

# Introduction to the analysis package.

Yo Sato (Tohoku University)

Belle II Starter Kit workshop, KEK, January 2020



**HELMHOLTZ** RESEARCH FOR  
GRAND CHALLENGES

**DESY.**



Istituto Nazionale di Fisica Nucleare  
SEZIONE DI TORINO



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Physics quantities and the VariableManager

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Nomenclature

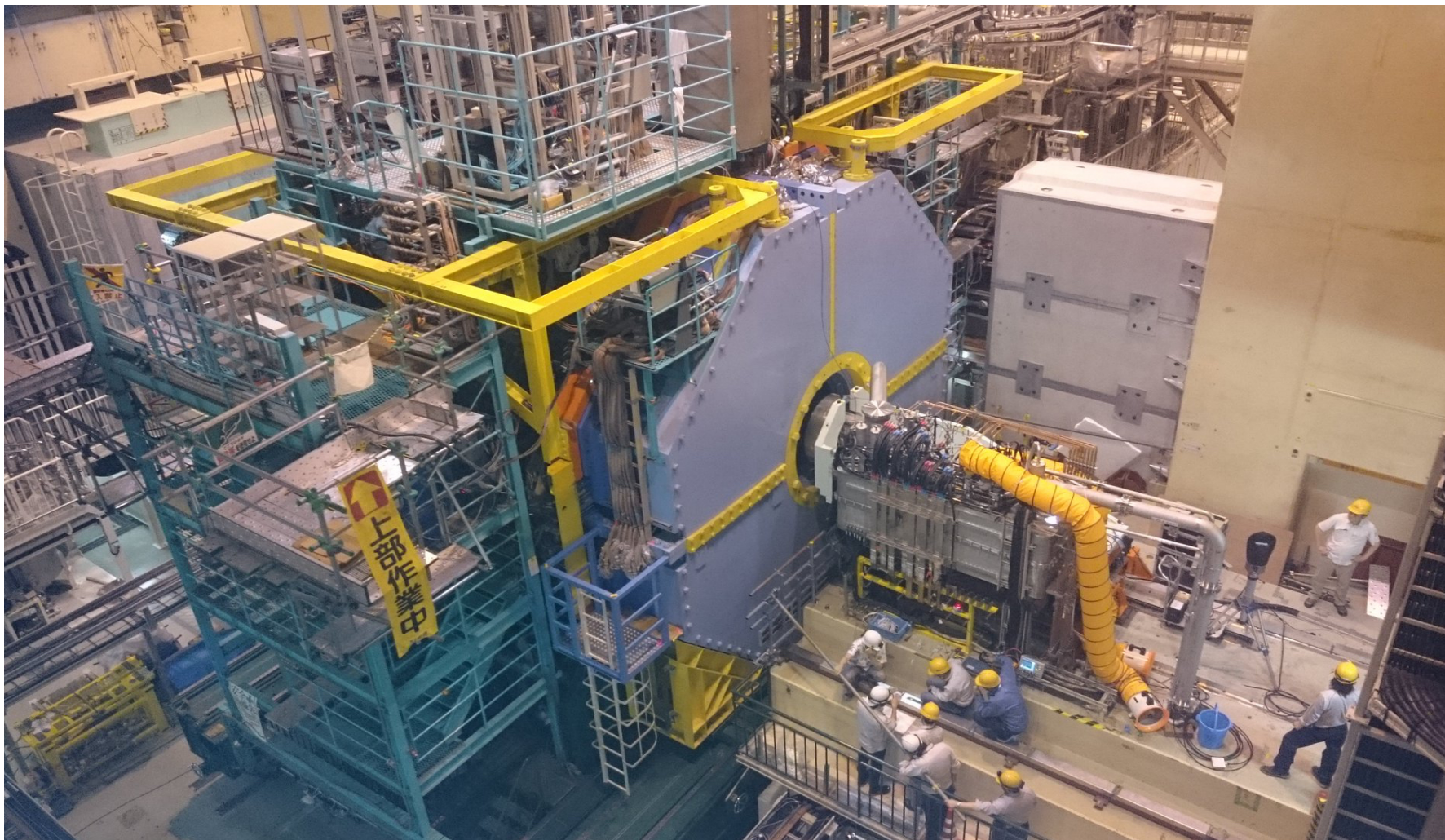
One final thing about Belle vs. Belle II



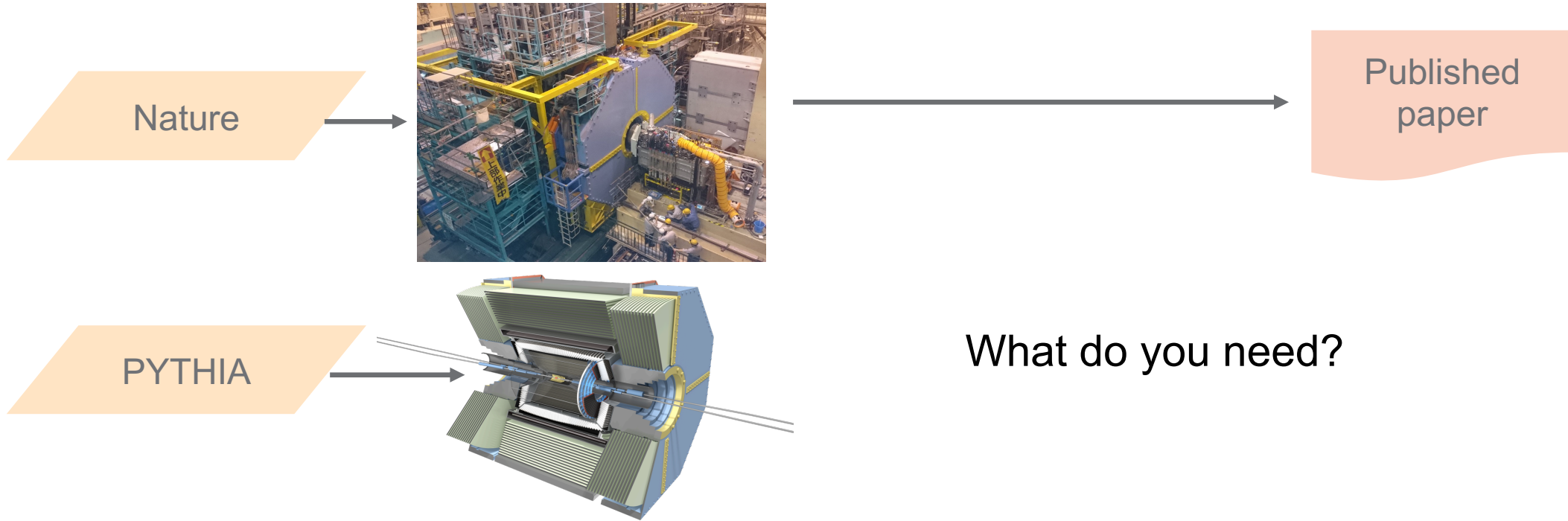
# An experiment's workflow

# The big picture

## Tsukuba Hall

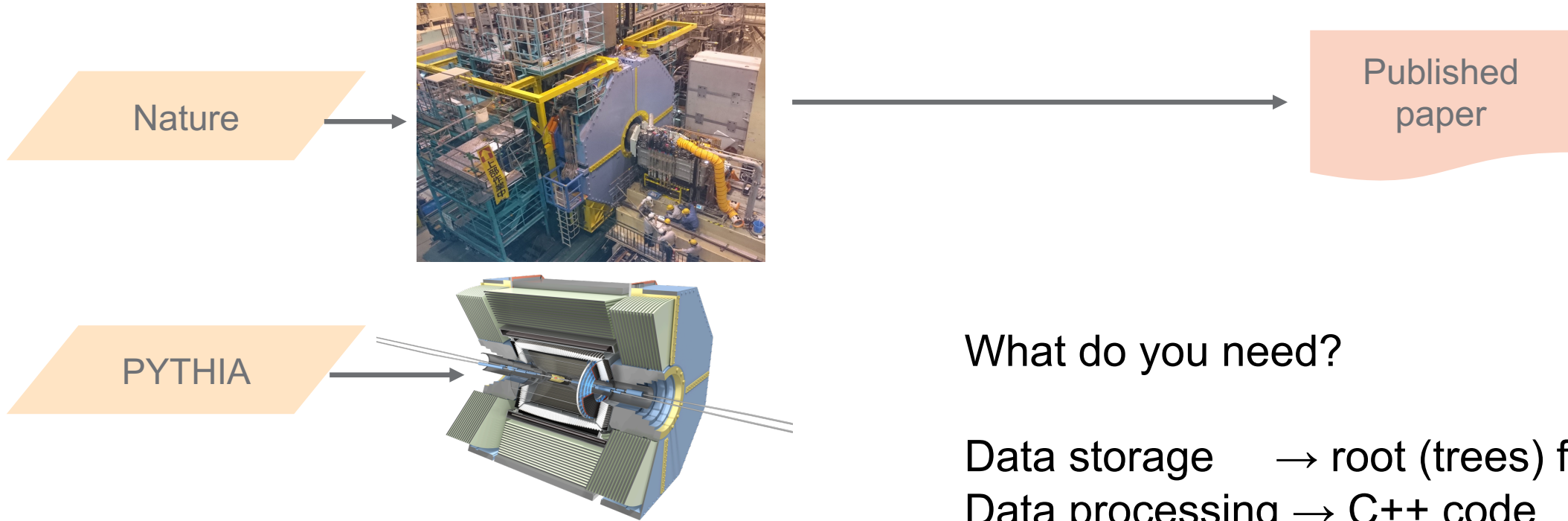


# The big picture





# The big picture



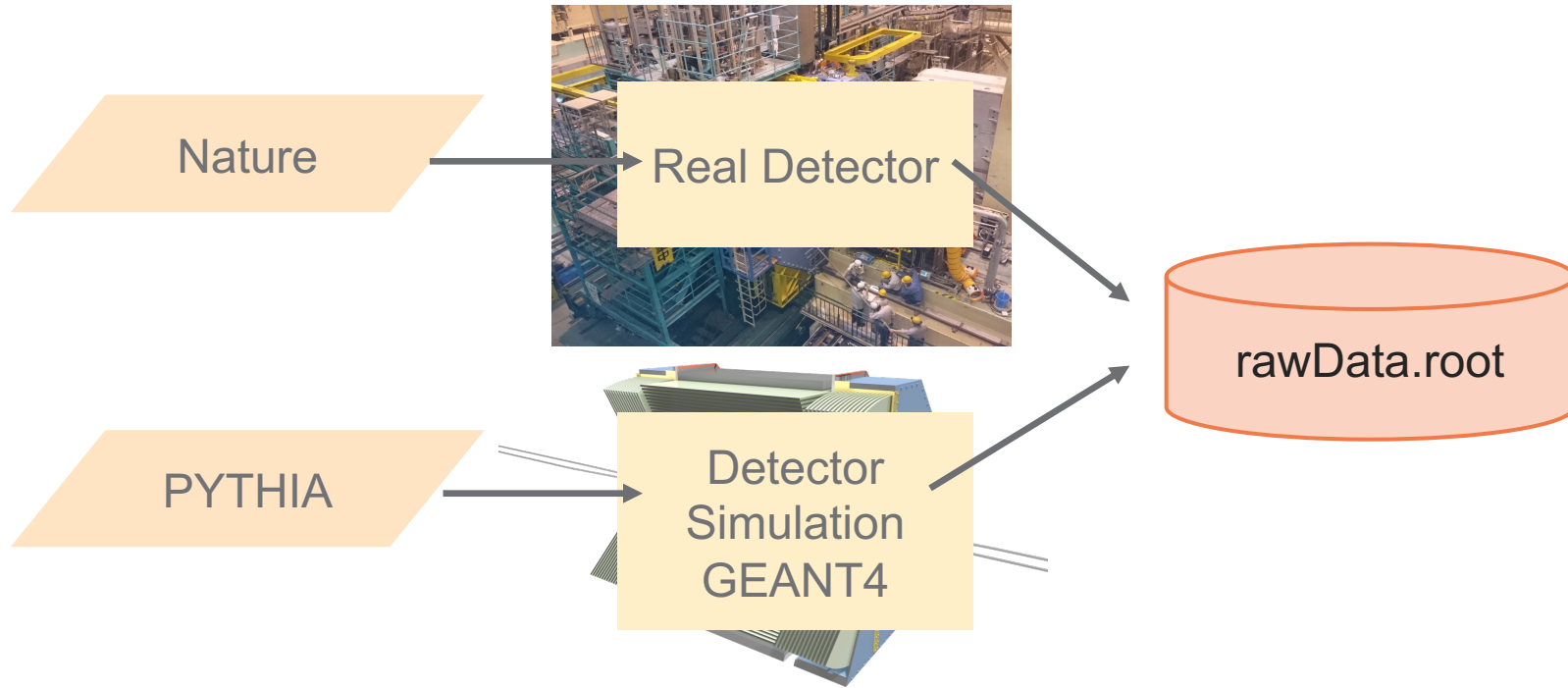
What do you need?

Data storage → root (trees) files

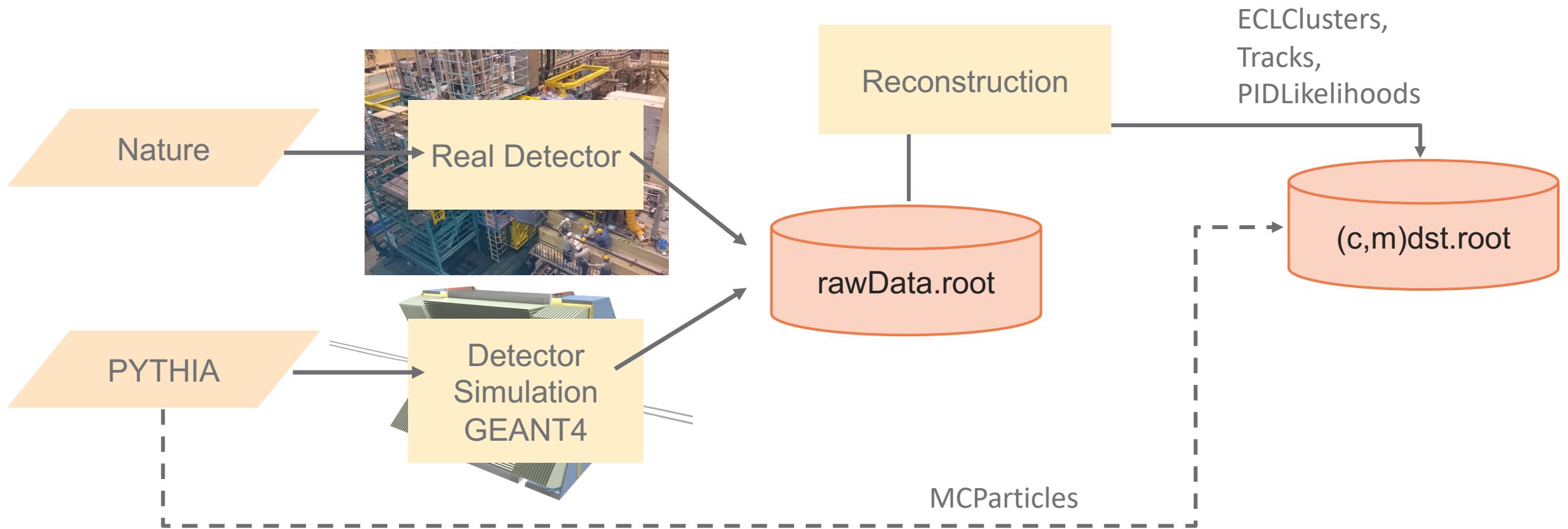
Data processing → C++ code

Scripting → Python

# The big picture

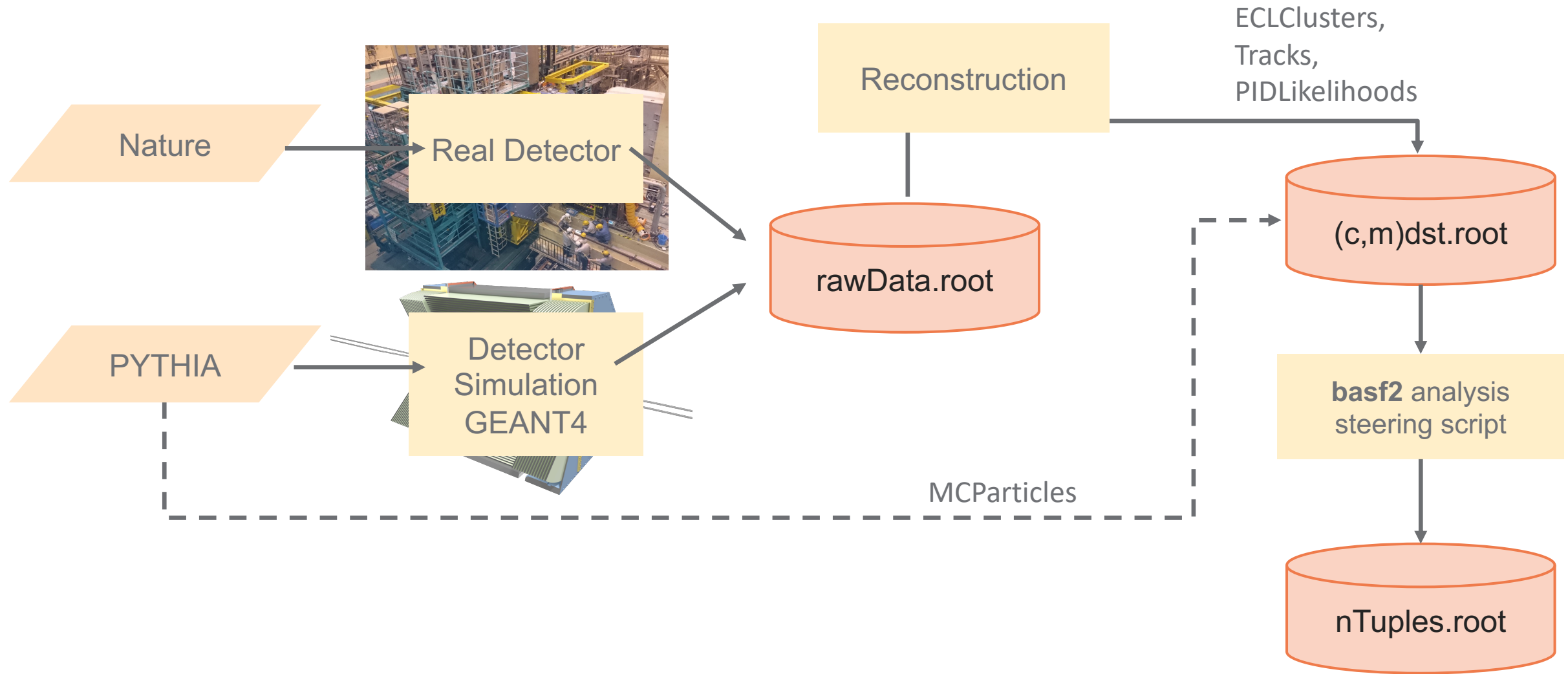


# The big picture

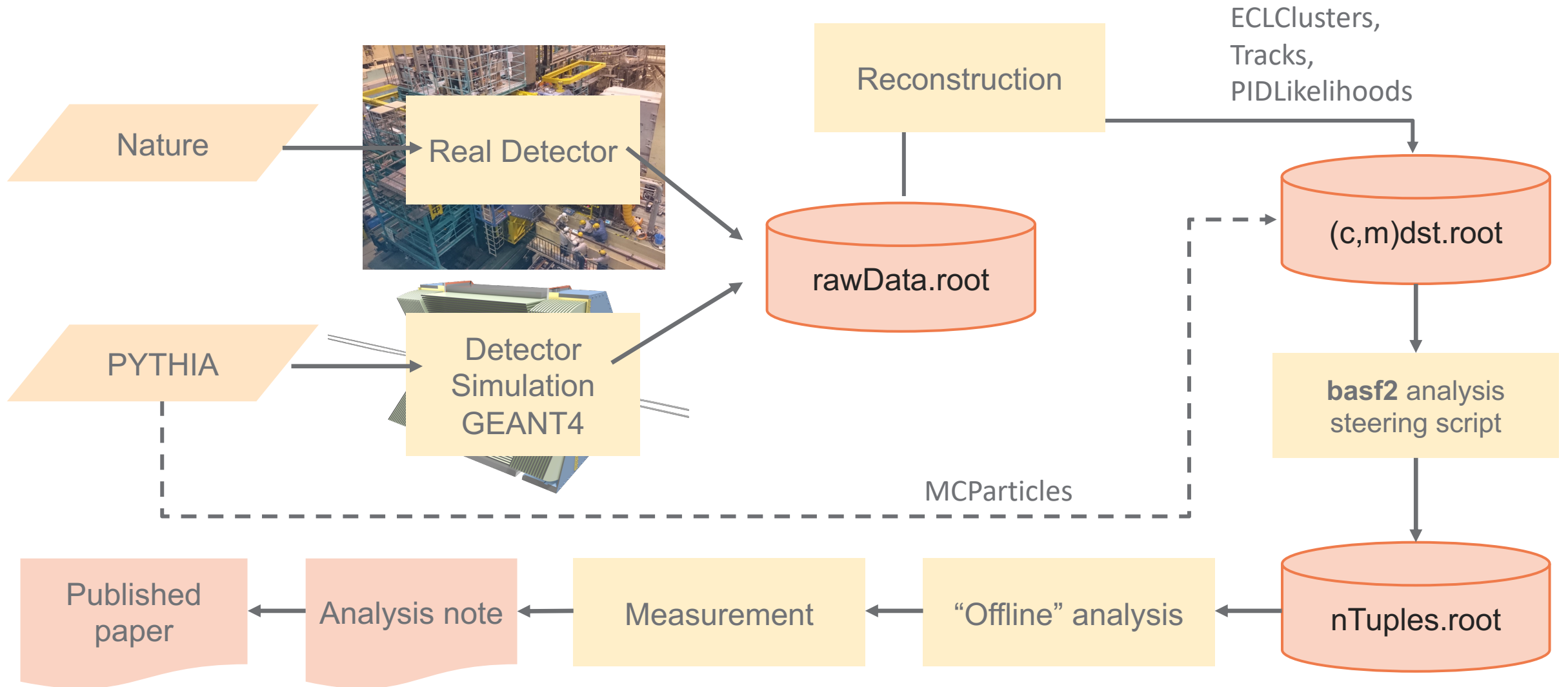




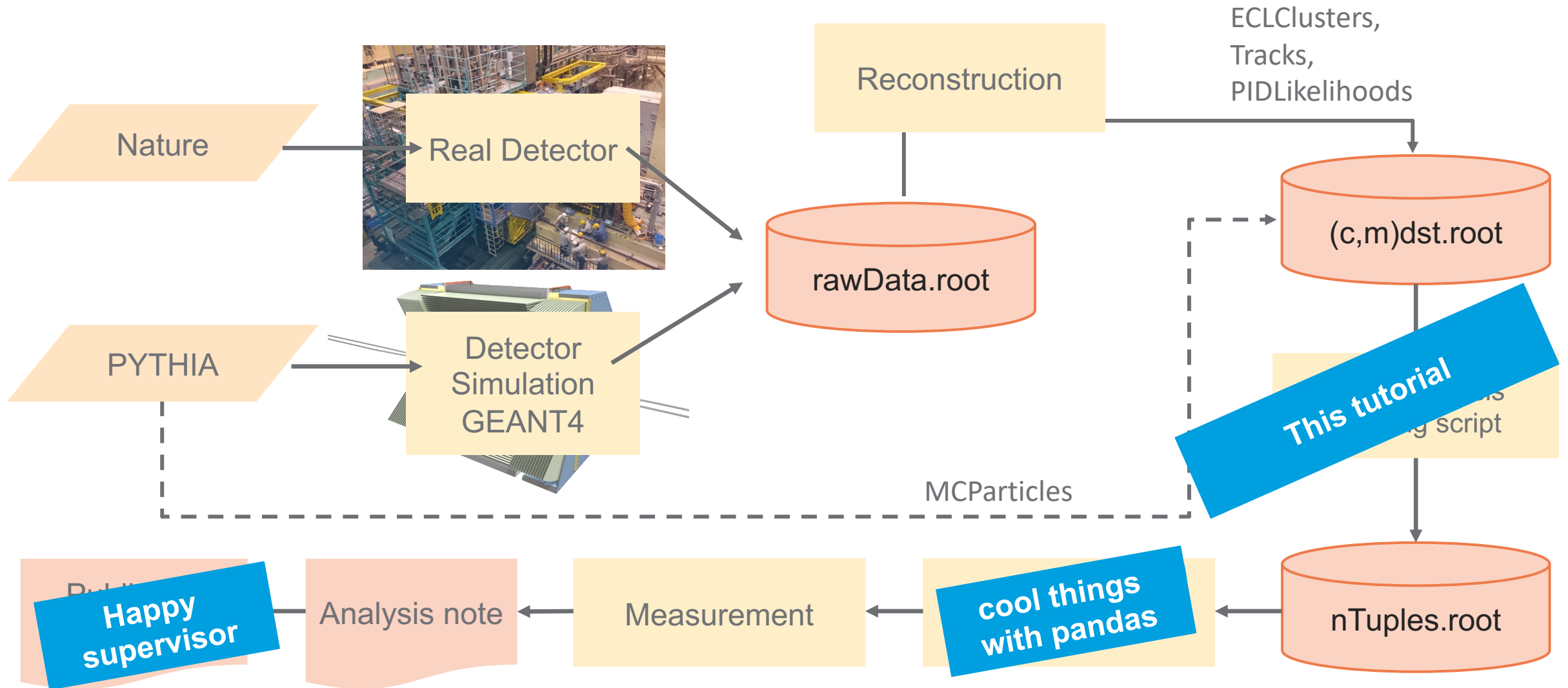
# The big picture



# The big picture



# The big picture



# What is basf2?

# The code

basf2 is C++14 "**under the hood**"

- Packages contain C++ **modules** to manipulate data.
- In analysis: we have code to build **particles** from primitive objects (like tracks and calorimeter clusters).
- We also calculate physics quantities, and apply cuts.

Python 3.6 code for **steering**

- Load and configure C++ modules
  - analysis modules and modules from other packages
- Also python does *some* high-level analysis tasks.
- You will write a fair bit of python during the workshop.



# The code

basf2 is C++14 "under the hood"

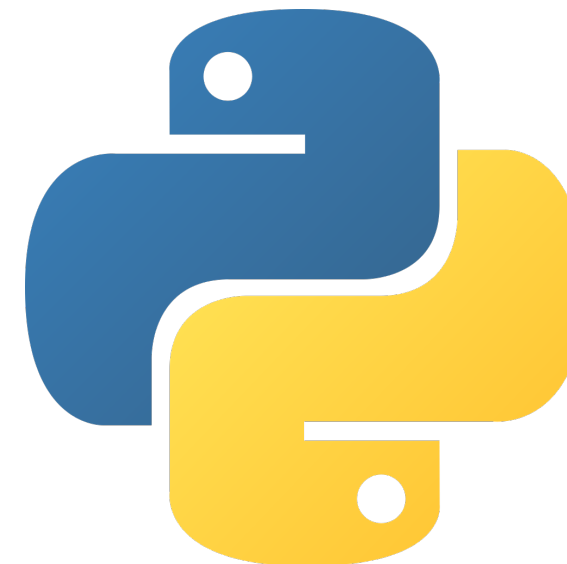
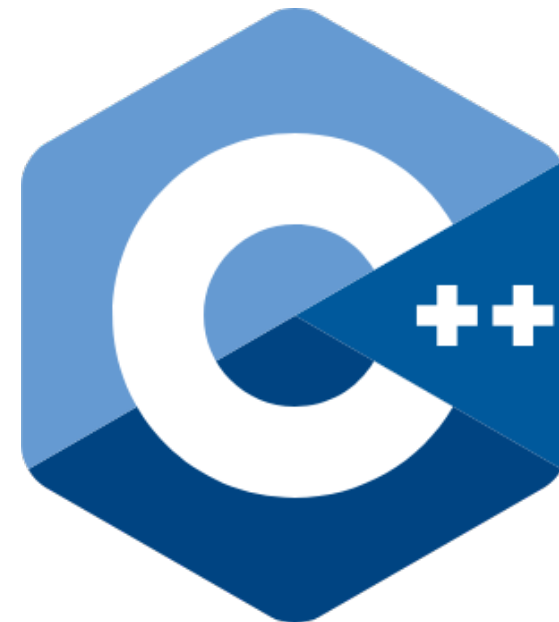
- Packages contain C++ **modules** to manipulate data.
- In analysis: we have **objects** from primitive objects (like tracks and calorimeters).
- We also calculate **observables** quantities, and apply cuts.

**Heavy lifting**

Python 3.6 code for **steering**

- Load and configure C++ modules
  - analysis modules and meta-packages
- Also python does **bookkeeping** tasks.
- You will write a **script** for python during the workshop.

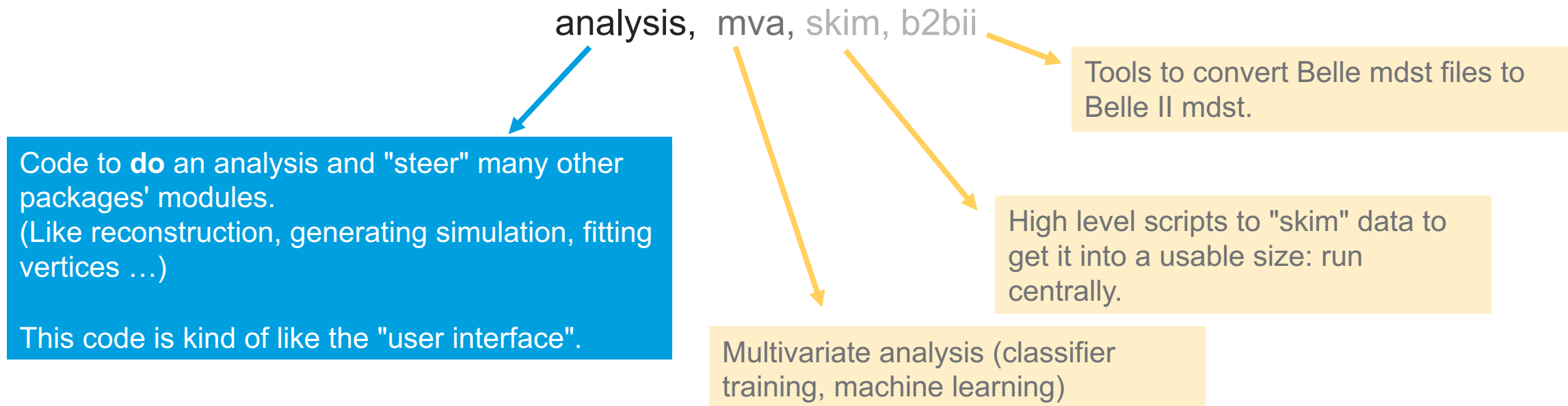
**Readable scripts**





# What is the analysis package?

- Our software is organised into "packages".
  - <https://stash.desy.de/projects/B2/repos/software/browse>
  - There are packages for subdetectors, tracking, simulation...
  - As a student/postdoc/collaborator you might work on some of them.
- BUT! When you want to do a physics measurement. You really only care about:



# I'm stuck, where do I go for help?

Probably the most important slide

## questions

for anything really, not just analysis package

<https://questions.belle2.org>

## documentation

there is fairly good documentation

<https://software.belle2.org>

## examples

so, for us <packagename> = analysis

`$BELLE2_RELEASE_DIR/<packagename>/examples`

## the code

<https://stash.desy.de/projects/B2/repos/software/browse>

`$BELLE2_RELEASE_DIR`

# I'm stuck, where do I go for help?

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so, for us <packagename> = analysis

`$BELLE2_RELEASE_DIR/<packagename>/examples`

## the code

[\\$BELLE2\\_RELEASE\\_DIR](https://stash.desy.de/projects/B2/repos/software/browse)



Go here first!

# Invoking basf2 and the equivalent of “Hello World”

# How to invoke basf2

## Belle analysis software framework 2

- Let's setup basf2 !
- But, first of all, you should know which release can be and should be used.

```
$ . /cvmfs/belle.cern.ch/tools/b2setup
$ b2setup --help # You can see available releases of basf2
...
  The following releases are available:
...
  release-04-01-00
  release-04-01-01
  light-1912-icarus

$ b2help-releases # or you can ask the recommended release
release-04-01-01
```

# How to invoke basf2

## Belle analysis software framework 2

Q. Which release should I use?

A. I really recommend you to use the latest `release-AA-BB-CC` or `light-YYMM-CODENAME`.

### `release-AA-BB-CC`

A full release, such as `release-04-01-01` and `release-03-02-04`.

If you execute ``b2help-releases``, it tells the latest full release.

### `light-YYMM-CODENAME`

A light release, such as `light-1912-icarus` and `light-1911-heracles`.

Made from only a few packages, analysis, mdst, skim, b2bii (from light-1912) etc..

They are suitable for doing analysis!



# How to invoke basf2

## Belle analysis software framework 2

- Hopefully you've seen this already.
- If not you will see it many times in this workshop.
- At the command line:

```
$ . /cvmfs/belle.cern.ch/tools/b2setup  
$ b2setup release-04-01-01  
$ basf2 --info
```

# How to invoke basf2

## Belle analysis software framework 2

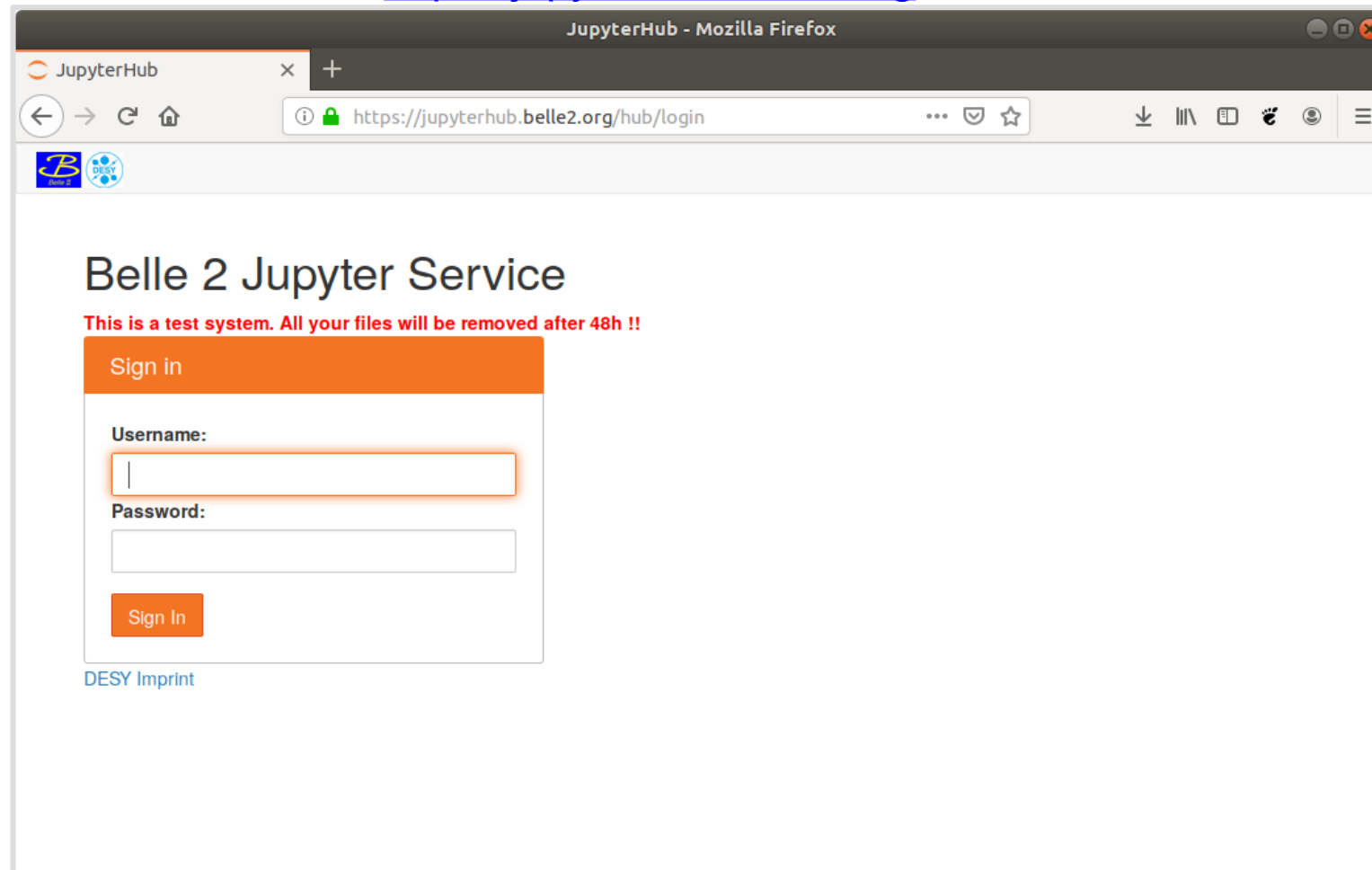
- Then Let's setup basf2 for the release!
- ``. /cvmfs/belle.cern.ch/tools/b2setup`` needs only once per session.

```
$  
$  
$ . /cvmfs/belle.cern.ch/tools/b2setup release-04-01-01  
$ basf2 --info
```

# Today we'll use jupyter

You will cover command line usage later in the workshop

<https://jupyterhub.belle2.org>



# Start a server, open the first exercise...

<https://jupyterhub.belle2.org>

B2T\_Basics\_1\_GettingHelp (autosaved)

File Edit View Insert Cell Kernel Widgets Help Not Trusted Python 3 (Belle2)

## BASF2 Basics

If you run things at the command line (for example at KEKCC or NAF) you will need to set up your environment and whatever release.

```
source /cvmfs/belle.cern.ch/tools/b2setup
b2setup release-03-01-01
```

Actually if you already know what release you want then you can do it all in once step:

```
source /cvmfs/belle.cern.ch/tools/b2setup release-03-01-01
```

To list all available releases:

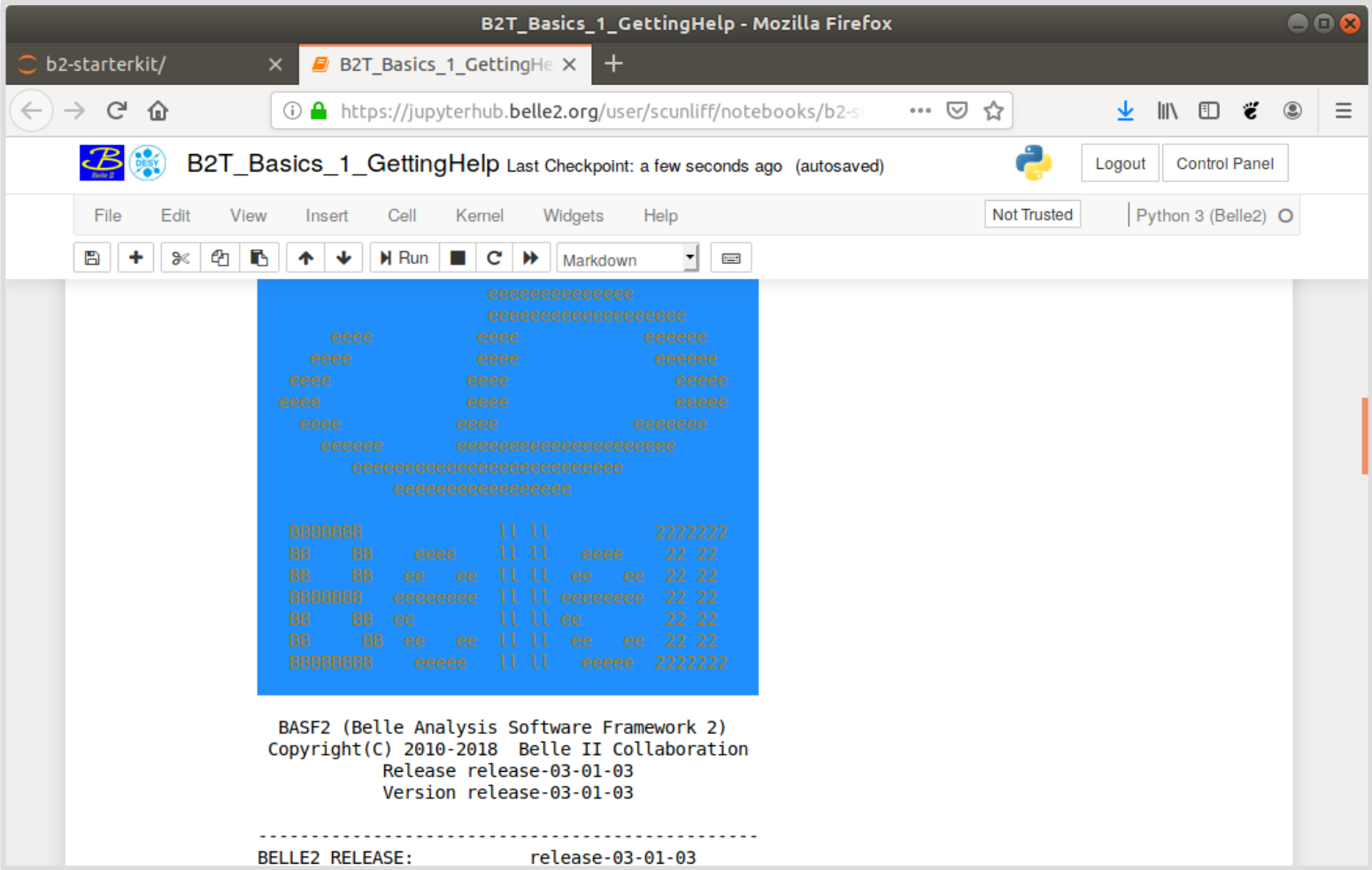
```
b2setup --help
```

You **don't** need to do that in this ipython/jupyter environment (we've set it all up for you). But please remember this, you'll need it later.

```
In [ ]: !basf2 --info
```

# basf2 --info

<https://jupyterhub.belle2.org>



# basf2 --info

Also works on a terminal to KEKCC or NAF

Source the logon script,  
pick a release

Execute this

```
Terminal
File Edit View Search Terminal Help
> source /cvmfs/belle.cern.ch/tools/b2setup release-03-01-03
Belle II software tools set up at: /cvmfs/belle.cern.ch/tools
Environment setup for release: release-03-01-03
Central release directory   : /cvmfs/belle.cern.ch/ubuntu1804/releases/release-03-01-03

> basf2 --info

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BASF2 (Belle Analysis Software Framework 2)
Copyright(C) 2010-2018 Belle II Collaboration
Release release-03-01-03
Version release-03-01-03

-----
BELLE2_RELEASE:      release-03-01-03
BELLE2_RELEASE_DIR:  /cvmfs/belle.cern.ch/ubuntu1804/releases/release-03-01-03
```



# basf2 --info

Also works on a terminal to KEKCC or NAF

Useful information

```
Terminal
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BASF2 (Belle Analysis Software Framework 2)
Copyright(C) 2010-2018 Belle II Collaboration
Release release-03-01-03
Version release-03-01-03

-----
BELLE2_RELEASE:      release-03-01-03
BELLE2_RELEASE_DIR:  /cvmfs/belle.cern.ch/ubuntu1804/releases/release-03-01-03
BELLE2_LOCAL_DIR:
BELLE2_SUBDIR:      Linux_x86_64/opt
BELLE2_EXTERNALS_VERSION: v01-07-01
BELLE2_ARCH:      Linux_x86_64
Default global tags:  release-03-01-00_rev3
Kernel version:      4.15.0-48-generic
Python version:      3.6.6
ROOT version:      6.14/06

basf2 module directories:
/home/scunliff
/cvmfs/belle.cern.ch/ubuntu1804/releases/release-03-01-03/modules/Linux_x86_64/opt
-----
>
```

# Modules, paths, the DataStore and how to steer them all

# Modules, paths, the DataStore

## What do we need to process the data?

Kuhr, Pulvermacher, Ritter, Hauth, Braun  
Comput. Softw. Big Sci. 3 (2019) no.1

1) A set of classes (modules) that process the data

→ **BASF2 module**

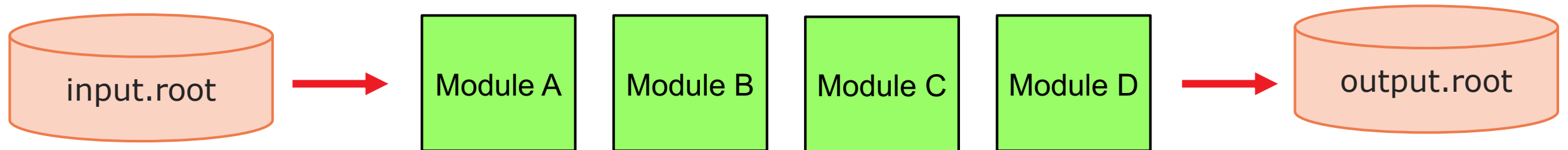
A module is written in C++ or Python and derived from a `Module` base class that defines the following interface methods:

- `initialize()`: called before the processing of events to initialize the module.
- `beginRun()`: called each time before a sequence of events of a new run is processed, e.g., to initialize run-dependent data structures like monitoring histograms.
- `event()`: called for each processed event.
- `endRun()`: called each time after a sequence of events of the same run is processed, e.g., to collect run-summary information.
- `terminate()`: called after the processing of all events.

# Modules, paths, the DataStore

What do we need to process the data?

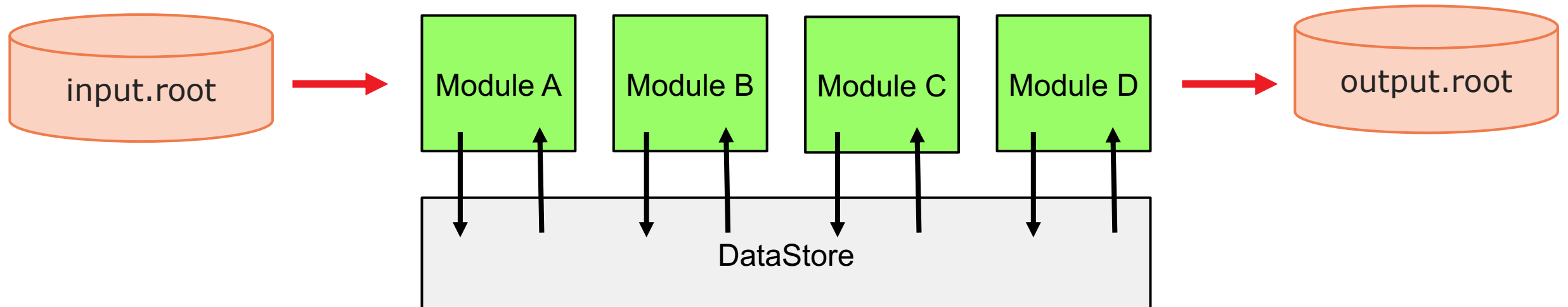
- 1) A set of classes (modules) that process the data  
→ **BASF2 module**



# Modules, paths, the DataStore

## What do we need to process the data?

- 1) A set of classes (modules) that process the data  
→ **BASF2 module**
- 2) A set of classes (dataobjects) that hold the data and allow module to pass thing one to the other  
→ **BASF2 dataStore**

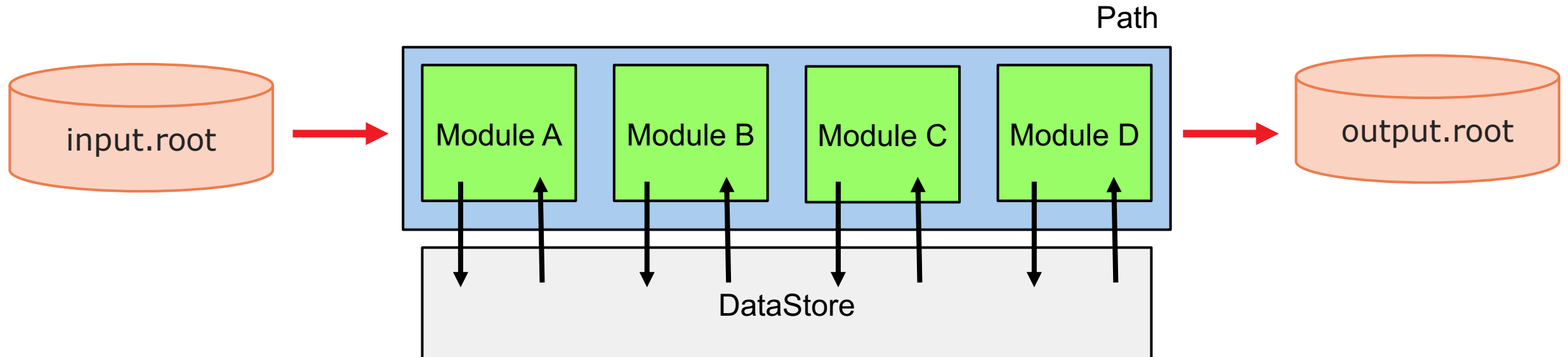


# Modules, paths, the DataStore

## What do we need to process the data?

1) A set of classes (modules) that process the data  
→ **BASF2 module**

2) A set of classes (dataobjects) that hold the data and allow module to pass thing one to the other  
→ **BASF2 dataStore**



3) An order in which the modules must be executed  
→ **BASF2 path**

# What does a steering file look like?

## How to I implement all this?

A steering file is a python script that runs

- the modules that you need
- in the order you need
- on the data you need

# A word about file types and why you should use the analysis package



# File types

## That basf2 can read and/or create

- A **dst** contains basf2 objects which will populate a **DataStore**.
  - **data summary table**
  - Basically: a special ROOT file.
- The data for physics analysis are "**mdst**"
  - **mini data summary table**.
  - Same structure of a **dst**, **but with much less information**
  - Input to your analysis package scripts
- The calibration & performance are "**cdst**"
  - **calibration data summary table**.
  - **mdst + digits**
- At the end of your analysis chain you will write out a "normal" root file containing a TTree, TNtuple, or histograms



A relevant question

<https://questions.belle2.org/question/219>

Objects allowed in an *mdst*:

<https://goo.gl/AB15Ud>

# Why use the analysis package?

## Can I read the mdst by myself?

mdst are basically root trees containing lists of:

- Track
- TrackFitResult
- V0
- PIDLikelihood
- ECLCluster
- KLMCluster
- KllD
- TRGSummary
- SoftwareTriggerResult
- (MCParticle)
- ...

The analysis package has modules to convert these into more friendly quantities like

- Particle
- ParticleList
- EventShapeContainer
- TagVertex
- ...

# Why use the analysis package?

Can I read the mdst by myself?

Can I open the mdst with my own, custom-made macro and run the analysis?

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Can I read the mdst by myself?

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The mdst contains also the relations between the objects stored in it, which are not trivially handled by a stand-alone root macro. **Use always basf2-based code.**

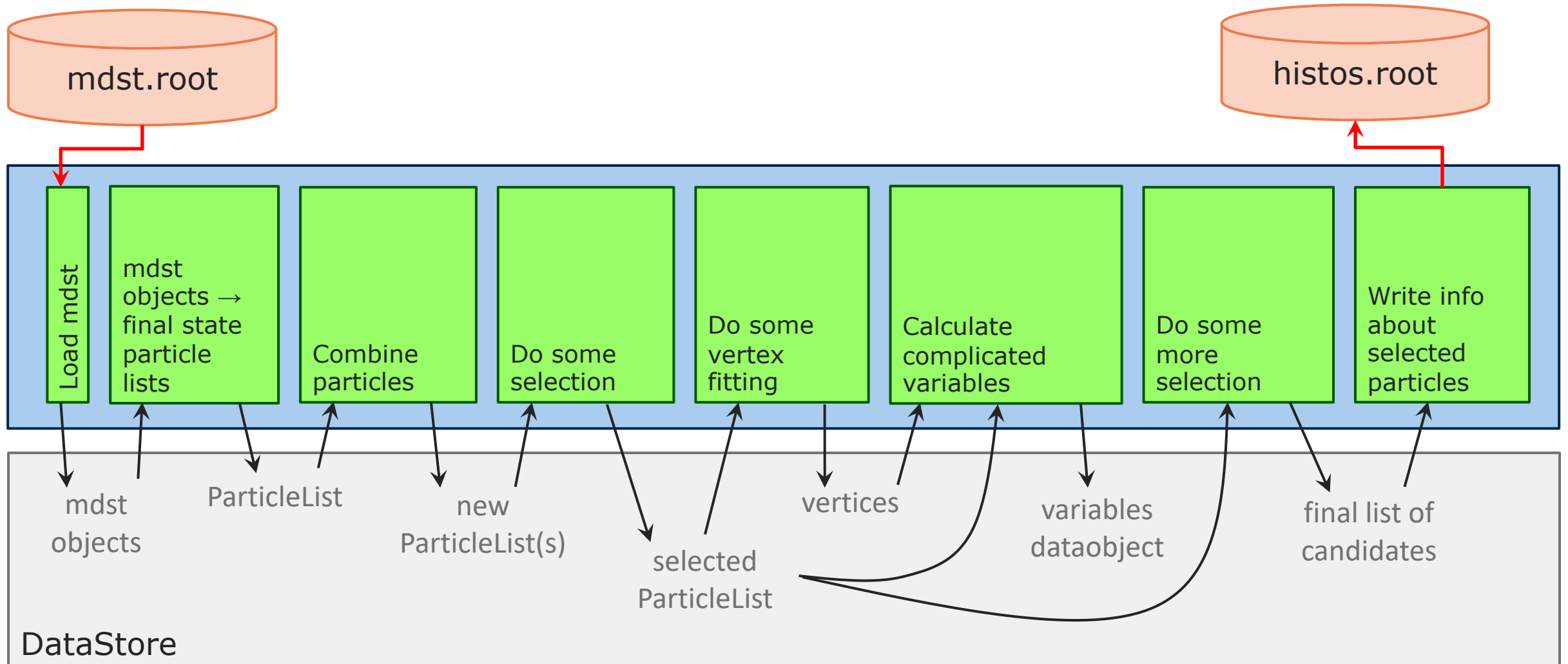
Should I write my own module that loops over reconstructed objects like the ECLClusters and do the analysis (i.e. Belle-style)?

**NO**

The relation between analysis object (particles) and the reconstructed objects is not always trivial.  
**One particle may have many trackFitResults**  
**The ECLClusters are not photons.**  
**Use the modules provided by a detector expert**

**Now let's step through an  
example**

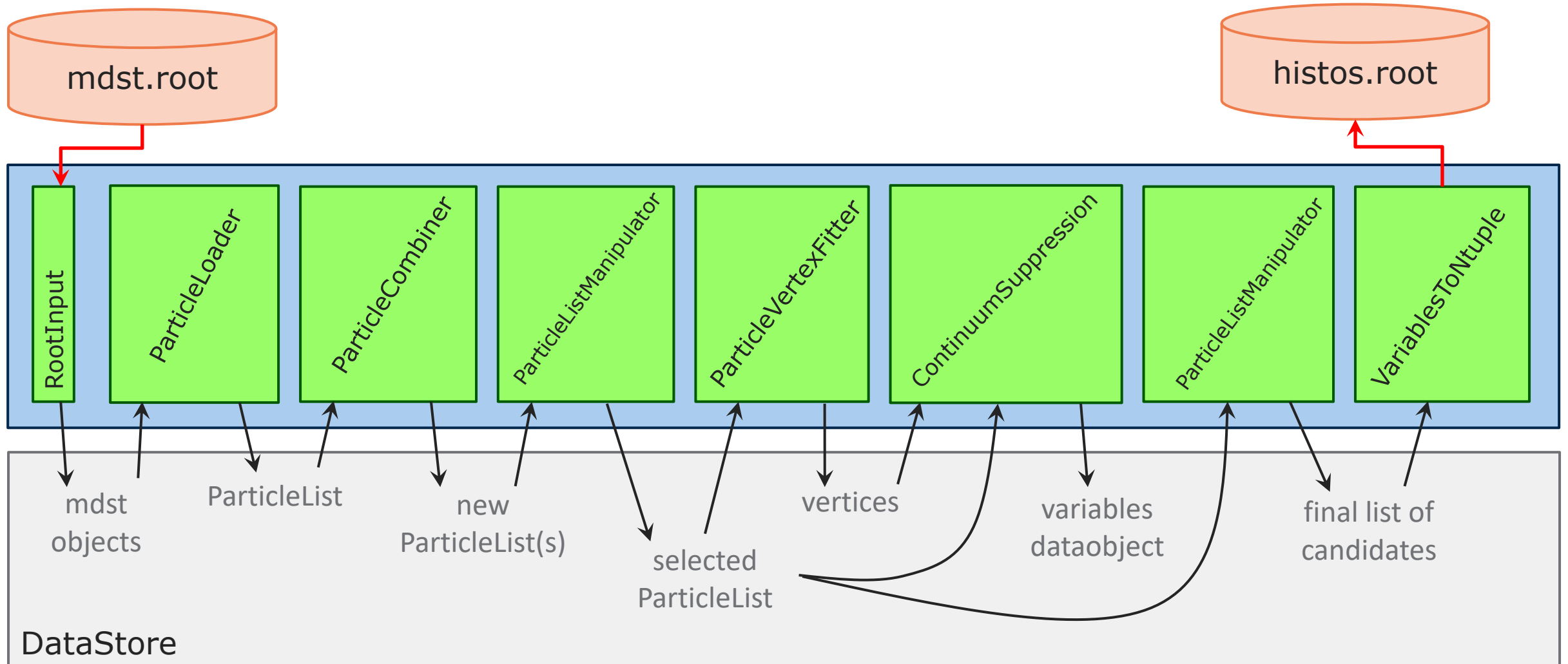
# A typical path for an analysis job





# A typical path for an analysis job

Now with the real names for the modules



# How to make a path and load a module

- First of all, you have make your path.

```
import basf2

mypath = basf2.Path()
# mypath basf2.create_path() # Both are OK!
```

# How to make a path and load a module

- Call a function to load a module and add it to your path.
- That is all you need to do!

```
from modularAnalysis import fillParticleList  
  
fillParticleList('pi+:highMom', 'p > 1', path=mypath)
```

Load a module to fill a list of particle, ParticleList.  
We'll see details of ParticleList later.

- What is happening in the function?

# The ParticleLoader

An analysis module which loads particles

```
fillParticleList('pi+:highMom', 'p > 1', path=mypath)
```



```
pload = register_module('ParticleLoader')  
pload.param('decayStringsWithCuts', [('pi+:highMom', 'p > 1')])  
mypath.add_module(pload)
```

- The function wraps these steps in a line.
- Of course, you can write them by hand.
  - **But quite quickly scripts become unreadable**

# The ParticleLoader

An analysis module which loads particles

- The procedure is same no matter what module you want in path.
  - "Register" the module you want
  - Set "parameters" to the module
  - "Add" the module to you path

```
pload = register_module('ParticleLoader')
pload.param('decayStringsWithCuts', [('pi+:highMom', 'p > 1')])
mypath.add_module(pload)
```

# How to find module documentation

## In general

- At the command line:

```
$ basf2 -m  
$ basf2 -m ParticleLoader
```

Take a look at  
B2T\_Basics\_1\_GettingHelp.ipynb  
on jupyterhub

Q: Can you find the **source code** for ParticleLoader?  
What do you notice about the C++ class name? How does this compare to  
the module name?

# Wrapper functions

- Functions such as `fillParticleList` are found in `<package>/scripts/*.py`
- The script that you will probably use a lot is **`analysis/scripts/modularAnalysis.py`**

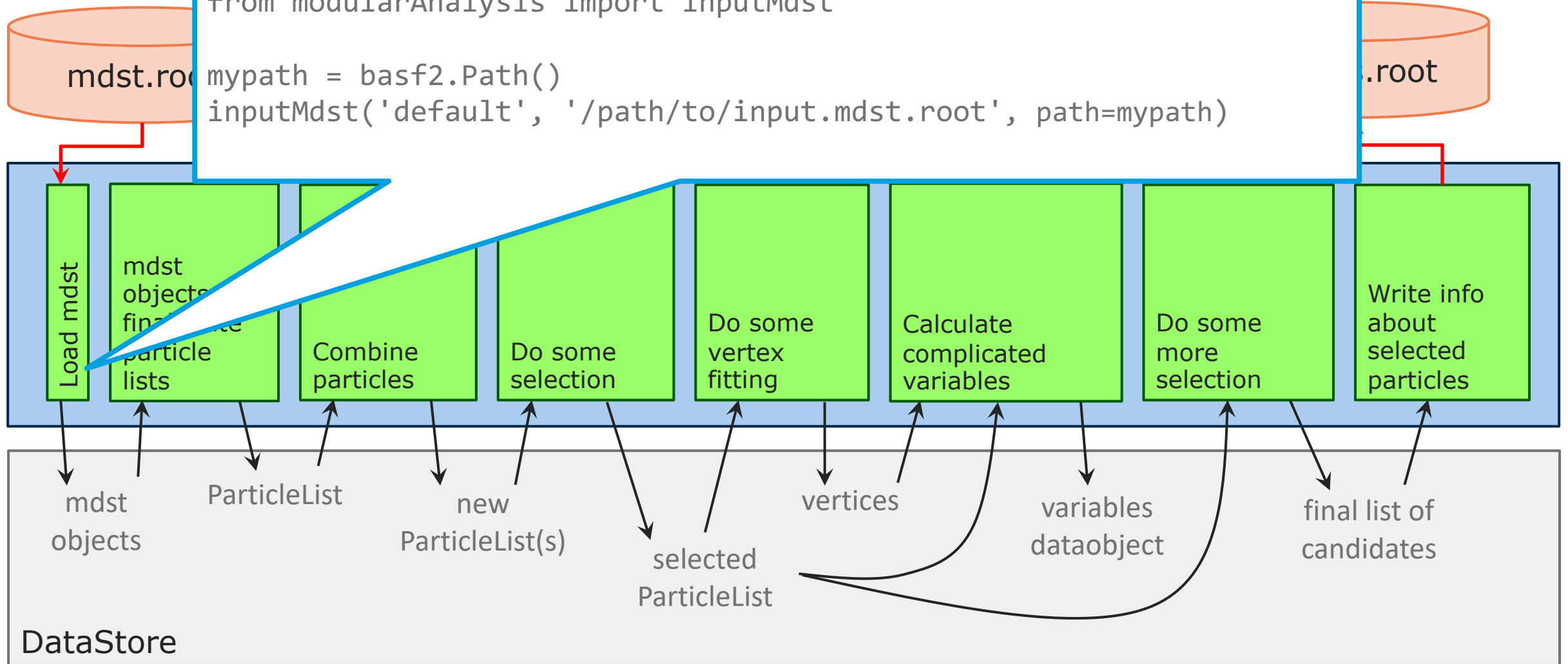
```
# analysis/scripts/modularAnalysis.py
from modularAnalysis import fillParticleList
fillParticleList('pi+:highMom', 'p > 1', path=mypath)
```

```
# analysis/scripts/vertex.py
from vertex import fitVertex
fitVertex('K*0:myKst', conf_level=0.0, path=mypath)
```

# A typical analysis workflow

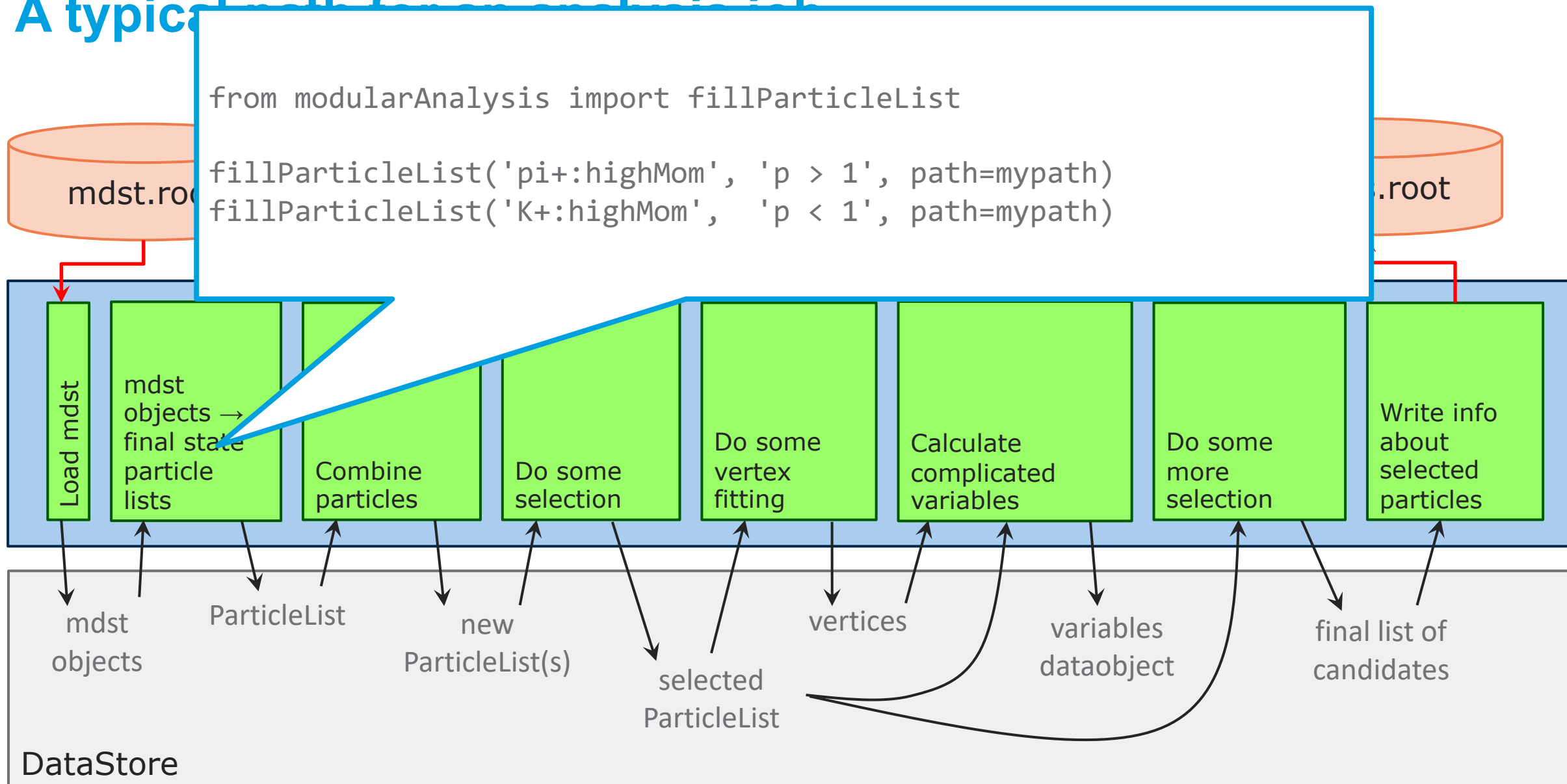
```
import basf2
from modularAnalysis import inputMdst

mdst.root mypath = basf2.Path()
inputMdst('default', '/path/to/input.mdst.root', path=mypath)
```



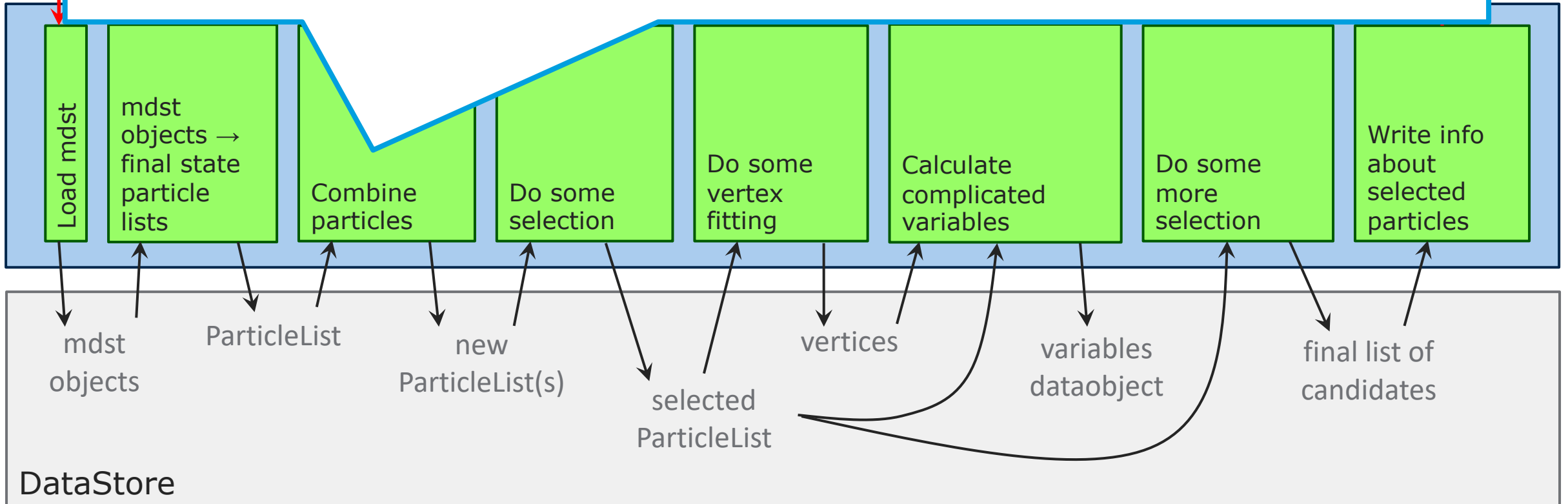


# A typical event analysis pipeline



A

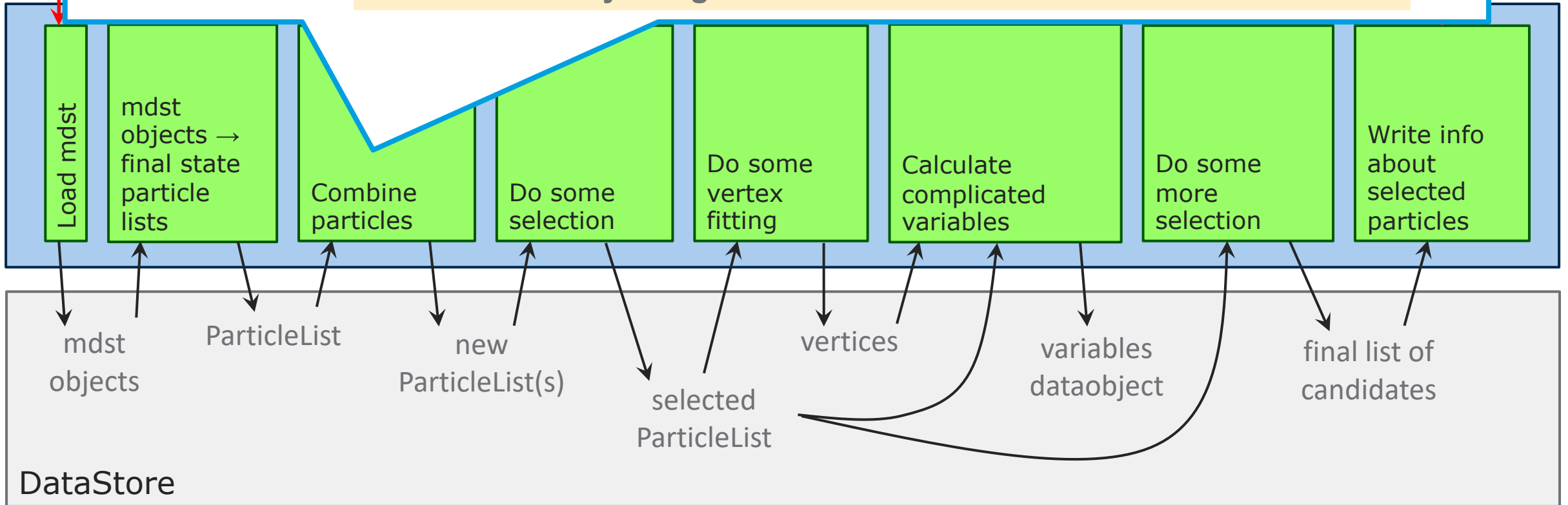
```
from modularAnalysis import reconstructDecay
reconstructDecay('K*0:myKst -> K+:highMom pi-:highMom', '0.6 < M < 1.0', path=mypath)
```



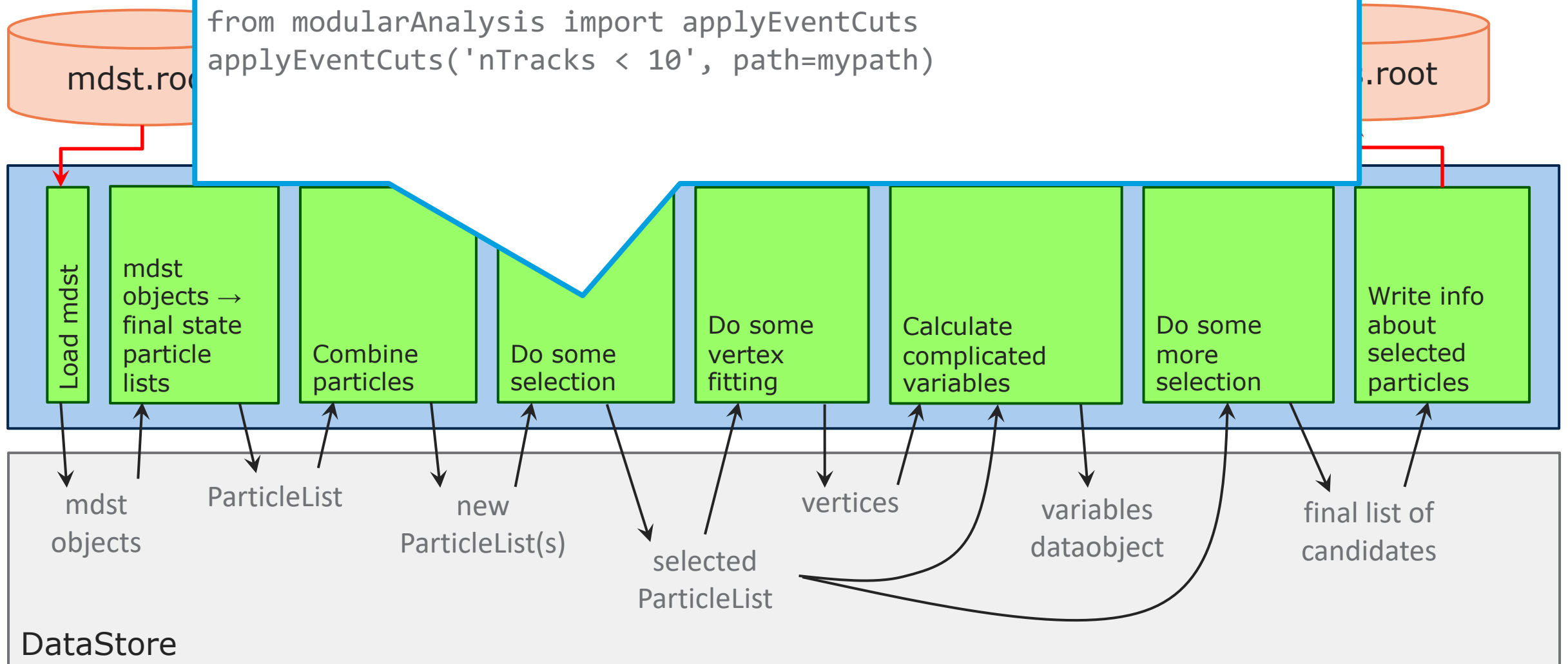
A

```
from modularAnalysis import reconstructDecay
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```

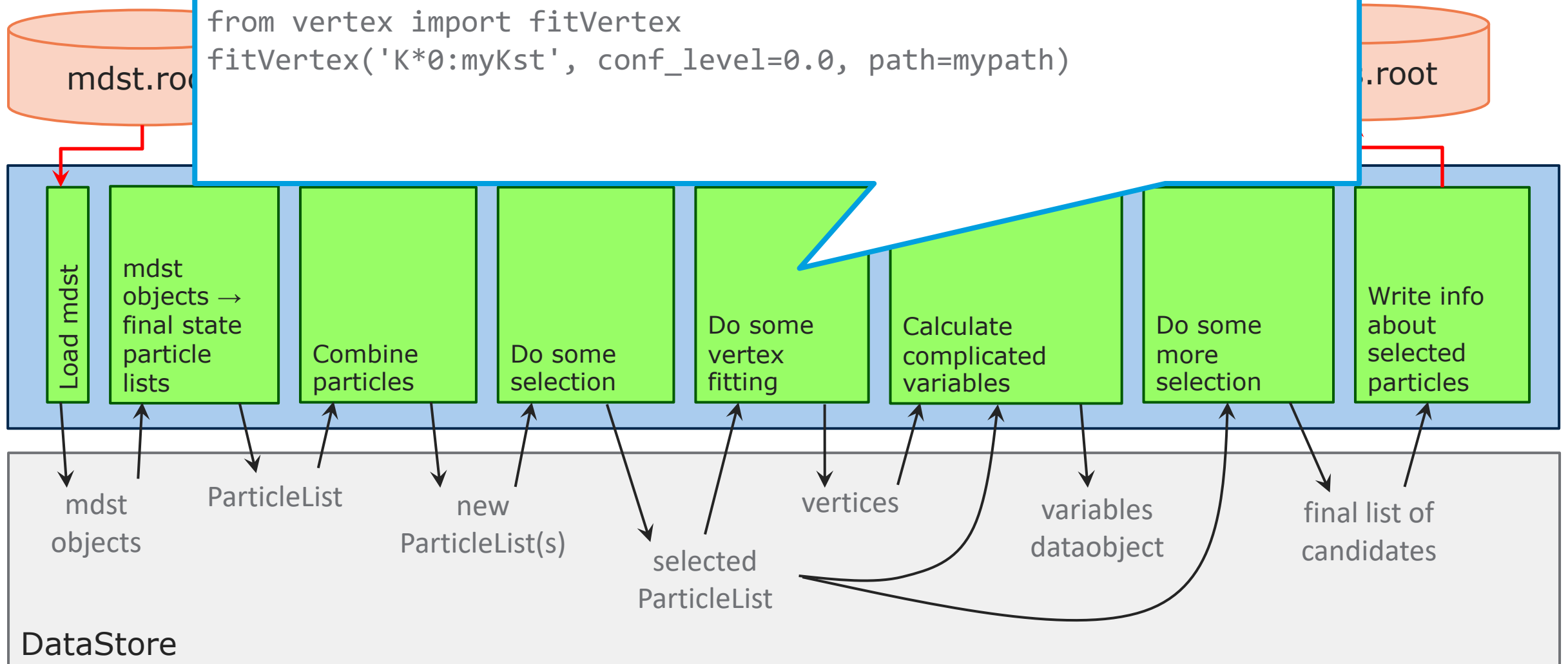
You can probably guess what this means already. But don't worry, we'll discuss the **DecayString** a bit later...



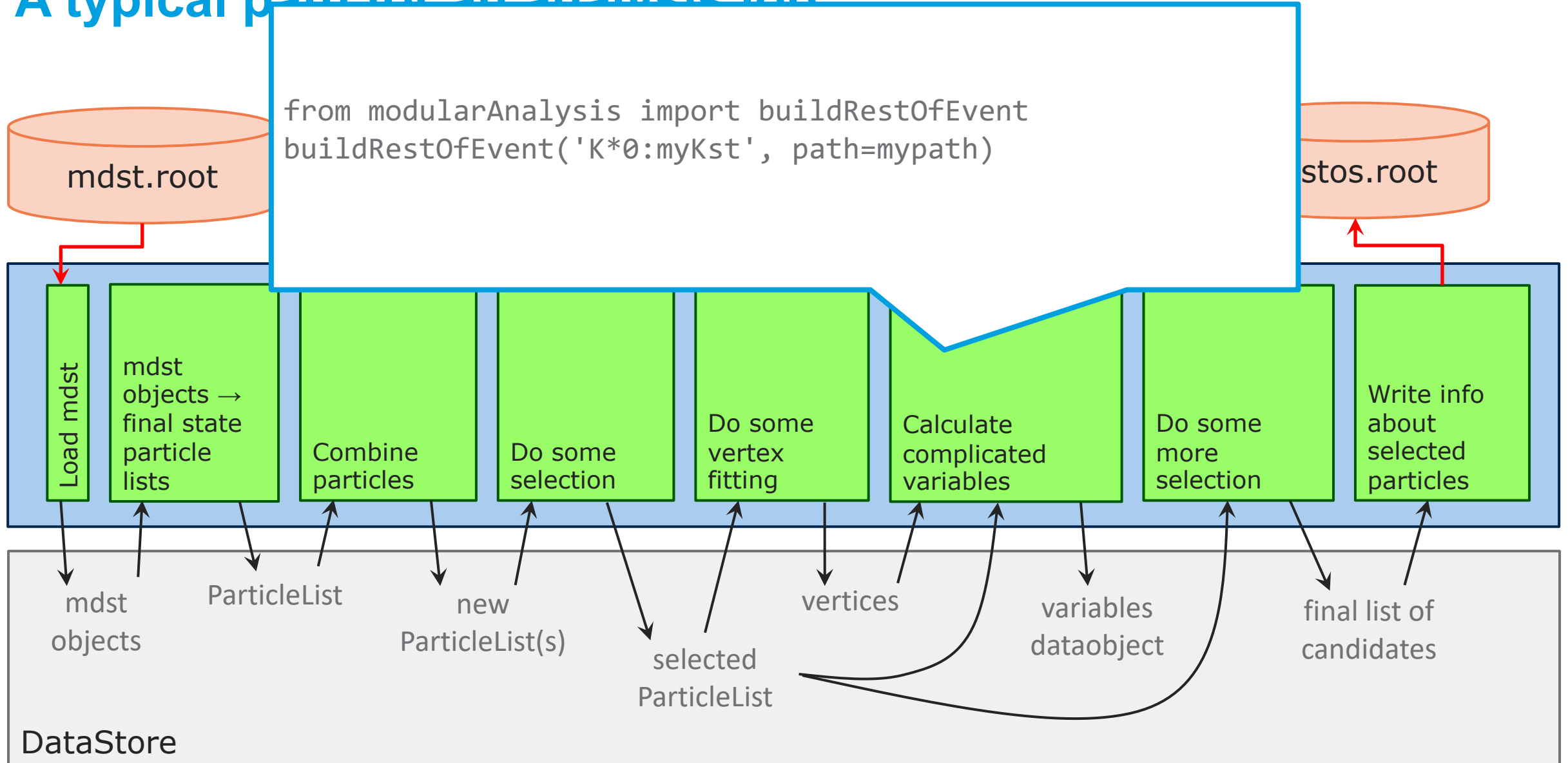
# A typical event analysis pipeline



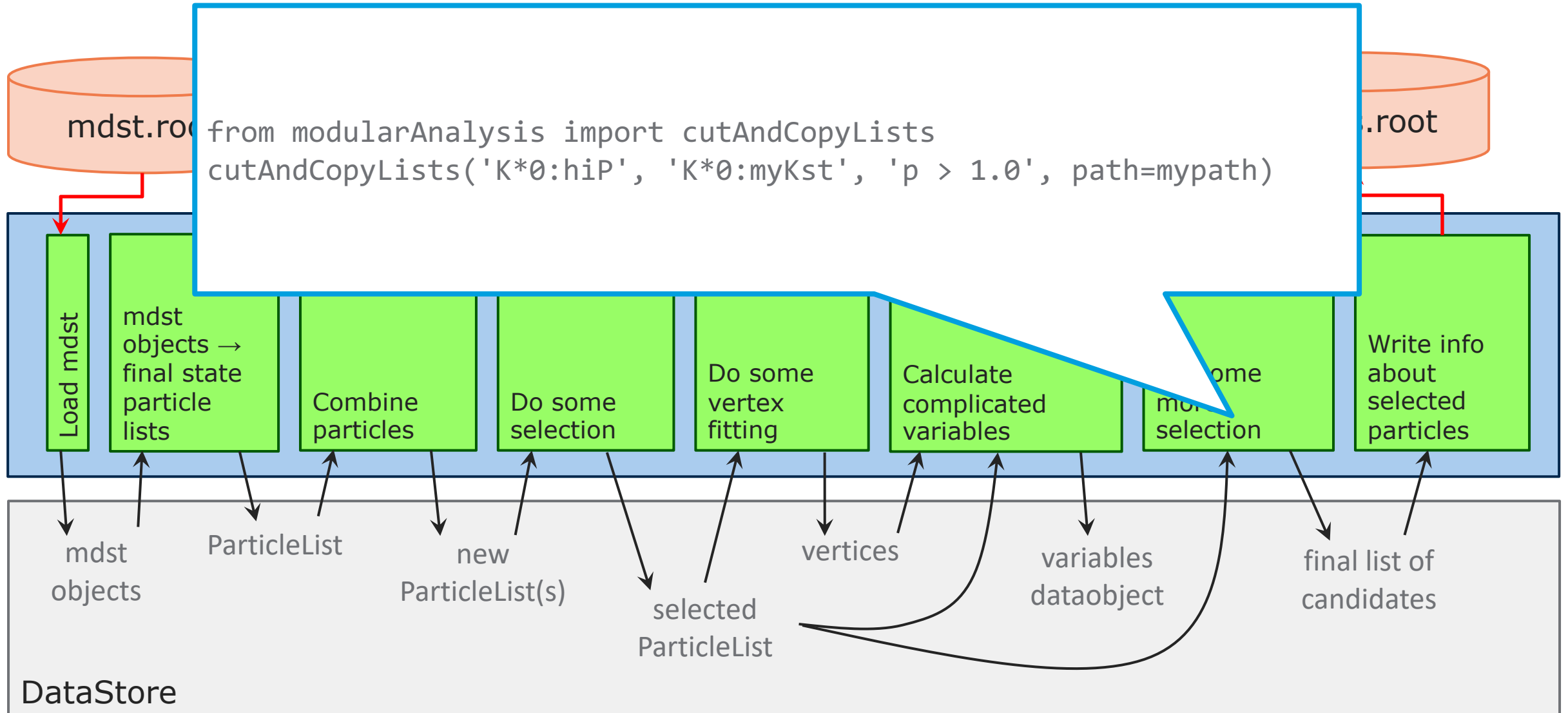
# A typical analysis workflow



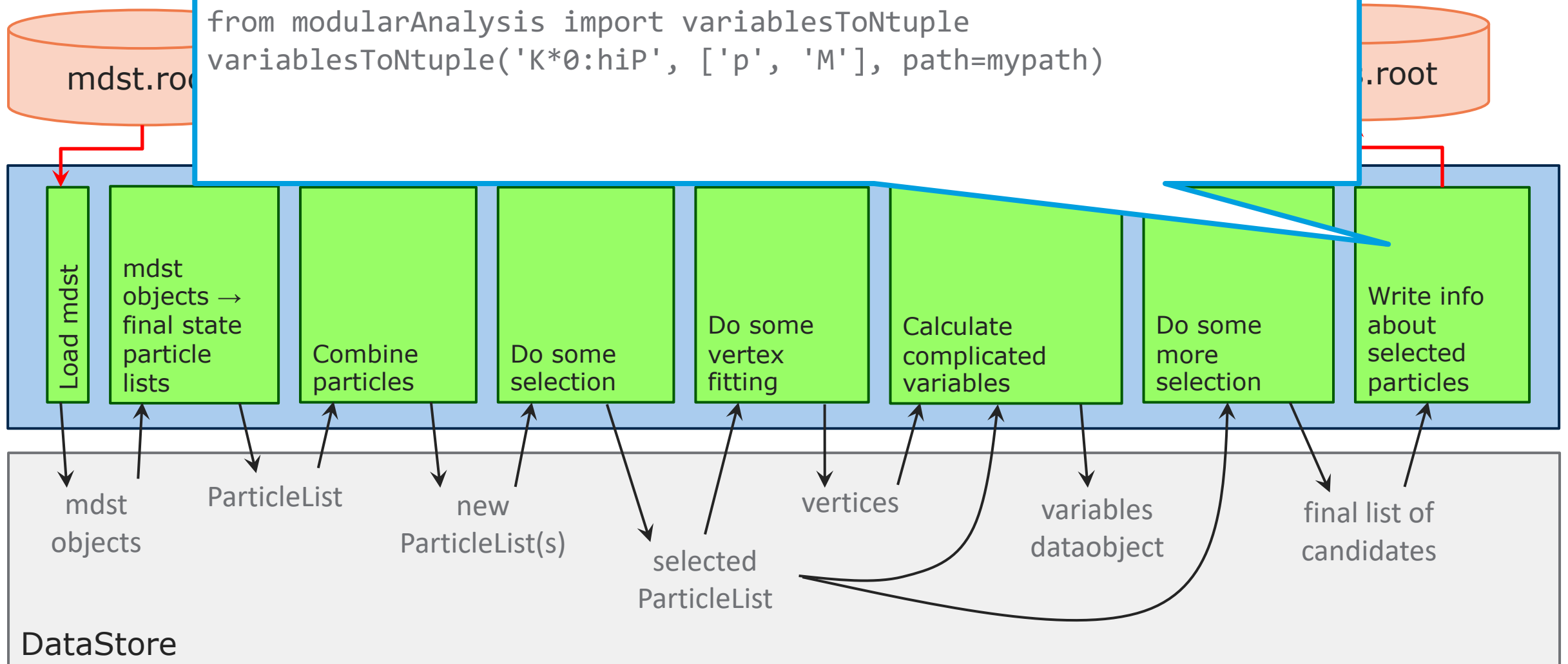
# A typical path for an analysis job



# A typical path for an analysis job



# A typical analysis workflow





# A template steering script

```
#!/usr/bin/env python3
# -*- coding: utf-8 -

import basf2
import modularAnalysis as ma

# create a path
mypath = basf2.Path()

# input mdst file
ma.inputMdst('default', 'inputMdstFile.root', path=mypath)

# PUT YOUR CODE HERE

# process the events
basf2.process(mypath)
# print out the summary
print(basf2.statistics)
```



# A template steering script

```
#!/usr/bin/env python3
# -*- coding: utf-8 -

import basf2
import modularAnalysis as ma

# create a path
mypath = basf2.Path()

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# PUT YOUR CODE HERE

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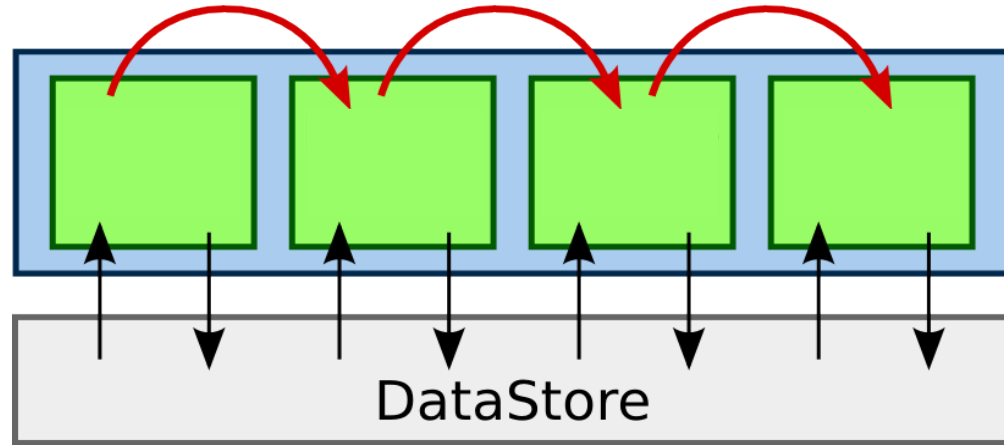


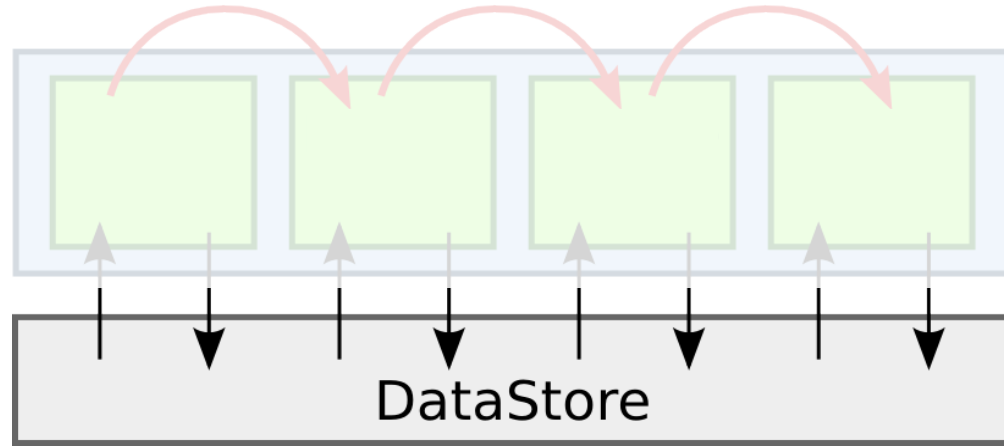
Q: what does 'default' mean?

What is the argument called?  
Can you find the documentation?  
Is there a questions post?

# Candidate/particle based analysis

# Before we get started, here's this diagram again





You can take a look at examples in `<package>/dataobjects`

- The datastore contains the **dataobjects**
- At the level of analysis, the main dataobject is: **ParticleList**

# Particle-based paradigms

In the wild

**Event based analysis  
frameworks**  
Babar, ATLAS, ILC

**Particle/candidate  
based analysis  
frameworks**  
LHCb, Belle II

# Particle-based analysis

- Take particle lists
- Build up decay parents from daughters
- Make ***candidates*** for your decay of interest
- Filter/cut/keep.
  
- You might have more than one candidate per event.
  - We deal with this after the fact.
  - This is fine. I promise.
  - <https://arxiv.org/abs/1703.01128>

# The Particle class

It's not crucial to understand the details

- A common representation of all particle types
  - Charged:  $e / \mu / \pi / K / p / d$  [built up from track + hypothesis]
  - $\gamma$  [built up from ECLClusters + !Track]
  - $K^0_L, n$  [built up from KLMLusters + ECLClusters + !Track ]
  - $K^0_S, \Lambda^0, \gamma$  [built up from V0 (2 tracks)]
  - Composite particles:  $\pi^0 / K^0_S / D / B$  [built up from combinations]
- Data members of the class are **common to all particle types**: mass, momentum, position, PDG code, ...
- Information which is only relevant to certain kinds of particle is saved in separate **analysis package dataobjects** and accessible by **relations**.
  - e.g. ContinuumSuppression
  - .... FlavorTaggerInfo



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Q: what is the name of the module to combine particles?

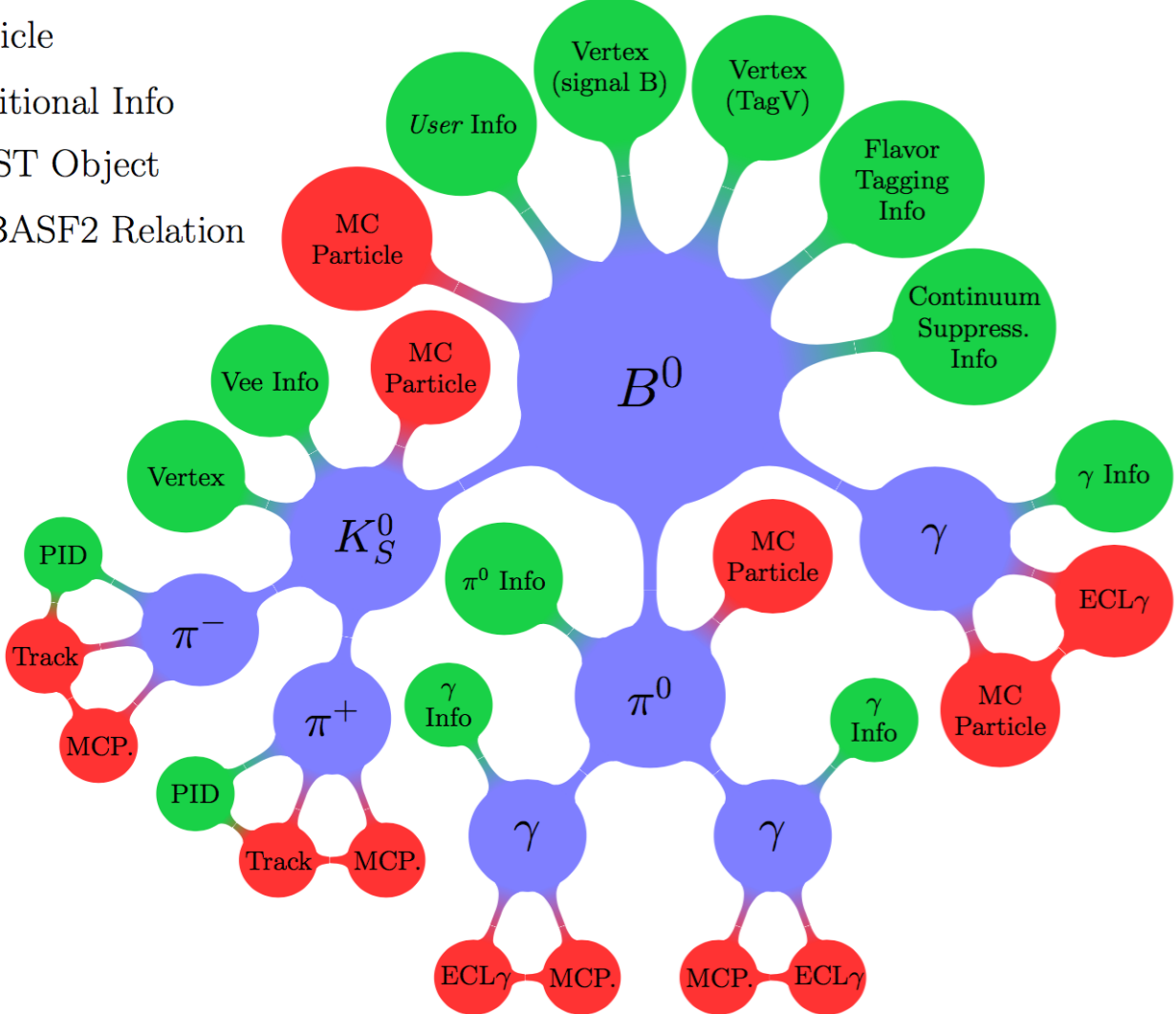
Q: which things on this slide are dataobjects?

# ParticleList

- A group of all particles and anti-particles that belong together logically.
  - e.g.  $K^{*0}$  s (decaying to  $K^\pm$  and  $\pi^\mp$  with invariant mass in a certain window)
- Can only store particles of the same PDG code (can be different decay modes).
- Doesn't have ownership of the **Particle** objects.
- **ParticleList** is the **dataobject** on which analysis modules operate.
  
- The physics-performance group provides **Standard Particle Lists** for which quality benchmarks exist, and systematics will be provided.
  - Currently you may need to optimize selection criteria of ParticleLists by yourself.
  - This will be really recommended to use in future.

# Some more details on Particles, ParticleLists and Relations

- Particle
- Additional Info
- MDST Object
- BASF2 Relation



- At each stage we build relations between the dataobjects
- Like vertex information, ContinuumSuppression → all related to Particles
- Particles themselves related to primitive mdst objects (clusters, tracks)

# Fill a ParticleList

There are two possible ways to fill a list of **stable\*** particles

1) fill it by hand

```
import modularAnalysis as ma

ma.fillParticleList('pi+:highMom', 'p > 1', path=mypath)
```

Q: What is a stable particle in this context?

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```

Particle name.

Tells basf2 if the list has to be created from tracks, ELCClusters, KLMClusters or V0

Must given in evt.pdl (or b2help-particles)

```
evt.pdl = $BELLE2_EXTERNALS_DIR/share/evtgen/evt.pdl
```

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Must given in evt.pdl (or b2help-particles)

Label name.

Used to distinguish different lists of the same particle type: you choose the name.

(But maybe call it something helpful)

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(But maybe call it something helpful)

ParticleList name

Same label name can be set to another particle name.  
e.g. pi+:highMom, K+:highMom

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Label name.

Used to distinguish different lists of the same particle type: you choose the name.  
(But maybe call it something helpful)

Selection criteria

Cuts used to select this list. **More on this later**

List name

Same label name can be set to another particle name.  
e.g. pi+:highMom, K+:highMom

```
evt.pdl = $BELLE2_EXTERNALS_DIR/share/evtgen/evt.pdl
```



# Fill a ParticleList

There are two possible ways to fill a list of **stable\* particles**

1) fill it by hand

```
import modularAnalysis as ma  
  
ma.fillParticleList('pi+:highMom', 'p > 1', path=mypath)
```

Once you call the function, pi-:highMom is created automatically.  
You **don't** have to call

```
ma.fillParticleList('pi-:highMom', 'p > 1', path=mypath)
```

```
evt.pdl = $BELLE2_EXTERNALS_DIR/share/evtgen/evt.pdl
```

# Fill a ParticleList

There are two possible ways to fill a list of **stable\*** particles

1) fill it by hand

```
import modularAnalysis as ma

ma.fillParticleList('pi+:highMom', 'p > 1', path=mypath)
```

2) Use the **standard particles lists** with pre-defined names and cuts

```
from stdPhotons import stdPhotons
from stdCharged import stdK

stdPhotons('all', path=mypath) # --> gamma:all
stdK('95eff', path=mypath)     # --> K+:95eff
```

First argument will be a label name of the particle list.

```
evt.pdl = $BELLE2_EXTERNALS_DIR/share/evtgen/evt.pdl
```

# DecayString

The last of the infrastructure things, I promise

- You've already seen the decay string in action.
- Now we can formally describe what this line is doing:

```
reconstructDecay('K*0:myKst -> K+:highMom pi-:highMom', '0.6 < M < 1.0', path=mypath)
```

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Decay string

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ParticleList ParticleList ParticleList

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```

Arrow  
(Indicate K\*0 decays to K+ pi-)

Decay string

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```
reconstructDecay('K*0:myKst -> K+:highMom pi-:highMom '0.6 < M < 1.0', path=mypath)
```

```
mother arrow daughter0 daughter1 ...
```

Decay string

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- Now we can formally describe what this line is doing:

```
reconstructDecay(K*0:myKst -> K+:highMom pi-:highMom' '0.6 < M < 1.0', path=mypath)
```

Particle name.  
Of your composite particle.  
Must given in evt.pdl  
(or b2help-particles)

Decay string

```
evt.pdl = $BELLE2_EXTERNALS_DIR/share/evtgen/evt.pdl
```



# DecayString

The last of the infrastructure things, I promise

- You've already seen the decay string in action.
- Now we can formally describe what this line is doing:

```
reconstructDecay(K*0:myKst -> K+:highMom pi-:highMom '0.6 < M < 1.0', path=mypath)
```

You can choose this.

List name that exist already

Decay string

- `reconstructDecay()` : Reconstruct mother particle from daughters with a given DecayString. List names for daughters must exist already. You can choose a list name of a mother.

You don't have to call

```
reconstructDecay('anti-K*0:myKst -> K-:95eff pi+:95eff', '0.6 < M < 1.0', path=mypath)
```

# DecayString

## One final thing...

- Sometimes you will need to select a particle within a decay string.
- This is done by "carat": ^

```
'K*0:myKst -> ^K+:highMom ^pi-:highMom'
```

- You will see this in action in the exercises later.

# Physics quantities and the VariableManager

# The VariableManager

## It manages variables

- VariableManager is a place in the analysis package to store variables
  - physics quantities: invariant mass, beam-constrained mass,  $E$ ,  $p$ ,  $p_T$ ,  $\theta$ ,  $\phi$ , highest energy in a cluster
  - counters: event\_number, nhits, i\_candidate
  
- Every variable takes at least a Particle\* as input and returns a double (even integer counters like event\_number where this doesn't make much sense)



# The VariableManager

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  - physics quantities: invariant mass, beam-constrained mass,  $E$ ,  $p$ ,  $p_T$ ,  $\theta$ ,  $\phi$ , highest energy in a cluster
  - counters: event\_number, nhits, i\_candidate
- Every variable takes at least a Particle\* as input and returns a double (even integer counters like event\_number where this doesn't make much sense)
- You've already seen it in **python**, when we used a cut on "M".
  - This comes from the VariableManager.

```
reconstructDecay('K*0:myKst -> K+:highMom pi-:highMom', '0.6 < M < 1.0', path=mypath)
```



# Where can I get the documentation?

- At the command line:

```
$ basf2 variables.py
```

- Online, the documentation is [software.belle2.org](https://software.belle2.org).

Take a look at  
B2T\_Basics\_1\_GettingHelp.ipynb on  
jupyterhub

Q: What is the beam-constrained mass,  $M_{bc}$   
called in the VariableManager?

# Have you been listening?

Q: Can you figure out the cut string and code to reconstruct  $B^0 \rightarrow K^0 \gamma$ ?  
Supposing you have already reconstructed  $K^0$ :myKst.

```
import modularAnalysis as ma

ma.fillParticleList('pi+:highMom', 'p > 1', path=mypath)
ma.fillParticleList('K+:highMom', 'p < 1', path=mypath)
ma.reconstructDecay('K*0:myKst -> K+:lowMom pi-:highMom', '0.6 < M < 1.0', path=mypath)
```

Q: How can we fill a particle list of gamma?

Q: How can we reconstruct  $B^0$  from  $K^0$  and gamma?

Q: Can you apply following selection on the  $B^0$  candidates?

$$M_{bc} > 5.2 \text{ GeV}/c^2$$

# Aliases

Are awesome and you should use them

- With more advanced variables, we deliberately give them verbose names in the VariableManager.
- It's important to be clear.
- Some examples:
  - `cosAngleBetweenMomentumAndVertexVector`
  - `totalPhotonEnergyOfEvent`
  - `trackFindingFailureFlag`
- You can also use `formula()`, `abs()`, `cos()` for simple math:
  - `formula(missingEnergyOfEventCMS/Ecms)`
  - ...



# Aliases

Are awesome and you should use them

- For your code, you can make short aliases.
- This makes your offline data more manageable, and can help with code readability.

```
from variables import variables as vm
vm.addAlias('cosPVtx',
           'cosAngleBetweenMomentumAndVertexVector')
```

There is no path here like path=myspath.  
The VariableManager exists alongside the path

# How to get data out

# Flat ntuples and histograms

## Getting output data

There are some modules to store variables.

### VariablesTo\*

- a) VariablesToNtuple
- b) VariablesToHistogram
- c) VariablesToEventBasedTree

# Flat ntuples and histograms

## Getting output data

There are some modules to store variables.

### VariablesTo\*

- a) VariablesToNtuple
- b) VariablesToHistogram
- c) VariablesToEventBasedTree

```
from modularAnalysis import variablesToNtuple  
from modularAnalysis import variablesToHistogram
```

# VariablesToNtuple

- You want to store physics quantities for one ParticleList from the VariableManager.
- Get an ttree (ntuple) of candidates (one row per candidate).

# VariablesToHistogram

- Perhaps you just want a quick histogram.
- Candidate information not preserved.

# VariablesToNtuple

```
# ... at the end of your script
from modularAnalysis import variablesToNtuple
variablesToNtuple('K*0:myKst', ['M', 'p', 'daughter(0,p)', 'daughter(1,p)'], path=mypath)

# process the events
basf2.process(mypath)
```

Daughter's variables can be stored with `daughter(i, variable)`.  
`i` is the daughter index, i.e. `0 = K+`, `1 = pi-`.

# VariablesToHistogram

```
# ... at the end of your script
from modularAnalysis import variablesToHistogram
variablesToHistogram('K*0:myKst',
                    [ ('M', 100, 0.7, 1.0),
                      ('p', 100, 0.1, 3.0) ], path=mypath)

# process the events
basf2.process(mypath)
```

# VariablesToNtuple

```
# ... at the end of your script
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                      ('p', 100, 0.1, 3.0) ], path=mypath)

# process the events
basf2.process(mypath)
```

Q: Can you guess what these numbers do?  
What is the (python) structure created by the parentheses?

# Utilities

## Helper functions to make daughter's aliases.

- You can store variables of daughters with `daughter(i, variable)` in a same ntuple.
- You may also want to short aliases for these variables.

```
reconstructDecay('K*0:myKst -> K+:highMom pi-:highMom', '0.6 < M < 1.0', path=mypath)

from variables import variables as vm
vm.addAlias('K_M', 'daughter(0,M)')
vm.addAlias('K_p', 'daughter(0,p)')
vm.addAlias('pi_M', 'daughter(1,M)')
vm.addAlias('pi_p', 'daughter(1,p)') # You may want to add more aliases

vars = ['M', 'p', 'K_M', 'K_p', 'pi_M', 'pi_p']
variablesToNTuple('K*0:myKst', vars, path=mypath)
```



# Utilities

## Helper functions to make daughter's aliases.

- A helper function creates aliases from a list of variables and a DecayString with carat '^'.

```
import variables.utils as vu
vars = vu.create_aliases_for_selected(['M', 'p', 'px', 'py', 'pz'],
                                     'K*0 -> ^K+ ^pi-')
# The function makes aliases for selected particles, for example,
# 'K_M' --> 'daughter(0,M)'
```

- All aliases can be checked with printAliases().

```
from variables import variables as vm
vm.printAliases()
# [INFO] =====
# [INFO] Following aliases exists:
# [INFO] 'K_M' --> 'daughter(0,M)'
# ...
```

# Nomenclature

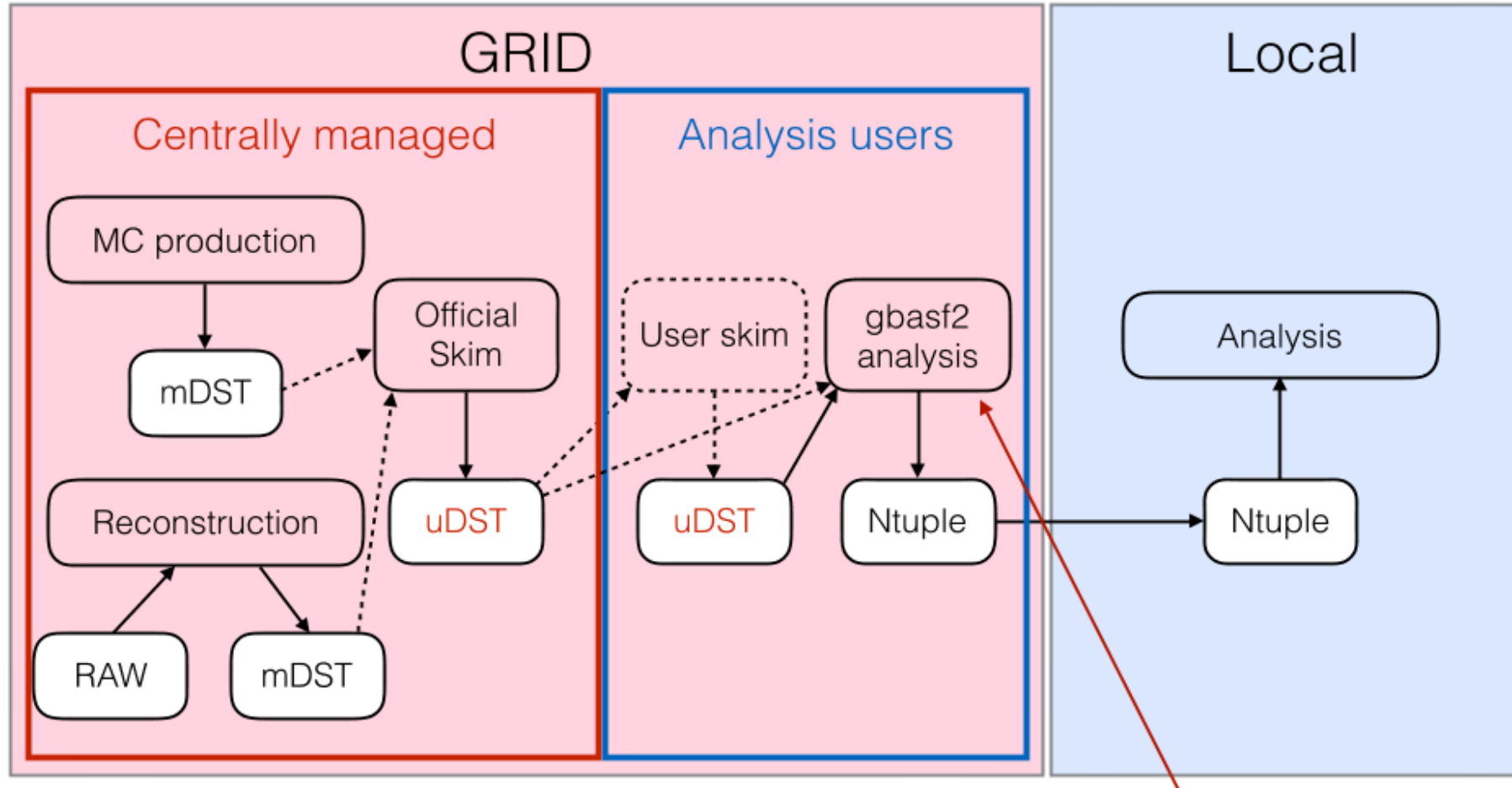
# Nomenclature

<https://confluence.desy.de/display/BI/Main+Glossary>

- **Experiment** (chunk of data-taking ~months).
- **Run** (chunk of data-taking w/ stable beams ~hours),
- Event.
  
- **TRG** the hardware trigger (group, device, DAQ)
- **L1** the hardware trigger (used interchangeably)
  
- SoftwareTrigger / **HLT** (the software trigger)
  
- **basf2** “Belle 2 analysis software framework” “the software”
- **gbasf2** “The grid job submission tool” “computing”

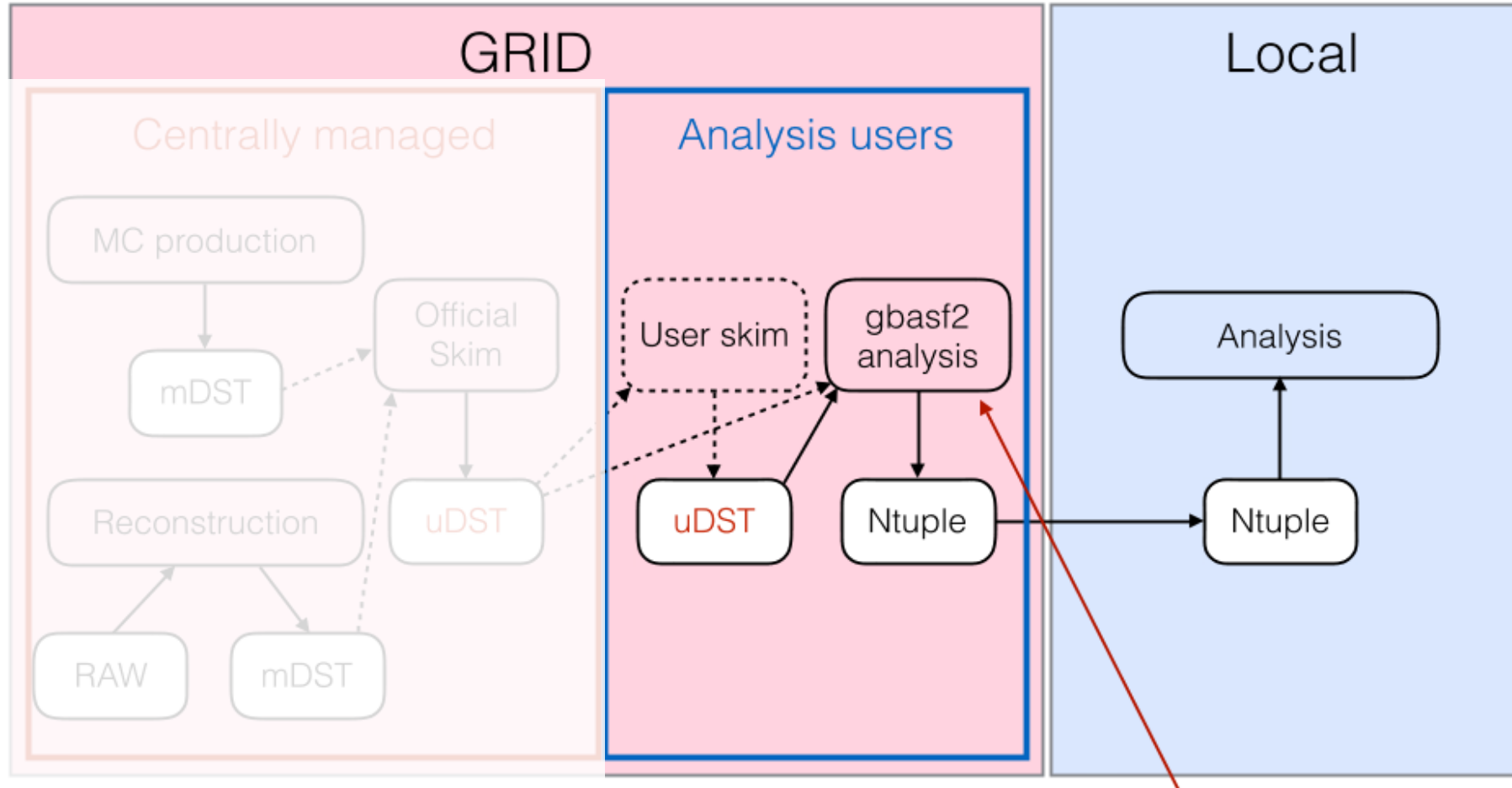
# One final thing about Belle vs. Belle II

# The Belle II analysis model



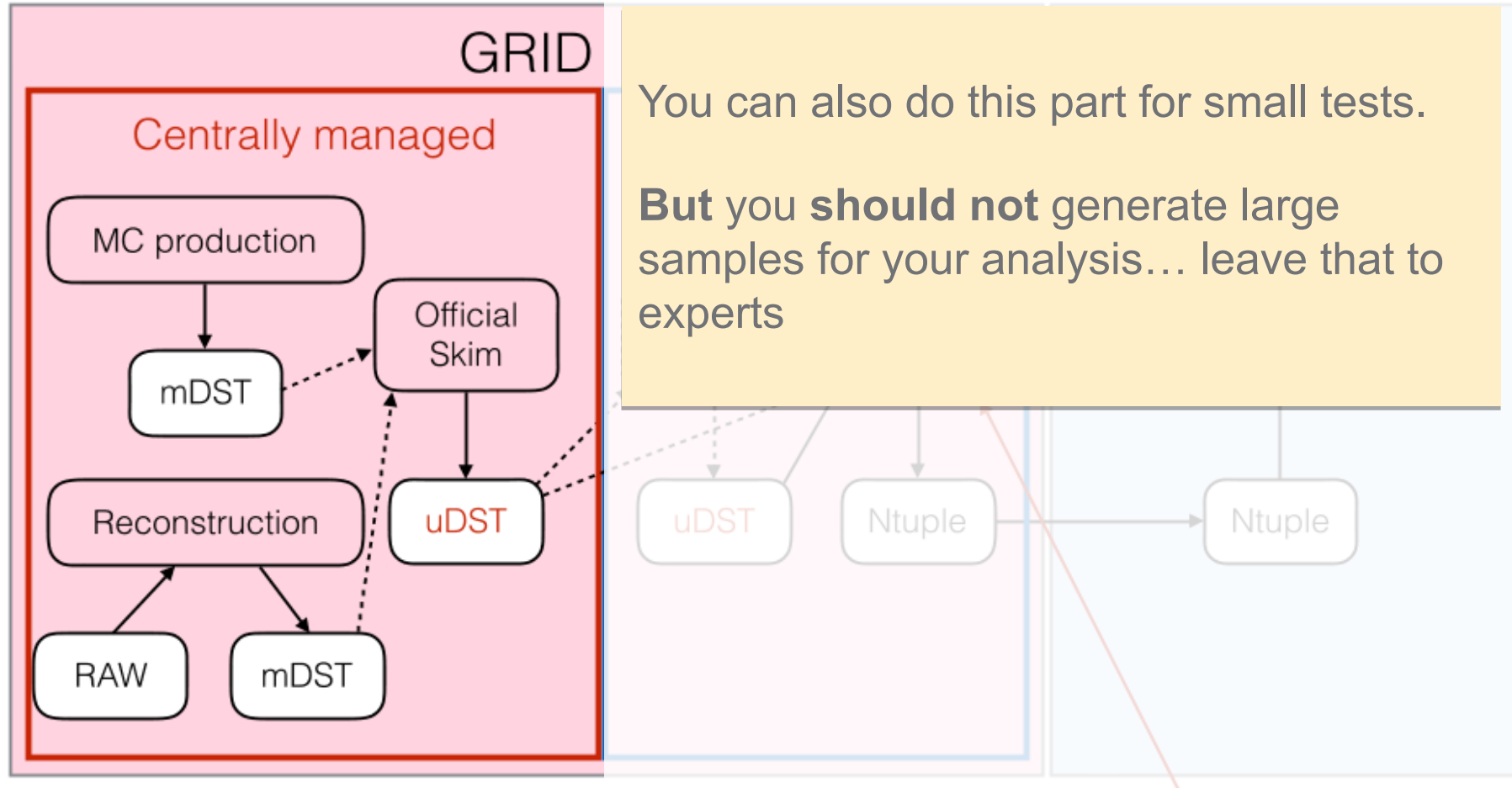
You will learn this tomorrow

# The Belle II analysis model



You will learn this tomorrow

# The Belle II analysis model



# Acknowledgements

- Much of the material for these slides (and ideas for how to present it) has been stolen from:

**Jake Bennett and Anže Zupanc**

## Versions of this tutorial material

- Version 1: Sam (February 2018).
- Version 2: Ilya and Sam (June 2018).
- Version 3: Sam (November 2018).
- Version 4: Umberto (February 2019)
- Version 5: Hannah and Sam (June 2019)
- Version 6: Yo and Ilya (January 2020)