# 5303hw11

# Jin Yao 2019/11/20

homework 11

# E 15.1

```
library(cfcdae)
library(conf.design)
gen1 <- rbind(c(0,1,1,1,0),c(1,0,1,1,0))
conf.design(gen1,2)
##
      Blocks T1 T2 T3 T4 T5
## 1
          00
             0
                0
                   0
                       0
## 2
          00
              1
                   1
                          0
## 3
          00
             1
## 4
          00
            0
                0
                   1
                          0
## 5
          00
              0
                0
                   0
                      0
## 6
          00
              1 1
                   1
                         1
## 7
          00
## 8
          00
             0 0
                   1
## 9
              1
                0
                   0
          01
## 10
          01
             0
                1
                         0
              0
## 11
          01
## 12
          01
              1
                0
                   1
                          0
## 13
          01
             1
                0
             0 1 1
## 14
          01
## 15
          01
             0
## 16
          01
             1
                0
                   1
## 17
          10
              0
## 18
          10
            1
## 19
          10
            1
                0
                  0
                       1
                          0
## 20
          10 0
                1
                   1
                          0
## 21
          10
             0 1
                   0
## 22
          10
             1 0
## 23
          10
             1
                0
                   0
                      1
## 24
              0
          10
## 25
             1
          11
## 26
          11
             0
## 27
          11
             0
                0
                   0
                      1
## 28
          11
              1
                1
                   1
## 29
          11
             1 1 0
## 30
          11
             0
                0 1
## 31
          11
             0
                0 0
                      1
## 32
          11
             1
                1
                   1
conf.design
```

```
## function (G, p, block.name = "Blocks", treatment.names = NULL)
## {
##
       if (!is.matrix(G))
           G <- rbind(G)
##
##
       if (is.null(treatment.names))
##
           treatment.names <- if (is.null(nam <- dimnames(G)[[2]]))</pre>
##
                .paste0("T", .zf(1:ncol(G)))
##
           else nam
##
       stopifnot(is.character(treatment.names), length(treatment.names) ==
           ncol(G), is.numeric(G), all(G >= 0), all(G\%1 == 0),
##
##
            is.numeric(p), length(p) == 1, p > 0, p\%1 == 0, p %in%
##
                primes(p), is.character(block.name), length(block.name) ==
                1)
##
       D <- as.matrix(expand.grid(rep(list(0:(p - 1)), ncol(G))))</pre>
##
##
       B <- .listMat((D %*% t(G))%%p, format)</pre>
##
       o <- do.call(order, B)
##
       B <- do.call(".paste0", B)</pre>
##
       D <- cbind(B, format(D))[o, ]</pre>
##
       dimnames(D) <- list(B[o], c(block.name, treatment.names))</pre>
##
       data.frame(.listMat(D, factor))
## }
## <bytecode: 0x7fce5d883818>
## <environment: namespace:conf.design>
```

# I use a,b,c,d to respectively represent T1,T2,T3,T4,T5

block 1:1,abc,abd,cd,e,ace,abde,cde block 2:a,bc,bd,acd,ae,bce,bde,acde block 3:b,ac,ad,bcd,be,ace,ade,bcde block 4:ab,c,d,abcd,abe,ce,de,abcde

### E 15.3

```
gen2 = rbind(c(1,1,1,0,0,1,0,0), c(1,1,0,1,1,0,0,0), c(1,0,1,1,1,0,0,0), c(0,1,1,1,0,0,0,1))
conf.set(gen2,2)
##
          [,1] [,2] [,3] [,4] [,5] [,6] [,7] [,8]
##
    [1,]
              1
                   1
                         1
                               0
                                     0
                                           1
##
   [2,]
              1
                   1
                         0
                                           0
                                                 0
                                                      0
                               1
                                     1
##
   [3,]
              0
                   0
                               1
                                     1
                                           1
                                                 0
                                                      0
   [4,]
                   0
                                           0
##
              1
                                                 0
                                                      0
                         1
                               1
                                     1
##
    [5,]
              0
                   1
                         0
                               1
                                     1
                                           1
##
   [6,]
              0
                   1
                         1
                               0
                                     0
                                           0
                                                 0
                                                      0
   [7,]
                   0
                                     0
                                           1
              1
                               0
##
   [8,]
              0
                   1
                         1
                                     0
                                           0
                                                 0
                                                      1
                               1
## [9,]
              1
                         0
                                           1
                   0
                               1
                                     0
                                                 0
                                                      1
## [10,]
              1
                   0
                                           0
                                                 0
                         1
                               0
                                     1
                                                      1
## [11,]
              0
                   1
                         0
                               0
                                     1
                                           1
                                                 0
## [12,]
              1
                   1
                         0
                               0
                                     1
                                           0
                                                 0
                                                      1
## [13,]
              0
                   0
                         1
                               0
                                     1
                                           1
                                                 0
                                                      1
                         0
                                     0
                                           0
                                                 0
## [14,]
              0
                   0
                               1
                                                      1
## [15,]
                         1
                                     0
                                           1
                                                 0
                                                      1
              1
                   1
                               1
```

#### # Use ABCDEFGH respective

counfounded:ABCF,ABDE,CDEF,ACDE,BDEF,BC,AF,BCDH,ADFH,ACEH,BEFH,ABEH,CEFH,DH,ABCDFH

### P 15.1

- (a) RCB I can split the grassland into 50 parts, so I can use RCB.
- (b) PBIBD 16 units, there are 8 treatment combinations, at about 4 blocks(4 runs per day), so lambda = 6/7
- (c) BIBD 12 units, there are 4 treatments(stints), so 3 units each group, lambda = 6
- (d) RCB there are 8 treatment combinations, because only 8 per block, so it is RCB
- (e) PBIBD there are 8 treatment combinations(A,B,C each 2), 4 units within each block, lambda = 6/7

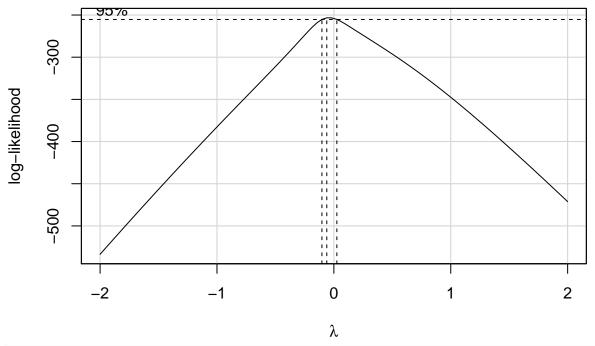
## P 15.2

- (a) youden square, there are two blocks, df(doctor)=63, df(drug)=1, df(comb)=1, df(error)=62
- (b) BIBD there are 6 treatments, 4 units within each block, df(brands)=5, df(day)=14, df(error)=40
- (c) Confounding ab, df(protein) = 1, df(choline) = 1, df(row) = 3, df(error) = 34
- (d) latin square, because each subject received drugs and each drug was used equal numbers df(subject)=29, df(sessision)=2, df(drug)=1, df(error)=28
- (e) BIBD there are 15 combinations of treatment, df(days)=2, df(charge)=2, df(additive)=4, because there are two treatments, df(charge:add)=8, df(error)=13

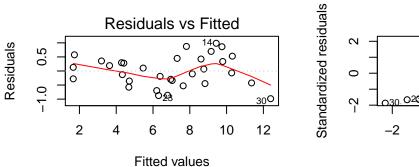
## P 15.3

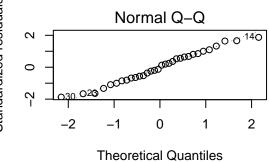
```
data("OdorIntensity")
summary(OdorIntensity)
##
                              В
                                        С
                                                   D
                                                          concentration
        judge
                    Α
##
    1
           :4
                high:16
                           high:16
                                     high:16
                                                high:16
                                                          Min.
##
    2
           :4
                low :16
                           low :16
                                     low :16
                                                low :16
                                                          1st Qu.:
                                                                      93
    3
##
           :4
                                                          Median:
                                                                     800
##
    4
           :4
                                                          Mean
                                                                  :11832
##
    5
           :4
                                                          3rd Qu.:11068
    6
           :4
                                                                  :90293
##
                                                          Max.
    (Other):8
head(OdorIntensity)
                         С
                              D concentration
##
     judge
                    В
## 1
         1
            low low low low
                                         8389
## 2
            low high high low
                                           816
## 3
         1 high high low high
                                            4
## 4
         1 high low high high
                                            46
         2 high low low low
## 5
                                          4351
## 6
         2 high high low
OdorIntensity$A =as.factor(OdorIntensity$A)
OdorIntensity$B =as.factor(OdorIntensity$B)
OdorIntensity$C =as.factor(OdorIntensity$C)
```

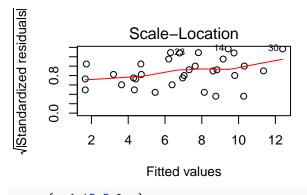
```
OdorIntensity$D =as.factor(OdorIntensity$D)
mod.15.3 <- lm(concentration~judge+A*B*C*D, data = OdorIntensity)</pre>
par(mfrow=c(2,2))
plot(mod.15.3)
                                                    Standardized residuals
                 Residuals vs Fitted
                                                                        Normal Q-Q
     20000
                                                                  270
                                                         \alpha
                               0
Residuals
     -20000
                                                         0
                0
                     20000
                                    60000
                                                                -2
                                                                                0
                                                                                               2
                      Fitted values
                                                                      Theoretical Quantiles
(Standardized residuals)
                                                                    Constant Leverage:
                                                    Standardized residuals
                   Scale-Location
                                                                 Residuals vs Factor Levels
                     0 011270
                               0 0
                                                         ^{\circ}
                                                                                  170
     0.8
                                                                      0 000
                                                                           0
                                                                                  ၀ ၀၀
                                                                 000
                                                                              o ထ
                                                                                       000
     0.0
                                                         7
                     20000
                                    60000
                                                                                              8
                0
                      Fitted values
                                                                   Factor Level Combinations
# Needs transformation
par(mfrow=c(1,1))
library(car)
## Loading required package: carData
boxCox(mod.15.3)# log
```

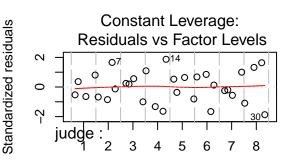


mod.15.3.log = lm(log(concentration)~judge+A\*B\*C\*D, data = OdorIntensity)
par(mfrow=c(2,2))
plot(mod.15.3.log)









**Factor Level Combinations** 

anova(mod.15.3.log)

## Analysis of Variance Table
##

```
## Response: log(concentration)
##
            Df Sum Sq Mean Sq F value
                                         Pr(>F)
## judge
            7 60.710
                       8.673 11.7590 0.0001586 ***
## A
             1 98.026 98.026 132.9079 7.548e-08 ***
## B
             1 74.362 74.362 100.8229 3.426e-07 ***
             1 0.862
## C
                       0.862
                              1.1686 0.3009266
## D
             1 10.788 10.788 14.6273 0.0024193 **
## A:B
             1 2.350
                       2.350
                               3.1863 0.0995380 .
## A:C
             1 1.125
                       1.125
                               1.5251 0.2404777
## B:C
            1 0.663
                      0.663
                               0.8986 0.3618536
## B:D
            1 0.002 0.002
                               0.0029 0.9582659
## C:D
             1 2.166
                        2.166
                               2.9364 0.1122962
                       1.175
                              1.5936 0.2307968
## A:B:D
             1 1.175
                      0.006
## A:C:D
            1 0.006
                               0.0082 0.9293287
## A:B:C:D 1 0.407
                        0.407
                               0.5520 0.4717972
## Residuals 12 8.851
                        0.738
## Signif. codes: 0 '***' 0.001 '**' 0.01 '*' 0.05 '.' 0.1 ' ' 1
\# the judge, A ,B, D are significant, the A:B is marginally significant
# I did not find A:D,A:B:C,B:C:D in the table, so it not very good.
```

### P 15.4

```
data("Milfoil")
Milfoil$A =as.factor(Milfoil$A)
Milfoil$B =as.factor(Milfoil$B)
Milfoil$C =as.factor(Milfoil$C)
mod.15.4 = lm(biomass~tank+A*B*C,data = Milfoil)
par(mfrow=c(2,2))
plot(mod.15.4)
```

