

Supplementary appendix 2. The WinBUGS code for random effects model.

The WinBUGS code for random effects model.

```
# Binomial likelihood, cloglog link
# Random effects model for multi-arm trials
model{                                     # *** PROGRAM STARTS
for(i in 1:ns){                           # LOOP THROUGH STUDIES
  w[i,1] <- 0    # adjustment for multi-arm trials is zero for control arm
  delta[i,1] <- 0          # treatment effect is zero for control arm
  mu[i] ~ dnorm(0,.0001)      # vague priors for all trial baselines
  for (k in 1:na[i]) {        # LOOP THROUGH ARMS
    r[i,k] ~ dbin(p[i,k],n[i,k]) # Binomial likelihood
  }
  # model for linear predictor
  cloglog(p[i,k]) <- log(time[i]) + mu[i] + delta[i,k]
  rhat[i,k] <- p[i,k] * n[i,k] # expected value of the numerators
}
#Deviance contribution
dev[i,k] <- 2 * (r[i,k] * (log(r[i,k])-log(rhat[i,k]))
+ (n[i,k]-r[i,k]) * (log(n[i,k]-r[i,k]) - log(n[i,k]-rhat[i,k])))
}
# summed residual deviance contribution for this trial
resdev[i] <- sum(dev[i,1:na[i]])
for (k in 2:na[i]) {                      # LOOP THROUGH ARMS
}
# trial-specific LOR distributions
delta[i,k] ~ dnorm(md[i,k],taud[i,k])
# mean of LOR distributions, with multi-arm trial correction
md[i,k] <- d[t[i,k]] - d[t[i,1]] + sw[i,k]
# precision of LOR distributions (with multi-arm trial correction)
taud[i,k] <- tau *2*(k-1)/k
# adjustment, multi-arm RCTs
w[i,k] <- (delta[i,k] - d[t[i,k]] + d[t[i,1]])
# cumulative adjustment for multi-arm trials
sw[i,k] <- sum(w[i,1:k-1])/(k-1)
}
}
totresdev <- sum(resdev[])           #Total Residual Deviance
d[1]<-0      # treatment effect is zero for reference treatment
# vague priors for treatment effects
for (k in 2:nt){ d[k] ~ dnorm(0,.0001) }
sd ~ dunif(0,2)      # vague prior for between-trial SD
tau <- pow(sd,-2)    # between-trial precision = (1/between-trial variance)

for (c in 1:(nt -1)) {
  for (k in (c+1):nt) {
    or[c,k] <- exp(d[k] - d[c])
    lor[c,k] <- (d[k] -d[c])
  }
}
```

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```
}

#Ranking of treatments#

for(k in 1:nt) {
    order[k]<- rank(d[],k)
    # this is when the outcome is positive - omit 'nt+1' when the outcome is negative
    most.effective[k]<-equals(order[k],1)

    for(j in 1:nt) {
        effectiveness[k,j]<- equals(order[k],j)
    }
}

for(k in 1:nt) {
    for(j in 1:nt) {
        cumeffectiveness[k,j]<- sum(effectiveness[k,1:j])
    }
}

#SUCRAS#

for(k in 1:nt) {
    SUCRA[k]<- sum(cumeffectiveness[k,1:(nt-1)]) /(nt-1)
}

# ranking on relative scale
for (k in 1:nt) {
    # rk[k] < - nt+1 -rank(d[],k) # assumes events are "good"
    rk[k] < - rank(d[],k) # assumes events are "bad"
    best[k] < - equals(rk[k],1) #calculate probability that treat k is best
}
}

# *** PROGRAM ENDS

#Data
# ns= number of studies; nt=number of treatments
list(ns=36, nt=15)
t[,1] r[,1] n[,1]      t[,2] r[,2] n[,2]      t[,3] r[,3] n[,3]      t[,4] r[,4] n[,4]      na[] time[]

...
...
.
.
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END
```