical. Return the linear models along with the adjusted image.

**Details**

The function consists of three main steps: First, it calculates pixel-wise similarity between the two rasters and identifies pseudo-invariant pixels based on a similarity threshold. In the second step the values of the pseudo-invariant pixels are regressed against each other in a linear model for each layer. Finally the linear models are applied to all pixels in the `img`, thereby matching it to the reference scene.

Pixel-wise similarity can be calculated using one of three methods: euclidean distance (`method = "ed"`), spectral angle ("`sam`") or pearsons correlation coefficient ("`cor`"). The threshold is defined as a similarity quantile. Setting `quantile=0.95` will select all pixels with a similarity above the 95% quantile as pseudo-invariant features.

Model fitting is performed with simple linear models ([lm](#)); fitting one model per layer.

**Value**

Returns a List with the adjusted image and intermediate products (if requested). #'

- `img`: the adjusted image
- `simMap`: pixel-wise similarity map (if `returnSimMap = TRUE`)
- `pifMap`: binary map of pixels selected as pseudo-invariant features (if `returnPifMap = TRUE`)
- `models`: list of linear models; one per layer (if `returnModels = TRUE`)

**Examples**

```
library(raster)

## Import Landsat example data
data(lsat)

## Create fake example data
## In practice this would be an image from another acquisition date
lsat_b <- log(lsat)

## Run pifMatch and return similarity layer, invariant features mask and models
lsat_b_adj <- pifMatch(lsat_b, lsat, returnPifMap = TRUE,
```

```

returnSimMap = TRUE, returnModels = TRUE)

## Pixelwise similarity
ggR(lsat_b_adj$simMap, geom_raster = TRUE)

## Pseudo invariant feature mask
ggR(lsat_b_adj$pifMap)

## Histograms of changes
par(mfrow=c(1,3))
hist(lsat_b[[1]], main = "lsat_b")
hist(lsat[[1]], main = "reference")
hist(lsat_b_adj$img[[1]], main = "lsat_b adjusted")

## Model summary for first band
summary(lsat_b_adj$models[[1]])

```

---

predict.superClass      *Predict a raster map based on a superClass model fit.*

---

## Description

useful to separate model fitting from spatial prediction, which can take some time.

## Usage

```

## S3 method for class 'superClass'
predict(object, img, predType = "raw",
        filename = NULL, datatype = "INT2U", ...)

```

## Arguments

object	superClass object
img	Raster object. Layer names must correspond to layer names used to train the superClass model, i.e. layer names in the original raster image.
predType	Character. Type of the final output raster. Either "raw" for class predictions or "prob" for class probabilities. Class probabilities are not available for all classification models ( <a href="#">predict.train</a> ).
filename	Character or NULL. Filename for output raster file.
datatype	Datatype of output raster file.
...	Further arguments passed to writeRaster.

**Examples**

```
## Load training data
data(rlogo)
train <- readRDS(system.file("external/trainingPoints.rds", package="RStoolbox"))

## Fit classifier
SC <- superClass(rlogo, trainData = train, responseCol = "class",
                 model = "rf", tuneLength = 1, predict = FALSE)

map <- predict(SC, rlogo)
```

---

predict.unsuperClass *Predict a raster map based on a unsuperClass model fit.*

---

**Description**

applies a kmeans cluster model to all pixels of a raster. Useful if you want to apply a kmeans model of scene A to scene B.

**Usage**

```
## S3 method for class 'unsuperClass'
predict(object, img, ...)
```

**Arguments**

object	unsuperClass object
img	Raster object. Layernames must correspond to layernames used to train the superClass model, i.e. layernames in the original raster image.
...	Further arguments passed to writeRaster.

**Examples**

```
## Load training data
data(rlogo)

## Perform unsupervised classification
uc <- unsuperClass(rlogo, nClasses = 10)

## Apply the model to another raster
map <- predict(uc, rlogo)
```

**Description**

Implements several different methods for radiometric calibration and correction of Landsat data. You can either specify a metadata file, or supply all necessary values manually. With proper parametrization `apref` and `sdos` should work for other sensors as well.

**Usage**

```
radCor(img, metaData, method = "apref", bandSet = "full", hazeValues,
       hazeBands, atmosphere, darkProp = 0.01, clamp = TRUE, verbose)
```

**Arguments**

<code>img</code>	raster object
<code>metaData</code>	object of class <code>ImageMetaData</code> or a path to the meta data (MTL) file.
<code>method</code>	Radiometric conversion/correction method to be used. There are currently four methods available (see Details): "rad", "apref", "sdos", "dos", "costz".
<code>bandSet</code>	Numeric or character. original Landsat band numbers or names in the form of ("B1", "B2" etc). If set to 'full' all bands in the solar (optical) region will be processed.
<code>hazeValues</code>	Numeric. Either a vector with dark DNs per hazeBand (method = 'sdos'); possibly estimated using <code>estimateHaze</code> . Or the 'starting haze value' (DN) for the relative scattering models in method = 'dos' or 'costz'. If not provided, hazeValues will be estimated in an automated fashion for all hazeBands. Argument only applies to methods 'sdos', 'dos' and 'costz'.
<code>hazeBands</code>	Character or integer. Bands corresponding to hazeValues (method = 'sdos') or band to select starting haze value from ('dos' or 'costz').
<code>atmosphere</code>	Character. Atmospheric characteristics. Will be estimated if not explicitly provided. Must be one of "veryClear", "clear", "moderate", "hazy" or "veryHazy".
<code>darkProp</code>	Numeric. Estimated proportion of dark pixels in the scene. Used only for automatic guessing of hazeValues (typically one would choose 1 or 2%).
<code>clamp</code>	Logical. Enforce valid value range. By default reflectance will be forced to stay within [0,1] and radiance $\geq 0$ by replacing invalid values with the corresponding boundary, e.g. -0.1 will become 0.
<code>verbose</code>	Logical. Print status information.

## Details

The atmospheric correction methods (sdos, dos and costz) apply to the optical (solar) region of the spectrum and do not affect the thermal band.

Dark object subtraction approaches rely on the estimation of atmospheric haze based on \*dark\* pixels. Dark pixels are assumed to have zero reflectance, hence the name. It is then assumed further that any radiation originating from such \*dark\* pixels is due to atmospheric haze and not the reflectance of the surface itself.

The following methods are available:

rad	Radiance
apref	Apparent reflectance (top-of-atmosphere reflectance)
dos	Dark object subtraction following Chavez (1989)
costz	Dark object subtraction following Chavez (1996)
sdos	Simple dark object subtraction. Classical DOS, Lhaze must be estimated for each band separately.

If either "dos" or "costz" are selected, radCor will use the atmospheric haze decay model described by Chavez (1989). Depending on the atmosphere the following coefficients are used:

veryClear	$\lambda^{-4.0}$
clear	$\lambda^{-2.0}$
moderate	$\lambda^{-1.0}$
hazy	$\lambda^{-0.7}$
veryHazy	$\lambda^{-0.5}$

For Landsat 8, no values for extra-terrestrial irradiation (esun) are provided by NASA. These are, however, necessary for DOS-based approaches. Therefore, these values were derived from a standard reference spectrum published by Thuillier et al. (2003) using the Landsat 8 OLI spectral response functions (for details, see <http://bleutner.github.io/RStoolbox/r/2016/01/26/estimating-landsat-8-esun-values>).

The implemented sun-earth distances neglect the earth's eccentricity. Instead we use a 100 year daily average (1979-2070).

## Value

RasterStack with top-of-atmosphere radiance ( $W/(m^2 * sr * \mu m)$ ), at-satellite brightness temperature (K), top-of-atmosphere reflectance (unitless) corrected for the sun angle or at-surface reflectance (unitless).

## Note

This was originally a fork of randcorr() function in the landsat package. This version works on Raster\* objects and hence is suitable for large rasters.

## References

S. Goslee (2011): Analyzing Remote Sensing Data in R: The landsat Package. Journal of Statistical Software 43(4).

G. Thuillier et al. (2003) THE SOLAR SPECTRAL IRRADIANCE FROM 200 TO 2400 nm AS MEASURED BY THE SOLSPEC SPECTROMETER FROM THE ATLAS AND EURECA MISSIONS. Solar Physics 214(1): 1-22 (

### Examples

```
library(raster)
## Import meta-data and bands based on MTL file
mtlFile <- system.file("external/landsat/LT52240631988227CUB02_MTL.txt",
  package="RStoolbox")
metaData <- readMeta(mtlFile)
lsat <- stackMeta(mtlFile)

## Convert DN to top of atmosphere reflectance and brightness temperature
lsat_ref <- radCor(lsat, metaData = metaData, method = "apref")

## Correct DN to at-surface-reflecatance with DOS (Chavez decay model)
lsat_sref <- radCor(lsat, metaData = metaData, method = "dos")

## Correct DN to at-surface-reflecatance with simple DOS
## Automatic haze estimation
hazeDN <- estimateHaze(lsat, hazeBands = 1:4, darkProp = 0.01, plot = TRUE)
lsat_sref <- radCor(lsat, metaData = metaData, method = "sdos",
  hazeValues = hazeDN, hazeBands = 1:4)
```

---

 rasterCVA

*Change Vector Analysis*


---

### Description

Calculates angle and magnitude of change vectors. Dimensionality is limited to two bands per image.

### Usage

```
rasterCVA(x, y, tmf = 2, ...)
```

### Arguments

x	RasterBrick or RasterStack with two layers. This will be the reference/origin for the change calculations. Both rasters (y and y) need to correspond to each other, i.e. same resolution, extent and origin.
y	RasterBrick or RasterStack with two layers. Both rasters (y and y) need to correspond to each other, i.e. same resolution, extent and origin.
tmf	Numeric. Threshold median factor. Used to calculate a threshold magnitude for which pixels are considered stable, i.e. no change. Defaults to 2 times the median non-zero magnitude. Calculated as $tmf * median(magnitude[magnitude > 0])$
...	further arguments passed to writeRaster

**Details**

Change Vector Analysis (CVA) is used to identify spectral changes between two identical scenes which were acquired at different times. CVA is limited to two bands per image. For each pixel it calculates the change vector in the two-dimensional spectral space. For example for a given pixel in image A and B for the red and nir band the change vector is calculated for the coordinate pairs: (red\_A | nir\_A) and (red\_B | nir\_B).

The coordinate system is defined by the order of the input bands: the first band defines the x-axis and the second band the y-axis, respectively. Angles are returned *\*in degree\** beginning with 0 degrees pointing 'north', i.e. the y-axis, i.e. the second band.

**Value**

Returns a RasterBrick with two layers: change vector angle and change vector magnitude

**Examples**

```
library(raster)
## Create example data
data(lsat)
pca <- rasterPCA(lsat)$map

## Do change vector analysis
cva <- rasterCVA(pca[[1:2]], pca[[3:4]])
cva
plot(cva)
```

---

rasterEntropy

*Multi-layer Pixel Entropy*


---

**Description**

Shannon entropy is calculated for each pixel based on it's layer values. To be used with categorical / integer valued rasters.

**Usage**

```
rasterEntropy(img, ...)
```

**Arguments**

```
img          RasterStack or RasterBrick
...          additional arguments passed to writeRaster
```

**Details**

Entropy is calculated as  $-\sum(p \log(p))$ ; p being the class frequency per pixel.



**Value**

RasterLayer "entropy"

**Examples**

```
data(rlogo)
re <- rasterEntropy(rlogo)
ggR(re, geom_raster = TRUE)
```

---

rasterPCA

*Principal Component Analysis for Rasters*


---

**Description**

Calculates R-mode PCA for RasterBricks or RasterStacks and returns a RasterBrick with multiple layers of PCA scores.

**Usage**

```
rasterPCA(img, nSamples = NULL, nComp = nlayers(img), spca = FALSE,
           maskCheck = TRUE, ...)
```

**Arguments**

<code>img</code>	RasterBrick or RasterStack.
<code>nSamples</code>	Integer or NULL. Number of pixels to sample for PCA fitting. If NULL, all pixels will be used.
<code>nComp</code>	Integer. Number of PCA components to return.
<code>spca</code>	Logical. If TRUE, perform standardized PCA. Corresponds to centered and scaled input image. This is usually beneficial for equal weighting of all layers. (FALSE by default)
<code>maskCheck</code>	Logical. Masks all pixels which have at least one NA (default TRUE is recommended but introduces a slow-down, see Details when it is wise to disable maskCheck). Takes effect only if nSamples is NULL.
<code>...</code>	further arguments to be passed to <a href="#">writeRaster</a> , e.g. filename.

**Details**

Internally rasterPCA relies on the use of [princomp](#) (R-mode PCA). If nSamples is given the PCA will be calculated based on a random sample of pixels and then predicted for the full raster. If nSamples is NULL then the covariance matrix will be calculated first and will then be used to calculate princomp and predict the full raster. The latter is more precise, since it considers all pixels, however, it may be slower than calculating the PCA only on a subset of pixels.

Pixels with missing values in one or more bands will be set to NA. The built-in check for such pixels can lead to a slow-down of rasterPCA. However, if you make sure or know beforehand that

all pixels have either only valid values or only NAs throughout all layers you can disable this check by setting `maskCheck=FALSE` which speeds up the computation.

Standardised PCA (SPCA) can be useful if imagery or bands of different dynamic ranges are combined. SPC uses the correlation matrix instead of the covariance matrix, which has the same effect as using normalised bands of unit variance.

## Value

RasterBrick

## Examples

```
library(ggplot2)
library(reshape2)
data(rlogo)
ggRGB(rlogo, 1,2,3)

## Run PCA
set.seed(25)
rpc <- rasterPCA(rlogo)
rpc
summary(rpc$model)

ggRGB(rpc$map,1,2,3, stretch="lin", q=0)
if(require(gridExtra)){
plots <- lapply(1:3, function(x) ggR(rpc$map, x, geom_raster = TRUE))
grid.arrange(plots[[1]],plots[[2]], plots[[3]], ncol=2)
}
```

---

readEE

*Tidy import tool for EarthExplorer .csv export files*

---

## Description

Imports and tidies CSV files exported from EarthExplorer into data.frames and annotates missing fields.

## Usage

```
readEE(x)
```

## Arguments

x Character, Character or list. One or more paths to EarthExplorer export files.

**Details**

The **EarthExplorer** CSV file can be produced from the search results page. Above the results click on 'export results' and select 'comma (,) delimited'.

Note that only a subset of columns is imported which was deemed interesting. Please contact the maintainer if you think an omitted column should be included.

**Value**

data.frame

**Examples**

```
library(ggplot2)
ee <- readEE(system.file("external/EarthExplorer_LS8.txt", package = "RStoolbox"))

## Scenes with cloud cover < 20%
ee[ee$Cloud.Cover < 20,]

## Available time-series
ggplot(ee) +
  geom_segment(aes(x = Date, xend = Date, y = 0, yend = 100 - Cloud.Cover,
                  col = as.factor(Year))) +
  scale_y_continuous(name = "Scene quality (% clear sky)")
```

---

readMeta

*Read Landsat MTL metadata files*

---

**Description**

Reads metadata and deals with legacy versions of Landsat metadata files and where possible adds missing information (radiometric gain and offset, earth-sun distance).

**Usage**

```
readMeta(file, raw = FALSE)
```

**Arguments**

file	path to Landsat MTL file (...MTL.txt)
raw	Logical. If TRUE the full raw metadata will be returned as a list. if FALSE (the default) all important metadata are homogenized into a standard format (ImageMetaData) and some information is added.

**Value**

Object of class ImageMetaData

### Examples

```
## Example metadata file (MTL)
mtlFile <- system.file("external/landsat/LT52240631988227CUB02_MTL.txt", package="RStoolbox")

## Read metadata
metaData <- readMeta(mtlFile)

## Summary
summary(metaData)
```

---

readSLI	<i>Read ENVI spectral libraries</i>
---------	-------------------------------------

---

### Description

read/write support for ENVI spectral libraries

### Usage

```
readSLI(path)
```

### Arguments

path                    Path to spectral library file with ending .sli.

### Details

ENVI spectral libraries consist of a binary data file (.sli) and a corresponding header (.hdr, or .sli.hdr) file.

### Value

The spectral libraries are read into a data.frame. The first column contains the wavelengths and the remaining columns contain the spectra.

### See Also

[writeSLI](#)

### Examples

```
## Example data
sliFile <- system.file("external/vegSpec.sli", package="RStoolbox")
sliTmpFile <- paste0(tempdir(), "/vegetationSpectra.sli")

## Read spectral library
sli <- readSLI(sliFile)
```

```

head(sli)
plot(sli[,1:2], col = "orange", type = "l")
lines(sli[,c(1,3)], col = "green")

## Write to binary spectral library
writeSLI(sli, path = sliTmpFile)

```

rescaleImage

*Linear Image Rescaling***Description**

performs linear shifts of value ranges either to match min and max of another image (y) or to any other min and max value (ymin and ymax).

**Usage**

```
rescaleImage(x, y, xmin, xmax, ymin, ymax, forceMinMax = FALSE)
```

**Arguments**

x	Raster* object. Image to normalise.
y	Raster* object. Reference image. Optional. Used to extract min and max values if ymin or ymax are missing.
xmin	Numeric. Min value of x. Either a single value or one value per layer in x. If xmin is not provided it will be extracted from x.
xmax	Numeric. Max value of x. Either a single value or one value per layer in x. If xmax is not provided it will be extracted from x.
ymin	Numeric. Min value of y. Either a single value or one value per layer in x. If ymin is not provided it will be extracted from y.
ymax	Numeric. Max value of y. Either a single value or one value per layer in x. If ymax is not provided it will be extracted from y.
forceMinMax	Logical. Forces update of min and max data slots in x or y.

**Details**

Providing xmin and xmax values manually can be useful if the raster contains a variable of a known, fixed value range, e.g. NDVI from -1 to 1 but the actual pixel values don't encompass this entire range. By providing xmin = -1 and xmax = 1 the values can be rescaled to any other range, e.g. 1 to 100 while comparability to other rescaled NDVI scenes is retained.

**Value**

Returns a Raster\* object of the same dimensions as the input raster x but shifted and stretched to the new limits.

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

```
## Create example data
data(lsat)
lsat2 <- lsat - 1000
lsat2

## Rescale lsat2 to match original lsat value range
lsat2_rescaled <- rescaleImage(lsat2, lsat)
lsat2_rescaled

## Rescale lsat to value range [0,1]
lsat2_unity <- rescaleImage(lsat2, ymin = 0, ymax = 1)
lsat2_unity
```

---

rlogo

*Rlogo as RasterBrick*

---

**Description**

Tiny example of raster data used to run examples.

**Usage**

```
data(rlogo)
```

**Examples**

```
data(rlogo)
ggRGB(rlogo, r = 1, g = 2, b = 3)
```

---

rsOpts

*Set global options for RStoolbox*

---

**Description**

shortcut to options(RStoolbox.\*)

**Usage**

```
rsOpts(verbose)
```

## Arguments

`verbose` Logical. If TRUE many functions will print status messages about the current processing step. By default verbose mode is disabled.

## Examples

```
# rsOpts(verbose=TRUE)
```

---

RStoolbox

*RStoolbox: A Collection of Remote Sensing Tools*

---

## Description

The RStoolbox package provides a set of functions which simplify performing standard remote sensing tasks in R. Most functions have built-in parallel support. All that is required is to run `beginCluster` beforehand.

## Data Import and Export

- `readMeta`: import Landsat metadata from MTL or XML files
- `stackMeta`: load Landsat bands based on metadata
- `readSLI` & `writeSLI`: read and write ENVI spectral libraries
- `saveRSTBX` & `readRSTBX`: save and re-import RStoolbox classification objects (model and map)
- `readEE`: import and tidy EarthExplorer search results

## Data Pre-Processing

- `radCor`: radiometric conversions and corrections. Primarily, yet not exclusively, intended for Landsat data processing. DN to radiance to reflectance conversion as well as DOS approaches
- `topCor`: topographic illumination correction
- `cloudMask` & `cloudShadowMask`: mask clouds and cloud shadows in Landsat or other imagery which comes with a thermal band
- `classifyQA`: extract layers from Landsat 8 QA bands, e.g. cloud confidence
- `rescaleImage`: rescale image to match min/max from another image or a specified min/max range
- `normImage`: normalize imagery by centering and scaling
- `histMatch`: matches the histograms of two scenes
- `coregisterImages`: co-register images based on mutual information
- `panSharpen`: sharpen a coarse resolution image with a high resolution image (typically panchromatic)

## Data Analysis

- `spectralIndices`: calculate a set of predefined multispectral indices like NDVI
- `tasseledCap`: tasseled cap transformation
- `sam`: spectral angle mapper
- `rasterPCA`: principal components transform for raster data
- `rasterCVA`: change vector analysis
- `unsupervisedClass`: unsupervised classification
- `superClass`: supervised classification
- `fCover`: fractional cover of coarse resolution imagery based on high resolution classification

## Data Display

- `ggR`: single raster layer plotting with `ggplot2`
- `ggRGB`: efficient plotting of remote sensing imagery in RGB with `ggplot2`

---

sam

*Spectral Angle Mapper*

---

## Description

Calculates the angle in spectral space between pixels and a set of reference spectra (endmembers) for image classification based on spectral similarity.

## Usage

```
sam(img, em, angles = FALSE, ...)
```

## Arguments

<code>img</code>	RasterBrick or RasterStack. Remote sensing imagery (usually hyperspectral)
<code>em</code>	Matrix or data.frame with endmembers. Each row should contain the endmember spectrum of a class, i.e. columns correspond to bands in <code>img</code> . It is recommended to set the rownames to class names.
<code>angles</code>	Logical. If TRUE a RasterBrick containing each one layer per endmember will be returned containing the spectral angles.
<code>...</code>	further arguments to be passed to <code>writeRaster</code>

## Details

For each pixel the spectral angle mapper calculates the angle between the vector defined by the pixel values and each endmember vector. The result of this is one raster layer for each endmember containing the spectral angle. The smaller the spectral angle the more similar a pixel is to a given endmember class. In a second step one can go ahead and enforce thresholds of maximum angles or simply classify each pixel to the most similar endmember.



**Value**

RasterBrick or RasterLayer If `angles = FALSE` a single Layer will be returned in which each pixel is assigned to the closest endmember class (integer pixel values correspond to row order of em).

**Examples**

```
library(raster)
library(ggplot2)
## Load example data-set
data(lsat)

## Sample endmember spectra
## First location is water, second is open agricultural vegetation
pts <- data.frame(x = c(624720, 627480), y = c(-414690, -411090))
endmembers <- extract(lsat, pts)
rownames(endmembers) <- c("water", "vegetation")

## Calculate spectral angles
lsat_sam <- sam(lsat, endmembers, angles = TRUE)
plot(lsat_sam)

## Classify based on minimum angle
lsat_sam <- sam(lsat, endmembers, angles = FALSE)

ggR(lsat_sam, forceCat = TRUE, geom_raster=TRUE) +
  scale_fill_manual(values = c("blue", "green"), labels = c("water", "vegetation"))
```

---

 saverSTBX

*Save and Read RStoolbox Classification Results*


---

**Description**

Saves objects of classes `unsuperClass`, `superClass`, `rasterPCA` and `fCover` to file. Useful to archive the fitted models.

**Usage**

```
saveRSTBX(x, filename, format = "raster", ...)

readRSTBX(filename)
```

**Arguments**

<code>x</code>	RStoolbox object of classes <code>c("fCover", "rasterPCA", "superClass", "unsuperClass")</code>
<code>filename</code>	Character. Path and filename. Any file extension will be ignored.
<code>format</code>	Character. Driver to use for the raster file
<code>...</code>	further arguments passed to <code>writeRaster</code>

**Value**

The output of writeRSTBX will be at least two files written to disk: a) an .rds file containing the object itself and b) the raster file (depending on the driver you choose this can be more than two files).

**Functions**

- saveRSTBX: Save RStoolbox object to file
- readRSTBX: Read files saved with saveRSTBX

**Note**

All files must be kept in the same directory to read the full object back into R by means of readRSTBX. You can move them to another location but you'll have to move *\*all\** of them (just like you would with Shapefiles). In case the raster file(s) is missing, readRSTBX will still return the object but the raster will be missing.

writeRSTBX and readRSTBX are convenience wrappers around saveRDS, readRDS. This means you can read all files created this way also with base functionality as long as you don't move your files. This is because x\$map is a Raster\* object and hence contains only a static link to the file on disk.

**Examples**

```
## Not run:
input <- brick(system.file("external/rlogo.grd", package="raster"))
## Create filename
file <- paste0(tempdir(), "/test", runif(1))
## Run PCA
rpc <- rasterPCA(input, nSample = 100)
## Save object
saveRSTBX(rpc, filename=file)
## Which files were written?
list.files(tempdir(), pattern = basename(file))
## Re-read files
re_rpc <- readRSTBX(file)
## Remove files
file.remove(list.files(tempdir(), pattern = basename(file), full = TRUE))

## End(Not run)
```

---

spectralIndices

*Spectral Indices*


---

**Description**

Calculate a suite of multispectral indices such as NDVI, SAVI etc. in an efficient way.

**Usage**

```
spectralIndices(img, blue = NULL, green = NULL, red = NULL,
  nir = NULL, swir1 = NULL, swir2 = NULL, swir3 = NULL,
  scaleFactor = 1, skipRefCheck = FALSE, indices = NULL,
  index = NULL, maskLayer = NULL, maskValue = 1, coefs = list(L =
  0.5, G = 2.5, L_evi = 1, C1 = 6, C2 = 7.5, s = 1, swir2ccc = NULL,
  swir2coc = NULL), ...)
```

**Arguments**

<code>img</code>	Raster* object. Typically remote sensing imagery, which is to be classified.
<code>blue</code>	Character or integer. Blue band.
<code>green</code>	Character or integer. Green band.
<code>red</code>	Character or integer. Red band.
<code>nir</code>	Character or integer. Near-infrared band (700-1100nm).
<code>swir1</code>	[temporarily deprecated]
<code>swir2</code>	Character or integer. Short-wave-infrared band (1400-1800nm).
<code>swir3</code>	Character or integer. Short-wave-infrared band (2000-2500nm).
<code>scaleFactor</code>	Numeric. Scale factor for the conversion of scaled reflectances to [0,1] value range (applied as reflectance/scaleFactor) Necessary for calculating EVI/EVI2 with scaled reflectance values.
<code>skipRefCheck</code>	Logical. When EVI/EVI2 is to be calculated there is a rough heuristic check, whether the data are inside [0,1] +/- 0.5 (after applying a potential scaleFactor). If there are invalid reflectances, e.g. clouds with reflectance > 1 this check will result in a false positive and skip EVI calculation. Use this argument to skip this check in such cases *iff* you are sure the data and scaleFactor are valid.
<code>indices</code>	Character. One or more spectral indices to calculate (see Details). By default (NULL) all implemented indices given the spectral bands which are provided will be calculated.
<code>index</code>	Character. Alias for indices.
<code>maskLayer</code>	RasterLayer containing a mask, e.g. clouds, for which pixels are set to NA. Alternatively a layername or -number can be provided if the mask is part of img.
<code>maskValue</code>	Integer. Pixel value in maskLayer which should be masked in output, i.e. will be set to NA in all calculated indices.
<code>coefs</code>	List of coefficients (see Details).
<code>...</code>	further arguments such as filename etc. passed to <a href="#">writeRaster</a>

**Details**

`spectralIndices` calculates all indices in one go in C++, which is more efficient than calculating each index separately (for large rasters). By default all indices which can be calculated given the specified indices will be calculated. If you don't want all indices, use the `indices` argument to specify exactly which indices are to be calculated. See the table below for index names and required bands.

Index values outside the valid value ranges (if such a range exists) will be set to NA. For example a pixel with NDVI > 1 will be set to NA.

Index	Description	Source	Bands	Formula
CTVI	Corrected Transformed Vegetation Index	Perry1984	red, nir	$(NDVI + 0.5) / (s * nir - red)$
DVI	Difference Vegetation Index	Richardson1977	red, nir	$G * ((nir - red) / (nir^2 - red^2))$
EVI	Enhanced Vegetation Index	Huete1999	red, nir, blue	$G * ((nir - red) / (nir^2 - red^2))$
EVI2	Two-band Enhanced Vegetation Index	Jiang 2008	red, nir	$G * ((nir - red) / (nir^2 - red^2))$
GEMI	Global Environmental Monitoring Index	Pinty1992	red, nir	$(nir - green) / (nir + green - swir2)$
GNDVI	Green Normalised Difference Vegetation Index	Gitelson1998	green, nir	$(green - swir2) / (nir + green - swir2)$
MNDWI	Modified Normalised Difference Water Index	Xu2006	green, swir2	$nir + 0.5 - (0.5 * (nir + 1) - (nir - swir3) / (nir + swir3))$
MSAVI	Modified Soil Adjusted Vegetation Index	Qi1994	red, nir	$(2 * (nir + 1) - (nir - swir3) / (nir + swir3)) / (nir + swir3)$
MSAVI2	Modified Soil Adjusted Vegetation Index 2	Qi1994	red, nir	$(nir - swir3) / (nir + swir3)$
NBRI	Normalised Burn Ratio Index	Garcia1991	nir, swir3	$(nir - red) / (nir + red)$
NDVI	Normalised Difference Vegetation Index	Rouse1974	red, nir	$(nir - red) / (nir + red)$
NDVIC	Corrected Normalised Difference Vegetation Index	Nemani1993	red, nir, swir2	$(nir - red) / (nir + red)$
NDWI	Normalised Difference Water Index	McFeeters1996	green, nir	$(green - nir) / (green + nir)$
NDWI2	Normalised Difference Water Index	Gao1996	nir, swir2	$(nir - swir2) / (nir + swir2)$
NRVI	Normalised Ratio Vegetation Index	Baret1991	red, nir	$(red/nir - 1) / (red/nir + 1)$
RVI	Ratio Vegetation Index		red, nir	$red/nir$
SATVI	Soil Adjusted Total Vegetation Index	Marsett2006	red, swir2, swir3	$(swir2 - red) / (nir - red) * (1 - L)$
SAVI	Soil Adjusted Vegetation Index	Huete1988	red, nir	$nir / (red + swir2)$
SLAVI	Specific Leaf Area Vegetation Index	Lymburger2000	red, nir, swir2	$nir/red$
SR	Simple Ratio Vegetation Index	Birth1968	red, nir	$sqrt((nir - red) / (nir + red))$
TVI	Transformed Vegetation Index	Deering1975	red, nir	$sqrt(abs((nir - red) / (nir + red)))$
TTVI	Thiam's Transformed Vegetation Index	Thiam1997	red, nir	$nir - s * red$
WDVI	Weighted Difference Vegetation Index	Richardson1977	red, nir	

Some indices require additional parameters, such as the slope of the soil line which are specified via a list to the `coefs` argument. Although the defaults are sensible values, values like the soil brightness factor `L` for SAVI should be adapted depending on the characteristics of the scene. The coefficients are:

Coefficient	Description	Affected Indices
<code>s</code>	slope of the soil line	DVI, WDVI
<code>L_evi</code> , <code>C1</code> , <code>C2</code> , <code>G</code>	various	EVI
<code>L</code>	soil brightness factor	SAVI, SATVI
<code>swir2ccc</code>	minimum <code>swir2</code> value (completely closed forest canopy)	NDVIC
<code>swir2coc</code>	maximum <code>swir2</code> value (completely open canopy)	NDVIC

The wavelength band names are defined following Schowengertd 2007, p10. The last column shows exemplarily which Landsat 5 TM bands correspond to which wavelength range definition.

Band	Description	Wavl_min	Wavl_max	Landsat5_Band
vis	visible	400	700	1,2,3
nir	near infra-red	700	1100	4
swir1	short-wave infra-red	1100	1351	-
swir2	short-wave infra-red	1400	1800	5
swir3	short-wave infra-red	2000	2500	7
mir1	mid-wave infra-red	3000	4000	-

mir2	mid-wave infra-red	45000	5000	-
tir1	thermal infra-red	8000	9500	-
tir2	thermal infra-red	10000	140000	6

**Value**

RasterBrick or a RasterLayer if length(indices) == 1

**Examples**

```
library(ggplot2)
library(raster)
data(lsat)

## Calculate NDVI
ndvi <- spectralIndices(lsat, red = "B3_dn", nir = "B4_dn", indices = "NDVI")
ndvi
ggR(ndvi, geom_raster = TRUE) +
  scale_fill_gradientn(colours = c("black", "white"))

## Calculate all possible indices, given the provided bands
## Convert DNS to reflectance (required to calculate EVI and EVI2)
mtlFile <- system.file("external/landsat/LT52240631988227CUB02_MTL.txt", package="RStoolbox")
lsat_ref <- radCor(lsat, mtlFile, method = "apref")

SI <- spectralIndices(lsat_ref, red = "B3_tre", nir = "B4_tre")
plot(SI)
```

---

srtm

*SRTM Digital Elevation Model*


---

**Description**

DEM for the Landsat example area taken from SRTM v3 tile: s04\_w050\_1arc\_v3.tif

**Usage**

```
data(srtm)
```

**Examples**

```
data(srtm)
ggR(srtm)
```

stackMeta

*Import separate Landsat files into single stack***Description**

Reads Landsat MTL or XML metadata files and loads single Landsat Tiffs into a rasterStack. Be aware that by default stackMeta() does NOT import panchromatic bands nor thermal bands with resolutions != 30m.

**Usage**

```
stackMeta(file, quantity = "all", category = "image",
          allResolutions = FALSE)
```

**Arguments**

file	Character. Path to Landsat MTL metadata (*.MTL.txt) file or an Landsat CDR xml metadata file (*.xml).
quantity	Character vector. Which quantity should be returned. Options: digital numbers ('dn'), top of atmosphere reflectance ('tre'), at surface reflectance ('sre'), brightness temperature ('bt'), spectral index ('index'), all quantities ('all').
category	Character vector. Which category of data to return. Options 'image': image data, 'pan': panchromatic image, 'index': multiband indices, 'qa' quality flag bands, 'all': all categories.
allResolutions	Logical. if TRUE a list will be returned with length = unique spatial resolutions.

**Value**

Returns one single RasterStack comprising all requested bands. If allResolutions = TRUE \*and\* there are different resolution layers (e.g. a 15m panchromatic band along wit 30m imagery) a list of RasterStacks will be returned.

**Note**

Be aware that by default stackMeta() does NOT import panchromatic bands nor thermal bands with resolutions != 30m. Use the allResolutions argument to import all layers. Note that nowadays the USGS uses cubic convolution to resample the TIR bands to 30m resolution.

**Examples**

```
library(rgdal)
## Example metadata file (MTL)
mtlFile <- system.file("external/landsat/LT52240631988227CUB02_MTL.txt", package="RStoolbox")

## Read metadata
metaData <- readMeta(mtlFile)
summary(metaData)
```

```
## Load rasters based on metadata file
lsat <- stackMeta(mtlFile)
lsat
```

---

superClass                      *Supervised Classification*

---

## Description

Supervised classification both for classification and regression mode based on vector training data (points or polygons).

## Usage

```
superClass(img, trainData, valData = NULL, responseCol = NULL,
  nSamples = 1000, polygonBasedCV = FALSE, trainPartition = NULL,
  model = "rf", tuneLength = 3, kfold = 5, minDist = 2,
  mode = "classification", predict = TRUE, predType = "raw",
  filename = NULL, verbose, overwrite = TRUE, ...)
```

## Arguments

img	Raster* object. Typically remote sensing imagery, which is to be classified.
trainData	SpatialPolygonsDataFrame or SpatialPointsDataFrame containing the training locations.
valData	SpatialPolygonsDataFrame or SpatialPointsDataFrame containing the validation locations (optional).
responseCol	Character or integer giving the column in trainData, which contains the response variable. Can be omitted, when trainData has only one column.
nSamples	Integer. Number of samples per land cover class.
polygonBasedCV	Logical. If TRUE model tuning during cross-validation is conducted on a per-polygon basis. Use this to deal with overfitting issues. Does not affect training data supplied as SpatialPointsDataFrames.
trainPartition	Numeric. Partition (polygon based) of trainData that goes into the training data set between zero and one. Ignored if valData is provided.
model	Character. Which model to use. See <a href="#">train</a> for options. Defaults to random-Forest ('rf'). In addition to the standard caret models, a maximum likelihood classification is available via model = 'mlc'.
tuneLength	Integer. Number of levels for each tuning parameter (see <a href="#">train</a> for details).
kfold	Integer. Number of cross-validation resamples during model tuning.
minDist	Numeric. Minimum distance between training and validation data, e.g. minDist=1 clips validation polygons to ensure a minimal distance of one pixel (pixel size according to img) to the next training polygon. Requires all data to carry valid projection information.



mode	Character. Model type: 'regression' or 'classification'.
predict	Logical. Produce a map (TRUE, default) or only fit and validate the model (FALSE).
predType	Character. Type of the final output raster. Either "raw" for class predictions or "prob" for class probabilities. Class probabilities are not available for all classification models ( <a href="#">predict.train</a> ).
filename	Path to output file (optional). If NULL, standard raster handling will apply, i.e. storage either in memory or in the raster temp directory.
verbose	Logical. prints progress and statistics during execution
overwrite	logical. Overwrite spatial prediction raster if it already exists.
...	further arguments to be passed to <a href="#">train</a>

## Details

SuperClass performs the following steps:

1. Ensure non-overlap between training and validation data. This is necessary to avoid biased performance estimates. A minimum distance (`minDist`) in pixels can be provided to enforce a given distance between training and validation data.
2. Sample training coordinates. If `trainData` (and `valData` if present) are `SpatialPolygons-DataFrames` `superClass` will calculate the area per polygon and sample `nSamples` locations per class within these polygons. The number of samples per individual polygon scales with the polygon area, i.e. the bigger the polygon, the more samples.
3. Split training/validation If `valData` was provided (reccomended) the samples from these polygons will be held-out and not used for model fitting but only for validation. If `trainPartition` is provided the trainingPolygons will be divided into training polygons and validation polygons.
4. Extract raster data The predictor values on the sample pixels are extracted from `img`
5. Fit the model. Using `caret::train` on the sampled training data the model will be fit, including parameter tuning (`tuneLength`) in `kfold` cross-validation. `polygonBasedCV=TRUE` will define cross-validation folds based on polygons (reccomended) otherwise it will be performed on a per-pixel basis.
6. Predict the classes of all pixels in `img` based on the final model.
7. Validate the model with the independent validation data.

## Value

A list containing `[[1]]` the model, `[[2]]` the predicted raster and `[[3]]` the class mapping

## See Also

[train](#)

**Examples**

```

library(caret)
library(randomForest)
library(e1071)
library(raster)
data(rlogo)
train <- readRDS(system.file("external/trainingPoints.rds", package="RStoolbox"))

## Plot training data
olpar <- par(no.readonly = TRUE) # back-up par
par(mfrow=c(1,2))
colors <- c("yellow", "green", "deppink")
plotRGB(rlogo)
plot(train, add = TRUE, col = colors[train$class], pch = 19)

## Fit classifier (splitting training into 70% training data, 30% validation data)
SC <- superClass(rlogo, trainData = train, responseCol = "class",
model = "rf", tuneLength = 1, trainPartition = 0.7)
SC

## Plots
plot(SC$map, col = colors, legend = FALSE, axes = FALSE, box = FALSE)
legend(1,1, legend = levels(train$class), fill = colors, title = "Classes",
horiz = TRUE, bty = "n")
par(olpar) # reset par

```

---

tasseledCap

*Tasseled Cap Transformation*


---

**Description**

Calculates brightness, greenness and wetness from multispectral imagery. Currently implemented Landsat 4 TM, Landsat 5 TM, Landsat 7ETM+, Landsat 8 OLI, MODIS, QuickBird, Spot5 and RapidEye.

**Usage**

```
tasseledCap(img, sat, ...)
```

**Arguments**

img	RasterBrick or RasterStack. Input image. Band order must correspond to sensor specifications (see Details and Examples)
sat	Character. Sensor; one of: c("Landsat4TM", "Landsat5TM", "Landsat7ETM", "Landsat8OLI", "MODIS", "QuickBird", "Spot5", "RapidEye"). Case is irrelevant.
...	Further arguments passed to writeRaster.

## Details

Currently implemented: Landsat 4 TM, Landsat 5 TM, Landsat 7ETM+, Landsat 8 OLI, MODIS, QuickBird, Spot5, RapdiEye. Input data must be in top of atmosphere reflectance. Moreover, bands must be provided in ascending order as listed in the table below. Irrelevant bands, such as Landsat Thermal Bands or QuickBird/Spot5 Panchromatic Bands must be omitted. Required bands are:

sat	bands	coefficients	data unit
Landsat4TM	1,2,3,4,5,7	Crist 1985	reflectance
Landsat5TM	1,2,3,4,5,7	Crist 1985	reflectance
Landsat7ETM	1,2,3,4,5,7	Huang 2002	reflectance
Landsat8OLI	2,3,4,5,6,7	Baig 2014	reflectance
MODIS	1,2,3,4,5,6,7	Lobser 2007	reflectance
QuickBird	2,3,4,5	Yarbrough 2005	reflectance
Spot5	2,3,4,5	Ivtis 2008	reflectance
RapidEye	1,2,3,4,5	Schoenert 2014	reflectance

## Value

Returns a RasterBrick with the three bands: brightness, greenness, and (soil) wetness.

## References

- Crist (1985) "A TM Tasseled Cap Equivalent Transformation for Reflectance Factor Data." *Remote Sensing of Environment* 17 (3): 301-306
- Huang et al. (2002) "Derivation of a Tasseled Cap Transformation Based on Landsat 7 At-Satellite Reflectance." *International Journal of Remote Sensing* 23 (8): 1741-1748
- Baig et al. (2014) "Derivation of a Tasseled Cap Transformation Based on Landsat 8 At-Satellite Reflectance." *Remote Sensing Letters* 5 (5): 423-431.
- Lobser et al. (2007) "MODIS Tasseled Cap: Land Cover Characteristics Expressed through Transformed MODIS Data." *International Journal of Remote Sensing* 28 (22): 5079-5101.
- Yarbrough et al. (2005) "QuickBird 2 tasseled cap transform coefficients: a comparison of derivation methods." *Pecora 16 Global Priorities in Land Remote Sensing*: 23-27.
- Ivtis et al. (2008) "Orthogonal transformation of segmented SPOT5 images." *Photogrammetric Engineering & Remote Sensing* 74 (11): 1351-1364.
- Schoenert et al. (2014) "Derivation of tasseled cap coefficients for RapidEye data." *Earth Resources and Environmental Remote Sensing/GIS Applications V (9245)*: 92450Qs.

## Examples

```
library(raster)
data(lsat)

## Run tasseled cap (exclude thermal band 6)
lsat_tc <- tasseledCap(lsat[[c(1:5,7)]], sat = "Landsat5TM")
lsat_tc
plot(lsat_tc)
```

---

topCor *Topographic Illumination Correction*

---

### Description

account and correct for changes in illumination due to terrain elevation.

### Usage

```
topCor(img, dem, metaData, solarAngles = c(), method = "C",
       stratImg = NULL, nStrat = 5, illu, ...)
```

### Arguments

img	Raster*. Imagery to correct
dem	Raster*. Either a digital elevation model as a RasterLayer or a RasterStack/Brick with pre-calculated slope and aspect (see <a href="#">terrain</a> ) in which case the layers must be named 'slope' and 'aspect'. Must have the same dimensions as img.
metaData	Character, ImageMetaData. Either a path to a Landsat meta-data file (MTL) or an ImageMetaData object (see <a href="#">readMeta</a> )
solarAngles	Numeric vector containing sun azimuth and sun zenith (in radians and in that order). Not needed if metaData is provided
method	Character. One of c("cos", "avgcos", "minnaert", "C", "stat", "illu"). Choosing 'illu' will return only the local illumination map.
stratImg	RasterLayer to define strata, e.g. NDVI. Or the string 'slope' in which case stratification will be on nStrat slope classes. Only relevant if method = 'minnaert'.
nStrat	Integer. Number of bins or quantiles to stratify by. If a bin has less than 50 samples it will be merged with the next bin. Only relevant if method = 'minnaert'.
illu	Raster*. Optional pre-calculated illumination map. Run topCor with method="illu" to calculate an illumination map
...	arguments passed to <a href="#">writeRaster</a>

### Details

For detailed discussion of the various approaches please see Riano et al. (2003).

The minnaert correction can be stratified for different landcover characteristics. If stratImg = 'slope' the analysis is stratified by the slope, i.e. the slope values are divided into nStrat classes and the correction coefficient k is calculated and applied separately for each slope class. An alternative could be to stratify by a vegetation index in which case an additional raster layer has to be provided via the stratImg argument.

### References

Riano et al. (2003) Assessment of different topographic correction in Landsat-TM data for mapping vegetation types. IEEE Transactions on Geoscience and Remote Sensing.

**Examples**

```
## Load example data
metaData <- system.file("external/landsat/LT52240631988227CUB02_MTL.txt", package="RStoolbox")
metaData <- readMeta(metaData)
lsat      <- stackMeta(metaData)
data(srtm)

## Minnaert correction, solar angles from metaData
lsat_minnaert <- topCor(lsat, dem = srtm, metaData = metaData, method = "minnaert")

## C correction, solar angles provided manually
lsat_C <- topCor(lsat, dem = srtm, solarAngles = c(1.081533, 0.7023922), method = "C")
```

---

unsuperClass

*Unsupervised Classification*


---

**Description**

Unsupervised clustering of Raster\* data using kmeans clustering

**Usage**

```
unsuperClass(img, nSamples = 10000, nClasses = 5, nStarts = 25,
             nIter = 100, norm = FALSE, clusterMap = TRUE,
             algorithm = "Hartigan-Wong", ...)
```

**Arguments**

img	Raster* object.
nSamples	Integer. Number of random samples to draw to fit cluster map. Only relevant if clusterMap = TRUE.
nClasses	Integer. Number of classes.
nStarts	Integer. Number of random starts for kmeans algorithm.
nIter	Integer. Maximal number of iterations allowed.
norm	Logical. If TRUE will normalize img first using <a href="#">normImage</a> . Normalizing is beneficial if your predictors have different scales.
clusterMap	Logical. Fit kmeans model to a random subset of the img (see Details).
algorithm	Character. <a href="#">kmeans</a> algorithm. One of c("Hartigan-Wong", "Lloyd", "MacQueen")
...	further arguments to be passed to <a href="#">writeRaster</a> , e.g. filename

## Details

Clustering is done using `kmeans`. This can be done for all pixels of the image (`clusterMap=FALSE`), however this can be slow and is not memory safe. Therefore if you have large raster data (> memory), as is typically the case with remote sensing imagery it is advisable to choose `clusterMap=TRUE` (the default). This means that a `kmeans` cluster model is calculated based on a random subset of pixels (`nSamples`). Then the distance of *\*all\** pixels to the cluster centers is calculated in a stepwise fashion using `predict`. Class assignment is based on minimum euclidean distance to the cluster centers.

The solution of the `kmeans` algorithm often depends on the initial configuration of class centers which is chosen randomly. Therefore, `kmeans` is usually run with multiple random starting configurations in order to find a convergent solution from different starting configurations. The `nStarts` argument allows to specify how many random starts are conducted.

## Examples

```
library(raster)
input <- brick(system.file("external/rlogo.grd", package="raster"))

## Plot
olpar <- par(no.readonly = TRUE) # back-up par
par(mfrow=c(1,2))
plotRGB(input)

## Run classification
set.seed(25)
unC <- unsuperClass(input, nSamples = 100, nClasses = 5, nStarts = 5)
unC

## Plots
colors <- rainbow(5)
plot(unC$map, col = colors, legend = FALSE, axes = FALSE, box = FALSE)
legend(1,1, legend = paste0("C",1:5), fill = colors,
      title = "Classes", horiz = TRUE, bty = "n")

par(olpar) # reset par
```

---

validateMap

*Map accuracy assessment*

---

## Description

validate a map from a classification or regression model. This can be useful to update the accuracy assessment after filtering, e.g. for a minimum mapping unit.

## Usage

```
validateMap(map, valData, responseCol, nSamples = 500,
  mode = "classification", classMapping = NULL)
```

**Arguments**

map	RasterLayer. The classified map.
valData	SpatialPolygonsDataFrame or SpatialPointsDataFrame with validation data.
responseCol	Character. Column containing the validation data in attribute table of valData.
nSamples	Integer. Number of pixels to sample for validation (only applies to polygons).
mode	Character. Either 'classification' or 'regression'.
classMapping	optional data.frame with columns 'class' and 'classID' defining the mapping from raster integers to class names.

**Note**

Polygons, which are smaller than the map resolution will only be considered if they overlap with a pixel center coordinate, otherwise they will be ignored.

**Examples**

```
## Not run:
library(caret)
library(raster)

## Training data
data(lsat)
poly <- readRDS(system.file("external/trainingPolygons.rds", package="RStoolbox"))

## Split training data in training and validation set (50%-50%)
splitIn <- createDataPartition(poly$class, p = .5)[[1]]
train <- poly[splitIn,]
val <- poly[-splitIn,]

## Classify (deliberately poorly)
sc <- superClass(lsat, trainData = train, responseCol = "class", nSamples = 50, model = "mlc")

## Polish map with majority filter

polishedMap <- focal(sc$map, matrix(1,3,3), fun = modal)

## Validation
## Before filtering
val0 <- validateMap(sc$map, valData = val, responseCol = "class",
                    classMapping = sc$classMapping)

## After filtering
val1 <- validateMap(polishedMap, valData = val, responseCol = "class",
                    classMapping = sc$classMapping)

## End(Not run)
```

---

writeSLI                      *Write ENVI spectral libraries*

---

### Description

Writes binary ENVI spectral library files (sli) with accompanying header (.sli.hdr) files OR ASCII spectral library files in ENVI format.

### Usage

```
writeSLI(x, path, wavl.units = "Micrometers", scaleF = 1,
         mode = "bin")
```

### Arguments

x	data.frame with first column containing wavelengths and all other columns containing spectra.
path	path to spectral library file to be created.
wavl.units	wavelength units. Defaults to Micrometers. Nanometers is another typical option.
scaleF	optional reflectance scaling factor. Defaults to 1.
mode	character string specifying output file type. Must be one of "bin" for binary .sli files or "ASCII" for ASCII ENVI plot files.

### Details

ENVI spectral libraries with ending .sli are binary arrays with spectra saved in rows.

### See Also

[readSLI](#)

### Examples

```
## Example data
sliFile <- system.file("external/vegSpec.sli", package="RStoolbox")
sliTmpFile <- paste0(tempdir(), "/vegetationSpectra.sli")

## Read spectral library
sli <- readSLI(sliFile)
head(sli)
plot(sli[,1:2], col = "orange", type = "l")
lines(sli[,c(1,3)], col = "green")

## Write to binary spectral library
writeSLI(sli, path = sliTmpFile)
```



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