

# Particles and decays in the Scikit-HEP project

Eduardo Rodrigues, Henry Schreiner  
University of Cincinnati , Princeton University

LHCb Week, 18<sup>th</sup> June 2019

---



PDG particle data  
and identification codes



Parse decay files, describe and  
convert particle decays between  
digital representations

# The grand picture – the Scikit-HEP project



- ❑ Create an ecosystem for particle physics data analysis in Python
- ❑ Initiative to improve the interoperability between HEP tools and the scientific ecosystem in Python
  - Expand the typical toolkitset for particle physicists
  - Set common APIs and definitions to ease “cross-talk”
- ❑ Initiative to build a community of developers and users
  - Community-driven and community-oriented project
- ❑ Effort to improve discoverability of relevant tools



**Collaboration**



**Reproducibility**



**Interoperability**



**Sustainability**

# Who uses (some of) Scikit-HEP ?

## Software projects



[Coffea](#) - a prototype [Analysis System](#) incorporating Scikit-HEP packages to provide a light-weight, scalable, portable, and user-friendly interface for columnar analysis of HEP data. Some of the sub-packages of Coffea may become Scikit-HEP packages as development continues.

The [zfit](#) project - it provides a model fitting library based on TensorFlow and optimised for simple and direct manipulation of probability density functions.

## Experiment collaborations



[BelleII](#) - the Belle II experiment at KEK, Japan.



[CMS](#) - the Compact Muon Solenoid experiment at CERN, Switzerland.

## Phenomenology projects



[flavio](#) - flavour physics phenomenology in the Standard Model and beyond.



***PDG particle data  
and identification codes***

identification codes  
particle data

# Particle package – motivation

---

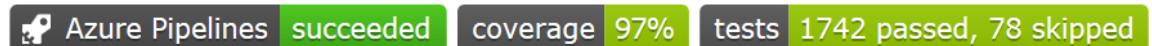
- The [Particle Data Group \(PDG\)](#) provides a downloadable table of particle masses, widths, charges and Monte Carlo particle ID numbers (PDG IDs)
  - Most recent file [here](#)
- It also provided an experimental file with extended information (spin, quark content, P and C parities, etc.) until 2008 *only*, see [here](#) (not widely known!)
- But **anyone** wanting to use these data, the only readily available, has to parse the file programmatically
- Why not make a Python package to deal with all these data, for everyone?
- The C++ HepPID and HepPDT libraries provide functions for processing particle ID codes in the standard particle (aka PDG) numbering scheme
- Different event generators have their separate set of particle IDs: Pythia, EvtGen, etc.
- Again, why not make a package providing all functionality/conversions, Python-ically, for everyone?



# Particle package – overview

---

- Pythonic interface to
  - PDG particle data table
  - and
  - particle identification codes
- With extra goodies
- 2 separate submodules
- Comprehensive documentation (docstrings)
- Continuous Integration (CI): extensive tests for excellent test coverage
  - In packages such as these, tests should target both the code itself but also the physics it deals with!
- We use [Azure DevOps](#)
  - Seamlessly test on Linux, macOS and Windows



# Particle package – PDG IDs module overview

## ❑ Process and query PDG IDs, and more – no look-up table needed

- Current version of package reflects the latest version of the HepPID & HepPDT utility functions defined in the [C++ HepPID and HepPDT versions 3.04.01](#)
- It contains more functionality than that available in the C++ code ... and minor fixes too

## ❑ Definition of a PDGID class, PDG ID literals, and set of standalone HepPID functions to query PDG IDs (is\_meson, has\_bottom, j\_spin, charge, etc.)

- All PDGID class functions are available standalone

## ❑ PDG ID queries also available on the command line

### ❑ PDGID class

- Wrapper class for PDG IDs
- Behaves like an int, with extra goodies
- Large spectrum of properties and methods, i.e. the functions defined in the HepPID and HepPDT C++ libraries, with a Pythonic interface, and yet more
- To print them all:

```
In [8]: print(PDGID(2212).info())
```

A	1
C	None
J	0.5

```
In [1]: from particle import PDGID  
  
pid = PDGID(211)  
pid
```

```
Out[1]: <PDGID: 211>
```

```
In [2]: pid.is_meson
```

```
Out[2]: True
```

```
In [3]: pid = PDGID(99999999)  
pid
```

```
Out[3]: <PDGID: 99999999 (is_valid==False)>
```

```
In [4]: from particle.pdgid import is_meson  
  
is_meson(211)
```

```
Out[4]: True
```

# Particle package – particle module overview

- Simple and natural API to deal with the PDG particle data table, with powerful look-up and search utilities
- Definition of a Particle class and particle name literals
  - Typical queries should be, and are, 1-liners

```
In [7]: from particle import Particle, SpinType  
  
Particle.findall(lambda p: p.pdgid.is_meson and p.pdgid.has_charm and p.spin_type==SpinType.PseudoScalar)  
  
Out[7]: [<Particle: name="D+", pdgid=411, mass=1869.65 ± 0.05 MeV>,  
          <Particle: name="D-", pdgid=-411, mass=1869.65 ± 0.05 MeV>,  
          <Particle: name="D0", pdgid=421, mass=1864.83 ± 0.05 MeV>,  
          <Particle: name="D~0", pdgid=-421, mass=1864.83 ± 0.05 MeV>,  
          <Particle: name="D(s)+", pdgid=431, mass=1968.34 ± 0.07 MeV>,  
          <Particle: name="D(s)-", pdgid=-431, mass=1968.34 ± 0.07 MeV>,  
          <Particle: name="eta(c)(1S)", pdgid=441, mass=2983.9 ± 0.5 MeV>,  
          <Particle: name="B(c)+", pdgid=541, mass=6274.9 ± 0.8 MeV>,  
          <Particle: name="B(c)-", pdgid=-541, mass=6274.9 ± 0.8 MeV>,  
          <Particle: name="eta(c)(2S)", pdgid=100441, mass=3637.6 ± 1.2 MeV>]
```

- Advanced usage: ability to specify or build a particle data table, conversion tools
- Particle / PDG ID searches available on the command line too

# Particle package – data files

---

## All data files stored under `particle/data/`

## PDG particle data files

- Original PDG data files, which are in a fixed-width format
- Code uses “digested forms” of the PDG files, stored as CSV, for optimised querying
- Latest PDG data used by default (2019 at present)
- Advanced usage: user can load older PDG table, load a “user table” with new particles, append to default table

ID	Mass	MassUpper	MassLower	Width	WidthUpper	WidthLower	I	G	P	C	Anti	Charge	Rank	Status	Name	Quarks
...																
441	2983.9	0.5	0.5	32	0.8	0.8	0	1	-1	1	0	0	0	0	eta(c)(1S)	cC
443	3096.9	0.006	0.006	0.0929	0.0028	0.0028	0	-1	-1	-1	0	0	0	0	J/psi(1S)	cC
445	3556.17	0.07	0.07	1.97	0.09	0.09	0	1	1	1	0	0	0	0	chi(c2)(1P)	cC
...																

## Other data files

- CSV file for mapping of PDG IDs to particle LaTeX names

# Particle package – particle look-up

## ❑ Particle class

- Standard look-up via `from_pdgid(...)`
- Various other `from_X(...)` methods exist

## ❑ - Large spectrum of properties and methods:

- Get particle properties
- Deal with underlying particle table
- Powerful search engine ...

```
In [1]: from particle import Particle
```

```
Particle.from_pdgid(-14122)
```

```
Out[1]:  $\bar{\Lambda}_c(2593)^-$ 
```

```
In [2]: Particle.from_string('pi-')
```

```
Out[2]:  $\pi^-$ 
```

```
In [3]: from IPython.core.display import display, HTML, Latex
```

```
print(Particle.from_pdgid(-10311).__repr__())
display(HTML('\t HTML name: {0}'.format(Particle.from_pdgid(-10311).html_name)))
display(Latex('\t LaTeX name: ${0}$.format(Particle.from_pdgid(-10311).latex_name)))
```

```
<Particle: name="K(0)*(1430)~0", pdgid=-10311, mass=1430 ± 50 MeV>
```

```
HTML name:  $\bar{K}_0^*(1430)^0$ 
```

```
LaTeX name:  $\bar{K}_0^*(1430)^0$ 
```

```
In [11]: from particle.particle import literals as lp
print(lp.K_0st_1430_0_bar.pdgid)
print(Particle.from_pdgid(-10311).programmatic_name)
<PDGID: -10311>
K_0st_1430_0_bar
```

## ❑ Particle literals

- Easily recognizable names for manipulations, e.g. in plots

# Particle package – powerful particle search

- ❑ `Particle.find(...)` – search a single match (exception raised if multiple particles match the search specifications)
- ❑ `Particle.findall(...)` – search a list of candidates

- ❑ Powerful search methods  
that can query any particle property!
- ❑ One-line queries

```
In [16]: print(Particle.find('J/psi').describe())
```

```
Name: J/psi(1S) ID: 443 Latex: $J/\psi(1S)$
Mass = 3096.900 ± 0.006 MeV
Lifetime = 7.09e-12 ± 2.2e-13 ns
Q (charge) = 0 J (total angular) = 1.0 P (space parity) = -
C (charge parity) = - I (isospin) = 0 G (G-parity) = -
SpinType: SpinType.Vector
Quarks: cC
Antiparticle name: J/psi(1S) (antiparticle status: Same)
```

- ❑ E.g., trivially find all pseudoscalar charm mesons:

```
In [7]: from particle import Particle, SpinType
Particle.findall(lambda p: p.pdgid.is_meson and p.pdgid.has_charm and p.spin_type==SpinType.PseudoScalar)
```

```
Out[7]: [<Particle: name="D+", pdgid=411, mass=1869.65 ± 0.05 MeV>,
<Particle: name="D-", pdgid=-411, mass=1869.65 ± 0.05 MeV>,
<Particle: name="D0", pdgid=421, mass=1864.83 ± 0.05 MeV>,
<Particle: name="D~0", pdgid=-421, mass=1864.83 ± 0.05 MeV>,
<Particle: name="D(s)+", pdgid=431, mass=1968.34 ± 0.07 MeV>,
<Particle: name="D(s)-", pdgid=-431, mass=1968.34 ± 0.07 MeV>,
<Particle: name="eta(c)(1S)", pdgid=441, mass=2983.9 ± 0.5 MeV>,
<Particle: name="B(c)+", pdgid=541, mass=6274.9 ± 0.8 MeV>,
<Particle: name="B(c)-", pdgid=-541, mass=6274.9 ± 0.8 MeV>,
<Particle: name="eta(c)(2S)", pdgid=100441, mass=3637.6 ± 1.2 MeV>]
```

# Particle package – future directions & developments

---

- **Addition of particle IDs and names relevant to other MC programs**

- (Yes, not consistent across programs!)
- Useful IDs such as those used in PYTHIA, Geant and EvtGen

- **Bring in other communities where Particle is / can be relevant**

- Ongoing discussions with astroparticle physics community
- Particle IDs used in EPOS, CORSIKA, DpmJet, QGSJet, Sybill, UrQMD, ...

- **Ongoing discussions with PDG group**

- Provide the right tool
- Can we provide more?
- Stay tuned ...



*Parse decay files, describe and  
convert particle decays  
between digital representations*

parse decay files, describe and  
convert particle decays between digital representations

parse decay files, describe and  
convert particle decays between digital representations

parse decay files, describe and  
convert particle decays between digital representations

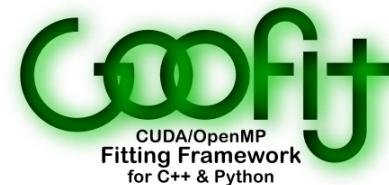
# DecayLanguage package – motivation and overview

## Motivation

- ❑ Ability to describe decay-tree-like structures
- ❑ Provide a translation of decay amplitude models from AmpGen to GooFit
  - Idea is to generalise this to other decay descriptions
- ❑ Any experiment uses event generators which, among many things, need to describe particle decay chains
- ❑ Programs such as EvtGen rely on so-called .dec decay files
- ❑ Many experiments need decay data files
- ❑ Why not make a Python package to deal with decay files, for everyone?



Library and set of applications for fitting and generating multi-body particle decays using the isobar model



## Overview

- ❑ Tools to parse decay files and programmatically manipulate them, query, display information
  - Descriptions and parsing built atop the Lark parser
- ❑ Tools to translate decay amplitude models from AmpGen to GooFit, and manipulate them



# DecayLanguage package – decay files

## “Master file” DECAY.DEC

- Gigantic file defining decay modes for all relevant particles, including decay model specifications
- LHCb example:
  - ~ 450 particle decays, thousands of decay modes, over 11k lines in total

```
Define dm 0.507e12
...
Alias      B0sig          B0
Alias      anti-B0sig     anti-B0
ChargeConj B0sig         anti-B0sig
...

Decay pi0
0.988228297  gamma   gamma
0.011738247  e+      e-      gamma
0.000033392  e+      e+      e-      e-
0.000000065  e+      e-
Enddecay
...
CDecay tau+
...
```

## User .dec files

- Needed to produce specific MC samples
- Typically contain a single decay chain (except if defining inclusive samples)

```
# Decay file for [B_c+ -> (B_s0 -> K+ K-) pi+]cc

Alias      B_c+sig        B_c+
Alias      B_c-sig        B_c-
ChargeConj B_c+sig       B_c-sig
Alias      MyB_s0          B_s0
Alias      Myanti-B_s0     anti-B_s0
ChargeConj MyB_s0         Myanti-B_s0

Decay B_c+sig
1.000    MyB_s0        pi+      PHOTOS PHSP;
Enddecay
CDecay B_c-sig
Decay MyB_s0
1.000    K+      K-      SSD_CP 20.e12 0.1 1.0 0.04 9.6 -0.8 8.4 -0.6;
Enddecay
CDecay Myanti-B_s0
```

# DecayLanguage package – decay file parsing and display

## ❑ Parsing should be simple

- Expert users can configure parser choice and settings, etc.

## ❑ Parsing should be (reasonably) fast

- Example of LHCb's master DECAY.DEC file:

Over 11k lines in total, ~ 450 particle decays, ~60 charge-conjugate decays created on-the-fly ('CDecay' statements), thousands of decay modes

```
p = DecFileParser(DIR / 'decaylanguage/data/DECAY_LHCB.DEC')
%timeit p.parse()
2.07 s ± 182 ms per loop (mean ± std. dev. of 7 runs, 1 loop each)

p.number_of_decays
506
```

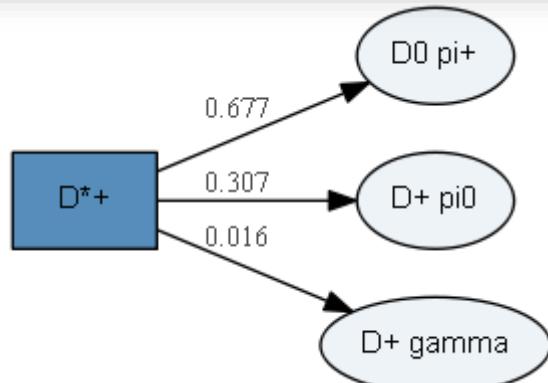
## ❑ After parsing, many queries are possible

## ❑ One can also visualise decay chains ... ☺

# Decay chain – simplest view with no sub-decays shown

```
from decaylanguage import DecFileParser  
  
p = DecFileParser('Dst.dec')  
p.parse()  
p  
  
<DecFileParser: decfile='Dst.dec', n_decays=5>
```

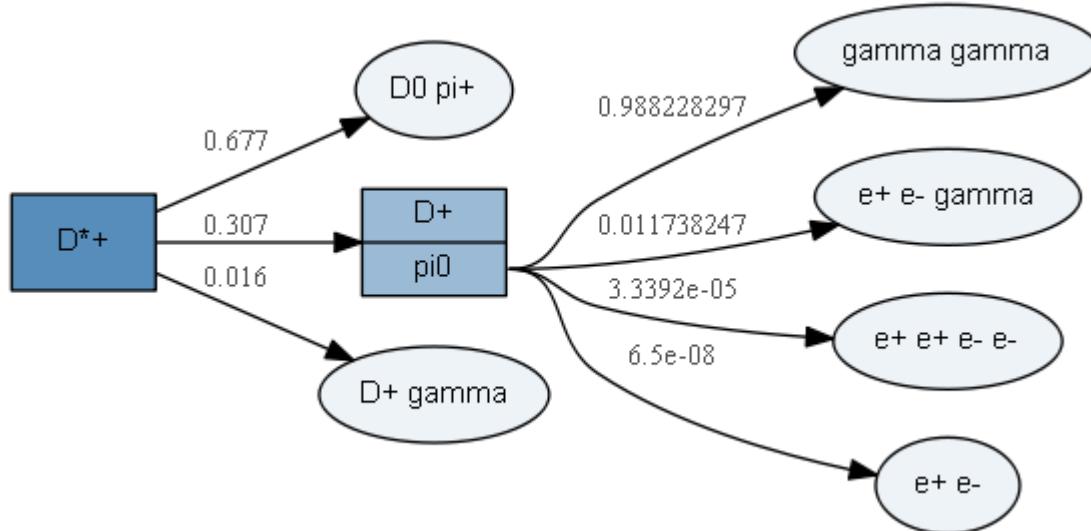
```
from decaylanguage.decay.viewer import DecayChainViewer  
  
chain = p.build_decay_chain('D*+', stable_particles=['D0', 'D+', 'pi0'])  
  
dcv = DecayChainViewer(chain)  
dcv
```



```
Decay D*+  
0.6770 D0 pi+  
0.3070 D+ pi0  
0.0160 D+ gamma  
Enddecay  
  
Decay D*-  
0.6770 anti-D0 pi-  
0.3070 D- pi0  
0.0160 D- gamma  
Enddecay  
  
Decay D0  
1.0 K- pi+  
Enddecay  
  
Decay D+  
1.0 K- pi+ pi+ pi0 PHSP;  
Enddecay  
  
Decay pi0  
0.988228297 gamma gamma PHSP;  
0.011738247 e+ e- gamma PI0_DALITZ;  
0.000033392 e+ e+ e- e- PHSP;  
0.000000065 e+ e- PHSP;  
Enddecay
```

(Considered by itself, this file is in fact incomplete, as there are no instructions on how to decay the anti- $D^0$  and the  $D^-$ . Good enough for illustration purposes, though.)

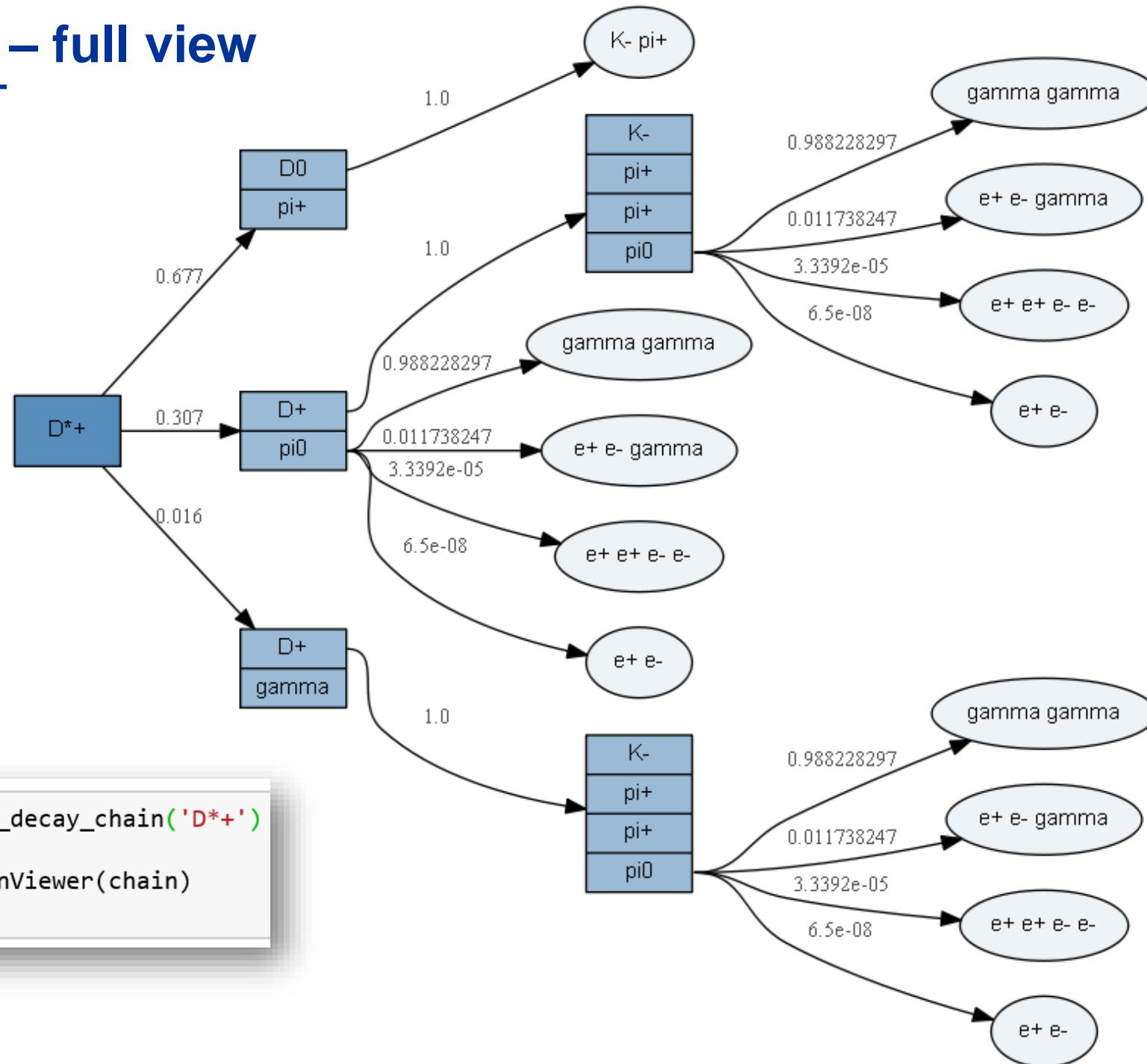
# Decay chain – $\pi^0$ decays displayed



```
chain = p.build_decay_chain('D*+', stable_particles=['D0', 'D+'])

dcv = DecayChainViewer(chain)
dcv
```

# Decay chain – full view



```
chain = p.build_decay_chain('D*+')  
  
dcv = DecayChainViewer(chain)  
dcv
```

# DecayLanguage package – conversion of decay models / representations

## □ Decay chains

- A universal modelling of decay chains would profit many use cases,  
e.g. description of components for amplitude analyses

## □ Present code understands AmpGen syntax and can generate code for the GooFit fitter

### □ Note:

makes use of the Particle package

```
[2]: lines, parameters, constants, states = AmplitudeChain.read_ampgen(text='''  
EventType D0 K- pi+ pi+ pi-  
  
D0[D]{K*(892)bar0{K-,pi+},rho(770)0{pi+,pi-}} 0 1 0.1 0 1 0.1  
  
K(1460)bar-_mass 0 1460 1  
K(1460)bar-_width 0 250 1  
  
a(1)(1260)+::Spline::Min 0.18412  
a(1)(1260)+::Spline::Max 1.86869  
a(1)(1260)+::Spline::N 34  
'''')  
  
[3]: lines[0].all_particles  
  
[3]: {<Particle: name="rho(770)0", pdgid=113, mass=775.3 ± 0.2 MeV>,  
<Particle: name="pi+", pdgid=211, mass=139.57061 ± 0.00024 MeV>,  
<Particle: name="pi-", pdgid=-211, mass=139.57061 ± 0.00024 MeV>,  
<Particle: name="K*(892)~0", pdgid=-313, mass=895.55 ± 0.20 MeV>,  
<Particle: name="K-", pdgid=-321, mass=493.677 ± 0.016 MeV>,  
<Particle: name="D0", pdgid=421, mass=1864.83 ± 0.05 MeV>}
```

# DecayLanguage package – future directions & developments

---

## *Decay files*

### Streamline and enhance the .dec parser

- Ex.: syntax such as

```
p.find_decay_chains(final_state=['K+', 'K-', 'pi+', 'pi-'], extra_particles=['pi0'])
```

could be a neat/trivial way to query the master DECAY.DEC and

“find all decay chains leading to either ‘K+ K- pi+ pi-’ or ‘K+ K- pi+ pi- pi0’”

### Provide a universal description and visualisation of decay trees (a lot done on this in the last week ...)

- We already have customers interested, e.g. visualisation of decays in [pyhepmc](#)

## *Decay models / representations*

### Implement more backend formats: GooFit in Python, etc.

### Longer term – implement decay logic inside model descriptions

- Provide a reference for other packages

# Interested ? Want to try it ?

---

## *Particle*

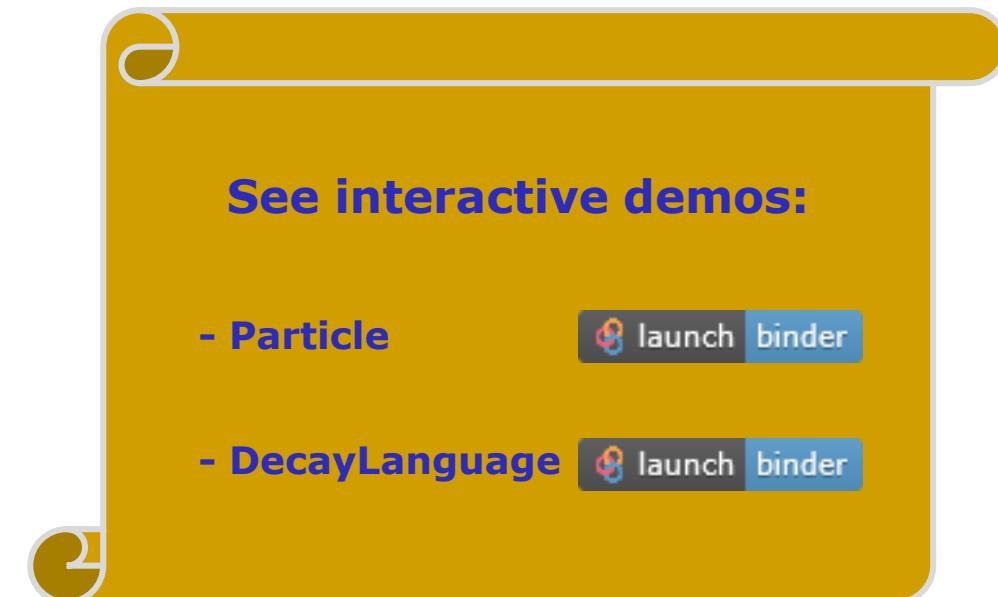
- GitHub: <https://github.com/scikit-hep/particle/>
- Releases: [PyPI](#) 
- Kindly recognise software work – cite us: DOI [10.5281/zenodo.2552429](https://doi.org/10.5281/zenodo.2552429)

## *DecayLanguage*

- GitHub: <https://github.com/scikit-hep/decaylanguage>
- Releases: [PyPI](#) 

## *Scikit-HEP project*

- GitHub: <https://github.com/scikit-hep/>
- Website: <http://scikit-hep.org/>
- Get in touch: <http://scikit-hep.org/get-in-touch.html>



***Back-up***

**Back-up**

# Particle package – PDG identification code literals

- ❑ **Literals:** handy way to manipulate things with human-readable names

- ❑ **PDGID literals**

- Provide (PDGID class) aliases  
for the most common particles,  
with easily recognisable names

- ❑ All is consistent. Ex.:

```
In [5]: from particle.pdgid import literals as lid
lid.pi_plus
Out[5]: <PDGID: 211>

In [6]: from particle.pdgid.literals import Lambda_b_0
Lambda_b_0
Out[6]: <PDGID: 5122>

In [7]: Lambda_b_0.has_bottom
Out[7]: True
```

```
In [8]: Particle.from_pdgid(-10311).pdgid == literals.K_0st_1430_0_bar
Out[8]: True
```

# DecayLanguage package – Lark parser grammar for decay files

- Decay file parser grammar:  
`decfile.lark !`

- This file is enough to parse and understand decay files

```
start : _NEWLINE? (line _NEWLINE)+ ("End" _NEWLINE)?
?line : define | pythia_def | alias | chargeconj | commands | decay | cdecay | setlspw

pythia_def : "PythiaBothParam" LABEL ":" LABEL "=" (LABEL | SIGNED_NUMBER)
setlspw : "SetLineshapePW" label label label value
cdecay : "CDecay" label
define : "Define" label value

alias : "Alias" label label
chargeconj : "ChargeConj" label label

?commands : global_photos
global_photos : boolean_photos
boolean_photos : "yesPhotos" -> yes
| "noPhotos" -> no

decay : "Decay" particle _NEWLINE decayline+ "Enddecay"
decayline : value particle* photos? model _NEWLINE // There is always a ; here
value : SIGNED_NUMBER
photos : "PHOTOS"

label : LABEL
particle : LABEL // Add full particle parsing here

model : MODEL_NAME model_options?
model_options : (value | LABEL)+

%import common.WS_INLINE
%import common.SIGNED_NUMBER

// New lines filter our comments too, and multiple new lines
_NEWLINE: ( /\r?\n[\t ]*/ | COMMENT )+

MODEL_NAME.2 : "BaryonPCR" | "BTO3PI_CP" | "BTOSLLALI" | "BTOSLLBALL" | "BTOXSGAMMA" | "BTOXSL" | ...
LABEL : /[a-zA-Z0-9\/-+*()']+
COMMENT : [;#][^\n]*/

// We should ignore comments
-ignore COMMENT

// Disregard spaces in text
-ignore WS_INLINE
```