# Lecture II: The Background

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struct background ba;
with fields ba.blabla, or through the pointer pba:
struct background * pba;
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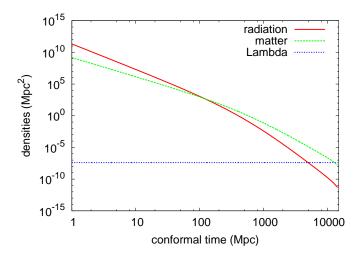
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struct background * pba;
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- the goal of this module is to solve the background evolution and store the
  results in a table. It should provide a function able to interpolate within this
  table at any value of time.
- other modules should be able to know all background quantities (densities, pressures, Hubble rate, angular/luminosity distances, etc.) at any given time or redshift.

• at the end of this lecture, we also see who to plot the background evolution.



### Units

Units assume c=1 and all quantities in  $\mathrm{Mpc}^n$ 

• times and distances are in Mpc: conformal time  $\tau$  in Mpc,  $H = \frac{a'}{a^2}$  in Mpc<sup>-1</sup>.

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- all densities and pressures appearing in the code are in fact some rescaled variables:

$$\rho = \frac{8\pi G}{3} \rho_{\rm physical}, ~~ p = \frac{8\pi G}{3} p_{\rm physical}, ~~ \text{in Mpc}^{-2}. \label{eq:rho_physical}$$

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So the Friedmann equation reads

$$H = \left(\sum_{i} \rho_i - \frac{K}{a^2}\right)^{1/2}$$

with the curvature K also in  $\mathrm{Mpc}^{-2}$ .



# The function background\_functions()

Most quantities can be immediately inferred from a given value of  $\boldsymbol{a}$  without integrating any differential equations:

$$\bullet$$
  $\rho_i = \Omega_i^0 H_0^2 \left(\frac{a}{a_0}\right)^{-3(1+w_i)}$ 

$$\bullet$$
  $p_i = w_i \rho_i$ 

$$\bullet \ H = \left(\sum_{i} \rho_{i} - \frac{K}{a^{2}}\right)^{1/2}$$

$$\bullet H' = \left(-\frac{3}{2}\sum_{i}(\rho_i + p_i) + \frac{K}{a^2}\right)a$$

$$\bullet$$
  $\rho_{\rm crit} = H^2$ 

• 
$$\Omega_i = \rho_i/\rho_{\rm crit}$$

These quantities are all returned by a function background\_functions(pba,a,...)

But to get them as a function of time we need to integrate one differential equation:

$$a' = a^2 H$$

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- comoving sound horizon:  $r_s'=c_s$ , since  $r_s=\int_{\tau_{\rm ini}}^{\tau_0}c_sd\tau$ , with a squared sound speed in the photon+baryon+electron fluid

$$c_s^2 = \frac{\delta p_\gamma}{\delta \rho_\gamma + \delta \rho_b} = \frac{1}{3(1 + [3\rho_b/4\rho_\gamma])}.$$



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• linear growth factor of density perturbations in minimal  $\Lambda$ CDM model (filled with dust),  $D'=1/(aH^2)$  (such that  $\delta_m \propto D$ ).



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Exemple of extended cosmology with quintessence  $\phi$  (see Thomas's lecture):

- $\{A\} = \{\rho_i, p_i, H, ..., V_{\phi}, \rho_{\phi}, p_{\phi}\}$
- $\bullet \ \{B\} = \{a, \phi, \phi'\}$

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Exemple of decaying dark matter with non-trivial differential equation giving  $\rho_{dm}(t)$ :

- $\{A\} = \{\rho_i, p_i, H, ...\}$
- $\bullet \{B\} = \{a, \rho_{dm}\}$



Reflected by arguments of

background\_functions(pba, pvecback\_B, format, pvecback)

# Input background parameters

In the \*.ini file, the user may pass:

- Hubble: h or HO
- Photons: T\_cmb or Omega\_g or omega\_g
- Ultra-relativvistic relics: N\_ur or Omega\_ur or omega\_ur
- CDM: Omega\_cdm or omega\_cdm
- Non-cold DM: ncdm: lots of possible input, see dedicated lectures
- Decaying CDM plus its relat. decay product: Omega\_dcdmdr or omega\_dcdmdr
- Curvature: Omega\_k
- Cosmological constant: Omega\_Lambda
- $\bullet$  Fluid: Omega\_fld, w0\_fld, wa\_fld, cs2\_fld (assuming CLP:  $w=w_0+w_a(1-a/a_0)$  and  $\delta p=c_s^2\delta\rho)$

# Input background parameters

#### Just one convention to remember !!!!

One of Omega\_Lambda or Omega\_fld must be left unspecified, to let the code match with  $H_0$ :

$$H_0^2 = \sum_i \rho_i^0 - K/a_0^2$$

If the two are passed, there is an error message.

All details on the syntax and on these rules are explicitly written in the comments of explanatory.ini

# Remark on the component called "fluid"

#### Remark on the fluid:

- $\rho_i = \Omega_i^0 H_0^2 \left(\frac{a}{a_0}\right)^{-3(1+w_i)}$  is only valid when  $w_i = \text{constant}$ .
- for  $w=w_0+w_a(1-a/a_0)$ , an analytic integration of the energy conservation equation gives

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Finally there is a parameter background\_verbose=... (one verbose parameter for each module). 0 gives no output at all, 1 the standard output that you see by default, etc.

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- ullet background\_tau\_of\_z(pba, z, &tau) returns au(z), can be useful just before calling background\_at\_tau(pba, tau,...)

- background\_indices(...) allocates dynamically all indices (see next slides)
- background\_functions(pba, pvecback\_B,...) gets  $\{B\}$  and returns  $\{A\}$ .
- background\_solve(...) integrates the differential system for  $\{B\}$ ,  $\{C\}$  variables
- ullet background\_initial\_conditions(...) assigns initial conditions  $\{B_{\mathrm{ini}}\}$ ,  $\{C_{\mathrm{ini}}\}$  to this system, e.g.  $a( au_{\mathrm{ini}})$  assuming pure radiation domination since Big Bang
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#### These functions call auxiliary functions in tools/:

- for array interpolation, tools/array.c (contains functions with all operations on arrays: interpolation, extrapolation, integration, derivation, smoothing, ...)
- for the integration of Ordinary Differential Equations, generic\_integrator(), which can be set either to rkck or ndf15 (see dedicated lecture). We will find this feature anytime we need to integrate ODE (also in thermodynamics and in perturbations).

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- Then we declare all possible indices index\_bg\_<blai> in common/background.h (more precisely, inside the structure background, because these indices are necessary for manipulating the background table).

# Dynamical indexing

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- Then we declare all possible indices index\_bg\_<blabla> in common/background.h (more precisely, inside the structure background, because these indices are necessary for manipulating the background table).
- We also declare flags saying whether these indices need to be defined or not.

# Dynamical indexing

#### In include/background.h:

```
struct background {
    /** input parameters with assigned in the input module*
    double OmegaO_cdm;
    /** flags and indices **/
    int has_cdm; // can take values _TRUE_ or _FALSE_
    . . . .
    int index_bg_rho_cdm;
    . . .
    int bg_size;
    /** interpolation table **/
    double * background_table;
```

# Dynamical indexing

In source/background.c, the function background\_indices() called at the beginning of background\_init() assigns numerical value to indices, that the user will never need to know (quantities always written symbolically as y[pba->index\_bg\_rho\_cdm])

```
int background_indices(pba,...) {
    /* initialize all flags */
    if (pba->0mega0_cdm != 0.)
       pba->has_cdm = _TRUE_;
    /* initialize all indices */
    index_bg=0;
    class_define_index(pba->index_bg_rho_cdm,
                       pba->has_cdm,
                       index_bg,
                       1):
    class_define_index(pba->index_bg_rho_fld,
                       pba->has_fld,
                       index_bg,
                       1):
    pba->bg_size = index_bg;
```

### Getting background quantities from outside

Calling background quantitites from another module is then very simple: e.g., in the perturbation module:

```
double * pvecback;
class_alloc(pvecback,
            pba->bg_size*sizeof(double),
            ppt->error_message);
class_call(background_at_tau(pba,tau,...,...,pvecback),
                             pba->error_message,
                             ppt->error_message);
/* We want here to compute the total background density*/
if (pba->has_cdm == _TRUE_) {
   rho_tot += pvecback[pba->index_bg_rho_cdm];
if (pba->has_fld == _TRUE_) {
   rho_tot += pvecback[pba->index_bg_rho_fld];
   p_tot += pvecback[pba->index_bg_p_fld];
. . .
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   p_tot += pvecback[pba->index_bg_p_fld];
```

5th argument can be: pba->long\_info, pba->normal\_info or pba->short\_info

# Full list of coded background quantities

Currently, the list of all available background quantities is:

```
index_bg_a
short info
                                               a.
                                               H
                 index bg H
                                               H'
                 index_bg_H_prime
normal_info
                index_bg_rho_<i>
                                               \rho for _b, _g,_cdm,_ur,_fld,_lambda
                variables for ncdm
                                               see dedicated lecture
                 index_bg_Omega_r
                                               \Omega_{\rm radiation}
                 index_bg_rho_crit
 long_info
                                               \rho_{\rm crit}
                 index_bg_Omega_m
                                               Ω<sub>matter</sub>
                 index_bg_conf_distance
                                               \tau_0 - \tau = \gamma
                 index_bg_ang_distance
                                              d A = a r
                                               d_{L} = (1+z)^{2} d_{A}
                 index_bg_lum_distance
                                               proper time t
                 index_bg_time
                 index_bg_rs
                                               conformal sound horizon r_s
                 index_bg_D
                                               density growth factor of \Lambda CDM
                                               velocity growth factor of \LambdaCDM
                 index bg f
```

(with 
$$r = \chi$$
, or  $\sin(\sqrt{K}\chi)/\sqrt{K}$ , or  $\sinh(\sqrt{-K}\chi)/\sqrt{-K}$ )



# To conclude on dynamical indexing in CLASS

These rules for flags and indices are followed everywhere in the code, with maybe 50 different lists of indices index\_ab\_blabla. In the background module, two such lists:

• for all variables in the table ( $\{A\}$ ,  $\{B\}$ ,  $\{C\}$ ),

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int index_bg_a;
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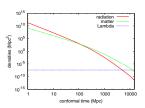
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• for the ODE dy[i] = f(y[j]), i.e. for variables  $\{B\}$ ,  $\{C\}$ ,

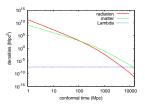
```
int index_bi_a;
int index_bi_time;
int index_bi_rs;
int index_bi_growth;
int bi_size;
```

declared in include/background.h outside the background structure, and erased/forgotten after the execution of background\_init().

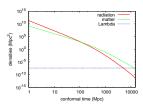
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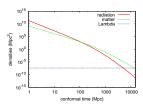
Two ways to produce such a plot with the quantitates of your choice (maybe customised to your own model):



The output module will also write a file output/toto\_background.dat with header:

- # Table of selected background quantitites
- # All densities are mutiplied by (8piG/3)
- # z, proper time [Gyr], conformal time \* c [Mpc], H/c [1/Mpc] (etc.)

Two ways to produce such a plot with the quantitates of your choice (maybe customised to your own model):

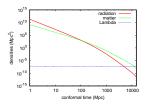


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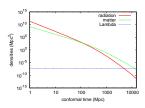
Output easy to customise in output.c, by editing: output\_one\_line\_of\_background(...) for the quantities to plot in each line output\_open\_background\_file(...) for the header (description of columns)

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2. good to know: directory test/ contains several test codes executing only part of the main(...) function.

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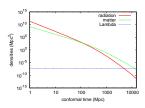


2. good to know: directory test/ contains several test codes executing only part of the main(...) function.

For instance: test/test\_background.c only executes input\_init(...), background\_init(...), and then outputs a table of all background quantities. Useful also for quick debugging!
Usage:

- > make test\_background
- > ./test\_background myinput.ini

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- Usage:
- > make test\_background
- > ./test\_background myinput.ini

Same with test\_thermodynamics.c, test\_perturbations.c, test\_nonlinear.c, test\_transfer.c...



# Background exercises (see exercise sheet for more details)

#### Exercise IIa

Reproduce this plot from the Dodelson book *Modern Cosmology*, using the plotting software of your choice.

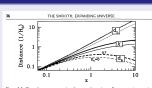


Figure 2.3. Three distance measures in a flat sepanding universe. From top to bottom, the luminosity distance, the comoving distance, and the angular diameter distance. The pair of lines in each case is for a flat universe with matter only (light curves) and 70% cosmological constant A (heavy curves). In a A-dominated universe, distances out to fixed redshift are larger than in a matter-dominated universe.

#### Exercise IIb

Add a new matter species with an equation of state  $p=w\rho$ . Visualise its evolution with time. Provoke an error on purpose, to check the error format.

By following a few general rules, we get automatically some very informative error messages like:

```
Error in thermodynamics_init
=>thermodynamics_init(L:292) :error in
    thermodynamics_helium_from_bbn(ppr,pba,pth);
=>thermodynamics_helium_from_bbn(L:1031) :condition (omega_b
    > omegab[num_omegab-1]) is true; You have asked for an
    unrealistic high value omega_b = 7.350000e-02. The
    corrresponding value of the primordial helium fraction
    cannot be found in the interpolation table. If you
    really want this value, you should fix YHe to a given
    value rather than to BBN
```

We only wrote the piece starting with "You have asked...". All the rest was generated automatically by the code. This follows from following everywhere 5 rules.

#### Rule 1:

All functions are of type int, and return either \_SUCCESS\_ or \_FAILURE\_ (defined internally in include/common.h: #define \_SUCCESS\_ 0 , #define \_FAILURE\_ 1 )

```
int function(input, &output) {
    ...
    if (something goes wrong) return _FAILURE_;
    ...
    return _SUCCESS_;
}
```

#### Rule 2:

All functions are called with the macro class\_call(.,.,.) (all macros class\_xxx(...) are defined in include/common.h):

#### Rule 3:

Each of the 9 main structures xx has a field called error\_message. Any function in the module xxx.c is called xxx\_something() and writes its error message in xx.error\_message (if pxx is a pointer to xx, in pxx->error\_message).

So if we are in perturb\_init() and we call perturb\_indices() we write:

But if we are in perturb\_init() and we call background\_at\_tau() we write:

#### Rule 4:

Whenever an error could occur, we first write a test with the macro class\_test(.,,,):

#### Example:

In the text, no need to say in which function we are, or to write that the number of points is zero, or to put a \n, all this is done automatically.

#### Rule 5:

Always allocate memory with the macros class\_alloc(), class\_calloc(), class\_realloc().

Instead of

```
malloc(parray, N*sizeof(double));
```

use

```
class_alloc(parray, N*sizeof(double), pxx->error_message);
```

If allocation fails (N too big, null or negative), the function will automatically return a <code>\_FAILURE\_</code> and the code will return an appropriate error message:

```
Error running background_init
=>background_init(L:537):error in background_solve(ppr,pba);
=>background_solve(L:1303):could not allocate pvecback with
    size -8
```

Final remark: in main/class.c there is no "higher level" so the 10 initialisation functions are called like e.g.:

But when CLASS is called as a function from another code (e.g. Monte Python or test/test\_loops.c) we can use the standard way:

Then, CLASS never crashes (in principle...), it only returns class(...) == \_FAILURE\_