exploratory

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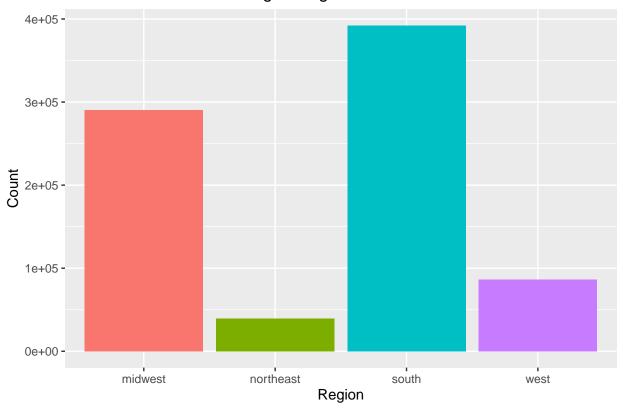
```
#Libraries
library(mosaic)
library(tidyverse)
library(usmap)
#importing in full dataset
import_df <- read_csv("data/chemical_data.csv")</pre>
## Parsed with column specification:
## cols(
##
     state = col_character(),
##
     site = col character(),
##
     disposal.area = col_character(),
##
     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

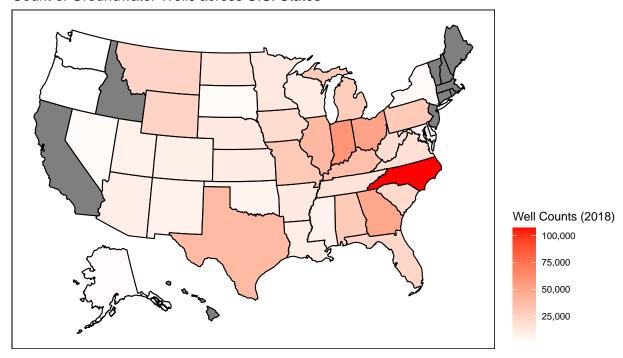


```
midwest_n <- midwest %>%
  group_by(state) %>%
  summarize(n = n()) %>%
  arrange(desc(n))
```

```
## '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
## 5 CCR Multiunit 2 (West Ash Basin, East and West FGD Settling Ponds, FGD~
                                                                                  6120
## 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 type
NC %>%
  group_by(type) %>%
  summarize(n = n()) \%
  arrange(desc(n))
## 'summarise()' ungrouping output (override with '.groups' argument)
## # A tibble: 3 x 2
     type
               n
##
     <chr> <int>
## 1 SI
           57084
## 2 M
           26309
## 3 L
           23741
```

In the case for the NC wells, there are 57,084 SI wells, 26,306 M wells, and 23,741 L 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)

Could the type of disposal unit affect the concentration (low/medium/high) of these contaminants?

```
NC_subset <- NC %>%
filter(contaminant %in% c("Arsenic, dissolved", "Arsenic, total",
```

```
"Cadmium, dissolved", "Cadmium, total",
                             "Chromium, total", "Selenium, Dissolved",
                             "Selenium, Total"))
NC_subset2 <- NC_subset %>%
  group_by(contaminant, type, measurement.unit, below.detection) %>%
  summarize(n = n())
## 'summarise()' regrouping output by 'contaminant', 'type', 'measurement.unit' (override with '.groups
NC_subset3 <- NC_subset2[-c(27), ] #removing strange sole observation
NC_subset4 <- NC_subset3 %>%
  group_by(contaminant, type, measurement.unit) %>%
  summarize(prop = n/(sum(n))) %>%
  mutate(below.detection = case_when(
    row_number() \ \%\% \ 2 == \ 1 \ \sim \ "<", \ \#odd
    row_number() %% 2 == 0 ~ "NA")) %>% #even
  filter(below.detection %in% "<") %>%
  arrange(desc(prop))
```

"Lead, total", "Mercury, total",

'summarise()' regrouping output by 'contaminant', 'type', 'measurement.unit' (override with '.groups
knitr::kable(NC_subset4)

contaminant	type	measurement.unit	prop	below.detection
Mercury, total	Μ	ug/l	0.6861314	<
Cadmium, total	\mathbf{M}	ug/l	0.5842788	<
Mercury, total	L	ug/l	0.4650767	<
Lead, total	${f M}$	ug/l	0.4505673	<
Cadmium, total	L	ug/l	0.4327087	<
Lead, total	L	ug/l	0.3637138	<
Mercury, total	SI	ug/l	0.3383978	<
Cadmium, total	SI	ug/l	0.3173343	<
Lead, total	SI	ug/l	0.2924724	<
Arsenic, total	\mathbf{M}	ug/l	0.1742301	<
Arsenic, total	L	ug/l	0.1533220	<
Chromium, total	\mathbf{M}	ug/l	0.1296596	<
Arsenic, total	SI	ug/l	0.1201657	<
Chromium, total	L	ug/l	0.1081772	<
Chromium, total	SI	ug/l	0.0949586	<