FCCAnalyses (for the casesstudies but not only)

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Outline

- Introduction
- Overall structure
- Workflow
- Examples
 - Higgs Recoil
 - Flavour physics at Z pole



Introduction

FCCAnalyses: https://github.com/HEP-FCC/FCCAnalyses/

- Originally developed for FCC-hh physics studies
 - Was based on heppy (full python analysis framework from CMS)
 - There is a heppy-legacy branch for nostalgic people
 - Was used to produced FCC-hh physics analyses published in the CDR

- Transitioned to RDataFrame
 - Heppy was slow
 - Achieved much higher performances
 - Edm4hep aware

Common software?

- <u>FCCAnalyses</u>
 - Is a common tool for analyzing large amount of data using RootDataFrame and produce flat ntuple
 - Is composed of a library of C++ analysers and python configurations
 - C++ analysers are developed in common
 - Python code specific to the analysis to define the analysers, output variables, input samples
 - Use those flat ntuples
 - To produce final variables for analysis, for MVA training, and for plotting
 - To run decay selector for flavour physics
 - Is a-priori transparent to the input format
- EDM4Hep
 - EDM4Hep is the common event data model
 - Is the only supported input format at the moment but this can evolve
 - Will be used for the very long term for any simulation (so better to start using it now!)

- Set of tools to help processing the output of 'simulation'
 - Agnostic to the type of simulation but specific reader functions are required
 - Build a common set of utility functions, algorithms for common use
 - Still possible for users to test their algorithms locally before publishing them
- How is FCCAnalyses structured up to now (all in the same repository)

Analysis configuration 4 python scripts to configure:

- 4 python scripts to configure:
- 1. Samples to run over (common production, EDM4Hep)
- 2. Functions/algorithm to call to produce variables of interests in a flat ntuple
- 3. Event selection and histograms definition using the flat ntuples
- 4. Plotting configuration

Common utility functions, algorithm, etc... C++ library Common interface code Sample database, RdataFrame, plotting **Python**

• How is FCCAnalyses will be structured for the case studies (two separate repositories)

Analysis configuration

4 python scripts to configure:

- 1. Samples to run over (common production, EDM4Hep)
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https://github.com/HEP-FCC/HEP-FCC/FCCeePhysicsPerformance/ Analysis configuration part close to the case studies and analysis specific code (like fitting, specific calculations, etc...)

Common utility functions, algorithm, etc... C++ library Common interface code Sample database, RdataFrame, plotting **Python** https://github.com/HEP-FCC/FCCAnalyses/ Centrally installed on cvmfs and accessible by common setup. But totally OK to use private version.

• How is FCCAnalyses will be structured for the case studies (two separate repositories)

3° master - FCCeePhysicsPerformance / case-stud	Go to file Add file -	
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<u>.</u>		
FCCAnalyses-config	change path	10 days ago
analysis	new configuration skim	13 days ago
images	new configuration skim	13 days ago
🗅 README.md	new configuration skim	13 days ago

• Each case study highly encouraged to have it's own FCCAnalyses-config

Analysis configuration

- 4 python scripts to configure:
- 1. Samples to run over (common production, EDM4Hep)
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• How is FCCAnalyses will be structured for the case studies (two separate repositories)

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analysis	new configuration skim	13 days ago
images	new configuration skim	13 days ago
🗅 README.md	new configuration skim	13 days ago

• But still possible to build and use custom algorithm provided they are properly linked

Common and user specific utility functions, algorithm, etc... C++ library

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Delphes FCCee Physic events tmp (IDEA with Track Covariance)

Q 240

NO	NAME	NEVENTS	NWEIGHTS	NFILES	NBAD	NEOS	SIZE (GB)	OUTPUT PATH	MAIN PROCESS	FINAL STATES	CROSS SECTION (PB)	K-FACTOR	MATCHING EFF
35	p8_ee_ZH_ecm240	9,920,000	0	992	0	992	90.31	<pre>/eos/experiment/fcc/ee/generation/DelphesEvents/fcc_tmp /p8_ee_ZH_ecm240/</pre>	ZH ecm=240GeV	inclusive decays	0.201037	1.0	1.0
36	p8_ee_ZZ_ecm240	9,980,000	0	998	0	998	67.76	/eos/experiment/fcc/ee/generation/DelphesEvents/fcc_tmp /p8_ee_ZZ_ecm240/	ZZ ecm=240GeV	inclusive decays	1.35899	1.0	1.0
37	p8_ee_WW_ecm240	10,000,000	0	1000	0	1000	65.44	/eos/experiment/fcc/ee/generation/DelphesEvents/fcc_tmp /p8_ee_WW_ecm240/	WW ecm=240GeV	inclusive decays	16.4385	1.0	1.0

Common samples are available on eos and can be accessed through a database But possible to by pass this and run with your own files

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Generalities

Each analysis is hosted in a single directory, for example FCCeeAnalyses/ZH_Zmumu/ and contains the same kind of files, please use the same naming convention for all analysis.

- 1. analysis.py : This class that is used to define the list of analysers and filters to run on as well as the output variables.
- 2. presel.py : This configuration file is used to define how to run the analysis.py . It contains the list of samples, the number of CPUs, the fraction of the original sample to process and the base directory for the yaml files (that contains the informations about the samples). This will run the analysis.py with a common code bin/runDataFrame.py (this last file is common to all analyses and should not be touched).
- 3. finalSel.py : This configuration file contains the final selections and it runs over the locally produced flat ntuples from the preSel.py . It contains a link to the procDict.json for getting cross section etc...(this might be removed later to include everything in the yaml, closer to the sample), the list of processes, the number of CPU, the cut list, and the variables (that will be both written in a TTree and in the form of TH1 properly normalised to an integrated luminosity of 1pb-1.
- 4. plots.py : This configuration files is used to select the final selections from running finalsel.py to plot. Informations about how to merge processes, write some extra text, normalise to a given integrated luminosity etc... For the moment it is possible to only plot one signal at the time, but several backgrounds.

python myAnalysis/preSel.py

will run analysis.py on the datasets defined in preSel.py and available in the common database

python myAnalysis/analysis.py /myfiles/file.root or /myfiles/*.root

will run analysis.py on all the input files specified (user responsibility to validate that the files are OK





ZH threshold



- Used the common delphes IDEA track covariance samples I produced
- Processed them through FCCAnalyses machinery with a Z->mumu selection
- And within very little time make this recoil plot
- Anybody with read access to eos can redo those plots in a matter of minutes

new configuration skim

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...

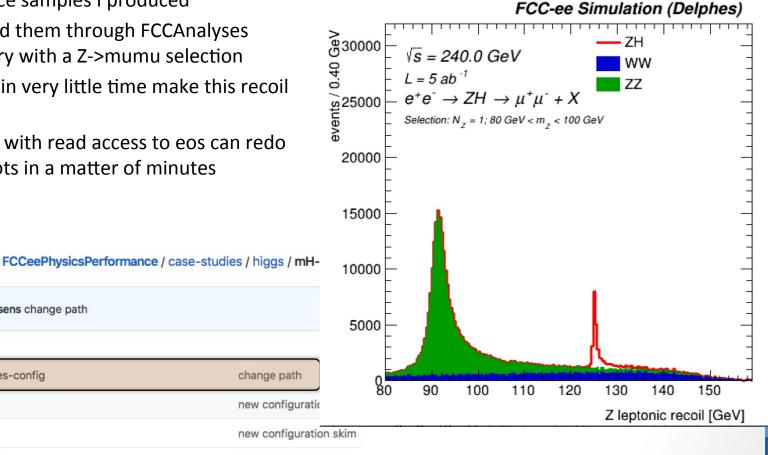
clementhelsens change path

FCCAnalyses-config

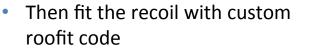
analysis

images

README.md



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- this part has to be written by • analysers and goes in the analysis
- Fitted mass: .

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100

...

125.07 +/- 0.00357 GeV ۲

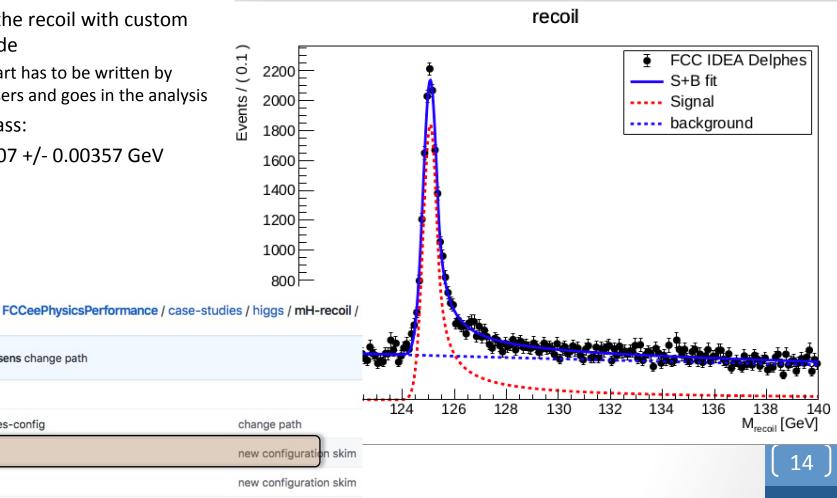
clementhelsens change path

FCCAnalyses-config

analysis

images

README.md



new configuration skim

Flavour physics at the Z pole



Define variables and filters: MC

analysis.py

.Alias("MCRecoAssociations0", "MCRecoAssociations#0.index") .Alias("MCRecoAssociations1", "MCRecoAssociations#1.index") .Alias("Particle0", "Particle#0.index")

```
.Filter("filterMC_pdgID(541, true)(Particle)==true")
.Filter("filterMC_pdgID(541, true)(Particle)==false")
```

```
.Define("MC px",
                         "getMC px(Particle)")
.Define("MC py",
                         "getMC py(Particle)")
                         "getMC pz(Particle)")
.Define("MC pz",
.Define("MC p",
                         "getMC p(Particle)")
.Define("MC e",
                         "getMC e(Particle)")
                         "getMC pdg(Particle)")
.Define("MC pdg",
.Define("MC charge",
                         "getMC charge(Particle)")
.Define("MC mass",
                         "getMC mass(Particle)")
.Define("MC_status",
                         "getMC genStatus(Particle)")
                         "getMC vertex x(Particle)")
.Define("MC vertex x",
.Define("MC vertex y",
                         "getMC vertex y(Particle)")
                         "getMC vertex z(Particle)")
.Define("MC_vertex_z",
```

Needed by RDF to understands #

Filter if event contains a Bc hadron or not, This will create an exclusive sample or inclusive -exclusive

Define the MC quantities to keep

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Define variables and filters: Reco

analysis.py

.Define("RP_p", .Define("RP_e", .Define("RP_px", .Define("RP_py", .Define("RP_pz", .Define("RP_charge", .Define("RP_mass",

"getRP_p(ReconstructedParticles)") "getRP_e(ReconstructedParticles)") "getRP_px(ReconstructedParticles)") "getRP_py(ReconstructedParticles)") "getRP_pz(ReconstructedParticles)") "getRP_charge(ReconstructedParticles)") "getRP_mass(ReconstructedParticles)")

Access reconstructed particles quantities

```
#.Define("RP_TRK_D0",
#.Define("RP_TRK_Z0",
```

```
"getRP2TRK_D0(ReconstructedParticles, EFlowTrack_1)")
"getRP2TRK_Z0(ReconstructedParticles, EFlowTrack_1)")
Can get the track associated to them
```

.Define('RP_MC_index',

"getRP2MC_index(MCRecoAssociations0,MCRecoAssociations1,ReconstructedParticles)")

Build a collection of size Reco that contains the index in the MC_* columns for the associated MC particle

.Define('RP_MC_parentindex', "getMC_parentid(RP_MC_index,Particle, Particle0)")

Use the column defined above and build a new column of MC parents index for each Reco particles

.Define('RP_MC_grandparentindex', "getMC_parentid(RP_MC_parentindex,Particle, Particle0)")

Use the column defined above and build a new column of MC grand parents index for each Reco particles

Define variables and filters: Event

'EVT thrust.at(0)')

'EVT thrust.at(5)')

.Define('EVT_thrust', 'minimize_thrust("Minuit2", "Migrad")(RP_px, RP_py, RP_pz)')

analysis.py

Call the minimisation for the thrust and sphericity and

define the relevant columns

.Define('EVT_sphericity', 'minimize_sphericity("Minuit2","Migrad")(RP_px, RP_py, RP_pz)') .Define('EVT_sphericity_val', 'EVT_sphericity.at(0)') .Define('EVT_sphericity_x', 'EVT_sphericity.at(1)') .Define('EVT_sphericity_x_err','EVT_sphericity.at(2)') .Define('EVT_sphericity_y', 'EVT_sphericity.at(3)') .Define('EVT_sphericity_y_err','EVT_sphericity.at(4)') .Define('EVT_sphericity_z', 'EVT_sphericity.at(5)') .Define('EVT_sphericity_z err','EVT_sphericity.at(6)')

EVT_thrust/sphericity is a vector of size 7 Value/x-axis/x-axis-error/y-axis/y-axis-error/z-axis/z-axis-error

.Define('EVT thrust val',

.Define('EVT thrust z',

.Define('EVT thrust x', 'EVT thrust.at(1)')

.Define('EVT thrust x err', 'EVT thrust.at(2)')

.Define('EVT_thrust_y', 'EVT_thrust.at(3)') .Define('EVT thrust y err','EVT thrust.at(4)')

.Define('EVT_thrust_z_err', 'EVT_thrust.at(6)')

Define variables and filters: Event

.Define('RP thrustangle', 'axisCosTheta(EVT thrust, RP px, RP py, RP pz)') cosTheta between an axis and the reconstructed particles .Define('RP sphericityangle', 'axisCosTheta(EVT sphericity, RP px, RP py, RP pz)')

.Define('EVT_thrusthemis0_q_kappal',	'getAxisCharge(0, 1)(RP_thrustangle,	, RP_charge, RP_px, RP_py, RP_pz)')	Charge of a given hemisphere
			(defined by the sign of cosTheta)

Number of particles (total/charged/neutral) .Define('EVT thrusthemis0 n', 'getAxisN(0)(RP thrustangle, RP charge)') .Define('EVT thrusthemisl n', 'getAxisN(1)(RP thrustangle, RP charge)') in each hemisphere

Energy sum of particles (total/charged/neutral) .Define('EVT thrusthemis0 e', 'getAxisEnergy(0)(RP thrustangle, RP charge, RP e)') 'getAxisEnergy(1)(RP_thrustangle, RP_charge, RP_e)') .Define('EVT_thrusthemisl_e', in each hemisphere

```
.Define('EVT thrutshemis0 ncharged',
                                     'EVT thrusthemis0 n.at(1)')
.Define('EVT thrutshemis0 nneutral',
                                     'EVT thrusthemis0 n.at(2)')
```

```
.Define('EVT thrutshemis0 echarged', 'EVT thrusthemis0 e.at(1)')
.Define('EVT thrutshemis0 eneutral', 'EVT thrusthemis0 e.at(2)')
```

```
.Define('EVT thrutshemis1 ncharged',
                                     'EVT thrusthemisl n.at(1)')
.Define('EVT_thrutshemis1_nneutral', 'EVT_thrusthemis1_n.at(2)')
```

.Define('EVT thrutshemis1 echarged', 'EVT thrusthemis1 e.at(1)') .Define('EVT thrutshemis1 eneutral', 'EVT thrusthemis1 e.at(2)')

Number of particles (charged/neutral) in hemisphere with cosTheta<0

Energy sum of particles (charged/neutral) in hemisphere with cosTheta<0

Number of particles (charged/neutral) in hemisphere with cosTheta>0

Energy sum of particles (charged/neutral) in hemisphere with cosTheta>0

.Define('EVT thrutshemis emax', 'if (EVT thrusthemis0 e.at(0)>EVT thrusthemis1 e.at(0)) return EVT thrusthemis0 e.at(0); else return EVT thrusthemis1 e.at(0); ') .Define('EVT thrutshemis emin', 'if (EVT thrusthemis0 e.at(0)>EVT thrusthemis1 e.at(0)) return EVT thrusthemis1 e.at(0); else return EVT thrusthemis0 e.at(0);')

Define the hemisphere with minimum/maximum energies

analysis.py

Define variables and filters: Run

analysis.py

select branches for output file branchList = ROOT.vector('string')() for branchName in [

"MC_px", "MC_py", "MC_pz", "MC_p", "MC_e", "MC_pdg", "MC_charge", "MC_mass", "MC_status", "MC_vertex_x", "MC_vertex_y", "MC_vertex_z",

And run

"EVT_thrust_x","EVT_thrust_y","EVT_thrust_z","EVT_thrust_val","EVT_thrusthemis0_q_kappal","EVT_thrusthemis1_q_kappal", "EVT_thrutshemis0_ncharged","EVT_thrutshemis1_ncharged","EVT_thrutshemis0_nneutral","EVT_thrutshemis1_nneutral", "EVT_thrutshemis0_echarged","EVT_thrutshemis1_echarged","EVT_thrutshemis0_eneutral","EVT_thrutshemis1_eneutral", "EVT_thrutshemis_emax","EVT_thrutshemis_emin",

```
"EVT_sphericity_x", "EVT_sphericity_y", "EVT_sphericity_z", "EVT_sphericity_val",
```

"RP_thrustangle", "RP_sphericityangle", "RP_p", "RP_px", "RP_py", "RP_pz", "RP_charge", "RP_mass",

```
"RP_MC_index", "RP_MC_parentindex", "RP_MC_grandparentindex",
```

```
]:
```

branchList.push_back(branchName)

Defined the columns to keep

```
#opts = ROOT.RDF.RSnapshotOptions()
#opts.fCompressionAlgorithm = ROOT.ROOT.kLZ4
#opts.fCompressionLevel = 3
#opts.fAutoFlush = -1024*1024*branchList.size()
#df2.Snapshot("events", self.outname, branchList, opts)
df2.Snapshot("events", self.outname, branchList)
```

preSel configuration

#python examples/FCCee/flavour/generic-analysis/preSel.py

```
from config.common_defaults import deffccdicts
import os
```

```
basedir=os.path.join(os.getenv('FCCDICTSDIR', deffccdicts), '') + "yaml/FCCee/fcc_tmp_v02/"
outdir="outputs/FCCee/flavour/generic-analysis/"
```

To run with the common production, files are validated etc...

preSel.py

```
import multiprocessing
```

```
NUM_CPUS = int(multiprocessing.cpu_count()-2)
process_list=[
    'p8_ee_Zbb_ecm91_EvtGen_Bc2TauNuTAUHADNU',
    'p8_ee_Zcc_ecm91',
    'p8_ee_Zuds_ecm91'
```

Define here the process list, names are in the Database

```
#output_list=['p8_ee_Zbb_ecm91']
output_list=[]
```

Possibility to change the ouptut name, can be used when filtering a more inclusive sample

.Filter("filterMC_pdgID(541, true)(Particle)==true") .Filter("filterMC_pdgID(541, true)(Particle)==false")

```
fraction=1.0
```

```
import config.runDataFrame as rdf
myana=rdf.runDataFrame(basedir,process_list, outlist=output_list)
myana.run(ncpu=NUM_CPUS,fraction=fraction,outDir=outdir)
```

```
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```

finalSel configuration

finaSel.py

###Input directory where the files produced at the pre-selection level are

- #baseDir = "outputs/FCCee/flavour/generic-analysis/"
- baseDir = "/eos/experiment/fcc/ee/tmp/flatntuples/Z_Zbb_Bc2TauNu/"

```
###Link to the dictonary that contains all the cross section informations etc...
#procDict = os.path.join(os.getenv('FCCDICTSDIR', deffccdicts), '') + "FCCee_procDict_fcc_tmp_v02.json"
#procDict = 'https://fcc-physics-events.web.cern.ch/fcc-physics-events/sharedFiles/FCCee_procDict_fcc_v02.json'
procDict = {
    "p8_ee_Zbb_ecm91": {"numberOfEvents": 19830000, "sumOfWeights": 19830000, "crossSection": 6645.46, "kfactor": 1.0, "matchingEfficiency": 1.0},
    "p8_ee_Zbb_ecm91_EvtGen_Bc2TauNuTAUHADNU": {"numberOfEvents": 14789, "sumOfWeights": 14789, "crossSection": 6645.46, "kfactor": 1.0, "matchingEfficiency": 1.0, "matchingEfficiency": 1.0,
    "p8_ee_Zbb_ecm91": {"numberOfEvents": 1000000, "sumOfWeights": 14789, "sumOfWeights": 12789, "crossSection": 6645.46, "kfactor": 1.0, "matchingEfficiency": 1.0,
    "p8_ee_Zcc_ecm91": {"numberOfEvents": 1000000, "sumOfWeights": 12789, "sumOfWeights": 12789, "crossSection": 6645.46, "kfactor": 1.0, "matchingEfficiency": 1.0,
    "p8_ee_Zcc_ecm91": {"numberOfEvents": 1000000, "sumOfWeights": 1000000, "crossSection": 5215.46, "kfactor": 1.0, "matchingEfficiency": 1.0,
    "p8_ee_Zuds_ecm91": {"numberOfEvents": 1000000, "sumOfWeights": 1000000, "crossSection": 5215.46, "kfactor": 1.0, "matchingEfficiency": 1.0},
    "p8_ee_Zuds_ecm91": {"numberOfEvents": 1000000, "sumOfWeights": 1000000, "crossSection": 18616.5, "kfactor": 1.0, "matchingEfficiency": 1.0},
    "p8_ee_Zuds_ecm91": {"numberOfEvents": 1000000, "sumOfWeights": 1000000, "crossSection": 18616.5, "kfactor": 1.0, "matchingEfficiency": 1.0}
}
process_list=['p8_ee_Zbb_ecm91_EvtGen_Bc2TauNuTAUHADNU', 'p8_ee_Zbb_ecm91', 'p8_ee_Zuds_ecm91']
define list={
```

"EVT_Ediff":"EVT_thrutshemis_emax-EVT_thrutshemis_emin",

```
"EVT_Echarged_max":"if (EVT_thrutshemis0_echarged>EVT_thrutshemis1_echarged) return EVT_thrutshemis0_echarged; else return EVT_thrutshemis1_echarged;",
"EVT_Echarged_min":"if (EVT_thrutshemis0_echarged>EVT_thrutshemis1_echarged) return EVT_thrutshemis1_echarged; else return EVT_thrutshemis0_echarged;",
"EVT_Eneutral_max":"if (EVT_thrutshemis0_eneutral>EVT_thrutshemis1_eneutral) return EVT_thrutshemis0_eneutral; else return EVT_thrutshemis1_eneutral;",
"EVT_Eneutral_max":"if (EVT_thrutshemis0_eneutral>EVT_thrutshemis1_eneutral) return EVT_thrutshemis0_eneutral; else return EVT_thrutshemis0_eneutral;",
```

"EVT_Ncharged_max":"if (EVT_thrutshemis0_ncharged>EVT_thrutshemis1_ncharged) return EVT_thrutshemis0_ncharged; else return EVT_thrutshemis1_ncharged;", "EVT_Ncharged_min":"if (EVT_thrutshemis0_ncharged>EVT_thrutshemis1_ncharged) return EVT_thrutshemis1_ncharged; else return EVT_thrutshemis0_ncharged;", "EVT_Nneutral_max":"if (EVT_thrutshemis0_nneutral>EVT_thrutshemis1_nneutral) return EVT_thrutshemis0_nneutral; else return EVT_thrutshemis1_nneutral;", "EVT_Nneutral_max":"if (EVT_thrutshemis0_nneutral>EVT_thrutshemis1_nneutral) return EVT_thrutshemis1_nneutral; else return EVT_thrutshemis0_nneutral;",

}

Using common procDict or custom Possibility to define new variables

```
R00T.gInterpreter.Declare("""
bool myFilter(R00T::VecOps::RVec<float> mass) {
    if (mass.size()<2) return false;
    for (size_t i = 0; i < mass.size(); ++i) {
        if (mass.at(i)<80. || mass.at(i)>100.)
            return false;
    }
    return true;
}
```



finalSel configuration

###Dictionary for the ouput variable/hitograms. The key is the name of the variable in the output files. "name" is the name of the v he histogram, "bin" the number of bins of the histogram, "xmin" the minimum x-axis value and "xmax" the maximum x-axis value. variables = {

"EVT_thrust_val":{"name":"EVT_thrust_val","title":"Event Thrust","bin":100,"xmin":0.4,"xmax":1.},
"EVT_thrusthemis_emax":{"name":"EVT_thrusthemis_emax","title":"Hemisphere energy (max) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_thrusthemis_emin":{"name":"EVT_thrusthemis_emin","title":"Hemisphere energy (min) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Ediff":{"name":"EVT_Ediff","title":"Hemisphere energy difference [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Echarged_max":{"name":"EVT_Echarged_max","title":"Hemisphere charged energy (max) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Echarged_min":{"name":"EVT_Echarged_max","title":"Hemisphere charged energy (max) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Eneutral_max":{"name":"EVT_Echarged_min","title":"Hemisphere charged energy (min) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Eneutral_max":{"name":"EVT_Echarged_min","title":"Hemisphere charged energy (min) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Eneutral_max":{"name":"EVT_Eneutral_max","title":"Hemisphere neutral energy (max) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Eneutral_max":{"name":"EVT_Eneutral_max","title":"Hemisphere neutral energy (min) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Eneutral_max":{"name":"EVT_Eneutral_min","title":"Hemisphere neutral energy (min) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Eneutral_min":{"name":"EVT_Eneutral_min","title":"Hemisphere neutral energy (min) [GeV]","bin":120,"xmin":0.,"xmax":60},
"EVT_Leneutral_max":{"name":"EVT_Eneutral_min","title":"Hemisphere neutral energy (min) [GeV]","bin":120,"xmin":0.,"xmax":25},
"EVT_Ncharged_max":{"name":"EVT_Ncharged_min","title":"Hemisphere charged multiplicity (max)","bin":25,"xmin":0.,"xmax":25},
"EVT_Nneutral_max":{"name":"EVT_Nneutral_max","title":"Hemisphere charged multiplicity (max)","bin":25,"xmin":0.,"xmax":25},
"EVT_Nneutral_min":{"name":"EVT_Nneutral_min","title":"Hemisphere neutral multiplicity (min)","bin":25,"xmin":0.,"xmax":25},"EVT_Nneutral_min":{"name":"EVT_Nneutral_min","title":"Hemisphere neutral

###Number of CPUs to use NUM_CPUS = 4 ###This part is standard to all analyses import config.runDataFrameFinal as rdf #myana=rdf.runDataFrameFinal(baseDir,procDict,process_list,cut_list,variables) myana=rdf.runDataFrameFinal(baseDir,procDict,process_list,cut_list,variables,defines=define_list) myana.run(ncpu=NUM_CPUS, doTree=False) #"")

```
ROOT.gInterpreter.Declare("""
bool myFilter(ROOT::VecOps::RVec<float> mass) {
    if (mass.size()<2) return false;
    for (size_t i = 0; i < mass.size(); ++i) {
        if (mass.at(i)<80. || mass.at(i)>100.)
            return false;
}
return true;
```

"sel3":"myFilter(zed_leptonic_m)"

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finaSel.py

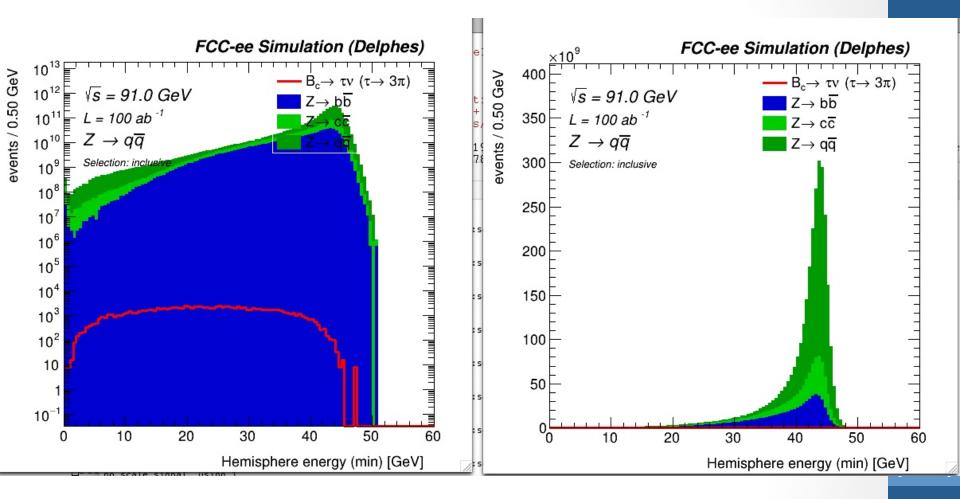
plots configuration

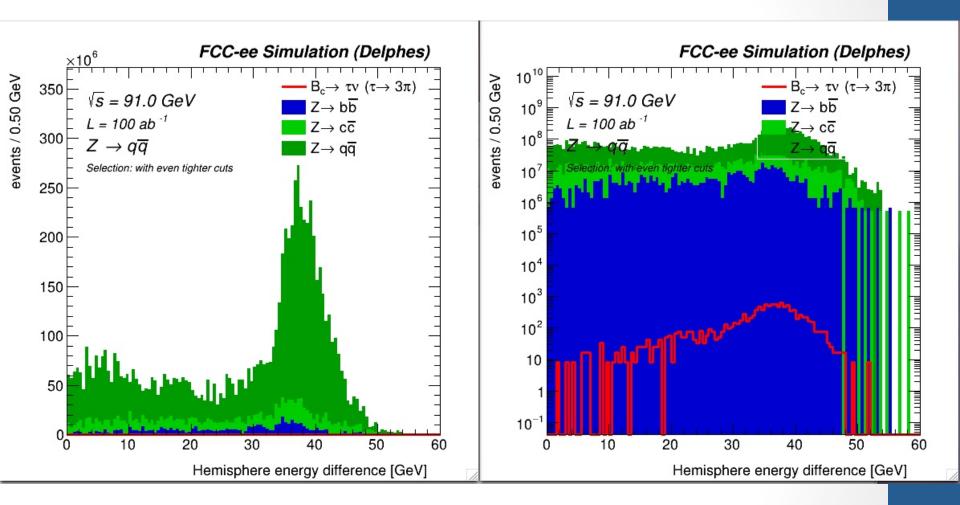
plots.py

```
###Dictonnary with the analysis name as a key, and the list of selections to be
                                                                            selections = {}
                                                                            selections['Flavour'] = ["sel0"."sel1"."sel2"."sel3"]
                                                                           extralabel = {}
# global parameters
                                                                            extralabel['sel0'] = "Selection: inclusive"
                = 100000000. #in pb-1
intLumi
                                                                            extralabel['sell'] = "Selection: with cuts"
               = "Z #rightarrow g#bar{g}"
ana tex
                                                                            extralabel['sel2'] = "Selection: with tighter cuts"
delphesVersion = "3.4.2"
                                                                            extralabel['sel3'] = "Selection: with even tighter cuts"
energy
               = 91.0
collider
               = "FCC-ee"
                                                                                                                                                         14/12
inputDir
               = "/eos/experiment/fcc/ee/tmp/flatntuples/Z Zbb Bc2TauNu/"
formats
                = ['png','pdf']
                = ['lin','log']
yaxis
                                                                          colors = {}
stacksig
                = ['stack', 'nostack']
                                                                          colors['Z flavour'] = ROOT.kRed
outdir
                = 'outputs/FCCee/flavour/generic-analysis/plots/'
                                                                          colors['Z bb'] = ROOT.kBlue+1
                                                                          colors['Z cc'] = R00T.kGreen+1
variables = [ "EVT_thrust_val",
                                                                          colors['Z uds'] = R00T.kGreen+2
               "EVT thrusthemis emax",
                                                                          plots = {}
               "EVT thrusthemis emin",
                                                                          plots['Flavour'] = {'signal':{'Z flavour':['p8 ee Zbb ecm91 EvtGen Bc2TauNuTAUHADNU']},
               "EVT Ediff",
                                                                                             'backgrounds':{'Z bb':['p8 ee Zbb ecm91'],
               "EVT Echarged max",
                                                                                                           'Z_cc':['p8_ee_Zcc_ecm91'],
               "EVT Echarged min",
                                                                                                           'Z uds':['p8 ee Zuds ecm91']}
               "EVT Eneutral max",
                                                                                     }
               "EVT Eneutral min".
                                                                          legend = \{\}
                                                                           legend['Z_flavour'] = 'B_{c}#rightarrow #tau#nu (#tau#rightarrow 3#pi)'
               "EVT Ncharged max",
                                                                          legend['Z bb'] = 'Z#rightarrow b#bar{b}'
               "EVT Ncharged min",
                                                                          legend['Z_cc'] = 'Z#rightarrow c#bar{c}'
               "EVT Nneutral max",
                                                                           legend['Z uds'] = 'Z#rightarrow q#bar{q}'
               "EVT Nneutral min"
```

]

Common plot style will be supported here. But users could write their own as well Plan to write a common plotting code independent of ROOT





FCCAnalysis: available algorithms

- Available functionality (see <u>doxygen</u>)
 - Thrust axis and thrust value
 - Sphericity axis and value
 - cos(theta) between an axis and particles
 - Weighted charge of all particles with cos(theta) axis > or <0
 - Preliminary jet clustering
 - Truth history from any MC particle of a given status
 - Tagging 2 body MC decays (for example H->gg, Z->bb)
 - Etc...
- Planned functionalities for the common analysers
 - Acceptance efficiencies
 - Parametric PID
 - Pi0 identification
 - Etc...
- Planning to have a dedicated uproot/awkward structure to easily run flavour exclusive decays, etc, etc...

Summary/Plans

- Easy to use analysis code to produce flat ntuples based on Root DataFrame ready for you to run analyses, please join and have fun
- Did not talked about FCC-hh but all this code can be used there as well (or in ATLAS/CMS?)
- Plans
 - Adding uproot and awkward from pure python or C++
 - Etc...