

# 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

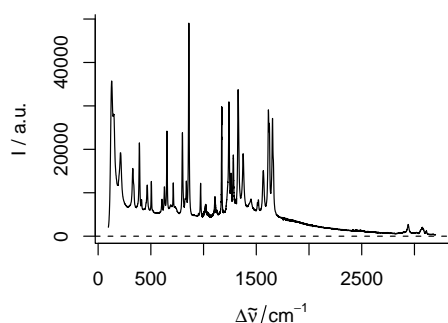
`levelplot` *hyperSpec* has a `levelplot` method

`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

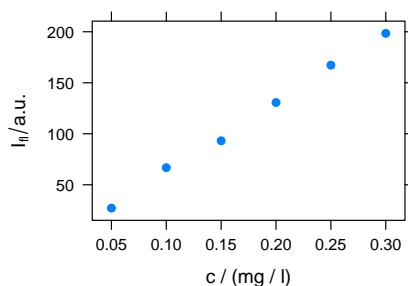
### plotspc



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

```
> plotspc (paracetamol)
```

### plotc

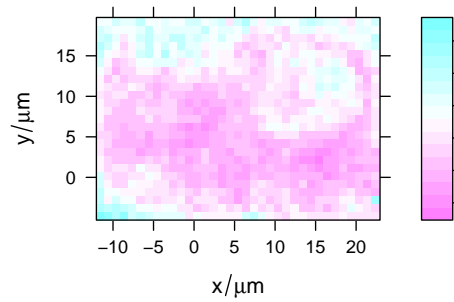


plots an intensity over a single other data column. e.g.

- time series
- calibration
- depth profile

```
> plotc (flu)
```

## levelplot

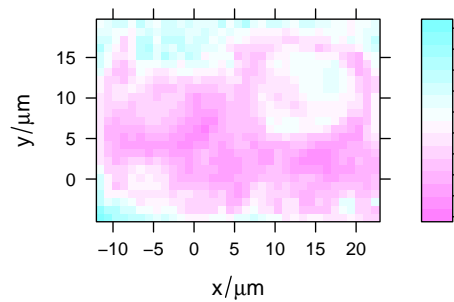


plots a false colour map, defined by a formula.

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

Warning: Only first wavelength is used for plotting

## plotmap



plotmap is a specialized version of levelplot. It uses a single value (e.g. average intensity or cluster membership) over two data columns (default  $x$  and  $y$ )

```
plotmap (chondro)
```

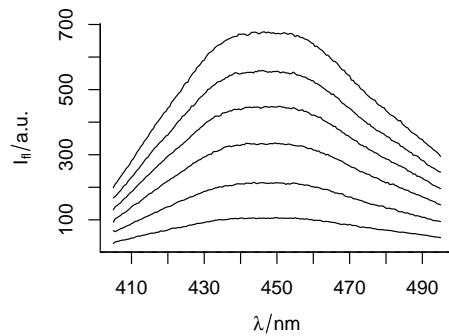
## 2 Arguments for plot

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.

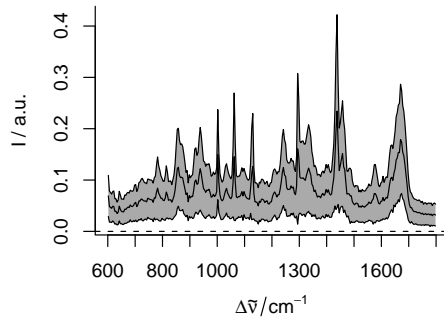
### plot (x, "spc")



is equivalent to `plotspc (flu)`

```
> plot (flu, "spc")
```

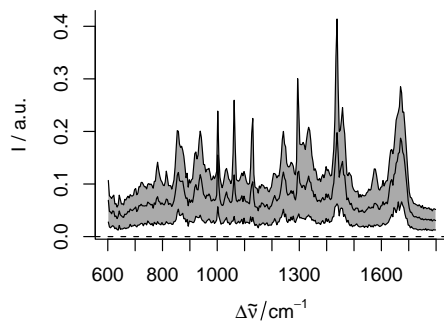
`plot (x, "spcmeansd")`



plots mean spectrum  $\pm 1$  standard deviation

`> plot (chondro.preproc, "spcmeansd")`

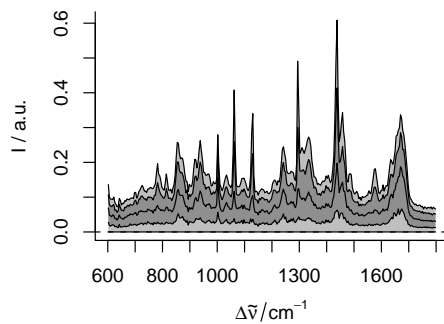
`plot (x, "spcprctile")`



plots median, 16<sup>th</sup> and 84<sup>th</sup> percentile for each wavelength. For Gaussian distributed data, 16<sup>th</sup>, 50<sup>th</sup> and 84<sup>th</sup> 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")`

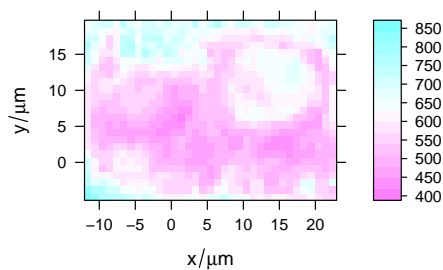
`plot (x, "spcprctl5")`



like "spcprctl" plus 5<sup>th</sup> and 95<sup>th</sup> percentile.

`> plot (chondro.preproc, "spcprctl5")`

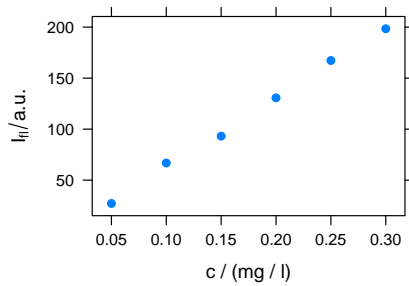
`plot (x, "map")`



is equivalent to `plotmap (chondro)`

`> plot (chondro, "map")`

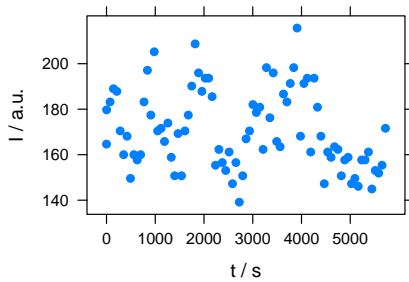
`plot (x, "c")`



is equivalent to `plotc (flu)`

`> plot (flu, "c")`

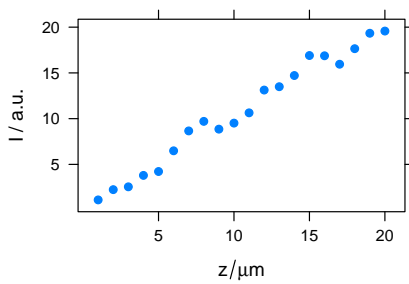
`plot (x, "ts")`



plots a time series plot, equivalent to `plotc (laser, spc ~ t)`

`> plot (laser, "ts")`

`plot (x, "depth")`

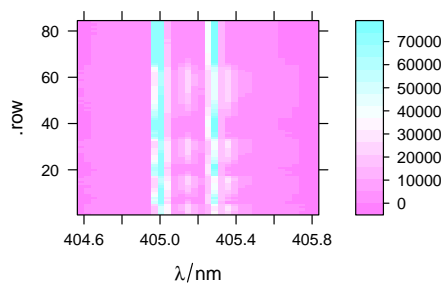


plots a depth profile plot, equivalent to `plotc (laser, spc ~ z)`

```
> depth.profile <- new ("hyperSpec",
+   spc = as.matrix (rnorm (20) + 1:20),
+   data = data.frame (z = 1 : 20),
+   label = list (spc = "I / a.u.",
+     z = expression (`/` (z, mu*m)),
+     .wavelength = expression (lambda)))
```

`> plot (depth.profile, "depth")`

`plot (x, "mat")`



plots the spectra matrix.

`> plot (laser, "mat")`

Equivalent to

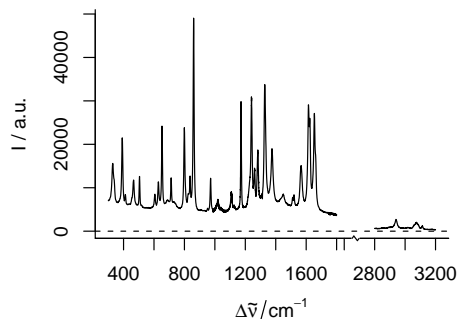
`> levelplot (spc ~ .wavelength * .row, laser)`

### 3 Spectra

plotspc

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

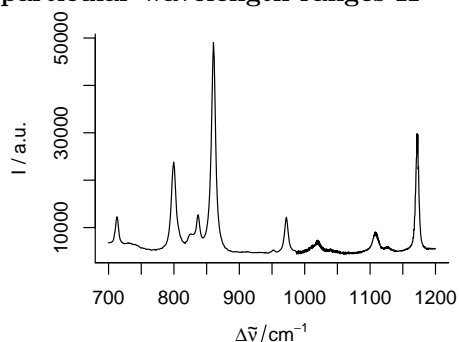
#### particular wavelength ranges



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`

```
> plotspc (paracetamol,  
+         wl.range = c (300 ~ 1800, 2800 ~ max),  
+         xoffset = 750)
```

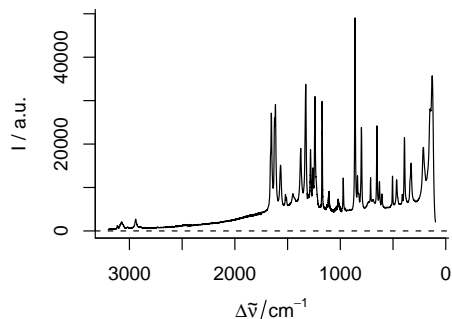
#### particular wavelength ranges II



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

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

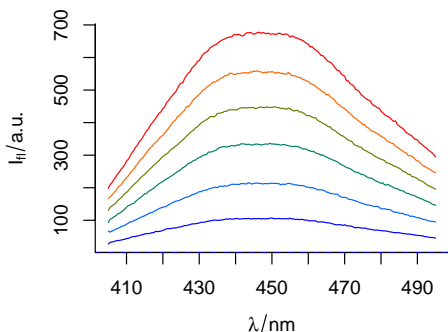
#### with reversed abscissa



use `wl.reverse = TRUE`

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

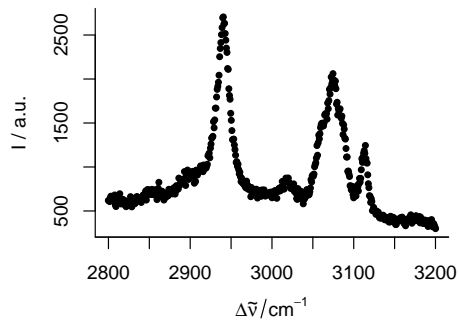
#### in different colours



use `col = vector.of.colours`

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

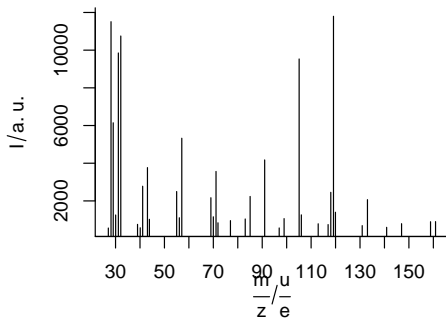
dots instead of lines



```
use lines.args = list (pch = 20, type = "p")

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

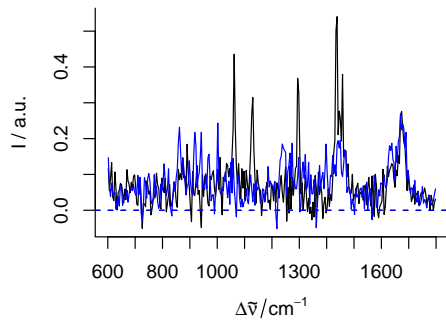
mass spectra



```
use lines.args = list (type = "h")

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

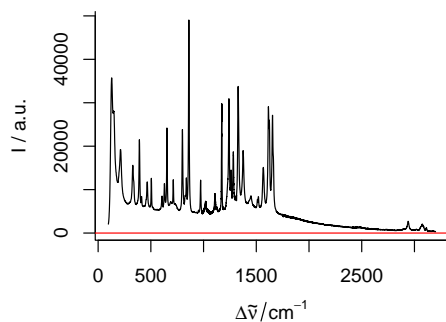
more spectra into an existing plot



```
use add = TRUE

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

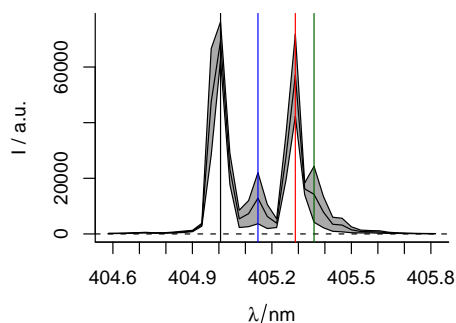
with different line at  $I = 0$



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

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

adding lines, etc.

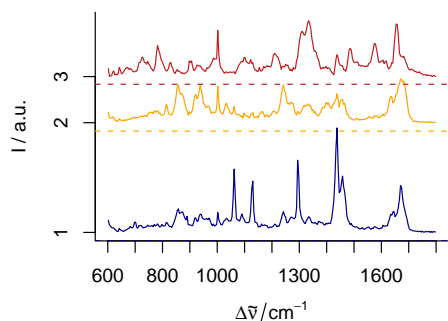


use `abline` for adding lines

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

### 3.1 Stacked spectra

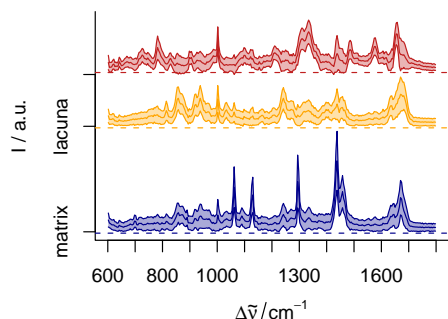
`stacked = TRUE`



use `stacked = TRUE`

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

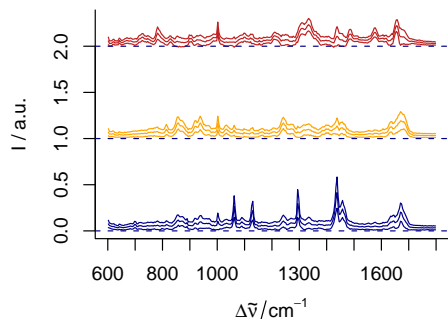
Stacking groups of spectra



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)
```

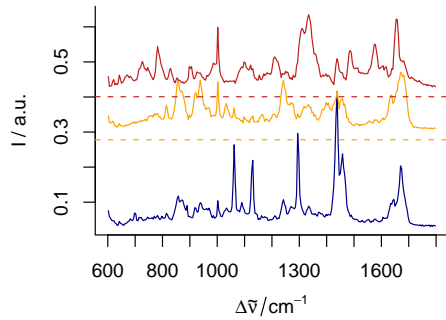
Manually giving yoffset



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))
```

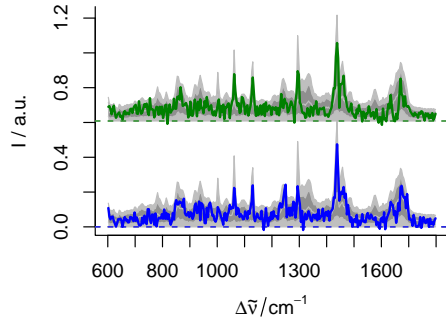




To obtain a denser stacking:

```
> ## 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:

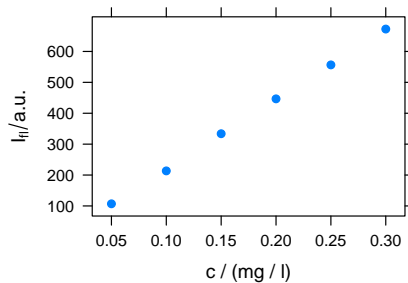
```
> ## coordinate system:
> yoffset <- apply (chondro.preproc, 2, quantile, c(0.05, 0.95))
> yoffset <- range (yoffset)
> plot(chondro.preproc[1],
+      plot.args = list (ylim = c (0, 2) * yoffset),
+      lines.args = list( type = "n")
+      )
> yoffset <- (0:1) * diff (yoffset)
> ## stacked spectra:
> for (i in 1 : 3){
+   plot(chondro.preproc, "spcprctl15",
+       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

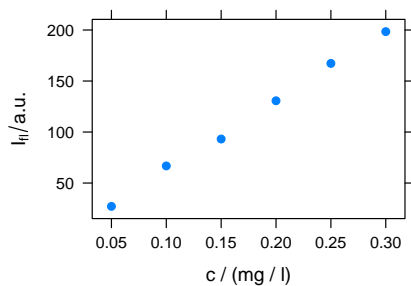
#### Intensities over concentration



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

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

## Integral Intensity over concentration

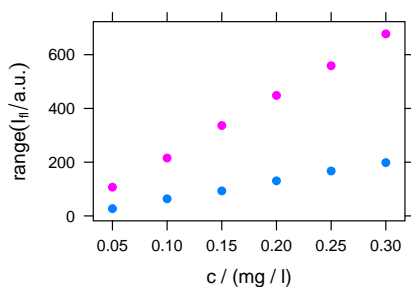


The default is to use the first intensity:

```
> plotc (flu)
```

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

## Other Summary Intensities over concentration

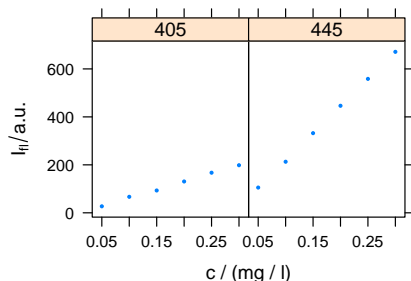


A function to compute a summary of the intensities before drawing can be used:

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

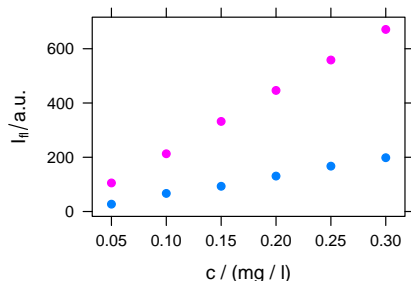
If `func` returns more than one value, the different results are accessible by `.wavelength`.

## Conditioning: plotting more traces separately



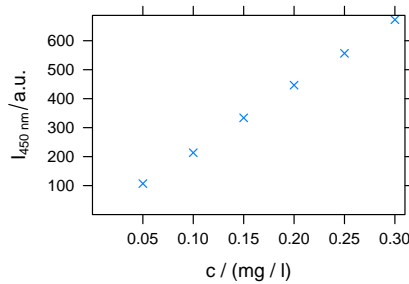
```
> plotc (flu [, c (405, 445)], spc ~ c | .wavelength,
+       cex = .3, scales = list (alternating = c(1, 1)))
```

## Grouping: plot more traces in one panel



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

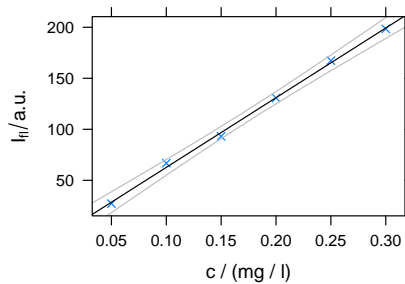
## Changing Axis Labels (and other parameters)



Arguments for `xyplot` can be given to `plotc`:

```
> plotc (flu [, 450],
+        ylab = expression (I ["450 nm"] / a.u.),
+        xlim = range (0, flu$c + .01),
+        ylim = range (0, flu$spc + 10),
+        pch = 4)
```

## Adding things to the plot: customized panel function



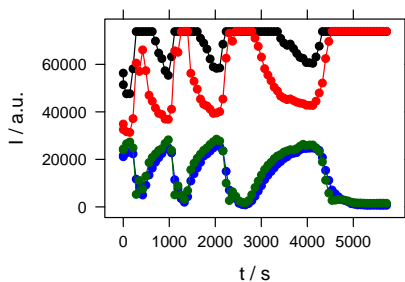
As `plotc` uses the *lattice* function `xyplot`, additions to the plot must be made via the panel function:

```
> panel.calibration <- function (x, y, ...,
+                               clim = range (x), level = .95) {
+   panel.xyplot (x, y, ...)
+   lm <- lm (y ~ x)
+   panel.abline (coef (lm), ...)
+   cx <- seq (clim [1], clim [2], length.out = 50)
+   cy <- predict (lm, data.frame (x = cx),
+                 interval = "confidence",
+                 level = level)
+   panel.lines (cx, cy [,2], col = "gray")
+   panel.lines (cx, cy [,3], col = "gray")
+ }
```

```
> plotc (flu [,405], panel = panel.calibration,
+        pch = 4, clim = c (0, 0.35), level = .99)
```

## 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:

```
> plotc (laser [, c (13, 17, 21, 23),
+               wl.index = TRUE],
+        spc ~ t,
+        groups = .wavelength,
+        type = "b",
+        col = c ("black", "blue", "red", "darkgreen"))
```

## 5 Levelplot

Levelplot can use two special column names:

`.wavelength` for the wavelengths

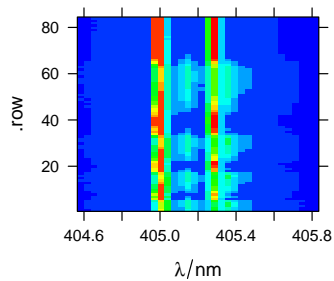
`.row` for the row index (i.e. spectrum number) in the data

## 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

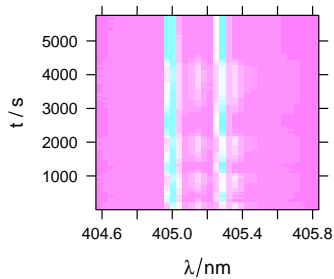


```
> plot (laser, "mat",  
+       col.regions = matlab.palette (20) )
```

is the same as

```
> levelplot (spc ~ .wavelength * .row,  
+            laser,  
+            col.regions = matlab.palette (20))
```

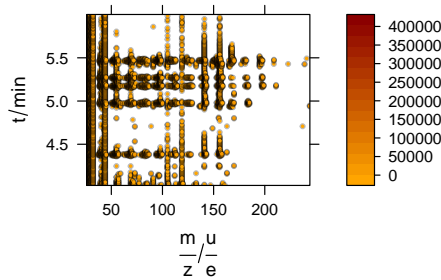
different y axis



Changing the y axis is only possible with `levelplot`:

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

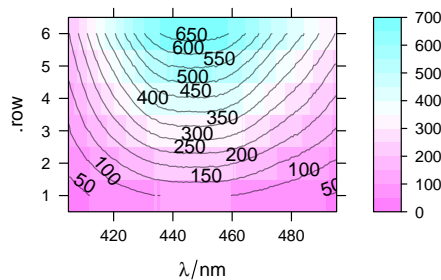
## different panel



```
> class (barbituates)
[1] "list"
> barbituates <- do.call (collapse, barbituates)
> barbituates <- orderwl (barbituates)

> levelplot (spc ~ .wavelength * z, barbituates,
+           panel = panel.levelplot.points,
+           cex = .5,
+           col.regions = colorRampPalette (c ("orange", "darkred"))
+           )
```

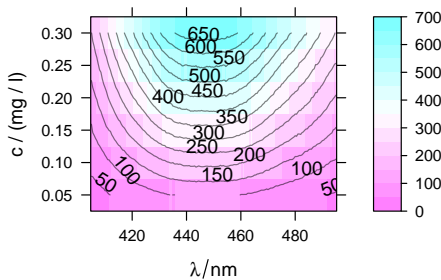
## contour lines with plot



Contourplots are possible with plot and levelplot:

```
> plot (flu,
+       "mat",
+       contour = TRUE,
+       labels = TRUE,
+       col = "#00000080",
+       at = seq (0, 700, by = 50))
```

## contour lines with levelplot



Contourplot with a different y axis:

```
> levelplot (spc ~ .wavelength * c,
+            flu,
+            contour = TRUE,
+            labels = TRUE,
+            col = "#00000080",
+            at = seq (0, 700, by = 50))
```

## 7 False-Colour Maps

plotmap

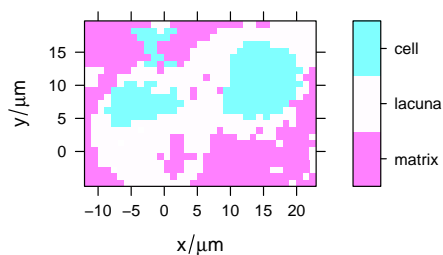
`plotmap` uses `levelplot`, a *lattice* function. Therefore, in loops, functions, Sweave chunks, etc. the lattice object needs to be printed explicitly by `print (plotmap (object))` ([R FAQ: Why do lattice/trellis graphics not work?](#)).

`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

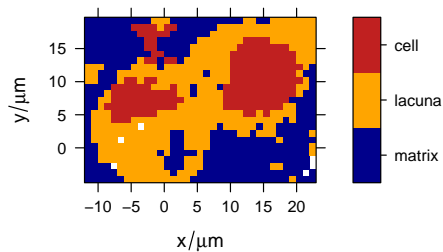
the function in argument *func*. The default setting is to plot the average intensity (no *z* and *func*=*mean*).

### plotting clusters



```
> plotmap (chondro, clusters ~ x * y)
```

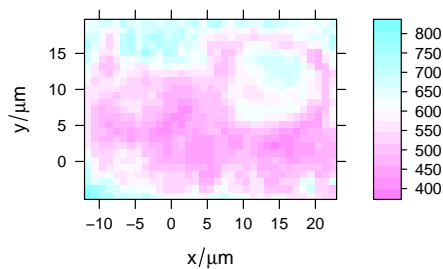
### different palette



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

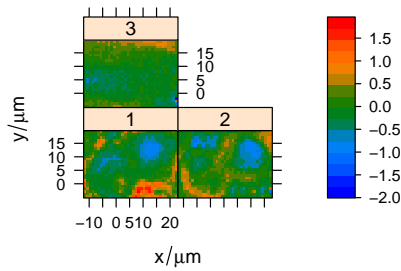
```
> cols <- c ("dark blue", "orange", "#C02020")
> print (plotmap (chondro,
+               clusters ~ x * y,
+               col.regions = cols))
```

### defined wavelengths



To plot a map with particular wavelengths use this:

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



Preprocessing of the data:

```
> baselines <- spc.fit.poly.below (chondro)
Fitting with npts.min = 15
> chondro <- chondro - baselines
> chondro <- sweep (chondro, 1, apply (chondro, 1, mean), "/")
> chondro <- sweep (chondro, 2, apply (chondro, 2, quantile,
+                                     0.05), "-")
> pca <- prcomp (~ spc, data = chondro$, center = TRUE)
> scores <- decomposition (chondro, pca$x,
+                           label.wavelength = "PC",
+                           label.spc = "score / a.u.")
```

The plot:

```
> levelplot (spc ~ x * y | as.factor(.wavelength),
+            scores [,1:3],
+            aspect = "iso",
+            col.regions = matlab.dark.palette(20))
```

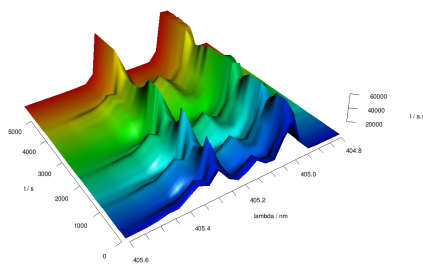
## Examples

**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 the 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)

> laser <- laser [,404.8 ~ 405.6]
> 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 / s", 'x+-', line = 2)
> mtext3d("lambda / nm", 'y--', line = 2)
> mtext3d("I / a.u.", edge = 'z--', line = 2.5)
```

## 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 *latticeist*. These packages allow easy customization of the plots and also identification of points.