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 Tuesday, 19 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] "/Users/yifeizhang/R/Environmental Data Analytics"
library(tidyverse)
## -- Attaching packages -
## v ggplot2 3.1.0
                             0.2.5
## v tibble 2.0.1
                             0.7.8
                    v dplyr
## v tidyr
           0.8.2
                    v stringr 1.3.1
## v readr
           1.3.1
                    v forcats 0.3.0
## -- Conflicts ------
## x dplyr::filter() masks stats::filter()
## x dplyr::lag()
                  masks stats::lag()
```

```
library(colormap)
PeterPaul <- read.csv("./Data/Processed/NTL-LTER_Lake_Nutrients_PeterPaul_Processed.csv")
PeterPaul_gathered <- read.csv("./Data/Processed/NTL-LTER_Lake_Nutrients_PeterPaulGathered_Processed.cs
USGS_Stream <- read.csv("./Data/Raw/USGS_Site02085000_Flow_Raw.csv")</pre>
```

Define your theme

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

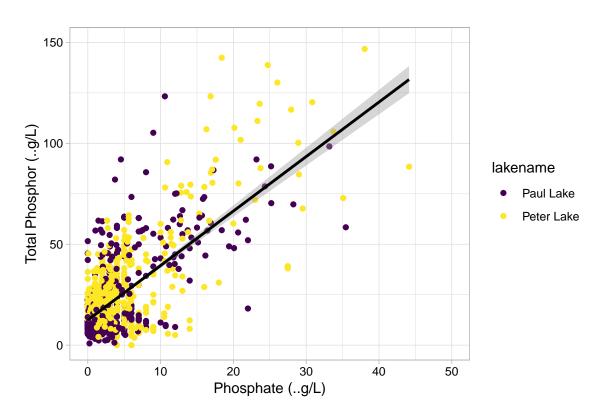
Create graphs

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.

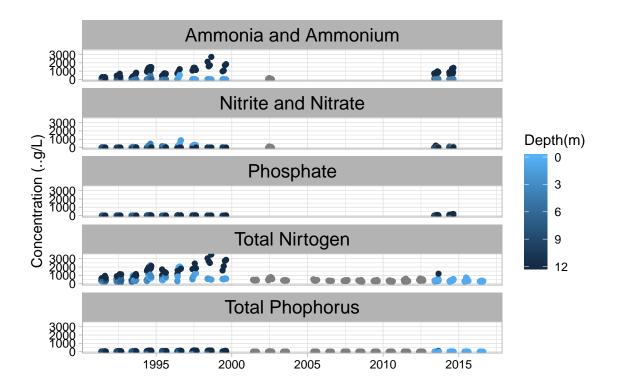
```
#4
plot4 <- ggplot(PeterPaul, aes(x = po4 , y= tp_ug, color = lakename) ) +
    geom_point() +
    geom_smooth(method = lm, color = "black") +
    Yifei_theme +
    labs(x = "Phosphate (\U0003BCg/L)" ,y = "Total Phosphor (\U0003BCg/L)") +
    xlim(0,50) +
    ylim(0,150)+
    scale_color_colormap(discrete = TRUE)</pre>
```



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

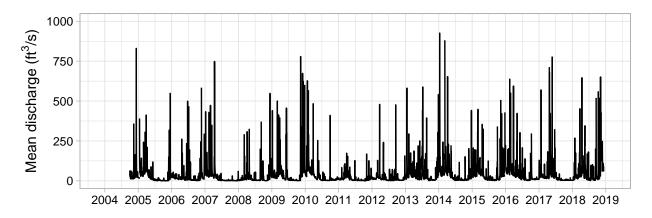
```
#5
plot5 <- ggplot(subset(PeterPaul_gathered, lakename == "Peter Lake"), aes(x = as.Date(sampledate), y =
    geom_point()+
    facet_wrap(vars(nutrient) , nrow = 5, labeller = as_labeller(c(`nh34` = "Ammonia and Ammonium" , `no23
    scale_color_continuous(trans = 'reverse')+
    labs(y = "Concentration (\U0003BCg/L)", x = "", color = "Depth(m)")+
    Yifei_theme

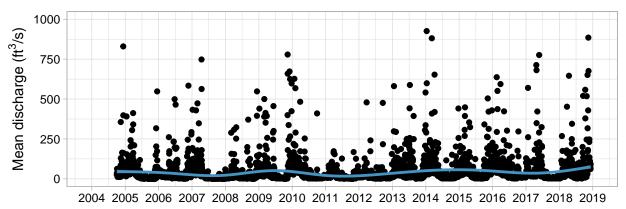
print(plot5)</pre>
```



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"). Place these graphs on the same plot (hint: ggarrange or something similar)

```
#6
#install.packages("ggpubr")
library(ggpubr)
plot6_1 <- ggplot(USGS_Stream, aes(y = discharge.mean, x = as.Date(datetime))) +</pre>
  geom_line()+
  scale_x_date(limits = as.Date(c("2004-01-01", "2018-12-31")),
 date_breaks = "1 year", date_labels = "%Y") +
 ylim(0,1000) +
  ylab(expression(paste("Mean discharge (ft"^3,"/s)"))) +
  xlab(expression(""))+
  Yifei_theme
#print(plot6_1)
plot6_2 <- ggplot(USGS_Stream, aes(y = discharge.mean, x = as.Date(datetime))) +
  geom_point()+
  geom_smooth(color= "#4393c3")+
  scale_x_date(limits = as.Date(c("2004-01-01", "2018-12-31")),
 date_breaks = "1 year", date_labels = "%Y") +
  ylim(0,1000) +
  ylab(expression(paste("Mean discharge (ft"^3,"/s)"))) +
  xlab(expression(""))+
  Yifei theme
#print(plot6_2)
ggarrange(plot6_1,plot6_2, nrow = 2)
```





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

Answer: The geom_line function just connects all the points together. The line shows the range of the discharge values. The geom_smooth function draws a line that shows the general trend of the discharge over the time.

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

```
#7
library(RColorBrewer)
plot7 <- ggplot(EPA_Ecotox, aes(x= Conc..Mean..Std.,color = Chemical.Name)) +
    geom_freqpoly() +
    Yifei_theme+
    scale_color_manual(values = c('#b2182b','#d6604d','#f4a582','#fddbc7','#f7f7f7','#d1e5f0','#92c5de','#
    labs(x="Concentration",y = "Count", color = "Chemical Name") +
        xlim(0,1000)

print(plot7)</pre>
```

