





### **Tutorial session: running EPOS and RIVET**

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# Program of this tutorial

In this tutorial, we will work with **EPOS 3.259** (the last tuned version of EPOS 3<sup>1</sup>) and **RIVET 3.1.5** (last version of RIVET<sup>2</sup> released until now).

#### What you will do:

- write files to parametrise  $e^+e^-$  and pp/pA/AA simulations with EPOS
- run simulations with EPOS
- see the content of the main output formats from EPOS
- use a (provided) HepMC data file from EPOS to analyse it with RIVET
- produce plots with RIVET

#### What you won't do:

- produce HepMC data file with EPOS (impossible with EPOS 3)
- use EPOS 4 3 (not publicly available yet)

All the instructions related to Docker will be written in grey, so that there's no confusion if you want to follow this tutorial with EPOS installed on you computer.

**IMPORTANT**: all words in blue in these slides are hyperlinks to the corresponding webpages

<sup>&</sup>lt;sup>1</sup>K. Werner et al., *Phys. Rev. C* **89** (2014), 064903

<sup>&</sup>lt;sup>2</sup>C. Bierlich et al., SciPost Phys. 8 (2020), 026

<sup>&</sup>lt;sup>3</sup>K. Werner, et al., *Phys. Atom. Nucl.* **84** (2021), 1026–1029

## Prerequisites

For simplicity, we will use **Docker** to run EPOS in a "container".

This way, you will benefit from a perfectly functional portable version, avoiding the necessity to install EPOS natively, with all its dependencies, on your computer.

Here are then the first necessary steps to follow, before to start this tutorial:

- Install Docker (for Linux, Windows or Mac) by following instructions from: https://docs.docker.com/engine/install/
- Download the EPOS 3 image and EPOS\_pp7T\_hepmc.tgz from : https://box.in2p3.fr/index.php/s/CgzXkqEDJPogcDc

You can then load this image, and check it worked (epos3259 should appear), with:

- > sudo docker load --input epos3259\_latest.tar.gz
- > sudo docker images

N.B.: normally, you would need to specify "sudo" before using any Docker command; to get rid of this, do:

- > sudo groupadd docker
- > sudo usermod -aG docker \$USER
- > newgrp docker (for Linux; otherwise disconnect+reconnect your session to apply changes)

(more information here : https://docs.docker.com/engine/install/linux-postinstall/)

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#### What is EPOS?

Event generators are programs made to compute models in order to simulate every step of a collision (e.g. EPOS, PYTHIA <sup>4</sup>...).

Advantages: - perfect detector, as final-state particles are all listed (no uncertainties)
- dynamical approach

(indeed, there's always a **shadow** in **the picture**: one has to be careful on the applicability, and phenomenological approaches generally requires parametrisation)

Energy conserving quantum mechanical approach, based on

Partons, parton ladders, strings,

Off-shell remnants, and

Saturation of parton ladders

Multi-purpose event generator based on parton-based Gribov-Regge Theory <sup>5</sup>. Can simulate with the same formalism any type of high-energy collision consistently:

$$e^{-} + e^{+}$$

$$e^- + p$$

$$q + q$$

$$p + A$$

$$A + A$$

N.B.: more details and references about EPOS in my talk on EPOS (Quarkonia As Tools 2022)

<sup>&</sup>lt;sup>4</sup>T. Sjöstrand et al., *Comput. Phys. Commun.* **191** (2015) 159-177

<sup>&</sup>lt;sup>5</sup>K. Werner et al., *Phys. Rep.* **350** (2001) 93-289

#### Simulation of $e^+e^-$ events with EPOS

To run simulations with EPOS, we need first to define a **.optns** file, containing all the information about the system.

First of all, create a dedicated directory in your folders (I will call it /Tuto\_EPOS).

Then, create a file "ee91.optns", and write in (comments are indeed not mandatory):

```
1 application ee
2 set engy 91.2 ! Center-of-mass energy of the collision (GeV)
3 set nevent 10 ! Number of events simulated
4 set modsho 1 ! Print a message every 'modsho' event(s) simulated
5 nodecays 130 -130 -20 end
6 ! Block the decays of particles (K+ K- and Klong here)
7 print * 2 ! Print the events history in a .check file
```

WARNING: never use tabs in .optns files!

To link your directory /Tuto\_EPOS to the one in the Docker image:

It will open a new shell, in which you can now execute (like if you had EPOS installed natively):

```
> ${EPO}epos ee91
```

### .check output

You should now have in your repertory /Tuto\_EPOS a new file called z-ee91.check.

```
        ior
        jor
        i
        ifr1
        ifr2
        id ist ity
        pt
        m
        E
        y

        0
        0
        1
        0
        0
        12
        1
        0
        0.000E+00
        0.511E-03
        0.456E+02
        0.121E+02

        0
        0
        2
        0
        0
        -12
        1
        0
        0.000E+00
        0.511E-03
        0.456E+02
        .121E+02

        1
        2
        3
        4
        9
        10
        1
        0
        0.000E+00
        0.912E+02
        0.912E+02
        0.000E+00

        3
        4
        4
        10
        0
        4
        21
        30
        0.442E+01
        0.000E+00
        0.242E+02
        .238E+01

        3
        4
        6
        10
        0
        9
        21
        30
        0.4373E+01
        0.000E+00
        0.175E+02
        .210E+01

        3
        -4
        7
        10
        0
        9
        21
        30
        0.173E+01
        0.000E+00
        0.750E+01
        .138E+01

        3
        -9
        8
        10
        0
        9
```

"ior"/"jor": parents of particle "i"
"ifr1"/"ifr2": children of particle "i"

You can clearly see the mechanism happening here : the colliding  $e^-$  and  $e^+$  annihilate into a photon, which is giving birth to a chain of partons  $c-g-...-g-\overline{c}$  forming a string.

N.B.: a lot of particle IDs in EPOS differ from PDG; look at idt.dt

### Simulation of pp/pA/AA events with EPOS

To run simulations for any hadronic system, you need a new file (called pp7T.optns here):

```
oplication hadron
set ecms 7000
                             ! Center-of-mass energy of the collision (GeV)
                            ! Atomic number Z of projectile
set laproj 1
                            ! Mass number A of projectile
set maproi 1
set latarg 1
                            ! Atomic number Z of target
set matarq 1
                            ! Mass number A of target
set ninicon 1
                            ! Number of initial conditions used for hydro
core off
                            ! Core-corona procedure off
hvdro x3ffoff
                            ! Hvdrodvnamics off
hacas off
                            ! Hadronic cascade (UrQMD) off
                            ! Number of full hydrodynamical simulation achieved
set nfull 10
set nfreeze 1
                            ! Number of freeze-out events per full hydro event
set modsho 1
                            ! Print a message every 'modsho' event(s) simulated
xinput KWt/icl70.optns
                            ! Input file containing centrality definitions
set centrality 0
                            ! Define centrality class (0 = Min.Bias.)
nodecays -20 end
                       ! Block the decays of particles (Klong here)
                            ! Record the events in Root files
```

WARNING: again, no tabs in .optns files!

If you want to activate hydro evolution of the core + final-state hadronic cascades : (hydro not recommended now, it would take a LOT of time)

```
core full ! Core-corona procedure ON
hydro x3ff ! Hydrodynamics ON
hacas full ! Hadronic cascade (UrQMD) ON
```

### **ROOT** output

In order to produce a **ROOT file** containing the simulated events (based on ROOT 5.8, you have to add the flag "-root" before the name of the file to execute:

```
> ${EPO}epos -root pp7T
```

Here is the list of variables stored in the Trees of ROOT files produced with EPOS 3:

```
versn = new int ;
                         // EPOS version number
                         // atomic number target
                         // mass number target
                         // number of particles in the event
                         // impact parameter (usually; other choices are possible)
us = new float [nmax];
x = new float [nmax]:
v = new float [nmax]:
z = new float [nmax];
                         // energy of particle
= new float [nmax];
= new float [nmax]:
                         // v component of formation point
= new float [nmax]:
= new float [nmax];
                         // particle id (see file "KWt/idt.dt")
d = new int [nmax]:
                         // particle status (hadron last generation (0) or not (1); other numbers refer to partons, Pomerons, etc)
                         // type of particle origin (20-29 from soft strings, 30-39 from hard strings, 40-59 from remnants, 60 from fluid
tv = new int [nmax]:
                         // index of mother (mother needed for exemple for strings; partons between ior and jor constitute the string)
```

#### And the list of variables added in EPOS 4:

```
signot = new float; // corresponding pp cross-section
whard = new int; // number of elementary hard parton-parton scatterings (set to 0)
npartproj = new int; // number of projectile's nucleons participants according to Glauber
npartarg = new int; // number of target's nucleons participants according to Glauber
nspecp = new int; // number of spectators protons according to Glauber
nspec = new int; // number of spectators returns according to Glauber
nspec = new int; // number of spectators returns according to Glauber
```

### HepMC output

With EPOS 4, it is possible to produce ASCII HepMC files (based on HepMC 2.06.09 below which contain some part of the simulated events, making it compatible with RIVET.

To produce such files with **EPOS 4**, you shall add "set ihepmc 1" in the .optns file and add the flag "-hepmc" before the name of the file to execute.

Detailed information about the structure of such files are given in HepMC 2 user manual (Sec.6).

All the details about the data recorded in EPOS .hepmc files (especially for heavy-ion collisions) are given in HepMC.txt (might evolve with later developments of EPOS 4).

In these files, we record the target + projectile, the final-state particles, as well as their parents (+grand-parents) when they decay with a lifetime  $\tau > 10^{-20} s$  (+ some exceptions, i.e.  $J/\Psi$ ,  $\Psi(2S/3770)$  and  $\Upsilon(1/2/3/4S)$ ).

⇒ enables to identify EM/weak feed-down contributions

N.B.: feature NOT available here as we use EPOS 3

<sup>&</sup>lt;sup>6</sup>https://hepmc.web.cern.ch/hepmc/

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## Prerequisites

For this tutorial, we will use the latest version released of RIVET 7, version 3.1.5.

To download a Docker image of RIVET version 3.1.5 (work the same for any version): > docker pull hepstore/rivet:3.1.5

Then, in order to link your /**Tuto\_EPOS** directory to the one in the RIVET image (like we did for EPOS):

Finally, you can test that RIVET works properly from the new shell opened with :

```
> rivet --version
```

N.B.: otherwise, you can define aliases in order to work from your directory, without opening a new shell, by following instructions from: https://gitlab.com/hepcedar/rivet/-/blob/release-3-1-x/doc/tutorials/docker.md

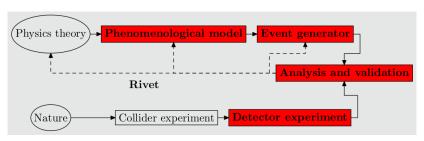
<sup>&</sup>lt;sup>7</sup>https://gitlab.com/hepcedar/rivet

## Robust Independant Validation of Experiment and Theory

RIVET is a "software" based on C++ libraries, installed with different packages :

- YODA: Python libraries and classes used for analyses and histogramming
- HepMC: simulations recording and reading for analyses
- FastJet: recombination algorithms, mainly used for jet analyses

<u>Purpose</u>: offer a simple and standardised tool to automatise comparison between event generators simulations and experimental data



(C. Bierlich, 2019)



# Robust Independant Validation of Experiment and Theory

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<u>Purpose</u>: offer a simple and standardised tool to automatise comparison between event generators simulations and experimental data

RIVET contains many analyses based on publications from many different experiments (data included), and develops thanks to contributions from the users community (from both experiments & theory).

#### Advantages:

- provides huge and constantly growing library of data and analyses
- easy to handle (a lot of documentation + helpful reactive developers)
- don't have to "think about" the analysis details anymore

⇒ RIVET is a very useful tool for us!

Deferences:

### Composition of an analysis

#### Each analysis composed of 4 files:

```
#include "AisotTrols/Percentile.bh" 
#include vospies.
#include vospies.
#include stringe
namespace Rivet {

class STM_ES_Centrality: public Single/alueProjection(

#INCLUDE STM_ES_CENTRALITY: public Single/alueProjec
```

#### ANALYSIS\_NAME.cc (analysis code)

```
Name: STAR_2017_PRCGG_644094
Year: 2017_Properties of the medium produced in Relativistic Heavy-Ion Collisic
Collider: BIOL
Co
```

#### ANALYSIS NAME info

(information about paper, beam, author...)

```
REGIN YODA SCATTERED V2 /REF/STAR 2017 PRC96 044984/d20-x04-v01
Variations: [""]
IsRef: 1
Path: /REF/STAR 2017 PRC96 044984/d20-x04-y81
Type: Scatter2D
# xval xerr- xerr+ vval
                                verr- verr+
1.398608e+81
               4.486080e+60
                               4.400800e+08
                                               4.418008e-81
                                                               6.080080e-02
                                                                               6.800808e-02
2.568888e+81
               7 1888888+88
                               7 189869e+98
                                               4 568888e-81
                                                               6.080080e-02
                                                                                6 888888e-82
4.478698e+81
                8.786986e+69
                               8.769869e+98
                                                4.918008e-81
                                                               6.480080e-02
                                                                                6.400000e-02
7.180000e+01
               1.020080e+01
                               1.020800e+01
                                               5.300000e-01
                                                               6.980080e-02
                                                                                6.900000e-02
1.099008e+82
               1.186086e+61
                               1.100800c+01
                                               5.858688e-81
                                                               7.480080e-02
                                                                                7.400800c-02
1.602008e+82
                               1.020800e+01
2.262008e+82
               7.986080e+60
                               7.900800e+08
                                               5.688008e-81
                                                               7.180980e-02
                                                                                7.100800e-02
2.984888e+82
               6.080080e+00
                               6.800800e+08
                                               5.918008e-81
                                                                7.580080e-02
                                                                                7.500000e-02
3.374898e+82
               2.186980c+69
                               2.100800c+00
                                               5.888888e-R1
                                                               7.580000e-02
                                                                                7.500000e-02
```

#### ANALYSIS\_NAME.yoda (experimental data)

```
BEGIN PLOT /STAR_2017 PROS0 044904/020-x04-y01
Title=9/p/pi* { NuA+0.7.0 PC/p()
XLabel=18/p,D1^+$
LogY=0
TWANN-0.8
TWANN-0.8
ComnectEdis=0
LegendY0x=0.8
LegendY0x=0.8
F any additional plot settings you might like, see make-plots documentation
```

END YOUA SCATTER2D V2

END PLOT

ANALYSIS\_NAME.plot (plotting options)



# Analyses in RIVET

#### Catalogue of the analyses available in RIVET: https://rivet.hepforge.org/analyses.html

#### WANTED: Analysis code

We need your analyses! Preserving analysis logic in a rerunnable, re-interpretable form is a key part of scientific reproducibility and impact at the LHC and other HEP experiments. If you are member of an experimental collaboration, please have a look at our wishilst and help us by providing us with Rivet analyses for your publications. This will also ensure that your measurements get used (and cited)!

#### Proposition of a list of relevant analyses to try during this tutorial:

- ALICE\_2010\_S8625980 (charged multiplicity)
- ALICE\_2015\_I1357424 (light hadrons)
- ALICE\_2016\_I1471838 (strange hadrons)
- ALICE\_2012\_I944757 (D mesons)
- ALICE\_2017\_I1645239 ( $\Lambda_c^+$ )

### Running analyses with RIVET

We will run the analyses using a set of 50k events of pp collisions at  $\sqrt{s}=7$  TeV, obtained with **EPOS 4.328**, stored in the file **EPOS\_pp7T\_50k-evts.hepmc**.

**IMPORTANT**: preliminary version of EPOS 4, tuning not complete yet

In order to run any analysis of the catalogue with RIVET, you need to use (you can "force" the analysis to run with an incompatible beam with "--ignore-beams"):

```
> rivet DATA_FILE.hepmc -a ANALYSIS_NAME -o ANALYSIS_NAME.yoda
```

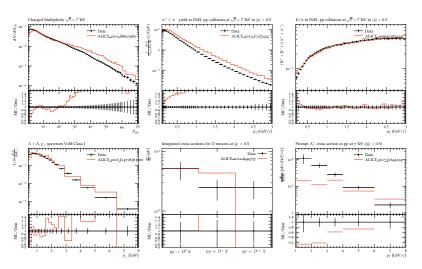
To get plots from the **.yoda** file obtained, you can type the following command: (it will create a folder called "ANALYSIS\_NAME", containing all plots in **.pdf** and **.png**, with a generated URL where all the figures can be displayed.)

```
> rivet-mkhtml ANALYSIS_NAME.yoda -o ANALYSIS_NAME
```

Finally, some analyses like **ALICE\_2016\_I1471838** require a centrality calibration. Hence, you need to run this calibration analysis (**ALICE\_2015\_PPCentrality** here) and call the obtained results after a "-p" flag when running your main analysis:

<u>N.B.:</u> for more information on how to use RIVET, look at RIVET's ReadMe or C. Bierlich's "Hands-on session" (COST workshop 2019)

### Some results you should obtain...



N.B.: these data are only used for illustration; they've been produced with a preliminary version of EPOS 4
(as already highlighted before), which explains why they don't fit well to the experimental results

# Thanks for your attention! Hope you enjoyed it:)



All remarks, suggestions and further expectations will be welcome!

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