

# Introduction to AlpsNMR (older API)

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**Abstract**

An introduction to the AlpsNMR package, showing the most relevant functions and a proposed workflow, using the older workflow.

**Package**

AlpsNMR 4.7.0

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## Introduction to AlpsNMR (older API)

The AlpsNMR package was written with two purposes in mind:

- to help **data analysts and NMR scientists** to work with NMR samples.
- to help **IT pipeline builders** implement automated methods for preprocessing.

Functions from this package written for data analysts and NMR scientists are prefixed with `nmr_`, while higher level functions written for IT pipeline builders are prefixed with `pipe_`. The main reason why all exported functions have a prefix is to make it easy for the user to discover the functions from the package. By typing `nmr_` RStudio will return the list of exported functions. In the R terminal, `nmr_` followed by the tab key (`<tab>`) twice will have the same effect. Other popular packages, follow similar approaches (e.g: `forcats`: `fct_*`, `stringr`: `str_*`).

This vignette is written for the first group. It assumes some prior basic knowledge of NMR and data analysis, as well as some basic R programming. In case you are interested in building pipelines with this package, you may want to open the file saved in this directory (run it on your computer):

```
pipeline_example <- system.file("pipeline-rmd", "pipeline_example.R", package = "AlpsNMR")
pipeline_example
```

```
library(BiocParallel)
library(AlpsNMR)
library(ggplot2)
```

## 1 Enable parallelization

This package is able to parallelize several functions through the use of the `BiocParallel` package. Whether to parallelize or not is left to the user that can control the parallelization registering backends. Please check the `BiocParallel` introduction for further details

```
library(BiocParallel)
#register(SerialParam(), default = TRUE) # disable parallelization
register(SnowParam(workers = 3, exportglobals = FALSE), default = TRUE) # enable parallelization with 4 workers
```

## 2 Data: The MeOH\_plasma\_extraction dataset

To explore the basics of the AlpsNMR package, we have included four NMR samples acquired in a 600 MHz Bruker instrument bundled with the package. The samples are pooled quality control plasma samples, that were extracted with methanol, and therefore only contain small molecules.

If you have installed this package, you can obtain the directory where the four samples are with the command

```
MeOH_plasma_extraction_dir <- system.file("dataset-demo", package = "AlpsNMR")
MeOH_plasma_extraction_dir
#> [1] "F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo"
```

The demo directory includes four samples (zipped) and a dummy Excel metadata file.

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```
fs::dir_ls(MeOH_plasma_extraction_dir)
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/10.zip
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/20.zip
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/30.zip
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/dummy_metadata.xlsx
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/README.txt
```

Given the name of the dataset, one may guess that the dataset was used to check the Methanol extraction in serum samples. The dummy metadata consists of dummy information, just for the sake of showing how this package can integrate external metadata. The excel file consists of two tidy tables, in two sheets.

```
MeOH_plasma_extraction_xlsx <- file.path(MeOH_plasma_extraction_dir, "dummy_metadata.xlsx")
exp_subj_id <- readxl::read_excel(MeOH_plasma_extraction_xlsx, sheet = 1)
subj_id_age <- readxl::read_excel(MeOH_plasma_extraction_xlsx, sheet = 2)
exp_subj_id
#> # A tibble: 3 x 3
#>   NMRExperiment SubjectID TimePoint
#>   <chr>         <chr>      <chr>
#> 1 10          Ana        baseline
#> 2 20          Ana        3 months
#> 3 30          Elia        baseline
subj_id_age
#> # A tibble: 2 x 2
#>   SubjectID Age
#>   <chr>     <dbl>
#> 1 Ana      29
#> 2 Elia     0
```

## 3 Loading samples

The function to read samples is called `nmr_read_samples`. It expects a character vector with the samples to load that can be paths to directories of Bruker format samples or paths to JDX files.

Additionally, this function can filter by pulse sequences (e.g. load only NOESY samples) or loading only metadata.

```
zip_files <- fs::dir_ls(MeOH_plasma_extraction_dir, glob = "*.zip")
zip_files
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/10.zip
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/20.zip
#> F:/biocbuild/bbs-3.20-bioc/tmpdir/RtmpgzB992/Rinst4468599d60ac/AlpsNMR/dataset-demo/30.zip
dataset <- nmr_read_samples(sample_names = zip_files)
dataset
#> An nmr_dataset (3 samples)
```

As we have not added any metadata to this dataset, the only column we see is the `NMRExperiment`:

```
nmr_meta_get(dataset, groups = "external")
#> # A tibble: 3 x 1
#>   NMRExperiment
#>   <chr>
#> 1 10
#> 2 20
#> 3 30
```

## 4 Adding metadata

Initially our dataset only has the NMRExperiment column:

```
nmr_meta_get(dataset, groups = "external")
#> # A tibble: 3 x 1
#>   NMRExperiment
#>   <chr>
#> 1 10
#> 2 20
#> 3 30
```

The exp\_subj\_id table we loaded links the NMRExperiment to the SubjectID.

As we already have the NMRExperiment column, we can use it as the merging column (note that both columns have the same column name to match the metadata such as group class, age, BMI...):

```
dataset <- nmr_meta_add(dataset, metadata = exp_subj_id, by = "NMRExperiment")
nmr_meta_get(dataset, groups = "external")
#> # A tibble: 3 x 3
#>   NMRExperiment SubjectID TimePoint
#>   <chr>          <chr>    <chr>
#> 1 10           Ana      baseline
#> 2 20           Ana      3 months
#> 3 30           Elia      baseline
```

If we have info from different files we can match them. For instance, now we have the SubjectID information so we can add the table that adds the SubjectID to the Age.

```
dataset <- nmr_meta_add(dataset, metadata = subj_id_age, by = "SubjectID")
nmr_meta_get(dataset, groups = "external")
#> # A tibble: 3 x 4
#>   NMRExperiment SubjectID TimePoint Age
#>   <chr>          <chr>    <chr>   <dbl>
#> 1 10           Ana      baseline 29
#> 2 20           Ana      3 months 29
#> 3 30           Elia      baseline 0
```

Now we have our metadata integrated in the dataset and we can make use of it in further data analysis steps.

## 5 Interpolation

1D NMR samples can be interpolated together, in order to arrange all the spectra into a matrix, with one row per sample. The main parameters we would need is the range of ppm values that we want to interpolate and the resolution.

We can see the ppm resolution by looking at the ppm axis of one sample:

```
ppm_res <- nmr_ppm_resolution(dataset)[[1]]
message("The ppm resolution is: ", format(ppm_res, digits = 2), " ppm")
#> The ppm resolution is: 0.00023 ppm
```

We can interpolate the dataset, obtaining an `nmr_dataset_1D` object:

```
dataset <- nmr_interpolate_1D(dataset, axis = c(min = -0.5, max = 10, by = 2.3E-4))
```

This operation changes the class of the object, as now the data is on a matrix. The dataset is now of class `nmr_dataset_1D`. The `axis` element is now a numeric vector and the `data_1r` element is a matrix.

## 6 Plotting samples

The AlpsNMR package offers the possibility to plot `nmr_dataset_1D` objects. Plotting many spectra with so many points is quite expensive so it is possible to include only some regions of the spectra or plot only some samples.

Use `?plot.nmr_dataset_1D` to check the parameters, among them:

- `NMRExperiment`: A character vector with the NMR experiments to plot
- `chemshift_range`: A ppm range to plot only a small region, or to reduce the resolution
- `interactive`: To make the plot interactive - `...``: Can be used to pass additional parameters such as `color = "SubjectID"` that are passed as aesthetics to ggplot.

```
plot(dataset, NMRExperiment = c("10", "30"), chemshift_range = c(2.2, 2.8))
```

### 6.1 Creating interactive plots

The option `interactive = TRUE` described above has some performance limitations. As high performance workaround, you can make many plots interactive with the function `plot_interactive`.

This function will use WebGL technologies to create a webpage that, once opened, allows you to interact with the plot.

Due to technical limitations, these plots need to be opened manually and can't be embedded in RMarkdown documents. Therefore, the function saves the plot in the directory for further exploration. Additionally, some old web browsers may not be able to display these interactive plots correctly.

```
plt <- plot(dataset, NMRExperiment = c("10", "30"), chemshift_range = c(2.2, 2.8))
plot_interactive(plt, "plot_region.html")
```

### 7 Exclude regions

---

Some regions can easily be excluded from the spectra with `nmr_exclude_region`. Note that the regions are fully removed and not zeroed, as using zeros complicates a lot the implementation<sup>1</sup> and has little advantages.

```
regions_to_exclude <- list(water = c(4.6, 5), methanol = c(3.33, 3.39))
dataset <- nmr_exclude_region(dataset, exclude = regions_to_exclude)
plot(dataset, chemshift_range = c(4.2, 5.5))
```

### 8 Filter samples

---

Maybe we just want to analyze a subset of the data, e.g., only a class group or a particular gender. We can filter some samples according to their metadata as follows:

```
samples_10_20 <- filter(dataset, SubjectID == "Ana")
nmr_meta_get(samples_10_20, groups = "external")
#> # A tibble: 2 x 4
#>   NMRExperiment SubjectID TimePoint   Age
#>   <chr>         <chr>    <chr>    <dbl>
#> 1 10           Ana      baseline    29
#> 2 20           Ana      3 months    29
```

### 9 Robust PCA for outlier detection

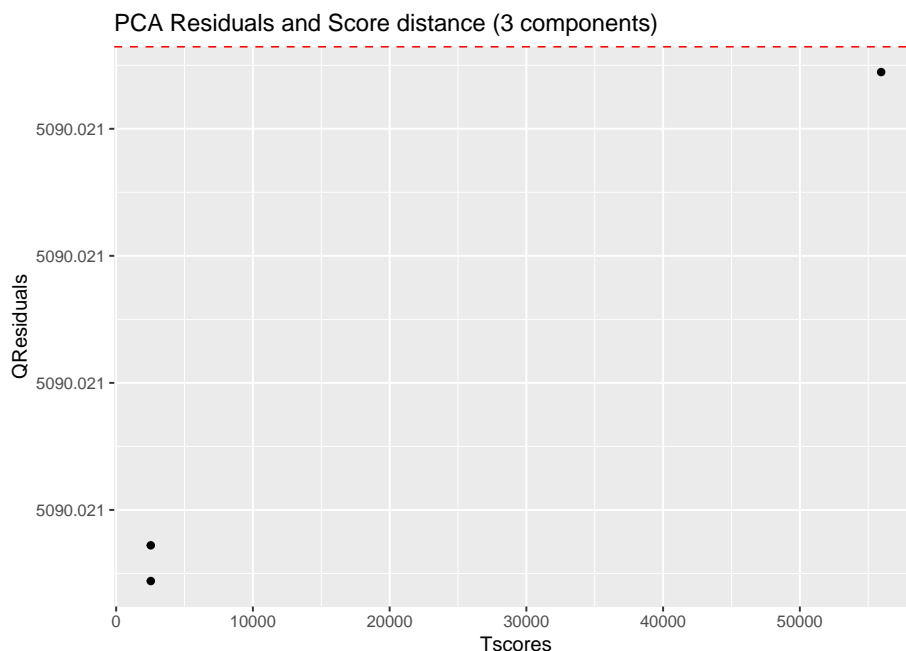
---

The AlpsNMR package includes robust PCA analysis for outlier detection. With such a small demo dataset, it is not practical to use, but check out the documentation of `nmr_pca_outliers_*` functions.

```
pca_outliers_rob <- nmr_pca_outliers_robust(dataset, ncomp = 3)
nmr_pca_outliers_plot(dataset, pca_outliers_rob)
```

---

<sup>1</sup>e.g. it can inadvertently distort the PQN normalization results



## 10 Baseline removal

Spectra may display an unstable baseline, specially when processing blood/fecal blood/fecal samples. If so, `nmr_baseline_removal` subtract the baseline by means of Asymmetric Least Squares method.

See before:

```
plot(dataset, chemshift_range = c(3.5,3.8))
```

And after:

```
dataset = nmr_baseline_removal(dataset, lambda = 6, p = 0.01)  
plot(dataset, chemshift_range = c(3.5,3.8))
```

## 11 Peak detection

The peak detection is performed on short spectra segments using a continuous wavelet transform. See `?nmr_detect_peaks` for more information.

Our current approach relies on the use of the baseline threshold (`baselineThresh`) automatic calculated (see `?nmr_baseline_threshold`) and the Signal to Noise Threshold (`SNR.Th`) to discriminate valid peaks from noise.

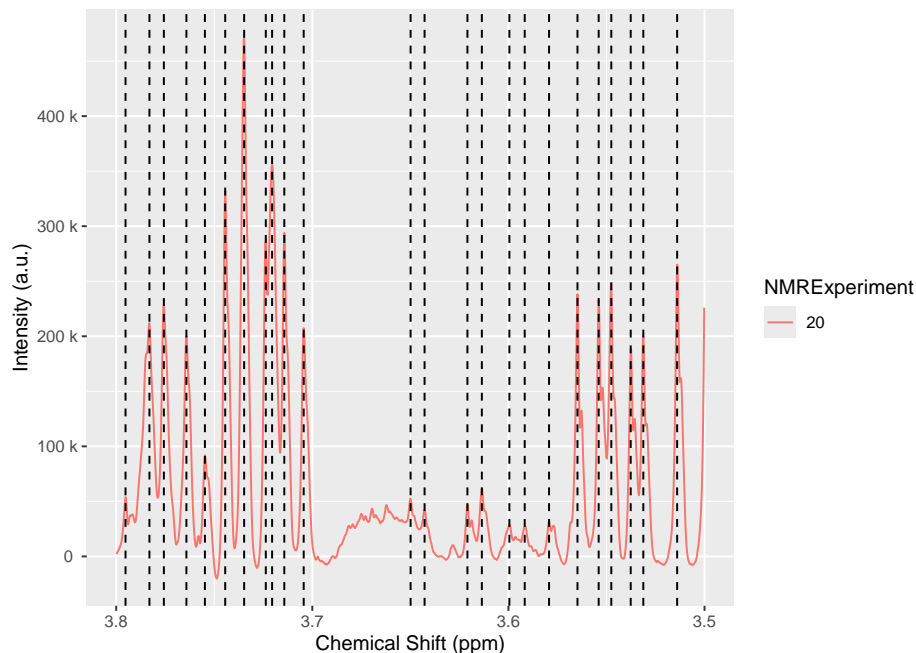
The combination of the `baselineThresh` and the `SNR.Th` optimizes the number of actual peaks from noise.

The advantage of the `SNR.Th` method is that it estimates the noise level on each spectra region independently, so in practice it can be used as a dynamic baseline threshold level.



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```
peak_table <- nmr_detect_peaks(dataset,
                              nDivRange_ppm = 0.1,
                              scales = seq(1, 16, 2),
                              baselineThresh = NULL, SNR.Th = 3)
NMExp_ref <- nmr_align_find_ref(dataset, peak_table)
message("Your reference is NMRExperiment ", NMExp_ref)
#> Your reference is NMRExperiment 30
nmr_detect_peaks_plot(dataset, peak_table, NMRExperiment = "20", chemshift_range = c(3.5, 3.8))
```



## 12 Spectra alignment

To align the sample, we use the `nmr_align` function, which in turn uses a hierarchical clustering method (see `?nmr_align` for further details).

The `maxShift_ppm` limits the maximum shift allowed for the spectra.

```
nmr_exp_ref <- nmr_align_find_ref(dataset, peak_table)
dataset_align <- nmr_align(dataset, peak_table, nmr_exp_ref, maxShift_ppm = 0.0015, acceptLostPeak = FALSE)

plot(dataset, chemshift_range = c(3.025, 3.063))
plot(dataset_align, chemshift_range = c(3.025, 3.063))
```

## 13 Normalization

There are multiple normalization techniques available. The most strongly recommended is the `pqn` normalization, but it may not be fully reliable when the number of samples is small, as it needs a computation of the median spectra. Nevertheless, it is possible to compute it:

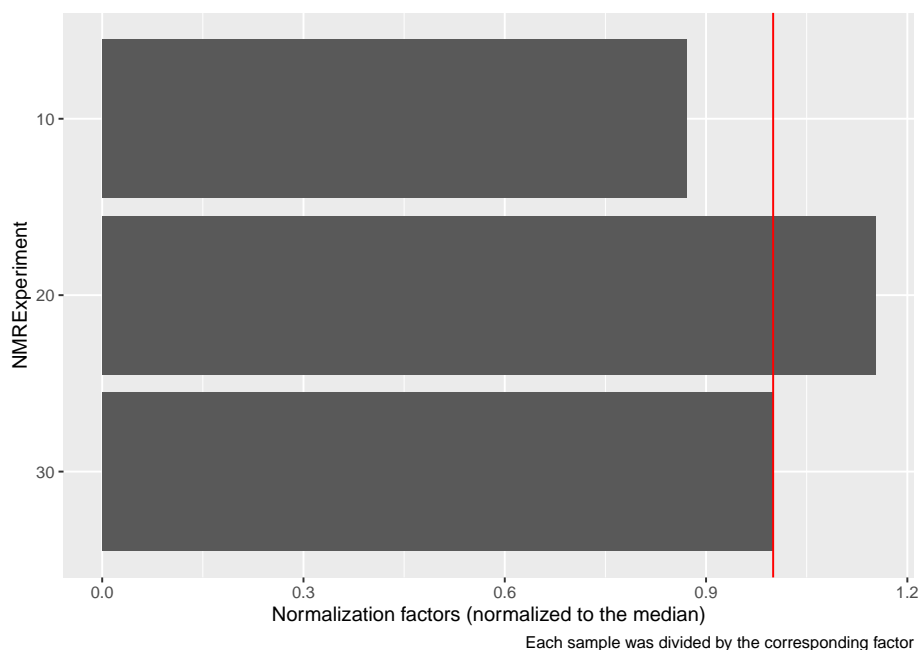
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```
dataset_norm <- nmr_normalize(dataset_align, method = "pqn")
#> Warning: There are not enough samples for reliably estimating the median spectra
#> i The Probabalistic Quotient Normalization requires several samples to compute the median spectra. Your n
#> i Review your peaks before and after normalization to ensure there are no big distortions
```

The AlpsNMR package offers the possibility to extract additional normalization information with `nmr_normalize_extra_info(dataset)`, to explore the normalization factors applied to each sample:

The plot shows the dispersion with respect to the median of the normalization factors, and can highlight samples with abnormally large or small normalization factors.

```
diagnostic <- nmr_normalize_extra_info(dataset_norm)
diagnostic$norm_factor
#>   NMRExperiment norm_factor norm_factor_norm
#> 1             10    0.8706322      0.8706322
#> 2             20    1.1523450      1.1523450
#> 3             30    1.0000000      1.0000000
diagnostic$plot
```



## 14 Peak integration

### 14.1 1. Integration based on peak center and width

If we want to integrate the whole spectra, we need ppm from the `peak_table`. See Peak detection section. The function `nmr_integrate_peak_positions` generates a new `nmr_dataset_1D` object containing the integrals from the `peak_table` (ppm values corresponding to detected peaks).

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```
peak_table_integration = nmr_integrate_peak_positions(  
  samples = dataset_norm,  
  peak_pos_ppm = peak_table$ppm,  
  peak_width_ppm = 0.006)
```

```
#> New names:
```

```
#> * `ppm_-0.0002` -> `ppm_-0.0002...6`  
#> * `ppm_0.8607` -> `ppm_0.8607...22`  
#> * `ppm_0.8655` -> `ppm_0.8655...23`  
#> * `ppm_0.8729` -> `ppm_0.8729...24`  
#> * `ppm_0.8784` -> `ppm_0.8784...25`  
#> * `ppm_0.8910` -> `ppm_0.8910...26`  
#> * `ppm_0.9032` -> `ppm_0.9032...27`  
#> * `ppm_0.9085` -> `ppm_0.9085...28`  
#> * `ppm_0.9196` -> `ppm_0.9196...30`  
#> * `ppm_0.9313` -> `ppm_0.9313...31`  
#> * `ppm_0.9437` -> `ppm_0.9437...32`  
#> * `ppm_0.9554` -> `ppm_0.9554...33`  
#> * `ppm_0.9658` -> `ppm_0.9658...34`  
#> * `ppm_0.9902` -> `ppm_0.9902...36`  
#> * `ppm_1.0019` -> `ppm_1.0019...37`  
#> * `ppm_1.0093` -> `ppm_1.0093...38`  
#> * `ppm_1.0208` -> `ppm_1.0208...39`  
#> * `ppm_1.0410` -> `ppm_1.0410...40`  
#> * `ppm_1.0716` -> `ppm_1.0716...43`  
#> * `ppm_1.0803` -> `ppm_1.0803...44`  
#> * `ppm_1.1505` -> `ppm_1.1505...46`  
#> * `ppm_1.1979` -> `ppm_1.1979...47`  
#> * `ppm_1.2084` -> `ppm_1.2084...48`  
#> * `ppm_1.2179` -> `ppm_1.2179...49`  
#> * `ppm_1.3262` -> `ppm_1.3262...54`  
#> * `ppm_1.3377` -> `ppm_1.3377...55`  
#> * `ppm_1.4424` -> `ppm_1.4424...57`  
#> * `ppm_1.4913` -> `ppm_1.4913...59`  
#> * `ppm_1.5693` -> `ppm_1.5693...66`  
#> * `ppm_1.6195` -> `ppm_1.6195...72`  
#> * `ppm_1.6443` -> `ppm_1.6443...75`  
#> * `ppm_1.6562` -> `ppm_1.6562...77`  
#> * `ppm_1.6799` -> `ppm_1.6799...80`  
#> * `ppm_1.6990` -> `ppm_1.6990...86`  
#> * `ppm_1.7351` -> `ppm_1.7351...92`  
#> * `ppm_1.7614` -> `ppm_1.7614...95`  
#> * `ppm_1.8131` -> `ppm_1.8131...97`  
#> * `ppm_1.8748` -> `ppm_1.8748...103`  
#> * `ppm_1.9210` -> `ppm_1.9210...106`  
#> * `ppm_1.9313` -> `ppm_1.9313...107`  
#> * `ppm_2.0026` -> `ppm_2.0026...114`  
#> * `ppm_2.0194` -> `ppm_2.0194...117`  
#> * `ppm_2.0307` -> `ppm_2.0307...118`  
#> * `ppm_2.0422` -> `ppm_2.0422...119`  
#> * `ppm_2.0806` -> `ppm_2.0806...123`  
#> * `ppm_2.1149` -> `ppm_2.1149...126`
```

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```
#> * `ppm_2.1192` -> `ppm_2.1192...127`  
#> * `ppm_2.1273` -> `ppm_2.1273...128`  
#> * `ppm_2.1324` -> `ppm_2.1324...129`  
#> * `ppm_2.1572` -> `ppm_2.1572...134`  
#> * `ppm_2.1655` -> `ppm_2.1655...135`  
#> * `ppm_2.2835` -> `ppm_2.2835...143`  
#> * `ppm_2.2950` -> `ppm_2.2950...144`  
#> * `ppm_2.3251` -> `ppm_2.3251...147`  
#> * `ppm_2.3451` -> `ppm_2.3451...149`  
#> * `ppm_2.3513` -> `ppm_2.3513...150`  
#> * `ppm_2.3575` -> `ppm_2.3575...151`  
#> * `ppm_2.3644` -> `ppm_2.3644...152`  
#> * `ppm_2.3711` -> `ppm_2.3711...153`  
#> * `ppm_2.4196` -> `ppm_2.4196...157`  
#> * `ppm_2.4912` -> `ppm_2.4912...166`  
#> * `ppm_2.4953` -> `ppm_2.4953...167`  
#> * `ppm_2.5284` -> `ppm_2.5284...171`  
#> * `ppm_2.6216` -> `ppm_2.6216...177`  
#> * `ppm_2.7577` -> `ppm_2.7577...188`  
#> * `ppm_2.9332` -> `ppm_2.9332...199`  
#> * `ppm_3.0305` -> `ppm_3.0305...206`  
#> * `ppm_3.0404` -> `ppm_3.0404...207`  
#> * `ppm_3.0489` -> `ppm_3.0489...208`  
#> * `ppm_3.2081` -> `ppm_3.2081...218`  
#> * `ppm_3.2154` -> `ppm_3.2154...219`  
#> * `ppm_3.2320` -> `ppm_3.2320...220`  
#> * `ppm_3.2382` -> `ppm_3.2382...221`  
#> * `ppm_3.2672` -> `ppm_3.2672...223`  
#> * `ppm_3.2706` -> `ppm_3.2706...224`  
#> * `ppm_3.4031` -> `ppm_3.4031...228`  
#> * `ppm_3.4243` -> `ppm_3.4243...231`  
#> * `ppm_3.4829` -> `ppm_3.4829...238`  
#> * `ppm_3.4981` -> `ppm_3.4981...240`  
#> * `ppm_3.5135` -> `ppm_3.5135...241`  
#> * `ppm_3.5372` -> `ppm_3.5372...243`  
#> * `ppm_3.7042` -> `ppm_3.7042...254`  
#> * `ppm_3.7141` -> `ppm_3.7141...255`  
#> * `ppm_3.7200` -> `ppm_3.7200...256`  
#> * `ppm_3.7237` -> `ppm_3.7237...257`  
#> * `ppm_3.7348` -> `ppm_3.7348...258`  
#> * `ppm_3.7442` -> `ppm_3.7442...259`  
#> * `ppm_3.7757` -> `ppm_3.7757...262`  
#> * `ppm_3.8233` -> `ppm_3.8233...265`  
#> * `ppm_3.8438` -> `ppm_3.8438...268`  
#> * `ppm_3.8528` -> `ppm_3.8528...270`  
#> * `ppm_3.8886` -> `ppm_3.8886...273`  
#> * `ppm_3.8923` -> `ppm_3.8923...274`  
#> * `ppm_3.9091` -> `ppm_3.9091...275`  
#> * `ppm_3.9128` -> `ppm_3.9128...276`  
#> * `ppm_3.9333` -> `ppm_3.9333...277`  
#> * `ppm_3.9493` -> `ppm_3.9493...278`
```

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```
#> * `ppm_3.9917` -> `ppm_3.9917...282`  
#> * `ppm_4.0020` -> `ppm_4.0020...283`  
#> * `ppm_4.0108` -> `ppm_4.0108...284`  
#> * `ppm_4.1205` -> `ppm_4.1205...288`  
#> * `ppm_4.1320` -> `ppm_4.1320...289`  
#> * `ppm_4.1927` -> `ppm_4.1927...291`  
#> * `ppm_4.2281` -> `ppm_4.2281...293`  
#> * `ppm_4.2398` -> `ppm_4.2398...295`  
#> * `ppm_4.2566` -> `ppm_4.2566...299`  
#> * `ppm_4.4376` -> `ppm_4.4376...315`  
#> * `ppm_4.4531` -> `ppm_4.4531...317`  
#> * `ppm_4.5218` -> `ppm_4.5218...320`  
#> * `ppm_4.5221` -> `ppm_4.5221...321`  
#> * `ppm_4.5906` -> `ppm_4.5906...324`  
#> * `ppm_4.5991` -> `ppm_4.5991...325`  
#> * `ppm_5.2369` -> `ppm_5.2369...329`  
#> * `ppm_5.4154` -> `ppm_5.4154...336`  
#> * `ppm_6.1026` -> `ppm_6.1026...350`  
#> * `ppm_6.9009` -> `ppm_6.9009...361`  
#> * `ppm_6.9154` -> `ppm_6.9154...362`  
#> * `ppm_7.2894` -> `ppm_7.2894...372`  
#> * `ppm_7.3014` -> `ppm_7.3014...373`  
#> * `ppm_7.3409` -> `ppm_7.3409...376`  
#> * `ppm_7.3816` -> `ppm_7.3816...380`  
#> * `ppm_7.4881` -> `ppm_7.4881...387`  
#> * `ppm_7.5553` -> `ppm_7.5553...390`  
#> * `ppm_8.2458` -> `ppm_8.2458...407`  
#> * `ppm_8.3515` -> `ppm_8.3515...409`  
#> * `ppm_8.4608` -> `ppm_8.4608...411`  
#> * `ppm_-0.0002` -> `ppm_-0.0002...422`  
#> * `ppm_0.8110` -> `ppm_0.8110...438`  
#> * `ppm_0.8552` -> `ppm_0.8552...442`  
#> * `ppm_0.8607` -> `ppm_0.8607...443`  
#> * `ppm_0.8784` -> `ppm_0.8784...446`  
#> * `ppm_0.8910` -> `ppm_0.8910...447`  
#> * `ppm_0.9085` -> `ppm_0.9085...449`  
#> * `ppm_0.9196` -> `ppm_0.9196...450`  
#> * `ppm_0.9313` -> `ppm_0.9313...451`  
#> * `ppm_0.9437` -> `ppm_0.9437...452`  
#> * `ppm_0.9554` -> `ppm_0.9554...453`  
#> * `ppm_0.9658` -> `ppm_0.9658...454`  
#> * `ppm_1.0093` -> `ppm_1.0093...458`  
#> * `ppm_1.0205` -> `ppm_1.0205...459`  
#> * `ppm_1.0716` -> `ppm_1.0716...463`  
#> * `ppm_1.0803` -> `ppm_1.0803...464`  
#> * `ppm_1.1505` -> `ppm_1.1505...468`  
#> * `ppm_1.2084` -> `ppm_1.2084...470`  
#> * `ppm_1.2179` -> `ppm_1.2179...471`  
#> * `ppm_1.2294` -> `ppm_1.2294...472`  
#> * `ppm_1.3262` -> `ppm_1.3262...476`  
#> * `ppm_1.3377` -> `ppm_1.3377...477`
```

## Introduction to AlpsNMR (older API)

```
#> * `ppm_1.4308` -> `ppm_1.4308...478`
#> * `ppm_1.4424` -> `ppm_1.4424...479`
#> * `ppm_1.4794` -> `ppm_1.4794...480`
#> * `ppm_1.5693` -> `ppm_1.5693...487`
#> * `ppm_1.6195` -> `ppm_1.6195...491`
#> * `ppm_1.6319` -> `ppm_1.6319...492`
#> * `ppm_1.6441` -> `ppm_1.6441...494`
#> * `ppm_1.6560` -> `ppm_1.6560...495`
#> * `ppm_1.6562` -> `ppm_1.6562...496`
#> * `ppm_1.6675` -> `ppm_1.6675...498`
#> * `ppm_1.6747` -> `ppm_1.6747...499`
#> * `ppm_1.6799` -> `ppm_1.6799...501`
#> * `ppm_1.6990` -> `ppm_1.6990...503`
#> * `ppm_1.7075` -> `ppm_1.7075...504`
#> * `ppm_1.7614` -> `ppm_1.7614...511`
#> * `ppm_1.8752` -> `ppm_1.8752...522`
#> * `ppm_1.9812` -> `ppm_1.9812...530`
#> * `ppm_2.0422` -> `ppm_2.0422...537`
#> * `ppm_2.0923` -> `ppm_2.0923...542`
#> * `ppm_2.1192` -> `ppm_2.1192...545`
#> * `ppm_2.1273` -> `ppm_2.1273...546`
#> * `ppm_2.1324` -> `ppm_2.1324...547`
#> * `ppm_2.1572` -> `ppm_2.1572...550`
#> * `ppm_2.1655` -> `ppm_2.1655...551`
#> * `ppm_2.2515` -> `ppm_2.2515...558`
#> * `ppm_2.2717` -> `ppm_2.2717...561`
#> * `ppm_2.3451` -> `ppm_2.3451...570`
#> * `ppm_2.3513` -> `ppm_2.3513...571`
#> * `ppm_2.3575` -> `ppm_2.3575...572`
#> * `ppm_2.3644` -> `ppm_2.3644...573`
#> * `ppm_2.4063` -> `ppm_2.4063...577`
#> * `ppm_2.4196` -> `ppm_2.4196...578`
#> * `ppm_2.4456` -> `ppm_2.4456...581`
#> * `ppm_2.4953` -> `ppm_2.4953...586`
#> * `ppm_2.5284` -> `ppm_2.5284...590`
#> * `ppm_2.7994` -> `ppm_2.7994...606`
#> * `ppm_2.9339` -> `ppm_2.9339...615`
#> * `ppm_3.2154` -> `ppm_3.2154...638`
#> * `ppm_3.2706` -> `ppm_3.2706...643`
#> * `ppm_3.3930` -> `ppm_3.3930...646`
#> * `ppm_3.4185` -> `ppm_3.4185...648`
#> * `ppm_3.5793` -> `ppm_3.5793...663`
#> * `ppm_3.6428` -> `ppm_3.6428...668`
#> * `ppm_3.6499` -> `ppm_3.6499...669`
#> * `ppm_3.7237` -> `ppm_3.7237...673`
#> * `ppm_3.7348` -> `ppm_3.7348...674`
#> * `ppm_3.7642` -> `ppm_3.7642...677`
#> * `ppm_3.7757` -> `ppm_3.7757...678`
#> * `ppm_3.8438` -> `ppm_3.8438...683`
#> * `ppm_3.8923` -> `ppm_3.8923...689`
#> * `ppm_3.9128` -> `ppm_3.9128...691`
```

## Introduction to AlpsNMR (older API)

```
#> * `ppm_3.9917` -> `ppm_3.9917...698`  
#> * `ppm_4.0020` -> `ppm_4.0020...699`  
#> * `ppm_4.2281` -> `ppm_4.2281...712`  
#> * `ppm_4.2398` -> `ppm_4.2398...714`  
#> * `ppm_4.2566` -> `ppm_4.2566...717`  
#> * `ppm_4.5089` -> `ppm_4.5089...735`  
#> * `ppm_4.5092` -> `ppm_4.5092...736`  
#> * `ppm_4.5221` -> `ppm_4.5221...737`  
#> * `ppm_4.5991` -> `ppm_4.5991...742`  
#> * `ppm_5.0973` -> `ppm_5.0973...744`  
#> * `ppm_5.4154` -> `ppm_5.4154...753`  
#> * `ppm_5.4220` -> `ppm_5.4220...754`  
#> * `ppm_6.1026` -> `ppm_6.1026...771`  
#> * `ppm_6.9154` -> `ppm_6.9154...790`  
#> * `ppm_6.9626` -> `ppm_6.9626...791`  
#> * `ppm_7.2066` -> `ppm_7.2066...798`  
#> * `ppm_7.3014` -> `ppm_7.3014...803`  
#> * `ppm_7.3409` -> `ppm_7.3409...806`  
#> * `ppm_7.4881` -> `ppm_7.4881...816`  
#> * `ppm_7.5415` -> `ppm_7.5415...819`  
#> * `ppm_8.2458` -> `ppm_8.2458...839`  
#> * `ppm_8.4608` -> `ppm_8.4608...842`  
#> * `ppm_-0.0002` -> `ppm_-0.0002...853`  
#> * `ppm_0.8110` -> `ppm_0.8110...865`  
#> * `ppm_0.8552` -> `ppm_0.8552...869`  
#> * `ppm_0.8655` -> `ppm_0.8655...871`  
#> * `ppm_0.8729` -> `ppm_0.8729...872`  
#> * `ppm_0.9032` -> `ppm_0.9032...875`  
#> * `ppm_0.9085` -> `ppm_0.9085...876`  
#> * `ppm_0.9196` -> `ppm_0.9196...877`  
#> * `ppm_0.9313` -> `ppm_0.9313...878`  
#> * `ppm_0.9902` -> `ppm_0.9902...883`  
#> * `ppm_1.0019` -> `ppm_1.0019...884`  
#> * `ppm_1.0205` -> `ppm_1.0205...886`  
#> * `ppm_1.0208` -> `ppm_1.0208...887`  
#> * `ppm_1.0410` -> `ppm_1.0410...888`  
#> * `ppm_1.0716` -> `ppm_1.0716...891`  
#> * `ppm_1.0803` -> `ppm_1.0803...892`  
#> * `ppm_1.1505` -> `ppm_1.1505...896`  
#> * `ppm_1.1979` -> `ppm_1.1979...897`  
#> * `ppm_1.2084` -> `ppm_1.2084...898`  
#> * `ppm_1.2179` -> `ppm_1.2179...899`  
#> * `ppm_1.2294` -> `ppm_1.2294...900`  
#> * `ppm_1.3262` -> `ppm_1.3262...904`  
#> * `ppm_1.3377` -> `ppm_1.3377...905`  
#> * `ppm_1.4308` -> `ppm_1.4308...906`  
#> * `ppm_1.4424` -> `ppm_1.4424...907`  
#> * `ppm_1.4794` -> `ppm_1.4794...908`  
#> * `ppm_1.4913` -> `ppm_1.4913...909`  
#> * `ppm_1.6319` -> `ppm_1.6319...919`  
#> * `ppm_1.6441` -> `ppm_1.6441...920`
```

## Introduction to AlpsNMR (older API)

```
#> * `ppm_1.6443` -> `ppm_1.6443...921`  
#> * `ppm_1.6560` -> `ppm_1.6560...922`  
#> * `ppm_1.6562` -> `ppm_1.6562...923`  
#> * `ppm_1.6675` -> `ppm_1.6675...924`  
#> * `ppm_1.6747` -> `ppm_1.6747...925`  
#> * `ppm_1.6799` -> `ppm_1.6799...927`  
#> * `ppm_1.6990` -> `ppm_1.6990...930`  
#> * `ppm_1.7075` -> `ppm_1.7075...933`  
#> * `ppm_1.7351` -> `ppm_1.7351...937`  
#> * `ppm_1.7614` -> `ppm_1.7614...940`  
#> * `ppm_1.8131` -> `ppm_1.8131...945`  
#> * `ppm_1.8748` -> `ppm_1.8748...952`  
#> * `ppm_1.8752` -> `ppm_1.8752...953`  
#> * `ppm_1.9210` -> `ppm_1.9210...956`  
#> * `ppm_1.9313` -> `ppm_1.9313...957`  
#> * `ppm_1.9812` -> `ppm_1.9812...962`  
#> * `ppm_2.0026` -> `ppm_2.0026...964`  
#> * `ppm_2.0194` -> `ppm_2.0194...967`  
#> * `ppm_2.0307` -> `ppm_2.0307...968`  
#> * `ppm_2.0806` -> `ppm_2.0806...974`  
#> * `ppm_2.0923` -> `ppm_2.0923...975`  
#> * `ppm_2.1149` -> `ppm_2.1149...977`  
#> * `ppm_2.1192` -> `ppm_2.1192...978`  
#> * `ppm_2.1273` -> `ppm_2.1273...979`  
#> * `ppm_2.2515` -> `ppm_2.2515...992`  
#> * `ppm_2.2717` -> `ppm_2.2717...995`  
#> * `ppm_2.2835` -> `ppm_2.2835...997`  
#> * `ppm_2.2950` -> `ppm_2.2950...999`  
#> * `ppm_2.3251` -> `ppm_2.3251...1002`  
#> * `ppm_2.3513` -> `ppm_2.3513...1005`  
#> * `ppm_2.3575` -> `ppm_2.3575...1006`  
#> * `ppm_2.3711` -> `ppm_2.3711...1008`  
#> * `ppm_2.4063` -> `ppm_2.4063...1011`  
#> * `ppm_2.4196` -> `ppm_2.4196...1012`  
#> * `ppm_2.4456` -> `ppm_2.4456...1015`  
#> * `ppm_2.4912` -> `ppm_2.4912...1020`  
#> * `ppm_2.6216` -> `ppm_2.6216...1029`  
#> * `ppm_2.7577` -> `ppm_2.7577...1039`  
#> * `ppm_2.7994` -> `ppm_2.7994...1041`  
#> * `ppm_2.9332` -> `ppm_2.9332...1050`  
#> * `ppm_2.9339` -> `ppm_2.9339...1051`  
#> * `ppm_3.0305` -> `ppm_3.0305...1058`  
#> * `ppm_3.0404` -> `ppm_3.0404...1059`  
#> * `ppm_3.0489` -> `ppm_3.0489...1060`  
#> * `ppm_3.2081` -> `ppm_3.2081...1070`  
#> * `ppm_3.2320` -> `ppm_3.2320...1072`  
#> * `ppm_3.2382` -> `ppm_3.2382...1073`  
#> * `ppm_3.2672` -> `ppm_3.2672...1075`  
#> * `ppm_3.3930` -> `ppm_3.3930...1079`  
#> * `ppm_3.4031` -> `ppm_3.4031...1080`  
#> * `ppm_3.4185` -> `ppm_3.4185...1082`
```



## Introduction to AlpsNMR (older API)

```
#> * `ppm_3.4243` -> `ppm_3.4243...1083`  
#> * `ppm_3.4829` -> `ppm_3.4829...1090`  
#> * `ppm_3.4981` -> `ppm_3.4981...1092`  
#> * `ppm_3.5135` -> `ppm_3.5135...1093`  
#> * `ppm_3.5372` -> `ppm_3.5372...1095`  
#> * `ppm_3.5793` -> `ppm_3.5793...1099`  
#> * `ppm_3.6428` -> `ppm_3.6428...1104`  
#> * `ppm_3.6499` -> `ppm_3.6499...1105`  
#> * `ppm_3.7042` -> `ppm_3.7042...1106`  
#> * `ppm_3.7141` -> `ppm_3.7141...1107`  
#> * `ppm_3.7200` -> `ppm_3.7200...1108`  
#> * `ppm_3.7348` -> `ppm_3.7348...1110`  
#> * `ppm_3.7442` -> `ppm_3.7442...1111`  
#> * `ppm_3.7642` -> `ppm_3.7642...1113`  
#> * `ppm_3.8233` -> `ppm_3.8233...1117`  
#> * `ppm_3.8528` -> `ppm_3.8528...1121`  
#> * `ppm_3.8886` -> `ppm_3.8886...1124`  
#> * `ppm_3.9091` -> `ppm_3.9091...1126`  
#> * `ppm_3.9333` -> `ppm_3.9333...1128`  
#> * `ppm_3.9493` -> `ppm_3.9493...1129`  
#> * `ppm_4.0108` -> `ppm_4.0108...1136`  
#> * `ppm_4.1205` -> `ppm_4.1205...1141`  
#> * `ppm_4.1320` -> `ppm_4.1320...1142`  
#> * `ppm_4.1927` -> `ppm_4.1927...1145`  
#> * `ppm_4.2281` -> `ppm_4.2281...1148`  
#> * `ppm_4.2566` -> `ppm_4.2566...1154`  
#> * `ppm_4.4376` -> `ppm_4.4376...1166`  
#> * `ppm_4.4531` -> `ppm_4.4531...1168`  
#> * `ppm_4.5089` -> `ppm_4.5089...1171`  
#> * `ppm_4.5092` -> `ppm_4.5092...1172`  
#> * `ppm_4.5218` -> `ppm_4.5218...1173`  
#> * `ppm_4.5906` -> `ppm_4.5906...1178`  
#> * `ppm_5.0973` -> `ppm_5.0973...1184`  
#> * `ppm_5.2369` -> `ppm_5.2369...1187`  
#> * `ppm_5.4220` -> `ppm_5.4220...1195`  
#> * `ppm_6.9009` -> `ppm_6.9009...1225`  
#> * `ppm_6.9154` -> `ppm_6.9154...1226`  
#> * `ppm_6.9626` -> `ppm_6.9626...1227`  
#> * `ppm_7.2066` -> `ppm_7.2066...1234`  
#> * `ppm_7.2894` -> `ppm_7.2894...1237`  
#> * `ppm_7.3409` -> `ppm_7.3409...1241`  
#> * `ppm_7.3816` -> `ppm_7.3816...1243`  
#> * `ppm_7.5415` -> `ppm_7.5415...1250`  
#> * `ppm_7.5553` -> `ppm_7.5553...1251`  
#> * `ppm_8.3515` -> `ppm_8.3515...1273`  
#> * `ppm_8.4608` -> `ppm_8.4608...1274`
```

```
peak_table_integration = get_integration_with_metadata(peak_table_integration)
```

We can also integrate with a specific peak position and some arbitrary width:

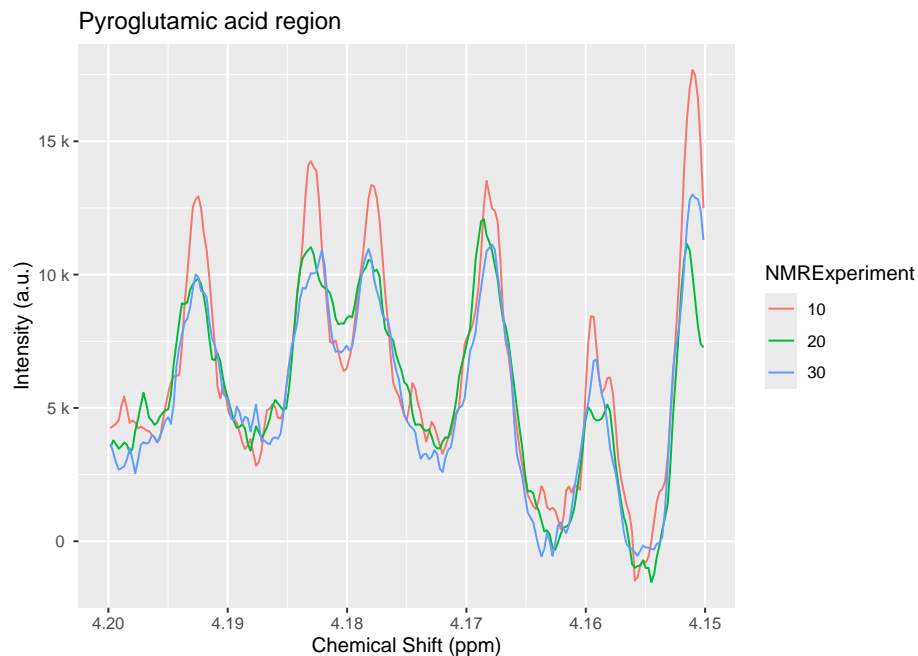
## Introduction to AlpsNMR (older API)

```
nmr_data(  
  nmr_integrate_peak_positions(samples = dataset_norm,  
                               peak_pos_ppm = c(4.1925, 4.183, 4.1775, 4.17),  
                               peak_width_ppm = 0.006)  
)  
#>      ppm_4.1925 ppm_4.1830 ppm_4.1775 ppm_4.1700  
#> 10      47.88901      53.53241      50.49386      45.62978  
#> 20      44.18572      51.07377      49.52798      44.53178  
#> 30      42.39196      47.99196      46.34812      39.07835
```

## 14.2 2. Integration based on peak boundaries

Imagine we only want to integrate the four peaks corresponding to the pyroglutamic acid:

```
pyroglutamic_acid_region <- c(4.15, 4.20)  
plot(dataset_norm, chemshift_range = pyroglutamic_acid_region) +  
  ggplot2::ggtitle("Pyroglutamic acid region")
```



We define the peak regions and integrate them. Note how we can correct the baseline or not. If we correct the baseline, the limits of the integration will be connected with a straight line and that line will be used as the baseline, that will be subtracted.

```
pyroglutamic_acid <- list(pyroglutamic_acid1 = c(4.19, 4.195),  
                           pyroglutamic_acid2 = c(4.18, 4.186),  
                           pyroglutamic_acid3 = c(4.175, 4.18),  
                           pyroglutamic_acid4 = c(4.165, 4.172))  
regions_base_corr_ds <- nmr_integrate_regions(dataset_norm, pyroglutamic_acid, fix_baseline = TRUE)  
regions_base_corr_matrix <- nmr_data(regions_base_corr_ds)  
regions_base_corr_matrix  
#>      pyroglutamic_acid1 pyroglutamic_acid2 pyroglutamic_acid3 pyroglutamic_acid4
```

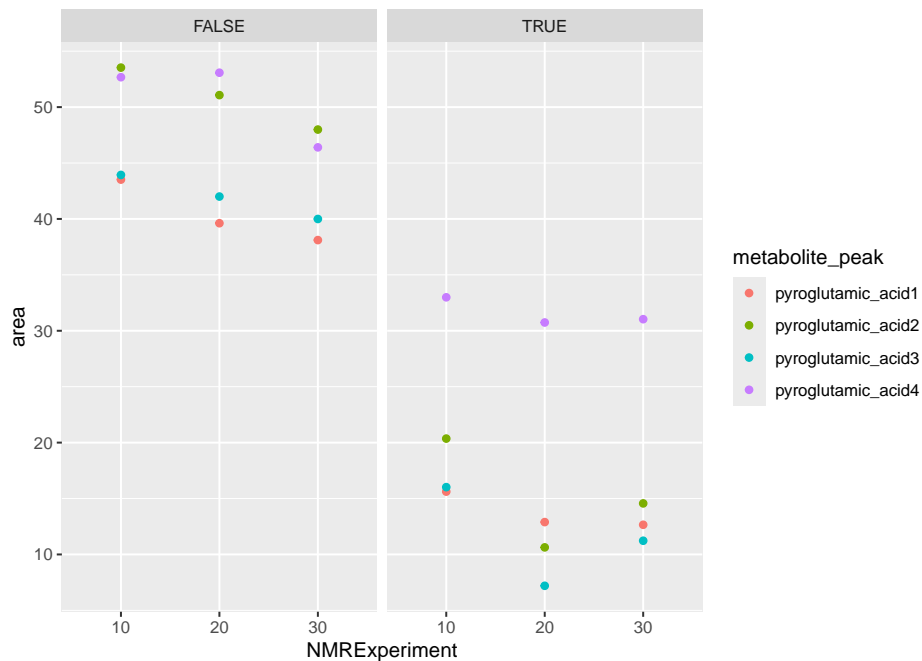
## Introduction to AlpsNMR (older API)

```
#> 10      15.61810      20.36002      16.018972      32.99148
#> 20      12.89344      10.63232       7.196728      30.74094
#> 30      12.65231      14.56388      11.226573      31.03655

regions_base_not_corr_ds <- nmr_integrate_regions(dataset_norm, pyroglutamic_acid, fix_baseline = FALSE)
regions_base_not_corr_matrix <- nmr_data(regions_base_not_corr_ds)
regions_base_not_corr_matrix
#>   pyroglutamic_acid1 pyroglutamic_acid2 pyroglutamic_acid3 pyroglutamic_acid4
#> 10      43.51763      53.53241      43.93533      52.67311
#> 20      39.61702      51.07377      42.00336      53.07528
#> 30      38.10972      47.99196      39.99623      46.40013
```

We may plot the integral values to explore variation based on the baseline subtraction.

```
dplyr::bind_rows(
  regions_base_corr_matrix %>%
    as.data.frame() %>%
    tibble::rownames_to_column("NMRExperiment") %>%
    tidyr::gather("metabolite_peak", "area", -NMRExperiment) %>%
    dplyr::mutate(BaselineCorrected = TRUE),
  regions_base_not_corr_matrix %>%
    as.data.frame() %>%
    tibble::rownames_to_column("NMRExperiment") %>%
    tidyr::gather("metabolite_peak", "area", -NMRExperiment) %>%
    dplyr::mutate(BaselineCorrected = FALSE)
) %>% ggplot() + geom_point(aes(x = NMRExperiment, y = area, color = metabolite_peak)) +
  facet_wrap(~BaselineCorrected)
```



## 15 Identification

After applying any feature selection or machine learning, Alps allows the identification of features of interest through `nmr_identify_regions_blood`. The function gives 3 possibilities sorted by the most probable metabolite (see `nmr_identify_regions_blood` for details).

```
ppm_to_assign <- c(4.060960203, 3.048970634, 2.405935596, 0.990616851, 0.986520147, 1.044258467)
identification <- nmr_identify_regions_blood(ppm_to_assign)
identification[!is.na(identification$Metabolite), ]
```

#>	Metabolite	HMDB_code	Shift_ppm	Type	J_Hz	Height
#> 1781	L-Valine	HMDB00883	0.991	d	7.01	1.0000
#> 1031	L-Alpha-aminobutyric acid	HMDB00452	0.997	t	7.58	1.0000
#> 178	L-Valine	HMDB00883	0.991	d	7.01	1.0000
#> 103	L-Alpha-aminobutyric acid	HMDB00452	0.997	t	7.58	1.0000
#> 179	L-Valine	HMDB00883	1.044	d	7.05	0.9614
#> 195	Pyroglutamic acid	HMDB00267	2.405	m		0.7466
#> 201	Succinic acid	HMDB00254	2.405	s		1.0000
#> 148	L-Lysine	HMDB00182	3.035	t		1.0000
#> 47	Creatinine	HMDB00562	3.045	s		1.0000
#> 44	Creatine	HMDB00064	3.035	s		1.0000
#> 46	Creatinine	HMDB00562	4.065	s		0.4374
#> 182	Myoinositol	HMDB00211	4.068	t	2.839	NA
#> 40	Choline	HMDB00097	4.071	ddd		0.0607

#>	Blood_concentration	n_reported_in_Blood	ppm_to_assign
#> 1781	179.04615	13	0.9865201
#> 1031	22.80000	6	0.9865201
#> 178	179.04615	13	0.9906169
#> 103	22.80000	6	0.9906169
#> 179	179.04615	13	1.0442585
#> 195	19.50000	2	2.4059356
#> 201	16.10000	8	2.4059356
#> 148	168.20000	14	3.0489706
#> 47	51.25167	14	3.0489706
#> 44	48.47000	6	3.0489706
#> 46	51.25167	14	4.0609602
#> 182	23.52500	5	4.0609602
#> 40	13.08000	5	4.0609602

## 16 Final thoughts

This vignette shows many of the features of the package, some features have room for improvement, others are not fully described, and the reader will need to browse the documentation. Hopefully it is a good starting point for using the package.

```
sessionInfo()
#> R version 4.4.0 RC (2024-04-16 r86468 ucrt)
#> Platform: x86_64-w64-mingw32/x64
#> Running under: Windows Server 2022 x64 (build 20348)
#>
#> Matrix products: default
```

## Introduction to AlpsNMR (older API)

```
#>
#>
#> locale:
#> [1] LC_COLLATE=C
#> [2] LC_CTYPE=English_United States.utf8
#> [3] LC_MONETARY=English_United States.utf8
#> [4] LC_NUMERIC=C
#> [5] LC_TIME=English_United States.utf8
#>
#> time zone: America/New_York
#> tzcode source: internal
#>
#> attached base packages:
#> [1] stats      graphics  grDevices  utils      datasets  methods   base
#>
#> other attached packages:
#> [1] AlpsNMR_4.7.0      future_1.33.2      BiocParallel_1.39.0
#> [4] readxl_1.4.3       ggplot2_3.5.1      dplyr_1.1.4
#> [7] BiocStyle_2.33.0
#>
#> loaded via a namespace (and not attached):
#> [1] baseline_1.3-5      gridExtra_2.3       rlang_1.1.3
#> [4] magrittr_2.0.3      MassSpecWavelet_1.71.0 matrixStats_1.3.0
#> [7] compiler_4.4.0      vctrs_0.6.5         reshape2_1.4.4
#> [10] RcppZigurat_0.1.6   quadprog_1.5-8      rvest_1.0.4
#> [13] stringr_1.5.1       pkgconfig_2.0.3     crayon_1.5.2
#> [16] fastmap_1.1.1       labeling_0.4.3      utf8_1.2.4
#> [19] promises_1.3.0      rmarkdown_2.26      ps_1.7.6
#> [22] itertools_0.1-3     tinytex_0.50        purrr_1.0.2
#> [25] xfun_0.43           Rfast_2.1.0         randomForest_4.7-1.1
#> [28] jsonlite_1.8.8      limSolve_1.5.7.1    later_1.3.2
#> [31] parallel_4.4.0      cluster_2.1.6       R6_2.5.1
#> [34] stringi_1.8.3       RColorBrewer_1.1-3  parallelly_1.37.1
#> [37] cellranger_1.1.0    Rcpp_1.0.12         bookdown_0.39
#> [40] iterators_1.0.14    knitr_1.46          snow_0.4-4
#> [43] Matrix_1.7-0        igraph_2.0.3        tidyselect_1.2.1
#> [46] yaml_2.3.8          websocket_1.4.1      codetools_0.2-20
#> [49] processx_3.8.4      listenv_0.9.1       doRNG_1.8.6
#> [52] lattice_0.22-6      tibble_3.2.1        plyr_1.8.9
#> [55] withr_3.0.0         rARPACK_0.11-0      evaluate_0.23
#> [58] signal_1.8-0        speaq_2.7.0         RcppParallel_5.1.7
#> [61] xml2_1.3.6          lpSolve_5.6.20      pillar_1.9.0
#> [64] BiocManager_1.30.22 rngtools_1.5.2      foreach_1.5.2
#> [67] ellipse_0.5.0       pcaPP_2.0-4         generics_0.1.3
#> [70] chromote_0.2.0      munsell_0.5.1       scales_1.3.0
#> [73] globals_0.16.3     glue_1.7.0          tools_4.4.0
#> [76] data.table_1.15.4   SparseM_1.81        RSpectra_0.16-1
#> [79] fs_1.6.4            mvtnorm_1.2-4       cowplot_1.1.3
#> [82] grid_4.4.0          impute_1.79.0       missForest_1.5
#> [85] tidyr_1.3.1         colorspace_2.1-0    cli_3.6.2
#> [88] fansi_1.0.6         mixOmics_6.29.0     corpcor_1.6.10
```

## Introduction to AlpsNMR (older API)

```
#> [91] doSNOW_1.0.20      gtable_0.3.5      digest_0.6.35  
#> [94] progressr_0.14.0   ggrepel_0.9.5     farver_2.1.1  
#> [97] htmltools_0.5.8.1 lifecycle_1.0.4    httr_1.4.7  
#> [100] MASS_7.3-60.2
```