# MCWrapper and Monte Carlo production





#### GlueX simulation chain

#### 1. Event generation

\* Your favourite event generator (gen\_amp, genr8, bggen, ...)

#### 2. Geant4 simulation

\* Use a Geant4 implementation of GlueX (hdgeant4) to simulate interactions inside the GlueX spectrometer

#### 3. Smearing

\* Correct for deficiencies in simulation ("make it look more like data")

treat like data

- 4. Reconstruction
- 5. Analysis (ReactionFilter)

### **MCW**rapper

- Most people doing simulations need to perform all these steps
  - Each step needs to be set up correctly
  - \* Environment variables, config files, ...
  - Time consuming and prone to errors
- \* Solution: Write some code to do most of the work for you MCWrapper, a one-stop shop for MC production
- \* A few scripts that help to generate MC and are controlled by config files
- It is not almighty: garbage in garbage out



## **MCW**rapper

- \* Two ways to generate MC with MCWrapper:
  - Locally



- Write a config file that controls MCWrapper
- Run either in shell or submit to batch farm

- \* OSG
  - \* Use a website to generate config file
  - \* Submit to the open science grid



 Recommended software setup: setenv MCWRAPPER\_CENTRAL /scigroup/mcwrapper/gluex\_MCwrapper/

# MCWrapper - locally

#### Usage:

```
ifarm1901.jlab.org> $MCWRAPPER CENTRAL/gluex MC.py
Usage: gluex MC.py config file Run Number/Range num events [all other options]
 where [all other options] are:
 variation=%s where %s is a valid jana calib context variation string (default is "mc")
 per file=%i where %i is the number of events you want per file/job (default is 10000)
 base file number=%i where %i is the starting number of the files/jobs (default is 0)
 numthreads=%i sets the number of threads to use to %i. Note that this will overwrite the NCORES set in MC.config
 generate=[0/1] where 0 means that the generation step and any subsequent step will not run (default is 1)
 geant=[0/1] where 0 means that the geant step and any subsequent step will not run (default is 1)
 mcsmear=[0/1] where 0 means that the mcsmear step and any subsequent step will not run (default is 1)
 recon=[0/1] where 0 means that the reconstruction step will not run (default is 1)
 cleangenerate=[0/1] where 0 means that the generation step will not be cleaned up after use (default is 1)
 cleangeant=[0/1] where 0 means that the geant step will not be cleaned up after use (default is 1)
 cleanmcsmear=[0/1] where 0 means that the mcsmear step will not be cleaned up after use (default is 1)
 cleanrecon=[0/1]where 0 means that the reconstruction step will not be cleaned up after running (default is 0)
 batch=[0/1/2] where 1 means that jobs will be submitted, 2 will do the same as 1 but also run the workflow in the case
of swif(2) (default is 0 [interactive])
logdir=[path] will direct the .out and .err files to the specified path for qsub
Options:
  -h, --help show this help message and exit
```

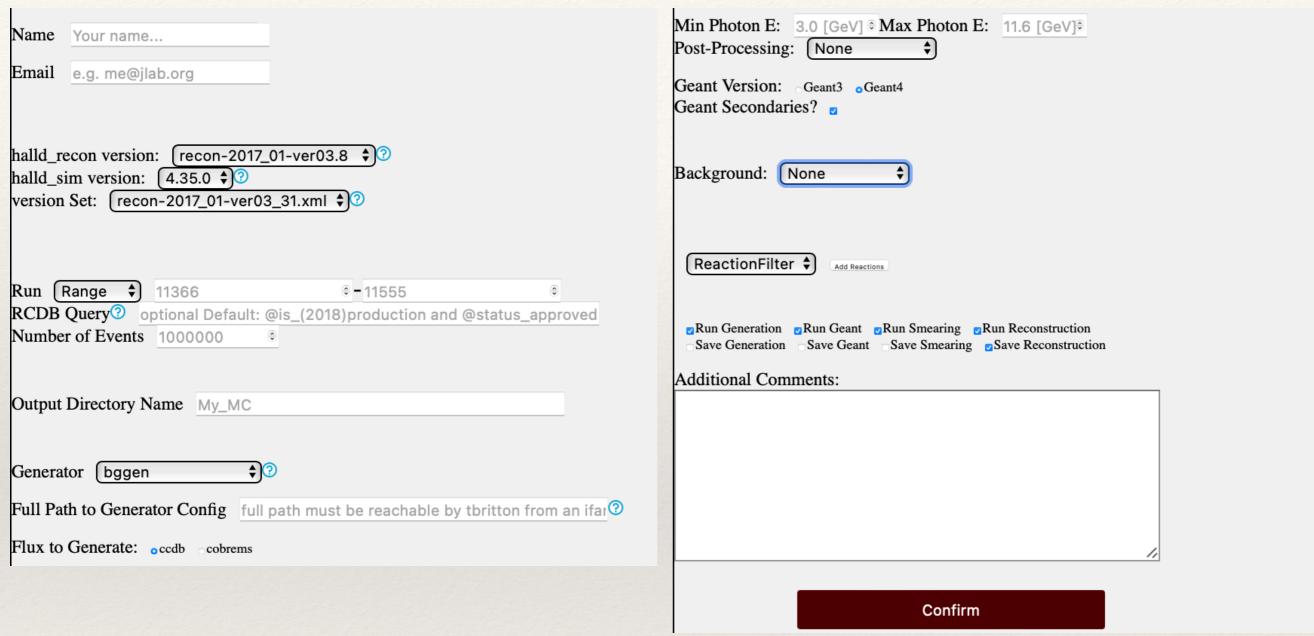
#### Need to provide config file

Example:

https://github.com/JeffersonLab/gluex\_MCwrapper/blob/master/examples/MC.config

### MCWrapper - OSG

https://halldweb.jlab.org/gluex\_sim/SubmitSim.html



- Drop-downs and text boxes to set up MC submission
- Click ? to access documentation
- Automatically tested and submitted to Open Science Grid

## MCWrapper - OSG

https://halldweb.jlab.org/gluex\_sim/Dashboard.html

- Dashboard lets you monitor progress (new improved performance!)
- \* Status:
  - Red = test no successful (check your emails!)
  - Green = tested successful, project running
- Left click more infos, right click options (cancel, declare complete)

Projects ? Show 10 entries										
Progress *	ID •	Email 💠	Status \$	RunNumLow \$	RunNumHigh \$	NumEvents \$	Generator \$	BKG \$	R	
2.7	2520	nwickjlb@jlab.org		50685	51768	250000000	gen_amp	Random:recon- 2018_08- ver02.2	@is_ @s bean	
0	2519	churaman@jlab.org		40856	42559	10000000	genEtaRegge	None		
30.79	2518	churaman@jlab.org		40856	42559	10000000	genEtaRegge	None		
29.41	2517	churaman@jlab.org		40856	42559	10000000	genEtaRegge	None		
0	2516	churaman@jlab.org		40856	42559	10000000	genEtaRegge	None		

## MCWrapper - OSG

https://halldweb.jlab.org/gluex\_sim/Records.html

- \* Records show all previously produced MC, worth checking before submitting new
- Contains all important information about projects

Projects Refresh tables Show 10 © entries Search:									
ID v	Email \$	user_id \$	Exp \$	Submit_Time \$	Tested \$	Is_Dispatched \$	Dispatched_Time \$	Completed_Time \$	RunNumLow \$
2510	zbaldwin@andrew.cmu.edu	51	GlueX	2022-05-11 08:53:40	1	1.0	2022-05-13 10:53:42		30274
2509	zbaldwin@andrew.cmu.edu	51	GlueX	2022-05-11 08:14:37	1	1.0	2022-05-13 10:53:33		30274
2508	zbaldwin@andrew.cmu.edu	51	GlueX	2022-05-09 14:22:00	4	1.0	2022-05-09 21:57:48	2022-05-11 09:09:26	30274
2506	gabyrod7@gmail.com	29	GlueX	2022-05-06 15:44:55	1	1.0	2022-05-09 21:55:37		30274
2505	jzarling@jlab.org	20	GlueX	2022-05-05 17:43:51	1	1.0	2022-05-09 21:55:14		51384
2504	jzarling@jlab.org	20	GlueX	2022-05-05 17:42:26	1	1.0	2022-05-09 21:56:26		51384

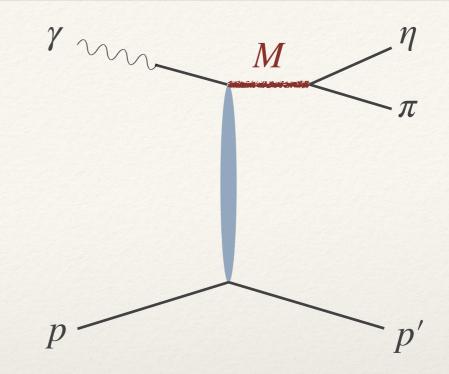
## Event generator

- Generic background samples:
  - \* bggen
- \* Simple phase space:
  - \* genr8
- Sophisticated amplitude models:
  - \* gen\_amp
  - \* gen\_vec\_ps
- For studies:
  - particle gun
  - geantBEAM

Decaying particles can either decay in hdgeant4 or you can specify the decay through generator post-processing, e.g. decay\_evtgen

#### Requirements

- We need flat MC to perform fits for PWA
- Use gen\_amp for both
  - First line in config file
     (commented out) contains
     switches for MCWrapper
  - Simple switch f to generate flat MC
- \* Force  $\eta \to \gamma \gamma$
- Generate for 2017-01 analysis ver52



```
session2b > ♣ EtaPiOFlat.cfg

1 # -f -u 3.0 -a 8.2 -b 8.8

2 reaction EtaPi0 Beam Proton Eta Pi0

3
```

```
session2b > ≡ EtaDecay.dec

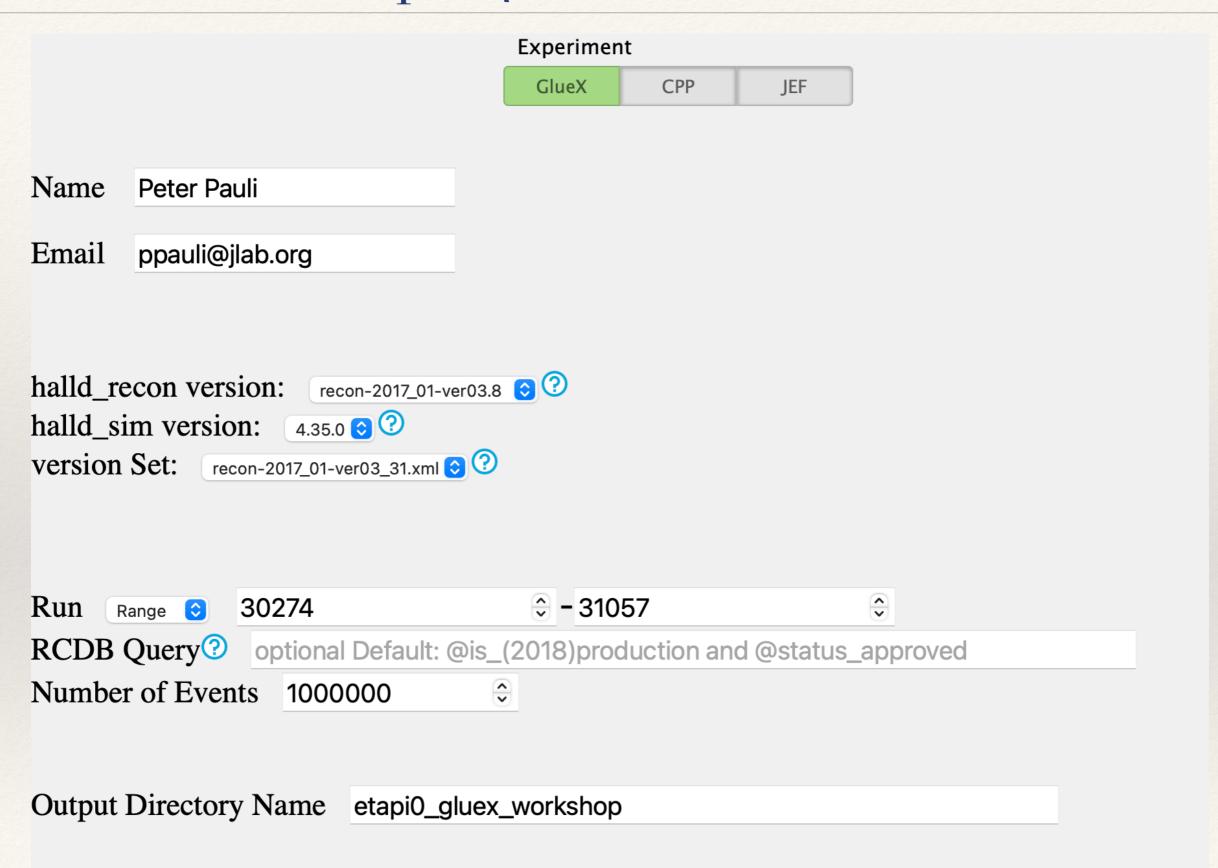
1 Decay eta
2 1.000 gamma gamma PHSP;
3 Enddecay
4
5 End
6
```

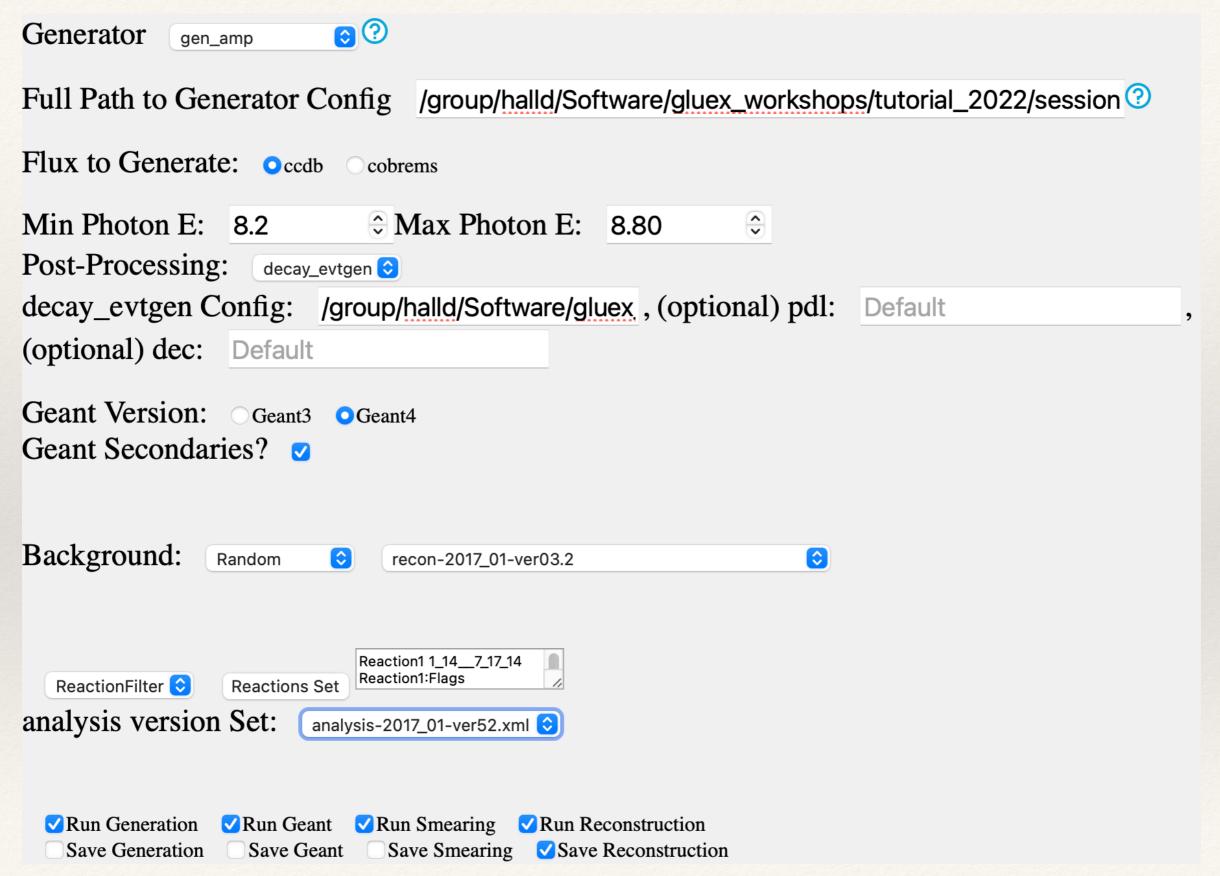
### Example: $\eta \pi$ simulations (locally)

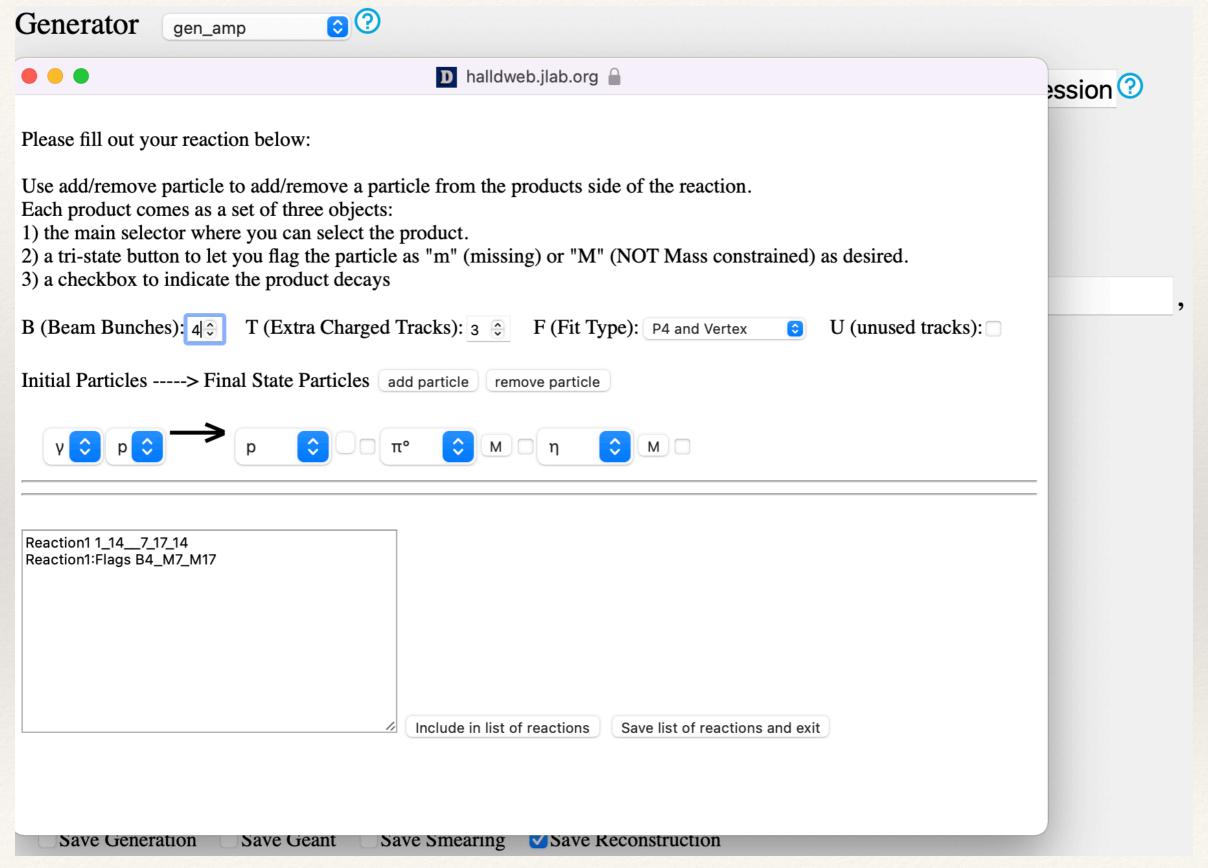
#### Example:

https://github.com/JeffersonLab/gluex\_MCwrapper/blob/master/examples/MC.config

```
session2b > A MC.config
      DATA_OUTPUT_BASE_DIR=/volatile/halld/home/ppauli/gluex_workshops/tutorial_2022/session2b/etapi_out/ #your desired output location
                    # Number of CPU threads to use or nodes:node-id:ppn or nodes:ppn depending on your system
      GENERATOR=gen amp #or you may specifile file:/.../file-to-use.hddm
      GENERATOR_CONFIG=/group/halld/Software/gluex_workshops/tutorial_2022/session2b/EtaPi0Flat.cfg
      GENERATOR_POSTPROCESS=decay_evtgen:/group/halld/Software/gluex_workshops/tutorial_2022/session2b/EtaDecay.dec #:pathToEVT.PDL:pathToDECAY.DEC # last two are optional
      #common parameters for generators
      GEN_MIN_ENERGY=8.2
      GEN_MAX_ENERGY=8.8
     GEANT_VERSION=4
      BKG=Random:recon-2017_01-ver03.2 #Use random trigger background for 2017-01
      ANA_ENVIRONMENT_FILE=/group/halld/www/halldweb/html/halld_versions/analysis-2017_01-ver52.xml #optional either a .(c)sh file to be sourced or .xml before the below pl
      #optional additional plugins that will be run along side danarest and hd_root. This should be a comma separated list (e.g. plugin1,plugin2)
      CUSTOM_PLUGINS=file:/group/halld/Software/gluex_workshops/tutorial_2022/session2b/jana_analysis.config #or file:/.../file-to-use which is a configuration file for jar
      #EVERYTHING BELOW FOR BATCH ONLY
      #VERBOSE=True
      BATCH_SYSTEM=swif2 #can be swif or swif2 or condor or osg or qsub adding : [name] will pass -q [name] into PBS.
      ENVIRONMENT_FILE=version.xml
                                       #change this to your own environment file
      WORKFLOW_NAME=FlatEtaPi0Tut22 #SWIF WORKFLOW NAME
 30 PROJECT = gluex
                               # http://scicomp.jlab.org/scicomp/#/projects
                               # https://scicomp.jlab.org/docs/batch_job_tracks
 31 TRACK= simulation
 32 # for swif2
 33 ACCOUNT = halld
                               # https://scicomp.jlab.org/scicomp/slurmJob/slurmAccount
                                    # https://scicomp.jlab.org/scicomp/slurmJob/slurmInfo
      PARTITION = production
      EXPERIMENT = GlueX
                               # GlueX (default) or CPP
     # RESOURCES for swif(2) jobs
 38 DISK=5GB
 39 RAM=5GB
                         # Max RAM usage
 40 TIMELIMIT=600minutes
                                # Max walltime. This may be of the form xx:xx:xx depending on your system
 41 OS=general
                        # Specify CentOS65 machines
```







Geant Version: ○Geant3 ○Geant4 Geant Secondaries? ✓							
Background: Random Corecon-2017_01-ver03.2							
Reaction 1 1_147_17_14 Reaction 1 Set:  Reaction 1 1_147_17_14 Reaction 1:Flags  analysis version Set:  analysis-2017_01-ver52.xml ©							
☑Run Generation ☑Run Geant ☑Run Smearing ☑Run Reconstruction ☑Save Generation ☑Save Geant ☑Save Smearing ☑Save Reconstruction Additional Comments:							
Confirm							

### Summary and Resources

- MCWrapper is a collection of scripts (not compiled!) that aids in producing Monte Carlo samples
- Use it locally (with config file) or submit to OSG
- \* Documentation:
  - \* <a href="https://github.com/JeffersonLab/gluex\_MCwrapper">https://github.com/JeffersonLab/gluex\_MCwrapper</a>
  - https://www.overleaf.com/read/bqynmnwstzfx
  - https://halldweb.jlab.org/wiki/index.php/Event\_generators
  - https://halldweb.jlab.org/wiki/index.php/
     HOWTO\_Use\_EvtGen\_to\_simulate\_particle\_decays\_in\_GlueX
  - https://halldweb.jlab.org/wiki/index.php/
     How to choose software versions on the MC submission form
- Need help? Email <u>ppauli@jlab.org</u> and <u>tbritton@jlab.org</u>