# Milestone 1: The background evolution of the universe

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**Abstract.** I set up the background cosmology of the universe.

### 1. Introduction

In this project I will follow the algorithm presented in Callin (2005)[1] for simulating the cosmic microwave background. This is part one of four for this project. The first part involves computing the expansion history of the universe, as well as looking at the evolution of the density of various matter and energy components. This sets up the background cosmology such that we in the next step can turn to perturbations. I have chosen to make this first part compatible with the inclusion of neutrinos. This can of course be removed by setting the neutrino density to zero and raising the radiation density accordingly.

To ease the development of this code, I have been provided with a skeleton code of the project. This code includes a variety of methods needed to solve the project. These methods will be explained as they are used throughout the project.

## 2. Equations

For the background cosmology I use the standard Friedmann-Lemaître-Robertson-Walker(FLRW) metric for flat space. This gives the line element.

$$ds^{2} = -dt^{2} + a^{2}(t)(r^{2}(d\theta^{2} + sin^{2}\theta d\phi^{2})$$
  
=  $a^{2}(\eta)(-d\eta^{2} + r^{2}(d\theta^{2} + sin^{2}\theta d\phi^{2}))$  (1)

where a(t) is the scale factor, and  $\eta$  conformal time. I also introduce a parameter x defined as

$$x = ln(a) \tag{2}$$

The redshift is defined as

$$1 + z = \frac{a_0}{a} \tag{3}$$

We assume that the universe consist of cold dark matter (CDM, m), baryons (b), radiation(r), neutrinos( $\nu$ ), and a cosmological constant ( $\Lambda$ ). With these components, the Hubble parameter H becomes

$$H = \frac{1}{a}\frac{da}{dt} = H_0\sqrt{(\Omega_m + \Omega_b)a^{-3} + (\Omega_r + \Omega_{\nu})a^{-4} + \Omega_{\Lambda}}.$$
 (4)

I also introduce a scaled H, namely

$$\mathcal{H} = \frac{1}{a} \frac{da}{d\eta} \equiv \frac{\dot{a}}{a} = aH$$

$$= H_0 \sqrt{(\Omega_m + \Omega_b)a^{-1} + (\Omega_r + \Omega_\nu)a^{-2} + \Omega_\Lambda a^2}.$$
(5)

Note that the dot means derivative with respect to conformal time. Why this is useful will become apparent later. We also want the derivative of this with respect to x.

$$\frac{d\mathcal{H}}{dx} = H_0 \frac{-(\Omega_m + \Omega_b)e^{-x} - 2(\Omega_r + \Omega_v)e^{-2x} + 2\Omega_\Lambda e^{2x}}{\sqrt{(\Omega_m + \Omega_b)e^{-x} + (\Omega_r + \Omega_v)e^{-2x} + \Omega_\Lambda e^{2x}}}.$$
 (6)

 $\Omega_x$  is the relative density of component x compared to the critical density  $\rho_c$  needed for the universe to be flat.

$$\Omega_x = \frac{\rho_x}{\rho_c} \tag{7}$$

$$\rho_c = \frac{3H^2}{8\pi G} \tag{8}$$

We want to keep track of the densities of each component.

$$\rho_m = \rho_{m,0} a^{-3} \tag{9}$$

$$\rho_b = \rho_{b,0} a^{-3} \tag{10}$$

$$\rho_r = \rho_{r,0} a^{-4} \tag{11}$$

$$\rho_{\nu} = \rho_{\nu,0} a^{-4} \tag{12}$$

$$\rho_{\Lambda} = \rho_{\Lambda,0} \tag{13}$$

We will also need to know the distance to the horizon at different times. This can be found by noting that

$$\frac{d\eta}{dt} = \frac{c}{a}$$

which can be rewritten such that

$$\frac{d\eta}{da} = \frac{c}{a\mathcal{H}} \tag{14}$$

There are two ways to solve this differential equation numerically. One can either intergrate directly, or one can use an ordinary differential equation solver(ODE solver). Since the skeleton code contains an ODE solver I have chosen the latter. (See section 3 for more information).

## 3. Implementation

The programming language of choice is Fortran90. This is chosen because of its speed, and because the skeleton code provided was written in it.

The first step is to set up arrays for a, x,  $\eta$ , all five  $\Omega_x$ , and  $\rho_x$ . We will also need arrays for  $\rho_c$ , H, and z.

We then set the first x value to correspond to  $a=10^{-10}$ , and the last to a=0 (today). I've chosen to use 1000 points for these arrays and make the steps equal in the x array. After that one computes the a, and z values for each of the points in this x array. Because of this we now have linear steps in x. With that done, one computes H, H,  $\rho_x$ , and  $\Omega_x$  for each x value.

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The last thing to find is  $\eta(x)$ . To do this we need to use an ODE solver on equation 14. For this we must have some initial value for the function. The equation we then want to solve is

$$\eta(a) = \int_0^a \frac{da'}{a'\mathcal{H}(a')} \tag{15}$$

This can be found by considering that  $a\mathcal{H} \to H_0 \sqrt{\Omega_r + \Omega_v}$  as  $a \to 0$ . This gives

$$\eta(a) = \frac{a}{H_0 \sqrt{\Omega_r + \Omega_v}} \tag{16}$$

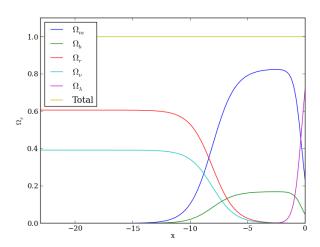
for small a. And thus one has an initial value.

Plugging this into the ODE solver returns  $\eta(x)$  for all values in the x array. Thus completing all calculations needed for the first part of the project. The ODE solver uses the Bulirsch–Stoer algorithm. The steplength is set to be one hundredth the length between two neighbouring x values.

As preparation for the next part I also spline the resulting  $\eta(x)$  values. These are then integrated and evaluated for a new set of x values not equal to but inside the x values used to make the spline. I have overplotted this in the plot for  $\eta(x)$  and the functions overlap.

## 4. Results

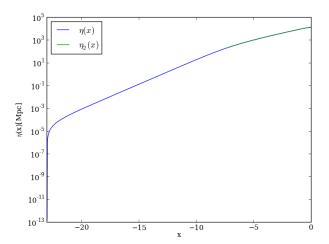
The  $\Omega_x$  values indicate the relative density of a given component compared to the critical density of the universe. The critical density being that which corresponds to the universe begin flat. Looking at the figure we see radiation dominating from the start together with neutrinos. This continues for some time until the dark matter component starts to rise. We see that this starts before the baryon component. This is good since that makes it possible for dark matter to form structures that baryons can later fall into. The radiation and neutrinos die out, and at a later point vacuum energy shoots up. At the end we end up with approximately 70% vacuum energy(dark energy?), 25% dark matter, and 5% baryons. This is exactly as it should be since that is what they were set to be at the present. See figure 1.



**Fig. 1.** The figure shows the evolution of the relative densities. As described they behave as they should.

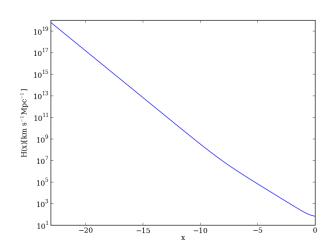
Conformal time, denoted  $\eta(x)$  measures the distance to the particle horizon for a given x value. A nice test of this is to insert

the x value of today. The answer should then be the radius of the observable universe. At the present this is measured to be approximately 14 billion parsecs (14Gpc)[2] The graph hits this value fairly well, indicating that everything is working properly so far. Note also that there are in fact two graphs in this figure, one which is calculated from the differential equation for  $\eta$ . And another one made by splining the first one and finding  $\eta$  values at arbitrary values between those of the first function.



**Fig. 2.** The conformal time hits the value of today fairly accurately. The splined function hits the computed values so well that they overlap. Here  $\eta(x)$  is the computed function, while  $\eta_2(x)$  is the splined one.

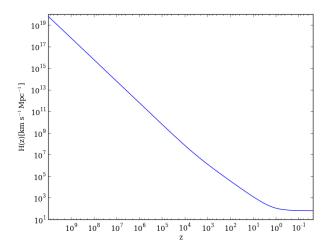
The Hubble parameter H is depicted both as a function of x, and z. Whether this is good or not is hard to say for the first part. We can at least put some faith in its precision by the fact that it ends at  $H_0 \approx 70 \,\mathrm{km \ s^{-1} Mpc^{-1}}$ , which is what  $H_0$  was set to be.



**Fig. 3.** The figure depicts the Hubble parameter as a function of x.

#### 5. Conclusions

The background cosmology of the universe is done and everything has worked the way it was expected to. There are some values that I do not have an intuition for what should be. Namely the Hubble parameter at times before today. However, I expect



**Fig. 4.** The figure depcits the Hubble parameter as a function of redshift *z*. Note that *z* increases to the left, such that time goes towards the right.

these to be correct since everything else has given results generally considered to be correct, and the Hubble parameter is computed directly from these values. With this, the code is ready for the introduction of perturbations in part two.

## 6. References

- [1] P. Callin, astro-ph/0606683
- [2] I. Bars and J. Terning, Extra Dimensions in Space and Time, Springer 2010

#### 7. Source code

The source code for the time\_mod file is included for inspection. Note that the code makes use of several different files, one with various parameters, as well as the ODE solver, and the spline.

```
module time_mod
 use healpix_types
 use params
 use spline_1D_mod
  use ode_solver
  implicit none
  integer(i4b)
                                                                ! Number of x-values
                                          :: n_t
  real(dp),
               allocatable, dimension(:) :: x_t
                                                                ! Grid of relevant x-values
                                                                ! Grid of relevant a-values
  real(dp),
               allocatable, dimension(:) :: a_t
 real(dp),
               allocatable, dimension(:) :: eta_t
                                                                ! Grid of relevant eta-values
                                                                ! Number of eta grid poins
  integer(i4b)
                                          :: n_eta
 real(dp),
               allocatable, dimension(:) :: z_eta !Grid points for eta
 real(dp),
               allocatable, dimension(:) :: x_eta
                                                                 ! Grid points for eta
 real(dp),
               allocatable, dimension(:) :: a_eta
                                                                ! Grid points for eta
                                                                ! Eta and eta'' at each grid point
 real(dp),
               allocatable, dimension(:) :: eta, eta2
               allocatable, dimension(:) :: dydx
  real(dp).
                :: rho_m0 !matter density today
 real(dp)
 real(dp)
             :: rho_b0 !baryong density today
  real(dp)
            :: rho_r0 !radiation density today
            :: rho_nu0 !neutrino density today
  real(dp)
  real(dp)
           :: rho_lambda0 !vacuum energy density today
  real(dp),
               allocatable, dimension(:) :: rho_m !matter density
  real(dp).
               allocatable, dimension(:) :: rho_b !baryong density
  real(dp),
               allocatable, dimension(:) :: rho_r !radiation density
  real(dp),
               allocatable, dimension(:) :: rho_nu !neutrino density
               allocatable, dimension(:) :: rho_lambda !vacuum energy density
  real(dp),
  real(dp),
               allocatable, dimension(:) :: Omega_mx !Relative densities
               allocatable, dimension(:) :: Omega_bx
  real(dp),
               allocatable, dimension(:) :: Omega_rx
 real(dp),
               allocatable, dimension(:) :: Omega_nux
 real(dp),
 real(dp),
               allocatable, dimension(:) :: Omega_lambdax
               allocatable, dimension(:) :: H !Huble constant as func of x
 real(dp),
contains
  subroutine initialize_time_mod
    implicit none
    integer(i4b) :: i, n, n1, n2
    real(dp)
                 :: z_start_rec, z_end_rec, z_0, x_start_rec, x_end_rec, x_0
                 :: dx, x_eta1, x_eta2, a_init,h1,eta_init,a_end,rho_crit0,rho_crit
    real(dp)
    real(dp)
                 :: eps,hmin,yp1,ypn
    ! Define two epochs, 1) during and 2) after recombination.
                = 200
                                             ! Number of grid points during recombination
   n1
                                             ! \ {\tt Number \ of \ grid \ points \ after \ recombination}\\
                = 300
   n2
                                             ! Total number of grid points
                = n1 + n2
   n_t
    z_start_rec = 1630.4d0
                                             ! Redshift of start of recombination
    z_{end}rec = 614.2d0
                                             ! Redshift of end of recombination
                = 0.d0
    z_0
                                             ! Redshift today
```

```
x_start_rec = -log(1.d0 + z_start_rec) ! x of start of recombination
                                         ! x of end of recombination
x_{end} = -\log(1.d0 + z_{end})
v_0
            = 0.d0
                                         ! x today
n_eta
           = 1000
                                         ! Number of eta grid points (for spline)
                                         ! Start value of a for eta evaluation
a_init
           = 1.d-10
           = 1.d0
a_end
          = log(a_init)
                                         ! Start value of x for eta evaluation
x_eta1
          = 0.d0
                                         ! End value of x for eta evaluation
x_eta2
eta_init = a_init/(H_0*sqrt(Omega_r+Omega_nu))
eps = 1.d-10
hmin = 0.d0
! Task: Fill in x and a grids ( These will be used in later milestones)
allocate(x_t(n_t))
do i = 0,n1-1! Fill interval during recombination
   x_t(i+1) = x_start_rec + i*(x_end_rec-x_start_rec)/(n1-1)
end do
do i = 1,n2 !Fill from end of recomb to today
    x_t(n1+i) = x_end_rec + (i)*(x_0-x_end_rec)/(n2)
!write(*,*) x_t !print x_t to terminal
allocate(a_t(n_t+1))
a_t = exp(x_t)!fill the a grid using the x grid
!write(*,*) a_t !print a_t to terminal
!Allocate and fill a,x, and z arrays
allocate(a_eta(n_eta))
allocate(x_eta(n_eta))
allocate(z_eta(n_eta))
x_{eta}(1) = x_{eta}(1)
do i = 1,n_eta-1
    x_{eta}(i+1) = x_{eta1} + i*(x_{eta2}-x_{eta1})/(n_{eta-1})
end do
a_{eta} = exp(x_{eta})
z_{eta} = 1.d0/a_{eta} - 1.d0
!write(*,*) z_eta
!write(*,*) size(z_eta)
!print *, "x"
!write(*,*) x_eta(1)
!write(*,*) x_eta(-1)
!print *, "a"
!write(*,*) a_eta(1)
!write(*,*) a_eta(-1)
!print *, "z"
!write(*,*) z_eta(1)
!write(*,*) z_eta(-1)
!Calculate the various densities for each scale factor
rho\_crit0 = 3.d0*H\_0**2.d0/(8.d0*pi*G\_grav)
rho_m0 = Omega_m *rho_crit0
rho_b0 = Omega_b
                       *rho_crit0
```

```
= Omega_r
    rho_r0
                          *rho_crit0
    rho_nu0 = Omega_nu
                          *rho_crit0
   rho_lambda0 = Omega_lambda*rho_crit0
    allocate(rho_m(n_eta))
    allocate(rho_b(n_eta))
    allocate(rho_r(n_eta))
    allocate(rho_nu(n_eta))
    allocate(rho_lambda(n_eta))
    allocate(Omega_mx(n_eta))
    allocate(Omega_bx(n_eta))
    allocate(Omega_rx(n_eta))
    allocate(Omega_nux(n_eta))
    allocate(Omega_lambdax(n_eta))
    allocate(H(n_eta))
   do i=1,n_{eta+1}
   H(i) = get_H(x_eta(i))
   Omega_rx(i) = Omega_r *H_0**2.d0/H(i)**2.d0 *a_eta(i)**-4.d0
    Omega_nux(i) = Omega_nu *H_0**2.d0/H(i)**2.d0 *a_eta(i)**-4.d0
    Omega_lambdax(i) = Omega_lambda *H_0**2.d0/H(i)**2.d0
    end do
    !End of density calculations
    allocate(eta(n_eta+1))
    eta(1) = eta_init !Start value of eta
   h1 = abs(1.d-2*(a_eta(1)-a_eta(2))) !Defines the steplength
    allocate(dydx(1))
    do i = 2, n_{eta+1}
eta(i) = eta(i-1)
       call odeint(eta(i:i),a_eta(i-1) ,a_eta(i), eps, h1, hmin, eta_derivs, bsstep, output)
    end do
    !write(*,*) eta !check that eta gives reasonable values
    !Spline eta and place the second derivative of
    !this function in eta2
    allocate(eta2(n_eta+1))
    yp1 = 1.d30
    ypn = 1.d30
    call spline(a_eta, eta, yp1, ypn, eta2)
    allocate(eta_t(n_t+1))
    do i=1,n_t+1
      eta_t(i) = get_eta(x_t(i))
    end do
  end subroutine initialize_time_mod
```

!Begin Stuff needed to make odeint work

```
subroutine eta_derivs(a, eta, dydx) !Define the derivative d/da(eta)
                     use healpix_types
                             implicit none
                             real(dp),
                                                                                                                    intent(in) :: a
                            real(dp), dimension(:), intent(in) :: eta
                            real(dp), dimension(:), intent(out) :: dydx
real(dp) :: H_p
real(dp) :: x
                            x = log(a)
H_p = get_H_p(x)
                            dydx = c/(a*H_p)
    end subroutine eta_derivs
    subroutine output(x, y)
                             use healpix_types
                             implicit none
                             real(dp),
                                                                                                                    intent(in) :: x
                            real(dp), dimension(:), intent(in) :: y
    end subroutine output
    !End Stuff needed to make odeint work
    ! Task: Write a function that computes H at given x
    function get_H(x)
           implicit none
          real(dp), intent(in) :: x
          real(dp)
                                                                                       :: get_H
          real(dp)
           a = exp(x)
          get_H = H_0*sqrt((Omega_b+Omega_m)*a**-3.d0 + (Omega_r+Omega_nu)*a**-4.d0 + Omega_lambda)
    end function get_H
    ! Task: Write a function that computes H' = a*H at given x
    function get_H_p(x)
           implicit none
          real(dp), intent(in) :: x
          real(dp)
                                                                                     :: get_H_p
          real(dp)
                                                   :: a
           a = exp(x)
           get_H_p = a*get_H(x)
    end function get_H_p
    ! Task: Write a function that computes dH'/dx at given x
    function get_dH_p(x)
           implicit none
          real(dp), intent(in) :: x
          real(dp)
                                                                                        :: get_dH_p
          \label{eq:get_dh_p} get_dH_p = H_0/2.d0*1/sqrt((Omega_m+Omega_b)*exp(-x)+Omega_r*exp(-2.d0*x) & \\ & (Omega_m+Omega_b)*exp(-x)+Omega_r*exp(-2.d0*x) & \\ & (Omega_m+Omega_b)*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*exp(-x)+Omega_r*e
          + Omega_lambda*exp(2.d0*x) * (-(Omega_m+Omega_b)*exp(-x)-2.d0*Omega_r*exp(-2.d0*x) & (-(Omega_m+Omega_b)*exp(-x)-2.d0*Omega_r*exp(-2.d0*x) & (-(Omega_m+Omega_b)*exp(-x)-2.d0*Omega_r*exp(-x)-2.d0*x) & (-(Omega_m+Omega_b)*exp(-x)-2.d0*Omega_r*exp(-x)-2.d0*x) & (-(Omega_m+Omega_b)*exp(-x)-2.d0*Omega_r*exp(-x)-2.d0*x) & (-(Omega_m+Omega_b)*exp(-x)-2.d0*Omega_r*exp(-x)-2.d0*x) & (-(Omega_m+Omega_b)*exp(-x)-2.d0*Omega_r*exp(-x)-2.d0*x) & (-(Omega_m+Omega_b)*exp(-x)-2.d0*x) & (-(Omega_m
          + 2.d0*Omega_lambda*exp(2.d0*x))
    end function get_dH_p
    ! Task: Write a function that computes eta(x), using the previously precomputed splined function
    function get_eta(x_in)
           implicit none
          real(dp), intent(in) :: x_in
          real(dp)
                                                                                      :: get_eta
          real(dp)
                                                  :: a_in
           a_{in} = exp(x_{in})
```

get\_eta = splint(a\_eta, eta, eta2, a\_in)
end function get\_eta

end module time\_mod