

SW Review (CMS top)

Status Report

4-MARCH-2022



Main version of the classes are ready

- Gitlab repo updated: https://gitlab.cern.ch/cdozen/nanoaodrdframe-ul/-/tree/UL_TTLFV
 - ➔ Before making any changes/modification try to compile and run it!
 - ❖ Detailed instruction: [NanoAODframework instruction](#)
- BaseAnalyser.cpp & BaseAnalyser.h —> Default version
 - User need to rename the source and header files according to the their analysis
 - BaseAnalyser.cpp —> TprimeAnalyser.cpp or SingletopAnalyser.cpp
 - BaseAnalyser.h —> TprimeAnalyser.h or SingletopAnalyser.h

Object ID definition

- In NanoAODAnalyzerrrdframe.cpp
- ElectronID added as an example. User need to define other object IDs (MuonID, JetID, etc.)

```
//cut-based ID Fall17 V2 (0:fail, 1:veto, 2:loose, 3:medium, 4:tight)
std::string NanoAODAnalyzerrrdframe::ElectronID(int cutbasedID){

if(cutbasedID<0 || cutbasedID>4){
    std::cout<< "Error Wrong Electron ID requested == " << cutbasedID << "!! Can't be applied" <<std::endl;
    std::cout<< "Please select number from 0 to 4 " <<std::endl;
}

//Rdataframe look for the variables in the input Ttree..
std::string output = Form("Electron_cutBased == %d ",cutbasedID);

return output;
}
```

Object ID definition

- In BaseAnalyser.cpp
- Call ElectronID

```
void BaseAnalyser::selectElectrons()
{
    cout << "select good electrons" << endl;
    //string veto_electron_str = Form ("Electron_pt>%f && abs(Electron_eta)<%f && %s", cutdb->Md_Electron_pT, cutdb->Md_Electron_eta, ElectronID(2).c_str());
    //_rlm = _rlm.Define("goodElectrons", veto_electron_str.c_str());

    _rlm = _rlm.Define("goodElectrons", ElectronID(2)); //without pt-eta cuts
    _rlm = _rlm.Define("goodElectrons_pt", "Electron_pt[goodElectrons]")
        .Define("goodElectrons_eta", "Electron_eta[goodElectrons]")
        .Define("goodElectrons_phi", "Electron_phi[goodElectrons]")
        .Define("goodElectrons_mass", "Electron_mass[goodElectrons]")
        .Define("goodElectrons_idx", ::good_idx, {"goodElectrons"})
        .Define("NgoodElectrons", "int(goodElectrons_pt.size())");

    //-----
    //generate electron 4vector
    //-----
    _rlm = _rlm.Define("goodElectron_4Vecs", ::generate_4vec, {"goodElectrons_pt", "goodElectrons_eta", "goodElectrons_phi", "goodElectrons_mass"});
}
```

You, last week • cleaned main class and added electron ID informat...

Running on a range of entries

- BaseAnalyser.cpp:
- In Rdataframe
 - Range() : filter based on entry number

```
//=====Event Selection=====//  
// Cuts to be applied in order  
//=====//  
void BaseAnalyser::defineCuts()  
{  
    // count some entries using ranges  
    auto Nentry = _rlm.Count();  
    // This is how you can express a range of the first 100 entries  
    _rlm = _rlm.Range(0, 100);  
    auto Nentry_100 = _rlm.Count();  
    // This is how you pick all entries from 10 onwards  
    _rlm = _rlm.Range(10, 0);  
    auto Nentry_10_end = _rlm.Count();  
    // We can use a stride too, in this case we pick an event every 3  
    _rlm = _rlm.Range(10, 0, 3);  
    auto Nentry_10_end_3 = _rlm.Count();  
  
    cout << "Usage of ranges:\n"  
        << " - All entries: " << *Nentry << endl  
        << " - Entries from 0 to 100: " << *Nentry_100 << endl  
        << " - Entries from 10 onwards: " << *Nentry_10_end << endl  
        << " - Entries from 100 onwards in steps of 3: " << *Nentry_10_end_3 << endl;
```

Running on Max number of files: - -nrootfiles

- Select number of files in the command line:
 - - -nrootfiles (or -N) 4 //4 files will be selected in the inputfile directory

```
time ./processnanoaod.py --allinone -Y 2018 --nrootfiles 4 --isDATA 0 roottestfilesUL18/ testBaseANA_UL18.root >testBaseUL.out
```

Running on Data / MC and UL / ReReco Selection

- DATA and MC selection can be done by user : on the command line
 - - -isDATA 0 : means MC
 - - -isDATA 1 : means DATA
 - - - isDATA -1 or nothing defined : it will be selected according to the input file “genweight” branch

```
time ./processnanoaod.py --allinone -Y 2018 --nrootfiles 4 --isDATA 0 roottestfilesUL18/ testBaseANA_UL18.root >testBaseUL.out
```

- UL and ReReco selection can be done by user : on the command line
 - - -runtime (or -R) UL
 - - - runtime (or -R) ReReco

```
time ./processnanoaod.py --allinone -Y 2018 --runtime UL --isDATA 0 roottestfilesUL18/ testBaseANA_UL18.root >testBaseUL.out
```

Running options

- Structure:
 - `./processnanoaod.py [option] [inputdir] [outputdir]`
 - `./processnanoaod.py [options] [inputDir] {$outputroot}` -> for “- - allinone” option
 - [option]:
 - **- - split** : How many jobs to split into. it will allow multithreading you to use multiple CPUs to process the files. (e.g. `—split=5`)
 - **[inputdir]**: Inputfiles directory : **roottestfilesUL18/**
 - **[outputdir]**: output directory : **testBaseANA_UL** : no need to create an output directory, it will be created if there isn't

```
./processnanoaod.py --split 4 -Y 2018 --runtype UL - -nrootfile 4 --isDATA 0 roottestfilesUL18/ testBaseANA_UL
```

- **--allinone** (-A) : will let you get a **single output root file** for all the files in the inputDir.
- **{outputroot}**: should be a root file name : **testBaseANA_UL18.root**

```
./processnanoaod.py --allinone -Y 2018 --runtype UL --nrootfiles 4 --isDATA 0 roottestfilesUL18/ testBaseANA_UL18.root
```

Bonus slide

Trello_NanoUL_cmstop_working_space

