4: Part 2 - Data Wrangling

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Objectives

- 1. Describe the usefulness of data wrangling and its place in the data pipeline
- 2. Wrangle datasets with dplyr functions
- 3. Apply data wrangling skills to a real-world example dataset

Set up your session

```
getwd()
## [1] "/home/guest/EDE_Fall2024"

library(plyr)
library(tidyverse)
library(lubridate)
NTL.phys.data.PeterPaul <- read.csv("./Data/Processed/NTL-LTER_Lake_ChemistryPhysics_PeterPaul_Processed/NTL.nutrient.data <- read.csv("./Data/Raw/NTL-LTER_Lake_Nutrients_Raw.csv",stringsAsFactors = TRUE)</pre>
```

Review of basic exploration and wrangling

```
# Data summaries for physical data
colnames(NTL.phys.data.PeterPaul)
   [1] "lakeid"
                       "lakename"
                                       "year4"
                                                       "daynum"
   [5] "month"
                       "sampledate"
                                       "depth"
##
                                                       "temperature_C"
   [9] "dissolvedOxygen" "irradianceWater" "irradianceDeck"
                                                       "comments"
dim(NTL.phys.data.PeterPaul)
## [1] 21613
str(NTL.phys.data.PeterPaul)
## 'data.frame':
                 21613 obs. of 12 variables:
## $ lakeid
                 : chr "L" "L" "L" "L" ...
## $ lakename
                  : chr "Paul Lake" "Paul Lake" "Paul Lake" "Paul Lake" ...
                  ## $ year4
```

```
## $ daynum
                  : int 148 148 148 148 148 148 148 148 148 1...
## $ month
                  : int 5555555555...
                  : chr "1984-05-27" "1984-05-27" "1984-05-27" "1984-05-27" ...
## $ sampledate
## $ depth
                  : num 0 0.25 0.5 0.75 1 1.5 2 3 4 5 ...
## $ temperature_C : num 14.5 NA NA NA 14.5 NA 14.2 11 7 6.1 ...
## $ dissolvedOxygen: num 9.5 NA NA NA 8.8 NA 8.6 11.5 11.9 2.5 ...
## $ irradianceWater: num 1750 1550 1150 975 870 610 420 220 100 34 ...
## $ comments
                  : chr NA NA NA NA ...
summary(NTL.phys.data.PeterPaul$comments)
##
     Length
               Class
                         Mode
##
      21613 character character
class(NTL.phys.data.PeterPaul$sampledate)
## [1] "character"
# Format sampledate as date
NTL.phys.data.PeterPaul$sampledate <- as.Date(NTL.phys.data.PeterPaul$sampledate, format = "%Y-%m-%d")
# Select Peter and Paul Lakes from the nutrient dataset
NTL.nutrient.data.PeterPaul <- filter(NTL.nutrient.data, lakename == "Paul Lake" | lakename == "Peter L
# Data summaries for nutrient data
colnames(NTL.nutrient.data.PeterPaul)
  [1] "lakeid"
                                          "daynum"
                                                      "sampledate"
                   "lakename"
                              "year4"
                                                      "nh34"
## [6] "depth id"
                   "depth"
                              "tn ug"
                                          "tp_ug"
## [11] "no23"
                  "po4"
                              "comments"
dim(NTL.nutrient.data.PeterPaul)
## [1] 2770
            13
str(NTL.nutrient.data.PeterPaul)
## 'data.frame':
                 2770 obs. of 13 variables:
## $ lakeid
            : Factor w/ 26 levels "B", "Berg", "Bolg", ...: 13 13 13 13 13 18 18 18 18 ...
\#\# $ lakename : Factor w/ 26 levels "Bergner Lake",..: 16 16 16 16 16 16 17 17 17 17 ...
             ## $ year4
              ## $ sampledate: Factor w/ 1294 levels "4/14/12","4/19/12",..: 31 31 31 31 31 31 31 31 31 31 ...
## $ depth_id : int 1 2 3 4 5 6 1 2 3 4 ...
              : num 0 0.85 1.75 3 4 6 0 1 2.25 3.5 ...
## $ depth
## $ tn_ug
              : num 538 285 399 453 363 583 352 356 364 582 ...
              : num 25 14 14 14 13 37 11 15 28 14 ...
## $ tp_ug
## $ nh34
              : num NA NA NA NA NA NA NA NA NA ...
              : num NA NA NA NA NA NA NA NA NA ...
## $ no23
              : num NA NA NA NA NA NA NA NA NA ...
## $ po4
## $ comments : Factor w/ 3 levels "", "sample missing", ..: 1 1 1 1 1 1 1 1 1 1 ...
```

```
summary(NTL.nutrient.data.PeterPaul$lakename)
##
         Bergner Lake
                                  Bog Pot
                                                   Bolger Bog
                                                                      Brown Lake
##
##
    Central Long Lake
                            Crampton Lake
                                               Cranberry Bog
                                                                  East Long Lake
##
                                                                                0
##
              Eds Bog Forest Service Bog
                                            Hummingbird Lake
                                                                     Inkpot Lake
##
                                                                                0
                    0
                                        0
##
                              Morris Lake
                                                                       Paul Lake
        Kickapoo Lake
                                              North Gate Bog
##
                                                                            1383
                                                            0
           Peter Lake
##
                                Plum Lake
                                              Raspberry Lake
                                                                 Reddington Lake
                 1387
##
                               Tender Bog
##
           Roach Lake
                                             Tenderfoot Lake
                                                                    Tuesday Lake
##
                    0
                                        0
                                                                                0
##
            Ward Lake
                           West Long Lake
##
                    0
# Notice that other lake names didn't go away, even though they have zero values
NTL.nutrient.data.PeterPaul <- droplevels(NTL.nutrient.data.PeterPaul)
summary(NTL.nutrient.data.PeterPaul$lakename)
    Paul Lake Peter Lake
##
         1383
                    1387
summary(NTL.nutrient.data.PeterPaul$comments)
##
## 2770
class(NTL.nutrient.data.PeterPaul$sampledate)
## [1] "factor"
NTL.nutrient.data.PeterPaul$sampledate <- as.Date(NTL.nutrient.data.PeterPaul$sampledate, format = "%m/
NTL.nutrient.data.PeterPaul <-
  NTL.nutrient.data.PeterPaul %>% #
  mutate(month = month(sampledate)) %>% #
  select(lakeid:daynum, month, sampledate:comments) %>% #
  drop_na(depth)
# Save processed nutrient file
write.csv(NTL.nutrient.data.PeterPaul, row.names = FALSE,
          file = "./Data/Processed/NTL-LTER_Lake_Nutrients_PeterPaul_Processed.csv")
# Remove columns that are not of interest for analysis
NTL.phys.data.PeterPaul.subset <- select(NTL.phys.data.PeterPaul,</pre>
                                          lakename:irradianceDeck)
NTL.nutrient.data.PeterPaul.subset <- select(NTL.nutrient.data.PeterPaul,
```

```
lakename, year4, daynum, month, sampledate, depth:po4)

# write a more succinct line of code to subset the nutrient dataset.
```

Gather and Spread

Gather nutrient data into one column

For most situations, data analysis works best when you have organized your data into a tidy dataset. A tidy dataset is defined as: * Each variable is a column * Each row is an observation * Each value is in its own cell

However, there may be situations where we want to reshape our dataset, for example if we want to facet numerical data points by measurement type (more on this in the data visualization unit). We can program this reshaping in a few short lines of code using the package tidyr, which is conveniently included in the tidyverse package.

We will work with the new functions from tidyr. For gather we will use pivot_longer and for spread we will use pivot_wider. Note that if you are familiar with gather and spread will may still use it, they are not going away, but they are not under active development.

```
#NTL.nutrient.data.PeterPaul.gathered <- gather(NTL.nutrient.data.PeterPaul.subset, "nutrient", "concen
# Gather nutrient data into one column using pivot_longer
NTL.nutrient.data.PeterPaul.gathered <- pivot_longer(NTL.nutrient.data.PeterPaul.subset,tn_ug:po4, name
#subset function does something similar to filtering, takes original data, replaces it with new data
NTL.nutrient.data.PeterPaul.gathered <- subset(NTL.nutrient.data.PeterPaul.gathered, !is.na(concentrati
count(NTL.nutrient.data.PeterPaul.gathered, nutrient)
## # A tibble: 5 x 2
##
    nutrient
##
     <chr>>
              <int>
## 1 nh34
              1130
## 2 no23
               1161
## 3 po4
               1186
## 4 tn_ug
               1425
## 5 tp_ug
               2279
write.csv(NTL.nutrient.data.PeterPaul.gathered, row.names = FALSE,
          file = "./Data/Processed/NTL-LTER_Lake_Nutrients_PeterPaulGathered_Processed.csv")
# Spread nutrient data into separate columns
#NTL.nutrient.data.PeterPaul.spread <- spread(NTL.nutrient.data.PeterPaul.gathered, nutrient, concentra
# Spread nutrient data into separate columns using pivot_wider
NTL.nutrient.data.PeterPaul.spread2 <- pivot_wider(NTL.nutrient.data.PeterPaul.gathered, names_from = n
# Split components of cells into multiple columns
# Opposite of 'separate' is 'unite'
NTL.nutrient.data.PeterPaul.dates <- separate(NTL.nutrient.data.PeterPaul.subset, sampledate, c("Y", "m
# I recommend using lubridate rather than separate and unite.
```

Combining multiple datasets

Join

In many cases, we will want to combine datasets into one dataset. If all column names match, the data frames can be combined with the rbind function. If some column names match and some column names don't match, we can combine the data frames using a "join" function according to common conditions that exist in the matching columns. We will demonstrate this with the NTL-LTER physical and nutrient datasets, where we have specific instances when physical and nutrient data were collected on the same date, at the same lake, and at the same depth.

In dplyr, there are several types of join functions:

- inner_join: return rows in x where there are matching values in y, and all columns in x and y (mutating join).
- semi_join: return all rows from x where there are matching values in y, keeping just columns from x (filtering join).
- left_join: return all rows from x, and all columns from x and y (mutating join).
- anti_join: return all rows from x where there are *not* matching values in y, keeping just columns from x (filtering join).
- full_join: return all rows and all columns from x and y. Returns NA for missing values (mutating join).

Let's say we want to generate a new dataset that contains all possible physical and chemical data for Peter and Paul Lakes. In this case, we want to do a full join.

```
NTL.phys.nutrient.data.PeterPaul <- full_join(NTL.phys.data.PeterPaul.subset, NTL.nutrient.data.PeterPaul
## Joining with 'by = join_by(lakename, year4, daynum, month, sampledate, depth)'
write.csv(NTL.phys.nutrient.data.PeterPaul, row.names = FALSE,</pre>
```

file ="./Data/Processed/NTL-LTER Lake Chemistry Nutrients PeterPaul Processed.csv")

rbind

Rbind stands for row bind. Combining datasets by row

The Niwot Ridge litter dataset, when downloaded from NEON, comes packaged with each month as a different .csv file. If we want to analyze the dataset as a single data frame, we need to combine each of these files.

```
Litter.June2016 <- read.csv("./Data/Raw/NIWO_Litter/NEON_NIWO_Litter_massdata_2016-06_raw.csv")
Litter.July2016 <- read.csv("./Data/Raw/NIWO_Litter/NEON_NIWO_Litter_massdata_2016-07_raw.csv")
Litter.August2016 <- read.csv("./Data/Raw/NIWO_Litter/NEON_NIWO_Litter_massdata_2016-08_raw.csv")
Litter.2019 <- rbind(Litter.June2016, Litter.July2016, Litter.August2016)
```

However, there are 20 months in this dataset, so importing all these files individually would be tedious to code. Here is a more efficient way to import and combine all files.

```
LitterFiles
    [1] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2016-06_raw.csv"
##
##
    [2] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2016-07_raw.csv"
   [3] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2016-08_raw.csv"
##
   [4] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2016-09_raw.csv"
##
   [5] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2016-10_raw.csv"
   [6] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2016-11_raw.csv"
##
##
   [7] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2017-07_raw.csv"
   [8] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2017-08_raw.csv"
##
##
  [9] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2017-10_raw.csv"
## [10] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2017-11_raw.csv"
## [11] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2018-06_raw.csv"
## [12] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2018-07_raw.csv"
## [13] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2018-08_raw.csv"
## [14] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2018-09_raw.csv"
## [15] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2018-10_raw.csv"
## [16] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2018-11_raw.csv"
## [17] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2019-06_raw.csv"
## [18] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2019-07_raw.csv"
## [19] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2019-08_raw.csv"
## [20] "./Data/Raw/NIWO_Litter//NEON_NIWO_Litter_massdata_2019-09_raw.csv"
#use a pipe, and for this object LitterFiles, read each element in file using read.csv file If we do th
Litter <- LitterFiles %>%
  plyr::ldply(read.csv)
```

LitterFiles = list.files(path = "./Data/Raw/NIWO_Litter/", pattern="*.csv", full.names=TRUE)

We also have information about individual traps, including the location and type of landcover. Let's join these two datasets. Note that "siteID", "plotID" and "trapID" exist in both datasets, and we can join them by these conditions. Notice the dimensions of the final dataset.

Split-Apply-Combine

dplyr functionality, combined with the pipes operator, allows us to split datasets according to groupings (function: group_by), then run operations on those groupings and return the output of those operations. There is a lot of flexibility in this approach, but we will illustrate just one example today.

```
NTL.PeterPaul.summaries <-
   NTL.phys.nutrient.data.PeterPaul %>%
   filter(depth == 0) %>%
   group_by(lakename, month) %>%
   filter(!is.na(temperature_C) & !is.na(tn_ug) & !is.na(tp_ug)) %>%
   summarise(meantemp = mean(temperature_C),
        sdtemp = sd(temperature_C),
        meanTN = mean(tn_ug),
        sdTN = sd(tn_ug),
        meanTP = mean(tp_ug),
        sdTP = sd(tp_ug))

## 'summarise()' has grouped output by 'lakename'. You can override using the
## '.groups' argument.

write.csv(NTL.PeterPaul.summaries, row.names = FALSE,
        file ="./Data/Processed/NTL-LTER_Lake_Summaries_PeterPaul_Processed.csv")
```

Alternative Methods for Data Wrangling

If you want to iteratively perform operations on your data, there exist several options. We have demonstrated the pipe as one option. Additional options include the apply function (https://www.rdocumentation.org/packages/base/versions/3.5.2/topics/apply) and for loops (https://swcarpentry.github.io/r-novice-inflammation/15-supp-loops-in-depth/). These options are good options as well (again, multiple ways to get to the same outcome). A word of caution: loops are slow. This may not make a difference for small datasets, but small time additions will make a difference with large datasets.