This is a guide for batch processing all tetrodes with mountainsort pipeline and converting the output to LFPO-readable form for post-processing. The mountainsort pipeline is based on instructions in Mountain Sort 25/10/19

Relevant Python/Bash scripts are located in /d1/data/bin/igor, which is normally part of the data user's PATH variable. If you run this pipeline as a different user, make the scripts available from your PATH. You need ms_run_and_adapt.sh, ms_lfpo_proc_tet.sh, ms_merge_clures.py, sgclust_to_lfpo.py, mount_sortb.sh, mountain_to_sgclustb.py, axtrk_to_whl.py, Dat_to_Mda.py, ws_interpolate, dat_to_mdab.sh.

Execute the following commands in the working directory containing dat and axtrk files.

- 1. Copy relevant files to working directory:
 - cp /d1/data/lfpo1/ms_and_lfpo/* .
- 2. Remove unwanted channels from *TEMPLATE.par*
- 3. Create a file named **BASELIST** with a list of basenames (1 per line) of *dat/axtrk* to be processed, in chronological order e.g.:

```
jc103-3005_01l
jc103-3005_02s
Jc103-3005_03l
```

To create one for all *dat* files in working dir, run (double check session order):

for f in *.dat; do printf '%s\n' "\${f%.dat}"; done > BASELIST

- 4. Run master script with following parameters: ANIMAL, DAY, DTHOLD (spike detection threshold for mountainsort), NCHAN (64 or 128) e.g.:
 ms run and adapt.sh 5401 07102020 5 34
- 6. [Optional] Perform the semi-automated cleaning steps (see below)
- 7. [Optional] Open clusters from all tetrodes together:
 - a. Merge clu/res: ms merge tetrodes.sh jc103 3005
 - b. [OPTIONAL] Adjust pf.sessions/pf.groups if you want to build rate maps
 - c. Open with config spike display mountainsort alltet.conf
- 8. [Optional] Adapt output to Jozsef's dataset format:
 - a. ms_to_sgclust.py sgclust/DRCCK15-11032018 ./ 0
 Parametres are: 1. basename for output files 2. Directory
 containing tet*/
 - 3. 1 if downsampling 24kHz->20kHz is needed
 - b. clures_merge_tetrodes_by_sessions.py DRCCK15-11032018-1724

 Run this in the output directory of previous step (sgclust/ in example above)
- Template display config file **spike_display_mountainsort_TEMPLATE.conf** is in /d1/data/lfpo1/Res/. It also relies on tetrode config files tetr_1tetr_* that are in /d1/data/lfpo1/Res/tetr/

• For post-processing, the most useful feature is displaying spikes that contaminate the autocorrelogram, which can be triggered with 'c'.

KNOWN ISSUES

- Pipeline doesn't work with symbolic links to *.dat
- Might not work with a single tetrode add a dummy tetrode in TEMPLATE.par to get over
- If a whole tetrode is broken, in the tetr configuration file just remove the line that contains the channels (instead of writing "0 [channels]")

SEMI-AUTOMATED CLUSTER CLEANING

By default, all scripts are run in the day directory (the one containing tet*/ directories)

This pipeline assumes you have first executed the mountainsort pipeline (see above).

- 1. [Pre-clean in lfp_online] Quickly go through all clusters obvious splits/merges, delete hopeless clusters
- 2. [FIT] Run cleaning script in the Fit mode nohup autoclean_alltet.sh &
- 3. [CLEAN] When fit ruIT mode takes few minutes per cluster: n is over, run clearing script in CLEAN mode:

run autoclean.sh

For every cluster, choose % of dirtiest spikes to be removed based on cleaning curve - click with middle mouse button or just close window to ignore the cluster.

Results are written to path with added 'clean' suffix after all clusters in the current tetrode have been cleaned.

- 4. [Replace clu/res] This will simply backup pre-clean clu/res and replace then with cleaned ones: copy_preclean.sh
- 5. [Post-clean] Open results in **lfp_online** to merge clusters and clean cross-contaminations.
- 6. [OPT] To redo cleaning, first restore backup from step 2 and run script in CLEAN mode

RUNNING SEMI-AUTOMATED CLUSTER CLEANING ON DIFFERENT MACHINE

Since FIT run takes time and resources, you can run it on a different machine. For a FIT run you only need fet, clu and res files. For a CLEAN run you then need the files in AUTOCLEAN/ directory that is created in every tet*/ directory. For example, in case of running mountainsort on pyramidal, doing pre-cleaning on home computer and want to run cleaning in FIT mode on mec, you need these commands to sync files:
on mec, in animal data directory

rsync -avP --include="*/" --include="*.fet.0" --exclude="*" igor@pyramidal:/hdr/data/bindata/\$ANIMAL/\$DAY

•

rsync -avP igor@pyramidal:/hdr/data/bindata/\$ANIMAL/\$DAY/TEMPLATE.par \$DAY/# on home, after pre-clean step
rsync -avP --include="*/" --include="*.clu" --include="*.res" --exclude="*" tet*
igor@mec:/temp_store/igor/\$ANIMAL/\$DAY

After FIT that, copy AUTOCLEAN/ directories back to *home* by running: # on home, in day directory (with tet*/ directories)

rsync_autoclean.sh igor@pyramidal:/hdr/data/bindata/\$ANIMAL/\$DAY

If you also do pre/post-cleaning on a machine different from the one where mountainsort has been running, you can use this command to avoid copying large spk files (it will interpolate the original waveshapes locally): rsync_msout.sh igor@pyramidal:/hdr/data/bindata/\$ANIMAL/\$DAY; run_ws_interporalte.sh

FEATURE REQUESTS FOR LFPO [OPTIONALLY]

[OPTIONALLY] - assign priority on scale 1-5

- (5) Add the ctrl+z or some command for undoing merging
 - o Implemented but needs testing
- In relation to the above, maybe also how many spikes fall within the refractory period (maybe when you press c to show them); so one can decide if they are a low percentage of the total...
 - Now can be done with clu_refractory_fraction.py (/d1/data/bin/igor)
- A way to see the inter spike interval of a specific cluster in more detail
 with a scale bar (e.g. so one can tell how wide the refractory period of a
 cell actually is, in ms)

HOW TO GET LATEST CLUSTERING SOFTWARE

If you don't have code, do git clone https://github.com/igridchyn/lfp_online.git

To update: git pull origin master

To compile: (while in lfp online/sdl example/Debug) make Ifp online

FOR YOAV

Merge whl before running ms_run_and_adapt.sh

It is not crucial to do it before, but it is easier.

Before merge, generate session_shifts.txt file. The DIV variable here should be equal <NUMBER OF CHANNELS X 2>, change it if number of channels is different from 34

DIV=68; ls -lU `awk '{print \$1".dat"}' BASELIST` | awk '{sum+=\$5/'\$DIV'; \$5=sum ; print \$5}' > session shifts.txt

merge_whl.py session_shifts.txt 4201-04082020.whl 4201-04082020_01.whl 4201-04082020_02.whl 4201-04082020_03.whl 4201-04082020_04.whl 4201-04082020_05.whl 4201-04082020_06.whl 4201-04082020_07.whl

For positrack, convert *.positrack files to whl:

positrack to whl.py 4201-04082020 01.positrack <SHIFT>

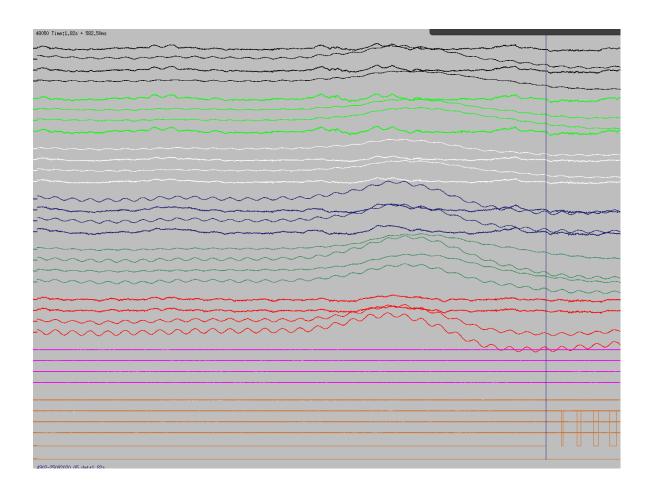
 $\label{thm:continuous} \mbox{Where} \mbox{\ensuremath{\mathsf{SHIFT}}\xspace} \mbox{\ensuremath{\mathsf{is}}} \mbox{\ensuremath{\mathsf{timestamp}}} \mbox{\ensuremath{\mathsf{of}}} \mbox{\ensuremath{\mathsf{the}}} \mbox{\ensuremath{\mathsf{first}}} \mbox{\ensuremath{\mathsf{tracking}}} \mbox{\ensuremath{\mathsf{TTL}}} \mbox{\ensuremath{\mathsf{sync}}} \mbox{\ensuremath{\mathsf{signal}}}, \\ \mbox{\ensuremath{\mathsf{signal}}} \mbox{\ensuremath{\mathsf{e}}} \mbox{$

which can be seen in **regaamc11** by putting pointer on it: (number is in the top left corner, pointer is invoked by 'p')

Command to run regaamc11:

regaamc11 4201-04082020 01.dat 34 50 ../regaa 34 chan.conf

The config file can be found in one of the data folders or in /d1/data/lfpol.



Copy clustering output to ca3. Change IP address if needed.

rsync -avP --exclude='*.mda' tet* guest03@10.4.44.175:/mnt/data_yoav/data/6001/02112020

Run lfp_online to do clustering

./lfp_online ../Res/spike_display_mountainsort_ca3.conf channel.num=64 animal=4001 day=25072020 tet=3 nchan=4

```
If d1 is not mounted ('ls /d1' shows empty folder), mount it with sudo mount alpha2:/d1 /d1

cd ~/lfp_online/sdl_example/Debug/

ms_to_sgclust.py sgclust/4001-24072020 ./ 0

cp ../4001-24072020.par .

clures_merge_tetrodes_by_sessions.py 4001-24072020

5.- Generate .light files for every session

$ dat2timestamps.py 4201-04082020_02.dat 34 33 20 4201-04082020_02.light
```