Causal Dynamic Resonance

Claudia Lainscsek

This software reproduces the results in the paper "Causal Dynamic Resonance", submitted to PNAS and can be found on Github (https://github.com/lclaudia/CDR). The code is written in JULIA and works on Linux, Mac, and Windows. For the epilepsy part we chose to include a shorter data file due to size and computation time. The Rössler part is written as a sequential code (calling julia) and for the epilepsy part we use parallel capabilities of JULIA (calling julia -pn, where n are the parallel nodes to be used).

The underlying C codes are compiled using cosmocc from https://github.com/jart/cosmopolitan.

LINUX: If you get the error "run-detectors: unable to find an interpreter", you can fix that by running these commands:

(see https://github.com/jart/cosmopolitan/blob/master/tool/cosmocc/README.md for more details).

```
sudo wget -0 /usr/bin/ape https://cosmo.zip/pub/cosmos/bin/ape-$(uname -m).elf
sudo chmod +x /usr/bin/ape
sudo sh -c "echo ':APE:M::MZqFpD::/usr/bin/ape:' >/proc/sys/fs/binfmt_misc/register"
sudo sh -c "echo ':APE-jart:M::jartsr::/usr/bin/ape:' >/proc/sys/fs/binfmt_misc/register"
```

WINDOWS: Microsoft might be seeing run_DDA_ASCII.exe as a virus and is deleting it. To fix this problem, turn the "Real-time protection" (temporarily) off to execute the codes.

1 Single Rössler system

Before introducing coupled Rössler systems the code to integrate a single system is presented. The equations for the Rössler system are

$$\dot{u}_1 = -u_2 - u_3
\dot{u}_2 = u_1 + a u_2
\dot{u}_3 = b - c u_3 + u_1 u_3$$
(1)

with a = 0.2 and c = 5.7 and $\delta t = 0.05$. This system can be encoded as

system	equation $\#$	variable		coefficients
$\dot{u}_1 = -u_2 - u_3$	0	0	2	-1
$\dot{u}_1 = -u_2 - \frac{u_3}{}$	0	0	3	-1
$\dot{u}_2 = \frac{u_1}{u_1} + a u_2$	1	0	1	1
$\dot{u}_2 = u_1 + \frac{a}{2} u_2$	1	0	2	a
$\dot{u}_3 = \frac{b}{c} - c u_3 + u_1 u_3$	2	0	0	b
$\dot{u}_3 = b - \frac{c}{c} \frac{u_3}{u_3} + u_1 u_3$	2	0	3	-c
$\dot{u}_3 = b - c u_3 + \mathbf{u_1} \mathbf{u_3}$	2	1	3	1

Note, that the equation numbers are (0,1,2) for the three equations. This defines DIM=3 (the number of equations). There are two "variable" columns which define the order of nonlinearity ODEorder=2. The numbers in the two columns are 1 for u_1 , 2 for u_2 , and 3 for u_3 . A line with only zeros denotes a constant term. All other entries are filled with zeros.

This encoding can be used to numerically integrate the Rössler system. The plots are shown in Fig. 1.

```
include("DDAfunctions.jl");
                                                                        # set of Julia functions
NrSyst=1;
                                                                        # 1 single system
ROS=[[0 0 2];
                                                                        # single Roessler system
     [0 0 3];
     [1 0 1];
     [1 0 2];
     [2 0 0];
     [2 0 3];
     [2 1 3]
    ];
 (MOD_nr,DIM,ODEorder,P) = make_MOD_nr(ROS,NrSyst);
                                                                         # encoding of the Roessler system
                                                                         # function defined in DDAfunctions.jl
a=.2; c=5.7;
dt=.05; X0=rand(DIM,1);
                                                                         # choice of parameters
L=10000; TRANS=5000;
                                                                         # integration length and transient
b=0.45;
                                                                         # chaotic attractor
MOD_par=[-1 -1 1 a b -c 1];
                                                                         # parameters
X = integrate_ODE_general_BIG(MOD_nr,MOD_par,dt,
                              L, DIM, ODEorder, XO, TRANS);
                                                                         # integrate system
                                                                         # function defined in DDAfunctions.jl
                                                                         # plot the attractor
plot(X[:,1],X[:,2],X[:,3],
     color=:blue,legend=false,
     xlabel=L"x",ylabel=L"y",zlabel=L"z")
plot!(size=(500,500))
display(current());
print("Make pdf file and continue? ");
readline()
savefig("Roessler_0.45.pdf")
b=1;
                                                                         # periodic attractor
MOD_par=[-1 -1 1 a b -c 1];
                                                                         # parameters
X = integrate_ODE_general_BIG(MOD_nr,MOD_par,dt,
                                                                         # integrate system
                              L, DIM, ODEorder, XO, TRANS);
                                                                         # function defined in DDAfunctions.jl
plot(X[:,1],X[:,2],X[:,3],
                                                                         # plot the attractor
     color=:blue,legend=false,
     xlabel=L"x", ylabel=L"y", zlabel=L"z")
plot!(size=(500,500))
display(current());
print ("Make pdf file and continue? ");
readline()
savefig("Roessler_1.pdf")
```

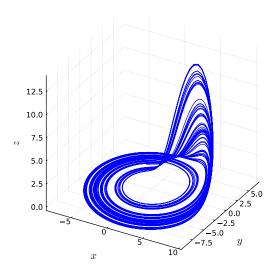
2 Coupled Rössler systems

We couple Rössler systems using diffusive coupling and consider here seven (coupled) Rössler systems

$$\dot{u}_{1,n} = -u_{2,n} - u_{3,n} + \sum_{j} \epsilon(u_{1,n} - u_{1,j})
\dot{u}_{2,n} = u_{1,n} + a_n u_{2,n}
\dot{u}_{3,n} = b_n + c_n u_{3,n} + u_{1,n} u_{3,n}$$
(2)

with n = 1, 2, ..., 7 and x_j is the u_1 -component of another system j. The values for a_n , b_n , and c_n are listed in Tab. 1. ϵ is either 0 or 0.15 depending on which systems are coupled.

We have 7 three-dimensional systems and therefore 21 variables. In the code we want to number them $x_1, x_2, \dots x_{21}$ and therefore need to make the following change in variables: $u_{1,n} \to x_{3n-2}, u_{2,n} \to x_{3n-1}$,



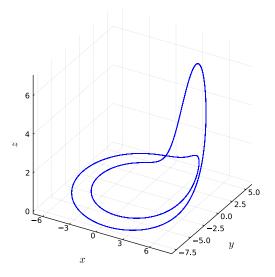


Figure 1: Rössler attractor with b = 0.45 (left) and b = 0.1 (right)

 $u_{3,n} \to x_{3n}$. In general, for a N-dimensional system we would have $u_{k,n} \to x_{Nn-(N-k)}$ This change of variables changes system (2) to

$$\dot{x}_{3n-2} = -x_{3n-1} - x_{3n} + \sum_{j} \epsilon(x_{3n-2} - x_{3j-2})
\dot{x}_{3n-1} = x_{3n-2} + a_n x_{3n-1}
\dot{x}_{3n} = b_n + c_n x_{3n} + x_{3n-2} x_{3n}$$
(3)

In the code we first encode the 7 systems without the coupling part:

```
a123=0.21;

a456=0.20;

a7 =0.18;

b1 = 0.2150;

b2 = 0.2020;

b3 = 0.2041;

b4 = 0.4050;

b5 = 0.3991;

b6 = 0.4100;
```

Table 1: Parameters of the seven Rössler systems

_#	a_n	b_n	c_n
1	0.21	0.21505	-4.5
2	0.21	0.20201	-4.5
3	0.21	0.20411	-4.5
4	0.20	0.40503	-4.5
5	0.20	0.39905	-4.5
6	0.20	0.41000	-4.5
7	0.18	0.50000	-6.8

```
b7 = 0.5000;

c =5.7;

c7=6.8;

MOD_par=[

-1 -1 1 a123 b1 -c 1

-1 -1 1 a123 b2 -c 1

-1 -1 1 a123 b3 -c 1

-1 -1 1 a456 b4 -c 1

-1 -1 1 a456 b5 -c 1

-1 -1 1 a456 b6 -c 1

-1 -1 1 a7 b7 -c7 1

];

MOD_par=reshape(MOD_par', size(ROS, 1) *NrSyst)';
```

2.1 First network example

For the first example the network of the seven Rössler systems in (2) reads as

$$\Xi \begin{cases}
\dot{x}_1 &= -y_1 - z_1 + 0.15(x_1 - x_7) \\
\dot{y}_1 &= x_1 + a_1 y_1 \\
\dot{z}_1 &= b_1 + c_1 z_1 + x_1 z_1 \\
\begin{cases}
\dot{x}_2 &= -y_2 - z_2 \\
\dot{y}_2 &= x_2 + a_2 y_2 \\
\dot{z}_2 &= b_2 + c_2 z_2 + x_2 z_2 \\
\begin{cases}
\dot{x}_3 &= -y_3 - z_3 \\
\dot{y}_3 &= x_3 + a_3 y_3 \\
\dot{z}_3 &= b_3 + c_3 z_3 + x_3 z_3
\end{cases}$$

$$\Xi \begin{cases}
\dot{x}_4 &= -y_4 - z_4 + 0.15(x_4 - x_7) \\
\dot{y}_4 &= x_4 + a_4 y_4 \\
\dot{z}_4 &= b_4 + c_4 z_4 + x_4 z_4 \\
\dot{z}_4 &= b_4 + c_4 z_4 + x_4 z_4 \\
\dot{z}_5 &= -y_5 - z_5 + 0.15(x_5 - x_7)
\end{cases}$$

$$\Xi \begin{cases}
\dot{x}_5 &= -y_5 - z_5 + 0.15(x_5 - x_7) \\
\dot{y}_5 &= x_5 + a_5 y_5 \\
\dot{z}_5 &= b_5 + c_5 z_5 + x_5 z_5 \\
\dot{z}_6 &= -y_6 - z_6 \\
\dot{y}_6 &= x_6 + a_6 y_6 \\
\dot{z}_6 &= b_6 + c_6 z_6 + x_6 z_6
\end{cases}$$

$$\Xi \begin{cases}
\dot{x}_7 &= -y_7 - z_7 + 0.15(x_7 - x_3) \\
\dot{y}_7 &= x_7 + a_7 y_7 \\
\dot{z}_7 &= b_7 + c_7 z_7 + x_7 z_7
\end{cases}$$

with the values for a_n , b_n , and c_n (n = (1, 2, ..., 7)) as listed in (1). (i), (ii), and (iii) denote the different dynamical regimes. The adjacency matrix and network motif of ground truth in (4) is shown in (2). The three different colors in the nodes of the motif denote the three dynamical regimes (i), (ii), (iii).

(4)

The encoding for the couplings in Fig. 2 is done in the following way:

from		to			
\underline{i}	Eq. #	variable	j	Eq. #	variable
3	0	0.1	7	0	0.1
7	0	0.1	1	0	0 1
7	0	0.1	4	0	0 1
7	0	0.1	5	0	0.1

```
FromTo=[
       [3 0 0 1 7 0 0 1];
       [7 0 0 1 1 0 0 1];
       [7 0 0 1 4 0 0 1];
       [7 0 0 1 5 0 0 1]
       ];
PF = "__FirstExample";  # parts of file names

II=make_MOD_nr_Coupling(FromTo,DIM,P);  # MOD_nr part for coupling epsilon=0.15;  # coupling strength
MOD_par_add=repeat([epsilon -epsilon],size(FromTo,1),1)'[:]';  # MOD_par for coupling part
```

We need to adjust the integration length according to the DDA parameters. For data of length L, the maximal delay TM, the number of data points for numerical integration dm, a window length WL, and a window shift WS the window number we loose dm + TM data points at the beginning of the time series and dm data points at the end. The number of windows WN of the DDA output is then WN = 1 + floor((L-WL-TM-2*dm)/WS): For anticipated 100 windows we then can compute the data length.

```
TAU=[32 9]; TM=maximum(TAU); dm=4; # DDA parameters
WL=4000;WS=2000; # window length and window shift for DDA
WN=100; # assign window number
LL=WS*(WN-1)+WL+TM+2*dm; # ajust integration length
```

The seven Rössler systems are integrated with a step size of 0.05 and down-sampled by a factor of two.

```
TRANS=20000;
                                                                        # transient
                                                                        # integration step size
dt = 0.05;
X0=rand(DIM*NrSyst,1);
                                                                        # initial conditions
DATA_DIR="DATA"; dir_exist(DATA_DIR);
                                                                        # DATA folder
noise="NoNoise";NOISE="NoNoise";
                                                                        # parts of file names
FN=@sprintf("%s%sCD_DDA_data_%s__WL%d_WS%d_WN%d%s.ascii",
            DATA_DIR, SL, noise, WL, WS, WN, PF);
                                                                        # noise free data file
CH_list=1:DIM:DIM*NrSyst;
                                                                        # only x
DELTA=2;
                                                                        # every second data point
if !isfile(FN)
   integrate_ODE_general_BIG([MOD_nr II],[MOD_par MOD_par_add],
                                                                        # encoding of the coupled systems
                                                                        # step size of num. integration
```

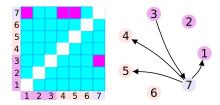


Figure 2: Adjacency matrix and network motif of ground truth. The connections are indicated as magenta boxes.

```
LL, # length
DIM*NrSyst,ODEorder,X0, # parameters
FN,
CH_list,DELTA,
TRANS);
end
```

We add noise to the data and run DDA

```
SNRadd_list= 20:-1:0;
                                                                          # add noise
MakeDataNoise(PF, noise, SNRadd_list);
DDA_DIR="DDA"; dir_exist(DDA_DIR);
                                                                         # DDA folder
nr_delays=2;
DDAmodel=[[0 0 1];
          [0 0 2];
          [1 1 1]];
(MODEL, L_AF, DDAorder) = make_MODEL(DDAmodel);
                                                                         # DDA model
NrCH=NrSyst; CH=collect(1:NrCH);
LIST=collect(combinations(CH,2));
LL1=vcat(LIST...)';
LIST=reduce(hcat, LIST)';
                                                                          # pairwise combinations
RunDDA(PF, NOISE, SNRadd_list);
                                                                          # run DDA
```

Then we plot the results:

```
(C,E) =makeCE(PF,NOISE,SNRadd_list);
                                                                                                                                                                                                                                                                                        # read outputs
 c=reshape(C[:,:,1],NrSyst^2); IDX=reverse(sortperm(c[:])); IDX=IDX[1:NrSyst^2-NrSyst];
 c=reshape(C,NrSyst^2,length(SNRadd_list)+1); c=c[IDX,:];
 e=reshape(E,NrSyst^2,length(SNRadd_list)+1); e=e[IDX,:];
 S=[repeat(1:NrSyst,NrSyst,1) repeat(1:NrSyst,1,NrSyst)'[:]][IDX,:];
 p1=Plots.palette(cgrad(:cool, scale=:log, rev=true), 42);
 p2=Plots.palette(cgrad(:cool,scale=:log),42);
 A = reshape(C[:,:,1], NrSyst^2); idx = reverse(sortperm(A[:])); A = A[idx[1:NrSyst^2-NrSyst]]; idx = reverse(sortperm(A[:])); A = A[idx[1:NrSyst]]; A = A[idx[1:NrSyst]]; A = A[idx[1:NrSyst]]; A = A[idx[1:NrSyst]]; A = A[i
 A=A .- A[end]; A=A ./ A[1] .* 1000; A=Int.(floor.(A)); A[A .== 0] .= 1;
 p1=Plots.palette(cgrad(:cool, scale=:log), 1000); p1=p1[A];
 pp=plot(size=(1000,550), margins=10*Plots.px, layout=@layout[a{0.8w} [b;c]]);\\
 \texttt{scatter!} (\texttt{pp,subplot=1,c[:,1:end]',e[:,1:end]',msw=0,palette=p1,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5,markersize=5
                                                xscale=:log10,yscale=:log10,label="")
 plot!(pp,subplot=1,c',e',msw=0,palette=p1,linewidth=0.2,xscale=:log10,yscale=:log10,label="",grid=false)
 plot!(pp, subplot=1, xlabel=L"$\mathcal{C}$");
 plot!(pp, subplot=1, ylabel=L"$\mathcal{E}$");
 for k=1:size(FromTo, 1)+3
                 txt=join([string(S[k,1]) " \u21FE " string(S[k,2])])
                annotate! (pp, subplot=1, c[k, 1], e[k, 1], Plots.text(txt, 18, p1[k]));
 end
 display(pp)
 cc=C[:,:,1]; cc=cc .- minimum(cc); cc=cc ./ maximum(cc);
 cc[diagind(cc)] .= NaN;
heatmap!(pp,subplot=2,cc,aspect_ratio=:equal,c=p2,colorbar=false,grid=false,axis=([], false))
```

```
display(pp)
GR.setarrowsize(0.5);
MS = [1,1,1,2,2,2,3];
colors = [colorant"plum2", colorant"mistyrose1", colorant"lavender"];
A=C[:,:,1];
A[A .== 0] .= 1;
A = A \cdot - \min(A);
A[A .== maximum(A)] .= 0;
A = A . / maximum(A);
A[A .< 0.33] .= 0;
graphplot!(pp,subplot=3,A,
              method=:circular,nodeshape=:circle,
              names=1:7,
              markersize=0.15,
              fontsize=20,
              linewidth=3,
              linealpha=1,
              markercolor = colors[MS],
              nodestrokecolor=colors[MS],
              arrow=arrow(:closed, 10),
display(pp)
```

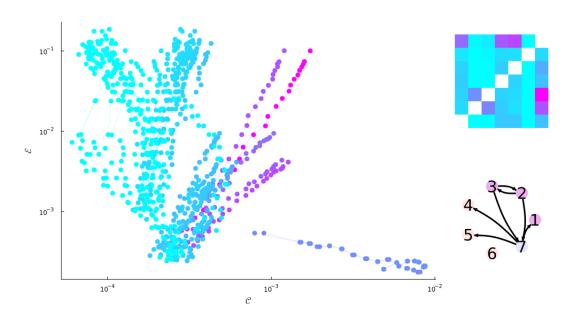


Figure 3: \mathcal{E} over \mathcal{C} ; numerical adjacency matrix and network motif. The colormap is on a log10 scale from cyan (low) to magenta (high). The colors in the adjacency matrix are the same as those in the \mathcal{C} - \mathcal{E} plot on the left (see code above).

We also plot the adjacency matrices for added noise:

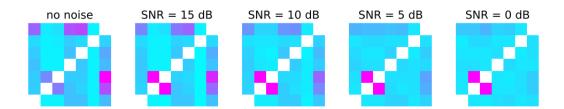


Figure 4: Adjacency matrices for added white noise.

2.2 Unconnected nodes

To change the network we have to change the following parameters in the code

```
FromTo=[];
PF = "__Empty";
```

and get the following results:

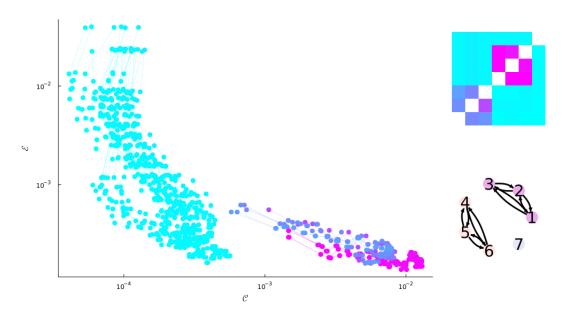


Figure 5: \mathcal{E} over \mathcal{C} ; numerical adjacency matrix and network motif. The colormap is on a log10 scale from cyan (low) to magenta (high). The colors in the adjacency matrix are the same as those in the \mathcal{C} - \mathcal{E} plot on the left (see code).

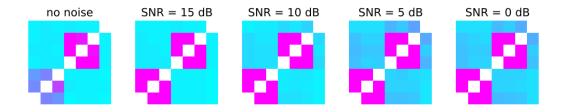


Figure 6: Adjacency matrices for added white noise.

3 Epilepsy

We here use a 4 minute data segment around the second seizure we show in the paper.

```
# julia -p10
@everywhere include("DDAfunctions.jl");
using JLD2
#import Pkg; Pkg.add("Distributed")
using Distributed
FN_part="S05__00_04";
NOISE_list=["NoNoise"; "add15dB"];
FN_NoNoise = @sprintf("%s__%s.ascii",FN_part,NOISE_list[1]);
FN_15dB = @sprintf("%s__%s.ascii",FN_part,NOISE_list[2]);
if !isfile(FN 15dB)
   X=readdlm(FN_NoNoise);
   NrCH=size(X, 2);
   SNR=15;
   Y=fill(NaN, size(X, 1), NrCH);
   for n_ch=1:NrCH
       Y[:, n_ch] = add_noise(X[:, n_ch], SNR);
   end
   fid=open(FN_15dB, "w");
   for k1=1:size(Y, 1)
       for k2=1:NrCH
          @printf(fid, "%13.10f ", Y[k1, k2]);
       @printf(fid, "\n");
   close(fid);
   X = nothing;
   Y = nothing; GC.gc();
end
NrCH=78:
CH=1:NrCH;
ONSET_CH = sort([45:52; 85:92; 61:68; 101:108] .- 30);
TAU=[7 \ 10]; TM = maximum(TAU); dm=4;
WL=500; WS=50;
nr_delays=2;
DDAmodel=[[0 0 0 1];
          [0 0 0 2];
          [1 1 1 1]];
(MODEL, L_AF, DDAorder) = make_MODEL(DDAmodel);
                                                                        # DDA model
DDA_DIR="DDA_Epilepsy"; dir_exist(DDA_DIR);
LIST=reduce(hcat, collect(combinations(CH, 2)))';
DELTA=20;
N=Int64(ceil(size(LIST, 1)/DELTA));
for n_FN=1:length(NOISE_list)
    FN_DATA = @sprintf("%s__%s.ascii",FN_part,NOISE_list[n_FN]);
    FN_DDA = @sprintf("%s%s%s__%s.DDA",DDA_DIR,SL,FN_part,NOISE_list[n_FN]);
    FN_ALL = @sprintf("%s%s%s__%s.jld2",DDA_DIR,SL,FN_part,NOISE_list[n_FN]);
   if !isfile(FN_ALL)
```

```
@printf("%s\n",FN_ALL);
if !isfile(join([FN_DDA,"_ST"]))
   if Sys.iswindows()
      if !isfile("run_DDA_ASCII.exe")
         run(`cp run_DDA_ASCII run_DDA_ASCII.exe`);
      end
      CMD=".\\run_DDA_ASCII.exe";
   else
     CMD="./run_DDA_ASCII";
   end
   CMD = "$CMD -ASCII";
   CMD = "$CMD -MODEL $(join(MODEL," "))"
   CMD = "$CMD -TAU $(join(TAU, " "))"
   CMD = "$CMD -dm $dm -order $DDAorder -nr_tau $nr_delays"
   CMD = "$CMD -DATA_FN $FN_DATA -OUT_FN $FN_DDA"
   CMD = "$CMD -WL $WL -WS $WS";
  CMD = "$CMD -SELECT 1 0 0 0";
                                                                 # ST-DDA
   if Sys.iswindows()
     run(Cmd(string.(split(CMD, " "))));
   else
     run(`sh -c $CMD`);
   rm(@sprintf("%s.info",FN_DDA));
end
@sync @distributed for n_N=1:N
    \label{eq:fndda} FN_DDAn = @sprintf("%s%s%s__%s__%03d.DDA",DDA_DIR,SL,FN_part,NOISE_list[n_FN],n_N); \\
    n=collect(1:DELTA) .+ (n_N-1)*DELTA; n=n[n.<=size(LIST,1)];
    LL1=LIST[n,:]; LL1=vcat(LL1'...)';
    if !isfile(join([FN_DDAn,"_CT"]))
       if Sys.iswindows()
          if !isfile("run_DDA_ASCII.exe")
             run(`cp run_DDA_ASCII run_DDA_ASCII.exe`);
          end
          CMD=".\\run_DDA_ASCII.exe";
       else
          CMD="./run_DDA_ASCII";
       end
       CMD = "$CMD -ASCII";
       CMD = "$CMD -MODEL $(join(MODEL," "))"
       CMD = "$CMD -TAU $(join(TAU, " "))"
       CMD = "$CMD -dm $dm -order $DDAorder -nr_tau $nr_delays"
       CMD = "$CMD -DATA_FN $FN_DATA -OUT_FN $FN_DDAn"
       CMD = "$CMD -WL $WL -WS $WS";
       CMD = "$CMD -SELECT 0 1 0 0"
                                                                # CT-DDA
       CMD = "$CMD -CT_CH_list $(join(LL1," "))";
                                                                # all pairwise channels
       CMD = "$CMD -WL_CT 2 -WS_CT 2";
       if Sys.iswindows()
         run(Cmd(string.(split(CMD, " "))));
       else
          run(`sh -c $CMD`);
       rm(@sprintf("%s.info",FN_DDAn));
    end
end
@sync @distributed for n_N=1:N
   FN_DDAn=@sprintf("%s%s%s__%s__%03d.DDA",DDA_DIR,SL,FN_part,NOISE_list[n_FN],n_N);
  n=collect(1:DELTA) .+ (n_N-1)*DELTA; n=n[n.<=size(LIST,1)];</pre>
```

```
LL1=LIST[n,:]; LL1=vcat(LL1'...)';
    if !isfile(join([FN_DDAn,"_CD_DDA_ST"]))
       if Sys.iswindows()
          if !isfile("run_DDA_ASCII.exe")
             run(`cp run_DDA_ASCII run_DDA_ASCII.exe`);
          end
          CMD=".\\run_DDA_ASCII.exe";
       else
          CMD="./run_DDA_ASCII";
       end
       CMD = "$CMD -ASCII";
       CMD = "$CMD -MODEL $(join(MODEL," "))"
       CMD = "$CMD -TAU $(join(TAU, " "))"
       \texttt{CMD} = \texttt{"$CMD} - \texttt{dm} \texttt{$dm} - \texttt{order $DDAorder -nr\_tau $nr\_delays"}
       CMD = "$CMD -DATA_FN $FN_DATA -OUT_FN $FN_DDAn"
       CMD = "$CMD -WL $WL -WS $WS";
       CMD = "$CMD -SELECT 0 0 1 0";
                                                                   # CD-DDA
       CMD = "$CMD -PAIRS $(join(LL1," "))";
                                                                   # all pairwise channels
       if Sys.iswindows()
          run(Cmd(string.(split(CMD, " "))));
       else
          run(`sh -c $CMD`);
       end
       rm(@sprintf("%s.info",FN_DDAn));
    end
end
######### stitch together outputs
ST=readdlm(join([FN_DDA,"_ST"]));
T=ST[:,1:2]; ST=ST[:,3:end];
WN=size(T, 1);
rhoS=ST[:,L_AF:L_AF:end];
ST = nothing; GC.gc();
E=fill(NaN, WN, NrCH, NrCH);
for n_N=1:N
    @printf("%3d ",n_N)
    FN_DDAn=@sprintf("%s%s%s__%s__%03d.DDA",DDA_DIR,SL,FN_part,NOISE_list[n_FN],n_N);
    n=collect(1:DELTA) .+ (n_N-1)*DELTA; n=n[n.<=size(LIST,1)];</pre>
    LL1=LIST[n,:] .- CH[1] .+ 1;
    CT=readdlm(join([FN_DDAn,"_CT"]));
    CT=CT[:,3:end];
    CT=CT[:,L_AF:L_AF:end];
    for l=1:size(LL1,1)
        ch1=LL1[1,1];ch2=LL1[1,2];
        E[:,ch1,ch2] = abs.( dropdims(mean(rhoS[:,[ch1,ch2]],dims=2),dims=2) ./ CT[:,1] .- 1 );
        E[:, ch2, ch1] = E[:, ch1, ch2];
    end
    CT = nothing; GC.gc();
end
@printf("\n");
C=fill (NaN, WN, NrCH, NrCH);
for n_N=1:N
    @printf("%3d ",n_N)
    FN_DDAn=@sprintf("%s%s%s__%s__%03d.DDA",DDA_DIR,SL,FN_part,NOISE_list[n_FN],n_N);
     n = collect (1:DELTA) .+ (n_N-1) *DELTA; n = n[n. <= size(LIST, 1)]; 
    LL1=LIST[n,:] .- CH[1] .+ 1;
    CD=readdlm(join([FN_DDAn,"_CD_DDA_ST"]));
```

```
CD=CD[:,3:end];
           CD=reshape(CD, WN, 2, size(LL1, 1));
           for l=1:size(LL1,1)
               ch1=LL1[1,1];ch2=LL1[1,2];
               C[:,ch1,ch2] = CD[:,2,1];
               C[:, ch2, ch1] = CD[:, 1, 1];
           end
           CD = nothing; GC.gc();
       end
       @printf("\n\n");
       @save FN_ALL C E rhoS T WN
       E = nothing; C = nothing; GC.gc();
    end
end
######### plot results
n_FN=1;
FN_ALL = @sprintf("%s%s%s__%s.jld2",DDA_DIR,SL,FN_part,NOISE_list[n_FN]);
@load FN_ALL C E T WN ;
C1 = C \cdot \star 1;
E1 = E \cdot \cdot \cdot 1;
E = nothing; C = nothing; GC.gc();
FN_ALL = @sprintf("%s%s%s__%s.jld2",DDA_DIR,SL,FN_part,NOISE_list[n_FN]);
@load FN_ALL C E T WN ;
C2 = C .* 1;
E2 = E \cdot \cdot 1;
E = nothing; C = nothing; GC.gc();
eLABEL=["LFP"; "LCG"; "LAT"; "LMT"; "LHP"; "LOC"; "LTH"; "LSU";
        "RFP"; "RCG"; "RAT"; "RMT"; "RHP"; "ROC"; "RTH"; "RSU"]
e_list = [
       [34:38; 15:22; 25:30],
                                    #LFP
       [31:33; 23; 24],
                                    #LCG
       [3:5; 10:14],
                                     #LAT
       Int[],
       [1:2; 7:9],
                                    #LHP
       Int[],
       Int[],
       Int[],
       [55:62; 71:78; 63; 65:70], #RFP
      Int[],
       [40:44; 50:52],
                                    #RAT
       Int[],
       [39; 47:49],
                                     #RHP
       Int[],
       Int[],
       [64]
                                    #RSU
       ];
e_NotZero = findall(x -> x == 1, length.(e_list)' .!= 0);
e_NotZero = [i[2] for i in e_NotZero];
SR = 500:
t = (T[:,1] .+ 1 .+ TM .+ dm) ./ SR ./ 60;
SEQ=1:length(e_NotZero);
CHs=vcat(e_list[e_NotZero][SEQ][:]...);
L_e_list=map(x -> length(e_list[e_NotZero][x]),SEQ);
IND=setdiff(1:length(CHs)^2, diagind(C1[1, CHs, CHs]));
c1=reshape(C1[:,CHs,CHs],WN,length(CHs)^2)[:,IND];
c2=reshape(C2[:,CHs,CHs],WN,length(CHs)^2)[:,IND];
el=reshape(El[:,CHs,CHs],WN,length(CHs)^2)[:,IND];
e2=reshape(E2[:,CHs,CHs],WN,length(CHs)^2)[:,IND];
```

```
c2[c2 .< 0.01] .= NaN;
c1[c1 .< 0.01] .= NaN;
BETA=asin.((c2 .- c1) ./ sqrt.( (c1 .- c2).^2 + (e1 .- e2).^2) );
SG = plot(size=(1000, 1000), layout=(2, 2));
heatmap! (SG, subplot=1,
        t,1:length(IND),c1',
        c=:jet,
        xtickfont=font(12), ytickfont=font(12),
        colorbar = true, clims=(0.1, 0.06)
Y = vcat(0, cumsum(L_e_list[:]) .* (length(CHs)-1));
Y = (Y .- [0; diff(Y)./2])[2:end];
heatmap!(SG, subplot=1, yticks=(Y, eLABEL[e_NotZero][SEQ]), xlabel="time [min]", clims=(0.01, 0.06))
display(SG)
heatmap! (SG, subplot=2,
        t,1:length(IND),c2',
        c=:jet,
        xtickfont=font(12), ytickfont=font(12),
        colorbar = true, clims=(0.1, 0.06)
hline!(SG, subplot=2, [cumsum(L_e_list[:]) .* (length(CHs)-1) .+ 0.5], legend=false, c=:black, linewidth=2)
\texttt{heatmap!} (\texttt{SG}, \texttt{subplot=2}, \texttt{yticks=(Y,eLABEL[e\_NotZero][SEQ])}, \texttt{xlabel="time [min]"}, \texttt{clims=(0.01,0.06)}) \\
display(SG)
ALPHA = BETA .* 1;
ALPHA[ALPHA .>= 0] .= 0;
ALPHA [ALPHA .< 0] .= 1;
ALPHA[ALPHA .== 0] .= NaN;
heatmap! (SG, subplot=3,
        t,1:length(IND),(c1 .* ALPHA)',
        xtickfont=font(12), ytickfont=font(12),
        colorbar = true, clims = (0.1, 0.06)
heatmap!(SG, subplot=3, yticks=(Y, eLABEL[e_NotZero][SEQ]), xlabel="time [min]", clims=(0.01, 0.06))
display(SG)
ALPHA = BETA .* 1;
ALPHA[ALPHA \cdot <= 0] \cdot = 0;
ALPHA [ALPHA .> 0] .= 1;
ALPHA[ALPHA .== 0] .= NaN;
heatmap! (SG, subplot=4,
        t,1:length(IND),(c1 .* ALPHA)',
        c=:jet,
        xtickfont=font(12), ytickfont=font(12),
        colorbar = true, clims = (0.1, 0.06)
hline!(SG, subplot=4, [cumsum(L_e_list[:]) .* (length(CHs)-1) .+ 0.5], legend=false, c=:black, linewidth=2)
heatmap!(SG, subplot=4, yticks=(Y, eLABEL[e_NotZero][SEQ]), xlabel="time [min]", clims=(0.01, 0.06))
display(SG)
```

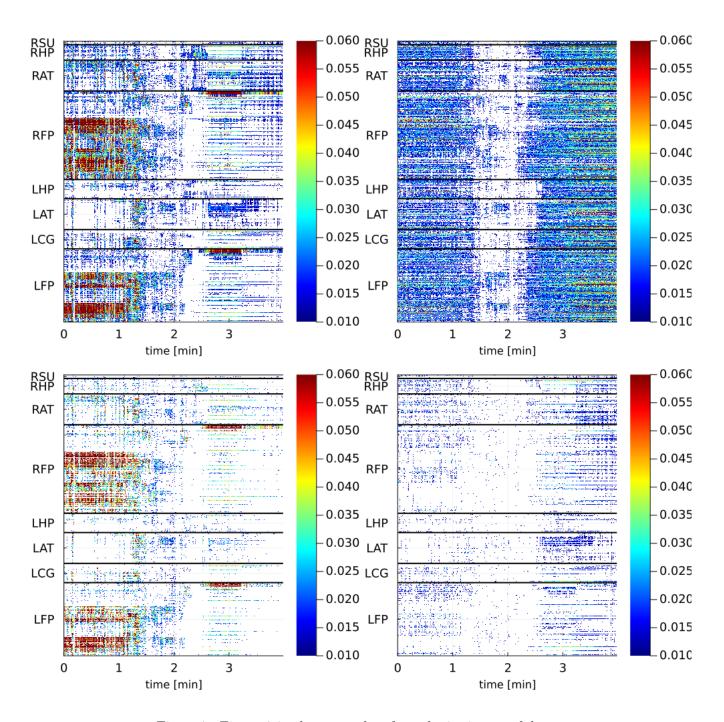


Figure 7: Figure 8 in the paper, but for only 4 minutes of data.