Final Workflow Assignment

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This project is based on my master's dissertation script. The original work is in .Rmd format. Here, as a follow-up to my previous *self-critique* assignment, I will try to improve my code, as well as illustrating some parts of my code that already follows the goals of this assignment (efficiency, fidelity, sharing/reproducibility)

Before starting the assignment, I need to briefly describe the "original files": thesisaltogether.Rmd is the main file in my original work. references.bib and Rreferences.bib are bibtex formats for managing my references. preamble-latex.tex is an addendum to my YAML section to address some technical issues with the formatting of figures, etc. Appendix, Figures, and Raw data contain several required files to execute the thesisaltogether.Rmd. However, the original Raw Data of my work is already posted in my GitHub repo.

I. Efficiency

Automation

##

##

psycho

TRUE

lmerTest

TRUE

rstanarm

TRUE

a function to optimize the package installation

In the original work, I inserted this function to reduce the amount of time re-installing/loading the packages.

```
required_packages <- function(pkg){</pre>
  new.pkg <- pkg[!(pkg %in% installed.packages()[, "Package"])]</pre>
  if (length(new.pkg))
    install.packages(new.pkg, dependencies = TRUE)
  sapply(pkg, require, character.only = TRUE)
}
packages <- c("rio", "readxl", "tidyverse", "devtools", "dplyr", "ggplot2", "magrittr", "Hmisc", "psycho", "lme</pre>
required_packages(packages)
                                                   dplyr
##
         rio
                 readxl tidyverse
                                     devtools
                                                            ggplot2
                                                                      magrittr
                                                                                    Hmisc
##
        TRUE
                   TRUE
                              TRUE
                                         TRUE
                                                    TRUE
                                                               TRUE
                                                                          TRUE
                                                                                     TRUE
```

TRUE

corrplot

TRUE

jtools bayesplot

TRUE

After examining my original work, I noticed that I had uploaded my "Raw Data" as an Excel spreadsheet. This means that unless the project is finished, every time I wanted to add an observation, I needed to replace the existing Excel file with a new one. This process can be improved if I insert the original works into a Google sheets file and R can have access to that file, which almost always contains the latest version of my raw data.

original piece from my work

```
# importing the main excel file from Github repository (mohdasti)
RAW_data <- rio::import('https://github.com/mohdasti/Queens-Thesis/blob/master/Raw%20data/RAW_data.xlsx
#View(RAW_data)</pre>
```

The improved version with access to the Google Sheets

I have been working on this piece of code for a while for another similar project. Here, I try to modify the code to show how it works. The code chunk is annotated.

googlesheets4::gs4_deauth()
}

if everything is fine, the the following function will run to read the google sheet. For the sake of

Here, I imported four columns from that 'raw google sheets' and renamed them in my raw spreadsheet in

```
knitr::opts_chunk$set(
  results='asis',
  echo = FALSE
)

# now, that I have assigned how I can detect the link and specified which columns to be selected and im
### the link is fake ###

library(magrittr) # For the pipe
Data <- create_data_object(
  data_location = "https://docs.google.com/spreadsheets/d/...",
  pdf_mode = params$pdf_mode
)
}</pre>
```

II. Fidelity

In terms of fidelity and implementing procedures that could reduce the additional manual copy/paste of different version of the data, the code above (using Google Sheets) can dramatically increase the fidelity of my work.

removing the hardcoding

Moreover, I went over my code and found that I used hard-coding approach to remove some observations.

```
#after importing the data from GitHub and renaming it RAW
RAW <- RAW_data
#creating subsets of data
MED <- RAW[which(RAW$Condition == 'MED'), ]</pre>
WAKE <- RAW[which(RAW$Condition == 'WAKE'), ]
NAP <- RAW[which(RAW$Condition == 'NAP'), ]
NAP_SWS <- RAW[which(RAW$percentSWS != 0), ]
NAP noSWS <- RAW[which(RAW$percentSWS == 0), ]
# Calculating Geometric Mean and Difference of Medians (two main measures of performance for the declar
# True Positive rate
RAW$TPR <-
  RAW$`Hit ratio` / (RAW$`Hit ratio` + RAW$`False Alarm ratio`)
#True Negative rate
RAW$TNR <-
  RAW$`Correct Rejection ratio` / (RAW$`Correct Rejection ratio` + RAW$`Miss ratio`)
#calculating the G-Mean
RAW$GMean <-
  sqrt(RAW$TPR * RAW$TNR)
#Excluding outliers - creating modified datasets
## any data points that is beyond ±2 SD - these exculusion are specific to each task (one participant m
## those who failed to meet the general criteria of their treatment condition (i.e. falling asleep in M
## those who their number of arousals were beyond ±2 SD
#MED$GMean[abs(scale(MED$GMean)) > 2]
```

```
#WAKE$GMean[abs(scale(WAKE$GMean)) > 2]
#NAP_SWS$GMean[abs(scale(NAP_SWS$GMean)) > 2]
#NAP_noSWS$GMean[abs(scale(NAP_noSWS$GMean)) > 2]
#creating two separate datasets based on the memory task
RAW_declr <- RAW[-c(34,39,25,13,46),]</pre>
```

So, here I would like to avoid hard-coding:

a simple plot of observations

Here, using ggplot package, I will plot the overall distribution

```
#overall pattern of distribution
NAP_dplyr <- rbind(NAP_SWS_dplyr, NAP_noSWS_dplyr)</pre>
NAP_GMean <-
  data.frame(Condition = "NAP", GeometricMean = NAP_dplyr$GMean)
MED_GMean <-
  data.frame(Condition = "MED", GeometricMean = MED dplyr$GMean)
WAKE GMean <-
  data.frame(Condition = "WAKE", GeometricMean = WAKE dplyr$GMean)
df_GMean <- rbind(NAP_GMean, MED_GMean, WAKE_GMean)</pre>
ggplot(df\_GMean, aes(x = Condition, y = GeometricMean, color = Condition)) +
  geom_point(size = 4,
             alpha = 0.7,
             position = position_jitter(w = 0.1, h = 0)) +
  stat_summary(
    fun.y = mean,
    geom = "point",
    shape = 23,
    color = "black",
    aes(fill = Condition),
    size = 4
  ) +
  stat_summary(
    fun.ymin = function(x)
      (mean(x) - sd(x)
    fun.ymax=function(x)(mean(x)+sd(x)),
    geom="errorbar", width=0.1) + xlab("Condition") + ylab("Geometric Mean") +
    theme_apa()
```

III. Sharing/Reproduciblity

As described in my last assignment, I had posted the original raw data in the Open Science Framework. Plus, I wrote the whole dissertation in RMarkdown and uploaded that in my GitHub Repository. You can also

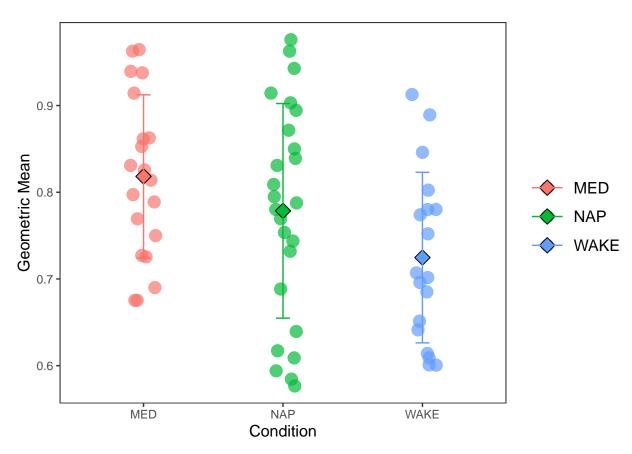


Figure 1: Overall distribution of Geometric Mean scores across the three experimental conditions (NAP; MED, meditation; WAKE). Error bars reflect standard deviation.

access the entire dissertation in this repository from thesisaltogether.Rmd.

The other issue that I briefly mentioned in my workflow critique was the usage of dev packages and how they might reduce the reproducibility. To avoid that, the groundhog package can bring a specific version of a package from a certain date. Below is a sample code chunk of groundhog—in case I needed to retrieve an older version of a package.

```
# install.packages("groundhog")
# library(groundhog)
# groundhog.library("psycho", "2018-12-26")
```