Assignment 5: Data Visualization

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OVERVIEW

This exercise accompanies the lessons in Environmental Data Analytics (ENV872L) on data wrangling.

Directions

- 1. Change "Student Name" on line 3 (above) with your name.
- 2. Use the lesson as a guide. It contains code that can be modified to complete the assignment.
- 3. Work through the steps, **creating code and output** that fulfill each instruction.
- 4. Be sure to **answer the questions** in this assignment document. Space for your answers is provided in this document and is indicated by the ">" character. If you need a second paragraph be sure to start the first line with ">". You should notice that the answer is highlighted in green by RStudio.
- 5. When you have completed the assignment, **Knit** the text and code into a single PDF file. You will need to have the correct software installed to do this (see Software Installation Guide) Press the **Knit** button in the RStudio scripting panel. This will save the PDF output in your Assignments folder.
- 6. After Knitting, please submit the completed exercise (PDF file) to the dropbox in Sakai. Please add your last name into the file name (e.g., "Salk_A04_DataWrangling.pdf") prior to submission.

The completed exercise is due on Thursday, 14 February, 2019 before class begins.

Set up your session

- Set up your session. Upload the NTL-LTER processed data files for chemistry/physics for Peter and Paul Lakes (tidy and gathered), the USGS stream gauge dataset, and the EPA Ecotox dataset for Neonicotinoids.
- 2. Make sure R is reading dates as date format, not something else (hint: remember that dates were an issue for the USGS gauge data).

```
#1
getwd()
## [1] "Y:/19spring/872/Environmental_Data_Analytics/Assignments"
setwd("Y:/19spring/872/Environmental_Data_Analytics/Assignments")
PPchemphy <- read.csv("../Data/Processed/NTL-LTER_Lake_Chemistry_Nutrients_PeterPaul_Processed.csv")
#Sorry, I didn't upload the gathered one.
USGSgage <- read.csv("../Data/Raw/USGS_Site02085000_Flow_Raw.csv")</pre>
ecotox <- read.csv(".../Data/Raw/ECOTOX_Neonicotinoids_Mortality_raw.csv")</pre>
library(tidyverse)
## Warning: package 'tidyverse' was built under R version 3.5.2
## -- Attaching packages ------
## v ggplot2 3.1.0
                               0.2.5
                      v purrr
## v tibble 1.4.2
                      v dplyr
                               0.7.8
## v tidyr
            0.8.2
                      v stringr 1.3.1
## v readr
            1.3.1
                      v forcats 0.3.0
## Warning: package 'ggplot2' was built under R version 3.5.2
## Warning: package 'tibble' was built under R version 3.5.2
```

```
## Warning: package 'tidyr' was built under R version 3.5.2
## Warning: package 'readr' was built under R version 3.5.2
## Warning: package 'purrr' was built under R version 3.5.2
## Warning: package 'dplyr' was built under R version 3.5.2
## Warning: package 'stringr' was built under R version 3.5.2
## Warning: package 'forcats' was built under R version 3.5.2
## -- Conflicts -----
## x dplyr::filter() masks stats::filter()
## x dplyr::lag()
                     masks stats::lag()
#2.
USGSgage$datetime <- as.Date(USGSgage$datetime, "%m/%d/%y")
USGSgage$datetime <- format(USGSgage$datetime,"%y%m%d")</pre>
creat.early.dates <- function(d) paste0(ifelse(d > 181231, "19", "20"),d)
USGSgage$datetime <- creat.early.dates(USGSgage$datetime)</pre>
USGSgage$datetime <- as.Date(USGSgage$datetime,"%Y%m%d")
PPchemphy$sampledate <- as.Date(PPchemphy$sampledate,"%Y-%m-%d")
```

Define your theme

3. Build a theme and set it as your default theme.

Create graphs

For numbers 4-7, create graphs that follow best practices for data visualization. To make your graphs "pretty," ensure your theme, color palettes, axes, and legends are edited to your liking.

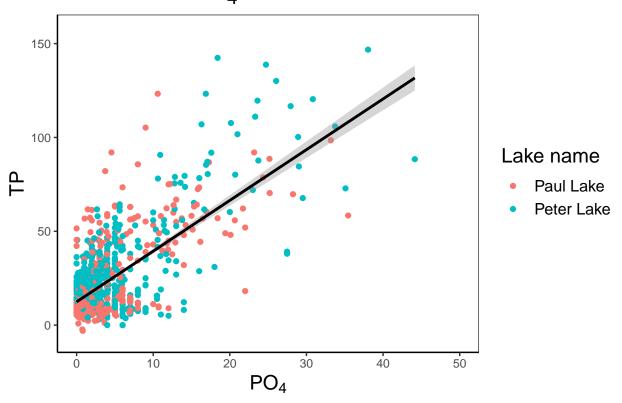
Hint: a good way to build graphs is to make them ugly first and then create more code to make them pretty.

4. [NTL-LTER] Plot total phosphorus by phosphate, with separate aesthetics for Peter and Paul lakes. Add a line of best fit and color it black.

```
#4
ggplot(PPchemphy,aes(x=po4,y=tp_ug,color=lakename)) +
    geom_point(size=1.5) +
    wang +
    xlim(c(0,50)) +
    labs(title = expression("TP-PO"[4]*"concentration"),x=expression("PO"[4]),y="TP",color="Lake name") +
    geom_smooth(aes(x=po4,y=tp_ug),method = lm,color="black")
## Warning: Removed 22310 rows containing non-finite values (stat_smooth).
```

Warning: Removed 22310 rows containing missing values (geom_point).

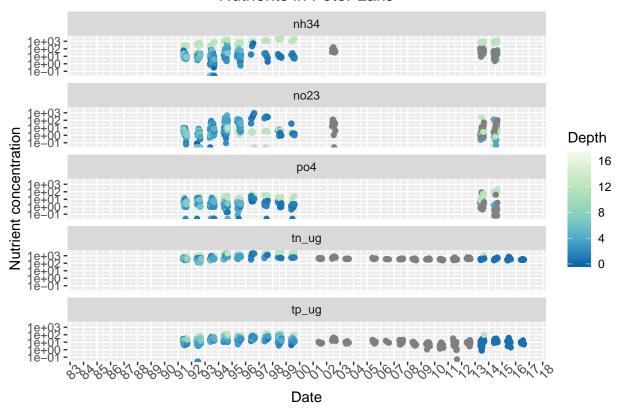
TP-PO₄concentration



5. [NTL-LTER] Plot nutrients by date for Peter Lake, with separate colors for each depth. Facet your graph by the nutrient type.

```
peter <- filter(PPchemphy,lakename=="Peter Lake")</pre>
## Warning: package 'bindrcpp' was built under R version 3.5.2
NTL.gathered <- gather(peter, "nutrient", "concentration", tn_ug:po4)
NTL.gathered$sampledate <- as.Date(NTL.gathered$sampledate,"%Y-%m-%d")
library(RColorBrewer)
## Warning: package 'RColorBrewer' was built under R version 3.5.2
ggplot(NTL.gathered,aes(x=sampledate,y=concentration,color=depth)) +
  geom_point(size=1.5) +
  facet_wrap(vars(nutrient),nrow = 5) +
  scale_color_distiller("Depth",palette = "GnBu",direction = -1) +
  scale_x_date(name = "Date",date_breaks = "1 year",date_labels = "%y") +
  scale_y_log10("Nutrient concentration") +
  labs(title = "Nutrients in Peter Lake") +
  theme(axis.text.x = element_text(size=10, angle = 45), plot.title=element_text(hjust = 0.5))
## Warning in self$trans$transform(x): NaNs produced
## Warning: Transformation introduced infinite values in continuous y-axis
## Warning: Removed 56671 rows containing missing values (geom point).
```

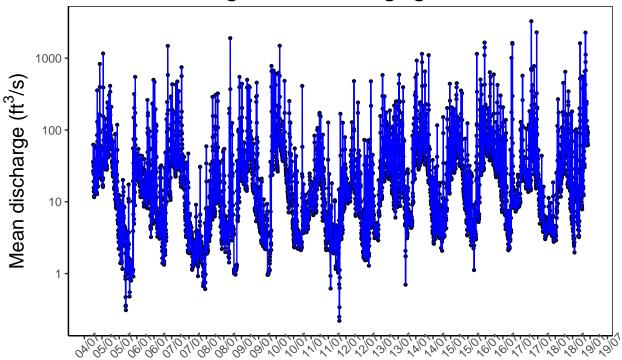
Nutrients in Peter Lake



6. [USGS gauge] Plot discharge by date. Create two plots, one with the points connected with geom_line and one with the points connected with geom_smooth (hint: do not use method = "lm").

```
#6
#I'm going to plot the mean discharge (84936_00060_00003).
discharge <- filter(USGSgage,X84936_00060_00003>0)
ggplot(discharge,aes(x=datetime,y=X84936_00060_00003)) +
    geom_point(size=0.8,na.rm = T) +
    geom_line(color="Blue") +
    scale_x_date(name = "Date",date_breaks = "6 months",date_labels = "%y/%m") +
    scale_y_log10(expression("Mean discharge (ft"^"3"*"/s)")) +
    labs(title = "Mean discharge for USGS gage site 02085000") +
    wang +
    theme(axis.text.x = element_text(angle = 45))
```

Mean discharge for USGS gage site 02085000

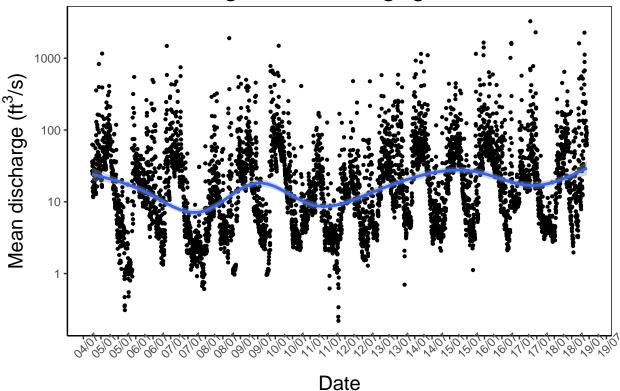


Date

```
ggplot(discharge,aes(x=datetime,y=X84936_00060_00003)) +
geom_point(size=0.8,na.rm = T) +
geom_smooth() +
scale_x_date("Date",date_breaks = "6 months",date_labels = "%y/%m") +
scale_y_log10(expression("Mean discharge (ft"^"3"*"/s)")) +
labs(title = "Mean discharge for USGS gage site 02085000") +
wang +
theme(axis.text.x = element_text(angle = 45))
```

$geom_smooth()$ using method = 'gam' and formula 'y ~ s(x, bs = "cs")'

Mean discharge for USGS gage site 02085000

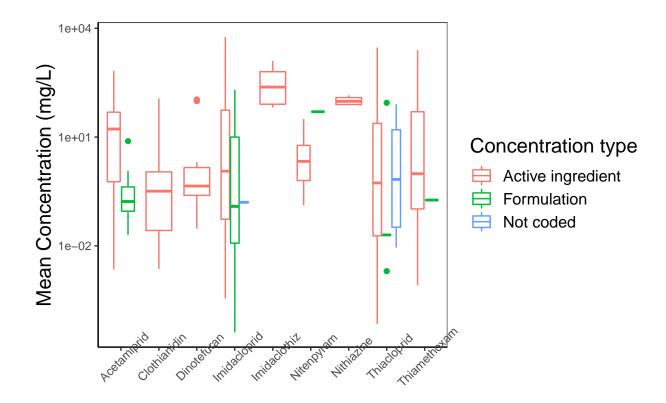


Question: How do these two types of lines affect your interpretation of the data?

Answer: The former one (geom_line) reflects the finest daily variation without any smoothing, while the latter (geom_smooth) shows the overall trend over time on a yearly temporal scale.

7. [ECOTOX Neonicotinoids] Plot the concentration, divided by chemical name. Choose a geom that accurately portrays the distribution of data points.

```
#7
attach(ecotox)
ecotox <- filter(ecotox,Conc..Units..Std.=="AI mg/L")
ecotox <- select(ecotox,Chemical.Name,Dur..Std.:Conc..Mean..Std.)
ggplot(ecotox,aes(x=Chemical.Name,y=Conc..Mean..Std.,color=Conc..Type)) +
    geom_boxplot() +
    scale_y_log10("Mean Concentration (mg/L)") + xlab("Chemical name") +
    wang +
    scale_color_discrete("Concentration type") +
    theme(axis.text.x = element_text(angle = 45))</pre>
```



Chemical name