

# Package ‘colorSpec’

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**Type** Package

**Title** Color Calculations with Emphasis on Spectral Data

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**Description** Calculate with spectral properties of light sources, materials, cameras, eyes, and scanners. Build complex systems from simpler parts using a spectral product algebra. For light sources, compute CCT, CRI, and SSI. For object colors, compute optimal colors and Logvinenko coordinates.

Work with the standard CIE illuminants and color matching functions, and read spectra from text files, including CGATS files. Estimate a spectrum from its response. A user guide and 8 vignettes are included.

**License** GPL (>= 3)

**LazyLoad** yes

**LazyData** yes

**Depends** R (>= 2.15)

**Imports** spacesXYZ, MASS

**Suggests** rootSolve, rgl, knitr, rmarkdown, spacesRGB, microbenchmark, arrangements

**NeedsCompilation** no

**Repository** CRAN

**VignetteBuilder** knitr

**BuildVignettes** yes

**ByteCompile** no

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colorSpec-package	<i>Package colorSpec - Color Calculations with Emphasis on Spectral Data</i>
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## Description

Package **colorSpec** is for working with spectral color data in R.

## Details

### Features:

1. a clear classification of the commonly seen spectra into 4 types - depending on the vector space to which they belong
2. flexible organization for the spectra in memory, using an S3 class - **colorSpec**
3. a product algebra for the **colorSpec** objects
4. uniform handling of biological eyes, electronic cameras, and general action spectra
5. a few advanced calculations, such as computing optimal colors (aka Macadam Limits)
6. inverse colorimetry, e.g. reflectance recovery from response
7. built-in essential tables, such as the CIE illuminants and color matching functions
8. a package logging system with log levels taken from the popular **Log4J**
9. support for reading a few spectrum file types, including CGATS
10. bonus files containing some other interesting spectra

11. minimal dependencies on other R packages

**Non-features:**

1. there is no support for many common 3D color spaces, such as CIELAB, HSL, etc.. For these spaces see packages **colorspace**, **colorscience**, **spacesRGB**, and **spacesXYZ**.
2. there are few non-linear operations
3. there is little support for scientific units; for these see package **colorscience**
4. photons are parameterized by wavelength in nanometers; other wavelength units (such as Angstrom and micron) and alternative parameterizations (such as wavenumber and electron-volt) are not available

Regarding the non-linear operations in 2, the only such operations are conversion of linear RGB to display RGB, conversion of absorbance to transmittance, and the reparameterized wavelength in [computeADL](#). The electronic camera model is purely linear with no dark current offset or other deviations.

Many ideas are taken from packages **hyperSpec**, **hsdar**, **pavo**, and **zoo**.

**Author(s)**

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**See Also**

[colorSpec](#) for the S3 class provided by this package.

**colorSpec User Guide**

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ABC

*Standard Illuminants A, B, and C (1931)*

---

**Description**

A. 1nm	standard Illuminant A, 360 to 780 nm at 1 nm intervals
B. 5nm	standard Illuminant B, 320 to 780 nm at 5 nm intervals
C. 5nm	standard Illuminant C, 320 to 780 nm at 5 nm intervals

**Format**

Each is a **colorSpec** object organized as a vector, with [quantity](#) equal to "energy".

**Details**

All of these have been divided by 100, to make the values at 560nm near 1 instead of 100.

**Source**

<http://www.cvr1.org>

**References**

Günther Wyszecki and W. S. Stiles. **Color Science : Concepts and Methods, Quantitative Data and Formulae**. Second Edition. Wiley-Interscience. 1982.

- A Table I(3.3.4) pp. 754-758.
- B Table II(3.3.4) pp. 759.
- C Table II(3.3.4) pp. 759.

**See Also**

[D50 D65](#)

**Examples**

```
summary(xyz1931.1nm)
white.point = product( D65.1nm, xyz1931.1nm, wave='auto' )
```

---

actinometric	<i>convert a colorSpec object to be actinometric</i>
--------------	--

---

**Description**

Convert a radiometric **colorSpec** object to have quantity that is actinometric (number of photons).  
Test an object for whether it is actinometric.

**Usage**

```
## S3 method for class 'colorSpec'
actinometric( x, multiplier=1, warn=FALSE )

## S3 method for class 'colorSpec'
is.actinometric( x )
```

**Arguments**

x	a <b>colorSpec</b> object
multiplier	a scalar which is multiplied by the output, and intended for unit conversion
warn	if TRUE and a conversion actually takes place, the a WARN message is issued. This makes the user aware of the conversion, so units can be verified. This can be useful when <code>actinometric()</code> is called from another <b>colorSpec</b> function.

## Details

If the `quantity` of `x` does not start with 'energy' then the quantity is not radiometric and so `x` is returned unchanged. Otherwise `x` is radiometric (energy-based), and must be converted.

If `type(x)` is 'light' then the most common radiometric energy unit is joule.

The conversion equation is:

$$Q = E * \lambda * 10^6 / (N_A * h * c)$$

where  $Q$  is the photon count,  $E$  is the energy of the photons,  $N_A$  is Avogadro's constant,  $h$  is Planck's constant,  $c$  is the speed of light, and  $\lambda$  is the wavelength. The output unit of photon count is ( $\mu$ mole of photons) = ( $6.02214 * 10^{17}$  photons). If a different unit for  $Q$  is desired, then the output should be scaled appropriately. For example, if the desired unit of photon count is exaphotons, then set `multiplier=0.602214`.

If the `quantity(x)` is 'energy->electrical', then the most common radiometric unit of responsivity to light is coulombs/joule (C/J) or amps/watt (A/W). The conversion equation is:

$$QE = R_e * ((h * c) / e) / \lambda$$

where  $QE$  is the quantum efficiency,  $R_e$  is the energy-based responsivity, and  $e$  is the charge of an electron (in C).

If the unit of `x` is not C/J, then `multiplier` should be set appropriately.

If the `quantity(x)` is 'energy->neural' or 'energy->action', the most common radiometric unit of energy is joule (J).

The conversion equation is:

$$R_p = R_e * 10^{-6} * (N_A * h * c) / \lambda$$

where  $R_p$  is the photon-based responsivity, and  $R_e$  is the energy-based responsivity, The output unit of photon count is ( $\mu$ mole of photons) = ( $6.02214 * 10^{17}$  photons). This essentially the reciprocal of the first conversion equation.

The argument `multiplier` is applied to the right side of all the above conversion equations.

## Value

`actinometric()` returns a `colorSpec` object with `quantity` that is actinometric (photon-based) and not radiometric (energy-based). If `type(x)` is a material type ('material' or 'responsivity.material') then `x` is returned unchanged.

If `quantity(x)` starts with 'photons', then `is.actinometric()` returns TRUE, and otherwise FALSE.

## Note

To log the executed conversion equation, execute `cs.options(loglevel='INFO')`.

## Source

Wikipedia. **Photon counting**. [http://en.wikipedia.org/wiki/Photon\\_counting](http://en.wikipedia.org/wiki/Photon_counting)

**See Also**

[quantity](#), [type](#), [cs.options](#), [radiometric](#)

**Examples**

```
colSums( solar.irradiance ) # step size is 1nm, from 280 to 1000 nm. organized as a matrix
# AirMass.0 GlobalTilt AirMass.1.5
# 944.5458 740.3220 649.7749 # irradiance, watts*m^{-2}

colSums( actinometric(solar.irradiance) )
# AirMass.0 GlobalTilt AirMass.1.5
# 4886.920 3947.761 3522.149 # photon irradiance, (umoles of photons)*sec^{-1}*m^{-2}

colSums( actinometric(solar.irradiance,multiplier=0.602214) )
# AirMass.0 GlobalTilt AirMass.1.5
# 2942.972 2377.397 2121.088 # photon irradiance, exaphotons*sec^{-1}*m^{-2}
```

---

appliespec

*apply a function to each spectrum in a colorSpec object*


---

**Description**

apply a function to each spectrum in a **colorSpec** object

**Usage**

```
## S3 method for class 'colorSpec'
appliespec( x, FUN, ... )
```

**Arguments**

x	a <b>colorSpec</b> object with N wavelengths
FUN	a function that takes an N-vector as argument and returns an N-vector
...	additional arguments passed to FUN

**Details**

appliespec() simply calls [apply\(\)](#) with the right MARGIN.

**Value**

a **colorSpec** object with the same dimensions, [wavelength](#), [quantity](#), and [organization](#). If FUN does not return an N-vector, it is an ERROR and appliespec() returns NULL.

**See Also**

[quantity](#), [wavelength](#), [linearize](#), [organization](#), [apply](#)

**Examples**

```
# convert absorbance to transmittance
path = system.file( "extdata/stains/Hematoxylin.txt", package='colorSpec' )
x = readSpectra( path )
x = applyspec( x, function(y) {10^(-y)} ) # this is what linearize(x) does
```

---

as.data.frame	<i>Convert a colorSpec Object to a data.frame</i>
---------------	---

---

**Description**

convert a **colorSpec** object to a data.frame

**Usage**

```
## S3 method for class 'colorSpec'
as.data.frame( x, row.names=NULL, optional=FALSE, organization='auto', ... )
```

**Arguments**

x	a <b>colorSpec</b> object
organization	The organization of the returned data.frame, which can be 'row', 'col', or 'auto'. If 'auto', then 'row' or 'col' is selected automatically, see <b>Details</b>
row.names	ignored
optional	ignored
...	extra arguments ignored

**Details**

If organization is 'auto', and the organization of x is 'df.row', then organization is set to 'row' and the returned data.frame has the spectra in the rows. Otherwise the returned data.frame has the spectra in the columns.

**Value**

If the returned data.frame has the spectra in the rows, then the spectra are in a matrix in the last column (with name spectra), and any existing extradata are also returned in the other columns. The wavelengths are only present in character form, as the colnames of the matrix. If the returned data.frame has the spectra in the columns, then the wavelengths are in the first column, and the spectra are in the other columns.

**See Also**

[as.matrix](#), [extradata](#)



---

atmosphere                      *atmospheric transmittance along a horizontal path*

---

### Description

Calculate transmittance along a horizontal optical path in the atmosphere, as a function of length (distance) and the molecular and aerosol properties. Because the path is horizontal, the atmospheric properties are assumed to be constant on the path. Only molecular scattering is considered. There is no modeling of molecular absorption; for visible wavelengths this is reasonable.

### Usage

```
atmosTransmittance( distance, wavelength=380:720,
                   molecules=list(N=2.547305e25,n0=1.000293),
                   aerosols=list(metrange=25000,alpha=0.8,beta=0.0001) )
```

### Arguments

distance	the length of the optical path, in meters. It can also be a numeric vector of lengths.
wavelength	a vector of wavelengths, in nm, for the transmittance calculations
molecules	a list of molecular properties, see <b>Details</b> . If this is NULL, then the molecular transmittance is identically 1.
aerosols	a list of aerosol properties, see <b>Details</b> . If this is NULL, then the aerosol transmittance is identically 1.

### Details

The list `molecules` has 2 parameters that describe the molecules in the atmosphere.  $N$  is the molecular density of the atmosphere at sea level, in *molecules/meter*<sup>3</sup>. The given default is the density at sea level.  $n_0$  is the refractive index of pure molecular air (with no aerosols). For the molecular attenuation, the standard model for Rayleigh scattering is used, and there is no modeling of molecular absorption.

The list `aerosols` has 3 parameters that describe the aerosols in the atmosphere. The standard Angstrom aerosol attenuation model is:

$$attenuation(\lambda) = \beta * (\lambda/\lambda_0)^{-\alpha}$$

$\alpha$  is the Angstrom exponent, and is dimensionless. *attenuation* and  $\beta$  have unit  $m^{-1}$ . And  $\lambda_0=550nm$ .

`metrange` is the *Meteorological Range* of the atmosphere in meters, as defined by *Koschmieder*. This is the distance at which the transmittance=0.02 at  $\lambda_0$ . If `metrange` is not NULL (the default is 25000) then both  $\alpha$  and  $\beta$  are calculated to achieve this desired `metrange`, and the supplied  $\alpha$  and  $\beta$  are ignored.  $\alpha$  is calculated from `metrange` using the *Kruse* model, see **Note**.  $\beta$  is calculated so that the product of molecular and aerosol transmittance yields the desired `metrange`. In fact:

$$\beta = -\mu_0 - \log(0.02)/V_r$$

where  $\mu_0$  is the molecular attenuation at  $\lambda_0$ , and  $V_r$  is the meteorological range. For a log message with the calculated values, execute `cs.options(loglevel='INFO')` before calling `atmosTransmittance()`.

### Value

`atmosTransmittance()` returns a **colorSpec** object with `quantity` equal to 'transmittance'. There is a spectrum in the object for each value in the vector distance. The specnames are set to `sprintf("dist=%gm", distance)`.

The final transmittance is the product of the molecular transmittance and the aerosol transmittance. If both molecules and aerosols are NULL, then the final transmittance is identically 1; the atmosphere has become a vacuum.

### Note

The Kruse model for  $\alpha$  as a function of  $V_r$  is defined in 3 pieces. For  $0 \leq V_r < 6000$ ,  $\alpha = 0.585 * (V_r/1000)^{1/3}$ . For  $6000 \leq V_r < 50000$ ,  $\alpha = 1.3$ . And for  $V_r \geq 50000$ ,  $\alpha = 1.6$ . So  $\alpha$  is increasing, but not strictly, and not continuously.  $V_r$  is in meters. See *Kruse* and *Kaushal*.

The built-in object `atmosphere2003` is transmittance along an optical path that is **NOT** horizontal, and extends to outer space. This is *much* more complicated to calculate.

### References

Angstrom, Anders. On the atmospheric transmission of sun radiation and on dust in the air. *Geogr. Ann.*, no. 2. 1929.

Kaushal, H. and Jain, V.K. and Kar, S. **Free Space Optical Communication**. Springer. 2017.

Koschmieder, Harald. Theorie der horizontalen Sichtweite. *Beitrag zur Physik der Atmosphere*. 1924. **12**. pages 33-53.

P. W. Kruse, L. D. McGlauchlin, and R. B. McQuistan. **Elements of Infrared Technology: Generation, Transmission, and Detection**. J. Wiley & Sons, New York, 1962.

### See Also

[solar.irradiance](#), [specnames](#)

### Examples

```
trans = atmosTransmittance( c(5,10,15,20,25)*1000 ) # 5 distances with atmospheric defaults

# verify that transmittance[550]=0.02 at dist=25000
plot( trans, legend='bottomright', log='y' )

# repeat, but this time assign alpha and beta explicitly
trans = atmosTransmittance( c(5,10,15,20,25)*1000, aero=list(alpha=1,beta=0.0001) )
```

---

bind *Combine colorSpec Objects*

---

### Description

Take a sequence of **colorSpec** objects and combine their spectra

### Usage

```
## S3 method for class 'colorSpec'  
bind( ... )
```

### Arguments

... **colorSpec** objects with the same **wavelength** and **quantity**, and with distinct **specnames** (no duplicates)

### Details

The **organization** of the returned object is the most complex of those in the inputs, where the order of complexity is:

```
'matrix' < 'df.col' < 'df.row'
```

If the selected organization is 'df.row' the extradata is combined in a way that preserves all the columns. Missing data is filled with NAs, analogous to `plyr::rbind.fill()`.

The **metadata** of the returned object is copied from the first object in the input list.

### Value

`bind()` returns a **colorSpec** object, or NULL in case of ERROR. If the bind is successful, the number of spectra in the output object is the sum of the number of spectra in the input objects.

### See Also

[wavelength](#), [quantity](#), [specnames](#), [organization](#), [extradata](#), [metadata](#), [rbind.fill](#)

### Examples

```
Rosco = readSpectra( system.file( 'extdata/objects/Rosco.txt', package='colorSpec' ) )  
Rosco = resample( Rosco, wave=wavelength(Hoya) )  
numSpectra(Hoya)      # prints 4  
numSpectra(Rosco)     # prints 42  
  
filters = bind( Hoya, Rosco )  
numSpectra(filters)   # prints 46  
  
colnames( extradata(Hoya) )  
## [1] "SAMPLE_NAME" "FresnelReflectance" "Thickness"
```

```

colnames( extradata(Rosco) )
## [1] "Model" "SampleID" "SAMPLE_NAME" "Substrate" "RefractiveIndex" "Thickness"

## The columns in common are "SAMPLE_NAME" and "Thickness"

colnames( extradata(filters) )
## [1] "FresnelReflectance" "Model" "RefractiveIndex" "SAMPLE_NAME"
## [5] "SampleID" "Substrate" "Thickness"
##
## "SAMPLE_NAME" and "Thickness" are combined in the usual way
## The other columns are present, and missing data is filled with NAs

```

---

calibrate

*make a linear modification to a colorSpec responder*


---

### Description

make a linear modification to a **colorSpec** responder with  $M$  spectra so a specific single spectrum (the stimulus) creates a given specific response. It is generalized *white balance*.

The options are complicated, but in all cases the returned object is `multiply(x,mat)` where `mat` is an internally calculated  $M \times M$  matrix. Stated another way, the spectra in the output are linear combinations of spectra in the input  $x$ .

In case of **ERROR**, a message is logged and the original  $x$  is returned.

### Usage

```

## S3 method for class 'colorSpec'
calibrate( x, stimulus=NULL, response=NULL, method=NULL )

```

### Arguments

<code>x</code>	a <b>colorSpec</b> responder with $M$ spectra. The type must be <code>'responsivity.light'</code> or <code>'responsivity.material'</code> .
<code>stimulus</code>	a <b>colorSpec</b> object with a single spectrum, with type <code>'light'</code> or <code>'material'</code> to match $x$ . The wavelength sequence of <code>stimulus</code> must be equal to that of $x$ . If <code>stimulus</code> is <code>NULL</code> , then an appropriate default is chosen, see <b>Details</b> .
<code>response</code>	an $M$ -vector, or a scalar which is replicated to length $M$ . All entries in <code>response</code> must be positive. If <code>response</code> is <code>NULL</code> , then an appropriate default <i>may be</i> chosen, see <b>Details</b> .
<code>method</code>	an $M \times M$ <i>adaptation matrix</i> . <code>method</code> can also be <code>'scaling'</code> and it is then set to the $M \times M$ identity matrix, which scales each responsivity spectrum in $x$ independently. If $M=3$ , <code>method</code> can also be <code>'Bradford'</code> , <code>'Von Kries'</code> , or <code>'MCAT02'</code> , and it is then set to the popular corresponding <i>chromatic adaptation matrix</i> . For these

special matrices, the spectra in `x` are **not** scaled independently; there is "cross-talk".

If `method` is NULL, then an appropriate default is chosen, see **Details**.

### Details

If `stimulus` is NULL, it is set to `illuminantE()` or `neutralMaterial()` to match `x`.

If `response` is NULL and the response of `x` is `electrical` or `action`, then `response` is set to an M-vector of all 1s. If `response` is NULL and the response of `x` is `neural`, then this is an ERROR and the user is prompted to supply a specific response.

If `method` is NULL and `M=3` and the response of `x` is `neural`, then the neural response is assumed to be human, and the method is set to the 3x3 Bradford matrix.

Otherwise `method` is set to the MxM identity matrix, which scales each responsivity spectrum in `x` independently. In cameras this is usually called *white balance*, and so `calibrate()` can be considered a generalization of white balance.

### Value

a `colorSpec` object equal to `multiply(x,mat)` where `mat` is an internally calculated MxM matrix. The `quantity` and `wavelength` are preserved.

Note that `mat` is not the same as the the MxM *adaptation matrix*. To inspect `mat` execute `summary` on the returned object. If `method` is `'scaling'` then `mat` is diagonal and the diagonal entries are the M gain factors needed to achieve the calibration.

Useful data is attached as attribute `"calibrate"`.

### Note

Chromatic adaptation transforms, such as `'Bradford'`, do not belong in the realm of spectra; this is not really a spectral calculation. For more about this subject see the explanation in *Digital Color Management*, Chapter 15 - Myths and Misconceptions. This adaptation option is provided in `calibrate` because it is possible and convenient.

### References

Edward J. Giorgianni and Thomas E. Madden. **Digital Color Management: Encoding Solutions**. 2nd Edition John Wiley. 2009. Chapter 15 - Myths and Misconceptions.

### See Also

`is.regular`, `quantity`, `wavelength`, `colorSpec`, `summary`, `illuminantE`, `neutralMaterial`, `product`

### Examples

```
# make an art gallery illuminated by illuminant A, and with tristimulus XYZ as output
gallery = product( A.1nm, 'artwork', xyz1931.1nm, wave='auto')
# chromatically adapt the output XYZs to D50 white point, using Bradford matrix
gallery.D50 = calibrate( gallery, response=spacesXYZ::standardXYZ('D50') )
```

```
# make an RGB flatbead scanner from illuminant F11 and a Flea2 camera
scanner = product( subset(Fs.5nm, 'F11'), 'paper', Flea2.RGB, wave='auto')
# adjust RGB gain factors (white balance) so the perfect reflecting diffuser yields RGB=(1,1,1)
scanner = calibrate( scanner )

# same flatbead scanner, but this time with some "white headroom"
scanner = product( subset(Fs.5nm, 'F11'), 'paper', Flea2.RGB, wave='auto' )
scanner = calibrate( scanner, response=0.95 )
scanner
```

---

chop

*chop spectra into low and high parts*


---

## Description

chop all spectra in a **colorSpec** object into low and high parts at a blending interval

## Usage

```
## S3 method for class 'colorSpec'
chop( x, interval, adj=0.5 )
```

## Arguments

x	a <b>colorSpec</b> object
interval	a numeric vector with length 2 giving the endpoints of the interval, in nm
adj	a number in [0,1] defining weights of low and high parts over the interval

## Details

For each spectrum, the low and high parts sum to the original spectrum. The low part vanishes on the right of the interval, and the high part vanishes on the left.

## Value

chop(x) returns a **colorSpec** object with twice the number of spectra in x and with [organization](#) equal to 'matrix'. The names of the new spectra are formed by appending ".lo" and ".hi" to the original spectrum names.

## See Also

[organization](#),

## Examples

```
# chop blue butane flame into diatomic carbon and hydrocarbon parts
path = system.file( "extdata/sources/BlueFlame.txt", package="colorSpec" )
blueflame = readSpectra( path, seq(375,650,0.5) )
plot( chop( blueflame, interval=c(432,435), adj=0.8 ) )

# chop 'white' LED into blue and yellow parts
path = system.file( "extdata/sources/Gepe-G-2001-LED.sp", package="colorSpec" )
LED = readSpectra( path )
plot( chop( LED, c(470,495) ) )
```

---

colorSpec

*constructing and testing colorSpec Objects*

---

## Description

The function `colorSpec()` is the constructor for **colorSpec** objects.

`is.colorSpec()` tests whether an object is a valid **colorSpec** object.

`as.colorSpec()` converts other variables to a **colorSpec** object, and is designed to be overridden by other packages.

## Usage

```
colorSpec( data, wavelength, quantity='auto', organization='auto', specnames=NULL )
```

```
is.colorSpec(x)
```

```
## Default S3 method:
```

```
as.colorSpec( ... )
```

## Arguments

<code>data</code>	a vector or matrix of the spectrum values. In case <code>data</code> is a vector, there is a single spectrum and the number of points in that spectrum is the length of the vector. In case <code>data</code> is a matrix, the spectra are stored in the columns, so the number of points in each spectrum is the number of rows in <code>data</code> , and the number of spectra is the number of columns in <code>data</code> . It is OK for the matrix to have only 0 or 1 column.
<code>wavelength</code>	a numeric vector of wavelengths for all the spectra, in nm. The length of this vector must be equal to <code>NROW(data)</code> . The sequence must be increasing. The <code>wavelength</code> vector can be changed after construction.
<code>quantity</code>	a character string giving the quantity of all spectra in <code>data</code> ; see <a href="#">quantity</a> for a list of possible values. In case of 'auto', a guess is made from the <code>specnames</code> . The <code>quantity</code> can be changed after construction.

organization	a character string giving the desired organization of the returned <b>colorSpec</b> object. In case of 'auto', the organization is 'vector' or 'matrix' depending on data. Other possible organizations are 'df.col' or 'df.row'; see <a href="#">organization</a> for discussion of all 4 possible organizations. The organization can be changed after construction.
specnames	a character vector with length equal to the number of spectra, and with no duplicates. If specnames is NULL and data is a vector, then specnames is set to <code>deparse(substitute(data))</code> . If specnames is NULL and data is a matrix, then specnames is set to <code>colnames(data)</code> . If specnames is <i>still</i> not a character vector with the right length, or if there are duplicate names, then specnames is set to 'S1', 'S2', ... with a warning message. The specnames vector can be changed after construction.
x	an R object to test for being a valid <b>colorSpec</b> object.
...	arguments for use in other packages.

### Details

Under the hood, a **colorSpec** object is either a vector, matrix, or data.frame. It is of S3 class 'colorSpec' with these extra attributes:

wavelength a numeric vector of wavelengths for all the spectra. If the organization of the object is 'df.col', then this is absent.

quantity a character string that gives the physical quantity of all spectra, see [quantity\(\)](#) for a list of possible values.

metadata a named list for user-defined data. The names 'path', 'header' and 'date' are already reserved; see [metadata\(\)](#).

step.wl step between adjacent wavelengths in nm. This is assigned only when the wavelengths are regular; see [is.regular\(\)](#).

specname only assigned when the organization is 'vector', in which case it is equal to the single character string name of the single spectrum. Note that the word specname is singular. Also see [specnames\(\)](#).

And there are a few more attributes that exist only in special cases; see the **colorSpec User Guide**.

### Value

`colorSpec()` returns a **colorSpec** object, or NULL in case of ERROR. Compare this function with `stats::ts()`.

`is.colorSpec()` returns TRUE or FALSE. It does more than check the class, and also checks wavelength, quantity, and organization. If FALSE, it logs (at `loglevel='DEBUG'`) the reason that x is invalid.

`as.colorSpec.default()` issues an ERROR message and returns NULL

### See Also

[wavelength](#), [quantity](#), [organization](#), [metadata](#), [step.wl](#), [specnames](#), [is.regular](#), [coredata](#)



**Examples**

```
# make a synthetic Gaussian bandpass filter

center = 600
wave   = 400:700
trans  = exp( -(wave-center)^2 / 20^2 )

filter.bp = colorSpec( trans, wave, quantity='transmittance', specnames='myfilter' )

organization( filter.bp ) # returns: "vector"

# and now plot it
plot( filter.bp )
```

---

computeADL

*compute ADL coordinates by ray tracing*


---

**Description**

Consider a **colorSpec** object  $x$  with type equal to `responsivity.material`. The set of all possible material reflectance functions (or transmittance functions) is convex, closed, and bounded (in any reasonable function space), and this implies that the set of all possible output responses from  $x$  is also convex, closed, and bounded. The latter set is called the *object-color solid* or *Rösch Farbkörper* for  $x$ . A color on the boundary of the *object-color solid* is called an *optimal color*. The special points **W** (the response to the perfect reflecting diffuser) and **0** are on the boundary of this set. The interior of the line segment of neutrals joining **0** to **W** is in the interior of the *object-color solid*. It is natural to parameterize this segment from 0 to 1 (from **0** to **W**). The solid is symmetrical about the neutral gray midpoint  $\mathbf{G}=\mathbf{W}/2$ .

Now suppose that  $x$  has 3 spectra (3 responses) and consider a color response **R** not equal to **G**. There is a ray based at **G** and passing through **R** that intersects the boundary of the *object-color solid* at an *optimal color* **B** on the boundary with Logvinenko coordinates  $(\delta, \omega)$ . If these 2 coordinates are combined with  $\alpha$ , where  $\mathbf{R} = (1 - \alpha)\mathbf{G} + \alpha\mathbf{B}$ , it yields the *Logvinenko coordinates*  $(\alpha, \delta, \omega)$  of **R**. These coordinates are also denoted by ADL; see **References**. A response is in the *object-color solid* iff  $\alpha \leq 1$ . A response is *optimal* iff  $\alpha = 1$ .

The coordinates of **0** are  $(\alpha, \delta, \omega)=(1,0,0)$ . The coordinates of **W** are  $(\alpha, \delta, \omega)=(1,1,0)$ . The coordinates of **G** are undefined.

**Usage**

```
## S3 method for class 'colorSpec'
computeADL( x, response )
```

**Arguments**

<code>x</code>	a <b>colorSpec</b> object with type equal to <code>responsivity.material</code> and 3 spectra
<code>response</code>	a numeric $N \times 3$ matrix with responses in the rows, or a numeric vector that can be converted to such a matrix, by row.

**Details**

For each response, a ray is computed and the ray tracing is done by `probeOptimalColors()`.

**Value**

`computeADL()` returns a data frame with a row for each response. The columns in the data frame are:

response	the input response vector
ADL	the computed ADL coordinates of the response vector
omega	the reparameterized $\lambda$ in the interval [0,1]; see <b>References</b>
lambda	lambda.1 and lambda.2 at the 2 transitions, in nm. lambda.1 < lambda.2 => bandpass, and lambda.1 > lambda.2 => bandstop.

If an individual ray could not be traced, or if the optimal spectrum has more than 2 transitions, the row contains NA in appropriate columns.

In case of global error, the function returns NULL.

**WARNING**

Since this function is really a simple wrapper around `probeOptimalColors()`, please see the performance warnings there.

**References**

Logvinenko, A. D. An object-color space. **Journal of Vision**. 9(11):5, 1-23, (2009). <https://jov.arvojournals.org/article.aspx?articleid=2203976>. doi:10.1167/9.11.5.

Godau, Christoph and Brian Funt. XYZ to ADL: Calculating Logvinenko's Object Color Coordinates. Proceedings Eighteenth IS&T Color Imaging Conference. San Antonio. Nov 2009.

**See Also**

`type()`, `probeOptimalColors()`, vignette **optimals**

**Examples**

```
D50.eye = product( D50.5nm, 'varmat', xyz1931.1nm, wave=seq(360,830,by=5) )
computeADL( D50.eye, c(30,50,70) )
##   response.X response.Y response.Z   ADL.alpha   ADL.delta  ADL.lambda   omega
## 1         30         50         70  0.7371475  0.5384104 473.3594572 0.3008817

## lambda.1 lambda.2
## 427.2011 555.5261

## since alpha < 1, XYZ=c(30,50,70) is *inside* the object-color solid of D50.eye
```

---

computeCCT

*Compute Correlated Color Temperature (CCT) of Light Spectra*


---

## Description

Compute the CCT, in K, of a **colorSpec** object with type equal to 'light'

## Usage

```
## S3 method for class 'colorSpec'
computeCCT( x, isotherms='robertson', locus='robertson', strict=FALSE )
```

## Arguments

x	an <b>colorSpec</b> R object with type equal to 'light', and M spectra
isotherms	A character vector whose elements match one of the available isotherm families: 'robertson', 'mccamy', and 'native'. Matching is partial and case-insensitive. When more than one family is given, a matrix is returned, see <b>Value</b> . When isotherms='native' the isotherms are defined implicitly as lines perpendicular to the locus, see <b>Details</b> in <code>spacesXYZ::CCTfromXYZ()</code> . The character NA (NA_character_) is taken as a synonym for 'native'.
locus	valid values are 'robertson' and 'precision', see above. Matching is partial and case-insensitive.
strict	The CIE considers the CCT of a chromaticity uv to be meaningful only if the distance from uv to the Planckian locus is less than or equal to 0.05 [in CIE UCS 1960]. If strict=FALSE, then this condition is ignored. Otherwise, the distance is computed along the corresponding isotherm, and if it exceeds 0.05 the returned CCT is set to NA.

## Details

In `computeCCT()`, for each spectrum, XYZ is computed using `xyz1931.1nm`, and the result passed to `spacesXYZ::CCTfromXYZ()`. If the quantity of x is 'photons' (actinometric) each spectrum is converted to 'energy' (radiometric) on the fly.

## Value

`computeCCT()` returns a numeric vector of length M, where M is the number of spectra in x. The vector's names is set to `specnames(x)`.  
If the type of x is not 'light', then a warning is issued and all values are NA\_real\_.

## References

- McCamy, C. S. *Correlated color temperature as an explicit function of chromaticity coordinates*. Color Research & Application. Volume 17. Issue 2. pages 142-144. April 1992.
- Robertson, A. R. Computation of correlated color temperature and distribution temperature. Journal of the Optical Society of America. 58. pp. 1528-1535 (1968).
- Wyszecki, Günther and W. S. Stiles. **Color Science: Concepts and Methods, Quantitative Data and Formulae, Second Edition**. John Wiley & Sons, 1982. Table 1(3.11). pp. 227-228.

## See Also

[type\(\)](#), [quantity\(\)](#), [xyz1931](#), [planckSpectra\(\)](#), [specnames\(\)](#), [spacesXYZ::CCTfromXYZ\(\)](#)

## Examples

```
computeCCT( D65.1nm )                # returns 6502.068
computeCCT( D65.1nm, isotherms='native' ) # returns 6503.323
computeCCT( A.1nm )                  # returns 2855.656
computeCCT( A.1nm, isotherms='native' ) # returns 2855.662
computeCCT( A.1nm, isotherms='mccamy' ) # returns 2857.188

moon = readSpectra( system.file( "extdata/sources/moonlight.txt", package='colorSpec' ) )
computeCCT( moon )                    # returns 4482.371
```

---

computeCRI

*Compute Color Rendering Index (CRI) of Light Spectra*

---

## Description

Compute the CIE 1974 color rendering index (CRI) of a light spectrum, called the *the test illuminant*.

From the given spectrum a *reference illuminant* is selected with the same CCT (Correlated Color Temperature). A selected set of 8 color samples is rendered in XYZ (1931) with both illuminants and 8 color differences are computed in a special **CIEUVW** color space. For each color difference a CRI is computed, where 100 is a perfect color match. The final CRI is the average of these 8 CRI values.

## Usage

```
## S3 method for class 'colorSpec'
computeCRI( x, adapt=TRUE, attach=FALSE, tol=5.4e-3 )
```

## Arguments

**x** an **colorSpec** R object with type equal to 'light', and exactly 1 spectrum

**adapt** if TRUE, then a special chromatic adaption is performed, see **Details**

attach	if TRUE, then a large list of intermediate calculations is attached to the returned number, as attribute data. This attached list includes data for all special 14 color samples, although the last 6 do not affect the returned CRI.
tol	for the CRI to be meaningful the chromaticities of the test and reference illuminants must be sufficiently close in the CIE 1960 uniform chromaticity space. If the tolerance is exceeded, the function returns NA. The default tol=5.4e-3 is the one recommended by the CIE, but the argument allows the user to override it.

### Details

The CCT of  $x$  is computed by `computeCCT()` with default options.

If `adapt` is TRUE the 8 test uv points are chromatically adapted from the test illuminant to the reference illuminant using a special von Kries type transformation; see *Oleari* and *Wikipedia*. The test UVW values are computed relative to the reference illuminant.

If `adapt` is FALSE the 8 test uv points are *not* chromatically adapted, and the test UVW values are computed relative to the test illuminant.

### Value

`computeCRI()` returns a single number  $\leq 100$ . In case of ERROR it returns NA. If `attach` is TRUE a large list of intermediate calculations is attached to the returned number.

### Source

The test color reflectance spectra are taken from:

[http://www.lrc.rpi.edu/programs/nlpip/lightinganswers/lightsources/scripts/NLPIP\\_LightSourceColor\\_Script.m](http://www.lrc.rpi.edu/programs/nlpip/lightinganswers/lightsources/scripts/NLPIP_LightSourceColor_Script.m)

### References

Oleari, Claudio, Gabriele Simone. **Standard Colorimetry: Definitions, Algorithms and Software**. John Wiley. 2016. pp. 465-470.

Günther Wyszecki and W. S. Stiles. **Color Science: Concepts and Methods, Quantitative Data and Formulae, Second Edition**. John Wiley & Sons, 1982. Table 1(3.11). p. 828.

Wikipedia. **Color rendering index**. [http://en.wikipedia.org/wiki/Color\\_rendering\\_index](http://en.wikipedia.org/wiki/Color_rendering_index)

Hunt, R. W. G. and M. R. Pointer. **Measuring Colour**. 4th edition. John Wiley & Sons. 2011. Appendix 7.

### See Also

[type](#), [xyz1931](#), [computeCCT](#)

### Examples

```
computeCRI( subset(Fs.5nm, 'F2') )      # returns 64.15195
computeCRI( subset(Fs.5nm, 'F4') )      # returns 51.36348
```

---

 computeSSI

*Compute the Spectrum Similarity Index of light spectra*


---

### Description

Compute the Spectrum Similarity Index (SSI), an index between 0 and 100, of a **colorSpec** object with type equal to 'light'. It compares a test spectrum with a reference spectrum (an ideal). The value 100 means a perfect match to the reference, and a smaller value mean a poorer match (similar to CRI). Only values in the interval [375,675] nm are used; for details see *Holm*.

### Usage

```
## S3 method for class 'colorSpec'
computeSSI( x, reference=NULL, digits=0, isotherms='mccamy', locus='robertson' )
```

### Arguments

x	a <b>colorSpec</b> object with type equal to 'light', and M test spectra
reference	a <b>colorSpec</b> object with type equal to 'light', and either 1 or M reference spectra. reference can also be NULL (the default), which means to generate each reference spectrum from the corresponding test spectrum.
digits	the number of digits after the decimal point in the returned vector. According to <i>Holm</i> the output should be rounded to the nearest integer, which corresponds to digits=0. To return full precision, set digits=Inf.
isotherms	this is only used when reference=NULL. It is passed to <code>computeCCT()</code> in order to compute the CCT of each test spectrum.
locus	this is only used when reference=NULL. It is passed to <code>computeCCT()</code> in order to compute the CCT of each test spectrum.

### Details

If reference contains a single spectrum, then each test spectrum is compared to that one reference spectrum. If reference contains M spectra, then the i'th test spectrum is compared to the i'th reference spectrum.

If reference=NULL then for each test spectrum the CCT is computed and used to compute a reference spectrum with the same CCT. It is either a Planckian (black-body) or daylight illuminant, see *Holm* for details. The test spectrum and auto-computed reference spectrum are then compared.

### Value

computeSSI() returns a numeric vector of length M, where M is the number of spectra in x. The vector's names is set from specnames(x) and a compact code for the corresponding reference spectrum.

If the type of x is not 'light', or reference is not valid, then the function returns NULL.

## References

J. Holm and T. Maier and P. Debevec and C. LeGendre and J. Pines and J. Erland and G. Joblove and S. Dyer and B. Sloan and J. di Gennaro and D. Sherlock. *A Cinematographic Spectral Similarity Index*. SMPTE 2016 Annual Technical Conference and Exhibition. pp. 1-36. (2016).

## See Also

[type\(\)](#), [computeCCT\(\)](#), [planckSpectra\(\)](#), [daylightSpectra\(\)](#), [specnames\(\)](#)

## Examples

```
computeSSI( planckSpectra( 1000*(2:6) ) )
## P2000_SSI[2027K] P3000_SSI[3057K] P4000_SSI[D4063] P5000_SSI[D5061] P6000_SSI[D6063]
##          99          98          93          92          92
```

---

convolvewith	<i>Convolve each spectrum in a colorSpec object with a kernel</i>
--------------	---

---

## Description

This function convolves each spectrum in a colorSpec object with a kernel of odd length. Its primary purpose is to correct raw spectrometer data (with positive instrumental bandwidth) to have bandwidth=0. Two popular correction kernels for this, with lengths 3 and 5, are built-in options, see **Details**.

## Usage

```
## S3 method for class 'colorSpec'
convolvewith( x, coeff )
```

## Arguments

x	a <b>colorSpec</b> object with N wavelengths
coeff	a convolution kernel of odd length. The center entry of this vector is taken as index 0 in the convolution. coeff can also be the string 'SS3' which means to apply the Stearns&Stearns bandwidth correction kernel $\text{coeff} = c(-1, 14, -1)/12$ , see <b>Details</b> . coeff can also be the string 'SS5' which means to apply the Stearns&Stearns bandwidth correction kernel $\text{coeff} = c(1, -12, 120, -12, 1)/98$ , see <b>Details</b> .

## Details

The built-in kernels, 'SS3' and 'SS5', were derived by *Stearns & Stearns* under specific hypotheses on the spectrometer profile, bandpass, and pitch; see **References**.

Missing values at both ends are filled by copying from the nearest valid value.

The function creates a function calling `stats::filter()` and passes that function to `applySpec()`.

**Value**

a **colorSpec** object with the same dimensions, [wavelength](#), [quantity](#), and [organization](#). If `coeff` is invalid it is an **ERROR** and `convolvewith()` returns **NULL**.

**References**

Stearns, E.I., Stearns R.E. An example of a method for correction radiance data for bandpass error. *Color Research and Application*. 13-4. 257-259. 1988.

Schanda, Janos. CIE Colorimetry, in *Colorimetry: Understanding the CIE System*. Wiley Interscience. 2007. p. 124.

Oleari, Claudio, Gabriele Simone. *Standard Colorimetry: Definitions, Algorithms and Software*. John Wiley. 2016. p. 309.

**See Also**

[quantity](#), [wavelength](#), [linearize](#), [appliespec](#), [organization](#)

---

coredata

*Extract the Core Data of a colorSpec Object*

---

**Description**

functions for extracting the core data contained in a **colorSpec** object.

**Usage**

```
## S3 method for class 'colorSpec'
coredata( x, forcemat=FALSE )
```

```
## S3 method for class 'colorSpec'
as.matrix( x, ... )
```

**Arguments**

<code>x</code>	a <b>colorSpec</b> object
<code>forcemat</code>	if <code>x</code> has only 1 spectrum, return a matrix with 1 column instead of a vector
<code>...</code>	extra arguments ignored

**Value**

<code>coredata</code>	If <code>x</code> has <code>organization</code> equal to <code>'vector'</code> then it returns <code>x</code> , unless <code>forcemat</code> is <b>TRUE</b> when it returns <code>x</code> as a matrix with 1 column. If <code>x</code> has any other <code>organization</code> then it returns a matrix with spectra in the columns. All of the <b>colorSpec</b> attributes are stripped except the column names, and the row names are set to <code>as.character(wavelength(x))</code> .
<code>as.matrix</code>	a wrapper for <code>coredata</code> with <code>forcemat=TRUE</code>



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

---

`cs.options`*Functions to set and retrieve colorSpec package options*

---

**Description**

**colorSpec** has a few options. The options are stored in the **R** global list and they are:

`colorSpec.loglevel`, `colorSpec.logformat`, and `colorSpec.stoponerror`

For details on what they do see [logging](#).

They can be set using the built-in function [options\(\)](#). When **R** starts up, an option can be set using a call to `options()` in the file **Rprofile.site**. If **colorSpec** is later loaded, the value of the option will not be changed. If an option has not been assigned, then it is created with a default value.

The function `cs.options()` makes setting the options a little easier in a few ways:

- it automatically prepends the string 'colorSpec.'
- partial matching of the option name is enabled
- a warning is issued when the option value has the wrong type

**Usage**

```
cs.options( ... )
```

**Arguments**

...            named arguments are set; unnamed arguments are ignored with a warning. See **Examples**.

**Value**

returns a list with all the **colorSpec** options.

**See Also**[logging](#), [options](#)**Examples**

```
cs.options( loglevel="DEBUG", stop=FALSE ) # 'stop' partially matches 'stoponerror'  
cs.options( stop='TRUE' )                 # warns that value has the wrong type  
cs.options( stop=FALSE, "DEBUG" )         # warns that the 2nd argument has no name  
cs.options( loglevel="WARN" )             # back to default
```

D50

*Standard Illuminant D50 (1964)***Description**

D50.5nm standard Illuminant D50, from 300 to 830 nm at 5 nm intervals.

**Format**

A **colorSpec** object organized as a vector, with 107 data points and **specnames** equal to 'D50'.

**Details**

This spectrum is not copied from a table from a CIE publication, though it does match such a table. It is computed using the function `daylightSpectra()` by following the special CIE recipe given in the **References**. The temperature is set to  $(14388/14380) * 5000 = 5002.781$  Kelvin. The coefficients of the daylight components  $S_0$ ,  $S_1$ , and  $S_2$  are rounded to 3 decimal places. This linear combination is computed at 10nm intervals and then linearly interpolated to 5nm intervals. The result is normalized to value 1 at 560nm (instead of the usual 100), and finally rounded to 5 decimal places. See **Examples**.

**References**

Günther Wyszecki and W.S. Stiles. **Color Science : Concepts and Methods, Quantitative Data and Formulae**. Second Edition. Wiley-Interscience. 1982. Table I(3.3.4) pp. 754-758

CIE 15: Technical Report: Colorimetry, 3rd edition. CIE 15:2004. Table T.1, pp 30-32, and Note 5 on page 69.

Schanda, Janos. CIE Colorimetry, in *Colorimetry: Understanding the CIE System*. Wiley Interscience. 2007. p. 42.

**See Also**

[ABC](#) , [D65](#) , [daylightSpectra](#)

**Examples**

```
# the CIE recipe for computing D50.5nm
correction = 14388 / 14380 # note 5, page 69 in CIE 15:2004
D50.10nm = daylightSpectra( correction*5000, wavelength=seq(300,830,by=10), roundMs=TRUE )
D50.5nm = resample( D50.10nm, seq(300,830,by=5), method='linear' )
D50.5nm = round( D50.5nm, 5 )

summary( D50.5nm )
white.point = product( D50.5nm, xyz1931.1nm, wave='auto' )
```

---

D65 *Standard Illuminant D65 (1964)*

---

### Description

D65.1nm standard Illuminant D65, 300 to 830 nm at 1 nm intervals  
 D65.5nm standard Illuminant D65, 380 to 780 nm at 5 nm intervals

### Format

Each is a **colorSpec** object organized as a vector, with `specnames` equal to 'D65'.

### Details

Both of these have been divided by 100, to make the values at 560nm equal to 1 instead of 100.

### Source

<http://www.cvrl.org>

### References

Günther Wyszecki and W.S. Stiles. **Color Science : Concepts and Methods, Quantitative Data and Formulae**. Second Edition. Wiley-Interscience. 1982. Table I(3.3.4) pp. 754-758

ASTM E 308-01. Standard Practice for Computing the Colors of Objects by Using the CIE System. Table 3. pages 3-4.

### See Also

[ABC](#), [D50](#), [daylightSpectra](#), [daylight](#)

### Examples

```
summary( D65.1nm )
white.point = product( D65.1nm, xyz1931.1nm, wave='auto' )
```

---

daylight *Standard Daylight Components*

---

### Description

daylight1964 spectral components  $S_0, S_1, S_2$ ; from 300 to 830 nm at 5 nm intervals  
 daylight2013 smoothed spectral components  $S_0, S_1, S_2$ ; from 300 to 830 nm at 1 nm intervals

## Format

Each is a **colorSpec** object organized as a matrix with 3 columns

S0 component 0, the mean power spectrum  
S1 component 1, the 1st characteristic spectrum  
S2 component 2, the 2nd characteristic spectrum

## Source

[http://www.cie.co.at/publ/abst/datatables15\\_2004/CIE\\_sel\\_colorimetric\\_tables.xls](http://www.cie.co.at/publ/abst/datatables15_2004/CIE_sel_colorimetric_tables.xls)

<http://vision.vein.hu/~schanda/CIE%20TC1-74/>

## References

Günther Wyszecki and W.S. Stiles. **Color Science : Concepts and Methods, Quantitative Data and Formulae**. Second Edition. Wiley-Interscience. 1982. Table V(3.3.4) p. 762.

Smoothing spectral power distribution of daylights. Zsolt Kosztyan and Janos Schanda. *Color Research & Application*. Volume 38, Issue 5, pages 316-321, October 2013.

CIE 15: Technical Report: Colorimetry, 3rd edition. CIE 15:2004. Table T.2, pp 33-34

JUDD, D.B., MACADAM, D.L. and WYSZECKI, G., with the collaboration of BUDDE, H.W, CONDIT, H.R, HENDERSON, S.T, and SIMONDS, J.L. Spectral distribution of typical daylight as a function of correlated color temperature. *J Opt. Soc. Am.* 54, 1031-1040, 1964.

Zsolt Kosztyan and Janos Schanda. Smoothing spectral power distribution of daylights. *Color Research & Application*. Volume 38, Issue 5, pages 316-321, October 2013.

## See Also

[D65](#), [D50](#), [daylightSpectra](#)

## Examples

```
summary( daylight1964 )
day1964 = daylightSpectra( c(5000,6500), comp=daylight1964 )
day2013 = daylightSpectra( c(5000,6500), comp=daylight2013 )

plot( day1964, col='black' )
plot( day2013, col='black', add=TRUE )
```

---

DisplayRGB

---

*Compute Display RGB from Linear RGB*


---

### Description

All RGB displays have a non-linear "gamma function" of some sort. This function converts from linear RGB to an RGB appropriate for the gamma function of the display; which is also called the *electro-optical conversion function* (EOCF).

### Usage

```
DisplayRGBfromLinearRGB( RGB, gamma='sRGB' )
```

### Arguments

RGB	linear RGB values organized as a vector or matrix of any size; all 3 channels are treated the same way so size does not matter
gamma	either the string 'sRGB' or a positive number giving the gamma of the display.

### Value

The function first clamps the input RGB to the interval [0,1]. If gamma='sRGB' (not case-sensitive) it then maps [0,1] to [0,1] using the special piecewise-defined sRGB function, see *Wikipedia*. In case gamma is a positive number, the function raises all values to the power 1/gamma. The dimensions and names of the input are copied to the output.

In case of error, the function returns the clamped input values.

### WARNING

This function is deprecated. New software should use `spacesRGB::SignalRGBfromLinearRGB()` instead.

### Source

Wikipedia. **sRGB**. <http://en.wikipedia.org/wiki/sRGB>

### See Also

[RGBfromXYZ](#)

### Examples

```
DisplayRGBfromLinearRGB( c(0.2, 0.5) )
# [1] 0.4845292 0.7353570      # this is display sRGB, in [0,1]

DisplayRGBfromLinearRGB( c(-0.1, 0.2, 0.5, 1), 2.2 )
# [1] 0.0000000 0.4811565 0.7297401 1.0000000      # gamma=2.2
```

```
x = seq( 0, 1, len=101)
plot( x, DisplayRGBfromLinearRGB(x), type='l' )
```

---

emulate	<i>modify a colorSpec responder to emulate (approximate) another responder</i>
---------	--

---

### Description

The two possible modifications are:

- pre-multiplication by a transmitting filter
- post-multiplication by a matrix

Both of these are optional. If neither of these modifications is enabled, the original `x` is returned.

### Usage

```
## S3 method for class 'colorSpec'
emulate( x, y, filter=FALSE, matrix=TRUE )
```

### Arguments

<code>x</code>	a <b>colorSpec</b> responder with <code>M</code> spectra, to be modified. The type must be 'responsivity.light' or 'responsivity.material'
<code>y</code>	a <b>colorSpec</b> responder with <code>N</code> spectra, to be emulated by a modified <code>x</code> . It must have the same type and wavelength vector as <code>x</code>
<code>filter</code>	enable filter pre-multiplication.
<code>matrix</code>	enable matrix post-multiplication. If <code>matrix=TRUE</code> then the computed matrix <code>A</code> is <code>MxN</code> .

### Details

If `filter=FALSE` and `matrix=TRUE` then the returned value is `multiply(x,A)`, where the matrix `A` is computed to minimize the difference with `y`, in the least squares sense (Frobenius matrix norm). The function `MASS::ginv()` is used here.

If `filter=TRUE` and `matrix=FALSE` then the returned value is `product(filter,x)`, where the object `filter` is computed to minimize the difference with `y`, in the least squares sense (Frobenius matrix norm). This calculation is fairly straightforward, but requires that the responsivity of `x` does not vanish at any wavelength. It also requires that `M=N`. The computed filter may be unrealistic, i.e. the transmittance may be  $> 1$ . If this happens a WARN message is issued.

If `filter=TRUE` and `matrix=TRUE` then the returned value is `product(filter,multiply(x,A))`, where `(filter,A)` are chosen with the above minimization criterion. If `N=1` then we must have `M=1`

as well; the calculation is trivial and the emulation is exact. If  $N \geq 2$ , the calculation is iterative - solving alternatively for `filter` and `A` until convergence. The function `MASS::ginv()` is used on each iteration. This is a bilinear optimization. If convergence fails, it is an error and the function returns `NULL`. If convergence succeeds, there is 1 degree of freedom in the `(filter,A)` pair. If one is scaled by a positive constant, the other can be scaled by the inverse, and the returned object is the same. The filter is scaled so the maximum transmittance is 1.

If `filter=FALSE` and `matrix=FALSE` then the original `x` is returned, with a `WARN` message.

### Value

a `colorSpec` object close to `y`, as in **Details**. The `quantity` is the same as `y`. The `specnames()` are the same as those of `y`, except that `".em"` is appended to each one. The function attaches attribute `"emulate"`, whose value is a list containing `filter` and/or `A` as appropriate.

### Examples

see the vignette **Emulation of one Camera by another Camera**

### See Also

[wavelength](#), [type](#), [quantity](#), [multiply](#), [product](#), [specnames](#)

---

extradata	<i>extradata of a colorSpec object</i>
-----------	--

---

### Description

Retrieve or set the extradata of a `colorSpec` object.

### Usage

```
## S3 method for class 'colorSpec'
extradata(x)

## S3 replacement method for class 'colorSpec'
extradata(x,add=FALSE) <- value
```

### Arguments

<code>x</code>	a <code>colorSpec</code> object with <code>M</code> spectra
<code>value</code>	a <code>data.frame</code> with <code>M</code> rows. It is OK for <code>value</code> to have 0 columns, and <code>value</code> can also be <code>NULL</code> ; see <code>add</code> .
<code>add</code>	If <code>add</code> is <code>FALSE</code> , then any existing extradata is discarded and replaced by <code>value</code> , except when <code>value</code> is <code>NULL</code> when <code>x</code> is left with no extradata. If <code>add</code> is <code>TRUE</code> , then <code>value</code> is appended to the existing extradata, except when <code>value</code> is <code>NULL</code> when <code>x</code> is left unchanged.

**Details**

If the organization of `x` is not `'df.row'`, then `extradata` cannot be stored in `x` and the assignment is ignored, with a warning. First change the `organization` to `'df.row'`, and *then* assign the `extradata`.

If the organization of `x` is `'df.row'`, but value does not have the right number of rows, the assignment is ignored, with a warning.

**Value**

`extradata(x)` returns a `data.frame` with `M` rows, where `M` is the number of spectra in `x`. The rownames are set to the `specnames` of `x`. If there is no extra data then the number of columns in this `data.frame` is 0.

**Note**

Do not confuse `extradata` and `metadata`.

`metadata` is unstructured data that is attached to the entire `colorSpec` object. `extradata` is structured data, with a row of data for each spectrum in the object.

**See Also**

[metadata](#), [organization](#)

---

F96T12

*Photon Irradiance of F96T12 Fluorescent Bulb*

---

**Description**

F96T12

Sylvania F96T12 CW/VHO 215-Watt fluorescent bulb photon irradiance, measured with a LI-COR LI-1800 spectroradiometer, from 300 to 900 nm at 1 nm intervals.

**Format**

A `colorSpec` object organized as a vector, with 601 data points and `specnames` equal to `'F96T12'`.

**Details**

The unit is  $(\mu\text{mole of photons}) * \text{sec}^{-1} * \text{m}^{-2} * \text{nm}^{-1}$ .

**Source**

Pedro J. Aphalo. <http://www.mv.helsinki.fi/aphalo/photobio/lamps.html>

**See Also**

[ABC](#), [D65](#), [daylightSpectra](#)



### Examples

```
sum( F96T12 )
# [1] 320.1132  photon irradiance, (micromoles of photons)*m^{-2}

sum( radiometric(F96T12) )
# [1] 68.91819  irradiance, watts*m^{-2}
```

---

Flea2.RGB

*Flea2 Camera FL2-14S3C from Point Grey*

---

### Description

Flea2.RGB an RGB responder to light, from 360 to 800 nm at 10 nm intervals

### Format

A **colorSpec** object with quantity equal to 'energy->electrical' and 3 spectra: Red, Green, and Blue.

### Details

This data is read from the file **Flea2-spectral.txt** which was digitized from the plot in **Flea2-spectral.png**.

### Source

<http://www.ptgrey.com/support/downloads/10117>

### See Also

[quantity](#), vignette [blueflame](#)

### Examples

```
# Make a scanner from a tungsten source and a Flea2 camera
Flea2.scanner = product( A.1nm, "VARMATERIAL", Flea2.RGB, wavelength=420:680 )
Flea2.scanner = calibrate( Flea2.scanner )
```

---

Fluorescents	<i>Standard series F Illuminants F1, F2, F3, F4, F5, F6, F7, F8, F9, F10, F11, and F12</i>
--------------	--

---

**Description**

Fs.5nm contains 12 CIE Fluorescent Illuminants, from 380 to 780 nm, at 5nm intervals.

**Format**

Fs.5nm is a **colorSpec** object with 12 spectra. It is organized as a data frame with `quantity` equal to "energy".

**Note**

The series F illuminants do not seem to be normalized in a consistent way.

**Source**

<http://www.rit-mcsl.org/UsefulData/Fluorescents.htm>

**See Also**

[ABC](#), [D50](#), [D65](#)

**Examples**

```
# plot only F4
plot( subset(Fs.5nm, "F4") )
```

---

HigherPasserines	<i>Cone Fundamentals for the Higher Passerines</i>
------------------	--

---

**Description**

HigherPasserines Tetrachromatic Cone Fundamentals of Higher Passerine Birds

**Format**

A **colorSpec** object organized as a matrix with the 4 spectra:

UV            the UV wavelength responsivity



**Source**

<http://www.hoyaoptics.com/>

**See Also**

[quantity](#)

**Examples**

```
# compute response of ACES scanner to the Hoya filters
product( Hoya, scanner.ACES, wave='auto' )
```

---

interpolate

*interpolate spectra*

---

**Description**

interpolate along a 1-parameter path of spectra

**Usage**

```
## S3 method for class 'colorSpec'
interpolate( x, p, pout, pname=deparse(substitute(p)) )
```

**Arguments**

x	a <b>colorSpec</b> object, typically with multiple spectra
p	a numeric vector with length(p)==numSpectra(x). The value p[i] is associated with the i'th spectrum in x.
pout	a numeric vector of parameter values at which interpolation of the spectra in x take place
pname	the name of the parameter p

**Details**

Each spectrum in x can be thought of as a point in a high-dimensional space, and each point has a real-valued parameter associated with it. The function performs natural spline interpolation on these points, one coordinate at a time. For each wavelength value it calls [spline](#) with method='natural'.

**Value**

interpolate(x) returns a **colorSpec** object y with a spectrum for each value in pout. The organization of y is 'df.row', and extradata(y) has a single column which is a copy of pout. The name of the column is pname. The names in specnames(y) are <pname>=<pout>. Other properties of y, e.g. [wavelength](#), [quantity](#), ..., are the same as x. In case of ERROR, the function returns NULL.

**See Also**

[organization](#), [wavelength](#), [extradata](#), [spline](#)

**Examples**

```
path = system.file( "extdata/stains/PhenolRed-Fig7.txt", package="colorSpec" )
wave = 350:650
phenolred = readSpectra( path, wavelength=wave )
pH = as.numeric( sub( '^[^0-9]+([0-9]+)$', '\\1', specnames(phenolred) ) )
pHvec = seq(min(pH),max(pH),by=0.05)
phenolinterp = interpolate( phenolred, pH, pHvec )
```

---

invert	<i>estimate spectra from responses, effectively inverting the operator from spectrum to response</i>
--------	--

---

**Description**

Given a light responder (e.g. an eye or a camera), two light spectra that produce the same response from the responder are called *metamers* for that responder. Similarly, given a material responder (e.g. a scanner), two materials that produce the same response from the responder are called *metamers* for that responder.

For a given responder and response, there are typically infinitely many *metamers*. The goal of the function `invert()` is to calculate a "good" metamer from a response. *Koenderink* calls this topic *inverse colorimetry*. In the case that the estimated spectrum is a reflectance spectrum, the topic is often called *reflectance estimation* or *reflectance recovery*, see *Bianco*.

The *centroid method*, which is the default and the featured method in this package, computes the centroid of the set of all the metamers (if any). The centroid is computed in an infinite-dimensional context and is expounded further in *Davis*. The second available method is the Hawkyard method, see *Hawkyard* and *Bianco*. The methods are quite similar.

The first argument to `invert()` is the responder `x`, and the second is the matrix response of responses (e.g. XYZs).

The goal is to return a "good" spectrum for each response so that:

$$\text{product}( \text{invert}(x, \text{response}), x ) \cong \text{response}$$

The error is returned as column `estim.precis`, see below.

First consider the case where `x` has type `type='responsivity.material'`. If the response is on the object-color boundary (an *optimal color*) or outside the boundary, the centroid method fails and the Hawkyard method returns a spectrum with an error that may be significant. The centroid method *may* also fail if the response is inside the object-color solid (the *Rösch Farbkörper*) and very close to the boundary.

Now consider the case where `x` has type `type='responsivity.light'`. The centroid method will fail if the response is too far from the response to Illuminant E. The Hawkyard method does not work in this case.

Equalization is controlled by the argument `alpha` and is enabled by default, see below. This means that if the response comes from a constant spectrum (a perfectly neutral gray material, or a multiple of Illuminant E), then the returned spectrum is that same constant spectrum (up to numerical precision). It is a complicated mechanism, for details see *Davis*.

NOTE: If the responder has only one output channel (e.g. a monochrome camera) and equalization is enabled, then *all* responses are inverted to a constant spectrum. This may or may not be desirable.

### Usage

```
## S3 method for class 'colorSpec'
invert( x, response, method='centroid', alpha=1 )
```

### Arguments

<code>x</code>	a <b>colorSpec</b> object with type equal to <code>'responsivity.material'</code> or <code>'responsivity.light'</code> and <code>M</code> responsivities
<code>response</code>	a numeric <code>NxM</code> matrix, or a numeric vector that can be converted to such matrix, by row. The <code>N</code> responses are contained in the rows. The <code>rownames(response)</code> are copied to the output <code>specnames</code> .
<code>method</code>	either <code>'centroid'</code> or <code>'Hawkyard'</code> . <code>'Hawkyard'</code> is only valid when <code>type(x)</code> is <code>'responsivity.material'</code> . Matching is partial and case-insensitive.
<code>alpha</code>	a vector of <code>M</code> weighting coefficients, or a single number that is replicated to length <code>M</code> . When <code>method='centroid'</code> , <code>alpha</code> is used for <i>equalizing</i> the responsivities, which is recommended. For <code>alpha</code> to be valid, the linear combination of the <code>M</code> responsivities, with coefficients <code>alpha</code> , must be positive. To disable equalization (not recommended) and use the original responsivities, set <code>alpha=NULL</code> . Similarly, when <code>method='Hawkyard'</code> , <code>alpha</code> is used for equalizing the responsivities, which is also recommended.

### Details

When `method='centroid'` the function calls the non-linear root-finder `rootSolve::multiroot()`.

### Value

If `type(x)='responsivity.material'` it returns a **colorSpec** object with `type = 'material'` (`quantity = 'reflectance'`).

If `type(x)='responsivity.light'` it returns a **colorSpec** object with `type = 'light'` (`quantity='energy'` or `quantity='photons'` depending on `quantity(x)`).

In either case, the returned object has `organization = 'df.row'` and the `extradata` is a `data.frame` with these columns:

<code>response</code>	the input matrix of desired responses
<code>estim.precis</code>	the difference between the desired response and actual response. It is the mean of the absolute value of the differences. See <code>rootSolve::multiroot()</code>
<code>time.msec</code>	the time to compute the spectrum, in msec. When <code>method='Hawkyard'</code> , all <code>N</code> spectra are computed at once, so all <code>N</code> spectra are assigned the same mean time.

iters	the number of iterations that were required to find the relevant root. This is present only when method='centroid'.
clamped	a logical indicating whether the reflectance was clamped to [0,1]. This is present only when method='Hawkyard'.

If a response could not be estimated, the row contains NA in appropriate columns.

In case of global error, the function returns NULL.

### Known Issues

If type(x)='responsivity.light' the centroid method may fail (not converge) if the response is too far from that of Illuminant E.

### References

Davis, Glenn. A Centroid for Sections of a Cube in a Function Space, with application to Colorimetry. <https://arxiv.org/abs/1811.00990>. [math.FA]. 2018.

Bianco, Simone. Reflectance spectra recovery from tristimulus values by adaptive estimation with metameric shape correction. vol. 27, no 8. *Journal of the Optical Society of America A*. pages 1868-1877. 2010 <https://www.osapublishing.org/josaa/abstract.cfm?uri=josaa-27-8-1868>.

Hawkyard, C. J. Synthetic reflectance curves by additive mixing. *Journal of the Society of Dyers and Colourists*. vol. 109. no. 10. Blackwell Publishing Ltd. pp. 323-329. 1993.

Koenderink, J.J. Color for the Sciences. MIT Press. 2010.

### See Also

[type\(\)](#), [quantity\(\)](#), [organization\(\)](#), [specnames\(\)](#), [product\(\)](#), [extradata\(\)](#), [rootSolve::multiroot\(\)](#), vignette **inversion**

### Examples

```

wave = 400:700
E.eye = product( illuminantE(1,wave), "material", xyz1931.1nm, wavelength=wave )
path = system.file( 'extdata/targets/CC_Avg30_spectrum_CGATS.txt', package='colorSpec' )
MacbethCC = readSpectra( path, wavelength=wave )
XYZ = product( MacbethCC, E.eye, wavelength=wave )
est.eq = invert( E.eye, XYZ, method='centroid', alpha=1 )
extra = extradata(est.eq)
range(extra$estim.precis)      # prints 0.000000e+00 3.191741e-08

```

---

lightResponsivitySpectra

*compute standard light responsivity spectra*

---

### Description

Some action spectra standards are defined by simple equations; the erythemal spectrum for human sunburn is one of them.

### Usage

```
erythemalSpectrum( wavelength=250:400 )
```

### Arguments

wavelength      a vector of wavelengths, in nm

### Details

This erythemal spectrum is defined in 4 pieces:  $\lambda \leq 298$ ,  $298 \leq \lambda \leq 328$ ,  $328 \leq \lambda \leq 400$ , and  $400 < \lambda$ . The unit is nm. The spectrum is used in the definition of the international standard **UV Index**.

### Value

For erythemalSpectrum()

A **colorSpec** object with **quantity** equal to 'energy->action'. The responsivity is 0 for  $\lambda > 400$  nm, so this putting this spectrum in the category of human vision is a bit of a stretch.

### Source

[http://en.wikipedia.org/wiki/Ultraviolet\\_index](http://en.wikipedia.org/wiki/Ultraviolet_index)

### References

McKinlay, A.F., and B.L. Diffey. A reference action spectrum for ultraviolet induced erythema in human skin. CIE Res. Note, 6(1), 17-22. (1987)

### See Also

[daylight](#), [quantity](#), [materialSpectra](#), [lightSpectra](#)



---

LightSpectra                      *compute standard light spectra*

---

### Description

Two families of standard illuminants that are parameterized by temperature are the Planckian spectra (black-body spectra), and daylight spectra. For the daylight spectra, a smoothed version is available. Illuminant E, a third and trivial spectrum, is also available.

### Usage

```
planckSpectra( temperature, wavelength=300:830, normalize=TRUE, c2=1.4388e-2 )
```

```
daylightSpectra( temperature, wavelength=NULL,
                 components=colorSpec::daylight1964, roundMs=FALSE )
```

```
illuminantE( energy=1, wavelength=380:780 )
```

### Arguments

temperature	a vector of temperatures, in Kelvin
wavelength	a vector of wavelengths. For <code>planckSpectra()</code> and <code>illuminantE()</code> this is required. For <code>daylightSpectra()</code> this is optional. The default <code>wavelength=NULL</code> means to use the wavelengths in <code>components</code> , and otherwise <code>components</code> is resampled at the given wavelength vector.
normalize	a logical value. If <code>TRUE</code> the Planck spectra are normalized to have value 1 at 560nm. If <code>FALSE</code> then the quantity returned is radiant exitance with unit $W * m^{-2} * nm^{-1}$ .
c2	the value of $hc/k$ in Planck's law. $h$ is the Planck constant; $c$ is the speed of light in $m/sec$ ; and $k$ is the Boltzmann constant. The default value of $1.4388e-2 m * K$ was recommended by the CIE in 2005; in 1986 the CIE recommended $c2=1.438e-2$ . If <code>c2='calc'</code> then <code>c2</code> is calculated directly from the 3 physical constants, as recommended by CODATA 2014.
components	a <b>colorSpec</b> object with the daylight components $S_0$ , $S_1$ , and $S_2$ . The default is <a href="#">daylight1964</a> and a smoothed version <a href="#">daylight2013</a> is also available.
roundMs	a logical value. The original CIE method for the daylight spectra requires rounding intermediate coefficients M1 and M2 to 3 decimal places. This rounding is necessary to reproduce the tabulated values in Table T.1 of the CIE publication in <b>References</b> .
energy	a vector of energy levels

### Details

For `planckSpectra()` the valid range of temperatures is 0 to  $\text{Inf} (\infty)$  K, but with exceptions at the endpoints. For a negative temperature the spectrum is set to all NAs.

If temperature=0 and normalize=TRUE, the spectrum is set to all NAs. If temperature=0 and normalize=FALSE, the spectrum is set to all 0s.

Conversely, if temperature=Inf and normalize=FALSE, the spectrum is set to all NAs. If temperature=Inf and normalize=TRUE, the spectrum is set to the pointwise limit  $(560/\lambda)^4$  (which appears blue).

For daylightSpectra() the valid range of temperatures is 4000 to 25000 K. For a temperature outside this range the spectrum is set to all NAs.

The equations for daylightSpectra() and planckSpectra() are complex and can be found in the **References**.

IlluminantE() is trivial - all constant energy.

### Value

For planckSpectra() and daylightSpectra() :

A **colorSpec** object with **quantity** equal to 'energy', and **organization** equal to 'matrix' or 'vector'. The specnames are PNNNN or DNNNN for planckSpectra() and daylightSpectra() respectively.

The number of spectra in the object is the number of temperatures = length(temperature).

For illuminantE() :

A **colorSpec** object with **quantity** equal to 'energy'.

The number of spectra in the object is the number of energy levels = length(energy).

### References

Günther Wyszecki and W.S. Stiles. **Color Science : Concepts and Methods, Quantitative Data and Formulae**. Second Edition. Wiley-Interscience. 1982. page 146.

CIE 15: Technical Report: Colorimetry, 3rd edition. CIE 15:2004. Table T.1, pp 30-32, and Note 5 on page 69.

Schanda, Janos. CIE Colorimetry, in *Colorimetry: Understanding the CIE System*. Wiley Interscience. 2007. p. 42.

### See Also

[daylight](#), [resample](#), [organization](#), [quantity](#), [materialSpectra](#)

---

linearize

*linearize a colorSpec object - to make it ready for colorimetric calculations*

---

### Description

linearize spectra and return modified object

### Usage

```
## S3 method for class 'colorSpec'
linearize( x )
```

**Arguments**

x                    a **colorSpec** object

**Details**

If the `quantity(x)` is not 'absorbance' then x is returned unchanged.

If the `quantity(x)` is 'absorbance' then absorbance is converted to transmittance using

$$transmittance = 10^{-absorbance}$$

Surprisingly, there does not seem to be a similar logarithmic version of reflectance. Plots with `log(responsivity)` is somewhat common, but does not seem to have a separate name. I have not seen `log(radiometric power)`.

**Value**

`linearize` returns a **colorSpec** object with linear quantities.

**See Also**

[quantity](#)

---

lms1971

*Cone Fundamentals - 2-degree (1971)*


---

**Description**

`lms1971.5nm` the Vos & Walraven (1971) 2° cone fundamentals from 380 to 780 nm, at 5nm intervals

**Format**

A **colorSpec** object organized as a matrix with 3 columns:

long	the long wavelength responsivity
medium	the medium wavelength responsivity
short	the short wavelength responsivity

**Source**

<http://www.cvr1.org/database/text/cones/vw.htm>

**References**

Vos, J. J. & Walraven, P. L. On the derivation of the foveal receptor primaries. **Vision Research**. 11 (1971) pp. 799-818.

**See Also**

[lms2000](#)

**Examples**

```
summary(lms1971.5nm)
white.point = product( D65.1nm, lms1971.5nm, wave='auto' )
```

---

lms2000

*Cone Fundamentals - 2-degree (2000)*

---

**Description**

lms2000.1nm the Stockman and Sharpe (2000) 2° cone fundamentals from 390 to 830 nm, at 1nm intervals

**Format**

A **colorSpec** object organized as a matrix with 3 columns:

long	the long wavelength responsivity
medium	the medium wavelength responsivity
short	the short wavelength responsivity

**Source**

<http://www.cvrl.org/cones.htm>

**References**

Stockman, A., Sharpe, L. T., & Fach, C. C. (1999). The spectral sensitivity of the human short-wavelength cones. **Vision Research**. 39, 2901-2927.

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.

**See Also**[lms1971](#)**Examples**

```
summary(lms2000.1nm)
white.point = product( D65.1nm, lms2000.1nm, wave='auto' )
```

logging

*Logging in colorSpec package***Description**

There is some flexibility in the **colorSpec** logging level and format. Logging output goes to `stderr()`, just like the message stream; but see [sink\(\)](#) (and the pitfalls of using it).

**Logging Options**

`colorSpec.loglevel` The levels are: "FATAL", "ERROR", "WARN", "INFO", "DEBUG", and "TRACE" - the usual ones from **Log4j**. The initial level is "WARN". A "FATAL" event usually means an internal package error. When setting `colorSpec.loglevel` an initial letter is sufficient.

`colorSpec.logformat` The format is given by a string with standard **Log4j** *conversion specifications*:

```
%t  the date/time of the logging event.  %t can be followed by standard strftime specs in braces; see example.
%l  the level of the logging event
%n  namespace where event occurred
%f  function where event occurred
%m  the message itself
```

`colorSpec.stoponerror` If the this option is TRUE (the default), a log event with level "ERROR" stops execution; otherwise, execution keeps going. For interactive use, TRUE is probably better. For long batch jobs, FALSE might be appropriate, since then a single error may not force a complete repeat. A "FATAL" event always stops execution.

**References**

Wikipedia. **Log4j**. <http://en.wikipedia.org/wiki/Log4j>

**See Also**

[options](#), [cs.options](#), [sink](#), [stderr](#)

**Examples**

```
options( colorSpec.logformat="%t{%H:%M:%OS3} %l %n::%f(). %m", colorSpec.stuponerror=TRUE )

# or equivalently
cs.options( logformat="%t{%H:%M:%OS3} %l %n::%f(). %m", stop=TRUE )
```

luminsivity

*Luminous Efficiency Functions (photopic and scotopic)***Description**

luminsivity.1nm Four luminous efficiency functions, from 360 to 830 nm, at 1nm step

**Format**

A **colorSpec** object, with quantity 'energy->neural', and with 4 spectra:

photopic1924 The luminous efficiency function adopted by the CIE in 1924, and defining the *standard photopic observer*. It is only to be used when light levels are high enough that the sensitivity of the eye is mediated by cones, and not rods. It is the same as the *y-bar* function in xyz1931.1nm. It is used to define the *candela* in the *International System* (SI) and is the only one of these functions to appear in the SI. It was downloaded from [http://www.cvrl.org/database/data/lum/vl1924e\\_1.csv](http://www.cvrl.org/database/data/lum/vl1924e_1.csv) where it is defined from 360 to 830 nm.

scotopic1951 The luminous efficiency function adopted by the CIE in 1951, and defining the *standard scotopic observer*. It is only to be used when light levels are low enough to exclude the activation of cones. It has no effective role in colorimetry. It was downloaded from [http://www.cvrl.org/database/data/lum/scvle\\_1.csv](http://www.cvrl.org/database/data/lum/scvle_1.csv) where it is defined from 380 to 780 nm. It has been padded with 0s to 360 to 830 nm.

photopic1978 The luminous efficiency function for photopic vision, with adjustments in the blue region by Judd (1951) and Vos (1978). It was published by the CIE in 1988. It was downloaded from [http://www.cvrl.org/database/data/lum/vme\\_1.csv](http://www.cvrl.org/database/data/lum/vme_1.csv) where it is defined from 380 to 780 nm. It has been padded with 0s to 360 to 830 nm.

photopic2008 The CIE (2008) *physiologically-relevant* luminous efficiency function for photopic vision, by Stockman, Jagle, Pirzer, & Sharpe. It was downloaded from [http://www.cvrl.org/database/data/lum/linCIE2008v2e\\_1.csv](http://www.cvrl.org/database/data/lum/linCIE2008v2e_1.csv) where it is defined from 390 to 830 nm. It has been padded with 0s to 360 to 830 nm.

**Note**

Luminsivity is a self-coined *portmanteau word*: luminsivity = luminous \* responsivity. The word is unrelated to *emissivity*. The term *luminous responsivity* is not common, but appears on page 15 of *Grum*. The term *luminous efficiency function* is standard, but too long. The term *luminosity*

*function* is common, but also appears in astronomy.  
The object `luminsivity.1nm` is used by the function `photometric()`.

### Source

Colour & Vision Research Laboratory. Institute of Ophthalmology. University College London. UK.  
<http://www.cvr1.org/>

### References

Grum, Franc and Richard J. Becherer. **Radiometry**. Optical Radiation Measurements, Volume 1. Academic Press. 1979.

Stockman, A., Jagle, H., Pirzer, M., & Sharpe, L. T. (2008). The dependence of luminous efficiency on chromatic adaptation. *Journal of Vision*, 8, 16:1, 1-26.

### See Also

[xyz1931.1nm](#), [photometric](#)

### Examples

```
summary(luminsivity.1nm)
product( D65.1nm, luminsivity.1nm, wave='auto' )
```

---

materialSpectra	<i>compute standard material spectra</i>
-----------------	--

---

### Description

Compute neutral gray material constant reflectance/transmittance, and rectangular spectra. Also compute absorbance of the human lens, as a function of age.

### Usage

```
neutralMaterial( gray=1, wavelength=380:780 )
rectangularMaterial( lambda, alpha=1, wavelength=380:780 )

lensAbsorbance( age=32, wavelength=400:700 )
```

### Arguments

gray	a numeric N-vector of gray levels, in the interval [0,1]. gray=1 represents the <i>Perfect Reflecting Diffuser</i> .
lambda	a numeric Nx2 matrix with wavelength pairs in the rows, or a vector that can be converted to such a matrix, by row. The two wavelengths are the two <i>transition wavelengths</i> of the returned spectrum, see <b>Details</b> .

alpha	a numeric N-vector of <i>chromatic amplitudes</i> in the interval [-1,1]. N must be equal to nrow(lambda). alpha can also be a single number, which is then replicated to length nrow(lambda). The <i>chromatic amplitude</i> is defined by <i>Logvinenko</i> and controls the size of both transitions, see <b>Details</b> .
age	a numeric N-vector of ages in years; all ages must be $\geq 20$ .
wavelength	a vector of wavelengths for the returned object

### Details

A *rectangular spectrum*, or *rectangular metamer*, is easiest to define when  $\alpha = 1$  and  $\lambda_1 < \lambda_2$ . In this case it is a band-pass filter with transmittance=1 for  $\lambda \in [\lambda_1, \lambda_2]$  and transmittance=0 otherwise. To create a long-pass filter, just set  $\lambda_2$  to Inf, or any large wavelength outside the spectrum range; and similarly for a short-pass filter.

When  $0 < \alpha < 1$  the spectrum is a weighted mixture of this band-pass filter with a perfect neutral gray filter with transmittance=0.5 at all  $\lambda$ , using  $\alpha$  and  $1 - \alpha$  as the two weights. The minimum transmittance is  $(1 - \alpha)/2$  and the maximum is  $(1 + \alpha)/2$ , and their difference, the *chromatic amplitude*, is  $\alpha$ . It is still a band-pass filter.

If  $\alpha = 0$  the spectrum is a perfect neutral with transmittance=0.5.

To "flip" the spectrum to its complement (change band-pass to band-stop, etc.), change  $\alpha$  to a negative number, or swap  $\lambda_1$  and  $\lambda_2$ . If  $\lambda_1 == \lambda_2$  then the spectrum is undefined and a warning is issued (unless  $\alpha = 0$ ).

### Value

neutralMaterial() returns a **colorSpec** object with **quantity** equal to 'reflectance'. The reflectance of each spectrum is constant and taken from gray. There are N spectra in the object - one for each gray level.

rectangularMaterial() returns a **colorSpec** object with **quantity** equal to 'transmittance'. The transmittance of each spectrum is a step function with 0, 1 or 2 transitions (jumps) defined by the corresponding row in lambda. If rownames(lambda) is not NULL, they are copied to specnames of the output. Otherwise the specnames are computed from the shape of the spectrum using these acronyms: LP (long-pass), SP (short-pass), BP (band-pass), BS (band-stop), and N (neutral, in case alpha==0).

lensAbsorbance() returns a **colorSpec** object with **quantity** equal to 'absorbance'. The absorbance model for the human lens is taken from *Pokorny*. There are N spectra in the object - one for each age (N=length(age)).

### Logvinenko

It is clear that there are 3 degrees-of-freedom in the spectra returned by rectangularMaterial(). *Logvinenko* shows that these spectra in fact form a 3D ball, which he calls the *rectangle color atlas*. He also shows that if a material responder satisfies the 2-transition condition, then these spectra uniquely generate *all* colors in the corresponding object color solid. For more on this, see the **inversion** vignette.

### Ostwald

Every spectrum returned by rectangularMaterial() is an Ostwald ideal spectrum. In Ostwald's terminology, the *color content* = *chromatic amplitude* =  $\alpha$ . And the *black content* = *white content*



$= (1 - \alpha)/2$ . Note that the sum of these 3 contents is 1. However, Ostwald allows *black content* and *white content* to be unequal, as long as the sum of the 3 contents is 1, and all are non-negative. Thus there is one extra degree-of-freedom for Ostwald's ideal spectra, for a total of 4 degrees-of-freedom. If an additional argument (or arguments) were added to `rectangularMaterial()`, then it could return all Ostwald ideal spectra.

## References

Foss, Carl E. and Dorothy Nickerson and Walter C. Granville. Analysis of the Ostwald Color System. *J. Opt. Soc. Am.*. vol. 34. no. 7. pp. 361-381. July, 1944.

Logvinenko, A. D. An object-color space. **Journal of Vision**. 9(11):5, 1-23, (2009). <https://jov.arvojournals.org/article.aspx?articleid=2203976>. doi:10.1167/9.11.5.

Pokorny, Joel, Vivianne C. Smith, and Margaret Lutze. Aging of the Human Lens. **Applied Optics**. Vol. 26, No. 8. 15 April 1987. Table I. Page 1439.

## See Also

[lightSpectra](#), [quantity\(\)](#), [specnames\(\)](#), [computeADL\(\)](#), vignette **inversion**

## Examples

```
# make a perfect reflecting diffuser (PRD)
prd = neutralMaterial( 1 )

# make a perfect transmitting filter (PTF)
ptf = prd
quantity(ptf) = 'transmittance'

# make a band-stop filter (for interval [500,550])
# with 1% transmittance in the band, and 99% outside the band
bs = rectangularMaterial( c(500,550), -0.98, 400:700 )
bs = rectangularMaterial( c(550,500), 0.98, 400:700 ) # equivalent to previous line

# compare transmittance at 3 ages: 20, 32, and 80 years
plot( linearize(lensAbsorbance( c(20,32,80) )), col='black', lty=1:3 )
```

---

mean	<i>calculate mean of multiple spectra</i>
------	---

---

## Description

compute mean of all spectra in a **colorSpec** object

## Usage

```
## S3 method for class 'colorSpec'
mean( x, ... )
```

**Arguments**

x                    a `colorSpec` object  
 ...                  further arguments ignored

**Details**

This function might be useful when capturing many spectra on a spectrometer and averaging them to reduce noise.

**Value**

a **colorSpec** object with single spectrum = average of all spectra in **colorSpec**.

---

metadata	<i>metadata of a colorSpec object</i>
----------	---------------------------------------

---

**Description**

Retrieve or set the metadata of a **colorSpec** object.

**Usage**

```
## S3 method for class 'colorSpec'
metadata(x, ...)

## S3 replacement method for class 'colorSpec'
metadata(x, add=FALSE ) <- value
```

**Arguments**

x                    a **colorSpec** R object  
 ...                  optional names of metadata to return  
 value                a named list. If `add` is `FALSE`, `value` replaces any existing metadata. If `add` is `TRUE`, `value` is appended to the existing list of metadata. If a name already exists, its value is updated using `modifyList()`. Unnamed items in `value` are ignored.  
 add                  if `add=FALSE`, any existing metadata is discarded. If `add=TRUE` then existing metadata is preserved, using `modifyList()`.

**Details**

The metadata list is stored as `attr(x, 'metadata')`. After construction this list is empty.

**Value**

`metadata(x)` with no additional arguments returns the complete named list of metadata. If arguments are present, then only those metadata items are returned.

**Note**

Do not confuse extradata and metadata.  
 metadata is unstructured data that is attached to the entire **colorSpec** object. extradata is structured data, with a row of data for each spectrum in the object.

**See Also**

[extradata](#), [modifyList](#)

**Examples**

```
## Not run:
# get list of *all* metadata
metadata(x)

# get just the file 'path'
metadata( x, 'path' )

# set the 'date'
metadata( x ) = list( date="2016-04-01" )

## End(Not run)
```

---

 multiply

---

*multiply a colorSpec object by scalar, vector, or matrix*


---

**Description**

multiply spectra by coefficients and return modified object

**Usage**

```
## S3 method for class 'colorSpec'
multiply( x, s )

## S3 method for class 'colorSpec'
normalize( x, norm='L1' )
```

**Arguments**

x	a <b>colorSpec</b> object with M spectra
s	a scalar, an M-vector, or an MxP matrix. In the case of a matrix, assigning colnames(s) is recommended; see <b>Details</b> .
norm	one of 'L1', 'L2', or 'Linf', specifying one of the standard vector norms $L^1$ , $L^2$ , or $L^{inf}$ . norm can also be a numeric wavelength (e.g. 560 nm), and then the spectrum is scaled to have value 1 at this wavelength. Of course, this is not a true vector norm.

**Details**

For `multiply()`:

If `s` is an  $M \times P$  matrix, say  $\mathbf{S}$ , and one thinks of the spectra as organized in an  $N \times M$  matrix  $\mathbf{X}$ , then the new spectra are defined by the matrix  $\mathbf{XS}$ , which is  $N \times P$ . If the  $P$  column names of `s` are set, then they are copied to the `specnames` of the output. Otherwise, default spectrum names are assigned as in `colorSpec()`, with a warning.

If `s` is an  $M$ -vector, then  $\mathbf{S}=\text{diag}(s)$  is computed and used in the previous sentence. This has the effect of multiplying spectrum  $i$  by `s[i]`.

If `s` is a scalar then every spectrum is multiplied by `s`.

The multiplication may produce negative entries, but no check is made for this.

WARNING: An  $M$ -vector and an  $M \times 1$  matrix may yield quite different results.

For `normalize()`:

`normalize()` calls `multiply()` with `s` = an  $M$ -vector. If the norm of a spectrum is 0, then it is left unchanged.

**Value**

`multiply` returns a **colorSpec** object with the matrix of spectra of `x` multiplied by `s`.

`normalize` returns a **colorSpec** object with each spectrum of `x` scaled to have given norm equal to 1.

In both functions, the `quantity` and `wavelength` are preserved.

**Note**

If `x` is organized as a matrix, and `s` is a scalar, the one can use the simpler and equivalent `s*x`.

**See Also**

`product()`, `quantity()`, `wavelength()`, `specnames()`, `colorSpec()`

---

officialXYZ

*Query the Official XYZ values for Standard Illuminants*

---

**Description**

In careful calculations with standard illuminants, it is often helpful to have the 'official' values of XYZ, i.e. with the right number of decimal places.

**Usage**

```
officialXYZ( name )
```

**Arguments**

`name` a subvector of `c('A', 'B', 'C', 'D50', 'D50.ICC', 'D55', 'D65', 'D75', 'E', 'F2', 'F7', 'F11')`, which are the names of some standard illuminants

**Details**

All XYZ values are taken from the ASTM publication in **References**, except B which is taken from *Wyszecki & Stiles* and D50.ICC which is taken from ICC publications. The latter is different than that of ASTM.

**Value**

An Mx3 matrix where M is the length of name. Each row filled with the official XYZ, but if the illuminant name is not recognized the row is all NAs. The matrix rownames are set to name, and colnames to c('X', 'Y', 'Z').

**WARNING**

This function is deprecated. New software should use `spacesRGB::standardXYZ()` instead.

**Note**

The input names are case-sensitive. The output XYZ is normalized so that Y=1.

**References**

ASTM E 308 - 01. Standard Practice for Computing the Colors of Objects by Using the CIE System. (2001).

Günther Wyszecki and W. S. Stiles. *Color Science: Concepts and Methods, Quantitative Data and Formulae*, Second Edition. John Wiley & Sons, 1982. Table I(3.3.8) p. 769.

**See Also**

[ABC](#), [D50](#), [D65](#), [Fluorescents](#), [illuminantE](#)

**Examples**

```
officialXYZ( c('A','D50','D50.ICC','D65') )
#           X Y       Z
# A       1.0985000 1 0.3558500
# D50     0.9642200 1 0.8252100
# D50.ICC 0.9642029 1 0.8249054
# D65     0.9504700 1 1.0888300
```

---

organization

*organization of a colorSpec object*

---

**Description**

Retrieve or set the organization of a **colorSpec** object.

**Usage**

```
## S3 method for class 'colorSpec'
organization(x)

## S3 replacement method for class 'colorSpec'
organization(x) <- value
```

**Arguments**

**x** a **colorSpec** R object

**value** a valid organization: 'vector', 'matrix', 'df.col', or 'df.row'.

**Details**

If `organization(x)` is "vector", then `x` is a vector representing a single spectrum. Compare this with `stats::ts()`.

If `organization(x)` is "matrix", then `x` is a matrix and the spectra are stored in the columns.

If `organization(x)` is "df.col", then `x` is a data.frame with `M+1` columns, where `M` is the number of spectra. The wavelengths are stored in column 1, and the spectra in columns 2:(`M+1`). This organization is good for printing to the console, and writing to files.

If the organization of `x` is "df.row", then `x` is a data.frame with `N` rows, where `N` is the number of spectra. The spectra are stored in the last column, which is a matrix with the name "spectra". The other columns preceding spectra (if present) contain extra data associated with the spectra; see [extradata](#).

**Value**

`organization(x)` returns a valid organization: 'vector', 'matrix', 'df.col', or 'df.row'.

**Note**

In `organization(x) <- value`

if `x` has more than 1 spectrum, then value equal to 'vector' is invalid and ignored.

If `organization(x)` is equal to 'df.row' and also has [extradata](#), then changing the organization silently discards the `extradata`.

**See Also**

[colorSpec](#); [extradata](#)

**Examples**

```
organization(Hoya)           # returns 'df.row'
organization(Hoya) = 'matrix' # extradata in Hoya is silently discarded
```

---

photometric                      *convert illuminant spectra to photometric units*

---

### Description

Convert radiometric units of power or energy to photometric units, using 4 standard photometric weighting curves. Actinometric units (number of photons) are converted to radiometric units (energy of photons) on-the-fly.

### Usage

```
## S3 method for class 'colorSpec'
photometric( x, photopic=683, scotopic=1700, multiplier=1 )
```

### Arguments

x	a <b>colorSpec</b> object with type equal to 'light', and with M spectra
photopic	the conversion factor for photopic vision, in lumen/watt. The CIE standard is 683, and another common value is 683.002.
scotopic	the conversion factor for scotopic vision, in lumen/watt. The CIE standard is 1700, and another common value is 1700.06.
multiplier	an conversion factor intended for conversion of units, and applied to both photopic and scotopic vision. For example if the input unit of x is $watt * sr^{-1}$ , and the desired output unit is <i>candlepower</i> , then set <code>multiplier=1/0.981</code> .

### Details

The function computes the product of x with `luminsivity.1nm`. This product is an Mx4 matrix, where M is the number of spectra in x. There are 3 columns for photopic vision, and 1 column for scotopic vision. These columns are multiplied by the appropriate conversion factors and the resulting Mx4 matrix is returned.

The 5 power-based input quantities and corresponding photometric outputs are:

<b>radiant power</b> [ <i>watt</i> ]	→	<b>luminous flux</b> [ <i>lumen</i> ]
<b>irradiance</b> [ $watt * m^{-2}$ ]	→	<b>illuminance</b> [ $lumen * m^{-2} = lux$ ]
<b>radiant exitance</b> [ $watt * m^{-2}$ ]	→	<b>luminous exitance</b> [ $lumen * m^{-2} = lux$ ]
<b>radiant intensity</b> [ $watt * sr^{-1}$ ]	→	<b>luminous intensity</b> [ $lumen * sr^{-1} = candela$ ]
<b>radiance</b> [ $watt * sr^{-1} * m^{-2}$ ]	→	<b>luminance</b> [ $candela * m^{-2} = nit$ ]

The 2 *common* energy-based input quantities and corresponding photometric outputs are:

<b>radiant energy</b> [ <i>joule</i> ]	→	<b>luminous energy</b> [ <i>talbot = lumen - second</i> ]
<b>radiant exposure</b> [ $joule * m^{-2}$ ]	→	<b>luminous exposure</b> [ $talbot * m^{-2} = lux - second$ ]

and there are 3 more obtained by integrating over time. For example "time-integrated radiance" —> "time integrated luminance". But I have not been able to find names for these 3. The *talbot* is the unofficial name for a lumen-second.

### Value

photometric() returns an Mx4 matrix, where M is the number of spectra in x. The rownames are specnames(x), and the colnames are specnames(luminsivity.1nm).  
In case of ERROR it returns NULL.

### Note

To get the right output quantity and units, the user must know the input quantity and units. If the units are different than those in the above list, then set multiplier appropriately.

It is up to the user to determine whether *photopic* or *scotopic* vision (or neither) is appropriate. The intermediate *mesopic* vision is currently a subject of research by the CIE, and might be added to this function in the future.

### References

Poynton, Charles. **Digital Video and HD - Algorithms and Interfaces**. Morgan Kaufmann. Second Edition. 2012. Appendix B, pp. 573-580.

### See Also

[quantity](#), [type](#), [luminsivity.1nm](#), [radiometric](#)

### Examples

```
photometric( solar.irradiance ) # unit is watt*m^{-2}

#           photopic1924 scotopic1951 photopic1978 photopic2008 # units are lux
# AirMass.0      133100.41    313883.2    133843.65    140740.3
# GlobalTilt    109494.88    250051.5    110030.31    115650.0
# AirMass.1.5   97142.25     215837.1    97571.57     102513.7
```

---

plot

*plot spectra*

---

### Description

plot the spectra in a **colorSpec** object as lines

### Usage

```
## S3 method for class 'colorSpec'
plot( x, color=NULL, subset=NULL, main=TRUE, legend=TRUE, CCT=FALSE, add=FALSE, ... )
```



**Arguments**

x	a colorSpec object
color	If color=NULL then colors are computed from the spectra themselves. If type(x) is 'material' the color is computed using illuminant D65.1nm and responder BT.709.RGB with no further normalization. Otherwise the spectrum color is faked by changing its quantity to 'energy' and taking the <code>product</code> with BT.709.RGB. The resulting RGBs are normalized to have a maximum of 1. This RGB normalization is done <i>before</i> processing the subset argument. If color='auto' then a suitable set of colors is generated using <code>colorRampPalette()</code> . Otherwise color is passed along to <code>lines()</code> as the col argument, e.g. col='black'.
subset	specifies a subset of x to plot; see <code>subset()</code> for acceptable arguments.
main	If main=TRUE then a main title is generated from the file 'path' in the metadata list, or from <code>deparse(substitute(x))</code> . If main=FALSE then no main title is displayed. And if main is a string then that string is used as the main title.
legend	If legend=TRUE then a pretty legend using <code>specnames(x)</code> is placed in the 'topright' corner of the plot. If legend is a string it is interpreted as naming a corner of the plot and passed as such to the function <code>legend</code> . If legend=FALSE then no legend is drawn.
CCT	If CCT=TRUE and the type of x is 'light' then the CCT of each spectrum is added to the legend; see <code>computeCCT()</code> .
add	If add=TRUE then the lines are added to an existing plot, and these arguments are ignored: main, ylab, xlim, ylim, and log; see <b>Details</b> .
...	other graphical parameters, see <b>Details</b>

**Details**

Commonly used graphical parameters are:

lty, lwd passed on to `lines()`, with no checks

pch If pch is an integer, it is passed on to `points()` which adds points on top of the lines

ylab If ylab is a string then it is passed on to `plot.default()`, otherwise suitable default string is generated.

xlim, ylim If xlim and ylim are 2-vectors, they are passed to `plot.default`. If one of the components is NA then a suitable default is supplied.

log passed on to `plot.default()`. Care must be taken for y because many spectra are 0 at some wavelengths, and even negative. Use ylim in such cases.

**Value**

TRUE or FALSE

**See Also**

`print`, `summary`, `colorRampPalette` in base.

**Examples**

```
plot( 100 * BT.709.RGB )
plot( xyz1931.1nm, add=TRUE, lty=2, legend=FALSE )
```

---

plotOptimals

*Plot Optimal Colors*


---

**Description**

Consider a **colorSpec** object *x* with type equal to 'responsivity.material' and 3 responsivity spectra. The function `plotOptimals3D()` makes a plot of the *object-color solid* for *x*. This solid is a *zonohedron* in 3D. The 3D drawing package **rgl** is required.

Consider a **colorSpec** object *x* with type equal to 'responsivity.material' and 2 responsivity spectra. The function `plotOptimals2D()` makes a plot of the *object-color solid* for *x*. This solid is a *zonogon* in 2D. The 3D drawing package **rgl** is *not* required.

The set of all possible material reflectance functions (or transmittance functions) is convex, closed, and bounded (in any reasonable function space), and this implies that the set of all possible output responses from *x* is also convex, closed, and bounded. The latter set is called the *object-color solid*, or *Rösch Farbkörper*, for *x*. A color on the boundary of the *object-color solid* is called an *optimal color*. For more discussion see `sectionOptimalColors()`.

**Usage**

```
## S3 method for class 'colorSpec'
plotOptimals3D( x, size=50, type='w', both=TRUE )
```

```
## S3 method for class 'colorSpec'
plotOptimals2D( x )
```

**Arguments**

<i>x</i>	a <b>colorSpec</b> object with type equal to 'responsivity.material' and 2 or 3 spectra, as appropriate.
<i>size</i>	an integer giving the number of wavelengths at which to resample <i>x</i> . To skip resampling, set <i>size</i> =NA.
<i>type</i>	<i>type</i> ='w' for a wireframe plot of the parallelogram faces. <i>type</i> ='p' for a point plot with points at the centers of the parallelograms.
<i>both</i>	the color solid is symmetric about its center, so only half of it must be computed. If <i>both</i> =TRUE it plots one half in black and the other half in red. If <i>both</i> =FALSE it only plots one half in black.

**Value**

The functions return TRUE or FALSE.

### Details for 3D

If  $n$  is the number of wavelengths, the number of parallelogram faces of the zonohedron is  $n*(n-1)$ . The time to compute these faces increase with  $n$  even faster, so that is why the default `size=50` is a fairly small number. It was chosen to be a reasonable compromise between detail and performance. In addition to the wireframe or points, it draws the box with opposite vertices at the "poles" **0** and **W** and the diagonal segment of neutral grays that connects **0** and **W**.

### Details for 2D

If  $n$  is the number of wavelengths, the number of edges in the zonogon is  $2*n$ . Computing these edges is fast and visualization is easy, so there are no plotting options at this time.

### Note

If all responsivity functions of  $x$  are non-negative, the *object-color solid* of  $x$  is inside the box. If the responsivity functions of  $x$  have negative lobes, the *object-color solid* of  $x$  extends outside the box. Indeed, the box may actually be *inside* the optimals.

### References

Centore, Paul. *A Zonohedral Approach to Optimal Colours*. **Color Research & Application**. Vol. 38. No. 2. pp. 110-119. April 2013.

Logvinenko, A. D. An object-color space. *Journal of Vision*. 9(11):5, 1-23, (2009). <https://jov.arvojournals.org/article.aspx?articleid=2203976>. doi:10.1167/9.11.5.

West, G. and M. H. Brill. Conditions under which Schrödinger object colors are optimal. **Journal of the Optical Society of America**. 73. pp. 1223-1225. 1983.

### See Also

`type()`, `probeOptimalColors()`, `sectionOptimalColors()`, vignette **optimals**

### Examples

```
human = product( D50.5nm, 'slot', xyz1931.5nm, wave=seq(400,770,by=5) )
plotOptimals3D( human )
```

```
plotOptimals2D( subset(human,2:3) )      # y and z only
```

```
scanner = product( D50.5nm, 'slot', BT.709.RGB, wave=seq(400,770,by=5) )
plotOptimals3D( scanner )
```

---

print                      *Convert colorSpec object to readable text*

---

### Description

display a **colorSpec** object as readable text. Output goes to `stdout()`.

### Usage

```
## S3 method for class 'colorSpec'  
print( x, ... )
```

```
## S3 method for class 'colorSpec'  
summary( object, long=TRUE, ... )
```

### Arguments

x	a colorSpec object
object	a colorSpec object
long	logical indicating whether to print metadata, calibrate, sequence, and emulate data
...	further arguments ignored

### Details

If `long=FALSE`, `summary()` prints a summary of the wavelength vector, and names of all spectra. For each spectrum it prints the range of values, `LambdaMax`, and `extradata` if any. If `long=TRUE` it also prints data listed above (if any).

The function `print()` simply calls `summary()` with `long=FALSE`.

### Value

Both functions return (invisibly) the character vector that was just printed to `stdout()`.

### See Also

[extradata](#), [print](#), [summary](#), [stdout](#)

### Examples

```
print( xyz1931.1nm )  
  
xyz1931.1nm    # same thing, just calls print()
```

---

 probeOptimalColors     *compute optimal colors by ray tracing*


---

### Description

Consider a **colorSpec** object  $x$  with type equal to 'responsivity.material'. The set of all possible material reflectance functions (or transmittance functions) is convex, closed, and bounded (in any reasonable function space), and this implies that the set of all possible output responses from  $x$  is also convex, closed, and bounded. The latter set is called the *object-color solid* or *Rösch Farbkörper* for  $x$ . A color on the boundary of the *object-color solid* is called an *optimal color*. The special points **W** (the response to the perfect reflecting diffuser) and **0** are on the boundary of this set. The interior of the line segment of neutrals joining **0** to **W** is in the interior of the *object-color solid*. It is natural to parameterize this segment from 0 to 1 (from **0** to **W**).

A ray  $r$  that is based at a point on the interior of the neutral line segment must intersect the boundary of the *object-color solid* in a unique optimal color. The purpose of the function probeOptimalColors() is to compute that intersection point.

Currently the function only works if the number of spectra in  $x$  is 3 (e.g. RGB or XYZ).

Before **colorSpec** v 0.8-1 this function used a 2D root-finding method that could only find optimal colors whose spectra contain 0, 1, or 2 transitions. But starting with v0.8-1, we have switched to zonohedral representation of the object-color solid, which makes it possible to discover more than 2 transitions. The inspiration for this change is the article by Centore. To *inspect* these computed spectra, the argument spectral must be set to TRUE.

### Usage

```
## S3 method for class 'colorSpec'
probeOptimalColors( x, gray, direction, aux=FALSE, spectral=FALSE, tol=1.e-6 )
```

### Arguments

x	a <b>colorSpec</b> object with type equal to 'responsivity.material' and 3 spectra
gray	vector of numbers in the open interval (0,1) that define neutral grays on the line segment from black to white; this neutral gray point is the basepoint of a probe ray
direction	a numeric Nx3 matrix with directions of the probe rays in the rows, or a numeric vector that can be converted to such a matrix, by row.
aux	a logical that specifies whether to return extra performance and diagnostic data; see <b>Details</b>
spectral	if TRUE, the function returns a <b>colorSpec</b> object with the optimal spectra, see <b>Value</b> .
tol	error tolerance for the intersection of probe and object-color boundary

## Details

Each gray level and each direction defines a ray. So the total number of rays traced is `length(gray) * nrow(direction)`. The 3 responsivities are regarded not as continuous functions, but as step functions. This implies that the color solid is a zonohedron. In the preprocessing phase the zonohedral representation is calculated. The faces of the zonohedron are either parallelograms, or *compound faces* that can be partitioned into parallelograms. The centers of all these parallelograms are computed, along with their normals and plane constants.

This representation of the color solid is very strict regarding the 2-transition assumption. During use, one can count on there being some spectra with more than two transitions. Forcing the best 2-transition spectrum is a possible topic for the future.

## Value

If argument `spectral=FALSE`, `probeOptimalColors()` returns a `data.frame` with a row for each traced ray. There are `length(gray) * nrow(direction)` rays. The columns in the output are:

<code>gray</code>	the graylevel defining the <i>basepoint</i> of the ray. $basepoint = gray * W$
<code>direction</code>	the <i>direction</i> of the ray
<code>s</code>	computed scalar so that $basepoint + s * direction$ is optimal
<code>optimal</code>	the optimal color on the boundary; $optimal = basepoint + s * direction$
<code>lambda</code>	<code>lambda.1</code> and <code>lambda.2</code> at the 2 transitions, in nm. <code>lambda.1 &lt; lambda.2</code> => bandpass, and <code>lambda.1 &gt; lambda.2</code> => bandstop. It will happen that the optimal spectrum has more than 2 transitions; in this case both lambdas are set to NA.
<code>dol</code>	delta and omega - the Logvinenko parameters $(\delta, \omega)$ for optimal colors, plus lambda ( $\lambda$ ) in nm. $\omega$ is the reparameterization of $\lambda$ ; see <i>Logvinenko</i> . If there are more than 2 transistions, these are set to NA.

If `aux` is TRUE, these auxiliary columns related to performance and diagnostics are added:

<code>timetrace</code>	time to trace the ray, in seconds
<code>parallelograms</code>	# of parallelograms in the (possibly compound) face. 1 means just a single parallelogram.
<code>tested</code>	# of parallelograms actually tested for ray intersection. This only has meaning for compound faces.
<code>alpha</code>	the 2 coordinates of the intersection point inside the parallelogram

If argument `spectral=TRUE`, `probeOptimalColors()` returns a **colorSpec** object with quantity 'reflectance'. This object contains the optimal spectra, and can be used to inspect the spectra with more than 2 transitions, which will happen. The above-mentioned `data.frame` can then be obtained by applying `extradata()` to the returned object.

If an individual ray could not be traced (which should be rare), the row contains NA in appropriate columns.

In case of global error, the function returns NULL.

**WARNING**

The preprocessing calculation of the zonohedron dominates the total time. And this time goes up rapidly with the number of wavelengths. We recommend using a wavelength step of 5nm, as in the **Examples**. For best results, batch a lot of rays into a single function call and then process the output.

Moreover, the preprocessing time is dominated by the partitioning of the compound faces into parallelograms. This is made worse by spectral responses with little overlap, as in `scanner.ACES`. In these cases, try a larger step size, and then reduce. Optimizing these compound faces is a possible topic for the future.

**References**

- Centore, Paul. *A zonohedral approach to optimal colours*. **Color Research & Application**. Vol. 38. No. 2. pp. 110-119. April 2013.
- Logvinenko, A. D. An object-color space. **Journal of Vision**. 9(11):5, 1-23, (2009). <https://jov.arvojournals.org/article.aspx?articleid=2203976>. doi:10.1167/9.11.5.
- Schrödinger, E. (1920). Theorie der Pigmente von grösster Leuchtkraft. **Annalen der Physik**. 62, 603-622.
- West, G. and M. H. Brill. Conditions under which Schrödinger object colors are optimal. **Journal of the Optical Society of America**. 73. pp. 1223-1225. 1983.

**See Also**

`type`, vignette `optimals`, `scanner.ACES`, `extradata()`

**Examples**

```

wave = seq(400,700,by=5)
D50.eye = product( D50.5nm, 'material', xyz1931.1nm, wavelength=wave )
probeOptimalColors( D50.eye, c(0.2,0.5,0.9), c(1,2,1, -1,-2,-1) )

##   gray direction.1 direction.2 direction.3      s optimal.1 optimal.2
## 1 0.2           1           2           1 32.306207 52.533143 85.612065
## 2 0.2          -1          -2          -1 8.608798 11.618138 3.782055
## 3 0.5           1           2           1 20.993144 71.560483 94.485416
## 4 0.5          -1          -2          -1 20.993144 29.574196 10.512842
## 5 0.9           1           2           1 4.333700 95.354911 103.165832
## 6 0.9          -1          -2          -1 35.621938 55.399273 23.254556

##   optimal.3 lambda.1 lambda.2   dol.delta   dol.omega   dol.lambda
## 1 49.616046 451.8013 598.9589 0.63409966 0.48287469 536.97618091
## 2 8.701041 636.3031 429.4659 0.08458527 0.99624955 674.30015903
## 3 64.267740 441.9105 615.0822 0.78101041 0.49048222 538.73234859
## 4 22.281453 615.0822 441.9105 0.21898959 0.99048222 662.20606601
## 5 82.227974 422.9191 648.7404 0.95800430 0.49825407 540.49590064
## 6 42.272337 593.2415 455.2425 0.42035428 0.97962398 650.57382749

# create a 0-1 spectrum with 2 transitions
rectspec = rectangularMaterial( lambda=c(579.8697,613.7544), alpha=1, wave=wave )

```

```

# compute the corresponding color XYZ
XYZ = product( rectspec, D50.eye )
XYZ
##                X                Y                Z
## BP_[579.87,613.754] 33.42026 21.96895 0.02979764

# trace a ray from middle gray through XYZ
white.XYZ = product( neutralMaterial(1,wave=wave), D50.eye )
direction = XYZ - white.XYZ/2

res = probeOptimalColors( D50.eye, 0.5, direction, aux=FALSE )
res$s
## 1.00004 the ray has gone past the original color to the boundary

res$optimal
##                X                Y                Z
## [1,] 33.41958 21.96774 0.02808178

res$lambda
## NA NA because there are more than 2 transitions in the true optimal

# since s=1.00004 > 1,
# XYZ is actually in the interior of the color solid, and not on the boundary.
# The boundary is a little-bit further along the ray,
# and the corresponding spectrum has more than 2 transitions.

```

---

product

*Compute the product of colorSpec objects*


---

## Description

Take a sequence of **colorSpec** objects and compute their product. Only certain types of sequences are allowed. The return value can be a new **colorSpec** object or a matrix; see **Details**.

## Usage

```
## S3 method for class 'colorSpec'
product( ... )
```

## Arguments

... unnamed arguments are **colorSpec** objects, and possibly a single character string, see **Details**. Possible named arguments are:

wavelength The default wavelength='identical' means that all the **colorSpec** objects must have the same wavelength sequence; if they do not it is an ERROR. wavelength can be a new wavelength sequence, and all the objects are then **resampled** at these new wavelengths. wavelength can



also be 'auto' or NULL which means to compute a suitable wavelength sequence from those of the objects, see **Details**. It is OK to abbreviate the string wavelength (e.g. to wave); see **Examples**. It is OK for the wavelength sequence to be irregular; when the return value is a matrix the integration weights the spectrum values appropriately.

method, span, extrapolation, clamp passed to `resample()` with no checking or changes

integration only applies when the return type is matrix. The default option is 'rectangular', which means to weight the spectrum value equally at all wavelengths; this is the ASTM E308-01 recommendation. The other option is 'trapezoidal', which means to give the 2 endpoint wavelength values 1/2 the weight of the others. Trapezoidal integration is provided mostly for compatibility with other software.

## Details

To explain the allowable product sequences it is helpful to introduce some simple notation for the objects:

notation	colorSpec type	description of the object
$L$	light	a light source
$M$	material	a material
$R_L$	responsivity.light	a light responder (aka detector)
$R_M$	responsivity.material	a material responder (e.g. a scanner)

It is also helpful to define a sequence of positive integers to be *conformable* iff it has at most one value greater than 1. For example, a sequence of all 1s is conformable. A sequence of all  $q$ 's is conformable. The sequences  $c(1, 3)$  and  $c(1, 1, 4, 1, 1, 4, 1)$  are conformable, but  $c(1, 1, 4, 1, 3, 4, 1)$  is not.

There are 6 types of sequences for which the product is defined:

$$1. M_1 * M_2 * \dots * M_m \mapsto M'$$

The product of  $m$  materials is another material. Think of a stack of  $m$  transmitting filters effectively forming a new filter. If we think of each object as a matrix (with the spectra in the columns), then the product is element-by-element using  $R$ 's  $*$  - the Hadamard product. The numbers of spectra in the terms must be conformable. If some objects have 1 spectrum and all the others have  $q$ , then the column-vector spectrums are repeated  $q$  times to form a matrix with  $q$  columns. If the numbers of spectra are not conformable, it is an ERROR and the function returns NULL.

As an example, suppose  $M_1$  has 1 spectrum and  $M_2$  has  $q$  spectra, and  $m = 2$ . Then the product is a material with  $q$  spectra. Think of an IR-blocking filter followed by the RGB filters in a 3-CCD camera.

$$2. L * M_1 * M_2 * \dots * M_m \mapsto L'$$

The product of a light source followed by  $m$  materials is a light source. Think of a light source followed by a stack of  $m$  transmitting filters, effectively forming a new light source. The numbers of spectra in the terms must be conformable as in sequence 1, and the matrices are multiplied element by element.

As an example, suppose  $L$  has 1 spectrum and  $M_1$  has  $q$  spectra, and  $m = 1$ . Then the product is a light source with  $q$  spectra. Think of a light source followed by a filter wheel with  $q$  filters.

$$3. M_1 * M_2 * \dots * M_m * R_L \mapsto R'_L$$

The product of  $m$  materials followed by a light responder, is a light responder. Think of a stack of  $m$  transmitting filters in front of a camera, effectively forming a new camera. The numbers of spectra in the terms must be conformable as in sequence 1, and the matrices are multiplied element by element.

As an example, suppose  $R_L$  has 1 spectrum and  $M_1$  has  $q$  spectra, and  $m = 1$ . Then the product is a light responder with  $q$  spectra. Think of a 3-CCD camera in which all 3 CCDs have exactly the same responsivity and so can be modeled with a single object  $R_L$ .

$$4. L * M_1 * \dots * \bullet * \dots * M_m * R_L \mapsto R'_M$$

This is the strangest product. The bullet symbol  $\bullet$  means that a variable material is inserted at that slot in the sequence (or light path). For each material spectrum inserted there is a response from  $R_L$ . Therefore the product of this sequence is a material responder  $R_M$ . Think of a light source  $L$  going through a transparent object  $\bullet$  on a flatbed scanner and into a camera  $R_L$ . For more about the mathematics of this product, see the [colorSpec-guide.pdf](#) in the doc directory. These material responder spectra are the same as the *effective spectral responsivities* in *Digital Color Management*. The numbers of spectra in the terms must be conformable as in sequence 1, and the product is a material responder with  $q$  spectra.

In the function `product()` the location of the  $\bullet$  is marked by any character string whatsoever - it's up to the user who might choose something that describes the typical material (between the light source and camera). For example one might choose:

```
scanner = product( A.1nm, 'photo', Flea2.RGB, wave='auto')
```

to model a scanner that is most commonly used to scan photographs. Other possible strings could be 'artwork', 'crystal', 'varmaterial', or even 'slot'. See the [gallery](#) vignette for a worked-out example.

$$5. L * M_1 * M_2 * \dots * M_m * R_L \mapsto matrix$$

The product of a light source, followed by  $m$  materials, followed by a light responder, is a matrix! The numbers of spectra in the terms must splittable into a conformable left part ( $L'$  from sequence 2.) and a conformable right part ( $R'_L$  from sequence 3.). There is a row for each spectrum in  $L'$ , and a column for each spectrum in  $R'_L$ . Suppose the element-by-element product of the left part is  $n \times p$  and the element-by-element product of the right part is  $n \times q$ , where  $n$  is the number of wavelengths. Then the output matrix is the usual matrix product `%%` of the transpose of the left part times and right part, which is  $p \times q$ .

As an example, think of a light source followed by a reflective color target with 24 patches followed by an RGB camera. The sequence of spectra counts is `c(1, 24, 3)` which is splittable into `c(1, 24)` and `c(3)`. The product matrix is  $24 \times 3$ . See the [gallery](#) vignette for a worked-out example.

Note that is OK for there to be no materials in this product; it is OK if  $m = 0$ . See the [blueflame](#)

vignette for a worked-out example.

6.  $M_1 * M_2 * \dots * M_m * R_M \mapsto matrix$

The product of  $m$  materials followed by a material responder, is a matrix ! The sequence of numbers of spectra must be splittable into left and right parts as in sequence 4, and the product matrix is formed the same way. One reason for computing this matrix in 2 steps is that one can [calibrate](#) the material responder separately in a customizable way. See the [gallery](#) vignette for a worked-out example with a flatbed scanner.

Note that sequences 5. and 6. are the only ones that use the usual matrix product `%*`. They may also use the Hadamard matrix product `*`, as in sequences 1 to 4.

The argument `wavelength` can also be 'auto' or NULL. In this case the intersection of all the wavelength ranges of the objects is computed. If the intersection is empty, it is an ERROR and the function returns NULL. The wavelength step `step.wl` is taken to be the smallest over all the object wavelength sequences. If the minimum `step.wl` is less than 1 nanometer, it is rounded off to the nearest power of 2 (e.g 1, 0.5, 0.25, ...).

### Value

`product()` returns either a **colorSpec** object or a matrix, see [Details](#).

If `product()` returns a **colorSpec** object, the [organization](#) of the object is 'matrix' or 'vector'; any [extradata](#) is lost. However, all terms in the product are saved in `attr(*, 'sequence')`. One can use `str()` to inspect this attribute.

If `product()` returns a matrix, this matrix can sometimes be ambiguous, see [Note](#).

All actinometric terms are converted to radiometric on-the-fly and the returned **colorSpec** object is also radiometric.

In case of ERROR it returns NULL.

### Note

The product for sequences 1, 2, and 3 is associative. After all matrices are filled out to have  $q$  columns, the result is essentially a Hadamard product of matrices, which is associative. Also note that a subsequence of sequences 2 and 3 might be sequence 1.

The product for sequence 4 is never associative, since subproducts that contain the variable  $\bullet$  are undefined. However the result is essentially a Hadamard product of matrices, and is unambiguous.

The product for sequence 5 is associative in special cases, but not in general. The problem is that the left and right splitting point is not unique. If all objects have only a single spectrum, then it `*is*` associative, and therefore unambiguous. If the left part has a different number of multiple spectra than the right part, then it is not associative in general since some ways of grouping the product may be undefined.

Moreover, in some cases the product can be ambiguous. Suppose that the vector of spectrum counts is `c(1, 3, 1)`; this could come from a single light source, followed by 3 filters (e.g. RGB), followed by a graylevel camera. There are 2 ways to split this: `"1|3,1"` and `"1,3|1"`. The first split is interpreted as the light source into a camera with 3 channels. The second split is interpreted as 3

colored light sources into a graylevel camera. In the first split the returned matrix is a 1x3 row vector. In the second split the returned matrix is a 3x1 column vector. For the vector "1, 3, 1", one can show that the computed components in the 2 vectors are equal, so the ambiguity is benign. But consider the longer sequence "1, 3, 3, 1". There are 3 ways to split this, and the returned matrices are 1x3, 3x3, and 3x1. So this ambiguity is obviously a problem. Whenever there is an ambiguity, the function chooses a splitting in which the left part is as long as possible, and issues a warning message. The user should inspect the result carefully. To avoid the ambiguity, the user should break the product into smaller pieces and call `product()` multiple times.

The product for sequence 6 is essentially the same as sequence 5, and the function issues a warning message when appropriate. Note that a subproduct is defined only if it avoids the final multiplication with  $R_M$ .

## References

Edward J. Giorgianni and Thomas E. Madden. **Digital Color Management: Encoding Solutions**. 2nd Edition John Wiley. 2009. Figure 10.11a. page 141.

Wikipedia. **Hadamard product (matrices)**. [http://en.wikipedia.org/wiki/Hadamard\\_product\\_%28matrices%29](http://en.wikipedia.org/wiki/Hadamard_product_%28matrices%29)

ASTM E308-01. Standard Practice for Computing the Colors of Objects by Using the CIE System. (2001).

## See Also

[wavelength](#), [type](#), [resample](#), [calibrate](#), [radiometric](#), [step.wl](#)

## Examples

```
# sequence 1.
path = system.file( "extdata/objects/Midwest-SP700-2014.txt", package='colorSpec' )
blocker.IR = readSpectra( path )
product( blocker.IR, Hoya, wave='auto' )

# sequence 2.
product( subset(solar.irradiance,1), atmosphere2003, blocker.IR, Hoya, wave='auto' )

# sequence 3.
plumbicon = readSpectra( system.file( "extdata/cameras/plumbicon30mm.txt", package='colorSpec' ) )
product( blocker.IR, subset(Hoya,1:3), plumbicon, wave='auto' )

# sequence 4.
# make an RGB scanner
bluebalancer = subset(Hoya,'LB')
# combine tungsten light source A.1nm with blue light-balance filter
# use the string 'artwork' to mark the variable material slot
scanner = product( A.1nm, bluebalancer, 'artwork', Flea2.RGB, wave='auto' )
```

```

# sequence 5.
product( D65.1nm, Flea2.RGB, wave='auto' ) # a 1x3 matrix, no materials
product( D65.1nm, neutralMaterial(0.01), Flea2.RGB, wave='auto' ) # a 1x3 matrix, 1 material
path = system.file( "extdata/sources/Lumencor-SpectraX.txt", package='colorSpec' )
lumencor = readSpectra( path, wave=340:660 )
product( lumencor, Flea2.RGB, wave='auto' ) # a 7x3 matrix, no materials

# sequence 6.
scanner = calibrate( scanner )
target = readSpectra( system.file( "extdata/targets/N130501.txt", package='colorSpec' ) )
product( target, scanner, wave='auto' ) # a 288x3 matrix

```

---

ptransform

*make a linear transformation to a colorSpec responder*


---

## Description

apply a linear transformation to a **colorSpec** responder with  $M$  spectra, so that multiples of  $M$  given *primary* vectors are transformed to the standard basis of  $R^M$ . And a given *white* vector is transformed to the  $M$ -vector of all 1s.

The returned object is always `multiply(x,A)` where  $A$  is an internally calculated  $M \times M$  matrix. The name `ptransform` is short for *projective transformation*.

In case of ERROR, a message is logged and NULL returned.

## Usage

```

## S3 method for class 'colorSpec'
ptransform( x, primary, white, digits=Inf )

```

## Arguments

<code>x</code>	a <b>colorSpec</b> responder with $M$ spectra. <code>type(x)</code> must be <code>'responsivity.light'</code> or <code>'responsivity.material'</code> .
<code>primary</code>	an $M \times M$ matrix whose rows define the $M$ primary vectors in the response space of $x$ . It is OK for each row to have a single value that is NA. In this case the NA value is changed so that the sum of the row is 1. This is done because typically the rows represent chromaticities in the response space of $x$ . After this change, the rows of <code>primary</code> must form a basis of $R^M$ . <code>rownames(primary)</code> must be defined; when $M=3$ they are typically <code>c('R', 'G', 'B')</code> .
<code>white</code>	an $M$ -vector in the response space of $x$ , that is typically the ideal white point of a color display. When <code>white</code> is expressed in the basis defined by <code>primary</code> , the coordinates must all be non-zero. <code>white</code> can also be a <b>colorSpec</b> object with a single spectrum suitable as stimulus for $x$ ; in this case the vector <code>white</code> is set to <code>product( white, x, wavelength='auto' )</code> .

`digits` if a positive integer, then `white` is rounded to this number of decimal digits, but in a non-standard way, see **Details**. This is typically done so the internally calculated MxM matrix `A` agrees with that from a color standard, see **Examples**.

### Details

The formal mathematical requirements for `primary` and `white` are:

- The rows of `primary` must form a basis of  $R^M$ . Equivalently, if  $P$  denotes the matrix `primary`, then  $P$  is invertible.
- The coordinates of `white` in this basis are all non-zero. Equivalently, if  $x$  is the solution of  $xP = \text{white}$ , then every component of  $x$  is non-zero.

Assuming both of these are true, then there is a unique matrix  $A$  so that

- $A$  transforms a multiple of the  $i$ 'th row of  $P$  to the  $i$ 'th standard basis vector of  $R^M$ .
- $A$  transforms `white` to the M-vector of all 1s.

This statement is essentially the fundamental theorem of (analytic) projective geometry; see *Bumcrot* page 87, and *Semple* page 398. The rows of  $P$  plus `white` define a *projective frame*; the former are the *fundamental points* and the latter is the *unit point*.

If `digits` is a positive integer, the chromaticity of `white` is computed by dividing `white` by `sum(white)`. The latter must be non-zero, or else it is an ERROR. This chromaticity is rounded to `digits` decimal digits, while preserving the sum of 1. This *rounded chromaticity* is non-zero, and defines a line through 0. The vector `white` is projected onto this line to get the new and rounded `white`, with the rounded chromaticity. See **Examples**.

### Value

a `colorSpec` object equal to `multiply(x,A)` where  $A$  is an internally calculated MxM matrix. The `quantity` and `wavelength` are preserved. The specnames of the returned object are set to `tolower(rownames(primary))`.

The user may want to change the `quantity` of the returned object; see **Examples**.

### References

Bumcrot, Robert J. **Modern Projective Geometry**. Holt, Rinehart, and Winston. 1969.

IEC 61966-2-1:1999. Multimedia systems and equipment - Colour measurement and management. Part 2-1: Colour management - Default RGB colour space - sRGB. <https://webstore.iec.ch/publication/6169>

Semple, J. G. and G. T. Kneebone. **Algebraic Projective Geometry**. Oxford. 1952.

### See Also

[quantity](#), [wavelength](#), [colorSpec](#), [multiply](#), [product](#)

**Examples**

```
##### Example for sRGB #####

# assign the standard sRGB primaries
P = matrix( c(0.64,0.33,NA, 0.3,0.6,NA, 0.15,0.06,NA ), 3, 3, byrow=TRUE )
rownames(P) = c('R','G','B')
P
#   [,1] [,2] [,3]
# R 0.64 0.33  NA
# G 0.30 0.60  NA
# B 0.15 0.06  NA

white = product( D65.1nm, xyz1931.1nm, wave='auto' )
white
#           X           Y           Z
# D65 100.437 105.6708 115.0574

white/sum(white)
#           X           Y           Z
# D65 0.3127269 0.3290232 0.3582499

# But the sRGB standard D65 is xy=(0.3127,0.3290)
# so when the next line is executed,
# the calculated 3x3 matrix will *NOT* agree with the sRGB standard
y = ptransform( xyz1931.1nm, P, white, digits=Inf )

product( D65.1nm, y, wave='auto' )
#   R G B
# D65 1 1 1   # this is exactly what we want, but the internal 3x3 matrix is a little off

# now repeat, but this time round the white chromaticity to
# xy=(0.3127,0.3290) in order to get the matrix right
y = ptransform( xyz1931.1nm, P, white, digits=4 )

rgb = product( D65.1nm, y, wave='auto' )
rgb
#           R           G           B
# D65 1.000238 1.000053 0.999835   # off in the 4'th digit (WARN: this is linear RGB)

255 * rgb
#           R           G           B
# D65 255.0607 255.0134 254.9579   # good enough for 8-bit RGB

65535 * rgb
#           R           G           B
# D65 65550.59 65538.44 65524.18   # NOT good enough for 16-bit RGB

# So for 16-bit RGB, one can get the white RGB right, or the 3x3 matrix right, but not both !

##### Example for ProPhoto RGB #####
```

```

# assign the standard ProPhoto RGB primaries
P = matrix( c(0.7347,0.2653,NA, 0.1596,0.8404,NA, 0.0366,0.0001,NA ), 3, 3, byrow=TRUE )
rownames(P) = c('R','G','B')
P
#      [,1] [,2] [,3]
# R 0.7347 0.2653  NA
# G 0.1596 0.8404  NA
# B 0.0366 0.0001  NA

white = product( D50.5nm, xyz1931.5nm, wave='auto' )
white
#           X           Y           Z
# D50 101.2815 105.042 86.67237

white / sum(white)
#           X           Y           Z
# D50 0.3456755 0.3585101 0.2958144

# but the ProPhoto RGB standard is xy=(0.3457,0.3585); proceed anyway
y = ptransform( xyz1931.5nm, P, white, digits=Inf )

product( D50.5nm, y, wave='auto' )
#      R G B
# D50 1 1 1 # this is exactly what we want, but the internal 3x3 matrix is a little off

# the following line is an equivalent way to compute y.
# pass D50.5nm directly as the 'white' argument
y = ptransform( xyz1931.5nm, P, D50.5nm )

```

---

quantity

*quantity of a colorSpec object*

---

## Description

Retrieve or set the quantity of a **colorSpec** object.

## Usage

```

## S3 method for class 'colorSpec'
quantity(x)

## S3 replacement method for class 'colorSpec'
quantity(x) <- value

## S3 method for class 'colorSpec'
type(x)

```



**Arguments**

x	a <b>colorSpec R</b> object
value	a valid quantity string; see Details.

**Details**

There are 4 valid types, which are further divided into 14 valid quantities. All of these are strings:

For type='light'

quantity = 'energy' (radiometric), or 'photons' (actinometric)

For type='responsivity.light'

quantity = 'energy->electrical', 'energy->neural', 'energy->action',  
'photons->electrical', 'photons->neural', or 'photons->action'

For type='material'

quantity = 'reflectance', 'transmittance', or 'absorbance'

For type='responsivity.material'

quantity = 'material->electrical', 'material->neural', or 'material->action'

Also see the **colorSpec User Guide**.

**Value**

quantity() returns the quantity of x

type() returns the type of x, which depends only on the quantity.

**Note**

The **colorSpec** quantity is more general than the physical SI quantity; for example quantity='energy' really includes 10 distinct SI quantities and maybe more. The unit is left arbitrary in most cases. Exceptions are reflectance, transmittance, and absorbance which are dimensionless.

Changing the quantity should only be done if one knows what one is doing. It does not change the underlying numbers. For example, changing photons to energy does not do numerical conversion. For this specific conversion, see [radiometric\(\)](#).

Similarly, see [linearize\(\)](#) for conversion from absorbance to transmittance.

**See Also**

[colorSpec](#), [radiometric](#), [linearize](#)

---

radiometric

*convert a colorSpec object from actinometric to radiometric*


---

### Description

Convert a **colorSpec** object to have quantity that is radiometric (energy of photons) - to prepare it for colorimetric calculations. Test an object for whether it is radiometric.

### Usage

```
## S3 method for class 'colorSpec'
radiometric( x, multiplier=1, warn=FALSE )
```

```
## S3 method for class 'colorSpec'
is.radiometric( x )
```

### Arguments

<code>x</code>	a <b>colorSpec</b> object
<code>multiplier</code>	a scalar which is multiplied by the output, and intended for unit conversion
<code>warn</code>	if TRUE and a conversion actually takes place, the a WARN message is issued. This makes the user aware of the conversion, so units can be verified. This can be useful when <code>radiometric()</code> is called from another <b>colorSpec</b> function.

### Details

If the `quantity` of `x` does not start with 'photons' then the quantity is not actinometric and so `x` is returned unchanged. Otherwise `x` is actinometric (photon-based).

If `type(x)` is 'light' then the most common actinometric unit of photon count is ( $\mu$ mole of photons) = ( $6.02214 \times 10^{17}$  photons). The conversion equation is:

$$E = Q * 10^{-6} * N_A * h * c / \lambda$$

where  $E$  is the energy of the photons,  $Q$  is the photon count,  $N_A$  is Avogadro's constant,  $h$  is Planck's constant,  $c$  is the speed of light, and  $\lambda$  is the wavelength in meters. The output energy unit is joule.

If the unit of  $Q$  is not ( $\mu$ mole of photons), then the output should be scaled appropriately. For example, if the unit of photon count is exaphotons, then set `multiplier=1/0.602214`.

If the `quantity(x)` is 'photons->electrical', then the most common actinometric unit of responsivity to light is quantum efficiency (QE). The conversion equation is:

$$R_e = QE * \lambda * e / (h * c)$$

where  $R_e$  is the energy-based responsivity,  $QE$  is the quantum efficiency, and  $e$  is the charge of an electron (in C). The output responsivity unit is coulombs/joule (C/J) or amps/watt (A/W).

If the unit of `x` is not quantum efficiency, then `multiplier` should be set appropriately.

If the `quantity(x)` is 'photons->neural' or 'photons->action', the most common actinometric unit of photon count is ( $\mu$ mole of photons) =  $(6.02214 \times 10^{17}$  photons). The conversion equation is:

$$R_e = R_p * \lambda * 10^6 / (N_A * h * c)$$

where  $R_e$  is the energy-based responsivity,  $R_p$  is the photon-based responsivity. This essentially the reciprocal of the first conversion equation.

The argument multiplier is applied to the right side of all the above conversion equations.

### Value

`radiometric()` returns a **colorSpec** object with `quantity` that is radiometric (energy-based) and not actinometric (photon-based). If `type(x)` is a material type ('material' or 'responsivity.material') then `x` is returned unchanged.

If `quantity(x)` starts with 'energy', then `is.radiometric()` returns TRUE, and otherwise FALSE.

### Note

To log the executed conversion equation, execute `cs.options(loglevel='INFO')`.

### Source

Wikipedia. **Photon counting**. [http://en.wikipedia.org/wiki/Photon\\_counting](http://en.wikipedia.org/wiki/Photon_counting)

### See Also

`quantity`, `type`, `F96T12`, `cs.options`, `actinometric`

### Examples

```
sum( F96T12 )      # the step size is 1nm, from 300 to 900nm
# [1] 320.1132  photon irradiance, (micromoles of photons)*m^{-2}*sec^{-1}

sum( radiometric(F96T12) )
# [1] 68.91819  irradiance, watts*m^{-2}
```

---

readCGATS

*read tables from files in ANSI/CGATS.17 format*

---

### Description

The CGATS text format supports a preamble followed by N tables, where  $N \geq 1$ . Each table can have a separate header. A table may or may not contain spectral data, see **Note**. The function converts each table to a `data.frame` with attributes; see **Details**.

### Usage

```
readCGATS( path, collapsesingle=FALSE )
```

## Arguments

- `path` the path name of a single file, in CGATS format
- `collapsesingle` If `path` has only one table ( $N=1$ ) and `collapsesingle` is `TRUE`, then return the single `data.frame` (instead of a list with 1 `data.frame`). If `path` has multiple tables ( $N \geq 2$ ), then `collapsesingle` is ignored.

## Details

The returned list is given attributes: "path", "preamble", and (if present) "date", "created", "originator", and "file\_descriptor". The attribute values are all character vectors. The value of attribute "path" is the argument `path`, and the other values are extracted from "preamble". The length of "preamble" is (typically) greater than 1, and the others have length 1. Each line of the preamble is a keyword-value pair. The keyword `ORIGINATOR` is converted to attribute "originator". The keyword `FILE_DESCRIPTOR` is converted to attribute "file\_descriptor". The keyword `CREATED` is converted to attributes "created" and "date". The list is also given names. If the keyword `TABLE_NAME` is present in the table header, then its value is used. Otherwise the names are "TABLE\_1", "TABLE\_2", ...

Each `data.frame` in the list is assigned attributes: "header", and (if present) "descriptor". The length of "header" is (typically) greater than 1, and "descriptor" has length 1. Each line of the table header is a keyword-value pair. The keywords `DESCRIPTOR` and `TABLE_DESCRIPTOR` are converted to attribute "descriptor".

For the lines between `BEGIN_DATA` and `END_DATA`, two conventions for separating the values are supported:

- In the standard convention, fields are separated by contiguous spaces or tabs, and character strings (which may have embedded spaces or even tabs) are enclosed by double-quotes. This is the convention in the CGATS standard. The function `scan()` is used here.
- In the non-standard convention, fields are separated by a *single* tab, and character strings (which may have embedded spaces but not tabs) are *not* enclosed by double-quotes. This convention is often easier to work with in spreadsheet software. The function `strsplit()` is used here.

The function `readCGATS()` selects the separation convention by examining the line after `BEGIN_DATA_FORMAT`. If this line is split by a single tab and the number of fields matches that given on the `NUMBER_OF_FIELDS` line, then the non-standard convention is selected; otherwise, the standard convention is selected.

## Value

`readCGATS()` returns a list of `data.frames` - one `data.frame` for each table found in `path`. The list and each individual `data.frame` have attributes, see **Details**.

If `path` has only a single table (the majority of files have only 1) and `collapsesingle` is `TRUE`, then the attributes of the list are copied to those of the `data.frame`, and the `data.frame` is then returned. The name of the table is lost.

If there is an error in any table, then the function returns `NULL`.

**Note**

In the BEGIN\_DATA\_FORMAT line(s), field names may not be quoted and may not have embedded spaces.

The CGATS standard allows duplicated field names, and readCGATS() returns them as they appear, with no attempt to append numbers in order to make them unique. Examples of field names which may be duplicated are: SPECTRAL\_NM, SPECTRAL\_DEC, and SPECTRAL\_PCT; for more on these see [readSpectraCGATS\(\)](#).

No attempt is made to recognize those tables that contain spectral data. For conversion of spectral data to **colorSpec** objects, see [readSpectraCGATS\(\)](#).

**References**

ANSI/CGATS.17. Graphic technology - Exchange format for colour and process control data using XML or ASCII text. <https://webstore.ansi.org/> 2009.

ISO/28178. Graphic technology - Exchange format for colour and process control data using XML or ASCII text. <https://www.iso.org/standard/44527.html>. 2009.

CGATS.17 Text File Format. [http://www.colorwiki.com/wiki/CGATS.17\\_Text\\_File\\_Format](http://www.colorwiki.com/wiki/CGATS.17_Text_File_Format).

**See Also**

[readSpectraCGATS](#), [scan](#), [strsplit](#), [names](#)

**Examples**

```
# read file with 2 tables of non-spectral data
A70 = readSpectra( system.file( "tests/A70.ti3", package='colorSpec' ) )
length(A70)          # [1] 2   # the file has 2 tables
ncol( A70[[1]] )    # [1] 7   # the 1st table has 7 columns
ncol( A70[[2]] )    # [1] 4   # the 2nd table has 4 columns
```

---

readSpectra                      *read colorSpec objects from files*

---

**Description**

These functions read **colorSpec** objects from files. In case of ERROR, they return NULL. There are 5 different file formats supported; see **Details**.

**Usage**

```
readSpectra( pathvec, ... )

readSpectraXY( path )
readSpectraSpreadsheet( path )
readSpectrumScope( path )

readSpectraCGATS( path )
```

```
readSpectraControl( path )
```

### Arguments

pathvec	a character vector to (possibly) multiple files. The file extension and a few lines from each file are read and a guess is made regarding the file format.
...	optional arguments passed on to <code>resample()</code> . The most important is <code>wavelength</code> . If these are missing then <code>resample()</code> is not called.
path	a path to a single file with the corresponding format: <code>XYZ</code> , <code>Spreadsheet</code> , <code>Scope</code> , <code>CGATS</code> , or <code>Control</code> . See <code>Details</code> . If the function cannot recognize the format, it returns <code>NULL</code> .

### Details

`readSpectra()` reads the first few lines of the file in order to determine the format, and then calls the corresponding format-specific function. If `readSpectra()` cannot determine the format, it returns `NULL`. The 5 file formats are:

#### XYZ

There is a column header line matching `^(wave|wv?l)` (not case sensitive) followed by the names of the spectra. All lines above this one are taken to be metadata. The separator on this header line can be space, tab, or comma; the line is examined and the separator found there is used in the lines with data below. The organization of the returned object is `'df.col'`. This is probably the most common file format; see the sample file `ciexyz31_1.csv`.

#### Spreadsheet

There is a line matching `"^(ID|SAMPLE|Time)"`. This line and lines below must be tab-separated. Fields matching `'^[A-Z]+([0-9.]+)nm$'` are taken to be spectral data and other fields are taken to be extradata. All lines above this one are taken to be metadata. The organization of the returned object is `'df.row'`. This is a good format for automated acquisition, using a spectrometer, of many spectra. See the sample file `N130501.txt` from **Wolf Faust**.

#### Scope

This is a file format used by **Ocean Optics** spectrometer software. There is a line

```
>>>>Begin Processed Spectral Data<<<<<
```

followed by wavelength and energy separated by a tab. There is only 1 spectrum per file. The organization of the returned object is `'vector'`. See the sample file `pos1-20x.scope`.

#### CGATS

This is a complex format that is best understood by looking at some samples, such as `extdata/objects/Rosco.txt`; see also the **References**. The function `readCGATS()` is first called to get all the tables, and then for each table the column names are examined. There are 2 conventions for presenting the spectral data:

- In the standard convention the fields `SPECTRAL_DEC` or `SPECTRAL_PCT` have the spectral values. The former is the true value, and the latter is the true value x 100. Each value column is preceded a corresponding wavelength column, which has the field name `SPECTRAL_NM`. Note

that these field names are highly duplicated. In principle, this convention allows each record in a CGATS table to have a different wavelength vector. However, this complication is rejected by `readSpectraCGATS()`, which treats it as an `ERROR`.

- In the non-standard convention the field names that match the pattern `"^(nm|SPEC_|SPECTRAL_)[_A-Z]*([0-9.]+"$` are considered to be spectral value data, and other fields are considered extradata. The wavelength is the numerical value of the 2nd parenthesized expression `([0-9.]+"$` in nanometers. Note that every record in this CGATS table has the same wavelength vector. Although this convention is non-standard, it appears in files from many companies, including X-Rite.

If a `data.frame` has spectral data, it is converted to a `colorSpec` object and placed in the returned list. The organization of the resulting `colorSpec` object is `'df.row'`. If the `data.frame` of extradata contains a column `SAMPLE_NAME`, `SAMPLE_ID`, `SampleID`, or `Name`, (examined in that order), then that column is taken to be the `specnames` of the object. If a table has no spectral data, then it is ignored. If the CGATS file has no tables with spectral data, then it is an `ERROR` and the function returns `NULL`.

#### Control

This is a personal format used for digitizing images of plots from manufacturer datasheets and academic papers. It is structured like a `.INI` file. There is a `[Control]` section establishing a simple linear map from pixels to the wavelength and spectrum quantities. Only 3 points are really necessary. It is OK for there to be a little rotation of the plot axes relative to the image. This is followed by a section for each spectrum, in `XY` pixel units only. Conversion to wavelength and spectral quantities is done on-the-fly after they are read. Extrapolation can be a problem, especially when the value is near 0. To force constant extrapolation (see `resample()`), repeat the control point (knot) at the endpoint. See the sample file `Lumencor-SpectraX.txt`. The organization of the returned objects is `'vector'`.

## Value

`readSpectra()` returns a single `colorSpec` object or `NULL` in case of `ERROR`. If there are multiple files in `pathvec` and they cannot be combined using `bind()` because their wavelengths are different, it is an `ERROR`. To avoid this `ERROR`, the `wavelength` argument can be used for resampling to common wavelengths. If there are multiple files, the `organization` of the returned object is `'df.row'` and the first column is the path from which the spectrum was read.

The functions `readSpectraXY()`, `readSpectraSpreadsheet()`, and `readSpectraScope()`, return a single `colorSpec` object, or `NULL` in case of `ERROR`.

The functions `readSpectraCGATS()` and `readSpectraControl()` are more complicated. These 2 file formats can contain multiple spectra with different wavelength sequences, so both functions return a *list* of `colorSpec` objects, even when that list has length 1. If no spectral objects are found, they return `NULL`.

If `readSpectra()` calls `readSpectraCGATS()` or `readSpectraControl()` and receives a list of `colorSpec` objects, `readSpectra()` attempts to `bind()` them into a single object. If they all have the same wavelength vector, then the `bind()` succeeds and the single `colorSpec` object is returned. Otherwise the `bind()` fails, and it is an `ERROR`. To avoid this error `readSpectra()` can be called with a `wavelength` argument. The multiple spectra are resampled using `resample()` and *then* combined using `bind()`, which makes it much more convenient to read such files.

**Note**

During import, each read function tries to guess the quantity from spectrum names or other cues. For example the first line in **N130501.txt** is IT8.7/1, which indicates that the quantity is 'transmittance' (a reflective target is denoted by IT8.7/2). If a confident guess cannot be made, it makes a wild guess and issues a warning. If the quantity is incorrect, one can assign the correct value after import. Alternatively one can add a line to the header part of the file with the keyword 'quantity' followed by some white-space and the correct value. It is OK to put the value in quotes. See example files under folder **extdata**.

**References**

CGATS.17 Text File Format. [http://www.colorwiki.com/wiki/CGATS.17\\_Text\\_File\\_Format](http://www.colorwiki.com/wiki/CGATS.17_Text_File_Format).

ANSI/CGATS.17. Graphic technology - Exchange format for colour and process control data using XML or ASCII text. <https://webstore.ansi.org/> 2009.

ISO/28178. Graphic technology - Exchange format for colour and process control data using XML or ASCII text. <https://www.iso.org/standard/44527.html>. 2009.

**See Also**

[wavelength](#), [quantity](#), [metadata](#), [resample](#), [bind](#), [readCGATS](#)

**Examples**

```
# read file with header declaring the quantity to be "photons->neural"
bird = readSpectra( system.file( "extdata/eyes/BirdEyes.txt", package='colorSpec' ) )
quantity(bird) # [1] "photons->neural"
```

---

resample

*resample a colorSpec Object to new wavelengths*

---

**Description**

interpolate or smooth to new wavelengths. Simple extrapolation and clamping is also performed.

**Usage**

```
## S3 method for class 'colorSpec'
resample( x, wavelength, method='auto', span=0.02, extrapolation='const', clamp='auto' )
```



**Arguments**

x	a <b>colorSpec</b> object
wavelength	vector of new wavelengths, in nanometers
method	interpolation methods available are 'sprague', 'spline', and 'linear'. Also available is 'auto' which means to use 'sprague' if x is regular, and 'spline' otherwise. An available smoothing method is 'loess'. See <b>Details</b> .
span	smoothing argument passed to <a href="#">loess()</a> during interpolation, and not used by other methods. The default value span=0.02 is suitable for .scope spectra but may be too small in many other cases.
extrapolation	extrapolation methods available are 'const' and 'linear'. These can be abbreviated to the initial letter. Also available is a numeric value, which is used for simple padding. See <b>Details</b> . Also available is a vector or list of length 2 that combines 2 of the above. The first item is used on the low side (shorter wavelengths), and the second item is used on the high side (longer wavelengths).
clamp	clamp methods available are 'auto', TRUE, and FALSE. Also available is a numeric vector of length 2, which defines the clamping interval. See <b>Details</b> .

**Details**

If method is 'sprague', the quintic polynomial in *De Kerf* is used. Six weights are applied to nearby data values: 3 on each side. The 'sprague' method is only supported when x is regular.

If method is 'spline', the function `stats::spline()` is called with `method='natural'`.

The 'spline' method is supported even when x is irregular.

If method is 'linear', the function `stats::approx()` is called. Two weights are applied to nearby data values: 1 on each side. The 'linear' method is supported even when x is irregular.

If method is 'loess', the function `stats::loess()` is called with the given span parameter. Smoothing is most useful for noisy data, e.g. raw data from a spectrometer. I have found that span=0.02 works well for Ocean Optics .scope files, but this may be too small in other cases, which triggers an error in `stats::loess()`. The 'loess' method is supported even when x is irregular.

If extrapolation is 'const', the extreme values at each end are simply extended. If extrapolation is 'linear', the line defined by the 2 extreme values at each end is used for extrapolation. If the ultimate and penultimate wavelengths are equal, then this line is undefined and the function reverts to 'const'.

If clamp is 'auto', output values are clamped to the physically realizable interval appropriate for x. This is the interval [0,1] when `quantity(x)` is 'reflectance' or 'transmittance', and the interval [0,∞) otherwise. Exception: If an input spectrum in x violates a limit, then clamping the output spectrum to this limit is *NOT* enforced. This happens most frequently for theoretical (or matrixed) cameras, such as [BT.709.RGB](#).

If clamp is TRUE, the result is the same as 'auto', but with no exceptions. If clamp is FALSE, then no clamping is done.

If clamp is a numerical interval, then clamping is done to that interval, with no exceptions. The two standard intervals mentioned above can be expressed in **R** as `c(0, 1)` and `c(0, Inf)` respectively.

**Value**

`resample(x)` returns a **colorSpec** object with the new wavelength. Other properties, e.g. [organization](#), [quantity](#), ..., are preserved.

In case of ERROR, the function returns NULL.

**References**

De Kerf, Joseph L. F. **The interpolation method of Sprague-Karup**. *Journal of Computational and Applied Mathematics*. volume I, no 2, 1975. equation (S).

**See Also**

[organization\(\)](#), [quantity\(\)](#), [wavelength\(\)](#), [is.regular\(\)](#), [theoreticalRGB](#), [spline\(\)](#), [approx\(\)](#), [loess](#)

**Examples**

```
path = system.file( "extdata/sources/pos1-20x.scope", package='colorSpec' )
y = readSpectra( path )
# plot noisy data in gray
plot( y, col='gray' )
# plot smoothed plot in black on top of the noisy one to check quality
plot( resample( y, 200:880, meth='loess', span=0.02 ), col='black', add=TRUE )
```

---

scanner

*standard RGB scanners*

---

**Description**

`scanner.ACES` is an RGB responder to material; an ACES/SMPTE standard for scanning RGB film. The 3 spectra are defined from 368 to 728 nm, at 2nm intervals.

**Format**

A **colorSpec** object with [quantity](#) equal to 'material->electrical' and 3 spectra: r, g, and b.

**Details**

The responsivities have been scaled (by [calibrate](#)) so the response to the *perfect transmitting filter* (PTF) is RGB=(1,1,1).

## References

Technical Bulletin TB-2014-005. Informative Notes on SMPTE ST 2065-2 - Academy Printing Density (APD). Spectral Responsivities, Reference Measurement Device and Spectral Calculation. SMPTE ST 2065-3 Academy Density Exchange Encoding (ADX). Encoding Academy Printing Density (APD) Values.

The Academy of Motion Picture Arts and Sciences. Science and Technology Council. Academy Color Encoding System (ACES) Project Committee. Version 1.0 December 19, 2014. Annex A Spectral Responsivities.

## See Also

[quantity, calibrate](#)

## Examples

```
# compute response of ACES scanner to the Hoya filters
product( Hoya, scanner.ACES, wave='auto' )

##           R           G           B
## R-60  0.902447043 2.022522e-05 0.00000000
## G-533 0.038450857 4.900983e-01 0.05431134
## B-440 0.008466317 1.686241e-02 0.42863320
## LB-120 0.184408941 3.264111e-01 0.53492533
```

---

sectionOptimalColors *compute sections of an optimal color surface by hyperplanes*

---

## Description

Consider a **colorSpec** object  $x$  with type equal to 'responsivity.material'. The set of all possible material reflectance functions (or transmittance functions) is convex, closed, and bounded (in fact they form a cube), and this implies that the set of all possible output responses from  $x$  is also convex, closed, and bounded. The latter set is called the *object-color solid* or *Rösch Farbkörper* for  $x$ . If the dimension of the response of  $x$  is 2, this solid is a convex polygon that is centrally symmetric - a *zonogon*. If the dimension of the response of  $x$  is 3 (e.g. RGB or XYZ), this solid is a special type of centrally symmetric convex polyhedron called a *zonohedron*, see *Centore*. This function only supports dimensions 2 and 3. Denote this object-color solid by  $\mathbf{Z}$ .

A color on the boundary of  $\mathbf{Z}$  is called an *optimal color*. Consider the intersection of a hyperplane with the boundary of  $\mathbf{Z}$ . Let the equation of the hyperplane be given by:

$$\langle v, normal \rangle = \beta$$

where *normal* is orthogonal to the hyperplane, and  $\beta$  is the plane constant, and  $v$  is a variable vector. The purpose of the function `sectionOptimalColors()` is to compute the intersection set.

In dimension 2 this hyperplane is a line, and the intersection is generically 2 points, and 1 point if the line only intersects the boundary (we ignore the special case when the intersection is an edge of the polygon).

In dimension 3 this hyperplane is a 2D plane, and the intersection is generically a polygon, and 1 point if the line only intersects the boundary (we ignore the special case when the intersection is a face of the zonohedron).

Of course, the intersection can also be empty.

### Usage

```
## S3 method for class 'colorSpec'
sectionOptimalColors( x, normal, beta )
```

### Arguments

x	a <b>colorSpec</b> object with type equal to 'responsivity.material' and M spectra, where M=2 or 3.
normal	a nonzero vector of dimension M, that is the normal to a hyperplane
beta	a vector of numbers of positive length. The number beta[k] defines the plane $\langle v, normal \rangle = beta[k]$ .

### Details

Consider first the case that the dimension of x is 3, so that **Z** is a zonohedron. In the preprocessing phase the zonohedral representation is calculated. The faces of **Z** are either parallelograms, or *compound faces* that are partitioned into parallelograms. The centers of all these parallelograms are computed, along with their extent in direction *normal*. For a given plane  $\langle v, normal \rangle = \beta$ , the parallelograms that intersect the plane are extracted. The boundary of each parallelogram intersects the plane in 2 points (in general) and one of those points is computed. The set of all these points is then sorted into proper order around the boundary.

In the case that the dimension of x is 2, so that **Z** is a zonogon, the parallelograms are replaced by line segments (edges), and the processing is much easier.

### Value

The function returns a list with an item for each value in vector beta. Each item in the output is a list with these items:

beta	the value of the plane constant $\beta$
section	an NxM matrix, where N is the number of points in the section, and M is the dimension of normal. If the intersection is empty, then N=0.

In case of global error, the function returns NULL.

### WARNING

The preprocessing calculation of the zonohedron dominates the total time. And this time goes up rapidly with the number of wavelengths. We recommend using a wavelength step of 5nm, as in the **Examples**. For best results, batch a lot of betas into a single function call and then process the output.

Moreover, the preprocessing time is dominated by the partitioning of the compound faces into parallelograms. This is made worse by an x whose spectral responses have little overlap, as in

`scanner.ACES`. In these cases, try a larger step size, and then reduce. Optimizing these compound faces is a possible topic for the future.

## References

Centore, Paul. *A Zonohedral Approach to Optimal Colours*. **Color Research & Application**. Vol. 38. No. 2. pp. 110-119. April 2013.

Logvinenko, A. D. An object-color space. **Journal of Vision**. 9(11):5, 1-23, (2009). <https://jov.arvojournals.org/article.aspx?articleid=2203976>. doi:10.1167/9.11.5.

## See Also

vignette `optimals`, `probeOptimalColors()`

## Examples

```

wave = seq(420,680,by=5)
Flea2.scanner = product( A.1nm, "material", Flea2.RGB, wavelength=wave )
seclist = sectionOptimalColors( Flea2.scanner, normal=c(0,1,0), beta=10 )
length( seclist[[1]]$section )
seclist[[1]]$section[ 1:5, ]
## [1] 207 # the polygon has 207 vertices, and the first 5 are:
##           Red Green   Blue
## [1,] 109.2756   10 3.5391342
## [2,] 109.5729   10 2.5403628
## [3,] 109.8078   10 1.7020526
## [4,] 109.9942   10 1.0111585
## [5,] 110.1428   10 0.4513051

```

---

solar.irradiance

*Standard Solar Irradiance - Extraterrestrial and Terrestrial*

---

## Description

solar.irradiance Three power spectra; from 280 to 1000 nm at 1 nm intervals. The unit is  $W * m^{-2} * nm^{-1}$ .

atmosphere2003 a transmittance spectrum = the quotient of 2 spectra from solar.irradiance

## Format

solar.irradiance is a **colorSpec** object with quantity equal to 'energy' and with 3 spectra:

AirMass.0 Extraterrestrial Radiation (solar spectrum at top of atmosphere) at mean Earth-Sun distance

GlobalTilt spectral radiation from solar disk plus sky diffuse and diffuse reflected from ground on south facing surface tilted 37 deg from horizontal

AirMass.1.5 the sum of Direct and Circumsolar irradiance, when the optical path is 1.5 times that of the sun at zenith, see **Details**

atmosphere2003 is a **colorSpec** object with quantity equal to 'transmittance' and with 1 spectrum:

AirMass.1.5 the quotient  $\text{AirMass.1.5} / \text{AirMass.0}$  from solar.irradiance

### Details

**Direct** is Direct Normal Irradiance Nearly parallel (0.5 deg divergent cone) radiation on surface with surface normal tracking (pointing to) the sun, excluding scattered sky and reflected ground radiation.

**Circumsolar** is Spectral irradiance within +/- 2.5 degree (5 degree diameter) field of view centered on the 0.5 deg diameter solar disk, but excluding the radiation from the disk.

### Note

The reference spectra in ASTM G173-03 are designed for Photovoltaic Performance Evaluation.

The original wavelength sequence in ASTM G173-03 is irregular. The interval is 0.5 nanometer from 280 to 400 nm, 1 nm from 400 to 1700 nm, an intermediate wavelength at 1702 nm, and 5 nm from 1705 to 4000 nm. To create the object solar.irradiance with a regular step size, the original was resampled from 280 to 1000 nm at 1nm intervals.

### Source

Reference Solar Spectral Irradiance: ASTM G-173. <http://rredc.nrel.gov/solar/spectra/am1.5/astmg173/astmg173.html>

### References

ASTM G173-03 Reference Spectra Derived from SMARTS v. 2.9.2.  
Standard Tables for Reference Solar Spectral Irradiances: Direct Normal and Hemispherical on 37-deg Tilted Surface (2003)

### See Also

[D65](#), [D50](#), [daylightSpectra](#), [resample](#), vignette **Blue Flame and Green Comet**

---

specnames

*specnames of a colorSpec object*

---

### Description

Retrieve or set the specnames of a **colorSpec** object. Retrieve the number of spectra.

**Usage**

```
## S3 method for class 'colorSpec'
specnames(x)

## S3 replacement method for class 'colorSpec'
specnames(x) <- value

## S3 method for class 'colorSpec'
numSpectra(x)
```

**Arguments**

x                    a **colorSpec** R object  
value                a character vector with length equal to the number of spectra in x.

**Details**

If the organization of x is "vector" then x is a vector and value is a single string, which is stored as attr(x, 'specname').

If the organization of x is "matrix", then x is a matrix and value is stored as colnames(x).

If the organization of x is "df.col", then x is a data.frame with N+1 columns, where N is the number of spectra. value is stored as colnames(x)[2:(N+1)].

If the organization of x is "df.row", then x is a data.frame and value is stored as row.names(x).

**Value**

specnames() returns a character vector with the names of the spectra.

numSpectra(x) is equal to length(specnames(x)) but much more efficient.

**See Also**

[rownames](#), [colnames](#)

---

standardRGB

*Convert from XYZ to some standard RGB spaces*

---

**Description**

To display an XYZ value, it typically must be converted to a standard RGB space. This is the function to do it.

**Usage**

```
RGBfromXYZ( XYZ, space )
```

**Arguments**

XYZ	a 3-vector, or a matrix with 3 columns with XYZs in the rows
space	the name of the RGB space - either 'sRGB' or 'Adobe RGB'. The match is case-insensitive, and spaces in the string are ignored.

**Details**

The input XYZ is multiplied by the appropriate 3x3 conversion matrix (for **sRGB** or **Adobe RGB**). These matrices are taken from *Lindbloom* and not from the corresponding *Wikipedia* articles; for the reason why see **Note**.

**Value**

An Mx3 matrix where M is the number of rows in XYZ, or M=1 if XYZ is a 3-vector. Each row of the returned matrix is filled with linear RGB in the appropriate RGB space. Values outside the unit cube are not clamped. To compute non-linear display RGB, see [DisplayRGBfromLinearRGB\(\)](#). In case of error the function returns NULL.

**WARNING**

This function is deprecated. New software should use `spacesRGB::RGBfromXYZ()` instead. The new function returns "signal RGB" instead of linear RGB.

**Note**

An RGB space is normally defined by the xy chromaticities of the 3 primaries and the white point. We follow *Lindbloom* in using the 'official' XYZ of the white point from ASTM E308. Using this XYZ of the white point makes the color space a little more consistent with other areas of color. For example, from IEC 61966-2-1 we have D65 xyY=(0.3127,0.3290,1) -> XYZ=(0.9504559,1,1.0890578). But from ASTM E308, D65 XYZ=(0.95047,1,1.08883), which is a little different.

**Source**

IEC 61966-2-1:1999. Multimedia systems and equipment - Colour measurement and management. Part 2-1: Colour management - Default RGB colour space - sRGB. <http://webstore.iec.ch/publication/6169>

Lindbloom, Bruce. RGB/XYZ Matrices. [http://brucelindbloom.com/index.html?Eqn\\_RGB\\_XYZ\\_Matrix.html](http://brucelindbloom.com/index.html?Eqn_RGB_XYZ_Matrix.html)

Wikipedia. **sRGB**. <http://en.wikipedia.org/wiki/SRGB>

Wikipedia. **Adobe RGB**. [http://en.wikipedia.org/wiki/Adobe\\_RGB\\_color\\_space](http://en.wikipedia.org/wiki/Adobe_RGB_color_space)

**See Also**

[D65](#), [officialXYZ](#), [DisplayRGBfromLinearRGB](#)



**Examples**

```

RGBfromXYZ( officialXYZ('D65'), 'sRGB' )
#      R G B
# [1,] 1 1 1   # not really 1s, but difference < 1.e-7

RGBfromXYZ( c(.3127,0.3290,0.3583)/0.3290, 'sRGB' )
#      R      G      B
# [1,] 0.9998409 1.000023 1.00024   difference > 1.e-5

```

---

subset	<i>extract a subset of a colorSpec Object</i>
--------	---

---

**Description**

extract a subset of the spectra in a **colorSpec** object.  
The subset can be specified by indexes, by a logical vector, or by a regular expression matching the [specnames](#)

**Usage**

```

## S3 method for class 'colorSpec'
subset( x, subset, ... )

```

**Arguments**

x	a <b>colorSpec</b> object
subset	an integer vector, a logical vector, or a regular expression
...	additional arguments ignored

**Details**

If subset is an integer vector, each integer must be between 1 and M, where M the number of spectra in x. No duplicates are allowed. The number of spectra returned is equal to length(subset). It is OK for the length to be 0, in which case the function returns the empty subset.

If subset is a logical vector, its length must be equal to M. The number of spectra returned is equal to the number of TRUEs in subset.

If subset is a regular expression, the number of spectra returned is equal to the number of specnames(x) matched by the expression.

**Value**

subset(x) returns a **colorSpec** object with the same [organization](#) as x. Exception: if the organization of x is 'vector' and the subset is empty, then the returned object is a matrix with 0 columns.

**Note**

subset() can also be used for re-ordering the spectra; just set argument subset to the desired permutation vector.

**See Also**

[organization](#)

**Examples**

```
tritanope = subset( lms2000.1nm, 1:2 ) # keep long and medium cone fundamentals, but drop the short
sml2000.1nm = subset( lms2000.1nm, 3:1 ) # reorder from short to long
```

---

theoreticalRGB	<i>Theoretical RGB Cameras - BT.709.RGB, Adobe.RGB, and ACES.RGB</i>
----------------	--

---

**Description**

These are 3 built-in **colorSpec** objects, with [quantity](#) equal to 'energy->electrical'.

BT.709.RGB a theoretical RGB responder to light. The 3 responsivity spectra are constructed so that the RGBs from this theoretical camera, when displayed on an sRGB display after correct EOTF adjustment, would emit light with the same XYZs as the captured scene (up to a constant multiple). All three responsivities have negative lobes.

Adobe.RGB a theoretical RGB responder to light. The 3 responsivity spectra are constructed so that the RGBs from this theoretical camera, when displayed on an Adobe RGB display after correct EOTF adjustment, would emit light with the same XYZs as the captured scene (up to a constant multiple). All three responsivities have negative lobes.

ACES.RGB a theoretical RGB responder to light. Unlike the two above cameras, the responsivities are non-negative and so this camera could be built, in principle. These are the ACES RICD (Reference Input Capture Device) spectral sensitivities.

**Format**

All are **colorSpec** objects with [quantity](#) equal to 'energy->electrical' and 3 spectra: r, g, and b. The wavelengths are 360 to 830 nm at 1 nm intervals.

**Details**

All responsivity spectra are linear combinations of the spectra in [xyz1931.1nm](#). These 3 theoretical cameras satisfy the *Maxwell-Ives criterion* by construction.

For BT.709.RGB and Adobe.RGB, the responsivities are scaled so the response to [D65.1nm](#) is RGB=(1,1,1). These responsivities have negative lobes.

The BT.709 primaries and white point are the same as those of sRGB (though the EOTF functions are different).

Adobe RGB and sRGB share the same Red, Blue, and White chromaticities, and only differ in the Green. This implies that for both cameras the Green output is 0 at Red and Blue, and 1 at White. This in turn implies that the Green output is identical for both cameras for all input spectra, and so the Green responsivity spectra are identical for both cameras.

For ACES.RGB the responsivities are area normalized as in Annex C of S-2008-001. They are scaled so that the response to Illuminant E is RGB=(1,1,1). For an example of white-balancing, as in Annex B, see the examples below.

## References

Poynton, Charles. **Digital Video and HD - Algorithms and Interfaces**. Morgan Kaufmann. Second Edition. 2012. Figure 26.5 on page 302.

**Academy Color Encoding Specification (ACES)**. S-2008-001. 2011. Annex B, pp. 23-25. Annex C, pp. 26-33.

## See Also

[quantity\(\)](#), [D65.1nm](#), [xyz1931.1nm](#), [ptransform\(\)](#), [calibrate\(\)](#), [vignette blueflame](#)

## Examples

```
##### BT.709.RGB is created using the following recipe #####
P = matrix( c(0.64,0.33,NA, 0.3,0.6,NA, 0.15,0.06,NA ), 3, 3, byrow=TRUE )
rownames(P) = c('R','G','B')
BT.709.RGB = ptransform( xyz1931.1nm, P, D65.1nm )
quantity(BT.709.RGB) = "energy->electrical"

##### Adobe.RGB recipe is the same, except for the matrix P #####
P = matrix( c(0.64,0.33,NA, 0.21,0.71,NA, 0.15,0.06,NA ), 3, 3, byrow=TRUE )
rownames(P) = c('R','G','B')
Adobe.RGB = ptransform( xyz1931.1nm, P, D65.1nm )
quantity(Adobe.RGB) = "energy->electrical"

##### white-balancing ACES.RGB for CIE Standard Illuminant D60 #####
# in a scene illuminated by daylight illuminant D6003,
# and with a perfect-reflecting-diffuser in that scene,
# object 'camera1' would have response RGB=(1,1,1) for that diffuser.
D6003 = daylightSpectra( 6000*1.4388/1.4380, wavelength=wavelength(ACES.RGB) )
camera1 = calibrate( ACES.RGB, D6003, 1, method='scaling' )
```

---

wavelength

*wavelength vector of a colorSpec object*

---

## Description

Retrieve or set the wavelengths of a **colorSpec** object. Retrieve the number of wavelengths, and whether the wavelength sequence is regular.

## Usage

```
## S3 method for class 'colorSpec'  
wavelength(x)  
  
## S3 replacement method for class 'colorSpec'  
wavelength(x) <- value  
  
## S3 method for class 'colorSpec'  
numWavelengths(x)  
  
## S3 method for class 'colorSpec'  
is.regular(x)  
step.wl(x)
```

## Arguments

x	a <b>colorSpec</b> R object
value	a numeric vector with length equal to the number of wavelengths in x. The wavelengths must be increasing. The unit must be nanometers.

## Details

If the organization of x is 'df.col', then x is a data.frame and the wavelength vector is stored in the first column of x.  
Otherwise, the wavelength vector is stored as attr(x, 'wavelength').

## Value

wavelength() returns a numeric vector with the wavelengths of the spectra, in nanometers.  
numWavelengths(x) is equal to length(wavelength(x)) but much more efficient.  
is.regular() returns TRUE or FALSE, depending on whether the step between consecutive wavelengths is a constant. A truncation error of 1.e-6 nm is tolerated here. For example, the X-Rite ColorMunki spectrometer in hi-res mode has a step of 3.33333nm, and it is considered regular.  
step.wl() returns the *mean* step in nm, whether the wavelengths are regular or not.

## See Also

[colorSpec](#)

## Description

xyz1931.1nm the 1931 2° functions from 360 to 830 nm, at 1nm intervals  
xyz1931.5nm the 1931 2° functions from 380 to 780 nm, at 5nm intervals

### Format

Each is a **colorSpec** object organized as a matrix with 3 variables.

x the x-bar responsivity function  
y the y-bar responsivity function  
z the z-bar responsivity function

### Source

<http://www.cvr1.org>

### References

Günther Wyszecki and W.S. Stiles. **Color Science : Concepts and Methods, Quantitative Data and Formulae**. Second Edition. Wiley-Interscience. 1982. Table I(3.3.1). pp. 723-735.

ASTM E 308 - 01. Standard Practice for Computing the Colors of Objects by Using the CIE System. Table 1

### See Also

[xyz1964](#)

### Examples

```
summary(xyz1931.1nm)
white.point = product( D65.1nm, xyz1931.1nm, wave='auto' )
```

---

xyz1964

*CIE Color Matching Functions - 10-degree (1964)*

---

### Description

xyz1964.1nm the 10° 1964 functions from 360 to 830 nm, at 1nm intervals  
xyz1964.5nm the 10° 1964 functions from 380 to 780 nm, at 5nm intervals

**Format**

Each is a **colorSpec** object organized as a matrix with 3 columns.

x	the x-bar responsivity function
y	the y-bar responsivity function
z	the z-bar responsivity function

**Source**

<http://www.cvrl.org>

**References**

Günther Wyszecki and W.S. Stiles. **Color Science : Concepts and Methods, Quantitative Data and Formulae**. Second Edition. Wiley-Interscience. 1982. Table I(3.3.1). pp. 723-735.

ASTM E 308 - 01. Standard Practice for Computing the Colors of Objects by Using the CIE System. Table 1

**See Also**

[xyz1931](#)

**Examples**

```
summary(xyz1964.1nm)
white.point = product( D65.1nm, xyz1964.1nm, wave='auto' )
```

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