

Using R for photobiological calculations

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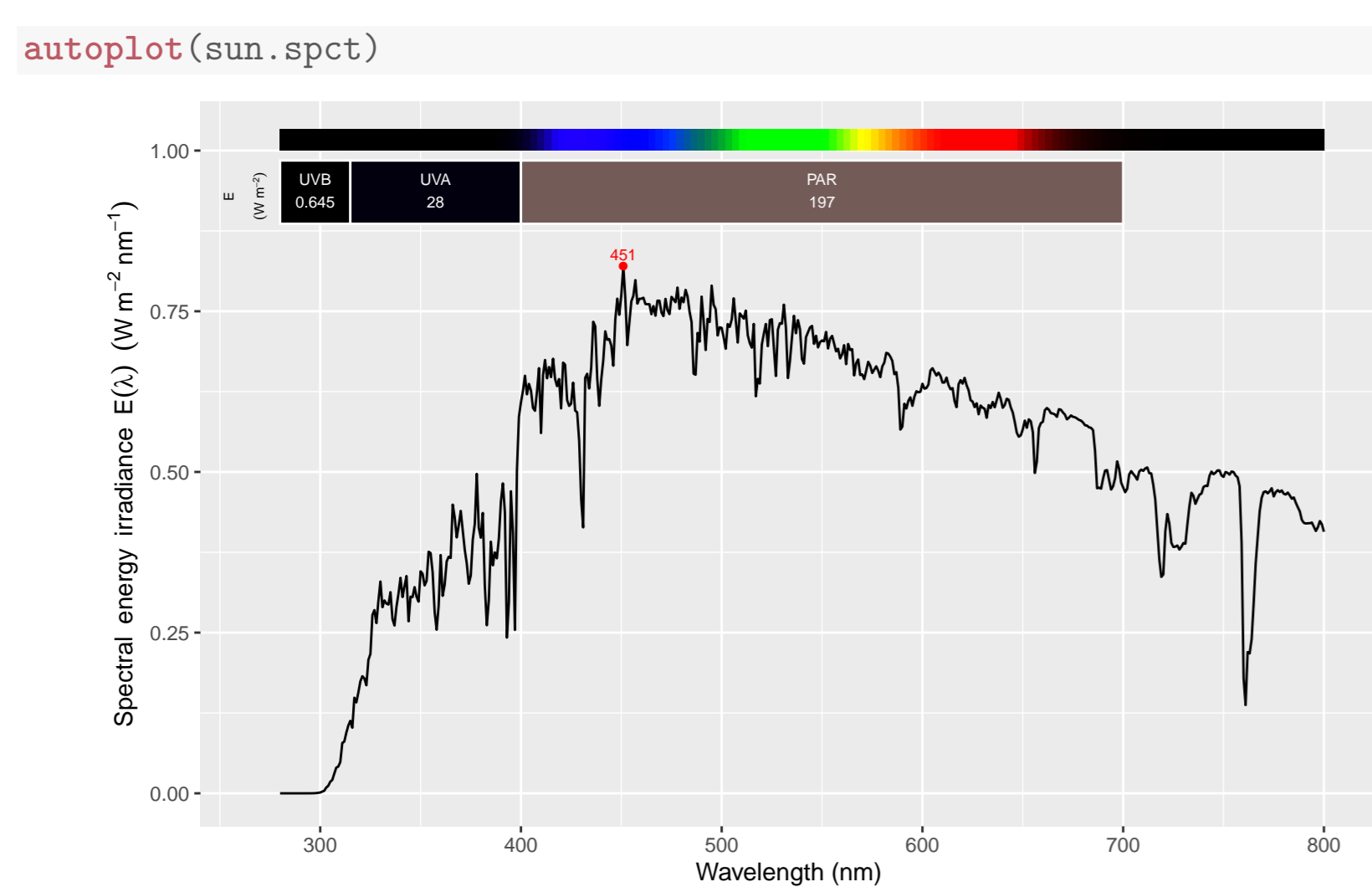
1. Introduction

- Here I showcase my R packages for photobiology.
- In the examples I also use data included in them.
- Several of the packages contain only data.
- Most of the packages are published in the CRAN repository, open-source and regularly updated.

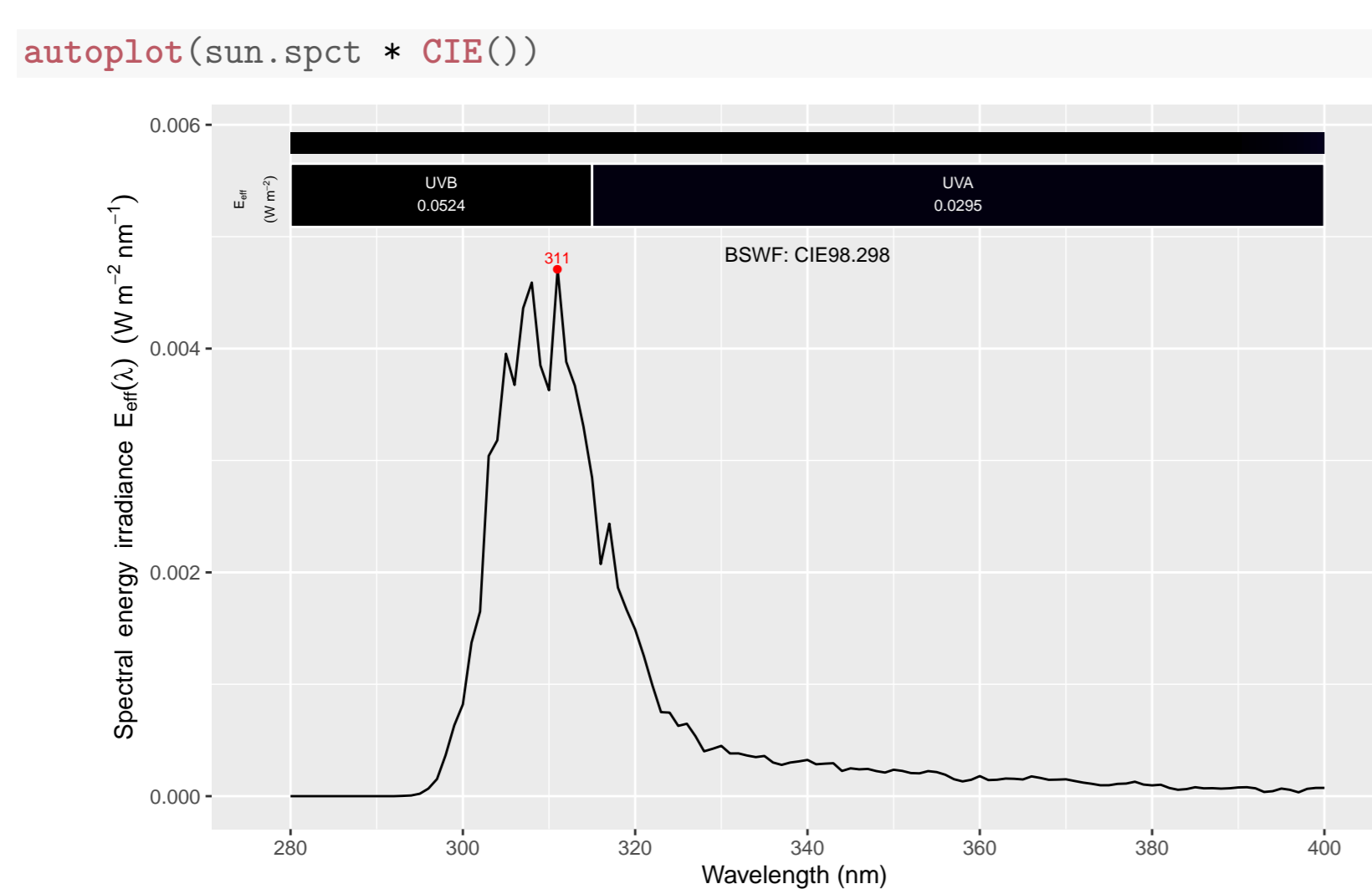
Repository	Package	Type	Contents
CRAN	photobiology	funs + classes	functions, class definitions and methods
CRAN	photobiologyInOut	functions	data import/export
CRAN	photobiologyWavebands	definitions	quantification of radiation
CRAN	ggspectra	methods	plotting of spectral data
CRAN	photobiologySun	data	solar and daylight
CRAN	photobiologyLamps	data	emission by light bulbs
CRAN	photobiologyLEDs	data	emission by LEDs
CRAN	photobiologyFilters	data	transmission of filters
CRAN	photobiologySensors	data	response of sensors
Bitbucket	photobiologyReflectors	data	reflection by materials
CRAN	photobiologyPlants	funs + data	plants' responses
Bitbucket	rOmniDriver	driver API	Ocean Optics spectrometers
Bitbucket	oacquire	data acquisition	Ocean Optics spectrometers

2. Plotting spectra

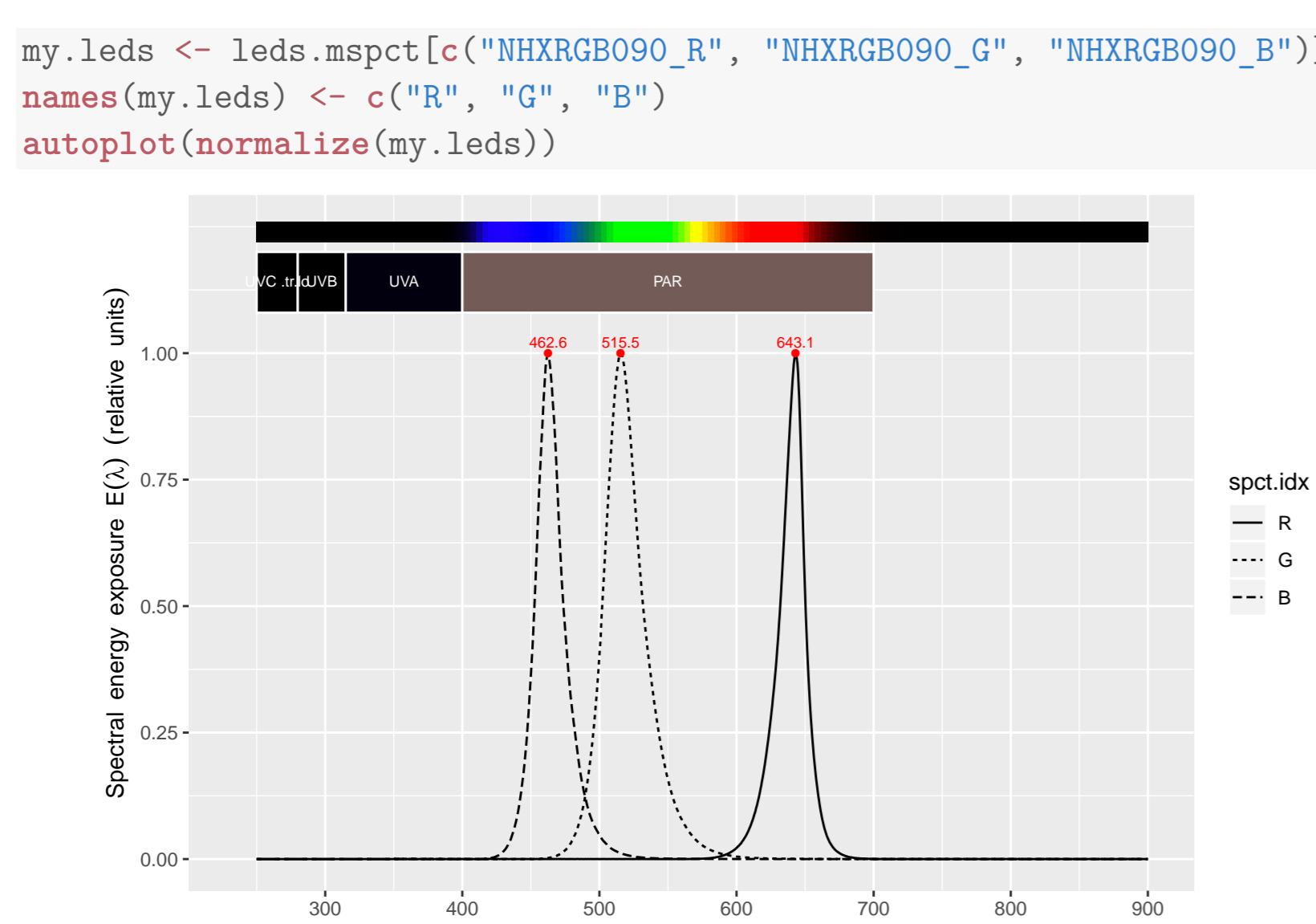
Units used when data is stored are consistent, and saved as metadata allowing their propagation to plot axis-labels and annotations.



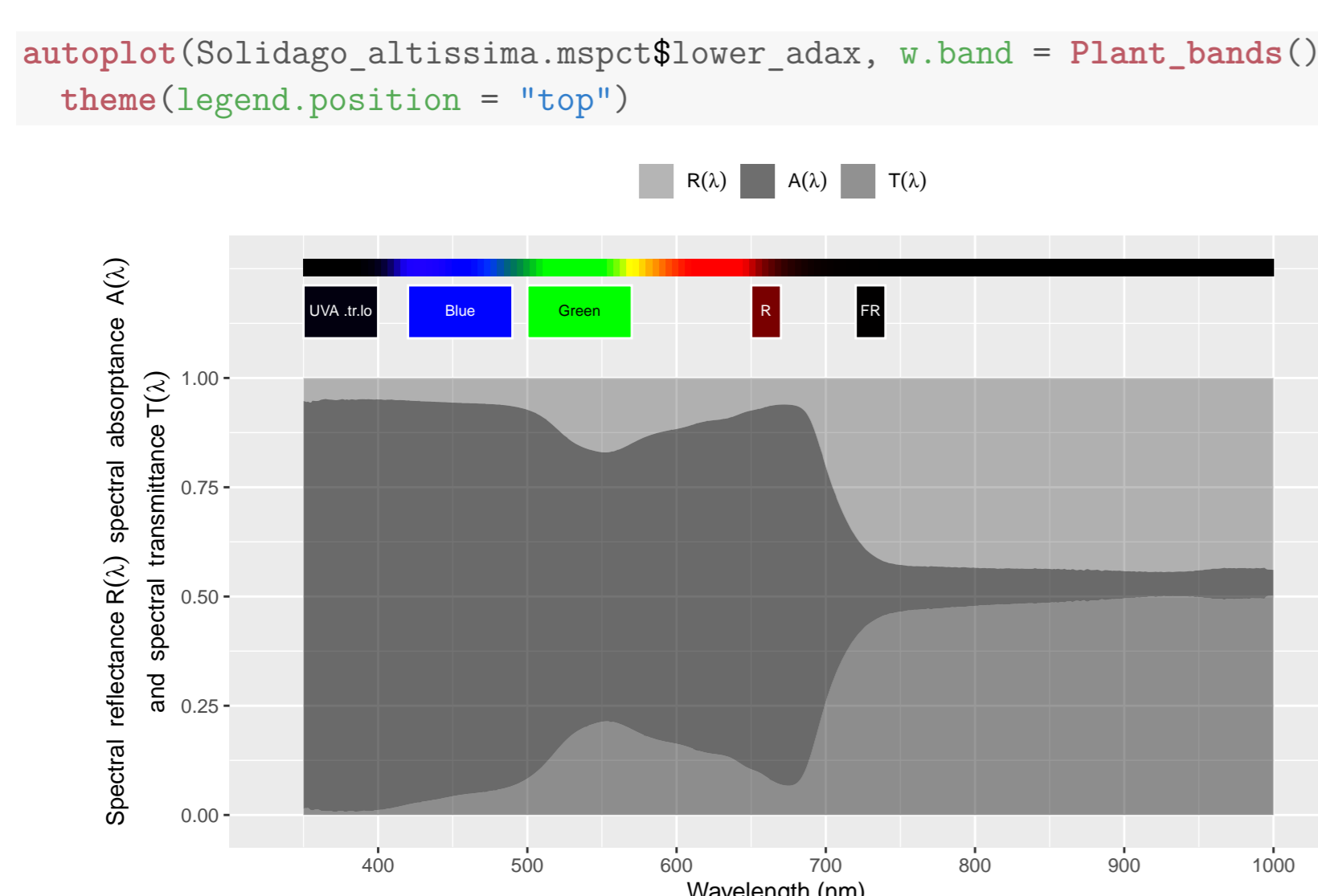
We convolute a radiation spectrum by multiplication with a weighting function.



Normalized spectral photon irradiance for the three channels of a LED array.



Spectral transmittance and reflectance for a leaf, with annotations corresponding to wavebands of interest in plant biology.



3. Irradiance

In this section we show examples of the calculation of energy and photon (quantum) irradiance, both weighted and unweighted, and for the whole spectrum or individual regions.

```
e_irrad(sun.spct)
## Total
## 269.1249
## attr(,"time.unit")
## [1] "second"
## attr(,"radiation.unit")
## [1] "energy irradiance total"
```

```
e_irrad(sun.spct, CIE())
## CIE98.298.tr.lo
## 0.08181583
## attr(,"time.unit")
## [1] "second"
## attr(,"radiation.unit")
## [1] "energy irradiance total"
```

```
e_irrad(sun.spct, VIS_bands())
## Purple.ISO Blue.ISO Green.ISO Yellow.ISO Orange.ISO Red.ISO
## 47.75529 37.55207 49.26860 13.67971 12.00432 79.38159
## attr(,"time.unit")
## [1] "second"
## attr(,"radiation.unit")
## [1] "energy irradiance total"
```

In the packages, S.I. units without 'multipliers' are used for radiation quantities. In this case we multiply the returned value, which is expressed in $\text{mol m}^{-2} \text{s}^{-1}$, by 10^6 to obtain the more familiar units $\mu\text{mol m}^{-2} \text{s}^{-1}$.

```
q_irrad(sun.spct, PAR()) * 1e6
## PAR
## 894.1352
## attr(,"time.unit")
## [1] "second"
## attr(,"radiation.unit")
## [1] "photon irradiance total"
```

4. Fluence

```
fluence(sun.spct, exposure.time = duration(12, "minutes"))
## Total
## 193769.9
## attr(,"radiation.unit")
## [1] "energy fluence (J m⁻²)"
## attr(,"exposure.duration")
## [1] "720s (~12 minutes)"
```

5. Ratios

Photon ratios can be calculated for any pair of arbitrary regions of the spectrum. Here we use predefined wavebands, but the user can easily define new ones as needed.

```
q_ratio(sun.spct, UVB(), PAR())
## UVB.ISO: PAR(q:q)
## 0.001873724
## attr(,"radiation.unit")
## [1] "q:q ratio"
```

Light source spectra can be convoluted with transmittance spectra as a simple multiplication operation. Interpolation is automatic when the wavelength values are not matched between the two spectra.

```
q_ratio(sun.spct * polyester.spct, UVB(), PAR())
## UVB.ISO: PAR(q:q)
## 5.142362e-05
## attr(,"radiation.unit")
## [1] "q:q ratio"
```

For R:FR ratio a convenience function is available.

```
R_FR(sun.spct)
## Red.Smith10: FarRed.Smith10(q:q)
## 1.266704
## attr(,"radiation.unit")
## [1] "q:q ratio"
```

Photon : energy ratios are also easily calculated. As the result is in mol J^{-1} we multiply it by 10^6 to obtain $\mu\text{mol J}^{-1}$.

```
qe_ratio(sun.spct, list(Blue(), Red())) * 1e6
## q:e( Blue.ISO) q:e( Red.ISO)
## 3.968591 5.682783
## attr(,"radiation.unit")
## [1] "q:e ratio"
```

6. Collections of spectra

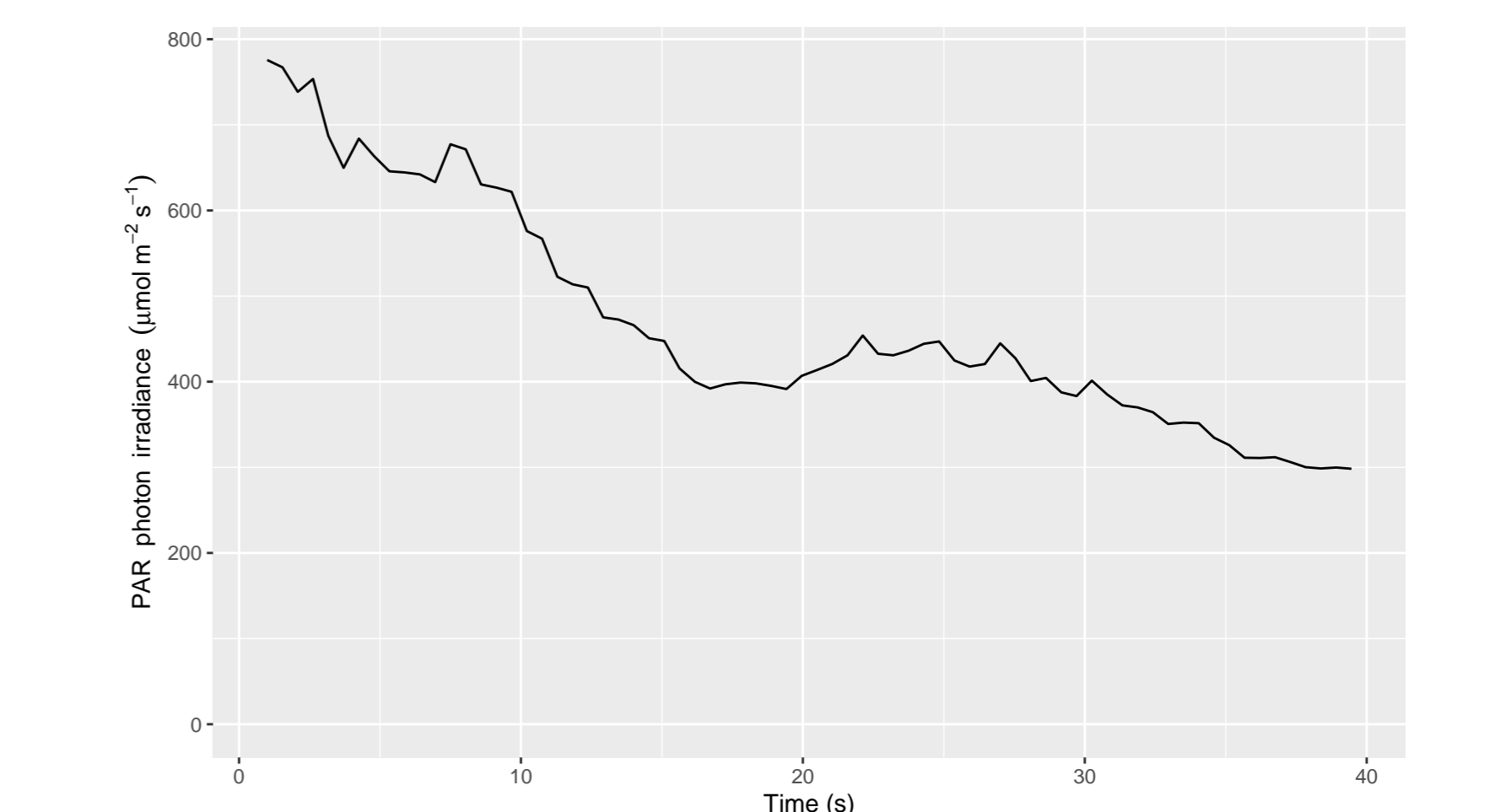
Collections of spectra, can be used to store groups of related spectra. Here we compare some Schott filters.

```
my.filters <- filters.mspect[intersect(schott_filters, uv_filters)]
summ.tb <- transmittance(my.filters, UV_bands())
# shorten column names
colnames(summ.tb) <- gsub("transmittance", "Tfr", colnames(summ.tb))
summ.tb

## # A tibble: 8 x 4
##   spct.idx   Tfr_UVC.ISO.tr.lo Tfr_UVB.ISO Tfr_UVA.ISO
##   <fct>      <dbl>      <dbl>      <dbl>
## 1 Schott_UG1      NA          0.153      0.659
## 2 Schott_UG5      0.540      0.957      0.924
## 3 Schott_UG11     0.150      0.851      0.694
## 4 Schott_N_WG280  0.0742     0.795     0.986
## 5 Schott_N_WG295  0.00266    0.529     0.978
## 6 Schott_N_WG305  0.0000290  0.305     0.961
## 7 Schott_N_WG320  0.0000100  0.0533    0.887
## 8 Schott_UG2A     NA          0.0472    0.660
```

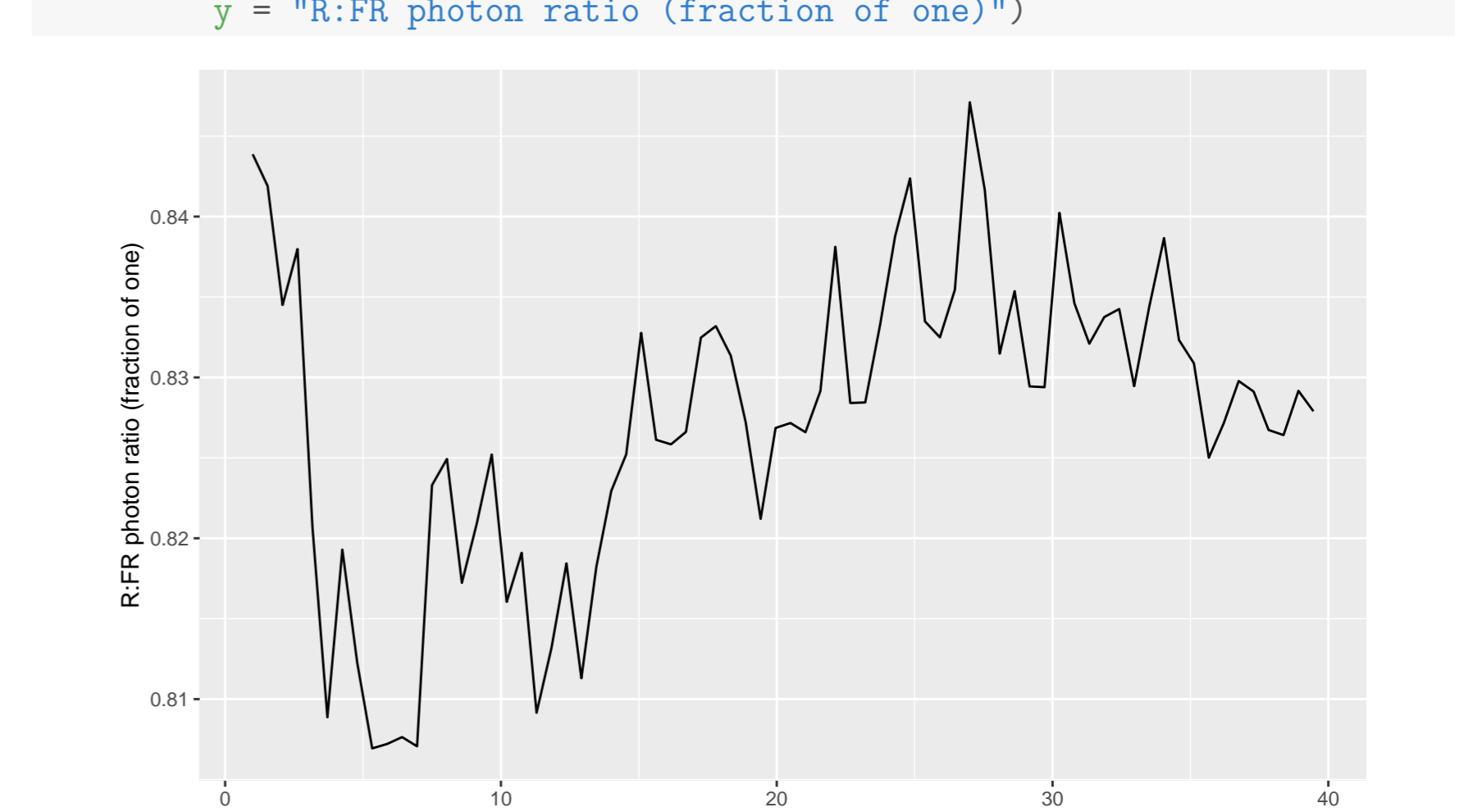
Collections of spectra can be also used to store time series of spectral measurements. We next use a time series of spectra measured with an array spectrometer in April in Helsinki, under silver birch trees. In this we use R package ggplot2 for plotting the summaries against time.

```
ppfd.df <- q_irrad(gap.mspect, PAR())
ppfd.df$time <- seq(from = 1, by = attr(gap.mspect, "time.step"),
  length.out = nrow(ppfd.df))
ggplot(ppfd.df, aes(time, q_irrad_PAR * 1e6)) +
  geom_line() + ylim(0, NA) +
  labs(x = "Time (s)",
  y = expression(PAR~photon~irradiance~(mu~mol~m^-2~s^-1)))
```



The time course of red:far-red photon ratio for the same time series of spectra.

```
R_FR.df <- q_ratio(gap.mspect, Red("Smith10"), Far_red("Smith10"))
names(R_FR.df)[2] <- "R_FR"
R_FR.df$time <- seq(from = 1, by = attr(gap.mspect, "time.step"),
  length.out = nrow(R_FR.df))
ggplot(R_FR.df, aes(time, R_FR)) +
  geom_line() +
  labs(x = "Time (s)",
  y = "R:FR photon ratio (fraction of one)")
```



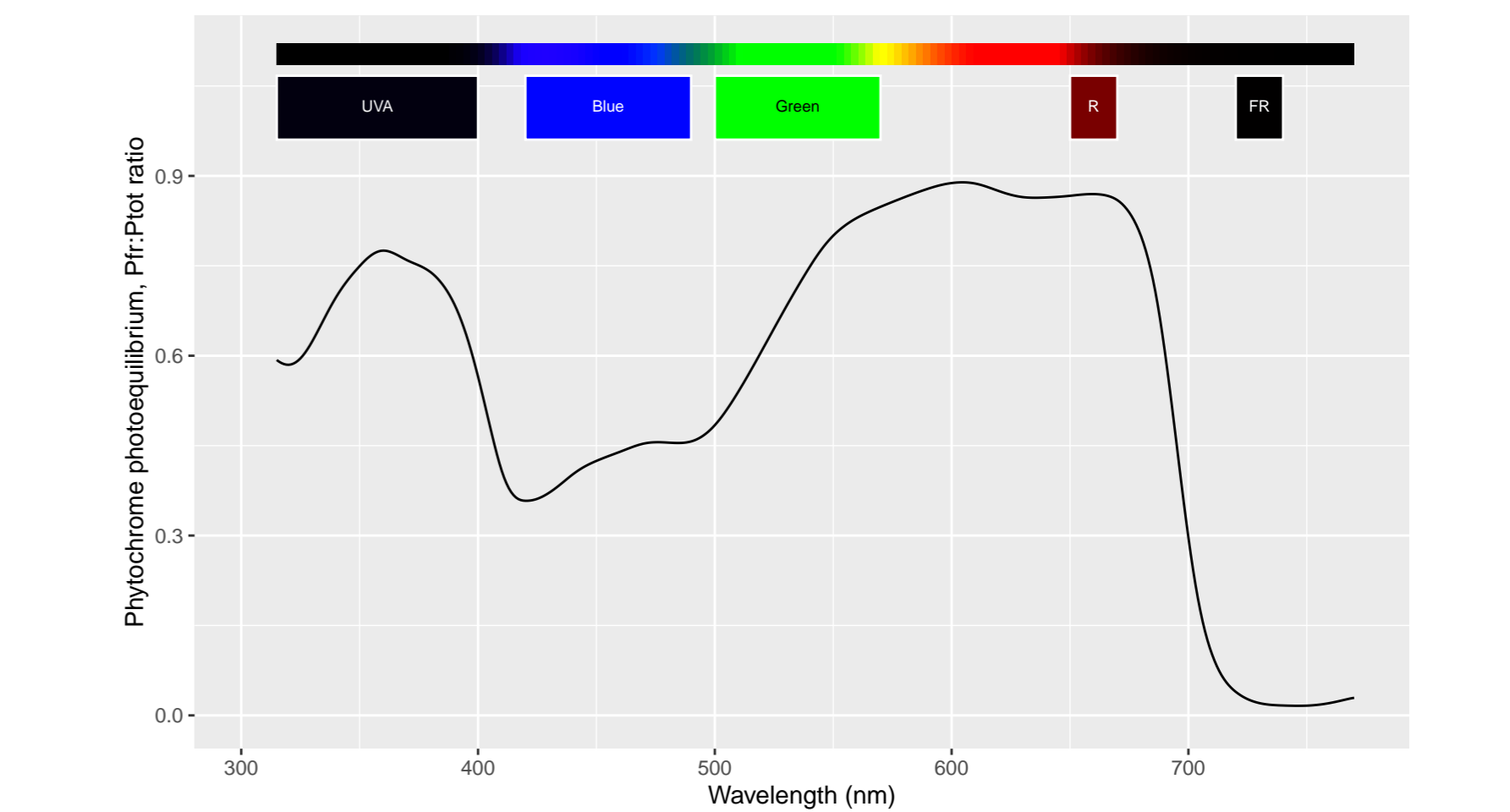
7. Photoreceptors

Phytochrome photoequilibrium from spectral data.

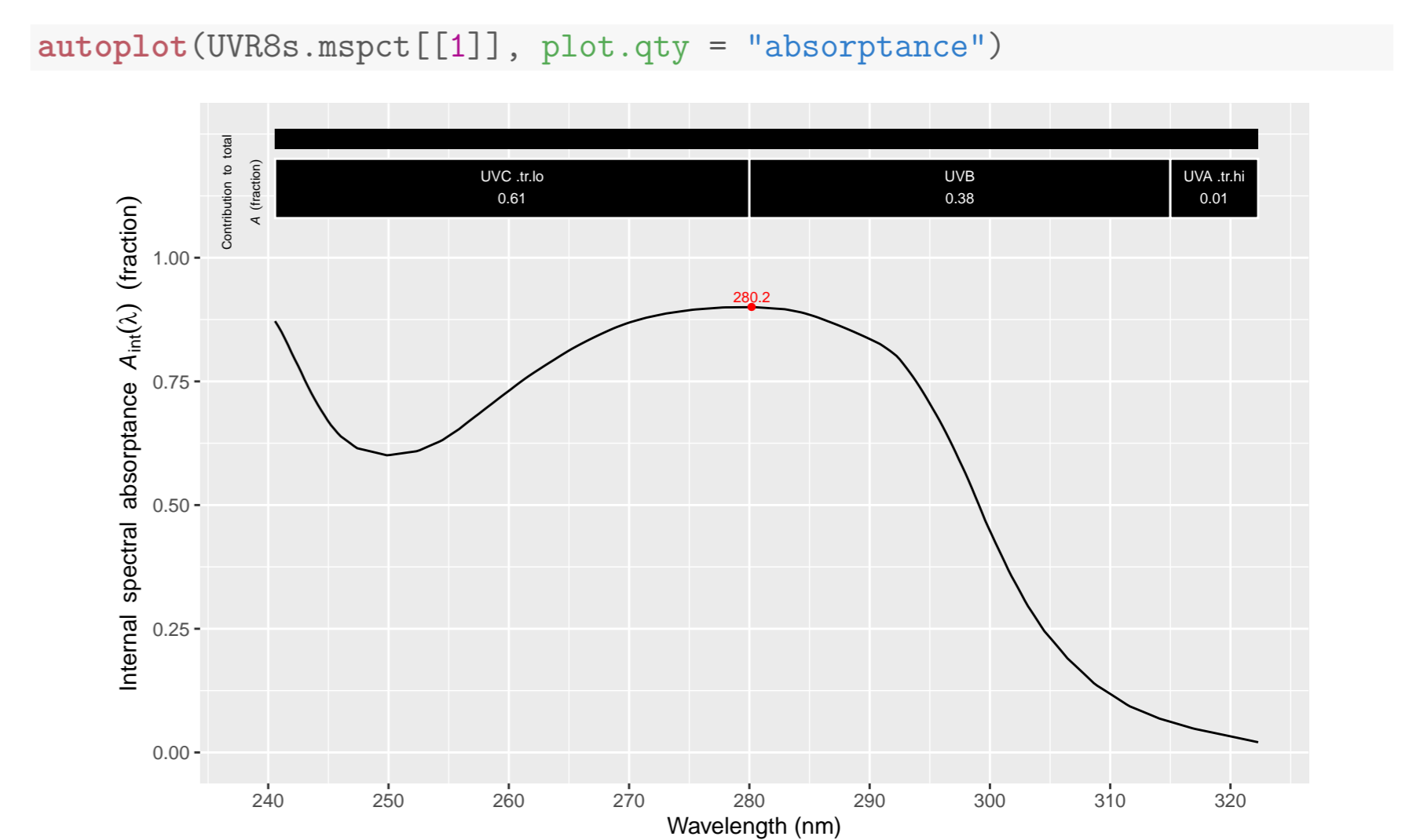
```
Pfr_Ptot(sun.spct)
## [1] 0.68341
```

Plot of Pfr:Ptot as a function of wavelength (nm) of monochromatic light.

```
autoplot(Pfr_Ptot(315:770), norm = NULL, unit.out = "photon",
  w.band = Plant_bands(),
  annotations = c("colour-guide", "labels", "boxes")) +
  labs(y = "Phytochrome photoequilibrium, Pfr:Ptot ratio")
```



The absorbance spectrum of uvr8.



<https://www.r4photobiology.info/>

More information is available at <https://www.r4photobiology.info/> where installation instructions, documentation and announcements of updates are published. Each package includes one or more user guides and help pages containing additional examples and explanations.