

Milestone 1 Code

Daniel Herman

February 2018

1 Code

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1 module time_mod
2   use healpix_types
3   use params
4   use spline_1D_mod
5   use ode_solver
6   implicit none
7
8   integer(i4b)                                :: n_t                !
9   Number of x-values
10  real(dp), allocatable, dimension(:) :: x_t                !
11  Grid of relevant x-values
12  real(dp), allocatable, dimension(:) :: a_t                !
13  Grid of relevant a-values
14  real(dp), allocatable, dimension(:) :: eta_t              !
15  Grid of relevant eta-values
16
17  integer(i4b)                                :: n_eta            !
18  Number of eta grid points
19  real(dp), allocatable, dimension(:) :: x_eta              !
20  Grid points for eta
21  real(dp), allocatable, dimension(:) :: z_eta              !
22  Grid points for z_eta
23  real(dp), allocatable, dimension(:) :: a_eta              !
24  Grid points for a_eta
25  real(dp), allocatable, dimension(:) :: eta, eta2          !
26  Eta and eta'' at each grid point
27  real(dp), allocatable, dimension(:) :: dydx
28
29  real(dp)                                :: rho_crit0        !
30  critical density today
31  real(dp)                                :: rho_m0            !
32  matter density today
33  real(dp)                                :: rho_b0            !
34  baryon density today
35  real(dp)                                :: rho_r0            !
36  radiation density today
37  real(dp)                                :: rho_nu0           !
38  neutrino density today
39  real(dp)                                :: rho_lambda0       !
40  vacuum energy density today
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26  real(dp),    allocatable , dimension(:) :: rho_m          !
    matter density grid
27  real(dp),    allocatable , dimension(:) :: rho_b          !
    baryon density grid
28  real(dp),    allocatable , dimension(:) :: rho_r          !
    radiation density grid
29  real(dp),    allocatable , dimension(:) :: rho_nu         !
    neutrino density grid
30  real(dp),    allocatable , dimension(:) :: rho_lambda     !
    vacuum energy density grid
31  real(dp),    allocatable , dimension(:) :: Omega_mx        !
    Relative densities
32  real(dp),    allocatable , dimension(:) :: Omega_bx        ! ''
33  real(dp),    allocatable , dimension(:) :: Omega_rx        ! ''
34  real(dp),    allocatable , dimension(:) :: Omega_nux       ! ''
35  real(dp),    allocatable , dimension(:) :: Omega_lambdax    ! ''
36  real(dp),    allocatable , dimension(:) :: Hx              !
    Hubble constant as a func of x
37  real(dp)                                :: m2mpc           !
    Value for changing m to Mpc
38  real(dp)                                :: kmmpc
39
40  contains
41
42  subroutine initialize_time_mod
43      implicit none
44
45      integer(i4b) :: i, n, n1, n2
46      real(dp)      :: z_start_rec, z_end_rec, z_0, x_start_rec,
    x_end_rec, x_0
47      real(dp)      :: dx, x_eta1, x_eta2, a_init, a_end, eta_init
48      real(dp)      :: eps, hmin, ypl, ypn, h1
49
50      ! Define two epochs, 1) during and 2) after recombination.
51
52      !-----
53      ! Initialize the values for variables that are used during the
    calculations
54
55      n1          = 200                                ! Number of grid points
    during recombination
56      n2          = 300                                ! Number of grid points
    after recombination
57      n_t         = n1 + n2                            ! Total number of grid
    points
58
59      z_start_rec = 1630.4d0                            ! Redshift of start of
    recombination
60      z_end_rec   = 614.2d0                              ! Redshift of end of
    recombination
61      z_0         = 0.d0                                ! Redshift today
62
63      x_start_rec = -log(1.d0 + z_start_rec)             ! x of start of
    recombination
64      x_end_rec   = -log(1.d0 + z_end_rec)               ! x of end of
    recombination
65      x_0         = 0.d0                                ! x today

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66      n_eta      = 1000                      ! Number of eta grid
67      points (for spline)
68      a_init     = 1.d-10                    ! Start value of a for
eta evaluation
69      a_end      = 1.d0
70      x_eta1     = log(a_init)                ! Start value of x for
eta evaluation
71      x_eta2     = 0.d0                      ! End value of x for
eta evaluation

72
73      eps = 1.d-10
74      hmin = 0.d0
75
76      m2mpc = 3.2408d-23                      ! 3.086*10^22m = 1Mpc
77      kmpc = 3.086d19                         ! ''
78
79      !-----
80      ! Task: Fill in x and a grids
81
82      allocate(x_t(n_t))
83      do i = 0,n1-1                            ! Fill recombination
interval
84          x_t(i+1) = x_start_rec + i*(x_end_rec-x_start_rec)/(n1-1)
85      end do
86
87      do i = 1,n2                            ! Fill post recombination
interval
88          x_t(n1+i) = x_end_rec + i*(x_0-x_end_rec)/(n2)
89      end do
90
91      allocate(a_t(n_t))
92      a_t = exp(x_t)                            ! Fill the a_t-grid using
the x_t values
93
94      ! Task: 1) Compute the conformal time at each eta time step
95      !         2) Spline the resulting function, using the provided "
spline" routine in spline_1D_mod.f90
96      allocate(x_eta(n_eta))
97      allocate(a_eta(n_eta))
98      allocate(z_eta(n_eta))
99      allocate(eta2(n_eta))
100
101      ! Compute the x_eta values
102      x_eta(1) = x_eta1
103      do i = 1,n_eta-1
104          x_eta(i+1) = x_eta1 + i*(x_eta2-x_eta1)/(n_eta-1)
105      end do
106
107      ! Compute the a_eta and z_eta values
108      a_eta = exp(x_eta)
109      z_eta = 1.d0/a_eta - 1.d0
110
111      !-----
112      ! Density calculations
113
114      rho_crit0 = 3.d0*H_0**2.d0/(8.d0*pi*G_grav)

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115 rho_m0      = Omega_m      * rho_crit0
116 rho_b0      = Omega_b      * rho_crit0
117 rho_r0      = Omega_r      * rho_crit0
118 rho_nu0     = Omega_nu     * rho_crit0
119 rho_lambda0 = Omega_lambda * rho_crit0
120
121 allocate(rho_m(n_eta))
122 allocate(rho_b(n_eta))
123 allocate(rho_r(n_eta))
124 allocate(rho_nu(n_eta))
125 allocate(rho_lambda(n_eta))
126 allocate(Omega_mx(n_eta))
127 allocate(Omega_bx(n_eta))
128 allocate(Omega_rx(n_eta))
129 allocate(Omega_nux(n_eta))
130 allocate(Omega_lambdax(n_eta))
131 allocate(Hx(n_eta))
132
133 do i=1,n_eta
134     Hx(i) = get_H(x_eta(i))*kmmpc
135     Omega_mx(i) = Omega_m      *H_0**2.d0/Hx(i)**2.d0 *
a_eta(i)**-3.d0
136     Omega_bx(i) = Omega_b      *H_0**2.d0/Hx(i)**2.d0 *
a_eta(i)**-3.d0
137     Omega_rx(i) = Omega_r      *H_0**2.d0/Hx(i)**2.d0 *
a_eta(i)**-4.d0
138     Omega_nux(i) = Omega_nu     *H_0**2.d0/Hx(i)**2.d0 *
a_eta(i)**-4.d0
139     Omega_lambdax(i) = Omega_lambda*H_0**2.d0/Hx(i)**2.d0
140 end do
141
142 ! Print the density values to file
143
144 open(50, file='densities.dat')
145 do i=1,n_eta
146     write(50, '(5(E17.8))') Hx(i),Omega_mx(i),Omega_bx(i),
Omega_rx(i),Omega_lambdax(i)
147 end do
148 close(50)
149
150 !-----
151 ! Calculate the eta values, initializing eta(1) using the
assumption that the universe
! is radiation dominated before recombination
152
153
154 allocate(eta(n_eta))
155 eta(1) = 1.d-10/(H_0*sqrt(Omega_r+Omega_nu)) ! early universe
is rad/neutrino dominated
156 h1 = abs(1.d-2*(a_eta(1)-a_eta(2)))
157 allocate(dydx(1))
158
159 do i =2,n_eta
160     eta(i) = eta(i-1)
161     call odeint(eta(i:i),a_eta(i-1),a_eta(i),eps,h1,hmin,
eta_derivs,bsstep,output)
162 end do
163

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164     eta = c*eta                                ! eta is defined by c/a
165     eta = m2mpc*eta                            ! Put eta in units of cMpc
166
167     ! Print x,a, and eta values to file for plotting
168     open(54, file='data.dat')
169     do i=1,n_eta
170         write(54,'(4(E17.8))') x_eta(i),a_eta(i),eta(i),z_eta(i)
171     end do
172     close(54)
173
174 end subroutine initialize_time_mod
175
176 !-----
177
178 subroutine eta_derivs(a,eta,dydx) ! Defining the eta deriv
179
180 use healpix_types
181 implicit none
182
183 real(dp), intent(in) :: a
184 real(dp), dimension(:), intent(in) :: eta
185 real(dp), dimension(:), intent(out) :: dydx
186 real(dp) :: H_p
187 real(dp) :: x
188
189 x = log(a)
190 H_p = get_H_p(x)
191 dydx = c/(a*H_p)
192
193 end subroutine eta_derivs
194
195 subroutine output(x,y)
196 use healpix_types
197 implicit none
198
199 real(dp), intent(in) :: x
200 real(dp), dimension(:), intent(in) :: y
201
202 end subroutine output
203
204 !-----Done with subroutines-----
205 !-----Define Functions-----
206
207 ! Task: Write a function that computes H at given x
208 function get_H(x)
209 implicit none
210
211 real(dp), intent(in) :: x
212 real(dp) :: get_H
213 real(dp) :: a
214 a = exp(x)
215
216 get_H = H_0*sqrt((Omega_b+Omega_m)*a**3 + (Omega_r +
217 Omega_nu)*a**4 + Omega_lambda)
218
219 end function get_H

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220 ! Task: Write a function that computes  $H' = a \cdot H$  at given x
221 function get_H_p(x)
222     implicit none
223
224     real(dp), intent(in) :: x
225     real(dp)              :: get_H_p
226     real(dp)              :: a
227     a = exp(x)
228
229     get_H_p = a*get_H(x)
230
231 end function get_H_p
232
233 ! Task: Write a function that computes  $dH'/dx$  at given x
234 function get_dH_p(x)
235     implicit none
236
237     real(dp), intent(in) :: x
238     real(dp)              :: get_dH_p
239     get_dH_p = H_0/2.d0*1/sqrt((Omega_m+Omega_b)*exp(-x)+Omega_r*
240     exp(-2.d0*x) &
241     + Omega_lambda*exp(2.d0*x)) * (-(Omega_m+Omega_b)*exp(-x)-2.d0*
242     Omega_r*exp(-2.d0*x) &
243     + 2.d0*Omega_lambda*exp(2.d0*x))
244
245 end function get_dH_p
246
247 ! Task: Write a function that computes eta(x), using the
248 ! previously precomputed splined function
249 function get_eta(x_in)
250     implicit none
251
252     real(dp), intent(in) :: x_in
253     real(dp)              :: get_eta
254     real(dp)              :: a_in
255     a_in = exp(x_in)
256     get_eta = splint(a_eta,eta,eta2,a_in)
257
258 end function get_eta
259
260 end module time_mod

```