

# Package ‘pavo’

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**Description** A cohesive framework for parsing, analyzing and organizing color from spectral data.

**Title** Perceptual Analysis, Visualization and Organization of Spectral Color Data in R

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**URL** <http://rafaelmaia.net/pavo/>, <https://github.com/rmaia/pavo/>

**BugReports** <https://github.com/rmaia/pavo/issues>

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<i>adjacent</i>	<i>Run an adjacency and boundary strength analysis</i>
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### Description

Calculate summary variables from the adjacency (Endler 2012) and boundary-strength (Endler et al. 2018) analyses, along with overall pattern contrast (Endler & Mielke 2005).

### Usage

```
adjacent(classimg, xpts = NULL, xscale = NULL, bkgID = NULL,
         polygon = NULL, exclude = c("none", "background", "object"),
         coldists = NULL, hsl = NULL, cores = getOption("mc.cores", 2L))
```

### Arguments

- |                       |  |
|-----------------------|--|
| <code>classimg</code> | (required) an xyz matrix, or list of matrices, in which x and y correspond to spatial (e.g. pixel) coordinates, and z is a numeric code specifying a colour-class. Preferably the result of <code>classify</code> , or constructed from grid-sampled spectra that have been visually modelled and clustered (as per Endler 2012).  |
| <code>xpts</code>     | an integer specifying the number of sample points along the x axis, from which the evenly-spaced sampling grid is constructed (if required). Defaults to the smallest dimension of <code>classimg</code> , though this should be carefully considered.   |
| <code>xscale</code>   | (required) an integer specifying the true length of the x-axis, in preferred units. Not required, and ignored, only if image scales have been set via <code>procimg</code> .   |
| <code>bkgID</code>    | an integer or vector specifying the colour-class ID number(s) of pertaining to the background alone, for relatively homogeneous and uniquely-identified backgrounds (e.g. the matte background of pinned specimens). Examine the attributes of, or call <code>summary</code> on, the result of <code>classify</code> to visualise the RGB values corresponding to colour-class ID numbers for classified images. Ignored if the focal object and background has been identified using <code>procimg</code> . |
| <code>polygon</code>  | a data.frame of x-y coordinates delineating a closed polygon that separates the focal object from the background. Not required, and ignored, if the focal object outline is specified using <code>procimg</code> .   |
| <code>exclude</code>  | the portion of the scene to be excluded from the analysis, if any. <ul style="list-style-type: none"> <li>'none': default</li> <li>'background': exclude everything <i>outside</i> the closed polygon specified using <code>procimg</code>, or the argument <code>polygon</code>. Alternatively, if the background is relatively homogeneous the colour-class ID(s) uniquely corresponding to the background can be specified via <code>bkgID</code>, and subsequently excluded.</li> </ul>                  |

	<ul style="list-style-type: none"> <li>• 'object': exclude everything <i>inside</i> the closed polygon specified using <a href="#">procimg</a>, or the argument polygon.</li> </ul>
coldists	a data.frame specifying the visually-modelled chromatic (dS) and/or achromatic (dL) distances between colour-categories. The first two columns should be named 'c1' and 'c2', and specify all possible combinations of numeric colour-class ID's (viewable by calling <code>summary(image, plot = TRUE)</code> on a colour classified image), with the remaining columns named dS (for chromatic distances) and/or dL (for achromatic distances). See <a href="#">vismodel</a> and <a href="#">colspace</a> for visual modelling with spectral data.
hsl	data.frame specifying the hue, saturation, and luminance of color patch elements, as might be estimated via <a href="#">vismodel</a> and <a href="#">colspace</a> . The first column, named 'patch', should contain numeric color category IDs, with the remaining columns specifying one or more of 'hue' (angle, in radians), 'sat', and/or 'lum'.
cores	number of cores to be used in parallel processing. If 1, parallel computing will not be used. Defaults to <code>getOption("mc.cores", 2L)</code> . Not available on Windows.

## Value

a data frame of summary variables:

- 'k': The number of user-specified colour and/or luminance classes.
- 'N': The grand total (sum of diagonal and off-diagonal) transitions.
- 'n\_off': The total off-diagonal transitions.
- 'p\_i': The overall frequency of colour class *i*.
- 'q\_i\_j': The frequency of transitions between *all* colour classes *i* and *j*, such that  $\sum(q_{i_j}) = 1$ .
- 't\_i\_j': The frequency of off-diagonal (i.e. class-change transitions) transitions *i* and *j*, such that  $\sum(t_{i_j}) = 1$ .
- 'm': The overall transition density (mean transitions), in units specified in the argument `xscale`.
- 'm\_r': The row-wise transition density (mean row transitions), in user-specified units.
- 'm\_c': The column-wise transition density (mean column transitions), in user-specified units.
- 'A': The transition aspect ratio ( $< 1 = \text{wide}$ ,  $> 1 = \text{tall}$ ).
- 'Sc': Simpson colour class diversity,  $Sc = 1/(\sum(p_i^2))$ . If all colour and luminance classes are equal in relative area, then  $Sc = k$ .
- 'St': Simpson transition diversity,  $St = 1/\sum(t_{i_j}^2)$ .
- 'Jc': Simpson colour class diversity relative to its achievable maximum.  $Jc = Sc/k$ .
- 'Jt': Simpson transition diversity relative to its achievable maximum.  $Jt = St/(k*(k-1)/2)$ .
- 'B': The animal/background transition ratio, or the ratio of class-change transitions entirely within the focal object and those involving the object and background,  $B = \sum(O_{a_a} / O_{a_b})$ .
- 'Rt': Ratio of animal-animal and animal-background transition diversities,  $Rt = St_{a_a} / St_{a_b}$ .
- 'Rab': Ratio of animal-animal and background-background transition diversities,  $Rt = St_{a_a} / St_{b_b}$ .

- 'm\_dS', 's\_dS', 'cv\_dS': weighted mean, sd, and coefficient of variation of the chromatic boundary strength.
- 'm\_dL', 's\_dL', 'cv\_dL': weighted mean, sd, and coefficient of variation of the achromatic boundary strength.
- 'm\_hue', 's\_hue', 'var\_hue': circular mean, sd, and variance of overall pattern hue (in radians).
- 'm\_sat', 's\_sat', 'cv\_sat': weighted mean, sd, and coefficient variation of overall pattern saturation.
- 'm\_lum', 's\_lum', 'cv\_lum': weighted mean, sd, and coefficient variation of overall pattern luminance.

### Author(s)

Thomas E. White <thomas.white026@gmail.com>

### References

- Endler, J. A. (2012). A framework for analysing colour pattern geometry: adjacent colours. *Biological Journal Of The Linnean Society*, 107(2), 233-253.
- Endler, J. A., Cole G., Kranz A. (2018). Boundary Strength Analysis: Combining color pattern geometry and coloured patch visual properties for use in predicting behaviour and fitness. *Methods in Ecology and Evolution*, 9(12), 2334-2348.
- Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.

### Examples

```
## Not run:
# Set a seed, for reproducibility
set.seed(153)

# Single image
papilio <- getimg(system.file("testdata/images/papilio.png", package = "pavo"))
papilio_class <- classify(papilio, kcols = 4)
papilio_adj <- adjacent(papilio_class, xscale = 100)

# Single image, with (fake) color distances and hsl values
# Fake color distances
distances <- data.frame(
  c1 = c(1, 1, 1, 2, 2, 3),
  c2 = c(2, 3, 4, 3, 4, 4),
  dS = c(5.3, 3.5, 5.7, 2.9, 6.1, 3.2),
  dL = c(5.5, 6.6, 3.3, 2.2, 4.4, 6.6)
)

# Fake hue, saturation, luminance values
hsl_vals <- data.frame(
  patch = seq_len(4),
  hue = c(1.5, 2.2, 1.0, 0.5),
```

```

lum = c(10, 5, 7, 3),
sat = c(3.5, 1.1, 6.3, 1.3)
)

# Full analysis, including the white background's ID
papilio_adj <- adjacent(papilio_class,
  xscale = 100, bkgID = 1,
  coldists = distances, hsl = hsl_vals
)

# Multiple images
snakes <- getimg(system.file("testdata/images/snakes", package = "pavo"))
snakes_class <- classify(snakes, kcols = 3)
snakes_adj <- adjacent(snakes_class, xpts = 120, xscale = c(50, 55))

## End(Not run)

```

---

aggplot

*Plot aggregated reflectance spectra*


---

## Description

Combines and plots spectra (by taking the average and the standard deviation, for example) according to an index or a vector of identities.

## Usage

```
aggplot(rspeccdata, by = NULL, FUN.center = mean, FUN.error = sd,
  lcol = NULL, shadecol = NULL, alpha = 0.2, legend = FALSE, ...)
```

## Arguments

rspeccdata	(required) data frame containing the spectra to be manipulated and plotted.
by	(required) either a single value specifying the range of spectra within the data frame to be combined (for example, <code>by = 3</code> indicates the function will be applied to groups of 3 consecutive columns in the spectra data frame) or a vector containing identifications for the columns in the spectra data frame (in which case the function will be applied to each group of spectra sharing the same identification).
FUN.center	the function to be applied to the groups of spectra, calculating a measure of central tendency (defaults to <code>mean</code> ).
FUN.error	the function to be applied to the groups of spectra, calculating a measure of variation (defaults to <code>sd</code> ).
lcol	color of plotted lines indicating central tendency.
shadecol	color of shaded areas indicating variance measure.
alpha	transparency of the shaded areas.

legend            automatically add a legend.  
 ...              additional graphical parameters to be passed to plot.

### Value

Plot containing the lines and shaded areas of the groups of spectra.

### Author(s)

Rafael Maia <rm72@zips.uakron.edu>, Chad Eliason <cme16@zips.uakron.edu>

### References

Montgomerie R (2006) Analyzing colors. In: Hill G, McGraw K (eds) Bird coloration. Harvard University Press, Cambridge, pp 90-147.

### Examples

```
# Load reflectance data
data(sicalis)

# Create grouping variable based on spec names
bysic <- gsub("^ind[0-9].", "", names(sicalis)[-1])

# Plot using various error functions and options
aggplot(sicalis, bysic)
aggplot(sicalis, bysic, FUN.error = function(x) quantile(x, c(0.0275, 0.975)))
aggplot(sicalis, bysic, shadecol = spec2rgb(sicalis), lcol = 1)
aggplot(sicalis, bysic, lcol = 1, FUN.error = function(x) sd(x) / sqrt(length(x)))
```

---

aggspec

*Aggregate reflectance spectra*

---

### Description

Combines spectra (by taking the average, for example) according to an index or a vector of identities.

### Usage

```
aggspec(rspeccdata, by = NULL, FUN = mean, trim = TRUE)
```

## Arguments

<code>rspecdata</code>	(required) data frame, possibly of class <code>rspec</code> containing the spectra to be manipulated. If it contains a wavelength column named "wl", that column will be ignored.
<code>by</code>	(required) either a single value specifying the range of spectra within the data frame to be combined (for example, <code>by = 3</code> indicates the function will be applied to groups of 3 consecutive columns in the spectra data frame); a vector containing identifications for the columns in the spectra data frame (in which case the function will be applied to each group of spectra sharing the same identification); or a list of vectors, e.g., <code>by = list(sex, species)</code> .
<code>FUN</code>	the function to be applied to the groups of spectra. (defaults to <a href="#">mean</a> )
<code>trim</code>	logical. if TRUE (default), the function will try to identify and remove numbers at the end of the names of the columns in the new <code>rspec</code> object.

## Value

A data frame of class `rspec` containing the spectra after applying the aggregating function.

## Author(s)

Chad Eliason <cme16@zips.uakron.edu>

## References

Montgomerie R (2006) Analyzing colors. In: Hill G, McGraw K (eds) Bird coloration. Harvard University Press, Cambridge, pp 90-147.

## Examples

```
data(teal)

# Average every two spectra
teal.sset1 <- aggspec(teal, by = 2)
plot(teal.sset1)

# Create factor and average spectra by levels 'a' and 'b'
ind <- rep(c("a", "b"), times = 6)
teal.sset2 <- aggspec(teal, by = ind)

plot(teal.sset2)
```



---

as.rimg	<i>Convert data to an rimg object</i>
---------	---------------------------------------

---

**Description**

Converts an array containing RGB image data data to an rimg object.

**Usage**

```
as.rimg(object, name = "img")  
  
is.rimg(object)
```

**Arguments**

object	(required) a three-dimensional array containing RGB values.
name	the name(s) of the image(s).

**Value**

an object of class rimg for use in further pavo functions  
a logical value indicating whether the object is of class rimg

**Author(s)**

Thomas E. White <thomas.white026@gmail.com>

**Examples**

```
# Generate some fake image data  
fake <- array(c(  
  as.matrix(rep(c(0.2, 0.4, 0.6), each = 250)),  
  as.matrix(rep(c(0.4, 0.7, 0.8), each = 250)),  
  as.matrix(rep(c(0.6, 0.1, 0.2), each = 250))  
)  
  ,  
  dim = c(750, 750, 3)  
)  
  
# Inspect it  
head(fake)  
  
# Determine if is rimg object  
is.rimg(fake)  
  
# Convert to rimg object and check again  
fake2 <- as.rimg(fake)  
is.rimg(fake2)
```

---

`as.rspec`*Convert data to an rspec object*

---

**Description**

Converts data frames or matrices containing spectral data to rspec object

**Usage**

```
as.rspec(object, whichwl = NULL, interp = TRUE, lim = NULL,
         exceed.range = TRUE)
```

```
is.rspec(object)
```

**Arguments**

<code>object</code>	(required) a data frame or matrix containing spectra to process.
<code>whichwl</code>	a numeric or character vector specifying which column contains wavelengths. If NULL (default), function searches for column containing equally spaced numbers and sets it as wavelengths "wl". If no wavelengths are found or <code>whichwl</code> is not given, returns arbitrary index values.
<code>interp</code>	whether to interpolate wavelengths in 1-nm bins (defaults to TRUE).
<code>lim</code>	vector specifying wavelength range to interpolate over (e.g. <code>c(300, 700)</code> ).
<code>exceed.range</code>	logical. Should data be interpolated to the limits specified by <code>lim</code> if <code>lim</code> exceeds the range of the actual data? Useful, and relatively safe, when the data range falls slightly within <code>lim</code> (e.g. 300.1 - 699 nm), but will produce spurious results if <code>lim</code> far exceeds the range of input data. Defaults to TRUE.

**Value**

an object of class `rspec` for use in further pavo functions

a logical value indicating whether the object is of class `rspec`

**Author(s)**

Chad Eliason <cme16@zip.uakron.edu>

**Examples**

```
# Generate some fake reflectance data
fakedat <- data.frame(wl = 300:700, ref1 = rnorm(401), ref12 = rnorm(401))
head(fakedat)

# Determine if is rspec object
is.rspec(fakedat)
```

```
# Convert to rspec object
fakedat2 <- as.rspec(fakedat)
is.rspec(fakedat2)
head(fakedat2)
```

---

axistetra

*Plot reference axes in a static tetrahedral colourspace*


---

### Description

Plots reference x, y and z arrows showing the direction of the axes in a static tetrahedral colourspace plot.

### Usage

```
axistetra(x = 0, y = 1.3, size = 0.1, arrowhead = 0.05,
  col = par("fg"), lty = par("lty"), lwd = par("lwd"),
  label = TRUE, adj.label = list(x = c(0.003, 0), y = c(0.003, 0.003),
  z = c(0, 0.003)), label.cex = 1, label.col = NULL)
```

### Arguments

x, y	position of the legend relative to plot limits (usually a value between 0 and 1, but because of the perspective distortion, values greater than 1 are possible)
size	length of the arrows. Can be either a single value (applied for x, y and z) or a vector of 3 separate values for each axis.
arrowhead	size of the arrowhead.
col, lty, lwd	graphical parameters for the arrows.
label	logical, include x, y and z labels (defaults to TRUE).
adj.label	position adjustment for the labels. a list of 3 named objects for x, y and z arrows, each with 2 values for x and y adjustment.
label.cex, label.col	graphical parameters for the labels.

### Value

axistetra adds reference arrows showing the direction of the 3-dimensional axes in a static tetrahedral colourspace plot.

### Author(s)

Rafael Maia <rm72@zips.uakron.edu>

---

 bgandilum

*Default background and illuminant data*


---

**Description**

Default background and illuminant data

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**References**

Endler, J. (1993). The Color of Light in Forests and Its Implications. *Ecological Monographs*, 63, 1-27.

---

 bootcoldist

*Bootstrap colour distance confidence intervals*


---

**Description**

Uses a bootstrap procedure to generate confidence intervals for the mean colour distance between two or more samples of colours

**Usage**

```
bootcoldist(vismodeldata, by, boot.n = 1000, alpha = 0.95,
  cores = getOption("mc.cores", 2L), ...)
```

**Arguments**

vismodeldata	(required) quantum catch color data. Can be the result from vismodel, or colspace. Data may also be independently calculated quantum catches, in the form of a data frame with columns representing photoreceptors.
by	(required) a numeric or character vector indicating the group to which each row from the object belongs to.
boot.n	number of bootstrap replicates (defaults to 1000)
alpha	the confidence level for the confidence intervals (defaults to 0.95)
cores	number of cores to be used in parallel processing. If 1, parallel computing will not be used. Defaults to <code>getOption("mc.cores", 2L)</code>
...	other arguments to be passed to <code>coldist</code> . Must at minimum include <code>n</code> and <code>weber</code> . See <code>coldist</code> for details.

**Value**

a matrix including the empirical mean and bootstrapped confidence limits for dS (and dL if achro = TRUE).

**References**

Maia, R., White, T. E., (2018) Comparing colors using visual models. Behavioral Ecology, ary017  
doi: [10.1093/beheco/ary017](https://doi.org/10.1093/beheco/ary017)

**Examples**

```
## Not run:
data(sicalis)
vm <- vismodel(sicalis, achro = "bt.dc")
gr <- gsub("ind..", "", rownames(vm))
bootcoldist(vm, by = gr, n = c(1, 2, 2, 4), weber = 0.1, weber.achro = 0.1, cores = 1)

## End(Not run)
```

---

 classify

*Identify colour classes in an image for adjacency analyses*


---

**Description**

Use k-means clustering to classify image pixels into discrete colour classes.

**Usage**

```
classify(imgdat, kcols = NULL, refID = NULL, interactive = FALSE,
  plotnew = FALSE, col = "red", cores = getOption("mc.cores", 2L),
  ...)
```

**Arguments**

imgdat	(required) image data. Either a single image, or a series of images stored in a list. Preferably the result of <a href="#">getimg</a> .
kcols	the number of discrete colour classes present in the input image(s). Can be a single integer when only a single image is present, or if kcols is identical for all images. When passing a list of images, kcols can also be a vector the same length as imgdat, or a data.frame with two columns specifying image file names and corresponding kcols. This argument can optionally be disregarded when interactive = TRUE, and kcols will be inferred from the number of selections.
refID	either the numeric index or name of a 'reference' image, for use when passing a list of images. Other images will be k-means classified using centres identified in the single reference image, thus helping to ensure that homologous pattern elements will be reliably classified between images, if so desired.

<code>interactive</code>	interactively specify the colour-category 'centers', for k-means clustering. When <code>TRUE</code> , the user is asked to click a number of points (equal to <code>kcols</code> , if specified, otherwise user-determined) that represent the distinct colours of interest. If a reference image is specified, it will be the only image presented.
<code>plotnew</code>	Should plots be opened in a new window when <code>interactive = TRUE</code> ? Defaults to <code>FALSE</code> .
<code>col</code>	the color of the marker points, when <code>interactive = TRUE</code> .
<code>cores</code>	number of cores to be used in parallel processing. If 1, parallel computing will not be used. Defaults to <code>getOption("mc.cores", 2L)</code> . Not available on Windows.
<code>...</code>	additional graphical parameters when <code>interactive = TRUE</code> . Also see <a href="#">par</a> .

### Value

A matrix, or list of matrices, of class `ring` containing the colour class classifications ID at each pixel location. The RGB values corresponding to k-means centres (i.e. colour classes) are stored as object attributes.

### Note

Since the `kmeans` process draws on random numbers to find initial cluster centres when `interactive = FALSE`, use `set.seed` if reproducible cluster ID's are desired between runs.

### Author(s)

Thomas E. White <thomas.white026@gmail.com>

### Examples

```
# Single image
papilio <- getimg(system.file("testdata/images/papilio.png", package = "pavo"))
papilio_class <- classify(papilio, kcols = 4)

## Not run:
# Multiple images, with interactive classification and a reference image
snakes <- getimg(system.file("testdata/images/snakes", package = "pavo"))
snakes_class <- classify(snakes, refID = "snake_01", interactive = TRUE)

## End(Not run)
```

---

coldist	<i>Colour distances</i>
---------	-------------------------

---

### Description

Calculates colour distances. When data are the result of `vismodel`, it applies the receptor-noise model of Vorobyev et al. (1998) to calculate colour distances with noise based on relative photoreceptor densities. It also accepts `colspace` data from the hexagon, colour-opponent-coding, categorical, segment, and cielab models, in which case euclidean distances (hexagon, cielab, categorical, segment) or manhattan distances (coc) are returned.

### Usage

```
coldist(modeldata, noise = c("neural", "quantum"), subset = NULL,
        achromatic = FALSE, qcatch = NULL, n = c(1, 2, 2, 4),
        weber = 0.1, weber.ref = "longest", weber.achro = 0.1)
```

### Arguments

modeldata	(required) quantum catch colour data. Can be the result from <code>vismodel</code> , or <code>colspace</code> . Data may also be independently calculated quantum catches, in the form of a data frame with columns representing photoreceptors.
noise	how the noise will be calculated. (Ignored for <code>colspace</code> objects if model is not a receptor noise model (i.e. hexagon, colour-opponent-coding, categorical, segment, and cie models)): <ul style="list-style-type: none"> <li>• neural (default): noise is proportional to the Weber fraction and is independent of the intensity of the signal received (i.e. assumes bright conditions).</li> <li>• quantum: noise is the sum of the neural noise and receptor noise, and is thus proportional to the Weber fraction and inversely proportional to the intensity of the signal received (the quantum catches). Note that the quantum option will only work with objects of class <code>vismodel</code>.</li> </ul>
subset	If only some of the comparisons should be returned, a character vector of length 1 or 2 can be provided, indicating which samples are desired. The subset vector must match the labels of the input samples, but partial matching (and regular expressions) are supported.
achromatic	Logical. If TRUE, last column of the data frame is used to calculate the achromatic contrast, with noise based on the Weber fraction given by the argument <code>weber.achro</code> . If the data are from the hexagon model (i.e. <code>colspace(space = 'hexagon')</code> ), it instead returns simple long (or 'green') receptor contrast.
qcatch	if the object is of class <code>vismodel</code> or <code>colspace</code> , this argument is ignored. If the object is a data frame of quantal catches from another source, this argument is used to specify what type of quantum catch is being used, so that the noise can be calculated accordingly: <ul style="list-style-type: none"> <li>• Qi: Quantum catch for each photoreceptor</li> </ul>

- `fi`: Quantum catch according to Fechner’s law (the signal of the receptor channel is proportional to the logarithm of the quantum catch)

<code>n</code>	photoreceptor densities for the cones used in visual modeling. must have same length as number of columns (excluding achromatic receptor if used; defaults to the Pekin robin <i>Leiothrix lutea</i> densities: <code>c(1, 2, 2, 4)</code> ). Ignored for <code>colspace</code> objects if model is not a receptor noise model (i.e. hexagon, colour-opponent-coding, categorical, and cie models).
<code>weber</code>	The Weber fraction to be used (often also referred to as receptor noise, or $e$ ). The noise-to-signal ratio $v$ is unknown, and therefore must be calculated based on the empirically estimated Weber fraction of one of the cone classes. $v$ is then applied to estimate the Weber fraction of the other cones. by default, the value of 0.1 is used (the empirically estimated value for the LWS cone from <i>Leiothrix lutea</i> ). See Olsson et al. 2017 for a review of published values in the literature. Ignored for <code>colspace</code> objects if model is not a receptor noise model (i.e. hexagon, colour-opponent-coding, categorical, segment, and cie models).
<code>weber.ref</code>	the cone class used to obtain the empirical estimate of the Weber fraction used for the <code>weber</code> argument. By default, <code>n4</code> is used, representing the LWS cone for <i>Leiothrix lutea</i> . Ignored for <code>colspace</code> objects if model is not a receptor noise model (i.e. hexagon, colour-opponent-coding, categorical, segment, and cie models).
<code>weber.achro</code>	the Weber fraction to be used to calculate achromatic contrast, when <code>achromatic = TRUE</code> . Defaults to 0.1. Ignored for <code>colspace</code> objects if model is not a receptor noise model (i.e. hexagon, colour-opponent-coding, categorical, segment, and cie models).

### Value

A data frame containing up to 4 columns. The first two (`patch1`, `patch2`) refer to the two colors being contrasted; `dS` is the chromatic contrast (delta S) and `dL` is the achromatic contrast (delta L). Units are JND’s in the receptor-noise model, euclidean distances in the categorical and segment space, manhattan distances in the color-opponent-coding space, green-receptor contrast in the hexagon, and lightness (L) contrast in the cielab model.

### Note on previous versions

previous versions of `coldist` calculated receptor noise using the arguments  $v$  for the individual cone noise-to-signal ratio and `n1`, `n2`, `n3`, `n4` for the relative cone densities. These arguments have been replaced by `weber` and `n`, which takes a vector of relative cone densities. `weber.ref` allows the user to specify which receptor to use as the reference to obtain the desired Weber fraction, and `coldist` calculates internally the value of  $v$  to be used when calculating the Weber fraction for the remaining cones.

This allows a more explicit choice of Weber fraction, without the need to find the right value of  $v$  to use in order to obtain the desired signal-to-noise ratio. Furthermore, by allowing `n` to be entered as a vector, `coldist` can now handle visual systems with more than four photoreceptors.

In addition, the achromatic noise is calculated based on the `weber.achro` argument directly, and not based on  $v$  and `n4` as before.



**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**References**

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- Olsson, P., Lind, O., & Kelber, A. (2015) Bird colour vision: behavioural thresholds reveal receptor noise. *Journal of Experimental Biology*, 218, 184-193.
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- Olsson, P., Lind, O., & Kelber, A. (2017) Chromatic and achromatic vision: parameter choice and limitations for reliable model predictions. *Behavioral Ecology*, doi: [10.1093/beheco/arx133](https://doi.org/10.1093/beheco/arx133)

**Examples**

```
## Not run:
# Dichromat
data(flowers)
vis.flowers <- vismodel(flowers, visual = "canis", relative = FALSE)
didist.flowers <- coldist(vis.flowers, n = c(1, 2))

# Trichromat
vis.flowers <- vismodel(flowers, visual = "apis", relative = FALSE)
tridist.flowers <- coldist(vis.flowers, n = c(1, 2, 1))

# Trichromat, color-hexagon model (euclidean distances)
vis.flowers <- vismodel(flowers,
  visual = "apis", qcatch = "Ei",
  relative = FALSE, vonkries = TRUE, achro = "l", bkg = "green"
)
hex.flowers <- colspace(vis.flowers, space = "hexagon")
hexdist.flowers <- coldist(hex.flowers)

# Trichromat, color-opponent-coding model (manhattan distances)
vis.flowers <- vismodel(flowers, visual = "apis", qcatch = "Ei", relative = FALSE, vonkries = TRUE)
coc.flowers <- colspace(vis.flowers, space = "coc")
hexdist.flowers <- coldist(coc.flowers)

# Tetrachromat
data(sicalis)
vis.sicalis <- vismodel(sicalis, visual = "avg.uv", relative = FALSE)
tetradist.sicalis.n <- coldist(vis.sicalis)
```

```
# This will also work, but give you several warnings you shouldn't ignore!!
col.sicalis <- colspace(vis.sicalis)
tetradist.sicalis.n <- coldist(col.sicalis)

tetradist.sicalis.q <- coldist(vis.sicalis, noise = "quantum")

## End(Not run)
```

---

colspace

*Model spectra in a colorspace*


---

## Description

Models reflectance spectra in a colorspace. For information on plotting arguments and graphical parameters, see [plot.colspace](#).

## Usage

```
colspace(vismodeldata, space = c("auto", "di", "tri", "tcs", "hexagon",
  "coc", "categorical", "ciexyz", "cielab", "cielch", "segment"),
  qcatch = NULL)
```

## Arguments

vismodeldata	(required) quantum catch color data. Can be either the result from <a href="#">vismodel</a> or independently calculated data (in the form of a data frame with columns representing quantum catches).
space	Which colorspace/model to use. Options are: <ul style="list-style-type: none"> <li>• auto: if data is a result from <a href="#">vismodel</a>, applies di, tri or tcs if input visual model had two, three or four cones, respectively.</li> <li>• di: dichromatic colourspace. See <a href="#">dispace</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• tri: trichromatic colourspace (i.e. Maxwell triangle). See <a href="#">trispace</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• tcs: tetrahedral colourspace. See <a href="#">tcspace</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• hexagon: the trichromatic colour-hexagon of Chittka (1992). See <a href="#">hexagon</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• coc: the trichromatic colour-opponent-coding model of Backhaus (1991). See <a href="#">coc</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• categorical: the tetrachromatic categorical fly-model of Troje (1993). See <a href="#">categorical</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• ciexyz: CIEXYZ space. See <a href="#">cie</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• cielab: CIELAB space. See <a href="#">cie</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• cielch: CIELCh space. See <a href="#">cie</a> for details. (<a href="#">plotting arguments</a>)</li> <li>• segment: segment analysis of Endler (1990). See <a href="#">segospace</a> for details. (<a href="#">plotting arguments</a>)</li> </ul>
qcatch	Which quantal catch metric is being inputted. Only used when input data is NOT an output from <a href="#">vismodel</a> . Must be Qi, fi or Ei.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

Thomas White <thomas.white026@gmail.com>

**References**

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Chittka L. (1992). The colour hexagon: a chromaticity diagram based on photoreceptor excitations as a generalized representation of colour opponency. *Journal of Comparative Physiology A*, 170(5), 533-543.

Chittka L, Shmida A, Troje N, Menzel R. (1994). Ultraviolet as a component of flower reflections, and the colour perception of Hymenoptera. *Vision research*, 34(11), 1489-1508.

Troje N. (1993). Spectral categories in the learning behaviour of blowflies. *Zeitschrift fur Naturforschung C*, 48, 96-96.

Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.

Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.

Kelber A, Vorobyev M, Osorio D. (2003). Animal colour vision - behavioural tests and physiological concepts. *Biological Reviews*, 78, 81 - 118.

Backhaus W. (1991). Color opponent coding in the visual system of the honeybee. *Vision Research*, 31, 1381-1397.

Endler, J. A. (1990) On the measurement and classification of color in studies of animal color patterns. *Biological Journal of the Linnean Society*, 41, 315-352.

**Examples**

```
data(flowers)

# A dichromat in a segment colourspace
vis.flowers <- vismodel(flowers, visual = "canis")
di.flowers <- colspace(vis.flowers, space = "di")

# Honeybee in the colour hexagon
vis.flowers <- vismodel(flowers,
  visual = "apis", qcatch = "Ei", relative = FALSE,
  vonkries = TRUE, achro = "l", bkg = "green"
)
hex.flowers <- colspace(vis.flowers, space = "hexagon")

# A trichromat in a Maxwell triangle
vis.flowers <- vismodel(flowers, visual = "apis")
tri.flowers <- colspace(vis.flowers, space = "tri")
plot(tri.flowers)
```

```
# A tetrachromat in a tetrahedral colour space
vis.flowers <- vismodel(flowers, visual = "bluetit")
tcs.flowers <- colspace(vis.flowers, space = "tcs")

# A housefly in the categorical colour space
vis.flowers <- vismodel(flowers, visual = "musca", achro = "md.r1")
cat.flowers <- colspace(vis.flowers, space = "categorical")
```

---

explorespec

*Plot spectral curves*


---

### Description

Plots one or multiple spectral curves in the same graph to rapidly compare groups of spectra.

### Usage

```
explorespec(rspeccdata, by = NULL, scale = c("equal", "free"),
  legpos = "topright", ...)
```

### Arguments

rspeccdata	(required) a data frame, possibly an object of class <code>rspec</code> that has wavelength range in the first column, named 'wl', and spectral measurements in the remaining columns.
by	number of spectra to include in each graph (defaults to 1)
scale	defines how the y-axis should be scaled. 'free': panels can vary in the range of the y-axis; 'equal': all panels have the y-axis with the same range.
legpos	legend position control. Either a vector containing x and y coordinates or a single keyword from the list: "bottomright", "bottom", "bottomleft", "left", "topleft", "top", "topright", "right" and "center".
...	additional parameters to be passed to plot

### Value

Spectral curve plots

### Note

Number of plots presented per page depends on the number of graphs produced.

### Author(s)

Pierre-Paul Bitton <bittonp@uwindsor.ca>

**Examples**

```
data(sicalis)
explorespec(sicalis, 3)
explorespec(sicalis, 3, ylim = c(0, 100), legpos = c(500, 80))
```

---

flowers	<i>Reflectance spectra from a suite of native Australian flowers, collected around Cairns, Queensland.</i>
---------	--

---

**Description**

dataset containing reflectance measurements from 36 native Australian angiosperm species, indicated by column names.

**Author(s)**

Thomas White <thomas.white026@gmail.com>

**References**

Dalrymple, R. L., Kemp, D. J., Flores-Moreno, H., Laffan, S. W., White, T. E., Hemmings, F. A., Tindall, M. L., & Moles, A. T. (2015). Birds, butterflies and flowers in the tropics are not more colourful than those at higher latitudes. *Global Ecology and Biogeography*, 24(12), 1424-1432. doi: [10.1111/geb.12368](https://doi.org/10.1111/geb.12368)

White, T. E., Dalrymple, R. L., Herberstein, M. E., & Kemp, D. J. (2017). The perceptual similarity of orb-spider prey lures and flowers colours. *Evolutionary Ecology*, 31(1), 1-20. doi: [10.1007/s106820169876x](https://doi.org/10.1007/s106820169876x)

Dalrymple, R. L., Flores-Moreno, H., Kemp, D. J., White, T. E., Laffan, S. W., Hemmings, F. A., Tindall, M. L., & Moles, A. T. (2018). Abiotic and biotic predictors of macroecological patterns in bird and butterfly coloration. *Ecological Monographs*, 88(2), 204-224. doi: [10.1002/ecm.1287](https://doi.org/10.1002/ecm.1287)

---

getimg	<i>Import image data</i>
--------	--------------------------

---

**Description**

Finds and imports PNG, JPEG, and/or BMP images.

**Usage**

```
getimg(imgpath = getwd(), subdir = FALSE, subdir.names = FALSE,
max.size = 1, cores = getOption("mc.cores", 2L))
```

**Arguments**

<code>imgpath</code>	(required) either the full path to a given image (including extension), or the path to a folder in which multiple image files are located. Mixed file formats within a folder are accepted.
<code>subdir</code>	should subdirectories within the <code>imgpath</code> folder be included in the search? (defaults to FALSE).
<code>subdir.names</code>	should subdirectory path be included in the name of the images? (defaults to FALSE).
<code>max.size</code>	maximum size of all images to be allowed in memory, in GB. Defaults to 1.
<code>cores</code>	number of cores to be used in parallel processing. If 1, or if total image sizes exceed 200 mb in memory, parallel computing will not be used. Defaults to <code>getOption("mc.cores", 2L)</code> . Not available on Windows.

**Value**

a image, or list of images, of class `ring`, for use in further `pavo` functions.

**Author(s)**

Thomas E. White <thomas.white026@gmail.com>

**Examples**

```
# Single image
papilio <- getimg(system.file("testdata/images/papilio.png", package = "pavo"))

# Multiple images
snakes <- getimg(system.file("testdata/images/snakes", package = "pavo"))
```

---

getspec

*Import spectra files*

---

**Description**

Finds and imports spectra files from a folder. Currently works for reflectance files generated in Ocean Optics SpectraSuite (USB2000, USB4000 and Jaz spectrometers), CRAIC software (after exporting) and Avantes (before or after exporting).

**Usage**

```
getspec(where = getwd(), ext = "txt", lim = c(300, 700),
  decimal = ".", sep = NULL, subdir = FALSE, subdir.names = FALSE,
  cores = getOption("mc.cores", 2L), ignore.case = TRUE)
```

**Arguments**

where	(required) folder in which files are located.
ext	file extension to be searched for, without the "." (defaults to "txt").
lim	a vector with two numbers determining the wavelength limits to be considered (defaults to 300 and 700).
decimal	character to be used to identify decimal plates (defaults to ".").
sep	column delimiting characters to be considered in addition to the default (which are: tab, space, and ";")
subdir	should subdirectories within the where folder be included in the search? (defaults to FALSE).
subdir.names	should subdirectory path be included in the name of the spectra? (defaults to FALSE).
cores	Number of cores to be used. If greater than 1, import will use parallel processing (not available in Windows).
ignore.case	Logical. Should the extension search be case insensitive? (defaults to TRUE)

**Value**

A data frame, of class `rspec`, containing individual imported spectral files as columns. Reflectance values are interpolated to the nearest wavelength integer.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>  
Hugo Gruson <hugo.gruson+R@normalesup.org>

**References**

Montgomerie R (2006) Analyzing colors. In: Hill G, McGraw K (eds) Bird coloration. Harvard University Press, Cambridge, pp 90-147.

**Examples**

```
getspec(system.file("testdata", package = "pavo"), lim = c(400, 900))
getspec(system.file("testdata", package = "pavo"), ext = "tnt")
```

---

img_conversion	<i>Convert images between class ring and cimg</i>
----------------	---

---

**Description**

Conveniently convert single objects of class `ring` and `cimg` (from the package `imager`, which contains a suite of useful image-processing capabilities).

**Usage**

```
ring2cimg(image)
cimg2ring(image, name = "img")
```

**Arguments**

image	an object of class ring or cimg.
name	the name(s) of the image(s).

**Value**

an image of the specified class

**Note**

Attributes (e.g. scales, color-classes) will not be preserved following conversion from class ring, so it's best to use early in the analysis workflow.

**Author(s)**

Thomas E. White <thomas.white026@gmail.com>  
Thomas E. White <thomas.white026@gmail.com>  
Thomas E. White <thomas.white026@gmail.com>

**Examples**

```
## Not run:
papilio <- getimg(system.file("testdata/images/papilio.png", package = "pavo"))

# From class ring to cimg
papilio_cimg <- ring2cimg(papilio)
class(papilio_cimg)

# From class cimg to ring
papilio_ring <- cimg2ring(papilio_cimg)
class(papilio_ring)

## End(Not run)
```



---

irrad2flux	<i>Converts between irradiance and photon (quantum) flux</i>
------------	--

---

**Description**

Some spectrometers will give illuminant values in units of irradiance ( $\mu\text{Watt} \cdot \text{cm}^{-2}$ ), but physiological models require illuminants in units of photon (quantum) flux ( $\mu\text{mol} \cdot \text{s}^{-1} \cdot \text{m}^{-2}$ ). The functions `irrad2flux` and `flux2irrad` allows for easy conversion of `rspec` objects between these units.

**Usage**

```
irrad2flux(rspecdata)
```

```
flux2irrad(rspecdata)
```

**Arguments**

`rspecdata` (required) a `rspec` object containing illuminant values.

**Value**

a converted `rspec` object.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

---

<code>is.colospace</code>	<i>Test is object is of class 'colospace'</i>
---------------------------	---

---

**Description**

Test is object is of class 'colospace'

**Usage**

```
is.colospace(object)
```

**Arguments**

`object` an R object

**Value**

a logical value indicating whether the object is of class `colospace`

**See Also**[colspace](#)


---

<code>is.vismodel</code>	<i>Test is object is of class 'vismodel'</i>
--------------------------	--

---

**Description**

Test is object is of class 'vismodel'

**Usage**

```
is.vismodel(object)
```

**Arguments**

`object`            an R object

**Value**

a logical value indicating whether the object is of class `vismodel`

**See Also**[vismodel](#)


---

<code>jnd2xyz</code>	<i>Convert JND distances into perceptually-corrected Cartesian coordinates</i>
----------------------	--

---

**Description**

Converts a `coldist` output into Cartesian coordinates that are perceptually-corrected (i.e. Euclidean distances = JND distances)

**Usage**

```
jnd2xyz(coldistres, center = TRUE, rotate = TRUE,
        rotcenter = c("mean", "achro"), ref1 = "l", ref2 = "u",
        axis1 = c(1, 1, 0), axis2 = c(0, 0, 1))
```

**Arguments**

<code>coldistres</code>	(required) the output from a <code>coldist</code> call.
<code>center</code>	logical indicating if the data should be centered on its centroid (defaults to TRUE).
<code>rotate</code>	logical indicating if the data should be rotated (defaults to TRUE).
<code>rotcenter</code>	should the vectors for rotation be centered in the achromatic center ("achro") or the data centroid ("mean", the default)?
<code>ref1</code>	the cone to be used as a the first reference. May be NULL (for no first rotation in the 3-dimensional case) or must match name in the original data that was used for <code>coldist</code> . Defaults to 'l'.
<code>ref2</code>	the cone to be used as a the second reference. May be NULL (for no first rotation in the 3-dimensional case) or must match name in the original data that was used for <code>coldist</code> . Defaults to 'u'. (only used if data has 3 dimensions).
<code>axis1</code>	A vector of length 3 composed of 0's and 1's, with 1's representing the axes (x,y,z) to rotate around. Defaults to <code>c(1,1,0)</code> , such that the rotation aligns with the xy plane (only used if data has 2 or 3 dimensions). Ignored if <code>ref1</code> is NULL (in 3-dimensional case only)
<code>axis2</code>	A vector of length 3 composed of 0's and 1's, with 1's representing the axes (x,y,z) to rotate around. Defaults to <code>c(0,0,1)</code> , such that the rotation aligns with the z axis (only used if data has 3 dimensions). Ignored if <code>ref2</code> is NULL (in 3-dimensional case only)

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**References**

Pike, T.W. (2012). Preserving perceptual distances in chromaticity diagrams. *Behavioral Ecology*, 23, 723-728.

Maia, R., White, T. E., (2018) Comparing colors using visual models. *Behavioral Ecology*, ary017  
doi: [10.1093/beheco/ary017](https://doi.org/10.1093/beheco/ary017)

**Examples**

```
data(flowers)
vis.flowers <- vismodel(flowers)
cd.flowers <- coldist(vis.flowers)
jnd2xyz(cd.flowers)
```

---

legendtetra	<i>Add legend to a static tetrahedral colourspace</i>
-------------	---

---

**Description**

Adds a legend to a static tetrahedral colourspace plot.

**Usage**

```
legendtetra(x = 0.8, y = 1.2, ...)
```

**Arguments**

x, y	position of the legend relative to plot limits (usually a value between 0 and 1, but because of the perspective distortion, values greater than 1 are possible)
...	additional arguments passed to <a href="#">legend</a> .

**Value**

legendtetra adds a legend to a static tetrahedral colourspace plot. for additional information on which arguments are necessary and how they are used, see [legend](#).

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

---

merge.rspec	<i>Merge two rspec objects</i>
-------------	--------------------------------

---

**Description**

Merges two rspec or data.frame objects into a single rspec object.

**Usage**

```
## S3 method for class 'rspec'  
merge(x, y, ...)
```

**Arguments**

x, y	(required) rspec objects to merge.
...	additional class arguments.

**Value**

an object of class `rspec` for use with `pavo` functions. Will use `by = "wl"` if unspecified, or automatically append `wl` to the `by` argument if one is specified.

**Author(s)**

Chad Eliason <cme16@zips.uakron.edu>

**See Also**

[as.rspec](#), [aggspec](#)

**Examples**

```
# Load and split dataset into 2 sections
data(teal)
teal1 <- teal[, c(1, 3:5)]
teal2 <- teal[, c(1, 2, 6:12)]
teal.mer <- merge(teal1, teal2, by = "wl")
head(teal.mer)
par(mfrow = c(1, 2))
plot(teal.mer)
plot(teal)
```

---

peakshape

*Peak shape descriptors*

---

**Description**

Calculates height, location and width of peak at the reflectance midpoint (FWHM). Note: bounds should be set wide enough to incorporate all minima in spectra. Smoothing spectra using [prospec](#) is also recommended.

**Usage**

```
peakshape(rspecdata, select = NULL, lim = NULL, plot = TRUE,
  ask = FALSE, absolute.min = FALSE, ...)
```

**Arguments**

<code>rspecdata</code>	(required) a data frame, possibly an object of class <code>rspec</code> , with a column with wavelength data, named <code>'wl'</code> , and the remaining column containing spectra to process.
<code>select</code>	specification of which spectra to plot. Can be a numeric vector or factor (e.g., <code>sex == 'male'</code> ).
<code>lim</code>	a vector specifying the wavelength range to analyze.

plot	logical. Should plots indicating calculated parameters be returned? (Defaults to TRUE).
ask	logical, specifies whether user input needed to plot multiple plots when number of spectra to analyze is greater than 1 (defaults to FALSE).
absolute.min	logical. If TRUE, full width at half maximum will be calculated using the absolute minimum reflectance of the spectrum, even if that value falls outside the range specified by lim. (defaults to FALSE)
...	additional arguments to be passed to plot.

**Value**

a data frame containing column names (id); peak height (max value, B3), location (hue, H1) and full width at half maximum (FWHM), as well as half widths on left (HWHM.l) and right side of peak (HWHM.r). Incl.min column indicates whether user-defined bounds incorporate the actual minima of the spectra. Function will return a warning if not.

**Author(s)**

Chad Eliason <cme16@zips.uakron.edu>  
 Rafael Maia <rm72@zips.uakron.edu>  
 Hugo Gruson <hugo.gruson+R@normalesup.org>

**See Also**

[procspec](#)

**Examples**

```
data(teal)

peakshape(teal, select = 3)
peakshape(teal, select = 10)

# Use wavelength bounds to narrow in on peak of interest
peakshape(teal, select = 10, lim = c(400, 550))
```

---

plot.colspace                      *Plot spectra in a colourspace*

---

**Description**

Plots reflectance spectra in the appropriate colorspace.

**Usage**

```
## S3 method for class 'colspace'
plot(x, ...)
```

**Arguments**

- x (required) an object of class colspace.
- ... additional graphical options, which vary by modeled space. Refer to their individual documentation:
- `diplot`: dichromat space
  - `tripplot`: trichromat space
  - `tetraplot`: tetrahedral space
  - `catplot`: categorical space
  - `hexplot`: colour hexagon
  - `cocplot`: colour-opponent-coding space
  - `cieplot`: cie spaces
  - `segplot`: segment analysis space
  - `jndplot`: perceptual, 'noise corrected' space (for the results of `jnd2xyz`)
- Also see `par`.

**Value**

A colorspace plot appropriate to the input data.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>  
 Thomas White <thomas.white026@gmail.com>  
 Chad Eliason <cme16@zips.uakron.edu>

**References**

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- Troje N. (1993). Spectral categories in the learning behaviour of blowflies. *Zeitschrift fur Naturforschung C*, 48, 96-96.
- Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.
- Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.
- Kelber A, Vorobyev M, Osorio D. (2003). Animal colour vision - behavioural tests and physiological concepts. *Biological Reviews*, 78, 81 - 118.

Backhaus W. (1991). Color opponent coding in the visual system of the honeybee. *Vision Research*, 31, 1381-1397.

Endler, J. A. (1990) On the measurement and classification of color in studies of animal color patterns. *Biological Journal of the Linnean Society*, 41, 315-352.

### See Also

[plot](#)

### Examples

```
## Not run:
data(flowers)
data(sicalis)

# Dichromat
vis.flowers <- vismodel(flowers, visual = "canis")
di.flowers <- colspace(vis.flowers, space = "di")
plot(di.flowers)

# Colour hexagon
vis.flowers <- vismodel(flowers,
  visual = "apis", qcatch = "Ei", relative = FALSE,
  vonkries = TRUE, achro = "l", bkg = "green"
)
hex.flowers <- colspace(vis.flowers, space = "hexagon")
plot(hex.flowers, sectors = "coarse")

# Tetrahedron (static)
vis.sicalis <- vismodel(sicalis, visual = "avg.uv")
tcs.sicalis <- colspace(vis.sicalis, space = "tcs")
plot(tcs.sicalis)

# Tetrahedron (interactive)
vis.sicalis <- vismodel(sicalis, visual = "avg.uv")
tcs.sicalis <- colspace(vis.sicalis, space = "tcs")
tcsplot(tcs.sicalis, size = 0.005)

## Add points to interactive tetrahedron
patch <- rep(c("C", "T", "B"), 7)
tcs.crown <- subset(tcs.sicalis, "C")
tcs.breast <- subset(tcs.sicalis, "B")
tcsplot(tcs.crown, col = "blue")
tcspoints(tcs.breast, col = "red")

## Plot convex hull in interactive tetrahedron
tcsplot(tcs.sicalis, col = "blue", size = 0.005)
tcsvol(tcs.sicalis)

## End(Not run)
```



---

`plot.rimg`*Plot unprocessed or colour-classified images*

---

**Description**

Plot unprocessed or colour-classified image data. If the images are in a list, they will be stepped through one by one.

**Usage**

```
## S3 method for class 'rimg'  
plot(x, axes = TRUE, col = NULL, ...)
```

**Arguments**

<code>x</code>	(required) an image of class <code>rimg</code> , or list thereof.
<code>axes</code>	should axes be drawn? (defaults to <code>TRUE</code> )
<code>col</code>	optional vector of colours when plotting colour-classified images. Defaults to the mean RGB values of the k-means centres (i.e. the average 'original' colours).
<code>...</code>	additional graphical parameters. Also see <a href="#">par</a> .

**Value**

an image plot.

**Author(s)**

Thomas E. White <thomas.white026@gmail.com>

**Examples**

```
## Not run:  
papilio <- getimg(system.file("testdata/images/papilio.png", package = "pavo"))  
plot(papilio)  
papilio_class <- classify(papilio, kcols = 4)  
plot(papilio_class)  
  
# Multiple images  
snakes <- getimg(system.file("testdata/images/snakes", package = "pavo"))  
plot(snakes)  
snakes_class <- classify(snakes, kcols = 3)  
plot(snakes_class)  
  
## End(Not run)
```

---

plot.rspec

*Plot spectra*


---

### Description

Plots reflectance spectra in different arrangements.

### Usage

```
## S3 method for class 'rspec'
plot(x, select = NULL, type = c("overlay", "stack",
  "heatmap"), varying = NULL, n = 100, labels.stack = NULL, ...)
```

### Arguments

x	(required) a data frame, possibly an object of class <code>rspec</code> , with a column with wavelength data, named 'wl', and the remaining column containing spectra to plot.
select	specification of which spectra to plot. Can be a numeric vector or factor (e.g., <code>sex=='male'</code> )
type	what type of plot should be drawn. Possibilities are: <ul style="list-style-type: none"> <li>• <code>overlay</code> (default) for plotting multiple spectra in a single panel with a common y-axis.</li> <li>• <code>stack</code> for plotting multiple spectra in a vertical arrangement.</li> <li>• <code>heatmap</code> for plotting reflectance values by wavelength and a third variable (<code>varying</code>).</li> </ul>
varying	a numeric vector giving values for y-axis in <code>heatmap</code> .
n	number of bins with which to interpolate colors and <code>varying</code> for the <code>heatmap</code> .
labels.stack	a vector of labels for the stacked spectra when using <code>type = stack</code> . Defaults to the numeric column ID's.
...	additional arguments passed to <code>plot</code> (or <code>image</code> for 'heatmap').

### Author(s)

Chad Eliason <cme16@zips.uakron.edu>

Thomas White <thomas.white026@gmail.com>

### See Also

[spec2rgb](#), [image](#), [plot](#)

## Examples

```
## Not run:

data(teal)
plot(teal, type = "overlay")
plot(teal, type = "stack")
plot(teal, type = "heatmap")

## End(Not run)
```

---

plotsmooth

*Plot loess smoothed curves*

---

## Description

Plots spectral curves with various levels of loess smoothing to help decide which loess parameters are best for subsequently smoothing the data (e.g. via procspec).

## Usage

```
plotsmooth(rspeccdata, minsmooth = 0.05, maxsmooth = 0.2, curves = 5,
  specnum = 0, ask = TRUE)
```

## Arguments

rspeccdata	(required) a data frame, possibly of class rspec, which contains a column containing a wavelength range, named 'wl', and spectra data in remaining columns.
minsmooth	the minimum f value of the loess function to visualize (defaults to 0.05).
maxsmooth	the maximum f value of the loess function to visualize (defaults to 0.20).
curves	the number of curves to display on the same plot (defaults to 5).
specnum	the number of spectral curves, from the data frame, to visualize (defaults to ALL).
ask	logical. if TRUE, asks for user input before changing plot pages

## Value

Series of plot with curves processed with varying level of loess smoothing

## Author(s)

Pierre-Paul Bitton <bittonp@uwindsor.ca>

**Examples**

```
## Not run:
data(sicalis)
plotsmooth(sicalis, minsmooth = 0.05, maxsmooth = 0.1, curves = 7, specnum = 6)

## End(Not run)
```

---

points.colospace	<i>Plot points in a colourspace</i>
------------------	-------------------------------------

---

**Description**

Add points to a colourspace plot

**Usage**

```
## S3 method for class 'colospace'
points(x, ...)
```

**Arguments**

x	(required) an object of class colospace.
...	additional graphical options. See <a href="#">par</a> .

**Value**

points.colospace adds points to a colorspace plot. When space = 'tcs', it creates 3D points in a tetrahedral color space plot using functions of the package rgl, based on openGL capabilities.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

Thomas White <thomas.white026@gmail.com>

---

procimg	<i>Process images</i>
---------	-----------------------

---

**Description**

Specify scales, resize, and/or define focal objects within images.

**Usage**

```
procimg(image, resize = NULL, rotate = NULL, scaledist = NULL,
        outline = FALSE, reclass = NULL, smooth = FALSE, iterations = 1L,
        col = "red", plotnew = FALSE, ...)
```

**Arguments**

image	(required) image data. Either a single image array, or a number of images stored in a list. Preferably the result of <a href="#">getimg</a> .
resize	an integer specifying the scaling factor for linearly resizing images, if so desired. E.g. 0.5 to half the size of an image, or 2 to double it.
rotate	an integer specifying the angle of image rotation, in degrees. Images are rotated around the centre, and linearly interpolated.
scaledist	an integer, or numeric vector equal in length to the number of images, specifying the length of the scale in the image(s). Image(s) will then be presented, and the user asked to select either end of the scale corresponding to the input value.
outline	interactively specify the focal object in an image by clicking around its outline. The xy-coordinates of the resulting closed polygon are saved as an attribute, for use in generating a masking layer & separating animals/plants from backgrounds in further analyses. This is particularly useful when backgrounds are complex, such as in natural settings.
reclass	interactively specify an area on a colour-classified image that is to be reclassified as the numeric value provided. e.g. when <code>reclass = 1</code> , the user will be asked to select a polygon on the image, within which all colour-category values will be changes to 1.
smooth	should the polygon specified when <code>outline = TRUE</code> be smoothed using Chaikin's corner-cutting algorithm? Defaults to FALSE.
iterations	the number of smoothing iterations, when <code>smooth = TRUE</code> . Defaults to 1.
col	the color of the marker points and/or line, when using interactive options.
plotnew	should plots be opened in a new window? Defaults to FALSE.
...	additional graphical parameters. Also see <a href="#">par</a> .

**Value**

an image, or list of images, for use in further pavo functions.

**Author(s)**

Thomas E. White <thomas.white026@gmail.com>

**References**

Chaikin, G. 1974. An algorithm for high speed curve generation. *Computer Graphics and Image Processing* 3, 346-349.

**Examples**

```
## Not run:
# Single image
papilio <- getimg(system.file("testdata/images/papilio.png", package = "pavo"))
papilio <- procimg(papilio, scaledist = 10)

# Assign individual scales to each image, after slightly reducing their size.
snakes <- getimg(system.file("testdata/images/snakes", package = "pavo"))
snakes <- procimg(snakes, scaledist = c(10, 14), resize = 0.95)

## End(Not run)
```

---

procspec

*Process spectra*

---

**Description**

Applies normalization and/or smoothing to spectra for further analysis or plotting.

**Usage**

```
procspec(rspeccdata, opt = c("none", "smooth", "maximum", "minimum",
  "bin", "sum", "center"), fixneg = c("none", "addmin", "zero"),
  span = 0.25, bins = 20)
```

**Arguments**

rspeccdata	(required) a data frame, possibly an object of class <code>rspec</code> , with a column with wavelength data, named 'wl', and the remaining column containing spectra to process.
opt	what type of processing options to apply. User can select multiple options by providing a vector. Possibilities are: <ul style="list-style-type: none"> <li>• "none" does not perform any processing (default).</li> <li>• "smooth" applies LOESS smoothing to each spectrum using <code>loess.smooth</code>. Optimal smoothing parameter can be assessed by using <code>plotsmooth</code>.</li> <li>• "minimum" subtracts the minimum from each individual spectra.</li> <li>• "maximum" divides each spectrum by its maximum value.</li> </ul>

- "sum" divides each spectrum by summed values.
- "bin" bins each spectrum into the specified number of bins. bins argument must be set.
- "center" centers individual spectra by subtracting mean reflectance from all values.

fixneg            how to handle negative values. Possibilities are:

- "none" does not perform negative value correction (default).
- "zero" sets all negative values to zero.
- "addmin" adds the absolute value of the maximally negative values of each spectra to the reflectance at all other wavelengths (setting the minimum value to zero, but scaling other values accordingly).

span             sets the smoothing parameter used by loess.smooth.

bins             sets the number of equally sized wavelength bins for opt = "bin".

**Value**

A data frame of class rspec with the processed data.

**Author(s)**

Chad Eliason <cme16@zip.s.uakron.edu>

**References**

Cuthill, I., Bennett, A. T. D., Partridge, J. & Maier, E. 1999. Plumage reflectance and the objective assessment of avian sexual dichromatism. *The American Naturalist*, 153, 183-200.

Montgomerie R. 2006. Analyzing colors. In Hill, G.E, and McGraw, K.J., eds. *Bird Coloration. Volume 1 Mechanisms and measurements*. Harvard University Press, Cambridge, Massachusetts.

White, T. E., Dalrymple, R. L., Noble D. W. A., O'Hanlon, J. C., Zurek, D. B., Umbers, K. D. L. 2015. Reproducible research in the study of biological coloration. *Animal Behaviour*, 106, 51-57.

**See Also**

[loess.smooth](#), [plotsmooth](#)

**Examples**

```
data(teal)
plot(teal, select = 10)

# Smooth data to remove noise
teal.sm <- specsproc(teal, opt = "smooth", span = 0.25)
plot(teal.sm, select = 10)

# Normalize to max of unity
teal.max <- specsproc(teal, opt = c("max"))
plot(teal.max, select = 10)
```

---

`projplot`*2D projection of a tetrahedral colourspace*

---

**Description**

Produces a 2D projection plot of points in a tetrahedral colour space

Adds points to a tetrahedral colorspace projection

**Usage**

```
projplot(tcsdata, ...)
```

```
projpoints(tcsdata, ...)
```

**Arguments**

`tcsdata` (required) tetrahedral color space coordinates, possibly a result from `colspace`, containing values for the 'h.theta' and 'h.phi' coordinates as columns (labeled as such).

`...` additional parameters to be passed to the plotting of data points.

**Value**

`projplot` creates a 2D plot of color points projected from the tetrahedron to its encapsulating sphere, and is ideal to visualize differences in hue.

`projpoints` creates points in a projection color space plot produced by `projplot`.

**Note**

`projplot` uses the Mollweide projection, and not the Robinson projection, which has been used in the past. Among other advantages, the Mollweide projection preserves area relationships within latitudes without distortion.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**References**

Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.

Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.



## Examples

```
data(sicalis)
vis.sicalis <- vismodel(sicalis, visual = "avg.uv")
tcs.sicalis <- colspace(vis.sicalis, space = "tcs")
projplot(tcs.sicalis, pch = 16, col = setNames(rep(seq_len(3), 7), rep(c("C", "T", "B"), 7)))
```

---

sensdata

*Retrieve or plot in-built spectral sensitivity data*

---

## Description

Retrieve (as an rspec object) or plot pavo's in-built spectral sensitivity data.

## Usage

```
sensdata(visual = c("none", "all", "avg.uv", "avg.v", "bluetit",
  "ctenophorus", "star", "pfowl", "apis", "canis", "cie2", "cie10",
  "musca", "habronattus", "rhinecanthus"), achromatic = c("none", "all",
  "bt.dc", "ch.dc", "st.dc", "md.r1", "ra.dc", "cf.r"), illum = c("none",
  "all", "bluesky", "D65", "forestshade"), trans = c("none", "all",
  "bluetit", "blackbird"), bkg = c("none", "all", "green"),
  plot = FALSE, ...)
```

## Arguments

visual

visual systems. Options are:

- 'none': no visual sensitivity data.
- 'all': all visual sensitivity data.
- 'apis': Honeybee *Apis mellifera* visual system.
- 'avg.uv': average avian UV system.
- 'avg.v': average avian V system.
- 'bluetit': Blue tit *Cyanistes caeruleus* visual system.
- 'canis': Canid *Canis familiaris* visual system.
- 'cie2': 2-degree colour matching functions for CIE models of human colour vision. Functions are linear transformations of the 2-degree cone fundamentals of Stockman & Sharpe (2000), as ratified by the CIE (2006).
- 'cie10': 10-degree colour matching functions for CIE models of human colour vision. Functions are linear transformations of the 10-degree cone fundamentals of Stockman & Sharpe (2000), as ratified by the CIE (2006).
- 'ctenophorus': Ornate dragon lizard *Ctenophorus ornatus*.
- 'musca': Housefly *Musca domestica* visual system.
- 'pfowl': Peafowl *Pavo cristatus* visual system.
- 'star': Starling *Sturnus vulgaris* visual system.
- 'habronattus': Jumping spider *Habronattus pyrrithrix*.

	<ul style="list-style-type: none"> <li>• 'rhinecanthus': Triggerfish <i>Rhinecanthus aculeatus</i>.</li> </ul>
achromatic	<p>the sensitivity data used to calculate luminance (achromatic) receptor stimulation. Options are:</p> <ul style="list-style-type: none"> <li>• 'none': no achromatic sensitivity data.</li> <li>• 'all': all achromatic sensitivity data.</li> <li>• 'bt.dc': Blue tit <i>Cyanistes caeruleus</i> double cone.</li> <li>• 'ch.dc': Chicken <i>Gallus gallus</i> double cone.</li> <li>• 'st.dc': Starling <i>Sturnus vulgaris</i> double cone.</li> <li>• 'cf.r': Canid <i>Canis familiaris</i> rod</li> <li>• 'md.r1': Housefly <i>Musca domestica</i> R1-6 photoreceptor.</li> <li>• 'ra.dc': Triggerfish <i>Rhinecanthus aculeatus</i> double cone.</li> </ul>
illum	<p>illuminants. Options are:</p> <ul style="list-style-type: none"> <li>• 'none': no illuminant data.</li> <li>• 'all': all background spectral data.</li> <li>• 'bluesky' open blue sky.</li> <li>• 'D65': standard daylight.</li> <li>• 'forestshade' forest shade.</li> </ul>
trans	<p>Ocular transmission data. Options are:</p> <ul style="list-style-type: none"> <li>• 'none': no transmission data.</li> <li>• 'all': all transmission data.</li> <li>• 'bluetit': blue tit <i>Cyanistes caeruleus</i> ocular transmission (from Hart et al. 2000).</li> <li>• 'blackbird': blackbird <i>Turdus merula</i> ocular transmission (from Hart et al. 2000).</li> </ul>
bkg	<p>background spectra. Options are:</p> <ul style="list-style-type: none"> <li>• 'none': no background spectral data.</li> <li>• 'all': all background spectral data.</li> <li>• 'green': green foliage.</li> </ul>
plot	should the spectral data be plotted, or returned instead (defaults to FALSE)?
...	additional graphical options passed to <code>plot.rspec</code> when <code>plot = TRUE</code> .

**Value**

An object of class `rspec` (when `plot = FALSE`), containing a wavelength column 'w1' and spectral data binned at 1 nm intervals from 300-700 nm.

**Author(s)**

Thomas White <thomas.white026@gmail.com>

Rafael Maia <rm72@zips.uakron.edu>

## Examples

```
# Plot the honeybee's receptors
sensdata(visual = "apis", ylab = "Absorbance", plot = TRUE)

# Plot the average UV vs V avian receptors
sensdata(visual = c("avg.v", "avg.uv"), ylab = "Absorbance", plot = TRUE)

# Retrieve the CIE colour matching functions as an rspec object
ciedat <- sensdata(visual = c("cie2", "cie10"))
```

---

sensmodel

*Modeling spectral sensitivity*

---

## Description

Models spectral sensitivity (with oil droplets; optional) based on peak cone sensitivity according to the models of Govardovskii et al. (2000) and Hart & Vorobyev (2005).

## Usage

```
sensmodel(peaksens, range = c(300, 700), lambdacut = NULL,
          Bmid = NULL, oiltype = NULL, beta = TRUE, om = NULL,
          integrate = TRUE)
```

## Arguments

peaksens	(required) a vector with peak sensitivities for the cones to model.
range	a vector of length 2 for the range over which to calculate the spectral sensitivities (defaults to 300nm to 700nm).
lambdacut	a vector of same length as peaksens that lists the cut-off wavelength value for oil droplets. Needs either Bmid or oiltype to also be entered. See Hart and Vorobyev (2005).
Bmid	a vector of same length as peaksens that lists the gradient of line tangent to the absorbance spectrum of the oil droplets. See Hart and Vorobyev (2005).
oiltype	a list of same length as peaksens that lists the oil droplet types (currently accepts only "T", "C", "Y", "R", "P") when Bmid is not known. Calculates Bmid based on the regression equations found in Hart and Vorobyev (2005).
beta	logical. If TRUE the sensitivities will include the beta peak See Govardovskii et al.(2000) (defaults to TRUE).
om	a vector of same length as range1-range2 that contains ocular media transmission data. If included, cone sensitivity will be corrected for ocular media transmission. Currently accepts "bird" using values from Hart et al. (2005), or user-defined values.
integrate	logical. If TRUE, each curve is transformed to have a total area under the curve of 1 (best for visual models; defaults to TRUE). NOTE: integration is applied before any effects of ocular media are considered, for compatibility with visual model procedures.

**Value**

A data frame of class `rspec` containing each cone model as a column.

**Author(s)**

Pierre-Paul Bitton <bittonp@uwindsor.ca>, Chad Eliason <cme16@ziips.uakron.edu>

**References**

Govardovskii VI, Fyhrquist N, Reuter T, Kuzmin DG and Donner K. 2000. In search of the visual pigment template. *Visual Neuroscience* 17:509-528

Hart NS, and Vorobyev M. 2005. modeling oil droplet absorption spectra and spectral sensitivities of bird cone photoreceptors. *Journal of Comparative Physiology A*. 191: 381-392

Hart NS, Partridge JC, Cuthill IC, Bennett AT (2000) Visual pigments, oil droplets, ocular media and cone photoreceptor distribution in two species of passerine bird: the blue tit (*Parus caeruleus* L) and the blackbird (*Turdus merula* L). *J Comp Physiol A* 186:375-387

**Examples**

```
# Blue tit visual system based on Hart et al (2000)
bluesens <- sensmodel(c(371, 448, 502, 563),
  beta = FALSE,
  lambdacut = c(330, 413, 507, 572),
  oiltype = c("T", "C", "Y", "R"), om = TRUE
)

# Danio aequipinnatus based on Govardovskii et al. (2000)
daniosens <- sensmodel(c(357, 411, 477, 569))
```

---

sicalis

*Spectral curves from three body regions of stripe-tailed yellow finch  
(Sicalis citrina) males*

---

**Description**

dataset containing reflectance measurements from 3 body parts ("C": crown, "B": breast, "T": throat) from seven male stripe-tailed yellow finches

**Author(s)**

Rafael Maia <rm72@ziips.uakron.edu>

---

`spec2rgb`*Spectrum to rgb colour conversion*

---

**Description**

Calculates rgb values from spectra based on human colour matching functions.

**Usage**

```
spec2rgb(rspeccdata, alpha = 1)
```

**Arguments**

<code>rspeccdata</code>	(required) a data frame, possibly an object of class <code>rspec</code> that has wavelength range in the first column, named 'wl', and spectral measurements in the remaining columns.
<code>alpha</code>	alpha value to use for colors (defaults to 1, opaque).

**Value**

A character vector consisting of hexadecimal colour values for passing to further plotting functions.

**Author(s)**

Hugo Gruson <hugo.gruson+R@normalesup.org>

Chad Eliason <cme16@zip.s.uakron.edu>

**References**

CIE(1932). Commission Internationale de l'Eclairage Proceedings, 1931. Cambridge: Cambridge University Press.

**Examples**

```
data(teal)
spec2rgb(teal)

# Plot data using estimated perceived colour
plot(teal, col = spec2rgb(teal), type = "o")
```

---

subset.rspec                    *Subset rspec, vismodel, and colspace objects*

---

## Description

Subsets various object types based on a given vector or grep partial matching of data names.

## Usage

```
## S3 method for class 'rspec'  
subset(x, subset, ...)  
  
## S3 method for class 'colspace'  
subset(x, subset, ...)  
  
## S3 method for class 'vismodel'  
subset(x, subset, ...)
```

## Arguments

**x**                    (required) an object of class `rspec`, `vismodel`, or `colspace`, containing spectra, visual model output or colorspace data to subset.

**subset**                a string used for partial matching of observations.

**...**                 additional attributes passed to `grep`. Ignored if `subset` is logical.

## Value

a subsetted object of the same class as the input object.

## Note

if more than one value is given to `subset`, any spectra that matches *either* condition will be included. It's a union, not an intersect.

## Author(s)

Chad Eliason <cme16@zip.uakron.edu>

## Examples

```
data(sicalis)  
vis.sicalis <- vismodel(sicalis)  
tcs.sicalis <- colspace(vis.sicalis, space = "tcs")  
  
# Subset all 'crown' patches (C in file names)  
head(subset(sicalis, "C"))  
subset(vis.sicalis, "C")  
subset(tcs.sicalis, "C")[, seq_len(5)]
```

```
subset(sicalis, c("B", "C"))
subset(sicalis, "T", invert = TRUE)
```

---

summary.colospace      *Colourspace data summary*

---

## Description

Returns the attributes of colospace objects.

## Usage

```
## S3 method for class 'colospace'
summary(object, by = NULL, ...)
```

## Arguments

object	(required) a colospace object.
by	when the input is in tcs colourspace, by is either a single value specifying the range of colour points for which summary tetrahedral-colourspace variables should be calculated (for example, by = 3 indicates summary will be calculated for groups of 3 consecutive colour points (rows) in the quantum catch colour data frame) or a vector containing identifications for the rows in the quantum catch colour data frame (in which case summaries will be calculated for each group of points sharing the same identification). If by is left blank, the summary statistics are calculated across all colour points in the data.
...	class consistency (ignored).

## Value

returns all attributes of the data as mapped to the selected colourspace, including options specified when calculating the visual model. Also return the default data.frame summary, except when the object is the result of tcs, in which case the following variables are output instead:

centroid.u, .s, .m, .l the centroids of usm1 coordinates of points.

c.vol the total volume occupied by the points.

rel.c.vol volume occupied by the points relative to the tetrahedron volume.

colspan.m the mean hue span.

colspan.v the variance in hue span.

huedisp.m the mean hue disparity.

huedisp.v the variance in hue disparity.

mean.ra mean saturation.

max.ra maximum saturation achieved by the group of points.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**References**

Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.

Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.

**Examples**

```
# Colour hexagon
data(flowers)
vis.flowers <- vismodel(flowers,
  visual = "apis", qcatch = "Ei", relative = FALSE,
  vonkries = TRUE, bkg = "green"
)
flowers.hex <- hexagon(vis.flowers)
summary(flowers.hex)

# Tetrahedral model
data(sicalis)
vis.sicalis <- vismodel(sicalis, visual = "avg.uv")
csp.sicalis <- colspace(vis.sicalis)
summary(csp.sicalis, by = rep(c("C", "T", "B"), 7))
```

---

summary.rimg

*Image summary*

---

**Description**

Returns the attributes of, and optionally plots, an image.

**Usage**

```
## S3 method for class 'rimg'
summary(object, plot = FALSE, axes = TRUE, col = NULL,
  ...)
```

**Arguments**

object	(required) an image of class rimg, or list thereof.
plot	logical; plot the image and, if the image is color-classified, the colours corresponding to colour class categories side-by-side? Defaults to FALSE.
axes	should axes be drawn when plot = TRUE? (defaults to TRUE).



`col` optional vector of colours when plotting colour-classified images with `plot = TRUE`. Defaults to the mean RGB values of the k-means centres (i.e. the 'original' colours).

`...` additional graphical options when `plot = TRUE`. Also see [par](#).

### Value

Either the RGB values of the k-means centres from the colour-classified image, or a plot of both the image and specified colours (when `plot = TRUE`).

### Author(s)

Thomas E. White <thomas.white026@gmail.com>

### Examples

```
## Not run:
papilio <- getimg(system.file("testdata/images/papilio.png", package = "pavo"))
papilio_class <- classify(papilio, kcols = 4)
summary(papilio_class)

# Plot the colour-classified image alongside the colour class palette
summary(papilio_class, plot = TRUE)

# Multiple images
snakes <- getimg(system.file("testdata/images/snakes", package = "pavo"))
snakes_class <- classify(snakes, kcols = 3)
summary(snakes_class, plot = TRUE)

## End(Not run)
```

---

summary.rspec

*Colourimetric variables*

---

### Description

Calculates all 23 colourimetric variables reviewed in Montgomerie (2006).

### Usage

```
## S3 method for class 'rspec'
summary(object, subset = FALSE, wlmin = NULL,
        wlmax = NULL, ...)
```

**Arguments**

object	(required) a data frame, possibly an object of class <code>rspec</code> , with a column with wavelength data, named 'wl', and the remaining column containing spectra to process.
subset	Either FALSE (the default), TRUE, or a character vector. If FALSE, all variables calculated are returned. If TRUE, only a subset of the complete output (composed of B2, S8 and H1; the variables described in Andersson and Prager 2006) are returned. Finally, a user-specified string of variable names can be used in order to filter and show only those variables.
wlmin, wlmax	minimum and maximum used to define the range of wavelengths used in calculations (default is to use entire range in the <code>rspec</code> object).
...	class consistency (ignored)

**Value**

A data frame containing either 23 or 5 (`subset = TRUE`) variables described in Montgomerie (2006) with spectra name as row names. The colorimetric variables calculated by this function are described in Montgomerie (2006) with corrections included in the README CLR file from the May 2008 distribution of the CLR software. Authors should reference both this package, Montgomerie (2006), and the original reference(s). Description and notes on the measures:

**B1 (Total brightness):** Sum of the relative reflectance over the entire spectral range (area under the curve). Frequently used but should be discouraged because values are difficult to compare across studies (B2 is preferred). REF 1-3, 7, 9-11, 13

**B2 (Mean brightness):** Mean relative reflectance over the entire spectral range. This is preferred to B1 since values are easier to compare across studies. REF 4, 12

**B3 (Intensity):** Maximum relative reflectance (Reflectance at wavelength of maximum reflectance). Note that may be sensitive to noise near the peak. REF 1, 5, 6

**S1 (Chroma):** Relative contribution of a spectral range to the total brightness (B1) S1 is arbitrarily divided in 6 measures of chroma based on the wavelength ranges normally associated with specific hues. The values are calculated using the following ranges: S1U (UV, if applicable): lambda min-400nm; S1V (Violet) lambda min-415nm; S1B (Blue) 400nm-510nm; S1G (Green) 510nm-605nm; S1Y (Yellow) 550nm-625nm; S1R (Red) 605nm-lambda max. REF 2, 7, 8, 11-13

**S2 (Spectral saturation):**  $R_{max}/R_{min}$  This measure is sensitive to spectral noise. Proper interpretation of this value may be difficult for spectra with multiple peaks in the range of interest. REF 1

**S3 (Chroma):** Reflectance over the  $R_{max} \pm 50$ nm range divided by B1. Values for peaks within 50nm of either the minimum or maximum range of the data will not be comparable since the area under the curve for the area of interest will not always be based on the same wavelength range. Therefore, S3 should be interpreted with caution for peaks in the UV or Red range. REF 11

**S4 (Spectral purity):**  $lb_{maxnegl}$ , calculated by approximating the derivative of the spectral curve. As such, it is very sensitive to noise and should only be considered when data is adequately smoothed. NAs are returned for curves which do not, at any range of wavelength, decrease in intensity. Therefore, reflectance curves for brown and red surfaces, for example, should not generate a values. REF 1

S5 (Chroma): Similar in design to segment classification measures (see Montgomerie 2006 for details). REF 10

S6 (Contrast):  $R_{max} - R_{min}$ . Because it uses both  $R_{min}$  and  $R_{max}$ , this measure may be sensitive to spectral noise. REF 5, 6

S7 (Spectral saturation): Difference between the relative reflectance before and after the wavelength at which reflectance is halfway between its minimum ( $R_{min}$ ) and its maximum ( $R_{max}$ ). Somewhat sensitive to noise and can be misleading when more than one maxima and/or minima are present. REF 3, 9

S8 (Chroma):  $(R_{max} - R_{min})/B2$ . Because it uses both  $R_{min}$  and  $R_{max}$ , this measure may be sensitive to spectral noise. REF 3, 13

S9 (Carotenoid chroma):  $(R700 - R450)/R700$ . Should only be used when the color of the surface is clearly due to carotenoid pigmentation and  $R450$  is lower than  $R700$ . Could be sensitive to noise. REF 8

S10 (Peaky chroma):  $(R_{max} - R_{min})/B2 \times lb_{maxnegl}$ . Should be used with properly smoothed curves. REF 7

H1 (Peak wavelength, hue): Wavelength of maximum reflectance. May be sensitive to noise and may be variable if there is more than one maxima. REF 1, 2, 4, 6, 7, 10-13

H2 (Hue): Wavelength at  $b_{maxneg}$ . Should be calculated using smoothed data. REF 2, 13

H3 (Hue): Wavelength at  $R_{mid}$ . Sensitive to noisy spectra and may be variable if there are more than one maxima and minima. REF 3, 9, 13

H4 (Hue): Similar in design to segment classification measures see Montgomerie (2006) for details. REF 10

H5 (Hue): Wavelength at  $b_{max}$ . Sensitive to noise and may be variable if there is more than one maxima and minima. REF 5

### Note

If minimum wavelength is over 400, UV chroma is not computed.

Variables which compute  $b_{max}$  and  $b_{maxneg}$  should be used with caution, for they rely on smoothed curves to remove noise, which would otherwise result in spurious results. Make sure chosen smoothing parameters are adequate.

Smoothing affects only B3, S2, S4, S6, S10, H2, and H5 calculation. All other variables can be reliably extracted using non-smoothed data.

### Author(s)

Pierre-Paul Bitton <bittonp@windsor.ca>, Rafael Maia <rm72@zips.uakron.edu>

### References

Montgomerie R. 2006. Analyzing colors. In Hill, G.E, and McGraw, K.J., eds. Bird Coloration. Volume 1 Mechanisms and measurements. Harvard University Press, Cambridge, Massachusetts.

References describing variables:

1- Andersson, S. 1999. Morphology of uv reflectance in a whistling-thrush: Implications for the study of structural colour signalling in birds. *Journal of Avian Biology* 30:193-204.

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- 5- Keyser, A. and G. Hill. 1999. Condition-dependent variation in the blue-ultraviolet coloration of a structurally based plumage ornament. *Proceedings of the Royal Society B* 266:771-777.
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- 12- Siefferman, L. and G. Hill. 2005. UV-blue structural coloration and competition for nestboxes in male eastern bluebirds. *Animal Behaviour* 69:67-72.
- 13- Smiseth, P., J. Ornborg, S. Andersson, and T. Amundsen. 2001. Is male plumage reflectance correlated with paternal care in bluethroats? *Behavioural Ecology* 12:164-170.

## Examples

```
data(sicalis)
summary(sicalis)
summary(sicalis, subset = TRUE)
summary(sicalis, subset = c("B1", "H4"))
```

---

summary.vismodel

*Visual model summary*

---

## Description

Returns the attributes used when calculating a visual model using `vismodel`

**Usage**

```
## S3 method for class 'vismodel'  
summary(object, ...)
```

**Arguments**

```
object      (required) Results of vismodel  
...         class consistency (ignored)
```

**Value**

Returns all attributes chosen when calculating the visual model, as well as the default data.frame  
summary

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**References**

Vorobyev, M., Osorio, D., Bennett, A., Marshall, N., & Cuthill, I. (1998). Tetrachromacy, oil droplets and bird plumage colours. *Journal Of Comparative Physiology A-Neuroethology Sensory Neural And Behavioral Physiology*, 183(5), 621-633.

Hart, N. S. (2001). The visual ecology of avian photoreceptors. *Progress In Retinal And Eye Research*, 20(5), 675-703.

Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.

Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.

**Examples**

```
data(sicalis)  
vis.sicalis <- vismodel(sicalis, visual = "avg.uv")  
summary(vis.sicalis)
```

---

tcsplot

*Interactive plot of a tetrahedral colourspace*

---

**Description**

Produces an interactive 3D plot of a tetrahedral colourspace using OpenGL capabilities.

Plots points in a tetrahedral colour space

**Usage**

```
tcsplot(tcsdata, size = 0.02, alpha = 1, col = "black",
        vertexsize = 0.02, achro = TRUE, achrosize = 0.01,
        achrocol = "grey", lwd = 1, lcol = "lightgrey", new = FALSE,
        hspin = FALSE, vspin = FALSE, floor = TRUE)
```

```
tcspoints(tcsdata, size = 0.02, col = "black", alpha = 1)
```

```
tcsvol(tcsdata, col = "black", alpha = 0.2, grid.alpha = 1,
        grid = TRUE, fill = TRUE, lwd = 1)
```

**Arguments**

tcsdata	(required) a data frame, possibly a result from the <code>colspace</code> or <code>tetraspace</code> function, containing values for the 'x', 'y' and 'z' coordinates as columns (labeled as such).
size	size of the points in the plot (defaults to 0.02)
alpha	transparency of points (or volume fill in <code>tcsvol</code> )
col	colour of the points in the plot (defaults to black)
vertexsize	size of the points at the vertices
achro	logical. Should the achromatic center be plotted? (defaults to TRUE)
achrosize	size of the point in the achromatic centre
achrocol	colour of the point in the achromatic centre
lwd, lcol	graphical parameters for the edges of the tetrahedron.
new	should a new 3D plot be called (defaults to FALSE)?
hspin	if TRUE, the graphic will spin horizontally (around the 'z' axis)(defaults to FALSE)
vspin	if TRUE, the graphic will spin vertically (around the 'x' axis)(defaults to FALSE)
floor	if TRUE, a reference xy plane is plotted under the tetrahedron (defaults to TRUE)
grid.alpha	transparency of the volume polygon grid lines
grid	if TRUE, connects the polygon outlining the volume occupied by points (defaults to TRUE)
fill	if TRUE, fills the volume occupied by points (WARNING: transparency is not saved properly if exported using <code>rgl.postscript</code> )(defaults to TRUE).

**Value**

`tcsplot` creates a 3D plot using functions of the package `rgl`, based on OpenGL capabilities. Plot is interactive and can be manipulated with the mouse (left button: rotate along 'z' axis; right button: rotate along 'x' axis; third button: zoom). `tcsvol` creates polygon based on points, determining the volume occupied by them in the colorspace. `tcspoints` adds points to the plot. Points are currently plotted only as spheres to maintain export capabilities.

`tcspoints` creates 3D points in a tetrahedral colour space plot produced by `tcsplot` using functions of the package `rgl`, based on OpenGL capabilities.

`tcsvol` creates a 3D convex hull within a `tcsplot` object.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**References**

- Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.
- Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.

**See Also**

[spheres3d](#), [rgl.postscript](#), [rgl.snapshot](#), [rgl.material](#)

**Examples**

```
## Not run:
# For plotting
data(sicalis)
vis.sicalis <- vismodel(sicalis, visual = "avg.uv")
tcs.sicalis <- colspace(vis.sicalis, space = "tcs")
tcsplot(tcs.sicalis, size = 0.005)
rgl::rgl.postscript("testplot.pdf", fmt = "pdf")
rgl::rgl.snapshot("testplot.png")

# For adding points
patch <- rep(c("C", "T", "B"), 7)
tcs.crown <- subset(tcs.sicalis, "C")
tcs.breast <- subset(tcs.sicalis, "B")
tcsplot(tcs.crown, col = "blue")
tcspoints(tcs.breast, col = "red")

# For plotting convex hull
tcsplot(tcs.sicalis, col = "blue", size = 0.005)
tcsvol(tcs.sicalis)

## End(Not run)
```

---

teal

*Angle-resolved reflectance data for the iridescent wing patch of a male green-winged teal (Anas carolinensis)*

---

**Description**

dataset containing reflectance measurements from the wing patch of a single male at different incident angles (15-75 degrees in 5-degree increments).

**Author(s)**

Chad Eliason <cme16@zip.s.uakron.edu>

---

transmissiondata      *Default ocular transmission data*

---

**Description**

Default ocular transmission data

**Author(s)**

Rafael Maia <rm72@zip.s.uakron.edu>

**References**

Hart, N. S., Partridge, J. C., Cuthill, I. C., Bennett, A. T. D. (2000). Visual pigments, oil droplets, ocular media and cone photoreceptor distribution in two species of passerine

---

ttvertex      *vertex for the tetrahedral color space*

---

**Description**

internal data for plotting devices.

**Author(s)**

Rafael Maia <rm72@zip.s.uakron.edu>

**References**

Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.



vismodel

*Visual models***Description**

Calculates quantum catches at each photoreceptor. Both raw and relative values can be returned, for use in a suite of colourspace and non-colourspace models.

**Usage**

```
vismodel(rspeccdata, visual = c("avg.uv", "avg.v", "bluetit",
  "ctenophorus", "star", "pfowl", "apis", "canis", "cie2", "cie10",
  "musca", "segment", "habronattus", "rhinecanthus"),
  achromatic = c("none", "bt.dc", "ch.dc", "st.dc", "md.r1", "ra.dc",
  "cf.r", "m1", "l", "all"), illum = c("ideal", "bluesky", "D65",
  "forestshade"), trans = c("ideal", "bluetit", "blackbird"),
  qcatch = c("Qi", "fi", "Ei"), bkg = c("ideal", "green"),
  vonkries = FALSE, scale = 1, relative = TRUE)
```

**Arguments**

- |            |  |
|------------|--|
| rspeccdata | (required) a data frame, possibly an object of class <code>rspec</code> that has wavelength range in the first column, named 'wl', and spectral measurements in the remaining columns.   |
| visual     | the visual system to be used. Options are: <ul style="list-style-type: none"> <li>• a data frame such as one produced containing by <code>sensmodel</code>, containing user-defined sensitivity data for the receptors involved in colour vision. The data frame must contain a 'wl' column with the range of wavelengths included, and the sensitivity for each other cone as a column.</li> <li>• 'apis': Honeybee <i>Apis mellifera</i>.</li> <li>• 'avg.uv': average avian UV system (default).</li> <li>• 'avg.v': average avian V system.</li> <li>• 'bluetit': Blue tit <i>Cyanistes caeruleus</i>.</li> <li>• 'canis': Canid <i>Canis familiaris</i>.</li> <li>• 'cie2': 2-degree colour matching functions for CIE models of human colour vision. Functions are linear transformations of the 2-degree cone fundamentals of Stockman &amp; Sharpe (2000), as ratified by the CIE (2006).</li> <li>• 'cie10': 10-degree colour matching functions for CIE models of human colour vision. Functions are linear transformations of the 10-degree cone fundamentals of Stockman &amp; Sharpe (2000), as ratified by the CIE (2006).</li> <li>• 'ctenophorus': Ornate dragon lizard <i>Ctenophorus ornatus</i>.</li> <li>• 'musca': Housefly <i>Musca domestica</i>.</li> <li>• 'pfowl': Peafowl <i>Pavo cristatus</i>.</li> <li>• 'segment': Generic tetrachromat 'viewer' for use in the segment analysis of Endler (1990).</li> </ul> |

	<ul style="list-style-type: none"> <li>• 'star': Starling <i>Sturnus vulgaris</i>.</li> <li>• 'habronattus': Jumping spider <i>Habronattus pyrrithrix</i>.</li> <li>• 'rhinecanthus': Triggerfish <i>Rhinecanthus aculeatus</i>.</li> </ul>
achromatic	<p>the sensitivity data to be used to calculate luminance (achromatic) receptor stimulation. Either a vector containing the sensitivity for a single receptor, or one of the options:</p> <ul style="list-style-type: none"> <li>• 'none': no achromatic stimulation calculated (default).</li> <li>• 'bt.dc': Blue tit <i>Cyanistes caeruleus</i> double cone.</li> <li>• 'ch.dc': Chicken <i>Gallus gallus</i> double cone.</li> <li>• 'st.dc': Starling <i>Sturnus vulgaris</i> double cone.</li> <li>• 'md.r1': Housefly <i>Musca domestica</i> R1-6 photoreceptor.</li> <li>• 'ra.dc': Triggerfish <i>Rhinecanthus aculeatus</i> double cone.</li> <li>• 'cf.r': Canid <i>Canis familiaris</i> cone.</li> <li>• 'm1': the summed response of the two longest-wavelength photoreceptors.</li> <li>• 'l': the longest-wavelength photoreceptor.</li> <li>• 'all': the summed response of all photoreceptors.</li> </ul>
illum	<p>either a vector containing the illuminant, or one of the options:</p> <ul style="list-style-type: none"> <li>• 'ideal': homogeneous illuminance of 1 across wavelengths (default)</li> <li>• 'bluesky': open blue sky.</li> <li>• 'D65': standard daylight.</li> <li>• 'forestshade': forest shade.</li> </ul>
trans	<p>either a vector containing the ocular or environmental transmission spectra, or one of the options:</p> <ul style="list-style-type: none"> <li>• 'ideal': homogeneous transmission of 1 across all wavelengths (default)</li> <li>• 'bluetit': blue tit <i>Cyanistes caeruleus</i> ocular transmission (from Hart et al. 2000).</li> <li>• 'blackbird': blackbird <i>Turdus merula</i> ocular transmission (from Hart et al. 2000).</li> </ul>
qcatch	<p>Which quantal catch metric to return. Options are:</p> <ul style="list-style-type: none"> <li>• 'Qi': Quantum catch for each photoreceptor (default)</li> <li>• 'fi': Quantum catch according to Fechner's law (the signal of the receptor channel is proportional to the logarithm of the quantum catch)</li> <li>• 'Ei': Hyperbolic-transformed quantum catch, where <math>E_i = Q_i / (Q_i + 1)</math>.</li> </ul>
bkg	<p>background spectrum. Note that this will have no effect when <code>vonkries = FALSE</code>. Either a vector containing the spectral data, or one of the options:</p> <ul style="list-style-type: none"> <li>• 'ideal': homogeneous illuminance of 1 across all wavelengths (default).</li> <li>• 'green': green foliage.</li> </ul>
vonkries	<p>logical. Should the von Kries color correction transformation be applied? (defaults to FALSE).</p>
scale	<p>a value by which the illuminant will be multiplied. Useful for when the illuminant is a relative value (i.e. transformed to a maximum of 1 or to a percentage), and does not correspond to quantum flux units (<math>\text{\\$umol*s}^{-1}\text{*m}^{-2}</math>). Useful values are, for example, 500 (for dim light) and 10000 (for bright illumination). Note that if <code>vonkries = TRUE</code> this transformation has no effect.</p>

relative            should relative quantum catches be returned (i.e. is it a color space model? Defaults to TRUE).

### Value

An object of class `vismodel` containing the photon catches for each of the photoreceptors considered. Information on the parameters used in the calculation are also stored and can be called using the `summary.vismodel` function.

### Author(s)

Rafael Maia <rm72@zips.uakron.edu>

Thomas White <thomas.white026@gmail.com>

### References

Vorobyev, M., Osorio, D., Bennett, A., Marshall, N., & Cuthill, I. (1998). Tetrachromacy, oil droplets and bird plumage colours. *Journal Of Comparative Physiology A-Neuroethology Sensory Neural And Behavioral Physiology*, 183(5), 621-633.

Hart, N. S., Partridge, J. C., Cuthill, I. C., Bennett, A. T. D. (2000). Visual pigments, oil droplets, ocular media and cone photoreceptor distribution in two species of passerine bird: the blue tit (*Parus caeruleus* L.) and the blackbird (*Turdus merula* L.). *Journal of Comparative Physiology A*, 186, 375-387.

Hart, N. S. (2001). The visual ecology of avian photoreceptors. *Progress In Retinal And Eye Research*, 20(5), 675-703.

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Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.

Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.

Chittka L. (1992). The colour hexagon: a chromaticity diagram based on photoreceptor excitations as a generalized representation of colour opponency. *Journal of Comparative Physiology A*, 170(5), 533-543.

Stockman, A., & Sharpe, L. T. (2000). Spectral sensitivities of the middle- and long-wavelength sensitive cones derived from measurements in observers of known genotype. *Vision Research*, 40, 1711-1737.

CIE (2006). Fundamental chromaticity diagram with physiological axes. Parts 1 and 2. Technical Report 170-1. Vienna: Central Bureau of the Commission Internationale de l' Eclairage.

Neitz, J., Geist, T., Jacobs, G.H. (1989) Color vision in the dog. *Visual Neuroscience*, 3, 119-125.

### Examples

```
# Dichromat (dingo)
data(flowers)
vis.flowers <- vismodel(flowers, visual = "canis")
```

```

di.flowers <- colspace(vis.flowers, space = "di")

# Trichromat (honeybee)
data(flowers)
vis.flowers <- vismodel(flowers, visual = "apis")
tri.flowers <- colspace(vis.flowers, space = "tri")

# Tetrachromat (blue tit)
data(sicalis)
vis.sicalis <- vismodel(sicalis, visual = "bluetit")
tcs.sicalis <- colspace(vis.sicalis, space = "tcs")

# Tetrachromat (starling), receptor-noise model
data(sicalis)
vis.star <- vismodel(sicalis, visual = "star", achromatic = "bt.dc", relative = FALSE)
dist.star <- coldist(vis.star, achromatic = TRUE)

```

---

vissyst

*Animal visual systems data*


---

### Description

Internal data for visual model calculations.

### Author(s)

Rafael Maia <rm72@zips.uakron.edu>

### References

Endler, J. A., & Mielke, P. (2005). Comparing entire colour patterns as birds see them. *Biological Journal Of The Linnean Society*, 86(4), 405-431.

---

vol

*Plot a tetrahedral color space*


---

### Description

Produces a 3D convex hull in tetrahedral color space when plotting a non-interactive tetrahedral plot.

### Usage

```

vol(tcsdata, alpha = 0.2, grid = TRUE, fill = TRUE, new = FALSE,
    ...)

```

**Arguments**

tcsdata	(required) a data frame, possibly a result from the <code>colspace</code> or <code>tetraspace</code> function, containing values for the 'x', 'y' and 'z' coordinates as columns (labeled as such).
alpha	transparency of volume (if <code>fill = TRUE</code> ).
grid	logical. if TRUE (default), draws the polygon outline defined by the points.
fill	logical. if TRUE (default), fills the volume defined by the points.
new	logical. Should a new plot be started or draw over an open plot? (defaults to FALSE)
...	additional graphical options. See <a href="#">polygon</a> and <a href="#">tetraplot</a> .

**Value**

`vol` creates a 3D convex hull within a static tetrahedral plot.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>

**Examples**

```
# For plotting
data(sicalis)
vis.sicalis <- vismodel(sicalis, visual = "avg.uv")
tcs.sicalis <- colspace(vis.sicalis, space = "tcs")
plot(tcs.sicalis)
vol(tcs.sicalis)
```

---

voloverlap

*Color volume overlap*

---

**Description**

Calculates the overlap between the volumes defined by two sets of points in cartesian space.

**Usage**

```
voloverlap(tcsres1, tcsres2, plot = FALSE, interactive = FALSE,
  col = c("blue", "red", "darkgrey"), fill = FALSE, new = TRUE,
  montecarlo = FALSE, nsamp = 1000, psize = 0.001, lwd = 1, ...)
```

**Arguments**

<code>tcsres1, tcsres2</code>	(required) data frame, possibly a result from the <code>colspace</code> function, containing values for the 'x', 'y' and 'z' coordinates as columns (labeled as such)
<code>plot</code>	logical. Should the volumes and points be plotted? (defaults to FALSE)
<code>interactive</code>	logical. If TRUE, uses the <code>rgl</code> engine for interactive plotting; if FALSE then a static plot is generated.
<code>col</code>	a vector of length 3 with the colors for (in order) the first volume, the second volume, and the overlap.
<code>fill</code>	logical. should the two volumes be filled in the plot? (defaults to FALSE)
<code>new</code>	logical. Should a new plot window be called? If FALSE, volumes and their overlap are plotted over the current plot (defaults to TRUE).
<code>montecarlo</code>	logical. If TRUE, Monte Carlo simulation is used instead of exact solution (not recommended; defaults to FALSE)
<code>nsamp</code>	if <code>montecarlo = TRUE</code> , determines the number of points to be sampled.
<code>psize</code>	if <code>montecarlo = TRUE</code> and <code>plot = TRUE</code> , sets the size to plot the points used in the Monte Carlo simulation.
<code>lwd</code>	if <code>plot = TRUE</code> , sets the line width for volume grids.
<code>...</code>	additional arguments passed to the plot. See <a href="#">vol</a>

**Value**

Calculates the overlap between the volumes defined by two set of points in colorspace. The volume from the overlap is then given relative to:

- `vsmallest` the volume of the overlap divided by the smallest of that defined by the the two input sets of color points. Thus, if one of the volumes is entirely contained within the other, this overlap will be `vsmallest = 1`.
- `vboth` the volume of the overlap divided by the combined volume of both input sets of color points.

The Monte Carlo solution is available mostly for legacy and benchmarking, and is not recommended (see notes). If used, the output will be different:

- `s_in1, s_in2` the number of sampled points that fall within each of the volumes individually.
- `s_inboth` the number of sampled points that fall within both volumes.
- `s_ineither` the number of points that fall within either of the volumes.
- `psmallest` the proportion of points that fall within both volumes divided by the number of points that fall within the smallest volume.
- `pboth` the proportion of points that fall within both volumes divided by the total number of points that fall within both volumes.

If the Monte Carlo solution is used, a number of points much greater than the default should be considered (Stoddard & Stevens(2011) use around 750,000 points, but more or fewer might be required depending on the degree of overlap).

**Note**

Stoddard & Stevens (2011) originally obtained the volume overlap through Monte Carlo simulations of points within the range of the volumes, and obtaining the frequency of simulated values that fall inside the volumes defined by both sets of color points.

Here we present an exact solution based on finding common vertices to both volumes and calculating its volume. However, the Monte Carlo solution is available through the `montecarlo=TRUE` option.

Stoddard & Stevens (2011) also return the value of the overlap relative to one of the volumes (in that case, the host species). However, for other applications this value may not be what one expects to obtain if (1) the two volumes differ considerably in size, or (2) one of the volumes is entirely contained within the other. For this reason, we also report the volume relative to the union of the two input volumes, which may be more adequate in most cases.

**Author(s)**

Rafael Maia <rm72@zips.uakron.edu>, with code from Sebastien Villeger

**References**

Stoddard, M. C., & Prum, R. O. (2008). Evolution of avian plumage color in a tetrahedral color space: A phylogenetic analysis of new world buntings. *The American Naturalist*, 171(6), 755-776.

Stoddard, M. C., & Stevens, M. (2011). Avian vision and the evolution of egg color mimicry in the common cuckoo. *Evolution*, 65(7), 2004-2013.

Villeger, S., Novack-Gottshall, P. M., & Mouillot, D. (2011). The multidimensionality of the niche reveals functional diversity changes in benthic marine biotas across geological time. *Ecology Letters*, 14(6), 561-568.

Maia, R., White, T. E., (2018) Comparing colors using visual models. *Behavioral Ecology*, ary017 doi: [10.1093/beheco/ary017](https://doi.org/10.1093/beheco/ary017)

**Examples**

```
## Not run:
data(sicalis)
tcs.sicalis.C <- subset(colspace(vismodel(sicalis)), "C")
tcs.sicalis.T <- subset(colspace(vismodel(sicalis)), "T")
tcs.sicalis.B <- subset(colspace(vismodel(sicalis)), "B")
voloverlap(tcs.sicalis.T, tcs.sicalis.B)
voloverlap(tcs.sicalis.T, tcs.sicalis.C, plot = T)
voloverlap(tcs.sicalis.T, tcs.sicalis.C, plot = T, col = seq_len(3))

## End(Not run)
```

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