hyperSpec Plotting functions

Claudia Beleites < Claudia. Beleites @chemometrix.gmbh>

DIA Raman Spectroscopy Group, University of Trieste/Italy (2005 – 2008) Spectroscopy · Imaging, IPHT, Jena/Germany (2008 – 2017) ÖPV, JKI, Berlin/Germany (2017 – 2019)

Arbeitskreis Lebensmittelmikrobiologie und Biotechnologie, Hamburg University, Hamburg/German Chemometric Consulting and Chemometrix GmbH, Wölfersheim/Germany (since 201

February 13, 2020

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, and others invisibly at the beginning of the file in order to have the code as similar as possible to interactive sessions.

Contents

1	Predefined functions	2
2	Arguments for plot	5
3	Spectra 3.1 Stacked spectra	8 10
4	Calibration Plots, (Depth) Profiles, and Time Series Plots4.1Calibration plots4.2Time series and other Plots of the Type Intensity-over-Something	12 12 13
5	Levelplot	14
6	Spectra Matrix	14
7	False-Colour Maps: plotmap	16
8	3D plots (with rgl)	20
9	Using ggplot2 with hyperSpec objects	21
10	Troubleshooting 10.1 No output is produced	22 22

11	Interactive Graphics	22
	11.1 spc.identify: finding out wavelength, intensity and spectrum	23
	11.2 map.identify: finding a spectrum in a map plot	23
	11.3 map.sel.poly: selecting spectra inside a polygon in a map plot	23
	11.4 Related functions provided by base graphics and lattice	23

Suggested Packages

latticeExtra: available deldir: available rgl: available ggplot2: available

In addition tripack, and latticist are mentioned, but not used in this vignette.

Preliminary Calculations

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

```
> chondro.preproc <- chondro - spc.fit.poly.below (chondro)
> chondro.preproc <- chondro.preproc / rowMeans (chondro)
> chondro.preproc <- chondro.preproc - quantile (chondro.preproc, 0.05)
> cluster.cols <- c ("dark blue", "orange", "#CO2020")
> cluster.meansd <- aggregate (chondro.preproc, chondro$clusters, mean_pm_sd)
> cluster.means <- aggregate (chondro.preproc, chondro$clusters, mean)</pre>
```

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

1 Predefined functions

hyperSpec comes with 6 major predefined plotting functions.

plotspc plots spectra

plotmat plots the spectra matrix

plotc calibration plot, time series, depth profile

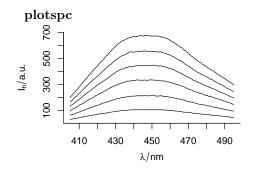
 ${ t plots}$ is a lattice function

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

plotmap is a lattice function

plotvoronoi more specialized version of plotmap that produces Voronoi tesselations. plotvoronoi is a *lattice* function

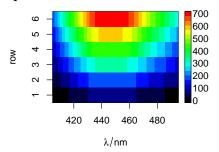
plotmap, plotvoronoi, and levelplot are *lattice* functions. Therefore, in loops, functions, Sweave chunks, etc. the lattice object needs to be printed explicitly by e.g. print (plotmap (object)) (R FAQ: Why do lattice/trellis graphics not work?).



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

> plotspc (flu)

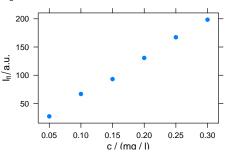
plotmat



plots the spectra, i.e. the colour coded intensities \$spc over the wavelengths @wavelength and the row number.

> plotmat (flu)

plotc

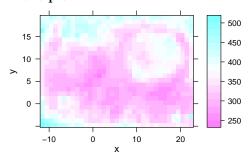


plots an intensity over a single other data column, e.g.

- calibration
- time series
- depth profile

> plotc (flu)

levelplot

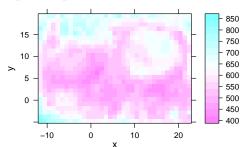


plots a false colour map, defined by a formula.

> levelplot (spc ~ x * y, chondro, aspect = "iso")

Warning: Only first wavelength is used for plotting

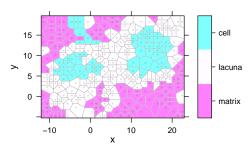
plotmap



 $\label{eq:plotmap} \begin{array}{l} \texttt{plotmap} \ \text{is a specialized version of levelplot}. \ \ \texttt{It} \ \text{uses a} \\ \texttt{single value} \ (\texttt{e.\,g.} \ \text{average intensity or cluster membership}) \\ \texttt{over two data columns} \ (\texttt{default \$x} \ \texttt{and \$y}) \end{array}$

> plotmap (chondro)

plotvoronoi

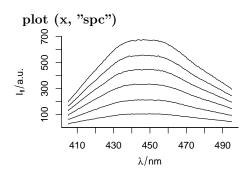


Function plotvoronoi is a special version of plotmap that produces Voronoi diagram of the hyperSpec object.

> plotvoronoi (sample (chondro, 300), clusters ~ x * y)

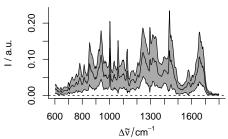
2 Arguments for plot

hyperSpec's plot method uses its second argument to determine which of the specialized plots to produce. This allows some handy abbreviations. All further arguments are handed over to the function actually producing the plot.



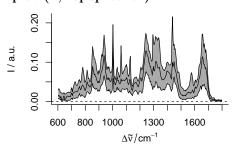
is equivalent to plotspc (flu)
> plot (flu, "spc")

plot (x, "spcmeansd")



plots mean spectrum ± 1 standard deviation
> plot (chondro.preproc, "spcmeansd")

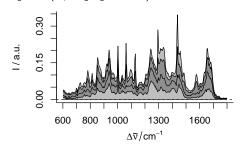
plot (x, "spcprctile")



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 "spcprct1" plus $5^{\rm th}$ and $95^{\rm th}$ percentile. > plot (chondro.preproc, "spcprct15")

plot (x, "c")

200

150

50

0.05

0.10

0.15

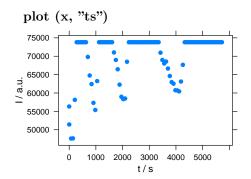
0.20

0.25

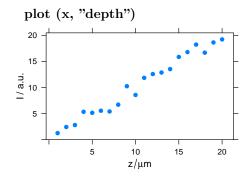
0.30

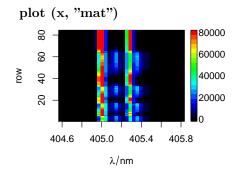
c / (ma / I)

> plot (flu, "c")
is equivalent to plotc (flu)

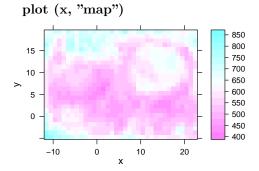


plots a time series plot
> plot (laser [,, 405], "ts")
equivalent to plotc (laser, spc ~ t)





plots the spectra matrix.
> plot (laser, "mat")
Equivalent to
> plotmat (laser)
A lattice alternative is:
> levelplot (spc ~ .wavelength * .row, laser)



is equivalent to plotmap (chondro)
> plot (chondro, "map")

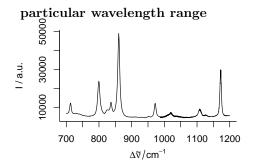
plot (x, "voronoi") -10

> plot (sample (chondro, 300), "voronoi") See plotvoronoi

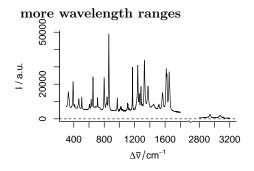
3 Spectra

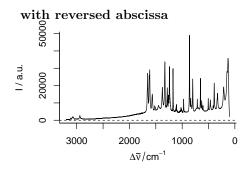
plotspc

 ${\tt plotspc}$ offers a variety of parameters for customized plots. To plot \dots



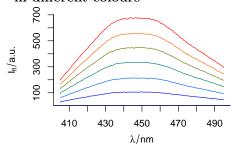
if only one wavelength range is needed, the extract command (see vignette ("introduction")) is handiest:
> plotspc (paracetamol [,, 700 ~ 1200])
wavelengths. If wl.range already contains indices use wl.index = TRUE.





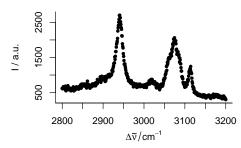
use wl.reverse = TRUE
> plotspc (paracetamol, wl.reverse = TRUE)

in different colours

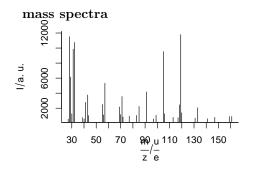


use col = vector.of.colours
> plotspc (flu, col = matlab.dark.palette (6))

dots instead of lines

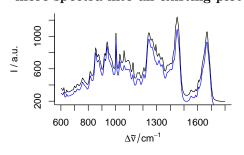


use lines.args = list (pch = 20, type = "p")
> plotspc (paracetamol [,, 2800 ~ 3200],
+ lines.args = list (pch = 20, type = "p"))



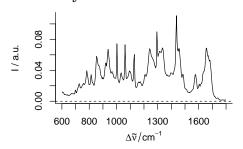
use lines.args = list (type = "h")
> plot (barbiturates [[1]], lines.args = list (type = "h"))

more spectra into an existing plot



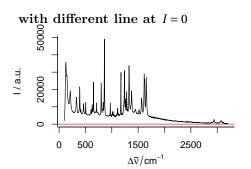
use add = TRUE
> plotspc (chondro [30,,])
> plotspc (chondro [300,,], add = TRUE, col = "blue")

Summary characteristics



 func may be used to calculate summary characteristics prior to plotting. To plot e.g. the standard deviation of the spectra, use:

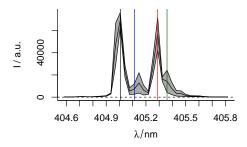
> plotspc (chondro.preproc, func = sd)



zero line takes a list with parameters to ${\tt abline}, \, {\tt NA} \, \, {\tt suppresses}$ the line.

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

adding to a spectra plot

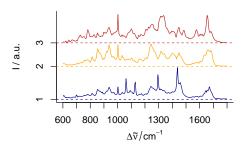


plotspc uses base graphics. After plotting the spectra, more content may be added to the graphic by abline, lines, points, etc.

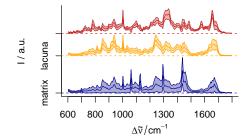
```
> plot (laser, "spcmeansd")
> abline (v = c(405.0063, 405.1121, 405.2885, 405.3591),
+ col = c("black", "blue", "red", "darkgreen"))
```

3.1 Stacked spectra

stacked



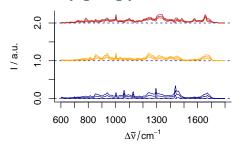
Stacking groups of spectra



The spectra to be stacked can be grouped: stacked = factor. Alternatively, the name of the grouping extra data column can be used:

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

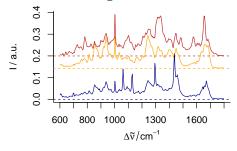
Manually giving yoffset



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

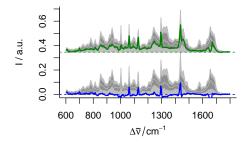
```
> plotspc (cluster.meansd,
+ yoffset = rep (0:2, each = 3),
+ col = rep (cluster.cols, each = 3))
```

Dense stacking



```
To obtain a denser stacking:
> yoffsets <- apply (cluster.means [[]], 2, diff)
> yoffsets <- - apply (yoffsets, 1, min)
> plot (cluster.means, yoffset = c (0, cumsum (yoffsets)),
+ col = cluster.cols)
```

Elaborate example



plotspc allows fine grained customization of almost all aspects of the plot. This is possible by

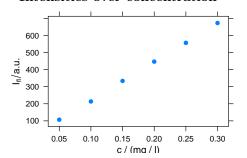
giving arguments to the functions that actually perform the plotting plot for setting up the plot area, lines for the plotting of the lines, axis for the axes, etc. The arguments for these functions should be given in lists as plot.args, lines.args, axis.args, etc.

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

plotc

4.1 Calibration plots

Intensities over concentration

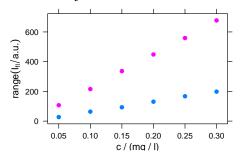


Plotting the Intensities of one wavelength over the concentration for univariate calibration:

> plotc (flu [,, 450])

The default is to use the first intensity only.

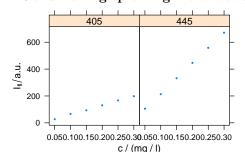
Summary Intensities over concentration



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

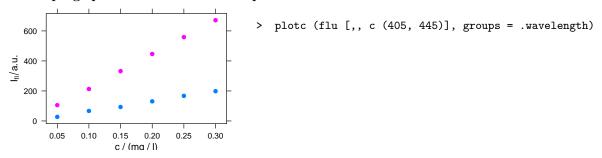
> plotc (flu, func = range, groups = .wavelength)
If func returns more than one value, the different results are accessible by .wavelength.

Conditioning: plotting more traces separately

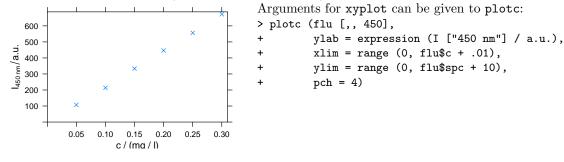


> plotc (flu [,, c (405, 445)], spc ~ c | .wavelength,
+ cex = .3, scales = list (alternating = c(1, 1)))

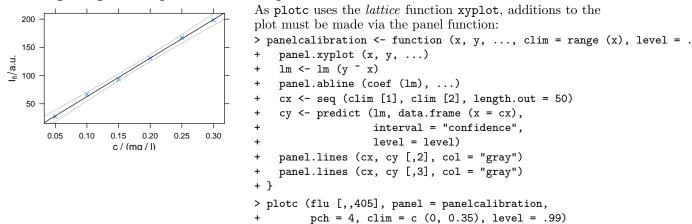
Grouping: plot more traces in one panel



Changing Axis Labels (and other parameters)

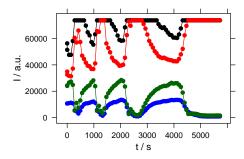


Adding things to the plot: customized panel function



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

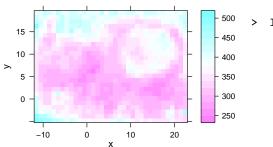
hyperSpec's levelplot can use two special column names:

.wavelength for the wavelengths

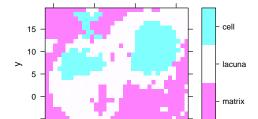
.row for the row index (i. e. spectrum number) in the data

Besides that, it behaves exactly like level plot. Particularly, the data is given as the second argument:

levelplot



> levelplot (spc ~ x * y, chondro)



10

20

If the colour-coded value is a factor, the display is adjusted to this fact:

> levelplot (clusters ~ x * y, chondro)

6 Spectra Matrix

-10

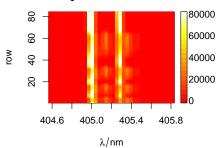
factors as z

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"). The actual plotting is done by image, but levelplot

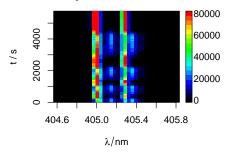
can produce spectra matrix plots as well and these plots can be grouped or conditioned.

different palette



```
> plot (laser, "mat", col = heat.colors (20))
is the same as
> plotmat (laser, col = heat.colors (20))
```

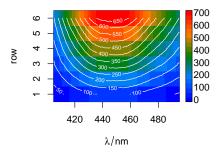
different y axis



Using a different extra data column for the y axis: > plotmat (laser, y = "t") alternatively, y values and axis label can be given separately.

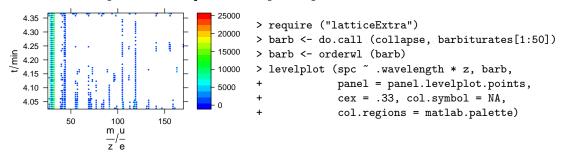
> plotmat (laser, y = laser\$t, ylab = labels (laser, "t"))

contour lines



```
Contour lines may be added:
> plotmat (flu, col = matlab.dark.palette (20))
> plotmat (flu, col = "white",
+ contour = TRUE, add = TRUE)
```

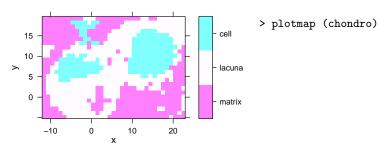
colour-coded points: levelplot with special panel function



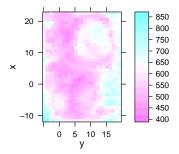
7 False-Colour Maps: plotmap

plotmap is a specialized version of levelplot. The spectral intensities may be summarized by a function before plotting (default: mean). The same scale is used for x and y axes (aspect = "iso").

plotting map

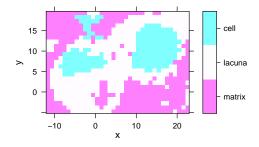


plotting maps with other than x and y



specify the colour-coded variable, abscissa and ordinate as
formula: colour.coded ~ abscissa * ordinate
> plotmap (chondro, spc ~ y * x)

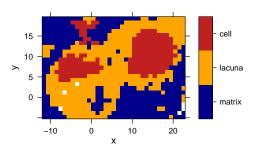
colour-coded factors



> plotmap (chondro, clusters ~ x * y)

If the colour-coded variable is a factor, each level gets its own colour, and the legend is labeled accordingly.

different palette



To plot with a different palette, use argument col.regions.

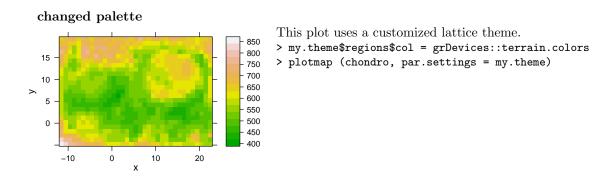
> plotmap (chondro, clusters ~ x * y,
+ col.regions = cluster.cols)

Fine tuning lattice parameters

The plotting of color maps is done via R package lattice (aka Trellis graphic approach), which is highly customizable. Use trellis.par.get and trellis.par.set to get/set the settings for the current graphics device.

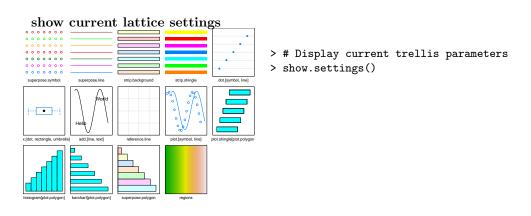
```
> my.theme = trellis.par.get()
> names(my.theme)
                           # note how many parameters are tunable
 [1] "grid.pars"
                          "fontsize"
                                               "background"
                                                                    "panel.background"
 [5] "clip"
                          "add.line"
                                                                    "plot.polygon"
                                               "add.text"
                          "box.rectangle"
 [9] "box.dot"
                                               "box.umbrella"
                                                                   "dot.line"
[13] "dot.symbol"
                          "plot.line"
                                               "plot.symbol"
                                                                    "reference.line"
[17] "strip.background"
                          "strip.shingle"
                                                                    "superpose.line"
                                               "strip.border"
[21] "superpose.symbol"
                                              "regions"
                                                                    "shade.colors"
                          "superpose.polygon"
                                                                    "layout.heights"
[25] "axis.line"
                          "axis.text'
                                               "axis.components
[29] "layout.widths"
                          "box.3d"
                                               "par.xlab.text"
                                                                    "par.ylab.text"
[33] "par.zlab.text"
                          "par.main.text"
                                               "par.sub.text"
```

Any of these parameters can be fine-tuned to produce the desired output. For example, parameter my.theme\$region is responsible for the appearance of color maps, and it contains elements \$alpha and \$col. By changing this parameters you can create your own theme for plotting and pass it to the plotting function via par.settings.



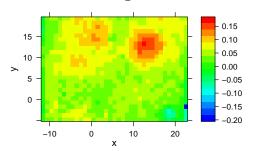
It is possible to **persistently** (i.e. inside of the current R session) set lattice parameters, so they would apply to all further plots. This is done via a call to trellis.par.set, for example trellis.par.set(my.theme).

The current settings can be visualized via a call to show.settings()



An overview of different color palettes, and ways to create your own, can be found in the **R** color cheatsheet.

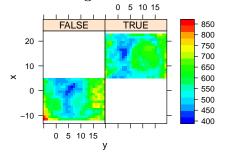
defined wavelengths



To plot a map of the average intensity at particular wavelengths use extraction:

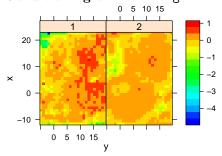
```
> plotmap (chondro.preproc [, , c(728, 782, 1098,
+ 1240, 1482, 1577)],
+ col.regions = matlab.palette)
```

Conditioning



```
> plotmap (chondro,
+            spc ~ y * x | x > 5,
+            col.regions = matlab.palette(20))
```

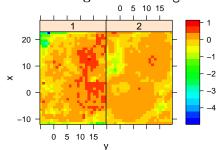
Conditioning on .wavelength



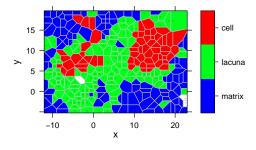
plotmap automatically applies the function in *func* before plotting. This defaults to the mean. In order to suppress this, use func = NULL. This allows conditioning on the wavelengths.

To plot e.g. the first two score maps of a principal component analysis:

Conditioning on .wavelength II



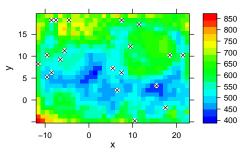
Voronoi plot



```
> plotvoronoi (sample (chondro, 300), clusters ~ x * y,
+ col.regions = matlab.palette(20))
```

Voronoi uses panel.voronoi from latticeExtra[2]. The tesselation is calculated by default using deldir[3], but tripack[4] can also be used. tripack seems to faster in general, but may "hang" with certain data sets (particularly regular grids with missing spectra as in this example). Furthermore, it is not FOSS (free and open source software), so users are kindly asked to review tripack's license before using it.

Mark missing spectra

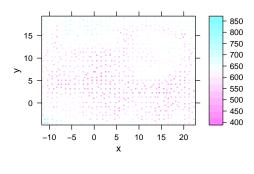


If the spectra come from a rectangular grid, missing positions can be marked with this panel function:

```
> mark.missing <- function (x, y, z, ...){
+ panel.levelplot (x, y, z, ...)
+
+ miss <- expand.grid (x = unique (x), y = unique (y))
+ miss <- merge (miss, data.frame (x, y, TRUE),
+ all.x = TRUE)
+ miss <- miss [is.na (miss[, 3]),]
+ panel.xyplot (miss [, 1], miss [, 2], pch = 4, ...)
+ }
> plotmap (sample (chondro, length(chondro) - 20),
+ col.regions = matlab.palette(20),
+ col = "black",
+ panel = mark.missing)
```

Unevenly spaced measurement grid

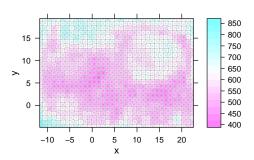
The panel function used by plotmap defaults to panel.levelplot.raster which assumes an evenly spaced measurement grid. Even if the spectra are measured on a nominally evenly spaced grid, the actual stage position may be slightly varying due to positioning inaccuracy and some manufacturers (e.g. Kaiser) record the position reported by the stage rather than the position requested by the stage control.



This leads to weird looking output with holes, and possibly wrong columns:

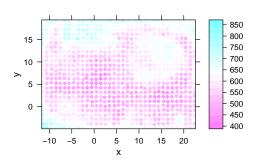
> plotmap (uneven)

The symptom of this situation are warnings about values in x and/or y not being equispaced; and that the output therefore may be wrong.



One possibility to obtain a correct map is using plotvoronoi instead which will construct a mosaic-like image with the respective "pixel" areas being centered around the actually recorded \$x\$ and \$y\$ position:

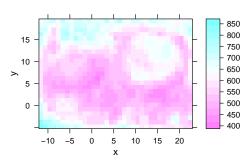
> plotvoronoi (uneven)



Another possibility that underlines a point shape of the measurements is switching to latticeExtra::panel.levelplot.points:

> plotmap (uneven, panel = panel.levelplot.points,

cex = 0.75, col.symbol=NA)



Alternatively, the measurement raster positions can be rounded to their nominal raster, e.g.:

- > rx <- makeraster (uneven\$x, start = -11.55, d = 1, tol = 0.3)
- > uneven\$x <- rx\$x
- > ry <- makeraster (uneven\$y, start = -4.77, d = 1, tol = 0.3)
- > uneven\$y <- ry\$x
- > plotmap (uneven)

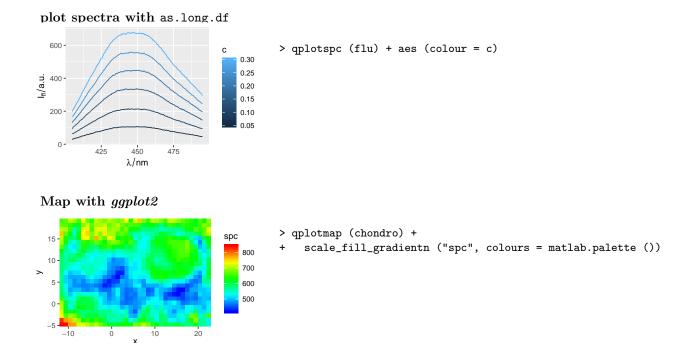
8 3D plots (with rgl)

```
rgl[5] offers fast 3d plotting in R. As rgl's axis annotations
are sometimes awkward, they may better be set manually:
> laser <- laser [,,404.8 ~ 405.6] / 10000
> laser$t <- laser$t / 3600
> cols <- rep (matlab.palette (nrow (laser)), nwl (laser))
> surface3d (y = wl (laser), x = laser$t,
+ z = laser$spc, col = cols)
> aspect3d (c (1, 1, 0.25))
> axes3d (c ('x+-', 'y--', 'z--'))
> axes3d ('y--', nticks = 25, labels= FALSE)
> mtext3d ("t / h", 'x+-', line = 2.5)
> mtext3d ("lambda / nm", 'y--', line = 2.5)
> mtext3d ("I / a.u.", edge = 'z--', line = 2.5)
```

Function not available package rgl needed.

9 Using ggplot2 with hyperSpec objects

hyperSpec objects do not yet directly support plotting with ggplot2 [6]. Nevertheless, ggplot2 graphics can easily be obtained, and qplot* equivalents to plotspc and plotmap are defined:

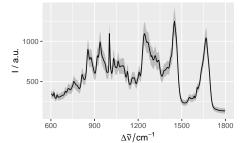


The two special columns .wavelength and .rownames contain the wavelength axis and allow to distinguish the spectra.

For more general plotting, as.long.df transforms a hyperSpec object into a long-form data.frame that is suitable for qplot, while as.t.df produces a data.frame where each spectrum is one column, and an additional first column gives the wavelength (see "plotting mean \pm sd" below for an example).

Long data.frames can be very memory consuming as they are of size $nrow \cdot nwl \times (ncol + 2)$ with respect to the dimensions of the hyperSpec object. Thus, e.g. the chondro data set (2 MB) as hyperSpec object) needs 28 MB as long-format data.frame. It is therefore highly recommended to calculate the particular data to be plotted beforehand.

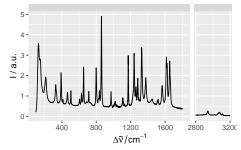
Mean \pm standard deviation with ggplot2



```
> qplotspc (mean (chondro)) +
+ geom_ribbon (aes (ymin = mean + sd,
+ ymax = mean - sd,
+ y = 0, group = NA),
+ alpha = 0.25,
+ data = as.t.df (mean_sd (chondro)))
```

Note that qpotspc specifies aesthetics y = spc and groups = .rownames, which do not have corresponding columns in the data.frame returned by as.t.df. These aesthetics must therefore be set manually in the aesthetics definition in geom_ribbon (or any other geom_ that uses as.t.df). Otherwise, errors occur that object spc (and/or .rownames) cannot be found.

Cut spectra with ggplot2



10 Troubleshooting

10.1 No output is produced

plotmap, plotvoronoi, levelplot, and plotc use *lattice* functions. 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?). The same holds for *ggplot2* graphics.

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

11 Interactive Graphics

hyperSpec offers basic interaction, spc.identify for spectra plots, and map.identify and map.sel.poly for maps. The first two identify points in spectra plots and map plots, respectively. map.sel.poly selects the part of a hyperSpec object that lies inside the user defined polygon.

11.1 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.

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

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

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

11.3 map.sel.poly: selecting spectra inside a polygon in a map plot

```
map.sel.poly returns a logical indicating which spectra are inside the polygon drawn by the user:
> map.sel.poly (chondro)
```

11.4 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 preceding 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.

References

- [1] Lemon J. Plotrix: a package in the red light district of r. R-News, 6(4):8–12, 2006.
- [2] Deepayan Sarkar and Felix Andrews. latticeExtra: Extra Graphical Utilities Based on Lattice, 2016. URL https://CRAN.R-project.org/package=latticeExtra. R package version 0.6-28.
- [3] Rolf Turner. deldir: Delaunay Triangulation and Dirichlet (Voronoi) Tessellation, 2019. URL https://CRAN.R-project.org/package=deldir. R package version 0.1-23.
- [4] Fortran code by R. J. Renka. R functions by Albrecht Gebhardt. With contributions from Stephen Eglen <stephen@anc.ed.ac.uk>, Sergei Zuyev, and Denis White. tripack: Triangulation of Irregularly Spaced Data, 2016. URL https://CRAN.R-project.org/package=tripack. R package version 1.3-8.
- [5] Daniel Adler, Duncan Murdoch, et al. rgl: 3D Visualization Using OpenGL, 2019. URL https://CRAN.R-project.org/package=rgl. R package version 0.100.30.
- [6] Hadley Wickham. ggplot2: Elegant Graphics for Data Analysis. Springer-Verlag New York, 2016. ISBN 978-3-319-24277-4. URL https://ggplot2.tidyverse.org.

Session Info

[49] tibble_2.1.3

[53] R6_2.4.1

[,1] "Linux" sysname "4.15.0-76-generic" release version "#86-Ubuntu SMP Fri Jan 17 17:24:28 UTC 2020" "cx17007" nodename "x86_64" machine "unknown" login "cb" user effective_user "cb" R version 3.6.2 (2019-12-12) Platform: x86_64-pc-linux-gnu (64-bit) Running under: Ubuntu 18.04.4 LTS Matrix products: default BLAS: /usr/lib/x86_64-linux-gnu/openblas/libblas.so.3 LAPACK: /usr/lib/x86_64-linux-gnu/libopenblasp-r0.2.20.so locale: [1] LC_CTYPE=de_DE.UTF-8 LC_NUMERIC=C LC_TIME=de_DE.UTF-8 [4] LC_COLLATE=C LC_MONETARY=de_DE.UTF-8 LC_MESSAGES=de_DE.UTF-8 [7] LC_PAPER=de_DE.UTF-8 LC_ADDRESS=C LC_NAME=C [10] LC_TELEPHONE=C LC_MEASUREMENT=de_DE.UTF-8 LC_IDENTIFICATION=C attached base packages: [1] tools grid stats graphics grDevices utils datasets methods base other attached packages: [1] latticeExtra_0.6-28 RColorBrewer_1.1-2 baseline_1.2-2 MASS_7.3-51.5 [5] hyperSpec_0.99-20200213 xml2_1.2.2 ggplot2_3.2.1 lattice_0.20-38 loaded via a namespace (and not attached): [1] tidyselect_0.2.5 xfun_0.11 reshape2_1.4.3 purrr_0.3.3 [5] colorspace_1.4-1 ${\tt testthat_2.3.1}$ miniUI_0.1.1.1 htmltools_0.4.0 [9] rlang_0.4.2 R.oo_1.23.0 manipulateWidget_0.10.0 pillar_1.4.2 [13] later_1.0.0 glue_1.3.1 withr_2.1.2 R.utils_2.9.2 [17] plyr_1.8.5 R.cache_0.14.0 lifecycle_0.1.0 stringr_1.4.0 [21] munsell_0.5.0 gtable_0.3.0 R.methodsS3_1.7.1 htmlwidgets_1.5.1 [25] mvtnorm_1.0-11 labeling_0.3 knitr_1.26 fastmap_1.0.1 [29] SparseM_1.78 httpuv_1.5.2 crosstalk_1.0.0 Rcpp_1.0.3 [33] xtable_1.8-4 scales_1.1.0 promises_1.1.0 plotrix_3.7-7 [37] webshot_0.5.2 jsonlite_1.6 deldir_0.1-23 ${\tt mime_0.7}$ [41] farver_2.0.1 R.rsp_0.43.2 digest_0.6.23 stringi_1.4.3 [45] dplyr_0.8.3 magrittr_1.5 lazyeval_0.2.2 shiny_1.4.0

crayon_1.3.4

compiler_3.6.2

pkgconfig_2.0.3

assertthat_0.2.1