## exploratory

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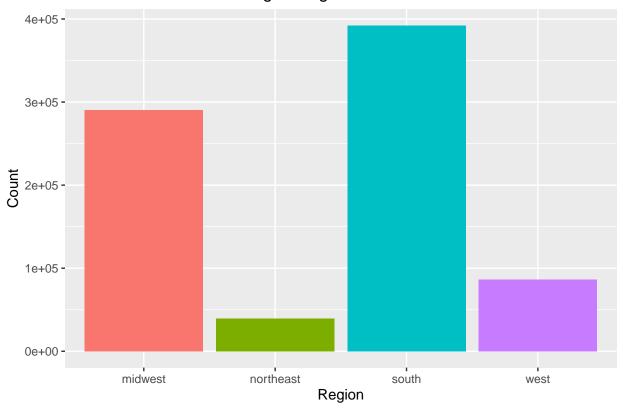
```
#Libraries
library(mosaic)
library(tidyverse)
library(usmap)
#importing in full dataset
import_df <- read_csv("data/chemical_data.csv")</pre>
## Parsed with column specification:
## cols(
##
     state = col_character(),
##
     site = col character(),
##
     disposal.area = col_character(),
##
     type = col_character(),
##
     well.id = col_character(),
##
     gradient = col_character(),
##
     samp.date = col_character(),
##
     contaminant = col_character(),
##
     measurement.unit = col_character(),
##
     below.detection = col_character(),
##
     concentration = col_double(),
##
     qualifier = col_character(),
     link = col_character()
##
#breaking apart into different datasets for each region
northeast <- import_df %>%
  filter(state %in% c("ME", "NH", "VT", "NY", "PA", "NJ", "MD",
                      "MA", "DE", "RI", "CT")) %>%
  mutate(region = "northeast")
midwest <- import_df %>%
  filter(state %in% c("OH", "IN", "MI", "IL", "WI", "MN", "IA",
                      "MO", "ND", "SD", "NE", "KS"))%>%
  mutate(region = "midwest")
west <- import_df %>%
  filter(state %in% c("WA", "MT", "OR", "ID", "WY", "CA", "NV",
                      "UT", "CO", "AZ", "NM", "AK", "HI")) %>%
  mutate(region = "west")
south <- import_df %>%
  filter(state %in% c("WV", "VA", "KY", "TN", "NC", "SC", "GA",
```

```
"FL", "MS", "AL", "LA", "AR", "OK", "TX", "PR")) %>%
mutate(region = "south")

#rejoin them back together for future ref. if needed
full <- list(northeast, midwest, west, south) %>%
    reduce(full_join)

## Joining, by = c("state", "site", "disposal.area", "type", "well.id", "gradient", "samp.date", "contat"
## Joining, by = c("state", "site", "disposal.area", "type", "well.id", "gradient", "samp.date", "contat"
## Joining, by = c("state", "site", "disposal.area", "type", "well.id", "gradient", "samp.date", "contat"
ggplot(full, aes(x = region)) +
    geom_bar(aes(fill = region), show.legend = FALSE) +
    ggtitle("Number of Wells according to Region") +
    xlab("Region") +
    ylab("Count")
```

## Number of Wells according to Region

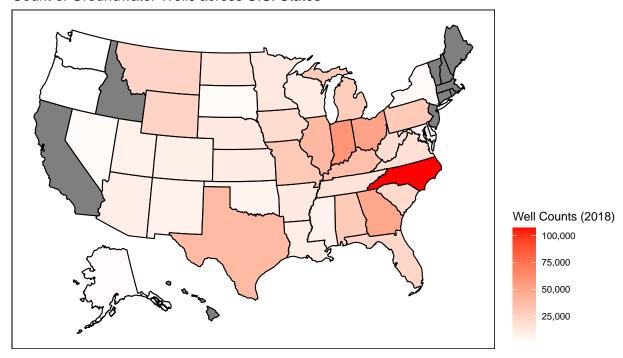


```
midwest_n <- midwest %>%
  group_by(state) %>%
  summarize(n = n()) %>%
  arrange(desc(n))
```

```
## 'summarise()' ungrouping output (override with '.groups' argument)
northeast_n <- northeast %>%
group_by(state) %>%
summarize(n = n()) %>%
```

```
arrange(desc(n))
## 'summarise()' ungrouping output (override with '.groups' argument)
south_n <- south %>%
  group_by(state) %>%
  summarize(n = n()) \%>\%
  arrange(desc(n))
## 'summarise()' ungrouping output (override with '.groups' argument)
west_n <- west %>%
  group_by(state) %>%
  summarize(n = n()) \%>\%
  arrange(desc(n))
## 'summarise()' ungrouping output (override with '.groups' argument)
states_n <- rbind(midwest_n, northeast_n, south_n, west_n)</pre>
idea: make colored map based on how many wells are in each state
state_name <- state.name</pre>
state_abb <- state.abb</pre>
states_map <- map_data("state")</pre>
plot_usmap(data = states_n, values = "n", regions = "states") +
  scale_fill_continuous(low = "white", high = "red",
                         name = "Well Counts (2018)",
                         label = scales::comma) +
  theme(legend.position = "right",
        panel.background = element_rect(color = "black",
                                         fill = "white")) +
  ggtitle("Count of Groundwater Wells across U.S. States")
```

## Count of Groundwater Wells across U.S. States



North Carolina has a significant number of wells amongst all states (over 100,000) compared to the next highest which is Indiana with around 58,000.

Let's focus in on North Carolina only for now!

```
NC <- south %>%
  filter(state %in% "NC")

#count of sites
NC %>%
  group_by(site) %>%
  summarize(n = n()) %>%
  arrange(desc(n))
```

## 'summarise()' ungrouping output (override with '.groups' argument)

```
## # A tibble: 13 x 2
##
     site
                                              n
##
      <chr>
                                          <int>
   1 "L.V. Sutton Energy Complex"
                                          15683
##
##
  2 "Belews Creek Steam Station"
                                          13414
  3 "Cliffside Steam Station"
##
                                          13362
##
  4 "Roxboro Steam Electric Plant"
                                          10320
## 5 "Allen Steam Station"
                                           9938
## 6 "Buck Steam Station"
                                           9813
## 7 "Dan River Steam Station"
                                           7885
## 8 "Mayo Steam Electric Plant"
                                           7560
## 9 "H.F. Lee Energy Complex"
                                           6120
```

```
## 10 "Marshall Steam Station"
                                            6047
## 11 "Asheville Steam Electric Plant"
                                            4115
## 12 "W.H. Weatherspoon Power Plant"
                                            2520
## 13 "Brickhaven No. 2 Mine Tract \"A\""
                                             357
There are 13 different "sites" in which the wells can belong to.
#count of disposal.area
NC %>%
  group by(disposal.area) %>%
  summarize(n = n()) \%>\%
  arrange(desc(n))
## 'summarise()' ungrouping output (override with '.groups' argument)
## # A tibble: 22 x 2
##
      disposal.area
                                                                                     n
##
      <chr>>
                                                                                 <int>
   1 Active Ash Basin
                                                                                 15434
  2 CCP Landfill
                                                                                 11572
## 3 1971 and 1984 Ash Basins
                                                                                 10620
## 4 Active Ash Basin, Retired Ash Basin, Retired Ash Basin Landfill
                                                                                  9938
                                                                                  6120
## 5 CCR Multiunit 2 (West Ash Basin, East and West FGD Settling Ponds, FGD~
## 6 Active Ash Basin and Industrial Landfill No. 1
                                                                                  6047
## 7 Craig Road Landfill
                                                                                  5156
## 8 Primary Pond (Ash Basin 2), Secondary Pond (Ash Basin 3)
                                                                                  5003
## 9 Additional Primary Pond (Ash Basin 1)
                                                                                  4810
## 10 CCR Multiunit 1 (East Ash Pond, Industrial Landfill)
                                                                                  4200
## # ... with 12 more rows
Within each well, there are multiple disposal areas also (total count of 22).
#count of gradient
NC %>%
  group_by(gradient) %>%
  summarize(n = n()) \%
  arrange(desc(n))
## 'summarise()' ungrouping output (override with '.groups' argument)
## # A tibble: 4 x 2
     gradient
                                     n
##
     <chr>
                                 <int>
## 1 Downgradient
                                 92169
## 2 Upgradient
                                 14063
## 3 Unknown
                                   629
## 4 Downgradient/Crossgradient
                                   273
```

In the case for the NC wells, there are 92.169 downgradient wells and 14,063 upgradient wells.

## Dangerous Toxins in Coal Ash

Some of the most dangerous contaminants often found in coal ash include: arsenic, lead, mercury, cadmium, chromium, and selenium (https://www.psr.org/wp-content/uploads/2018/05/coal-ash-toxics.pdf)

Are there different rates of censoring for different contaminants?

```
NC_subset <- NC %>%
  filter(contaminant %in% c("Arsenic, dissolved", "Arsenic, total",
                            "Lead, total", "Mercury, total",
                            "Cadmium, dissolved", "Cadmium, total",
                            "Chromium, total", "Selenium, Dissolved",
                            "Selenium, Total")) %>%
  #exclude crossgradient and unknown
  filter(gradient %in% c("Upgradient", "Downgradient"))
NC subset2 <- NC subset %>%
  group_by(contaminant, gradient, measurement.unit, below.detection) %>%
  summarize(n = n())
## 'summarise()' regrouping output by 'contaminant', 'gradient', 'measurement.unit' (override with '.gr
NC_subset3 <- NC_subset2[-c(27), ] #removing strange sole observation
## Warning: The 'i' argument of ''[.tbl_df'()' must lie in [-rows, 0] if negative, as of tibble 3.0.0.
## Use 'NA integer ' as row index to obtain a row full of 'NA' values.
## This warning is displayed once every 8 hours.
## Call 'lifecycle::last_warnings()' to see where this warning was generated.
NC subset4 <- NC subset3 %>%
  group_by(contaminant, gradient, measurement.unit) %>%
  summarize(prop = n/(sum(n))) \%>\%
  mutate(below.detection = case_when(
   row_number() %% 2 == 1 ~ "<", #odd
   row_number() %% 2 == 0 ~ "NA")) %>% #even
  filter(below.detection %in% "<") %>%
  arrange(order(contaminant))
```

## 'summarise()' regrouping output by 'contaminant', 'gradient', 'measurement.unit' (override with '.gr knitr::kable(NC\_subset4)

contaminant	gradient	measurement.unit	prop	below.detection
Arsenic, total	Downgradient	ug/l	0.1235191	<
Arsenic, total	Upgradient	ug/l	0.2423803	<
Cadmium, total	Downgradient	ug/l	0.3968846	<
Cadmium, total	Upgradient	ug/l	0.4702467	<
Chromium, total	Downgradient	ug/l	0.0976305	<
Chromium, total	Upgradient	ug/l	0.1567489	<
Lead, total	Downgradient	ug/l	0.3394032	<
Lead, total	Upgradient	ug/l	0.3889695	<
Mercury, total	Downgradient	ug/l	0.4429574	<
Mercury, total	Upgradient	mg/l	1.0000000	<
Mercury, total	Upgradient	ug/l	0.4912791	<

Generally, when considering only the wells that have measurements below detection – it seems like a higher proportion of the upgradient wells have measurements below detection when compared to their downgradient counterparts. This is to be expected – upgradient wells are used to measure water quality before it passes through a coal ash plant, downgradient wells to measure water quality after passing through the plant.

Ideally, we would like to know the average level of contamination (for a contaminant) with regards to upgradient and downgradient wells. However, due to the proportions of measurements which are below

the limit of detection being so high – it may pose challenges in our endeavor.

We could try to calculate averages without accounting for censoring to see what happens and then applying our methods and see if there are any differences. We have no way of knowing what the true averages will be due to so many being below the limit of detection – however, we can definitely look to see if there are differences.

@Kelderman2019 compiled a list of the "top 10 most contaminated sites in the U.S." which I wanted to delve further into:

```
top10 <- c("San Miguel Plant",</pre>
           "Allen Steam Station",
           "Jim Bridger Power Plant",
           "Naughton Power Plant",
           "New Castle Generating Station",
           "Allen Fossil Plant",
           "Brandywine Ash Management Facility",
           "Hunter Power Plant",
           "R.D. Morrow, Sr. Generating Station",
           "Ghent Generating Station")
top10df <- full %>%
  filter(site %in% top10)
top10df %>%
  group_by(site)%>%
  summarize(n = n()) \%>\%
  arrange(desc(n))
## 'summarise()' ungrouping output (override with '.groups' argument)
## # A tibble: 10 x 2
##
      site
                                               n
##
      <chr>
                                           <int>
## 1 Allen Steam Station
                                            9938
## 2 Jim Bridger Power Plant
                                            8932
## 3 Naughton Power Plant
                                            6485
## 4 San Miguel Plant
                                            4177
## 5 Ghent Generating Station
                                            3895
## 6 Hunter Power Plant
                                            2010
## 7 Brandywine Ash Management Facility
                                            1988
## 8 Allen Fossil Plant
                                            1952
## 9 New Castle Generating Station
                                            1749
## 10 R.D. Morrow, Sr. Generating Station 1701
/newpage
#how many obs are below detection in each site
top10df %>%
  group_by(site, below.detection) %>%
  summarize(n = n())
```

There is a significant number of observations below the limit of detection which was not discussed (of being accounted for) in @Kelderman2019's report. I'm curious to see if the top 10 most contaminated sites in the U.S. would change, if we start accounting for these below detection values.

How did the authors decide how to rank these top 10 sites by? They calculated a ratio of the "highest onsite average concentration of that pollutant to the health-based thresholds", and essentially the higher the ratio was – the more dangerous that contaminant was determined to be. A contaminant with the ratio "x200" had concentrations that were 200 times greater than the health-based threshold. It is also essential to note that the authors only looked at the downgradient wells to determine if a well was contaminated or not (when in reality, the upgradient *could* be contaminated also).

Table 5: The Ten Most Contaminated Sites in the Country

Rank	Site	Pollutants Exceeding Safe Levels
ı	San Miguel Plant (TX)	Arsenic (x7), Beryllium (x138), Boron (x23), Cadmium (x124), Cobalt (x522), Fluoride (x3), Lithium (x93), Mercury (x3), Radium 226+228 (x6), Selenium (x8), Sulfate (x20), Thallium (x9)
2	Allen Steam Station (NC)	Arsenic (x6), Beryllium (x6), Cadmium (x1), Cobalt (x532), Fluoride (x1), Lithium (x12), Selenium (x7), Sulfate (x3), Thallium (x1)
3	Jim Bridger Power Plant (WY)	Antimony (x1), Arsenic (x5), Boron (x6), Cadmium (x4), Cobalt (x96), Fluoride (x3), Lead (x5), Lithium (x170), Molybdenum (x12), Radium 226+228 (x2), Selenium (x116), Sulfate (x131), Thallium (x13)
4	Naughton Power Plant (WY)	Arsenic (x5), Beryllium (x2), Boron (x1), Cobalt (x3), Lead (x1), Lithium (x195), Radium 226+228 (x1), Selenium (x159), Sulfate (x65), Thallium (x14)
5	New Castle Generating Station (PA)	Arsenic (x372), Boron (x3), Cobalt (x5), Lithium (x54), Molybdenum (x1), Sulfate (x4)
6	Allen Fossil Plant (TN)	Arsenic (x350), Boron (x2), Fluoride (x1), Lead (x4), Molybdenum (x9)
7	Brandywine Ash Management Facility (MD)	Arsenic (x5), Beryllium (x2), Boron (x16), Cobalt (x47), Lithium (x222), Molybdenum (x111), Selenium (x9), Sulfate (x10)
8	Hunter Power Plant (UT)	Boron (x9), Cobalt (x26), Lithium (x228), Molybdenum (x11), Radium 226+228 (x2), Sulfate (x66)
9	R.D. Morrow, Sr. Generating Station (MS)	Arsenic (x3), Boron (x12), Lead (x1), Lithium (x193), Molybdenum (x171), Sulfate (x6), Thallium (x1)
10	Ghent Generating Station (KY)	Antimony (x2), Arsenic (x2), Beryllium (x2), Boron (x4), Chromium (x3), Cobalt (x12), Lead (x3), Lithium (x154), Molybdenum (x16), Radium 226+228 (x31), Sulfate (x3), Thallium (x2)

NOTE: The number that follows each pollutant is the ratio of the highest onsite average concentration of that pollutant to the health-based thresholds that we identify in Table B1.

Figure 1: Interval Censoring Example

We can apply the techniques in the simulation study (substitution, km, mle, ros) to estimate the average concentration of each contaminant in the downgradient wells across all sites, and compile a list of the top 10 wells, much like @Kelderman2019 did, and then compare to see whether or not if the list matches up – and if the ratio changes if we use these methods to handle left-censored values.