hyperSpec Plotting functions

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Vignette under Development

This file is currently undergoing a thorough revision. Changes may happen frequently. Even if the file is not yet nice to read, the shown code does work.

Reproducing the Examples in this Vignette

All spectra used in this manual are installed automatically with *hyperSpec*. Note that some definitions are executed in vignette.defs.

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For some plots of the chondro dataset, the pre-processed spectra are preferred, and their cluster averages \pm one standard deviation:

> chondro.preproc <- chondro - spc.fit.poly.below (chondro)

Fitting with npts.min = 15

- > chondro.preproc <- sweep (chondro.preproc, 1, mean, "/")
- > chondro.preproc <- sweep (chondro.preproc, 2, apply (chondro.preproc, 2, quantile, 0.05), "-")
 > cluster.cols <- c ("dark blue", "orange", "#C02020")
- > cluster.meansd <- aggregate (chondro.preproc, chondro\$clusters, mean_pm_sd)
- > cluster.means <- aggregate (chondro.preproc, chondro\$clusters, mean)

For details about the pre-processing, please refer to the vignette vignette ("chondro"), or the help ? chondro.

1 Predefined functions

hyperSpec comes with 5 major predefined plotting functions.

plot main switchyard for most plotting tasks

levelplot hyperSpec has a levelplot method

plotspc plots spectra

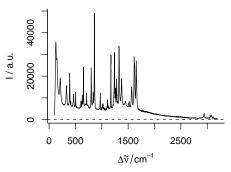
plotc calibration plot, time series, depth profile

plotc is a lattice function

plotmap more specialized version of levelplot for map or image plots.

plotmap is a *lattice* function

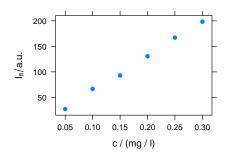
plotspc



plots the spectra, i.e. the intensities \$spc over the wavelengths @wavelength.

> plotspc (paracetamol)

plotc

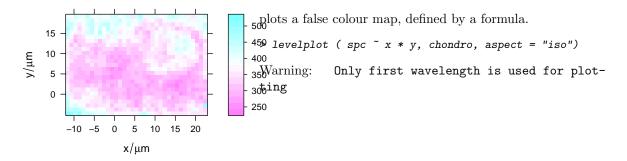


plots an intensity over a single other data column. e.g.

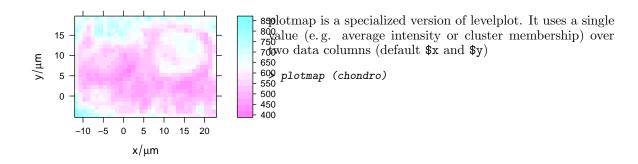
- time series
- calibration
- depth profile

> plotc (flu) 2

levelplot



plotmap

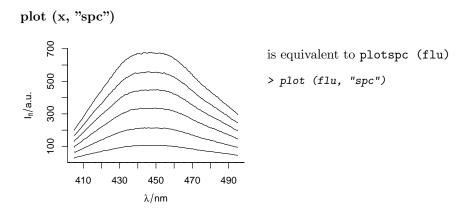


2 Arguments for plot

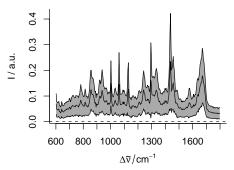
The three specialized functions are also accessible via plot:

hyperSpec's plot method uses the second argument to determine which of the three specialized plot functions to call. All further arguments are handed over to this function.

This allows a few more handy abbreviations.



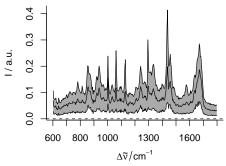
plot (x, "spcmeansd")



plots mean spectrum \pm 1 standard deviation

> plot (chondro.preproc, "spcmeansd")

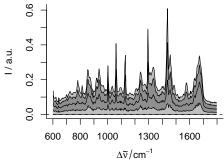
plot (x, "spcprctile")



plots median, $16^{\rm th}$ and $84^{\rm th}$ percentile for each wavelength. For Gaussian distributed data, $16^{\rm th}$, $50^{\rm th}$ and $84^{\rm th}$ percentile are equal to mean \pm standard deviation. Spectroscopic data frequently are not Gaussian distributed. The percentiles give a better idea of the true distribution. They are also less sensitive to outliers.

> plot (chondro.preproc, "spcprctile")

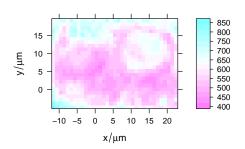
plot (x, "spcprctl5")



like "spcprctl" plus 5th and 95thpercentile.

> plot (chondro.preproc, "spcprct15")

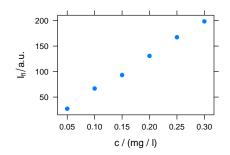
plot (x, "map")



is equivalent to plotmap (chondro)

> plot (chondro, "map")

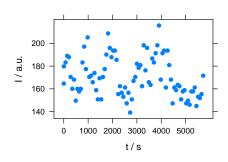
plot (x, "c")



is equivalent to plotc (flu)

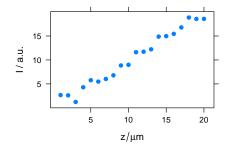
> plot (flu, "c")

plot (x, "ts")



> plot (laser, "ts")

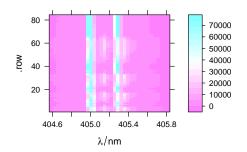
plot (x, "depth")



plots a depth profile plot, equivalent to plotc (laser, spc
~ z)

> plot (depth.profile, "depth")

plot (x, "mat")



plots the spectra matrix.

> plot (laser, "mat")

Equivalent to

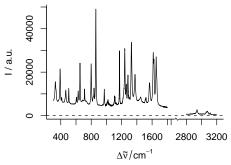
> levelplot (spc ~ .wavelength * .row, laser)

3 Spectra

plotspc

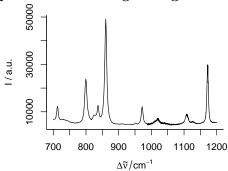
plotspc offers a variety of parameters for customized plots. To plot ...

particular wavelength ranges



use wl.range = list (600 ~ 1800, 2800 ~ 3100). If wl.range already contains indices: use wl.index = TRUE. Cut the wavelength axis appropriately with xoffset = 750

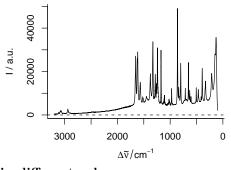
particular wavelength ranges II



if only one wavelength range is needed, the extract command is handier:

> plotspc (paracetamol[,, 700 ~ 1200])

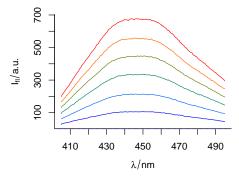
with reversed abscissa



use wl.reverse = TRUE

> plotspc (paracetamol, wl.reverse = TRUE)

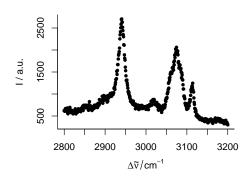
in different colours



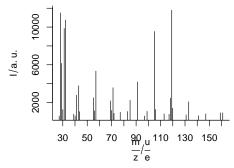
use col = vector.of.colours

> plotspc (flu, col = matlab.dark.palette(6))

dots instead of lines

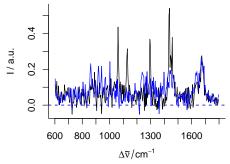


${\bf mass\ spectra}$



use lines.args = list (type = "h")
> plot (barbituates [[1]], lines.args = list (type = "h"))

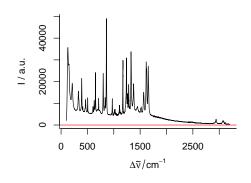
more spectra into an exsisting plot



use add = TRUE

> plotspc (chondro.preproc [30,,])
> plotspc (chondro.preproc [300,,], add = TRUE, col = "blue")

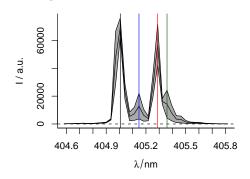
with different line at I=0



use zeroline = list.of.arguments.to.abline. NULL suppresses the line.

> plotspc (paracetamol,
+ zeroline = list (col = "red"))

adding lines, etc.

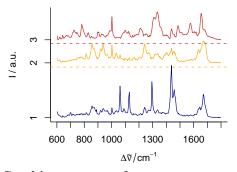


use abline for adding lines

```
> plot (laser, "spcmeansd")
> abline (v = wl (laser)[c (13, 17, 21, 23)],
+ col = c("black", "blue", "red", "darkgreen") )
```

3.1 Stacked spectra

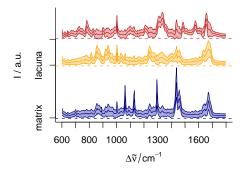
stacked = TRUE



use stacked = TRUE

```
> plotspc (cluster.means,
+ col = cluster.cols,
+ stacked = TRUE)
```

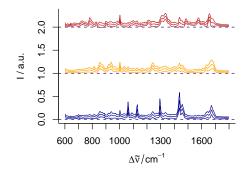
Stacking groups of spectra



The spectra to be stacked can be grouped: stacked = "grouping.column.name" The same applies to

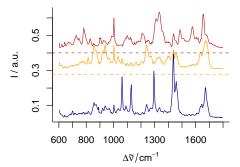
```
> plot (cluster.meansd,
+ stacked = ".aggregate",
+ fill = ".aggregate",
+ col = cluster.cols)
```

Manually giving yoffset



Stacking values can also be given manually as numeric values in yoffset:

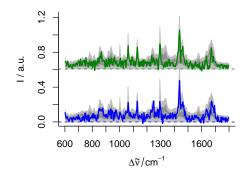
```
> plotspc (cluster.meansd,
+ yoffset = rep (0:2, each = 3),
+ col = rep (cluster.cols, each = 3))
```



To obtain a denser stacking:

```
> ## coordinate system:
> yoffsets <- apply (cluster.means [[]], 2, diff)
> yoffsets <- apply (yoffsets, 1, min)
> plot (cluster.means, yoffset = c (0, cumsum (yoffsets)), col = clust
```

Elaborate example



A more elaborate example of manual stacking:

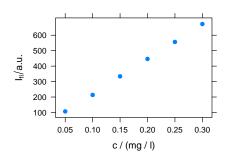
```
> ## coordinate system:
> yoffset <- apply (chondro.preproc, 2, quantile, c(0.05, 0.95))
 yoffset <- range (yoffset)</pre>
> plot(chondro.preproc[1],
       plot.args = list (ylim = c (0, 2) * yoffset),
       lines.args = list( type = "n")
> yoffset <- (0:1) * diff (yoffset)
> ## stacked spectra:
 for (i in 1 : 3){
   plot(chondro.preproc, "spcprct15",
         yoffset = yoffset [i],
         col = "gray", add = TRUE)
   plot (chondro.preproc [i],
          yoffset = yoffset [i],
          col = matlab.dark.palette (3) [i],
          add = TRUE,
          lines.args = list (lwd = 2))
```

4 Calibration Plots, (Depth) Profiles, and Time Series Plots

plotc

4.1 Calibration plots

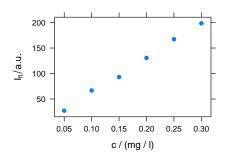
Intensities over concentration



Plotting the Intensities of one wavelength over the concentration for univariate calibration:

```
> plotc (flu [,, 450])
```

Integral Intensity over concentration

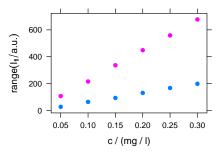


The default is to use the first intensity:

> plotc (flu)

Warning: In plotc(flu): Intensity at first wavelengh only is used.

Other Summary Intensities over concentration

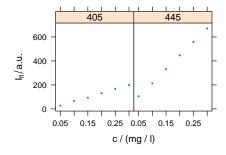


A function to compute a summary of the intensities before drawing can be used:

> plotc (flu, func = range, groups = .wavelength)

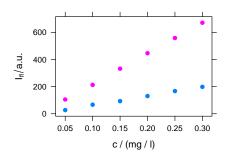
If func returns more than one value, the different results are accessible by .wavelength.

Conditioning: plotting more traces separately



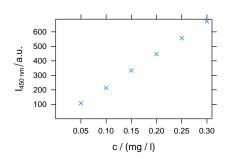
> plotc (flu [,, c (405, 445)], spc ~ c | .wavelength,
+ cex = .3, scales = list (alternating = c(1, 1)))

Grouping: plot more traces in one panel



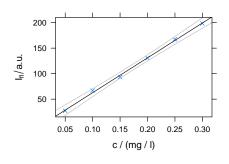
> plotc (flu [,, c (405, 445)], groups = .wavelength)

Changing Axis Labels (and other parameters)



Arguments for xyplot can be given to plotc:

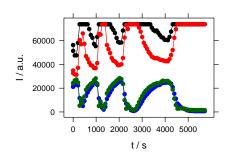
Adding things to the plot: customized panel function



As plotc uses the *lattice* function xyplot, additions to the plot must be made via the panel function:

4.2 Time series and other Plots of the Type Intensity-over-Something

Abscissae other than c



Other abscissae may be specified by explicitly giving the model formula:

5 Levelplot

Levelplot can use two special column names:

.wavelength for the wavelengths

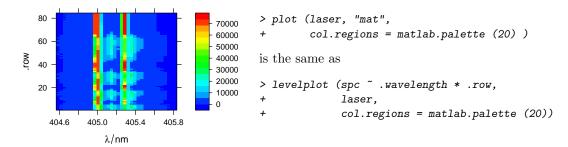
.row for the row index (i. e. spectrum number) in the data

6 Spectra Matrix

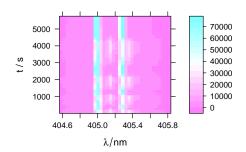
It is often useful to plot the spectra against an additional coordinate, e.g. the time for time series, the depth for depth profiles, etc.

This can be done by plot (object, "mat") or levelplot (model = spc ~.wavelength * other.data.column, object). The actual plotting is done by levelplot, so the plots can be grouped or conditioned.

different palette



different y axis



Changing the y axis is only possible with levelplot:

```
> levelplot (spc ~ .wavelength * t,
+ laser)
```

different panel

```
400000
                                  > class (barbituates)
                          350000
300000
                                  [1] "list"
                          250000
                                  > barbituates <- do.call (collapse, barbituates)
                          200000
                          150000
                                  > barbituates <- orderwl (barbituates)
                          100000
                          50000
                                  > levelplot (spc \tilde{} .wavelength * z, barbituates,
50
    100
        150
             200
                                                 panel = panel.levelplot.points,
      m .u
       z
                                                 cex = .5,
         e
                                                 col.regions = colorRampPalette (c ("orange", "darkred"))
```

Contourplot with a different y axis:

Contourplots are possible with plot and levelplot:

contour lines with plot

700 > plot (flu, 6 650 600 550 600 "mat", 5 500 contour = TRUE, 400 labels = TRUE, 300 col = "#00000080",200 2 at = seq(0, 700, by = 50))100 0 480 440 460 λ/nm

contour lines with levelplot

```
700
                   650
600<sub>550</sub>
                                                       levelplot (spc ~ .wavelength * c,
   0.30
                                             600
                                                                      flu,
   0.25
                                             500
c / (mg / I)
                                                                       contour = TRUE,
   0.20
                                             400
                                                                       labels = TRUE,
                                             300
   0.15
                                                                       col = "#00000080",
                                             200
   0.10
                                                                       at = seq(0, 700, by = 50))
          50
                                             100
   0.05
            420
                  440
                        460
                              480
                   \lambda/nm
```

7 False-Colour Maps

plotmap

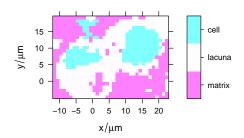
plotmap uses levelplot, a *lattice* function. Therefore, in loops, functions, Sweave chunks, etc. the lattice object needs to be printed explicitly by print (plotmap (object)) (R FAQ: Why do lattice/trellis graphics not work?).

plotmap produces a 3d plot, with the z axis colour-coded. plotmap's arguments x and y take the name of extra data columns.

The colour-coded axis. Also z can be used to select one column of the extra data by name. Alternatively, it may be a numeric or factor directly giving the values to be used. Each level of a factor will have one colour. It is also possible to plot a sum characteristic of the spectra: supply

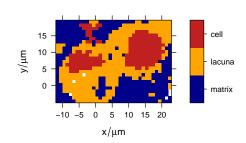
the function in argument func. The default setting is to plot the average intensity (no z and func=mean).

plotting clusters



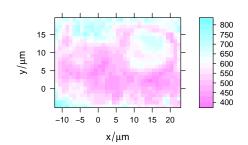
> plotmap (chondro, clusters ~ x * y)

different palette



To plot with a different palette, use *trellis.args=* list (col.regions = palette).

defined wavelengths



To plot a map with particular wavelengths use this:

```
> plotmap (chondro[, , c( 728, 782, 1098, 1240,
+ 1482, 1577)])
```

1.5 1.0 0.5 0.0.0

Preprocessing of the data:

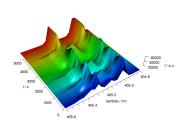
Examples

Conditioning. Lattice graphics have a concept of conditioning a plot. Instead of plotting all data in one diagram, a diagram is produced for each of the groups specified by the condition. plotmap's argument *cond* takes he name of the extra data column used for conditioning. This could e.g. be a column containing the sample number of a *hyperSpec* object that contains several samples.

Beispiele: voronoi

8 3 D

3D figures are possible with rgl



> library (rgl)

9 Troubleshooting

9.1 No output is produced

plotmap and plotc use levelplot, a *lattice* function. Therefore, in loops, functions, Sweave chunks, etc. the lattice object needs to be printed explicitly by print (plotmap (object)) (R FAQ: Why do lattice/trellis graphics not work?).

For suggestions how the lattice functions can be redefined so that the result is printed without external print command, see vignettes.defs.

10 Interactive Graphics

hyperSpec offers two basic interaction functions, spc.identify, and map.identify. They identify points in spectra plots and map plots, respectively.

10.1 spc.identify: finding out wavelength, intensity and spectrum

spc.identify allows to measure points in graphics produced by plotspc. It works correctly with reversed and cut wavelength axes.

```
> spc.identify (plotspc (paracetamol, wl.range = c (600 ~ 1800, 2800 ~ 3200), xoffset = 800))
```

The result is a data.frame with the indices of the spectra, the wavelength, and its intensity.

10.2 map.identify: finding a spectrum in a map plot

map.identify returns the spectra indices of the clicked points.

```
> map.identify (chondro)
```

10.3 Related functions provided by base graphics and lattice

For base graphics (as produced by plotspc), locator may be useful as well. It returns the clicked coordinates. Note that these are *not* transformed according to xoffset & Co.

For lattice graphics, grid.locator may be used instead. If it is not called in the panel function, a preceeding call to trellis.focus is needed:

```
> plot (laser, "mat")
> trellis.focus ()
> grid.locator ()
```

identify (or panel.identify for lattice graphics) allows to identify points of the plot directly. Note that the returned indices correspond to the plotted object.

10.4 Interactively changing graphics

hyperSpec's lattice functions work with playwith and latticist. These packages allow easy customization of the plots and also identification of points.