

Monte Python

the way back from cosmology to Fundamental Physics

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and

UC Berkeley



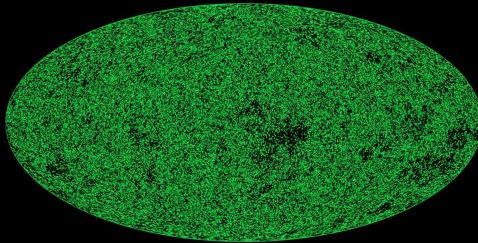
NORDITA



IFT School on Cosmology Tools

March 2017

Related Tools

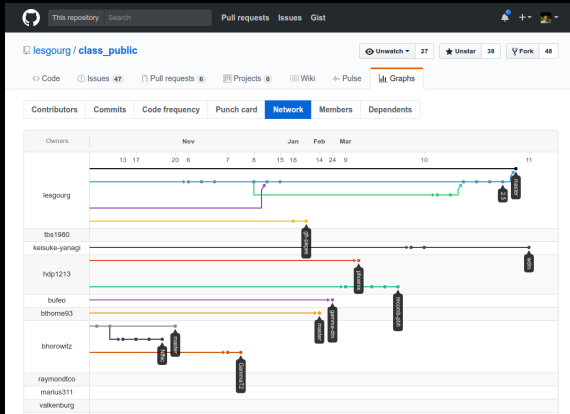


Version control and Python interfacing

PLANCK

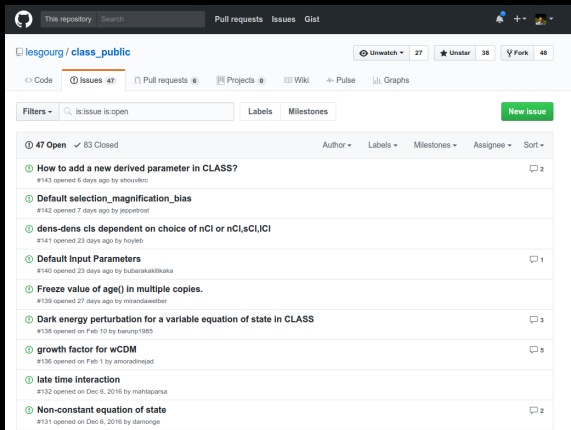
Version control: GIT + github

- GIT: access all previous versions, restore, compare, branch...
- Github: website + interface
 - Network: users, branches intermediate versions



Version control: GIT + github

- GIT: access all previous versions, restore, compare, branch...
- Github: website + interface
 - Network: users, branches intermediate versions
 - Issues: troubleshooting, forum/improvements



The screenshot shows the GitHub interface for the repository 'lesgourg / class_public'. The 'Issues' tab is selected, showing 47 open issues. The issues are listed in a table with columns for status, title, number, age, author, and comments. The first issue is 'How to add a new derived parameter in CLASS?' (#143) opened 6 days ago by shouvillec. Other issues include 'Default selection_magnification_bias' (#142), 'dens-dens cls dependent on choice of nCl or nCl,sCl,ICl' (#141), 'Default Input Parameters' (#140), 'Freeze value of age() in multiple copies.' (#139), 'Dark energy perturbation for a variable equation of state in CLASS' (#138), 'growth factor for wCDM' (#136), 'late time interaction' (#132), and 'Non-constant equation of state' (#131).

Status	Title	Number	Age	Author	Comments
Open	How to add a new derived parameter in CLASS?	#143	opened 6 days ago	shouvillec	2
Open	Default selection_magnification_bias	#142	opened 7 days ago	jospetrost	
Open	dens-dens cls dependent on choice of nCl or nCl,sCl,ICl	#141	opened 23 days ago	hoyleb	
Open	Default Input Parameters	#140	opened 23 days ago	bubarakakikaka	1
Open	Freeze value of age() in multiple copies.	#139	opened 27 days ago	mirandawebster	
Open	Dark energy perturbation for a variable equation of state in CLASS	#138	opened on Feb 10 by barump1985		3
Open	growth factor for wCDM	#136	opened on Feb 1 by amoradnejad		5
Open	late time interaction	#132	opened on Dec 6, 2016 by mahtaparsa		
Open	Non-constant equation of state	#131	opened on Dec 6, 2016 by damonge		2

Python interfacing with classy

- classy → use CLASS as a Python module
 - Required for MCMC (tomorrow!)
 - Useful for plotting

```
from classy import Class
import numpy as np
import matplotlib.pyplot as plt
cosmo = Class ()
cosmo.set ({'output': 'tCl, pCl, lCl', 'lensing': 'yes'})
cosmo.compute ()
l = np.array ( range (2 ,2501) )
factor = l*(l +1) /(2*np.pi )
lensed_cl = cosmo.lensed_cl (2500)
#then just plot lensed_cl...
```

Python interfacing with classy

- classy → use CLASS as a Python module
 - Required for MCMC (tomorrow!)
 - Useful for plotting
- IPython → Interactive Python frontend
 - TAB auto-completion

```
miguel@Goedel:~$ ipython
Python 2.7.6 (default, Oct 26 2016, 20:30:19)
Type "copyright", "credits" or "license()" for more information.

IPython 5.1.0 -- An enhanced Interactive Python.
?                -> Introduction and overview of IPython's features.
%quickref        -> Quick reference.
help             -> Python's own help system.
object?         -> Details about 'object', use 'object??' for extra details.
```

```
In [1]: from classy import Class
```

```
In [2]: c = Class()
```

```
In [3]: c.
c.age              c.h              c.Omega_nu
c.angular_distance c.Hubble         c.pars
c.baryon_temperature c.ionization_fraction c.pk
c.compute          c.lensed_cl      c.raw_cl
c.density_cl       c.luminosity_distance c.rs_drag
c.empty            c.n_s            c.set
c.get_background   c.Neff           c.set_default
c.get_current_derived_parameters c.nonlinear_method c.sigma8
c.get_perturbations c.nonlinear_scale c.state
c.get_pk           c.Omega0_m       c.struct_cleanup
c.get_prinordial   c.Omega_b        c.T_cmb
c.get_thermodynamics c.omega_b        c.z_of_r
c.get_transfer      c.Omega_m
```

Python interfacing with classy

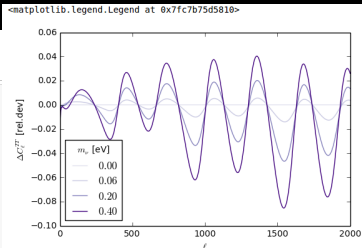
- classy → use CLASS as a Python module
 - Required for MCMC (tomorrow!)
 - Useful for plotting
- IPython → Interactive Python frontend
 - TAB auto-completion
- Jupyter → Notebook interface (Julia+Python+R)

Jupyter example

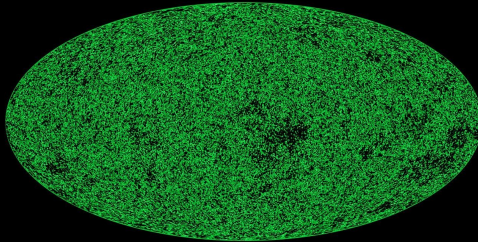
You can write \mathcal{L}_{eff} and awesome equations like $m_\nu \gtrsim 60$ meV. And plot the effect of m_ν in few lines

```
cm = plt.get_cmap('Purples')
c = Class()
cl = {} #dictionary for output
for m in [0.0, 0.06, 0.2, 0.4]:
    c.set({'N_ncdm':1, 'N_ur':2.033, 'm_ncdm':m, 'output':'tcl'})
    c.compute()
    cl[m] = c.raw_cl(2000)
    plt.plot(cl[m]['ell'][2:], cl[m]['tt'][2:]/cl[0]['tt'][2:]-1.,
             color=cm((m+0.1)/0.5), label=r'$%.2f$'%(m))
    c.empty()

plt.xlabel(r'$\ell$')
plt.ylabel(r'$\Delta C_{\ell}^{\text{TT}}$ [rel.dev]')
plt.legend(loc='lower left', fontsize = 12, title= r'$m_{\nu}$ [eV]')
```



Monte Python



from cosmology back to fundamental physics

PLANCK

DISCLAIMER: Short time!

$\lesssim 3h$ course \Rightarrow overview and basic usage

Learn further:

- MontePython slides by Sebastien Clesse ($\ll 1h?$)

https://lesgourg.github.io/class-tour/16.06.02_Lisbon_intro.pdf

- MontePython course ($\sim 5h$)

<https://lesgourg.github.io/class-tour-Tokyo.html>

- Links to extra resources in exercise sheet

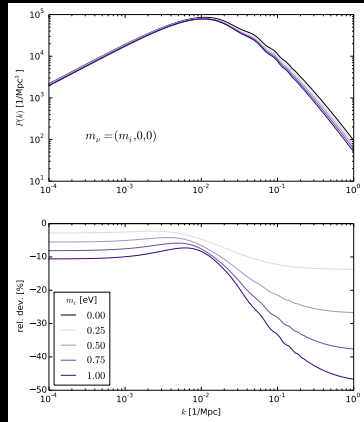
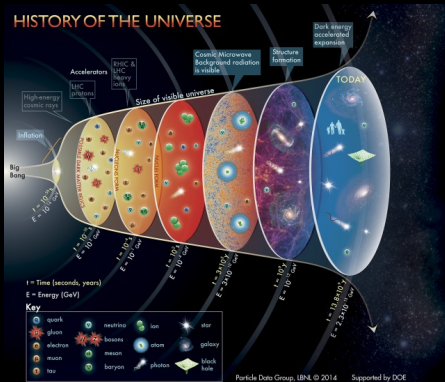
IMPORTANT DISCLAIMER:

↓ I'm mainly a user with little experience developing!

↑ Help from experts:

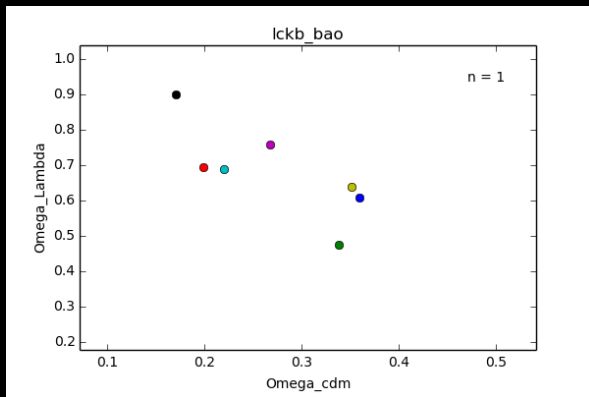
Thejs Brinckmann, Carlos Garcia, Deanna Hooper & Vivian Poulin

Fundamental physics and cosmology



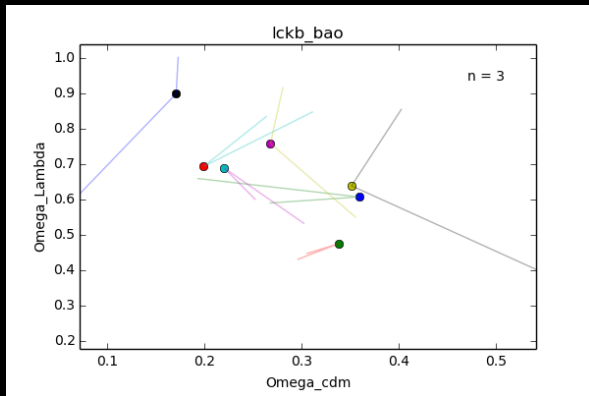
Initial conditions, Dark Matter, Neutrinos, Dark Energy, Gravity...

Scan space of parameters



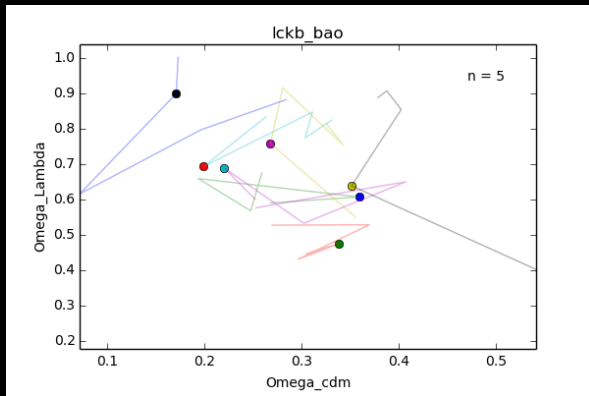
(recall lecture by Will Handley)

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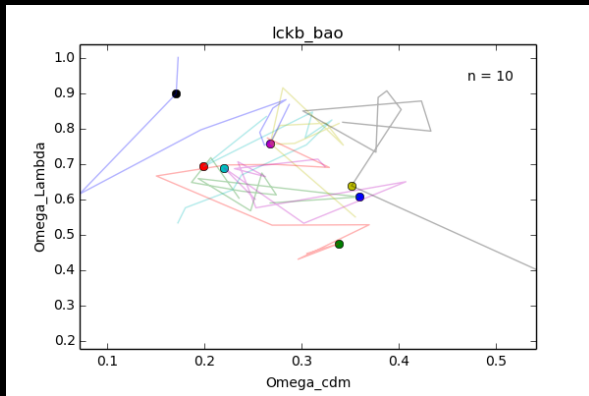
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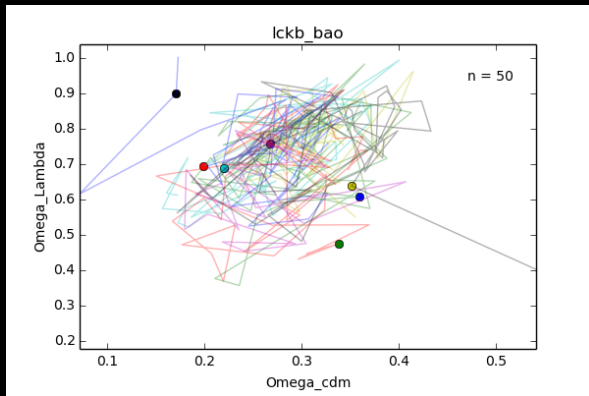
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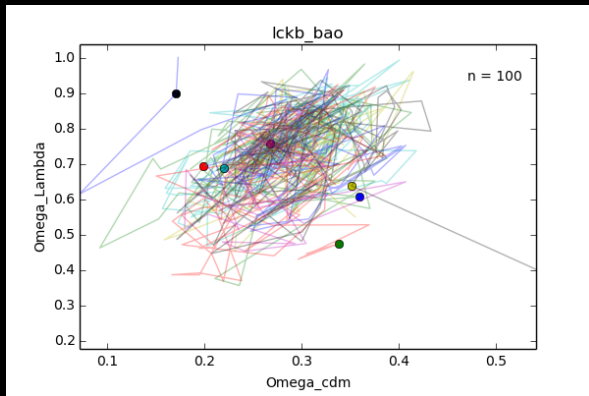
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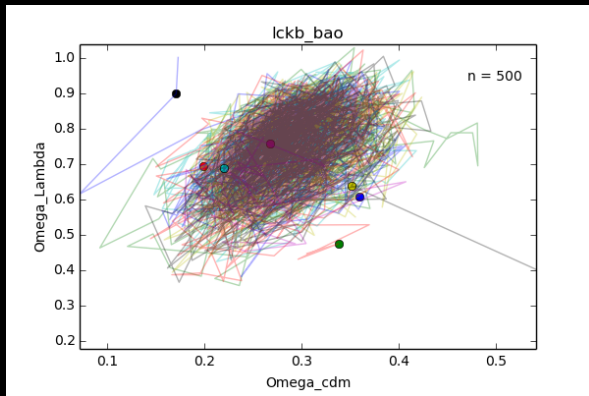
(recall lecture by Will Handley)

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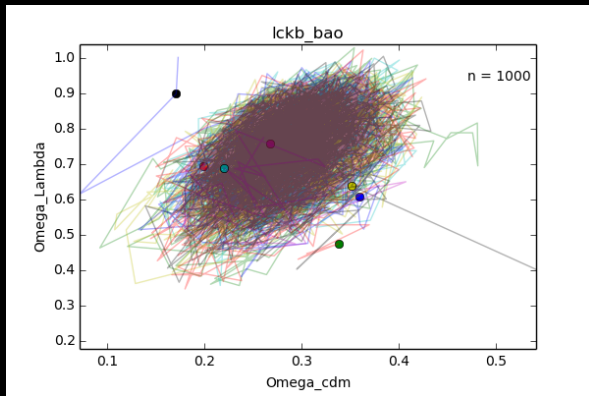
(recall lecture by Will Handley)

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MontePython

An Markov Chain Monte Carlo engine for parameter extraction:

Features

- Written in Python
 - Python is practically magic!
 - imports routines from `numpy` and `scipy`
 - useful outside academia, standard for Big Data

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- Modular, easy to add
 - likelihoods for new experiments
 - features for sampling, plotting...

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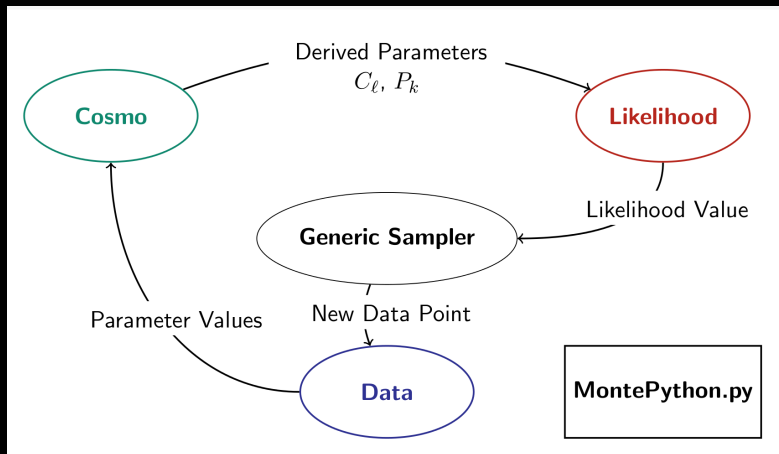
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- Modular, easy to add
 - likelihoods for new experiments
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- Easy to use, intensively documented
- Parallelization is optional
 - simpler to install
 - runs in old/separate computers, short queue

Modular Structure



(from B. Audren's Monte Python slides)

Defining a run (model.param)

- List of experiments

```
data.experiments=['Planck_highl','Planck_lowl','Planck_lensing']
```

Collected in montepython/likelihoods:

```
miguel@Goedel:~/code/montepython_zuma/montepython/likelihoods$ ls
acbar          clik_wmap_full      __init__.py         quad
bao            clik_wmap_lowl      __init__.pyc        sdss_lrgDR4
bao_boss       cosmic_clocks_BC03  JLA                  sn
bao_boss_aniso cosmic_clocks_BC03_all JLA_simple           spt
bao_boss_aniso_gauss_approx cosmic_clocks_MaStro lowlike              spt_2500
bao_known_rs   da_rec              Planck_actspt        test_gaussian
bicep          euclid_lensing      Planck_highl         test_nuisance1
bicep2         euclid_pk           Planck_highl_lite    test_nuisance2
boomerang      fake_desi           Planck_highl_TTTEEE  timedelay
cbl            fake_planck_bluebook Planck_lensing        WiggleZ
CFHTLens       gunn_peterson       Planck_lowl          WiggleZ_bao
CFHTLens_correlation hst                 Planck_SZ            wmap
clik_fake_planck igm_temperature     polarbear            wmap_9yr
```

Defining a run (model.param)

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- Cosmological parameters

```
                # [mean, (bounds) , SIGMA,scale, type ]  
data.parameters['n_s']=[0.96, None,None, 0.008, 1 , 'cosmo']
```

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Fixed values:

- set $SIGMA = 0$
- `data.cosmo_arguments['N_ncdm'] = 1`

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- Derived and Nuisance parameters

```
data.parameters['sigma8'] = [0, None, None, 0, 1, 'derived']  
data.parameters['A_cib_217'] = [61,0,200,7,1,'nuisance']
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```

- MCMC parameters: data.N=10, data.write_step=5 ...

Running Monte Python (run)

Single chain:

```
python montepython/MontePython.py run \
  -p model.param \
  -o output_directory (...)
```

Parallel run (4 chains):

```
mpirun -nc 4 python (...)
```

Some options:

```
#!/bin/sh
#
# Appending to an existing folder: using the log paran instead of
# input/lambda_bao_b.paran
Running Monte Python v2.2.2

with CLASS v2.4.5

Testing likelihoods for:
-> bao_boss, bao_boss_aniso

#!/bin/sh
#
# excluding isotropic CMASS measurement
Creating chains/lcdm_bao_b/2017-03-12_100__2.txt

Creating error file chains/lcdm_bao_b/error_log.txt

Deduced starting covariance matrix:
['omega_b', 'Omega_cdm', 'Omega_k']
[[ 0. 0. 0.]
 [ 0. 0. 0.]
 [ 0. 0. 0.]]

# -logLik 1e+02omega_b Omega_cdm Omega_k Omega_Lambda
3 11.9558 2.221100e+00 3.132032e-01 -1.692095e-02 6.553103e-01
1 12.7946 2.262336e+00 3.271086e-01 2.742679e-02 5.961614e-01
1 12.8483 2.256136e+00 2.494162e-01 1.271974e-01 5.742162e-01
3 11.6619 2.286885e+00 2.388487e-01 1.014554e-01 6.478604e-01
1 11.4992 2.275734e+00 2.188346e-01 5.439644e-02 6.851744e-01
1 11.1477 2.309970e+00 2.013242e-01 3.477642e-02 7.135600e-01
3 7.83276 2.297757e+00 2.751741e-01 9.733703e-02 5.774153e-01
14 6.96653 2.302347e+00 2.748664e-01 5.921469e-02 6.157454e-01
4 6.46121 2.372939e+00 2.523288e-01 2.276382e-02 6.731983e-01
13 6.9573 2.334726e+00 2.363992e-01 -3.071722e-02 7.434402e-01
4 8.883 2.350200e+00 2.572473e-01 -3.950319e-02 7.311015e-01
14 7.19446 2.380873e+00 2.210523e-01 -2.615132e-02 7.532173e-01
2 6.95177 2.359364e+00 2.629244e-01 7.590971e-02 6.097521e-01
6 7.18686 2.339764e+00 2.595452e-01 8.159985e-02 6.078676e-01
1 6.44771 2.353995e+00 2.392110e-01 3.775342e-02 6.717389e-01
4 8.22755 2.326585e+00 2.075475e-01 -2.491768e-03 7.442001e-01
1 7.89432 2.390601e+00 2.076928e-01 -2.671562e-02 7.669296e-01
6 7.21255 2.388188e+00 2.145373e-01 7.295908e-03 7.261260e-01
2 7.07413 2.397083e+00 2.685047e-01 6.233775e-02 6.169233e-01
1 7.37498 2.406336e+00 2.207678e-01 4.592795e-02 6.808096e-01
2 10.6892 2.439898e+00 1.946383e-01 5.075931e-02 7.014388e-01
3 11.7196 2.476394e+00 2.298767e-01 1.535277e-01 5.626362e-01
1 12.3995 2.441832e+00 3.239238e-01 2.005879e-01 4.222807e-01
3 10.2062 2.420720e+00 2.880782e-01 1.751385e-01 4.840349e-01
1 7.3809 2.439488e+00 2.699004e-01 7.482511e-02 6.021178e-01
3 7.25137 2.437977e+00 2.578438e-01 2.764830e-02 6.613929e-01
1 6.83719 2.459588e+00 2.214344e-01 4.100809e-02 6.839635e-01

# 100 steps done, acceptance rate: 0.26
```

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Single chain:

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python montepython/MontePython.py run \
  -p model.param \
  -o output_directory (...)
```

Parallel run (4 chains):

```
mpirun -nc 4 python (...)
```

Some options:

- `-N` → # points
- `-C` → covariance matrix
- `-r` → restart from last point of chain
- `--update` → update sampling + covariance

```
#!/usr/bin/env python
#
# Appending to an existing folder: using the log paran instead of
# input/lambda_bao_b.paran
Running Monte Python v2.2.2

with CLASS v2.4.5

Testing likelihoods for:
-> bao_boss, bao_boss_aniso

#!/usr/bin/env python
#
# excluding isotropic CMASS measurement
Creating chains/icdmk_bao_b/2017-03-12_100_2.txt
Creating error file chains/icdmk_bao_b/error_log.txt

Deduced starting covariance matrix:
['omega_b', 'Omega_cdm', 'Omega_k']
[[ 0.  0.  0.]
 [ 0.  0.  0.]
 [ 0.  0.  0.]]

# -LogLikl      1e+02omega_b      Omega_cdm      Omega_k      Omega_Lambda
3  11.9558      2.221100e+00      3.132032e-01      -1.692095e-02      6.553103e-01
1  12.7946      2.262336e+00      3.271086e-01      2.742679e-02      5.961614e-01
1  12.8483      2.255136e+00      2.494162e-01      1.271974e-01      5.742182e-01
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13  6.9573      2.334726e+00      2.363992e-01      -3.071722e-02      7.434402e-01
4  8.883      2.350200e+00      2.572473e-01      -3.950319e-02      7.311015e-01
14  7.19446      2.380873e+00      2.210523e-01      -2.615132e-02      7.532173e-01
2  6.95177      2.359364e+00      2.629244e-01      7.590971e-02      6.097521e-01
6  7.18686      2.339764e+00      2.595452e-01      8.159985e-02      6.078676e-01
1  6.44771      2.353985e+00      2.392110e-01      3.775342e-02      6.717309e-01
4  8.22755      2.326585e+00      2.075475e-01      -2.491768e-03      7.442001e-01
1  7.89432      2.390601e+00      2.076928e-01      -2.671562e-02      7.669296e-01
6  7.21255      2.388188e+00      2.145373e-01      7.295908e-03      7.261260e-01
2  7.07413      2.397083e+00      2.685047e-01      6.233775e-02      6.169233e-01
1  7.37498      2.406336e+00      2.207678e-01      4.592705e-02      6.808096e-01
2  10.6892      2.439808e+00      1.940383e-01      5.075931e-02      7.014388e-01
3  11.7196      2.476394e+00      2.298767e-01      1.535277e-01      5.626362e-01
1  12.3995      2.441832e+00      2.239238e-01      2.005879e-01      4.222807e-01
3  10.2062      2.420720e+00      2.880782e-01      1.751385e-01      4.840349e-01
1  7.3809      2.439488e+00      2.699004e-01      7.482511e-02      6.021178e-01
3  7.25137      2.437977e+00      2.578438e-01      2.764830e-02      6.613929e-01
1  6.83719      2.459508e+00      2.214344e-01      4.100809e-02      6.839635e-01

# 100 steps done, acceptance rate: 0.26
```

All options explained

```
python montepython/MontePython.py info --help
```

Analyzing results (info)

Single model/experiment:

```
python montepython/MontePython.py info \
    output_directory (...)
```

Comparing several runs:

```
python montepython/MontePython.py info \
    output_1 output_2 output_3 (...)
```

Configuring the output/analysis

```
miguel@Goedel:~/code/montepython_zuma/chains_itp$ python ../montepython/MontePython.py info lckb_bao/
Running Monte Python v2.2.2

--> Finding global maximum of likelihood
--> Removing burn-in
--> Scanning file lckb_bao/2017-03-12_100000_3.txt : Removed 16
6 non-markovian points, 0 points of burn-in, keep 10397 steps
    2017-03-12_10_1.txt : Removed 0
non-markovian points, 2 points of burn-in, keep 1 steps
    2017-03-12_100000_4.txt : Removed 57
non-markovian points, 0 points of burn-in, keep 10405 steps
    2017-03-12_100000_8.txt : Removed 0
non-markovian points, 2 points of burn-in, keep 6783 steps
    2017-03-12_100000_7.txt : Removed 0
non-markovian points, 4 points of burn-in, keep 13666 steps
    2017-03-12_100000_1.txt : Removed 92
non-markovian points, 0 points of burn-in, keep 7975 steps
    2017-03-12_100000_6.txt : Removed 0
non-markovian points, 4 points of burn-in, keep 20 steps
    2017-03-12_100000_5.txt : Removed 15
5 non-markovian points, 0 points of burn-in, keep 11158 steps
    2017-03-12_10_2.txt : Removed 0
non-markovian points, 2 points of burn-in, keep 1 steps
--> Computing mean values
--> Computing variance
--> Computing convergence criterium (Gelman-Rubin)
-> R-1 is 0.002068 for Omega_b
    0.001200 for Omega_cdm
    0.002191 for Omega_k
    0.002137 for Omega_Lambda
-----
-> Computing histograms for Omega_b
-> Computing histograms for Omega_cdm
-> Computing histograms for Omega_k
-> Computing histograms for Omega_Lambda
-----
--> Saving figures to .pdf files
--> Writing .info and .tex files
```


Analyzing results (info)

Single model/experiment:

```
python montepython/MontePython.py info \
    output_directory (...)
```

Comparing several runs:

```
python montepython/MontePython.py info \
    output_1 output_2 output_3 (...)
```

Configuring the output/analysis

- `--extra` → file with plot options
- `--bins` → # bins for posterior
- `--all` → plot every subplot separately
- `--no-mean` → only marginalized in 1D

```
miguel@Goedel:~/code/montepython_zuma/chains_itp$ python ../montepython/MontePython.py info lckb_bao/
Running Monte Python v2.2.2

--> Finding global maximum of likelihood
--> Removing burn-in
--> Scanning file lckb_bao/2017-03-12_100000_3.txt : Removed 16
6 non-markovian points, 0 points of burn-in, keep 10397 steps
2017-03-12_10_1.txt : Removed 0
non-markovian points, 2 points of burn-in, keep 1 steps
2017-03-12_100000_4.txt : Removed 57
non-markovian points, 0 points of burn-in, keep 10405 steps
2017-03-12_100000_8.txt : Removed 0
non-markovian points, 2 points of burn-in, keep 6783 steps
2017-03-12_100000_7.txt : Removed 0
non-markovian points, 4 points of burn-in, keep 13666 steps
2017-03-12_100000_1.txt : Removed 92
non-markovian points, 0 points of burn-in, keep 7975 steps
2017-03-12_100000_6.txt : Removed 0
non-markovian points, 4 points of burn-in, keep 20 steps
2017-03-12_100000_5.txt : Removed 15
5 non-markovian points, 0 points of burn-in, keep 11158 steps
2017-03-12_10_2.txt : Removed 0
non-markovian points, 2 points of burn-in, keep 1 steps
--> Computing mean values
--> Computing variance
--> Computing convergence criterion (Gelman-Rubin)
-> R-1 is 0.002068 for Omega_b
0.001280 for Omega_cdm
0.002191 for Omega_k
0.002137 for Omega_Lambda

-----
-> Computing histograms for Omega_b
-> Computing histograms for Omega_cdm
-> Computing histograms for Omega_k
-> Computing histograms for Omega_Lambda
-----
--> Saving figures to .pdf files
--> Writing .info and .tex files
```

All options explained

```
python montepython/MontePython.py info --help
```

A very minimal run

Write `lckb.param`:

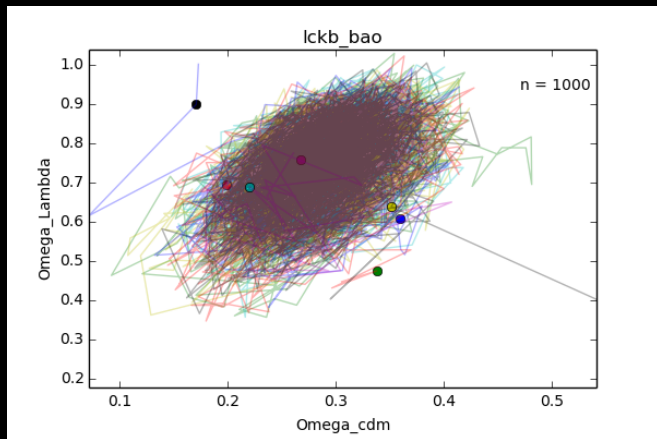
```
data.experiments=['bao_boss','bao_boss_aniso']
#Cosmo parameteress          [mean, min, max, sigma, scale, type]
data.parameters['Omega_b'] = [0.045,0.01, None,0.01,1,'cosmo']
data.parameters['Omega_cdm'] = [0.3, 0, None, 0.1, 1, 'cosmo']
data.parameters['Omega_k'] = [0.0, -0.5, 0.5, 0.1, 1, 'cosmo']
#Fixed parameters (sigma = 0)
data.parameters['H0'] = [67.8, None, None, 0, 1, 'cosmo']
data.cosmo_arguments['YHe'] = 0.24
#derived parameters
data.parameters['Omega_Lambda'] = [1,None,None,0,1,'derived']
#mcmc parameters
data.N=10
data.write_step=5
```

Run ~ 7 chains with

```
python montepython/MontePython.py run -o chains/lckb_bao \
-p lckb_param --update 300 -N 100000
```

A very minimal run

The 7 chains explore the parameter space



Chains named `yyyy-mm-dd_N_n.txt` (`date_points_id`)

A very minimal run

Analyze:

```
python montepython/MontePython.py info chains/lckb_bao
```

- lckb_bao.tex → table with MCMC results

Param	best-fit	mean $\pm\sigma$	95% lower	95% upper
Ω_b	0.03595	$0.03977^{+0.0095}_{-0.015}$	0.01662	0.06547
Ω_{cdm}	0.2931	$0.2872^{+0.049}_{-0.048}$	0.1892	0.3847
Ω_k	-0.1183	$-0.08087^{+0.11}_{-0.14}$	-0.3182	0.1755
Ω_Λ	0.7891	$0.7538^{+0.12}_{-0.091}$	0.542	0.9564

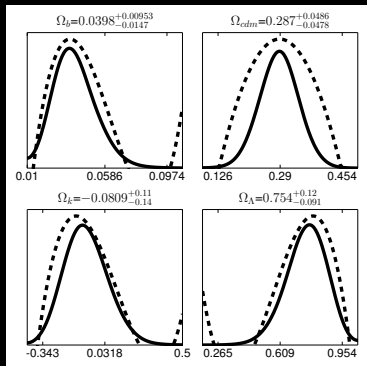
$-\ln \mathcal{L}_{\min} = 5.57269$, minimum $\chi^2 = 11.15$

- `lckb_bao.covmat` → covariance matrix
- `lckb_bao.bestfit` → best fit values

→ arguments for another run

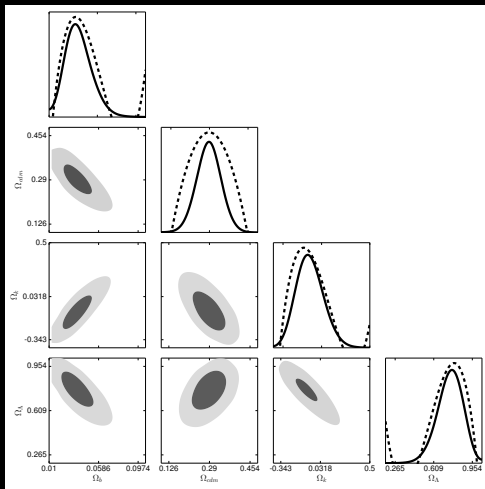
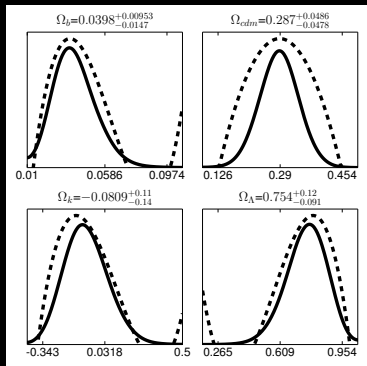
A very minimal run

In `lckb_bao/plots`:



A very minimal run

In `lckb_bao/plots`:

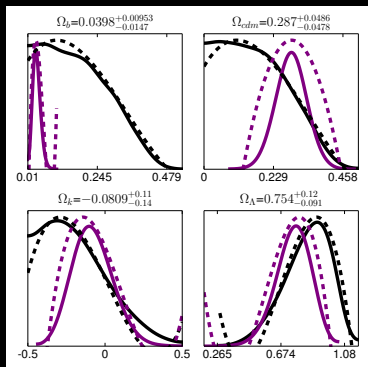


Comparing several runs

Run chains lckb_sne with `data.experiments=['sne']`

Analyze: `python ... info chains/lckb_sne chains/lckb_bao`

In lckb_sne/plots:

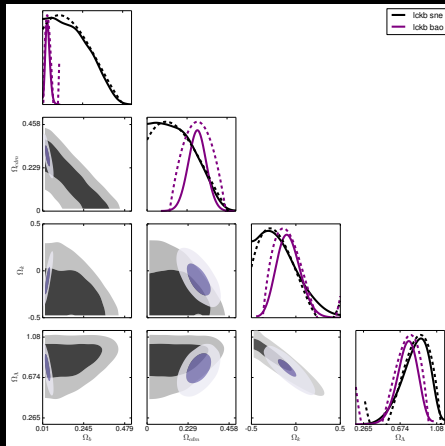
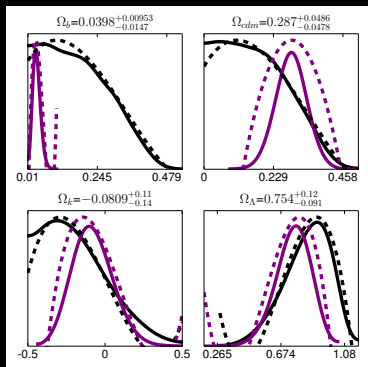


Comparing several runs

Run chains lckb_sne with `data.experiments=['sne']`

Analyze: `python ... info chains/lckb_sne chains/lckb_bao`

In `lckb_sne/plots`:



Experiments and Likelihoods

Simplest example ever: Prior on H_0 (1103.2976)

The Hubble Space Telescope measured $h_{obs} = 0.738 \pm 0.024$

$$\log(\mathcal{L}) = -\frac{1}{2} \frac{(h_{th} - h_{obs})^2}{\sigma_h^2}$$

Likelihood (montepython/likelihoods/hst/__init__):

```
from montepython.likelihood_class import Likelihood_prior
class hst(Likelihood_prior):
    def loglkl(self, cosmo, data):
        h = cosmo.h()
        loglkl = -0.5 * (h - self.h) ** 2 / (self.sigma ** 2)
        return loglkl
```

Data (montepython/likelihoods/hst/hst.data):

```
# Values for Hubble Space Telescope (following 1103.2976)
hst.h      = 0.738
hst.sigma  = 0.024
```

Likelihood rules

- Likelihoods in directory `montepython/likelihoods/l_name`
- Needed files: `__init__.py` and `l_name.data`
- `__init__.py` defines a class, inheriting from `Likelihood`
- Contains function `loglkl` $\rightarrow \log(\mathcal{L})$

Introducing your own Likelihoods

- Follow the above rules
- Inspire yourself with the examples
- \exists similar likelihood? \rightarrow you can inherit its methods!
- You can use additional python packages

(See also B. Audren's lecture on likelihoods)

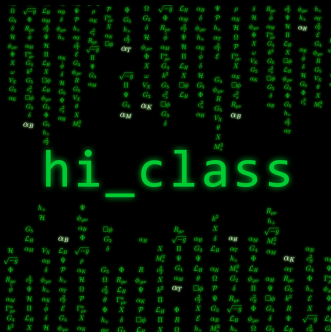
Conclusions

- Brings all the power of CLASS to Python
- Easy to run chains and analyze likelihoods
- Many available experiments
- Advantages from object oriented features in python
 - Add likelihoods
 - Add samplers or other features
- This just scratches the surface, many more options!

(See also B. Audren's slides)

The hi_class academy

Coming soon!



- Set of interrelated projects:
 - ★ Theory & model building
 - ★ Implementation and phenomenology
 - ★ Compare with data
- Collaboration → Publishable results
 - ★ Review of models
 - ★ Observational constraints
- Stay tuned for more info!

www.hiclass-code.net