# EZ Bayesian Hierarchical Drift Diffusion Model

Based on Joachim's python code

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Last time knitted: 25 September, 2023

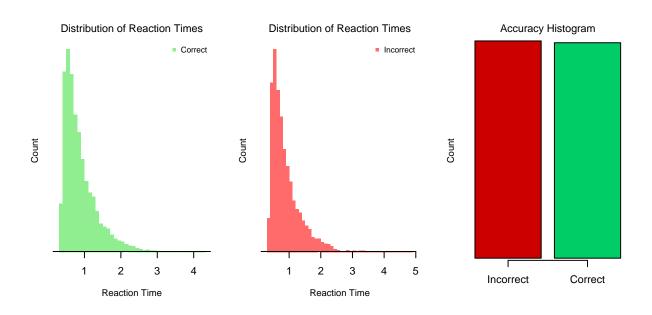
### Basic functions to generate DDM data

```
# Part 1: Simulate single trial outcome
simulate_ddm <- function(a, v, dt, max_steps){</pre>
     x <- 0
     random_dev <- rnorm(max_steps)</pre>
     # Scale step changes by dt
     noise <- random_dev * sqrt(dt)</pre>
     drift <- v * dt
     for(i in 2:max_steps){
        this_step = drift + noise[i]
        x = x + this_step
             if(abs(x)>=(a/2)){break}
     output \leftarrow list("RT" = (i+1)*dt, "C" = x)
     return(output)
}
# Part 2: Simulate over 'n' trials
wdmrnd <- function(a,v,t,n){</pre>
    dt = 0.001
    max_steps = 10 / dt
    rt = rep(NA,n)
    accuracy = rep(NA,n)
    for(i in 1:n){
        X <- simulate_ddm(a, v, dt, max_steps)</pre>
        rt[i] <- X$RT
        if(X$C>0){ accuracy[i] <- 1</pre>
                     accuracy[i] <- 0 }</pre>
          }else{
    output <- data.frame("RT" = rt + t, "accuracy" = accuracy)</pre>
    return(output)
}
```

# Example: Generate some data

```
a = 1.50
v = 0.00
t = 0.30
n = 10000

data <- wdmrnd(a, v, t, n)
rt <- data$RT
accuracy <- data$accuracy</pre>
```



#### Simulation Study environment and variables

#### Auxiliary functions

The code for the auxiliary functions listed below is hidden from this .pdf file (but can be checked on the .Rmd file).

- 1. design\_summary: A function to print the settings used in the simulation.
- 2. default\_priors: A function to load and print default prior values.
- 3. sample\_parameters: A function to sample true parameter values from the priors specified.
- 4. write\_JAGSmodel: A function to write the JAGS model using the prior values.
- 5. data\_toJAGS: A function to create a list with all the data objects to be passed to JAGS.
- 6. default\_inits: A function to create an object containing initial values for the drift.
- 7. extractSamples: A function to extract all samples associated with a parameter.name
- 8. plot.Chain: A function to plot the merging chains for hierarchical parameters
- 9. getError: A function to compute the difference between the true value and estimate retrieved for every parameter.
- 10. recoveryPlot : A function to create a quick recovery plot

#### Core functions

```
# Sample data using simulation settings and true parameter values sampled
sample_data <- function(settings, parameter_set){</pre>
  nObs <- settings$nPart*settings$nTrials
  data <- matrix(NA,ncol=3,nrow=nObs)</pre>
  data[,1] <- rep(1:settings$nPart, each=settings$nTrials)</pre>
                              # Get data for every Participant
  for(i in 1:settings$nP){
      this.sub <- which(data[,1]==i)
      accuracy = 0
      while(sum(accuracy)==0){
            temp <- wdmrnd(a = parameter_set$bound[i], v = parameter_set$drift[i],</pre>
                             t = parameter_set$nondt[i], n = settings$nTrials)
            accuracy = temp$accuracy
      }
      data[this.sub,3] <- accuracy</pre>
      data[this.sub,2] <- temp$RT</pre>
  }
  data <- as.data.frame(data)</pre>
  colnames(data) <- c("sub", "rt", "accuracy")</pre>
  return(data)
}
# Get individual statistics from raw data: mean accuracy and mean and variance of correct-RT
get_Statistics <- function(data){</pre>
  if(is.null(data$accuracy)|is.null(data$rt)){
        error.msg = "Data not available."
        return(print(error.msg))
    }
  subID <- unique(data$sub)</pre>
  sum_correct <- tapply(data$accuracy, data$sub, sum)</pre>
  # Remove participants with no correct answer
  always_0 <- which(sum_correct==0)</pre>
  if(length(always_0)!=0){
        bad_participants <- (data$sub %in% always_0)</pre>
        data <- data[-bad_participants,]</pre>
        sum_correct <- tapply(data$accuracy, data$sub, sum)</pre>
```

```
# Get proportion of correct responses
mean_accuracy <- tapply(data$accuracy, data$sub, mean)
# Get mean and variance of correct RT
keep.correct <- which(data$accuracy==1)
correct_only <- data[keep.correct,]
mean_rt_correct <- tapply(correct_only$rt, correct_only$sub, mean)
var_rt_correct <- tapply(correct_only$rt, correct_only$sub, var)
# Create a data.frame with just summary statistics
data_statistics <- cbind(subID, sum_correct, mean_accuracy, mean_rt_correct, var_rt_correct)
data_statistics <- as.data.frame(data_statistics)
colnames(data_statistics) = c("sub", "sum_correct", "meanAccuracy", "meanRT_correct", "varRT_correct")
return(data_statistics)
}</pre>
```

#### Main functions

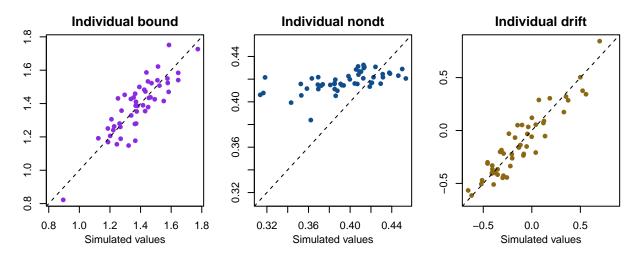
```
# A function to load priors and true values to use in simulation
Hddm_Parameter_Set <-function(nParticipants, nTrials, Show=TRUE){</pre>
    if(Show){ design_summary(nParticipants,nTrials) }
    prior <- default_priors(Show)</pre>
    settings <- list("nPart"= nParticipants, "nTrials"= nTrials, "prior"= prior)</pre>
    parameter_set <- sample_parameters(settings, Show)</pre>
    return(list("settings" = settings, "parameter_set" = parameter_set))
}
# A function to generate raw data, summary statistics and data to be passed on JAGS
Hddm_Data <- function(settings, parameter_set){</pre>
    rawData = sample data(settings,parameter set)
    sumData = get_Statistics(rawData)
    jagsData = data_toJAGS()
    return(list("rawData" = rawData, "sumData" = sumData, "jagsData" = jagsData))
}
# A function to run JAGS model
Hddm_runJAGS <- function(getData, settings, n.chains, modelFile="./EZHBDDM.bug", plot.Chains = FALSE){</pre>
    # Write model
    write_JAGSmodel(settings$prior)
    # Load settings
    parameters <- c("bound_mean", "drift_mean", "nondt_mean", "bound", "nondt",
                     "drift_sdev", "nondt_sdev", "bound_sdev", "drift")
    myinits
               <- default_inits(n.chains, settings$nPart)</pre>
    data <- getData$jagsData
    # Prepare data
            <- getData$sumData$sub</pre>
    sub
    correct <- getData$sumData$sum_correct</pre>
    varRT <- getData$sumData$varRT_correct</pre>
    meanRT <- getData$sumData$meanRT_correct</pre>
    nTrialsPerPerson <- as.numeric(unique(tapply(getData$rawData$accuracy,getData$rawData$sub,length)))
                     <- length(getData$sumData$sub)</pre>
    nParticipants
    # Run model and get samples
    suppressMessages(library(R2jags))
    suppressMessages(samples <- jags(data=data,</pre>
                                  parameters.to.save=parameters,
```

```
model=modelFile,
                                 n.chains=n.chains,
                                 n.iter=1000,
                                 n.burnin=200,
                                 n.thin=1,
                                 DIC=T,
                                 inits=myinits))
    if(plot.Chains){    plot.Chain(samples)
return(list("drift" = apply(extractSamples("drift", samples),3,mean),
            "drift_mean" = mean(extractSamples("drift_mean", samples)),
            "drift_sdev" = mean(extractSamples("drift_sdev", samples)),
            "bound" = apply(extractSamples("bound", samples),3,mean),
            "bound_mean" = mean(extractSamples("bound_mean", samples)),
            "bound sdev" = mean(extractSamples("bound sdev", samples)),
            "nondt" = apply(extractSamples("nondt", samples),3,mean),
            "nondt_mean" = mean(extractSamples("nondt_mean", samples)),
            "nondt_sdev" = mean(extractSamples("nondt_sdev", samples))))
}
# Main function: A function to run a complete simulation for nParticipants and nTrials per participant
Hddm_runSim <- function(nParticipants, nTrials, n.chains = 4, Show=TRUE){</pre>
      design.parameters <- Hddm_Parameter_Set(nParticipants,nTrials, Show=Show)
      settings <- design.parameters$settings</pre>
      parameter_set <- design.parameters$parameter_set</pre>
      getData <- Hddm_Data(settings,parameter_set)</pre>
      estimates <- Hddm_runJAGS(getData=getData, n.chains = n.chains, settings = settings)
      error <- getError(estimates, parameter_set)</pre>
return(list("trueValues" = parameter_set, "estValues" = estimates, "error" = error))
}
Hddm_simStudy <- function(nSim, nParticipants, nTrials, n.chains = 4,</pre>
                           outputFile = "./output_simStudy.RData", Show=FALSE, forceSim = FALSE){
    runSim <- TRUE
    if((!forceSim)&(file.exists(outputFile))){runSim <- FALSE}</pre>
    if(runSim){
      simOutput <- array(NA, dim=c(nSim,3,3))</pre>
      for(k in 1:nSim){
          set.seed(k)
          cat("Iteration", k, "of", nSim,"\n")
          tryCatch({
            sim <- Hddm_runSim(nParticipants = 50, nTrials = 150, Show = FALSE)</pre>
            simOutput[k,,1] <- c(sim$trueValues$drift_mean, sim$estValues$drift_mean, sim$error$drift_mean)</pre>
            simOutput[k,,2] <- c(sim$trueValues$bound_mean, sim$estValues$bound_mean, sim$error$bound_mean)
            simOutput[k,,3] <- c(sim$trueValues$nondt_mean, sim$estValues$nondt_mean, sim$error$nondt_mean)},
            error=function(e){cat("ERROR :",conditionMessage(e), "\n")})
      }
      simOutput <- simOutput[which(!is.na(simOutput[,1,1])),,]</pre>
      dimnames(simOutput) <- list(NULL, c("true", "est", "error"), c("drift_mean", "bound_mean", "nondt_mean"))
      save(simOutput,file=outputFile)
    }else{
              load(outputFile)
    return(simOutput)
```

### Run simple example

```
sim <- Hddm_runSim(nParticipants = 50, nTrials = 150)</pre>
```

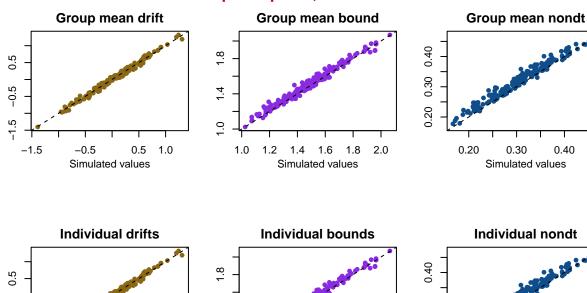
```
====== EZBHDDM Design Parameters: ======
## Number of Participants:
                             50
## Trials Per Person:
                             150
## ====== EZBHDDM Priors: =========
## Bound Mean Mean:
                    1.5
## Bound Mean Std Dev: 0.2
## Drift Mean Mean:
## Drift Mean Std Dev: 0.5
## Non-decision Time Mean Mean: 0.3
## Non-decision Time Mean Std: 0.06
## Bound Std Dev Shape: 0.1
## Bound Std Dev Scale: 0.2
## Drift Std Dev Shape: 0.2
## Drift Std Dev Scale: 0.4
## Non-decision Time Shape: 0.01
## Non-decision Time Scale: 0.05
## ==== EZBHDDM True Parameters: ========
## Bound Mean: 1.387905
## Bound SD:
                 0.1528105
## Drift Mean:
                -0.1150887
## Drift SD:
                0.3784838
## Non-decision Time Mean: 0.3935225
  Non-decision Time SD:
                         0.0320574
  _____
##
  Compiling model graph
     Resolving undeclared variables
     Allocating nodes
##
  Graph information:
     Observed stochastic nodes: 150
##
##
     Unobserved stochastic nodes: 156
##
     Total graph size: 1579
## Initializing model
```

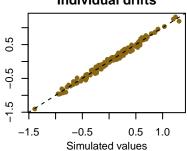


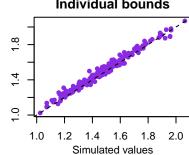
# Smulation study (150 trials - 200 repetitions)

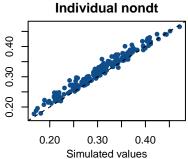
```
nT <- 150
nP <- 50
simStudy <- Hddm_simStudy(nSim = 200, nParticipants = nP, nTrials = nT, outputFile = "./sim-P50T150.RData")
showRecovery(simStudy, nP, nT)
```

### 50 participants, 150 trials each



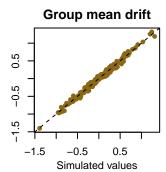


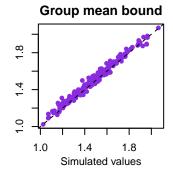


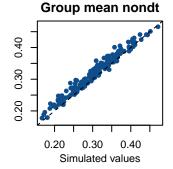


0.40

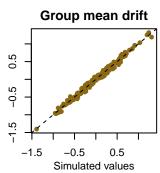
# 50 participants, 150 trials each

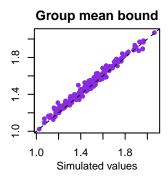


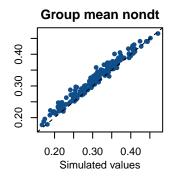




# 50 participants, 50 trials each







# 500 participants, 10 trials each

