## Interdependence between Drugs

```
library(BCor) # Install via remotes::install_github("jan-lukas-wermuth/BCor") if not done yet
# Package names
packages <- c("tidyverse",</pre>
              "mvtnorm",
              "readxl", # import from Excel
              "formatR", # 'double'-pipe
              "patchwork", # plot arrangement
              "Cairo") # save pdf (only necessary to have \phi properly illustrated with qqsave). This may
# Install packages
# Loop: checks if packages are already installed. If not, they will be installed.
# Otherwise, they will be updated.
for(i in packages){
  if(!require(i, character.only = T)){
    install.packages(i, dependencies = T)
    require(i, character.only = T)
  }
}
## Loading required package: tidyverse
## -- Attaching core tidyverse packages ------ tidyverse 2.0.0 --
## v dplyr
              1.1.4
                        v readr
                                     2.1.4
              1.0.0
## v forcats
                                     1.5.1
                        v stringr
## v ggplot2 3.4.4
                        v tibble
                                     3.2.1
## v lubridate 1.9.3
                         v tidyr
                                     1.3.0
## v purrr
               1.0.2
## -- Conflicts ------ tidyverse_conflicts() --
## x dplyr::filter() masks stats::filter()
## x dplyr::lag()
                     masks stats::lag()
## i Use the conflicted package (<a href="http://conflicted.r-lib.org/">http://conflicted.r-lib.org/</a>) to force all conflicts to become errors
## Loading required package: mvtnorm
## Loading required package: readxl
##
## Loading required package: formatR
##
## Loading required package: patchwork
##
## Loading required package: Cairo
```

## Preparation work

Load data in terms of p, q and r. The structure of pqr\_arr is the following: The first dimension marks the first drug and the second dimension the second drug. The third dimension marks p, q or r. p always refers to the relative consumption frequency of the first drug in the comparison and q to the relative consumption frequency of the second drug in the comparison. r denotes the joint relative consumption frequency.

load("~/Dropbox/DimitriadisPohleWermuth/Binary Correlation/replication\_BCor/application/data\_pqr/pqr\_arr.R

## Load functions:

- 1. comparison binary: Function yielding a list containing Cole's C, Yule's Q and Phi.
- 2. plot\_contour\_matrix: Function creating heatmap with rows/columns ordered according to skewness of the marginal distribution.

```
# Insert variable1 and variable2 as strings, e.g. 'MET'
comparison_binary <- function(variable1, variable2) {</pre>
    phi <- as.numeric(Phi(pgr arr[variable1, variable2, c("p",
        "q", "r")], alpha = FALSE))
    cole <- as.numeric(Cole(pqr_arr[variable1, variable2, c("p",</pre>
        "q", "r")], alpha = FALSE))
    yule <- as.numeric(YuleQ(pqr_arr[variable1, variable2, c("p",</pre>
        "q", "r")], alpha = FALSE))
    result <- list(Phi = phi, `Cole's C` = cole, `Yule's Q` = yule)
    return(result)
}
plot_contour_matrix <- function(mat, xlab, ylab, bins) {</pre>
    drug_d_sorted <- as.factor(c("ALC", "MAR", "CIG", "AMP",</pre>
        "TRQ", "NAR", "LSD", "PSY", "SED", "MDM", "COK", "CRA",
        "MET", "HER")) # drugs sorted according to their marginal relative consumption frequencies
    drug_d_sorted_desc <- as.factor(rev(drug_d_sorted))</pre>
    data <- expand_grid(x = drug_d_sorted, y = drug_d_sorted_desc)</pre>
    z <- rep(NA, dim(mat)[1]^2)
    for (i in 1:dim(mat)[1]) {
        for (j in 1:dim(mat)[1]) {
            z[(i-1) * dim(mat)[1] + j] \leftarrow mat[j, i]
    }
    data <- mutate(data, z)</pre>
    labels <- rep("", bins + 1)</pre>
    addlabels <- seq(from = -1, to = 1, length.out = 11)
    addlabelsat <- seq(from = 1, to = bins + 1, length.out = 11)
    labels[addlabelsat] <- addlabels</pre>
    g <- ggplot(data, aes(x = fct_inorder(x), y = fct_inorder(y),
        fill = z)) + geom_tile() + scale_fill_stepsn(n.breaks = bins,
        colors = hcl.colors(20, "Spectral"), limits = c(-1, 1),
        labels = labels) + guides(fill = guide_colourbar(title = NULL,
        ticks = FALSE, barheight = 10)) + xlab(xlab) + ylab(ylab)
    g
    # 'Green-Brown' or 'Blue-Red 3' instead of 'Spectral'
    # are colorblind-safe, but for smaller ranges of values
    # do not distinguish as well
}
```

Create sorted (according to skewness) drug name vector in order to have the correlation (Cole's C, Yule's Q, Phi) matrices in the order in which the heatmap function needs it to be.

```
drug_d_sorted_desc <- rev(drug_d_sorted)</pre>
```

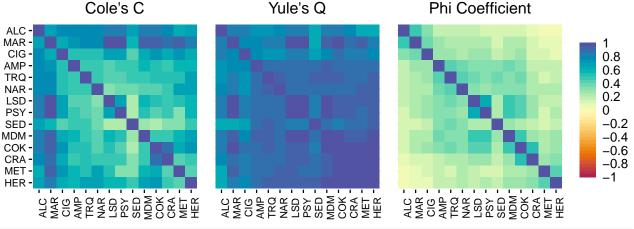
Create correlation matrices (Cole's C, Yule's Q, Phi):

```
# Matrix 1: Phi
df_phi <- matrix(NA, ncol = length(drug_d_sorted), nrow = length(drug_d_sorted))</pre>
colnames(df_phi) <- drug_d_sorted</pre>
rownames(df_phi) <- drug_d_sorted_desc</pre>
k \leftarrow 0
# Fill the data frame with the appropriate dependence
# measures for the appropriate comparison (sex vs.
# consumption of some drug)
for (i in drug_d_sorted_desc) {
    k < - k + 1
    1 <- 0
    for (j in drug_d_sorted) {
        1 <- 1 + 1
        df_phi[k, 1] <- comparison_binary(i, j)[[1]]</pre>
    }
}
# Matrix 2: Cole's C
# -----
df_cole <- matrix(NA, ncol = length(drug_d_sorted), nrow = length(drug_d_sorted))</pre>
colnames(df_cole) <- drug_d_sorted</pre>
rownames(df_cole) <- drug_d_sorted_desc</pre>
k < -0
# Fill the data frame with the appropriate dependence
# measures for the appropriate comparison (sex vs.
# consumption of some drug)
for (i in drug_d_sorted_desc) {
    k < - k + 1
    1 <- 0
    for (j in drug_d_sorted) {
        1 <- 1 + 1
        df_cole[k, 1] <- comparison_binary(i, j)[[2]]</pre>
    }
}
# Matrix 3: Yule's Q
df_yule <- matrix(NA, ncol = length(drug_d_sorted), nrow = length(drug_d_sorted))</pre>
colnames(df_yule) <- drug_d_sorted</pre>
rownames(df_yule) <- drug_d_sorted_desc</pre>
k < - 0
# Fill the data frame with the appropriate dependence
# measures for the appropriate comparison (sex vs.
# consumption of some drug)
for (i in drug_d_sorted_desc) {
    k <- k + 1
    1 <- 0
   for (j in drug_d_sorted) {
```

## **Including Plots**

Equipped with those three matrices, we can produce the respective heatmaps:

```
adj width <- 50
adj_height <- 96
size <- 14
plot_contour_matrix(df_cole, "", "", 40) + coord_fixed() + theme_minimal(base_size = size) +
    ggtitle("Cole's C") + theme(plot.title = element_text(hjust = 0.5)) +
    theme(axis.text.x = element text(color = "black", angle = 90),
        axis.text.y = element_text(color = "black"), axis.ticks = element_line(color = "black"),
        legend.text = element_text(size = size)) + plot_contour_matrix(df_yule,
    "", "", 40) + coord_fixed() + theme_minimal(base_size = size) +
    ggtitle("Yule's Q") + theme(plot.title = element_text(hjust = 0.5)) +
    theme(axis.text.x = element_text(color = "black", angle = 90),
        axis.ticks.x = element_line(color = "black"), axis.text.y = element_blank(),
        axis.ticks.y = element_blank(), axis.title.y = element_blank(),
        legend.text = element_text(size = size)) + plot_contour_matrix(df_phi,
    "", "", 40) + coord_fixed() + theme_minimal(base_size = size) +
    ggtitle("Phi Coefficient") + theme(plot.title = element_text(hjust = 0.5)) +
    theme(axis.text.x = element text(color = "black", angle = 90),
        axis.ticks.x = element_line(color = "black"), axis.text.y = element_blank(),
        axis.ticks.y = element_blank(), axis.title.y = element_blank(),
        legend.text = element_text(size = size)) + plot_layout(guides = "collect",
    ncol = 3)
```



```
# ggsave(filename = '', height = 200 - adj_height, width =
# (1+sqrt(5))/2*200 - adj_width, device = 'pdf', units =
# 'mm')
```

The last piece of code computes the confidence intervals of Cole's C, Phi and Yule's Q for all comparisons involving marijuana and meth.

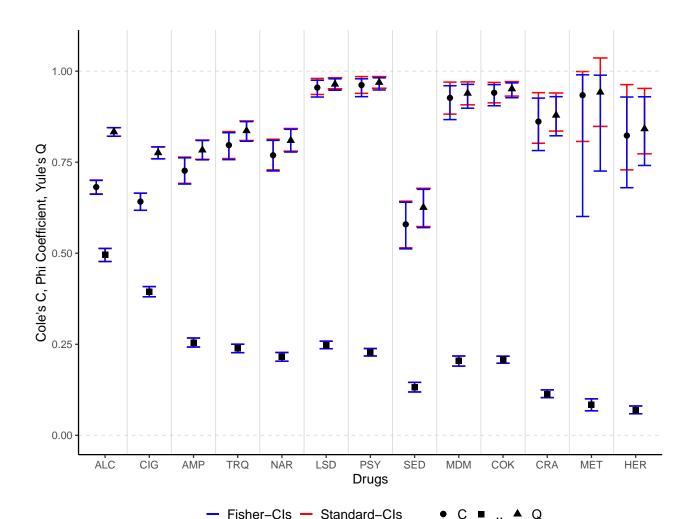
```
# Order drugs with respect to the frequency of use (except marijuana, which is our reference point)
  drugs_inference_sorted <- c("ALC", "CIG", "AMP", "TRQ", "NAR", "LSD", "PSY", "SED", "MDM", "COK", "CRA",
  drugs_inference_sorted_desc <- rev(drug_d_sorted)</pre>
marijuana_CI <- function(variable){</pre>
  C_est <- Cole(pqr_arr["MAR", variable, c("p", "q", "r")], alpha = FALSE)[[1]]</pre>
  Phi_est <- Phi(pqr_arr["MAR", variable, c("p", "q", "r")], alpha = FALSE)[[1]]
  Q_est <- YuleQ(pqr_arr["MAR", variable, c("p", "q", "r")], alpha = FALSE)[[1]]
  marijuana_CI_C <- c(Cole(pqr_arr["MAR", variable, c("p", "q", "r")], Fisher = FALSE, alpha = 0.9, n = pq
  marijuana_CI_CZ <- c(Cole(pqr_arr["MAR", variable, c("p", "q", "r")], alpha = 0.9, n = pqr_arr["MAR", va
  marijuana_CI_Q <- c(YuleQ(pqr_arr["MAR", variable, c("p", "q", "r")], Fisher = FALSE, alpha = 0.9, n = p
  marijuana_CI_phi <- c(Phi(pqr_arr["MAR", variable, c("p", "q", "r"))], Fisher = FALSE, alpha = 0.9, n = p
  marijuana_CI_phiZ <- c(Phi(pqr_arr["MAR", variable, c("p", "q", "r")], alpha = 0.9, n = pqr_arr["MAR", v</pre>
  res <- dplyr::tribble(~C_est, ~Phi_est, ~Q_est, ~marijuana_CI_C, ~marijuana_CI_CZ, ~marijuana_CI_Q, ~mari
                          #--|--|--|--|--
                         C_est, Phi_est, Q_est, marijuana_CI_C, marijuana_CI_CZ, marijuana_CI_Q, marijuan
  return(res)
}
Q_est <- NA
C_est <- NA
Phi_est <- NA
CI_lower_Q <- NA
CI_upper_Q <- NA
CI_lower_QZ <- NA
CI_upper_QZ <- NA
CI_lower_C <- NA
CI_upper_C <- NA
CI_lower_CZ <- NA
CI_upper_CZ <- NA
CI_lower_phi <- NA
CI_upper_phi <- NA
CI_lower_phiZ <- NA
CI_upper_phiZ <- NA
for (i in drugs_inference_sorted) {
  Q_est <- c(Q_est, marijuana_CI(i)$Q_est)</pre>
for (i in drugs_inference_sorted) {
  C_est <- c(C_est, marijuana_CI(i)$C_est)</pre>
for (i in drugs_inference_sorted) {
  Phi_est <- c(Phi_est, marijuana_CI(i)$Phi_est)</pre>
for (i in drugs_inference_sorted) {
  CI_lower_Q <- c(CI_lower_Q, marijuana_CI(i)$marijuana_CI_Q[[1]][1])</pre>
for (i in drugs_inference_sorted) {
  CI_upper_Q <- c(CI_upper_Q, marijuana_CI(i) marijuana_CI_Q[[1]][2])</pre>
}
for (i in drugs_inference_sorted) {
  CI_lower_QZ <- c(CI_lower_QZ, marijuana_CI(i)$marijuana_CI_QZ[[1]][1])
for (i in drugs_inference_sorted) {
```

```
CI_upper_QZ <- c(CI_upper_QZ, marijuana_CI(i)$marijuana_CI_QZ[[1]][2])</pre>
for (i in drugs_inference_sorted) {
  CI_lower_C <- c(CI_lower_C, marijuana_CI(i)$marijuana_CI_C[[1]][1])</pre>
for (i in drugs_inference_sorted) {
  CI_upper_C <- c(CI_upper_C, marijuana_CI(i)$marijuana_CI_C[[1]][2])</pre>
}
for (i in drugs_inference_sorted) {
  CI_lower_CZ <- c(CI_lower_CZ, marijuana_CI(i) marijuana_CI_CZ[[1]][1])
for (i in drugs_inference_sorted) {
  CI_upper_CZ <- c(CI_upper_CZ, marijuana_CI(i) marijuana_CI_CZ[[1]][2])
}
for (i in drugs_inference_sorted) {
  CI_lower_phi <- c(CI_lower_phi, marijuana_CI(i)$marijuana_CI_phi[[1]][1])</pre>
for (i in drugs inference sorted) {
  CI_upper_phi <- c(CI_upper_phi, marijuana_CI(i) marijuana_CI_phi[[1]][2])
for (i in drugs_inference_sorted) {
  CI_lower_phiZ <- c(CI_lower_phiZ, marijuana_CI(i) marijuana_CI_phiZ[[1]][1])
}
for (i in drugs_inference_sorted) {
  CI_upper_phiZ <- c(CI_upper_phiZ, marijuana_CI(i) marijuana_CI_phiZ[[1]][2])
}
Q_est <- as.numeric(na.omit(Q_est))</pre>
C_est <- as.numeric(na.omit(C_est))</pre>
Phi_est <- as.numeric(na.omit(Phi_est))</pre>
CI_lower_Q <- as.numeric(na.omit(CI_lower_Q))</pre>
CI_upper_Q <- as.numeric(na.omit(CI_upper_Q))</pre>
CI_lower_QZ <- as.numeric(na.omit(CI_lower_QZ))</pre>
CI_upper_QZ <- as.numeric(na.omit(CI_upper_QZ))</pre>
CI_lower_C <- as.numeric(na.omit(CI_lower_C))</pre>
CI_upper_C <- as.numeric(na.omit(CI_upper_C))</pre>
CI_lower_CZ <- as.numeric(na.omit(CI_lower_CZ))</pre>
CI_upper_CZ <- as.numeric(na.omit(CI_upper_CZ))</pre>
CI_lower_phi <- as.numeric(na.omit(CI_lower_phi))</pre>
CI_upper_phi <- as.numeric(na.omit(CI_upper_phi))</pre>
CI_lower_phiZ <- as.numeric(na.omit(CI_lower_phiZ))</pre>
CI_upper_phiZ <- as.numeric(na.omit(CI_upper_phiZ))</pre>
data_Q <- data.frame(drugs = fct_inorder(drugs_inference_sorted),</pre>
                    Q_est = Q_est,
                    C_est = C_est,
                    Phi_est = Phi_est,
                    CI_lower_Q = CI_lower_Q,
                    CI_upper_Q = CI_upper_Q,
                    CI_lower_QZ = CI_lower_QZ,
                    CI_upper_QZ = CI_upper_QZ,
                    CI_lower_C = CI_lower_C,
                    CI_upper_C = CI_upper_C,
                    CI_lower_CZ = CI_lower_CZ,
                    CI_upper_CZ = CI_upper_CZ,
                    CI_lower_phi = CI_lower_phi,
                    CI_upper_phi = CI_upper_phi,
```

```
CI_lower_phiZ = CI_lower_phiZ,
                   CI_upper_phiZ = CI_upper_phiZ)
options(ggplot2.discrete.colour= c("blue", "red"))
ggplot(data_Q, aes(drugs, Q_est)) +
    geom_errorbar(aes(ymin = CI_lower_Q, ymax = CI_upper_Q, color = "Standard-CIs"), linewidth = 0.8, widt
    geom_errorbar(aes(ymin = CI_lower_QZ, ymax = CI_upper_QZ, color = "Fisher-CIs"), linewidth = 0.8, widt
    geom_errorbar(aes(ymin = CI_lower_C, ymax = CI_upper_C, color = "Standard-CIs"), linewidth = 0.8, widt
    geom_errorbar(aes(ymin = CI_lower_CZ, ymax = CI_upper_CZ, color = "Fisher-CIs"), linewidth = 0.8, widt
    geom_errorbar(aes(ymin = CI_lower_phi, ymax = CI_upper_phi, color = "Standard-CIs"), linewidth = 0.8,
    geom_errorbar(aes(ymin = CI_lower_phiZ, ymax = CI_upper_phiZ, color = "Fisher-CIs"), linewidth = 0.8,
    geom_point(aes(drugs, Q_{est}, shape = "Q"), size = 3, position = position_nudge(x = 0.2)) +
    geom_point(aes(drugs, C_est, shape = "C"), size = 3, position = position_nudge(x = -0.2)) +
    geom_point(aes(drugs, Phi_est, shape = "\u03D5"), size = 3) +
    labs(x = "Drugs",
         y = "Cole's C, Phi Coefficient, Yule's Q") +
    geom_hline(yintercept = c(0,1), linetype = "dashed", color = "grey", alpha = 0.5) +
    geom_vline(xintercept = c(1.5, 2.5, 3.5, 4.5, 5.5, 6.5, 7.5, 8.5, 9.5, 10.5, 11.5, 12.5), color = "gre
    theme_classic(base_size = 15) +
    theme(legend.text = element text(size=15), legend.position = "bottom", axis.title=element text(size=15)
    scale_shape_manual(values = c(15, 16, 17)) +
    scale colour discrete(name = "") +
    scale_shape_discrete(name = "", breaks = c("C", "\u03D5", "Q")) +
    ylim(0, 1.06)
## Scale for shape is already present.
## Adding another scale for shape, which will replace the existing scale.
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <cf>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
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## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <cf>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <cf>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
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## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
## Warning in grid.Call.graphics(C_text, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <cf>
## Warning in grid.Call.graphics(C_text, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
```



# ggsave(filename = "", height = 200, width = (1+sqrt(5))/2\*200, device = cairo\_pdf, units = "mm")

```
options(ggplot2.discrete.colour = c("blue", "red", "green4"))
ggplot(data_Q, aes(drugs, Q_est)) + geom_errorbar(aes(ymin = CI_lower_QZ,
    ymax = CI_upper_QZ), color = "black", linewidth = 0.8, width = 0.3,
    position = position_nudge(x = 0.2)) + geom_errorbar(aes(ymin = CI_lower_CZ,
    ymax = CI_upper_CZ), color = "black", linewidth = 0.8, width = 0.3,
    position = position_nudge(x = -0.2)) + geom_errorbar(aes(ymin = CI_lower_phiZ,
    ymax = CI upper phiZ), color = "black", linewidth = 0.8,
    width = 0.3) + geom_point(aes(drugs, Q_est, shape = "Q",
    color = "Q"), size = 3, position = position_nudge(x = 0.2)) +
    geom_point(aes(drugs, C_est, shape = "C", color = "C"), size = 3,
       position = position_nudge(x = -0.2)) + geom_point(aes(drugs,
    Phi_est, shape = " ", color = " "), size = 3) + labs(x = "Drugs",
    y = "Cole's C, Phi Coefficient, Yule's Q") + geom_hline(yintercept = c(0,
    1), linetype = "dashed", color = "grey", alpha = 0.5) + geom_vline(xintercept = c(1.5,
    2.5, 3.5, 4.5, 5.5, 6.5, 7.5, 8.5, 9.5, 10.5, 11.5, 12.5),
    color = "grey", alpha = 0.5) + theme_classic(base_size = 15) +
   theme(legend.text = element_text(size = 15), legend.position = "bottom",
       axis.title = element text(size = 15)) + scale colour manual(values = c("blue",
    "red", "yellow")) + scale_shape_manual(values = c(15, 16,
    17)) + scale_colour_discrete(name = "", breaks = c("C", " ",
    "Q")) + scale_shape_discrete(name = "", breaks = c("C", " ",
    "Q")) + ylim(0, 1.06)
```

```
## Scale for colour is already present.
## Adding another scale for colour, which will replace the existing scale.
## Scale for shape is already present.
## Adding another scale for shape, which will replace the existing scale.
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <cf>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on '' in 'mbcsToSbcs': dot substituted for <cf>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
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## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
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## conversion failure on '' in 'mbcsToSbcs': dot substituted for <95>
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Cole's C, Phi Coefficient, Yule's Q
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                      AMP
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                                          LSD
                                                              MDM
                                                                            CRA
                                                                                   MET
                                                                                         HER
                                                Drugs
# ggsave(filename = '', height = 200, width
# (1+sqrt(5))/2*200, device = cairo_pdf, units = 'mm')
# Order drugs with respect to the frequency of use (except meth, which is our reference point)
  drugs_inference_sorted <- c("ALC", "MAR", "CIG", "AMP", "TRQ", "NAR", "LSD", "PSY", "SED", "MDM", "COK",
  drugs_inference_sorted_desc <- rev(drug_d_sorted)</pre>
meth_CI <- function(variable){</pre>
```

## Warning in grid.Call(C\_textBounds, as.graphicsAnnot(x\$label), x\$x, x\$y, :
## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <95>

## Warning in grid.Call(C\_textBounds, as.graphicsAnnot(x\$label), x\$x, x\$y, :

## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <cf>

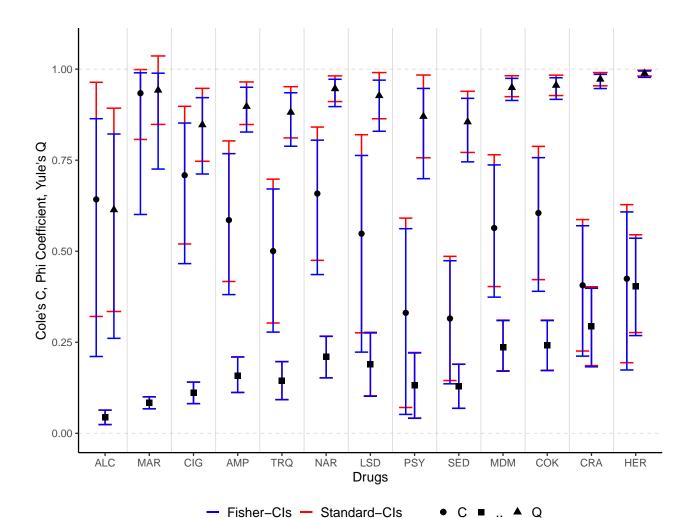
```
C_est <- Cole(pqr_arr["MET", variable, c("p", "q", "r")], alpha = FALSE)[[1]]</pre>
  Phi_est <- Phi(pqr_arr["MET", variable, c("p", "q", "r")], alpha = FALSE)[[1]]
  Q_est <- YuleQ(pqr_arr["MET", variable, c("p", "q", "r")], alpha = FALSE)[[1]]
  meth_CI_C <- c(Cole(pqr_arr["MET", variable, c("p", "q", "r")], Fisher = FALSE, alpha = 0.9, n = pqr_arr
  meth_CI_CZ <- c(Cole(pqr_arr["MET", variable, c("p", "q", "r")], alpha = 0.9, n = pqr_arr["MET", variable
  meth_CI_Q <- c(YuleQ(pqr_arr["MET", variable, c("p", "q", "r")], Fisher = FALSE, alpha = 0.9, n = pqr_ar
  meth_CI_QZ <- c(YuleQ(pqr_arr["MET", variable, c("p", "q", "r")), alpha = 0.9, n = pqr_arr["MET", variable
  meth_CI_phi <- c(Phi(pqr_arr["MET", variable, c("p", "q", "r")], Fisher = FALSE, alpha = 0.9, n = pqr_ar
  meth_CI_phiZ <- c(Phi(pqr_arr["MET", variable, c("p", "q", "r")], alpha = 0.9, n = pqr_arr["MET", variab</pre>
  res <- dplyr::tribble(~C_est, ~Phi_est, ~Q_est, ~meth_CI_C, ~meth_CI_CZ, ~meth_CI_Q, ~meth_CI_QZ, ~meth_
                           #--|--|--|--|--|--
                           C_est, Phi_est, Q_est, meth_CI_C, meth_CI_CZ, meth_CI_Q, meth_CI_QZ, meth_CI_phi
  return(res)
}
Q_est <- NA
C_est <- NA
Phi est <- NA
CI_lower_Q <- NA
CI_upper_Q <- NA
CI_lower_QZ <- NA
CI_upper_QZ <- NA
CI_lower_C <- NA
CI_upper_C <- NA
CI_lower_CZ <- NA
CI_upper_CZ <- NA
CI_lower_phi <- NA
CI_upper_phi <- NA
CI_lower_phiZ <- NA
CI_upper_phiZ <- NA
for (i in drugs_inference_sorted) {
  Q_est <- c(Q_est, meth_CI(i)$Q_est)</pre>
for (i in drugs_inference_sorted) {
  C_est <- c(C_est, meth_CI(i)$C_est)</pre>
for (i in drugs_inference_sorted) {
  Phi_est <- c(Phi_est, meth_CI(i)$Phi_est)</pre>
for (i in drugs_inference_sorted) {
  CI_lower_Q <- c(CI_lower_Q, meth_CI(i) meth_CI_Q[[1]][1])
for (i in drugs_inference_sorted) {
  CI_upper_Q <- c(CI_upper_Q, meth_CI(i)$meth_CI_Q[[1]][2])</pre>
}
for (i in drugs_inference_sorted) {
  CI_lower_QZ <- c(CI_lower_QZ, meth_CI(i) meth_CI_QZ[[1]][1])
for (i in drugs_inference_sorted) {
  CI_upper_QZ <- c(CI_upper_QZ, meth_CI(i)$meth_CI_QZ[[1]][2])</pre>
for (i in drugs_inference_sorted) {
  CI_lower_C <- c(CI_lower_C, meth_CI(i) meth_CI_C[[1]][1])</pre>
}
for (i in drugs_inference_sorted) {
  CI_upper_C <- c(CI_upper_C, meth_CI(i)$meth_CI_C[[1]][2])</pre>
```

```
for (i in drugs_inference_sorted) {
  CI_lower_CZ <- c(CI_lower_CZ, meth_CI(i)$meth_CI_CZ[[1]][1])</pre>
for (i in drugs_inference_sorted) {
  CI_upper_CZ <- c(CI_upper_CZ, meth_CI(i) meth_CI_CZ[[1]][2])</pre>
}
for (i in drugs_inference_sorted) {
  CI_lower_phi <- c(CI_lower_phi, meth_CI(i) meth_CI_phi[[1]][1])</pre>
for (i in drugs_inference_sorted) {
  CI_upper_phi <- c(CI_upper_phi, meth_CI(i) meth_CI_phi[[1]][2])</pre>
for (i in drugs_inference_sorted) {
  CI_lower_phiZ <- c(CI_lower_phiZ, meth_CI(i)$meth_CI_phiZ[[1]][1])</pre>
}
for (i in drugs_inference_sorted) {
  CI_upper_phiZ <- c(CI_upper_phiZ, meth_CI(i) $meth_CI_phiZ[[1]][2])</pre>
Q_est <- as.numeric(na.omit(Q_est))</pre>
C_est <- as.numeric(na.omit(C_est))</pre>
Phi_est <- as.numeric(na.omit(Phi_est))</pre>
CI_lower_Q <- as.numeric(na.omit(CI_lower_Q))</pre>
CI_upper_Q <- as.numeric(na.omit(CI_upper_Q))</pre>
CI_lower_QZ <- as.numeric(na.omit(CI_lower_QZ))</pre>
CI_upper_QZ <- as.numeric(na.omit(CI_upper_QZ))</pre>
CI_lower_C <- as.numeric(na.omit(CI_lower_C))</pre>
CI_upper_C <- as.numeric(na.omit(CI_upper_C))</pre>
CI_lower_CZ <- as.numeric(na.omit(CI_lower_CZ))</pre>
CI_upper_CZ <- as.numeric(na.omit(CI_upper_CZ))</pre>
CI_lower_phi <- as.numeric(na.omit(CI_lower_phi))</pre>
CI_upper_phi <- as.numeric(na.omit(CI_upper_phi))</pre>
CI_lower_phiZ <- as.numeric(na.omit(CI_lower_phiZ))</pre>
CI_upper_phiZ <- as.numeric(na.omit(CI_upper_phiZ))</pre>
data_Q <- data.frame(drugs = fct_inorder(drugs_inference_sorted),</pre>
                    Q_est = Q_est,
                    C_est = C_est,
                    Phi_est = Phi_est,
                    CI_lower_Q = CI_lower_Q,
                    CI_upper_Q = CI_upper_Q,
                    CI_lower_QZ = CI_lower_QZ,
                    CI_upper_QZ = CI_upper_QZ,
                    CI_lower_C = CI_lower_C,
                    CI_upper_C = CI_upper_C,
                    CI_lower_CZ = CI_lower_CZ,
                    CI_upper_CZ = CI_upper_CZ,
                    CI_lower_phi = CI_lower_phi,
                    CI_upper_phi = CI_upper_phi,
                    CI_lower_phiZ = CI_lower_phiZ,
                    CI_upper_phiZ = CI_upper_phiZ)
options(ggplot2.discrete.colour= c("blue", "red"))
ggplot(data_Q, aes(drugs, Q_est)) +
    geom_errorbar(aes(ymin = CI_lower_Q, ymax = CI_upper_Q, color = "Standard-CIs"), linewidth = 0.8, widt
    geom_errorbar(aes(ymin = CI_lower_QZ, ymax = CI_upper_QZ, color = "Fisher-CIs"), linewidth = 0.8, widt
    geom_errorbar(aes(ymin = CI_lower_C, ymax = CI_upper_C, color = "Standard-CIs"), linewidth = 0.8, widt
```

```
geom_errorbar(aes(ymin = CI_lower_CZ, ymax = CI_upper_CZ, color = "Fisher-CIs"), linewidth = 0.8, widt
    geom_errorbar(aes(ymin = CI_lower_phi, ymax = CI_upper_phi, color = "Standard-CIs"), linewidth = 0.8,
    geom_errorbar(aes(ymin = CI_lower_phiZ, ymax = CI_upper_phiZ, color = "Fisher-CIs"), linewidth = 0.8,
    geom_point(aes(drugs, Q_{est}, shape = "Q"), size = 3, position = position_nudge(x = 0.2)) +
    geom_point(aes(drugs, C_est, shape = "C"), size = 3, position = position_nudge(x = -0.2)) +
    geom_point(aes(drugs, Phi_est, shape = "\u03D5"), size = 3) +
    labs(x = "Drugs",
         y = "Cole's C, Phi Coefficient, Yule's Q") +
    geom_hline(yintercept = c(0,1), linetype = "dashed", color = "grey", alpha = 0.5) +
    geom_vline(xintercept = c(1.5, 2.5, 3.5, 4.5, 5.5, 6.5, 7.5, 8.5, 9.5, 10.5, 11.5, 12.5), color = "gre
    theme_classic(base_size = 15) +
    theme(legend.text = element_text(size=15), legend.position = "bottom", axis.title=element_text(size=15)
    scale_colour_manual(values = c("blue", "red")) +
    scale_shape_manual(values = c(15, 16, 17)) +
    scale_colour_discrete(name = "") +
    scale_shape_discrete(name = "", breaks = c("C", "\u03D5", "Q")) +
    ylim(0, 1.06)
## Scale for colour is already present.
## Adding another scale for colour, which will replace the existing scale.
## Scale for shape is already present.
## Adding another scale for shape, which will replace the existing scale.
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
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## conversion failure on ' ' in 'mbcsToSbcs': dot substituted for <cf>
## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
```

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## Warning in grid.Call(C_textBounds, as.graphicsAnnot(x$label), x$x, x$y, :
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```



#  $ggsave(filename = "", height = 200, width = (1+sqrt(5))/2*200, device = cairo_pdf, units = "mm")$ 

```
options(ggplot2.discrete.colour = c("blue", "red", "green4"))
ggplot(data_Q, aes(drugs, Q_est)) + geom_errorbar(aes(ymin = CI_lower_QZ,
    ymax = CI_upper_QZ), color = "black", linewidth = 0.7, width = 0.3,
   position = position_nudge(x = 0.2)) + geom_errorbar(aes(ymin = CI_lower_CZ,
   ymax = CI_upper_CZ), color = "black", linewidth = 0.7, width = 0.3,
    position = position_nudge(x = -0.2)) + geom_errorbar(aes(ymin = CI_lower_phiZ,
    ymax = CI upper phiZ), color = "black", linewidth = 0.7,
    width = 0.3) + geom_point(aes(drugs, Q_est, shape = "Q",
    color = "Q"), size = 3, position = position_nudge(x = 0.2)) +
    geom_point(aes(drugs, C_est, shape = "C", color = "C"), size = 3,
       position = position_nudge(x = -0.2)) + geom_point(aes(drugs,
    Phi_est, shape = " ", color = " "), size = 3) + labs(x = "Drugs",
    y = "Cole's C, Phi Coefficient, Yule's Q") + geom_hline(yintercept = c(0,
    1), linetype = "dashed", color = "grey", alpha = 0.5) + geom_vline(xintercept = c(1.5,
    2.5, 3.5, 4.5, 5.5, 6.5, 7.5, 8.5, 9.5, 10.5, 11.5, 12.5),
    color = "grey", alpha = 0.5) + theme_classic(base_size = 15) +
    theme(legend.text = element_text(size = 15), legend.position = "bottom",
       axis.title = element_text(size = 15)) + scale_colour_manual(values = c("blue",
    "red")) + scale_shape_manual(values = c(15, 16, 17)) + scale_colour_discrete(name = "",
    breaks = c("C", " ", "Q")) + scale_shape_discrete(name = "",
    breaks = c("C", "", "Q")) + ylim(0, 1.06)
```

## Scale for colour is already present.

```
## Adding another scale for colour, which will replace the existing scale.
## Scale for shape is already present.
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                      CİG
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                                    TRQ
                                          NAR
                                                                     MDM
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                                                Drugs
# ggsave(filename = '', height = 200, width
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# (1+sqrt(5))/2\*200, device = cairo\_pdf, units = 'mm')