



# Supplementary material: The rose of the Sainte-Chapelle in Paris: sophisticated stained glasses for late medieval painters

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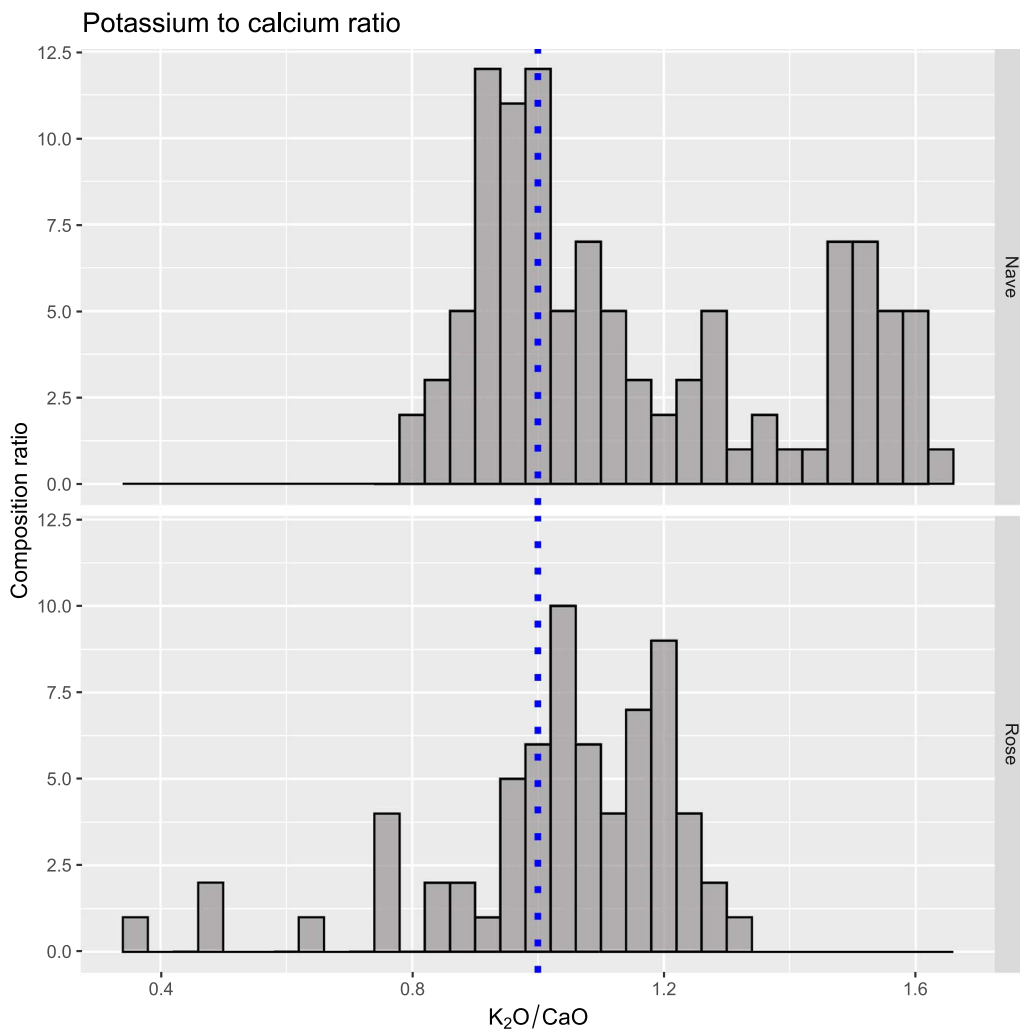
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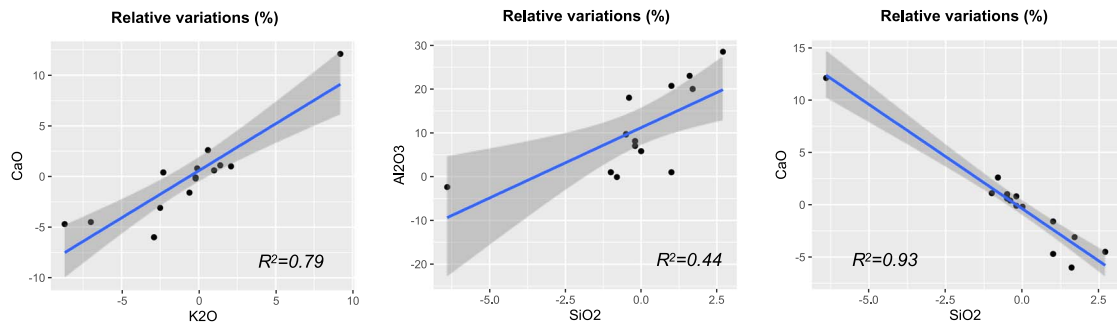
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\* Corresponding author.



**Supplementary Figure S1.** Composition ratio between potassium and calcium. We observe that the glasses from the Rose show a different pattern from the glasses from the Nave.





**Supplementary Figure S3.** Correlations between relative variations between interior and exterior side of non-flashed glasses. Linear Regression is indicated with the blue line, the r square-factor is given on each graph and the grey shadow indicates the interval of confidence with level of 0.95.

## R Code

```
##### R code to load and make the plots of the paper:
#####
#The Rose of the Sainte-Chapelle in Paris
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#####
##### HEADER PACKAGES TO LOAD #####@
#####statistics#####@
library(corrplot)
library(gridExtra)
library(FactoMineR)
library(PerformanceAnalytics)
##### plotting #####
library(ggdendro)
library(ggplot2)
library(gridExtra)
library(dendextend)
library(Hmisc)
library(reshape2)
library(dplyr)
library(ggtern)

##### HEADER PACKAGES TO LOAD #####@
## Set your working directory
#setwd(dir="~/XXXXXX")

## load data from NAVE ###
mydata_NEF <- read.table("SI_chemical_compositions_Sainte-Chapelle_Paris_nave.txt", header=TRUE,
  sep="\t", quote="", dec=".",row.names=1, na.strings = "NA") #na.string = "NULL"
mydata_NEF_sorted <- mydata_NEF[order(mydata_NEF $Na2O) ,]
mydata_NEF_ancient <- mydata_NEF_sorted[c(1:20,22:106),]
mydata_NEF_ancient_num <- mydata_NEF_ancient[,-c(1:6,41)]

## load data from ROSE ###
mydata_glass = read.table("SI_chemical_compositions_Sainte-Chapelle_Paris_rose.txt", header=TRUE,
  sep="\t", quote="", dec=".",row.names=1, na.strings = "NA") #na.string = "NULL"
mydata_glass_num <- mydata_glass[c(7:40)]

#####
#####Plots from main text #####
#####
# FIGURE 5 (Ternary diagrams)

ggtern(data = mydata_glass, aes(x =CaO , y = Na2O, z = K2O ))+
  geom_point(color = as.character(mydata_glass$Color))

colors_to_use <- as.character(mydata_glass$Color)
p5 <- ggtern(data = mydata_glass, aes(x =CaO , y = Na2O, z = K2O ))+
  geom_point(color =colors_to_use, shape= mydata_glass $Color_plaquette, labels=mydata_glass $Panel) +
  tern_limits(T=.2,L=0.6,R=0.6)

#####NAVE ternary
p <- ggtern(data = mydata_NEF_ancient, aes(x =CaO , y = Na2O, z = K2O )) +
  geom_point(color = "grey", pch = 19) + tern_limits(T=.2,L=0.6,R=0.6) + theme_light()

p <- p + theme_ggtern()
p

# FIGURE 6 a
#####
my.formula <- y ~ x
p6a <- ggplot(mydata_glass,aes(x= Al2O3, y=TiO2, color="white")) +
```

```

geom_point(color = 'white', pch=19)+
geom_smooth(method=lm,se=FALSE, color="black", formula = my.formula) +
#stat_poly_eq(formula = my.formula, aes(label = paste(.eq.label,..rr.label.., sep = "~~~")),
parse = TRUE) +
geom_point(data = mydata_NEF_ancient, aes(x = Al2O3, y = TiO2), color = 'grey', pch=19) +
  geom_point(data = mydata_glass, aes(x = Al2O3, y = TiO2), color =
as.character(mydata_glass$Color),pch= as.numeric(mydata_glass$Color_plaquette))

p6a + scale_x_continuous(limits = c(0,52000), expand = c(0,0)) +
  scale_y_continuous(limits = c(0,3800),expand = c(0,0)) +
  labs(x=expression("Al"[2]*"O"[3]* " (ppm)")) +
  labs(y=expression("Ti"[1]*"O"[2]* " (ppm)")) + theme_bw() + theme(aspect.ratio=1)

# FIGURE 6 b
#####
my.formula <- y ~ x
p6b <- ggplot(mydata_glass,aes(x= ZrO2, y=TiO2, color="white")) +
  geom_point(color = 'white', pch=19)+
  geom_smooth(method=lm,se=FALSE, color="black", formula = my.formula) +
#stat_poly_eq(formula = my.formula, aes(label = paste(.eq.label,..rr.label.., sep = "~~~")),
parse = TRUE) +
  geom_point(data = mydata_NEF_ancient, aes(x = ZrO2, y = TiO2), color = 'grey', pch=19) +
  geom_point(data = mydata_glass, aes(x = ZrO2, y = TiO2), color =
as.character(mydata_glass$Color),pch= as.numeric(mydata_glass$Color_plaquette))

p6b + scale_x_continuous(limits = c(0,350), expand = c(0,0)) +
  scale_y_continuous(limits = c(0,3800),expand = c(0,0)) +
  labs(x=expression("Zr"[1]*"O"[2]* " (ppm)")) +
  labs(y=expression("Ti"[1]*"O"[2]* " (ppm)")) + theme_bw() + theme(aspect.ratio=1)

#####
# FIGURE 7
#####
majors <- rose_alteration_2[c(5:10,13:15,31,38)]

major_Test <- melt(majors,id.vars='Plaquette',
  measure.vars=c('Na2O','MgO','Al2O3','SiO2','P2O5','K2O','CaO'), variable.name = "element",
  value.name="compo")

Rstd_FE <- data.frame(element=c('Na2O','K2O','CaO','Al2O3','MgO','P2O5','SiO2'), Rstd_FE
=c(4.6,1.0,1.9,1.8,2.6,2.4,0.9))
Rstd_FI <- data.frame(element=c('Na2O','K2O','CaO','Al2O3','MgO','P2O5','SiO2'),
  Rstd_FI=c(3.3,0.8,1.2,1.7,2.6,1.1,0.5))
Rstd_all <- data.frame(element=c('Na2O','K2O','CaO','Al2O3','MgO','P2O5','SiO2'),
  Rstd_all=c(3.9,0.9,1.5,1.7,2.6,1.8,0.7))

## Nouvel ordre des éléments chimiques
Rstd_FE <- data.frame(element=c('Na2O','MgO','Al2O3','SiO2','P2O5','K2O','CaO'), Rstd_FE
=c(4.6,2.6,1.8,0.9,2.4,1.0,1.9))
Rstd_FI <- data.frame(element=c('Na2O','MgO','Al2O3','SiO2','P2O5','K2O','CaO'),
  Rstd_FI=c(3.3,2.6,1.7,0.5,1.1,0.8,1.2))
Rstd_all <- data.frame(element=c('Na2O','MgO','Al2O3','SiO2','P2O5','K2O','CaO'),
  Rstd_all=c(3.9,2.6,1.7,0.7,1.8,0.9,1.5))

new_data_alteration <- merge(major_Test, Rstd_FE, by.x = "element", by.y = "element")
new_data_alteration <- merge(new_data_alteration, Rstd_FI, by.x = "element", by.y = "element")
new_data_alteration <- merge(new_data_alteration, Rstd_all, by.x = "element", by.y = "element")

p7 <- ggplot(new_data_alteration) +
  geom_boxplot(aes(x=element, y=compo))+ facet_wrap(~element2, scales="free", ncol=2,labeller =
  label_parsed) + labs(title="Effect of alteration",x="", y = "Relative variation (%)") +
  geom_hline(aes(yintercept = Rstd_all,color= 'red')) + geom_hline(aes(yintercept = -Rstd_all,color=
  'red')) +theme(axis.text.x=element_blank())

p7

#####
# FIGURE 8
#####
rose_diff_FE_FI = read.table("FE_FI_comparaison_RSTDev.txt", header=TRUE, sep="\t", quote="",
dec=".",row.names=1, na.strings = "NA") #na.string = "NULL"

rose_diff_FE_FI_Test <- melt(rose_diff_FE_FI,id.vars='Plaquette',
  measure.vars=c('Na2O','MgO','Al2O3','SiO2','P2O5','K2O','CaO'), variable.name = "element",

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value.name="compo")

workshop_variation <- data.frame(element=c('Na2O', 'MgO', 'Al2O3', 'SiO2', 'P2O5', 'K2O', 'CaO'),
  hline=c(14,4,6,1,3,4,7))

new_data <- merge(rose_diff_FE_FI_Test, workshop_variation, by.x = "element", by.y = "element")

element.labs <-
  c(expression("Na"[2]*"O"), expression(paste("MgO")), expression("Al"[2]*"O"[3]), expression("SiO"[2]),
  expression("P"[2]*"O"[5]), expression("K"[2]*"O"), expression("CaO"))
names(element.labs) <- c('Na2O', 'MgO', 'Al2O3', 'SiO2', 'P2O5', 'K2O', 'CaO')

new_data$element2 <- factor(new_data$element, labels =
  c("Na[2]*O", "MgO", "Al[2]*O[3]", "Si*O[2]", "P[2]*O[5]", "K[2]*O", "CaO"))

p8 <- ggplot(new_data) +
  geom_boxplot(aes(x=element, y=compo, color=Plaquege)) + facet_wrap(~element2, scales="free",
  ncol=2, labeller = label_parsed) + scale_x_discrete(breaks=NULL) + labs(title="Relative stDev
  between FE and FI", x="", y="Relative stDev (%)", col="Flashed glass") + geom_hline(aes(yintercept =
  hline))
p8

#####
# FIGURE 9
#####

data <- scale(mydata_glass_num[, c("CaO", "MgO", "K2O", "P2O5", "Na2O")]) #SiO2

d_scale <- dist(as.matrix(data))
hc_scale <- hclust(d_scale, method = "ward.D2") # methods : Ward.D2
dend_scale <- as.dendrogram(hc_scale)

dend= dend_scale

colors_to_use <- mydata_glass$Color
labels_to_use <- mydata_glass$Panel
leaves_symbols <- mydata_glass$Color_plaquege

dend %>% set("leaves_pch", 15) %>% # node point type
  set("leaves_cex", 1) %>% # node point size
  set("leaves_col", as.character(colors_to_use[order.dendrogram(dend)])) %>% # node point color
  #set("labels_col", c("green", "blue")) %>% # change label color
  set("labels_cex", 0.5) %>% # Change label size
  set("labels", as.character(labels_to_use[order.dendrogram(dend)])) %>% # change label names
  set("leaves_pch", leaves_symbols[order.dendrogram(dend)]) %>% ## set leaves
  plot(main = "Ca K Na P WARD.D2", xlab = "", ylab = "", sub = "")

#####
##### COLOR ANALYSES #####
#####

##### SPLIT COLORS #####

split_colors <- split(mydata_glass, mydata_glass$Color)

blue_glass <- split_colors$"blue"[,]
colorless_glass <- split_colors$"black"[,]
purple_glass <- split_colors$"hotpink"[,]
green_glass <- split_colors$"green"[,]
yellow_glass <- split_colors$"yellow"[,]
green_glass <- split_colors$"green"[,]
red_glass <- split_colors$"red"[,]
violet_glass <- split_colors$"violet"[,]

df <- split(purple_glass, purple_glass$Face)
new_df <- cbind(df$FI[c("MnO")], df$FE[c("MnO")], df$FI[c("Fe2O3")], df$FE[c("Fe2O3")])
names(new_df)[1] <- "MnO_1"
names(new_df)[2] <- "MnO_2"
names(new_df)[3] <- "Fe2O3_1"
names(new_df)[4] <- "Fe2O3_2"

test_purple <- mutate(purple_glass, paired = rep(1:4, 2))

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test_blue <- blue_glass[ order(row.names(blue_glass)), ]
test_blue <- mutate(test_blue, paired = rep(1:7,c(2,2,2,6,2,2,2)))

test_violet <- mutate(violet_glass, paired = rep(1:2,2))

##### Colorless + Purple GLASSES
#####
#FIGURE 11a
#####
p11a <- ggplot( test_purple) +
  geom_line(aes(x = Fe2O3, y = MnO,group = paired),color="grey", linetype = "dashed")+
  geom_point(data = mydata_glass, aes(x = Fe2O3, y = MnO), color = "grey", pch=
as.numeric(mydata_glass $Color_plaquette)) +
  geom_point(data = colorless_glass, aes(x =
Fe2O3, y = MnO), color = "black",pch= as.numeric(colorless_glass $Color_plaquette)) +
  geom_point(data= test_purple ,aes(x = Fe2O3, y = MnO), color = "hotpink",pch=
as.numeric(purple_glass $Color_plaquette))+
  geom_point(data = test_violet, aes(x = Fe2O3, y = MnO), color = "violet",pch=
as.numeric(test_violet$Color_plaquette) ) +
  geom_line(data = test_violet, aes(x = Fe2O3, y = MnO,group = paired),color="grey", linetype =
"dashed")

p11a <- p11a + coord_cartesian(xlim = c(0,18000), ylim = c(0,20000))+
  scale_x_continuous( expand = c(0,0)) +
  scale_y_continuous(expand = c(0,0)) +
  labs(x=expression("Fe"[2]*"O"[3]* " (ppm)")) +
  labs(y=expression("Mn"[1]*"O"[1]* " (ppm)")) + theme(aspect.ratio=1) + theme_bw()

p11a

#####
#FIGURE 11b
#####
p11b <- ggplot( test_purple) +
  geom_line(aes(x = BaO, y = MnO,group = paired),color="grey", linetype = "dashed")+
  geom_point(data = mydata_glass, aes(x = BaO, y = MnO), color = "grey", pch=
as.numeric(mydata_glass $Color_plaquette)) +
  geom_point(data = colorless_glass,
aes(x = BaO, y = MnO), color = "black",pch= as.numeric(colorless_glass $Color_plaquette)) +
  geom_point(data = purple_glass, aes(x = BaO, y = MnO), color = "hotpink",pch=
as.numeric(purple_glass $Color_plaquette))

p11b <- p11b + coord_cartesian(xlim = c(0,5200), ylim = c(0,20000))+
  scale_x_continuous(expand = c(0,0)) +
  scale_y_continuous(expand = c(0,0)) +
  labs(x=expression("Ba"[1]*"O"[1]* " (ppm)")) +
  labs(y=expression("Mn"[1]*"O"[1]* " (ppm)")) + theme(aspect.ratio=1) + theme_bw()

p11b

#####
#FIGURE 13
#####
# MoO3

p <- ggplot(test_blue) +
  geom_line(aes(x = CoO, y = MoO3,group = paired),color="grey", linetype = "dashed")+
  geom_point(data = mydata_glass, aes(x = CoO, y = MoO3), color = "grey", pch=
as.numeric(mydata_glass $Color_plaquette)) +
  geom_point(data = blue_glass, aes(x =
CoO, y = MoO3), color = "blue",pch= as.numeric(blue_glass $Color_plaquette)) +
  geom_point(data = test_violet, aes(x = CoO, y = MoO3), color = "violet",pch=
as.numeric(test_violet$Color_plaquette) ) +
  geom_line(data = test_violet, aes(x = CoO, y = MoO3,group = paired),color="grey", linetype =
"dashed")

p <- p + coord_cartesian(xlim = c(0,4200), ylim = c(0,150))+
  scale_x_continuous(expand = c(0,0)) +
  scale_y_continuous(expand = c(0,0)) +
  labs(x=expression("Co"[1]*"O"[1]* " (ppm)")) +
  labs(y=expression("Mo"[1]*"O"[3]* " (ppm)")) + theme(aspect.ratio=1) + theme_bw()

p
# NiO

p <- ggplot(test_blue) +
  geom_line(aes(x = CoO, y = NiO,group = paired),color="grey", linetype = "dashed")+
  geom_point(data = mydata_glass, aes(x = CoO, y = NiO), color = "grey", pch=
as.numeric(mydata_glass $Color_plaquette)) +
  geom_point(data = blue_glass, aes(x =
CoO, y = NiO), color = "blue",pch= as.numeric(blue_glass $Color_plaquette) ) +
  geom_point(data = test_violet, aes(x = CoO, y = NiO), color = "violet",pch=
as.numeric(test_violet$Color_plaquette) ) +

```



```

geom_line(data = test_violet, aes(x = CoO, y = NiO,group = paired),color="grey", linetype =
"dashed")

p <- p + coord_cartesian(xlim = c(0,4200), ylim = c(0,2200))+
  scale_x_continuous(expand = c(0,0)) +
  scale_y_continuous(expand = c(0,0)) +
  labs(x=expression("Co"[2]*"O")* " (ppm)") +
  labs(y=expression("Ni"[2]*"O")* " (ppm)") + theme(aspect.ratio=1) + theme_bw()
p

#####
#FIGURE 15b
#####
test_green <- mutate(green_glass, paired = rep(1:5,2))

test_red <- red_glass[ order(row.names(red_glass)), ]
test_red <- mutate(test_red, paired = rep(1:5,c(2,2,2,2,1)))

# ZnO

p <- ggplot(test_red) +
  geom_line(aes(x = CuO, y = ZnO,group = paired),color="grey", linetype = "dashed")+
  geom_point(data = mydata_glass, aes(x = CuO, y = ZnO), color = "grey", pch=
as.numeric(mydata_glass $Color_plaquette)) + geom_point(data = test_red, aes(x = CuO,
y = ZnO), color = "red",pch= as.numeric(test_red $Color_plaquette) ) +
  geom_point(data = test_green, aes(x = CuO, y = ZnO), color = "green",pch=
as.numeric(test_green $Color_plaquette) ) +
  geom_line(data= test_green, aes(x = CuO, y = ZnO,group = paired),color="grey", linetype = "dashed")

p <- p + coord_cartesian(xlim = c(0,15400), ylim = c(0,5000))+
  scale_x_continuous(expand = c(0,0)) +
  scale_y_continuous(expand = c(0,0)) +
  labs(x=expression("Cu"[2]*"O")* " (ppm)") +
  labs(y=expression("Zn"[2]*"O")* " (ppm)") + theme_bw() + theme(aspect.ratio=1)
p

#####
#####Plots from supporting information #####
#####
test <- cbind(mydata_NEF_ancient_num, new_col = "Nave")
test2 <- cbind(mydata_glass_num, new_col = "Rose")
test_merge <- rbind(test, test2)
K_Ca_tot <- test_merge[,c("K2O")]/test_merge[,c("CaO")]
Mg_K_tot <- test_merge[,c("MgO")]/test_merge[,c("K2O")]

test_merge <- cbind(test_merge, K_Ca_tot)
test_merge <- cbind(test_merge, Mg_K_tot)

# FIGURE S1
p1 <- ggplot(test_merge, aes(x= K_Ca_tot)) +
  geom_histogram(color = "black",fill="white", alpha=0.5, position="identity",binwidth=0.04) +
  labs(title="Potassium to calcium ratio",x=expression("K"[2]*"O"/CaO),y="Composition
ratio",col="")+
  facet_grid(new_col ~ .)+ geom_vline(xintercept = 1,linetype="dotted", color = "blue", size=1.5)
p1

# FIGURE S2
plot(test_merge[
c("CaO", "MgO", "K2O", "P2O5", "Na2O", "Al2O3", "TiO2", "ZrO2", "SiO2", "Fe2O3", "MnO", "BaO", "Rb2O", "Cl", "SrO
")],col = test_merge$new_col)

# FIGURE S3

```