

Milestone 3

Elisabeth Strøm¹
AST5220 16.04.2018

¹ Institute of Theoretical Astrophysics, University of Oslo

These results are in accordance with those obtained by Callin (2006).

1. Introduction

The full equation set outside the tight coupling regime, is as follows;

2. Method

Here we present the methods we have used to obtain the results in Section 3.

2.1. Theory

The initial conditions when not including polarization, are

$$\Phi = 1 \quad (1)$$

$$\delta = \delta_b = \frac{3}{2}\Phi \quad (2)$$

$$v = v_b = \frac{ck}{2\mathcal{H}}\Phi \quad (3)$$

$$\Theta_0 = \frac{1}{2}\Phi \quad (4)$$

$$\Theta_1 = -\frac{ck}{6\mathcal{H}}\Phi \quad (5)$$

$$\Theta_2 = -\frac{20ck}{45\mathcal{H}\tau'}\Theta_1 \quad (6)$$

$$\Theta_l = -\frac{l}{2l+1}\frac{ck}{\mathcal{H}\tau'}\Theta_{l-1} \quad (7)$$

$$(8)$$

$$R = \frac{4\Omega_r}{3\Omega_b a} \quad (10)$$

$$\Theta'_0 = -\frac{ck}{\mathcal{H}}\Theta_1 - \Phi', \quad (11)$$

$$\Theta'_1 = \frac{ck}{3\mathcal{H}}\Theta_0 - \frac{2ck}{3\mathcal{H}}\Theta_2 + \frac{ck}{3\mathcal{H}}\Psi + \tau' \left[\Theta_1 + \frac{1}{3}v_b \right], \quad (12)$$

$$\Theta'_2 = \frac{2ck}{5\mathcal{H}}\Theta_1 - \frac{3ck}{5\mathcal{H}}\Theta_3 + \frac{9\tau'}{10}\Theta_l, \quad (13)$$

$$\text{for } l < l_{\max}: \quad (14)$$

$$\Theta'_l = \frac{lck}{(2l+1)\mathcal{H}}\Theta_{l-1} - \frac{(l+1)ck}{(2l+1)\mathcal{H}}\Theta_{l+1} + \tau'\Theta_l, \quad (15)$$

$$\Theta_{l_{\max}} = \frac{ck}{\mathcal{H}}\Theta_{l_{\max}-1} - c\frac{l_{\max}+1}{\mathcal{H}\eta(x)}\Theta_{l_{\max}} + \tau'\Theta_{l_{\max}}, \quad (16)$$

$$\Phi' = \Psi - \frac{1}{3}\left(\frac{ck}{\mathcal{H}}\right)^2\Phi + \frac{1}{2}\left(\frac{H_0}{\mathcal{H}}\right)^2 \quad (17)$$

$$\cdot [\Omega_m a^{-1}\delta + \Omega_b a^{-1}\delta_b + 4\Omega_r a^{-2}\Theta_0], \quad (18)$$

$$\delta = \frac{ck}{\mathcal{H}}v - 3\Phi', \quad (19)$$

$$\delta'_b = \frac{ck}{\mathcal{H}}v_b - 3\Phi', \quad (20)$$

$$v' = -v - \frac{ck}{\mathcal{H}}\Psi, \quad (21)$$

$$v'_b = -v_b - \frac{ck}{\mathcal{H}}\Psi + \tau'R(3\Theta_1 + v_b). \quad (22)$$

$$(23)$$

When within the tight-coupling regime, there are certain changes, namely for $\Theta_{l>0}$ and Θ'_l . We also need to require the equation of baryon velocity, as for small values of

We also have a general algebraic expression for Ψ , which we can implement when needed,

$$\Psi = -\Psi - 12\left(\frac{H_0}{cka}\right)^2\Omega_r\Theta_2 \quad (9)$$

$$q = \frac{1}{(1+R)\tau' + \frac{\mathcal{H}'}{\mathcal{H}} - 1} \left[-[(1-2R)\tau' + (1+R)\tau''] \cdot (3\Theta_1 + v_b) - \frac{ck}{\mathcal{H}}\Psi' + \left(1 - \frac{\mathcal{H}'}{\mathcal{H}}\right)\frac{ck}{\mathcal{H}}(2\Theta_2 - \Theta_0) - \frac{ck}{\mathcal{H}}\Theta'_0 \right] \quad (24)$$

$$v'_b = \frac{1}{1+R} \left[-v_b - \frac{ck}{\mathcal{H}} \Psi + R \left(q + \frac{ck}{\mathcal{H}} (2\Theta_2 - \Theta_0) - \frac{ck}{\mathcal{H}} \Psi \right) \right], \quad (25)$$

$$\Theta'_1 = \frac{1}{3} (q - v'_b), \quad (26)$$

$$\Theta_2 = -\frac{20ck}{45\mathcal{H}\tau'} \Theta_1, \quad (27)$$

$$\Theta_l = -\frac{l}{2l+1} \frac{ck}{\mathcal{H}\tau'} \Theta_{l-1}. \quad (28)$$

2.2. Implementation

3. Results

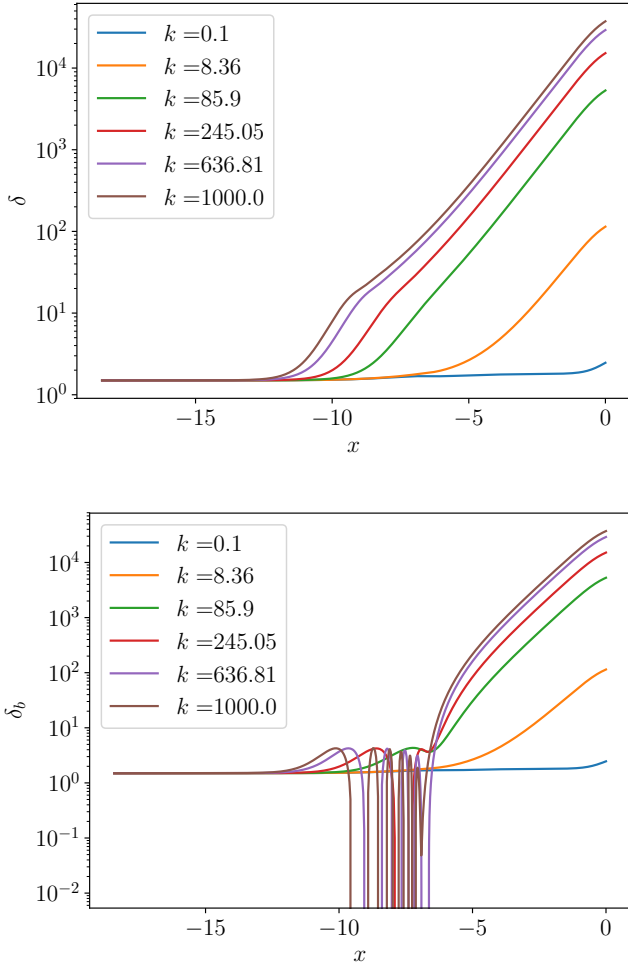


Fig. 1

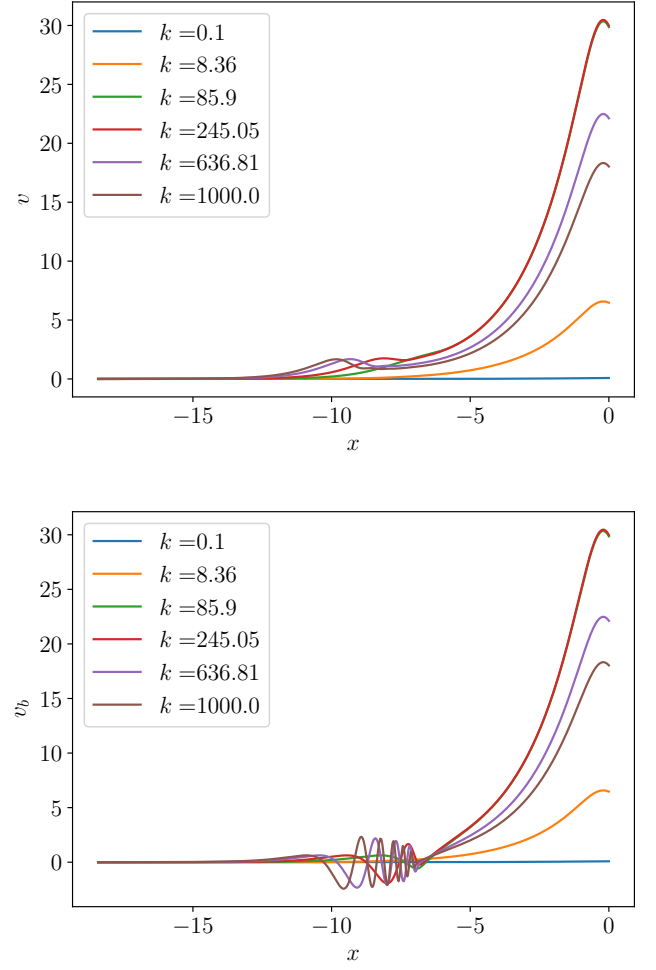


Fig. 2

4. Conclusions

References

Callin, P. 2006, ArXiv Astrophysics e-prints

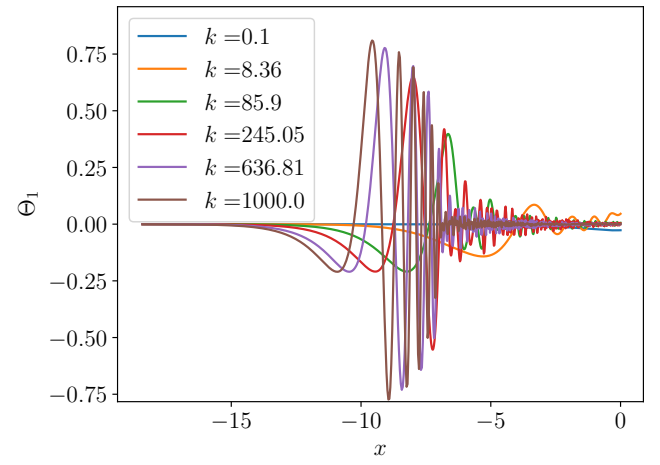
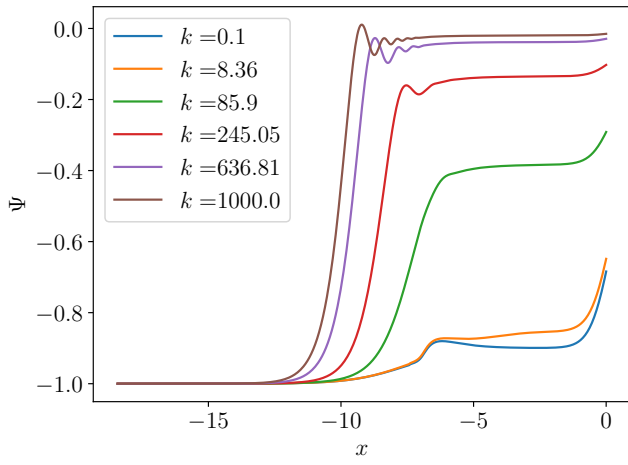
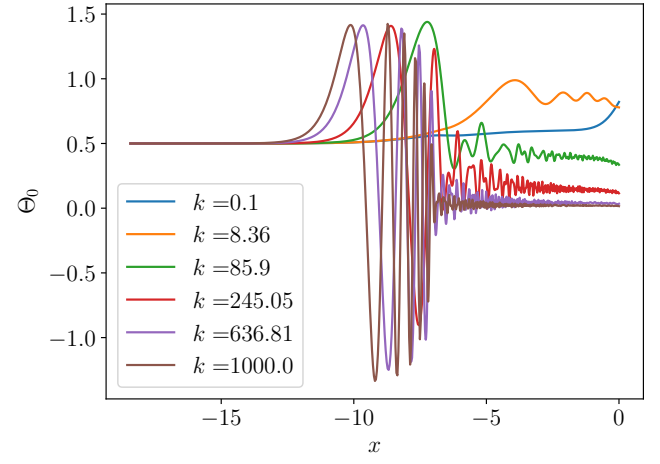
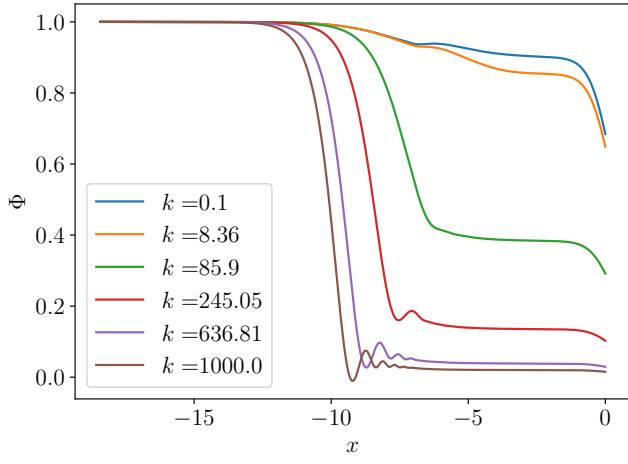


Fig. 3

Fig. 4

5. Appendix

Source code

Listing 1: C:/Users/elini/Documents/AST5220/Ast5220/src/evolution_mod.f90

```
module evolution_mod
  use healpix_types
  use params
  use time_mod
  use ode_solver
  use rec_mod
  implicit none

  ! Accuracy parameters
  real(dp), parameter, private :: a_init = 1.d-8
  real(dp), parameter, private :: x_init = log(a_init)
  real(dp), parameter, private :: k_min = 0.1d0 * H_0 / c
  real(dp), parameter, private :: k_max = 1.d3 * H_0 / c
  integer(i4b), parameter :: n_k = 100
  integer(i4b), parameter, private :: lmax_int = 6

  ! Perturbation quantities
  real(dp), allocatable, dimension(:, :, :) :: Theta
  real(dp), allocatable, dimension(:, :) :: delta
  real(dp), allocatable, dimension(:, :) :: delta_b
  real(dp), allocatable, dimension(:, :) :: Phi
  real(dp), allocatable, dimension(:, :) :: Psi
  real(dp), allocatable, dimension(:, :) :: v
  real(dp), allocatable, dimension(:, :) :: v_b
  real(dp), allocatable, dimension(:, :) :: dPhi
  real(dp), allocatable, dimension(:, :) :: dPsi
  real(dp), allocatable, dimension(:, :) :: dv_b
  real(dp), allocatable, dimension(:, :, :) :: dTheta

  ! Fourier mode list
  real(dp), allocatable, dimension(:) :: ks

  ! Book-keeping variables
  real(dp), private :: k_current
  integer(i4b), private :: npar = 6+lmax_int

  real(dp), private :: ck, H_p, ckH_p, dt

contains

  ! NB!!! New routine for 4th milestone only; disregard until then!!!
  subroutine get_hires_source_function(k, x, S)
    implicit none

    real(dp), pointer, dimension(:), intent(out) :: k, x
    real(dp), pointer, dimension(:, :), intent(out) :: S

    integer(i4b) :: i, j
    real(dp) :: g, dg, ddg, tau, dt, ddt, H_p, dH_p, ddHH_p, Pi, dPi, ddPi
    real(dp), allocatable, dimension(:, :) :: S_lores

    ! Task: Output a pre-computed 2D array (over k and x) for the
    !       source function, S(k,x). Remember to set up (and allocate) output
    !       k and x arrays too.
    !
    ! Substeps:
    ! 1) First compute the source function over the existing k and x
    !    grids
    ! 2) Then spline this function with a 2D spline
    ! 3) Finally, resample the source function on a high-resolution uniform
    !    5000 x 5000 grid and return this, together with corresponding
    !    high-resolution k and x arrays
```

```

end subroutine get_hires_source_function

! Routine for initializing and solving the Boltzmann and Einstein equations
subroutine initialize_perturbation_eqns
  implicit none

  integer(i4b) :: l, i

  ! Task: Initialize k-grid, ks; quadratic between k_min and k_max
  allocate(ks(n_k))
  do i = 1, n_k
    ks(i) = k_min + (k_max-k_min)*((i-1.d0)/(n_k-1.d0))**2.d0
  end do

  ! Allocate arrays for perturbation quantities
  allocate(Theta(0:n_t, 0:lmax_int, n_k))
  allocate(delta(0:n_t, n_k))
  allocate(delta_b(0:n_t, n_k))
  allocate(v(0:n_t, n_k))
  allocate(v_b(0:n_t, n_k))
  allocate(Phi(0:n_t, n_k))
  allocate(Psi(0:n_t, n_k))
  allocate(dPhi(0:n_t, n_k))
  allocate(dPsi(0:n_t, n_k))
  allocate(dv_b(0:n_t, n_k))
  allocate(dTheta(0:n_t, 0:lmax_int, n_k))
  ! Task: Set up initial conditions for the Boltzmann and Einstein equations
  !Theta(:, :, :) = 0.d0
  !dTheta(:, :, :) = 0.d0
  !dPhi(:, :) = 0.d0
  !dPsi(:, :) = 0.d0

  Phi(0, :) = 1.d0
  delta(0, :) = 1.5d0 * Phi(0, :)
  delta_b(0, :) = delta(0, :)
  Theta(0, 0, :) = 0.5d0*Phi(0, :)
  H_p = get_H_p(x_init)
  dt = get_dtau(x_init)

  do i = 1, n_k
    ckH_p = c*ks(i)/H_p

    v(0, i) = ckH_p/2.d0*Phi(0, i)
    v_b(0, i) = v(0, i)

    Theta(0, 1, i) = -ckH_p/6.d0*Phi(0, i)
    Theta(0, 2, i) = -20.d0/45.d0*ckH_p/(dt)*Theta(0, 1, i)
    do l = 3, lmax_int
      Theta(0, l, i) = - 1/(2.d0*l + 1.d0)*ckH_p/dt *Theta(0, l-1, i)
    end do
    Psi(0, i) = - Phi(0, i) - 12.d0*(H_0/(c*ks(i)*a_init))**2.d0*Omega_r*Theta(0, 2, i)
  end do

end subroutine initialize_perturbation_eqns

subroutine integrate_perturbation_eqns
  implicit none

  integer(i4b) :: i, j, k, l, i_tc
  real(dp) :: x1, x2
  real(dp) :: eps, hmin, h1, x_tc, t1, t2

  real(dp), allocatable, dimension(:) :: y, y_tight_coupling, dydx

  eps = 1.d-8
  hmin = 0.d0

```

```

h1      = 1.d-5

allocate(y(npar))
allocate(dydx(npar))
allocate(y_tight_coupling(7))

! Propagate each k-mode independently
do k = 1, n_k
  write(*,*) "starting k loop in integrate"
  k_current = ks(k) ! Store k_current as a global module variable
  ck = c*k_current

  ! Initialize equation set for tight coupling
  y_tight_coupling(1) = delta(0,k)
  y_tight_coupling(2) = delta_b(0,k)
  y_tight_coupling(3) = v(0,k)
  y_tight_coupling(4) = v_b(0,k)
  y_tight_coupling(5) = Phi(0,k)
  y_tight_coupling(6) = Theta(0,0,k)
  y_tight_coupling(7) = Theta(0,1,k)

  ! Find the time to which tight coupling is assumed,
  ! and integrate equations to that time
  write(*,*) "entering get_tight_coupling_time"
  x_tc = get_tight_coupling_time(k_current)
  write(*,*) "x_tc", x_tc
  write(*,*) "k", k
  ! Task: Integrate from x_init until the end of tight coupling, using
  !       the tight coupling equations
  write(*,*) "integrating tight coupling equations"
  i_tc = 1

  do while(x_t(i_tc) < x_tc)
    !write(*,*) "evol i_tc lopp!", i_tc
    ! Integration while tc
    call odeint(y_tight_coupling, x_t(i_tc-1), x_t(i_tc), eps, h1, hmin, dy_tc_dx, bsstep, output)
    ! some parameters
    ckH_p = ck*get_H_p(x_t(i_tc))
    dt     = get_dtau(x_t(i_tc))

    delta(i_tc,k) = y_tight_coupling(1)
    delta_b(i_tc,k) = y_tight_coupling(2)
    v(i_tc,k) = y_tight_coupling(3)
    v_b(i_tc,k) = y_tight_coupling(4)
    Phi(i_tc,k) = y_tight_coupling(5)
    Theta(i_tc,0,k) = y_tight_coupling(6)
    Theta(i_tc,1,k) = y_tight_coupling(7)
    Theta(i_tc,2,k) = -20.d0/45.d0*ckH_p/dt * Theta(i_tc,1,k)
    do l = 3, lmax_int
      Theta(i_tc,l,k) = - 1/(2.d0*1 + 1.d0)*ckH_p/dt *Theta(i_tc,l-1,k)
    end do
    Psi(i_tc,k) = - Phi(i_tc,k) - 12.d0*(H_0/(ck*a_t(i_tc)))*2.d0*Omega_r*Theta(i_tc,2,k)

    ! The store derivatives necessary here?
    call dy_tc_dx(x_t(i_tc), y_tight_coupling, dydx)
    dv_b(i_tc,k) = dydx(4)
    dPhi(i_tc,k) = dydx(5)
    dTheta(i_tc,0,k) = dydx(6)
    dTheta(i_tc,1,k) = dydx(7)
    dTheta(i_tc,2,k) = 2.d0/5.d0*ckH_p*Theta(i_tc,1,k) -
      3.d0/5.d0*ckH_p*Theta(i_tc,3,k)+dt*0.9d0*Theta(i_tc,2,k)

    do l=3,lmax_int-1
      dTheta(i_tc,l,k) = 1/(2.d0*1+1.d0)*ckH_p*dTheta(i_tc,l-1,k) -
        (1+1.d0)/(2.d0*1+1.d0)*ckH_p*dTheta(i_tc,l+1,k) + dt*Theta(i_tc,l,k)
    end do
    dPsi(i_tc,k) = -dPhi(i_tc,k) - 12.d0*(H_0/(ck*a_t(i_tc)))*2.d0
      *Omega_r*(-2.d0*Theta(i_tc,2,k)+dTheta(i_tc,2,k))
  end while
end do

```

```

        i_tc = i_tc+1
    end do ! end while do

    ! Task: Set up variables for integration from the end of tight coupling
    ! until today
    y(1:7) = y_tight_coupling(1:7)
    y(8) = Theta(i_tc-1,2,k)
    do l = 3, lmax_int
        y(6+l) = Theta(i_tc-1,l,k)
    end do

    write(*,*) "integrating non-tight coupling equations"
    do i = i_tc, n_t-1

        !write(*,*) "after tc loop", i
        ! Task: Integrate equations from tight coupling to today
        call odeint(y, x_t(i-1), x_t(i),eps, h1, hmin, dy_dx, bsstep, output)
        ! Task: Store variables at time step i in global variables
        !write(*,*) "made it through"
        delta(i,k) = y(1)
        delta_b(i,k) = y(2)
        v(i,k) = y(3)
        v_b(i,k) = y(4)
        Phi(i,k) = y(5)

        do l = 0, lmax_int
            Theta(i,l,k) = y(6+l)
        end do

        Psi(i,k) = - Phi(i,k) - 12.d0*(H_0/(ck*a_t(i)))**2.d0*Omega_r*Theta(i,2,k)

        ! Task: Store derivatives that are required for C_l estimation
        call dy_dx(x_t(i), y, dydx)
        dv_b(i,k) = dydx(4)
        dPhi(i,k) = dydx(5)
        do l=0, lmax_int
            dTheta(i,l,k) = dydx(6+l)
        end do
        dPsi(i,k) = -dPhi(i,k) - 12.d0*(H_0/(ck*a_t(i)))**2 * Omega_r*(dTheta(i,2,k)-2.d0*Theta(i,2,k))
    end do

end do

deallocate(y_tight_coupling)
deallocate(y)
deallocate(dydx)

end subroutine integrate_perturbation_eqns

! Task: Complete the following routine, such that it returns the time at which
!       tight coupling ends. In this project, we define this as either when
!        $d\tau < 10$  or  $c*k/(H_p*dt) > 0.1$  or  $x > x(\text{start of recombination})$ 
function get_tight_coupling_time(k)
    implicit none

    real(dp), intent(in) :: k
    real(dp) :: get_tight_coupling_time
    real(dp) :: x, x_start_rec, z_start_rec
    integer(i4b) :: i, n

    z_start_rec = 1630.4d0 ! Redshift of start of recombination
    x_start_rec = -log(1.d0 + z_start_rec) ! x of start of recombination

    n=1d4
    do i=0,n
        x = x_init + i*(0.d0- x_init)/n
        dt = get_dtau(x)
        H_p = get_H_p(x)
    end do
end function

```

```

    if (abs(dt) > 10.d0 .and. abs((c*k/H_p)/dt) <= 0.1d0 .and. x<=x_start_rec) then
        get_tight_coupling_time = x
    end if

end do

end function get_tight_coupling_time

subroutine dy_tc_dx(x, y, dydx)
! Tight coupling, only l=0,1 for dTheta
use healpix_types
implicit none
real(dp),          intent(in) :: x
real(dp), dimension(:), intent(in) :: y
real(dp), dimension(:), intent(out) :: dydx

real(dp) :: delta, delta_b, v, v_b, Phi, Theta0, Theta1, Theta2, Psi
real(dp) :: ddelta, ddelta_b, dv, dv_b, dPhi, dTheta0, dTheta1
real(dp) :: q, R, a, dH_p, ddt

delta      = y(1)
delta_b    = y(2)
v          = y(3)
v_b        = y(4)
Phi        = y(5)
Theta0     = y(6)
Theta1     = y(7)

a          = exp(x)
dt         = get_dtau(x)
ddt        = get_ddtau(x)
H_p        = get_H_p(x)
dH_p       = get_dH_p(x)
ckH_p      = ck/H_p

! Derivatives
Theta2     = - 20.d0*ckH_p/(45.d0*dt) * Theta1
R          = (4.d0*Omega_r)/(3.d0*Omega_b*a)
Psi        = - Phi - 12.d0*(H_0/(ck*a))**2.d0 * Omega_r * Theta2

dPhi       = Psi - ckH_p**2.d0/3.d0*Phi + 0.5d0*(H_0/H_p)**2.d0 * (Omega_m/a*delta + Omega_b/a*delta_b +
4.d0*Omega_r*Theta0/a**2.d0)
dv         = - v - ckH_p * Psi

ddelta     = ckH_p * v - 3.d0*dPhi
ddelta_b   = ckH_p * v_b - 3.d0*dPhi

dTheta0    = - ckH_p*Theta1 - dPhi
!----- special for tight coupling -----
q          = (-((1.d0-2.d0*R)*dt + (1.d0+R)*ddt)*(3.d0*Theta1 + v_b)- ckH_p*Psi +
(1.d0-(dH_p/H_p))*ckH_p*(-Theta0 + 2.d0*Theta2) - ckH_p*dTheta0)/((1.d0+R)*dt + (dH_p/H_p) -1.d0)

dv_b       = (1.d0/(1.d0 + R)) * (-v_b - ckH_p*Psi + R*(q + ckH_p*(-Theta0 + 2.d0*Theta2) - ckH_p*Psi))
dTheta1    = (1.d0/3.d0)*(q-dv_b)
! -----

! Final array
dydx(1) = ddelta
dydx(2) = ddelta_b
dydx(3) = dv
dydx(4) = dv_b
dydx(5) = dPhi
dydx(6) = dTheta0
dydx(7) = dTheta1

end subroutine dy_tc_dx

subroutine dy_dx(x, y, dydx)
! we define dy/dx

```



```

use healpix_types
implicit none
real(dp),          intent(in) :: x
real(dp), dimension(:), intent(in) :: y
real(dp), dimension(:), intent(out) :: dydx

integer(i4b) :: l
real(dp) :: delta, delta_b, v, v_b, Phi, Theta0, Theta1, Theta2, Psi
real(dp) :: ddelta, ddelta_b, dv, dv_b, dPhi, dTheta0, dTheta1
real(dp) :: q, R, a, eta

! what we take in, use in derivation
delta = y(1)
delta_b = y(2)
v = y(3)
v_b = y(4)
Phi = y(5)
Theta0 = y(6)
Theta1 = y(7)
Theta2 = y(8)
! Theta3-6: y(9)-y(12)

a = exp(x)
eta = get_eta(x)
dt = get_dtau(x)
H_p = get_H_p(x)
ckH_p = ck/H_p

! Derivatives
R = (4.d0*Omega_r)/(3.d0*Omega_b*a)
Psi = - Phi - 12.d0*(H_0/(ck*a))**2.d0 * Omega_r * Theta2
dPhi = Psi - ckH_p**2.d0/3.d0*Phi + 0.5d0*(H_0/H_p)**2.d0 * (Omega_m/a*delta + Omega_b/a*delta_b +
4.d0*Omega_r*Theta0/a**2.d0)

dv = - v - ckH_p*Psi
dv_b = - v_b - ckH_p*Psi + dt*R*(3.d0*Theta1 + v_b)

ddelta = ckH_p * v - 3.d0*dPhi
ddelta_b = ckH_p * v_b - 3.d0*dPhi

dTheta0 = - ckH_p*Theta1 - dPhi
dTheta1 = ckH_p/3.d0*Theta0 - 2.d0/3.d0*ckH_p*Theta2 + ckH_p/3.d0*Psi + dt*(Theta1 + v_b/3.d0)

! dTheta2 - dTheta5
do l = 2, lmax_int-1
    dydx(6+l) = 1/(2.d0*l+1.d0)*ckH_p*y(6+l-1) - (l+1.d0)/(2.d0*l + 1.d0)*ckH_p*y(6+l+1) + dt*(y(6+l) -
0.1d0*y(6+l)*abs(l==2))
end do

! Final array
dydx(1) = ddelta
dydx(2) = ddelta_b
dydx(3) = dv
dydx(4) = dv_b
dydx(5) = dPhi
dydx(6) = dTheta0
dydx(7) = dTheta1
! dTheta6
dydx(6+l) = ckH_p*y(6+l-1) - c*(l+1.d0)/(H_p*eta)*y(6+l) + dt*y(6+l)

end subroutine dy_dx

subroutine write_to_file_evolution_mod
    use healpix_types
    implicit none

    integer(i4b) :: i
    integer(i4b), dimension(6) :: k

```

```

write(*,*) "writing to file; evolution_mod"

k(1:6)=(/1, 10, 30, 50, 80, 100 /)
!k(1:6)=(/1, 2, 3, 4, 5, 10 /)

!----- write to file ---
write(*,*) "opening files "
open (unit=0, file = 'k_ks.dat', status='replace')
open (unit=1, file = 'x_t.dat', status='replace')
open (unit=2, file = 'Phi.dat', status='replace')
open (unit=3, file = 'Psi.dat', status='replace')
open (unit=4, file = 'delta.dat', status='replace')
open (unit=5, file = 'delta_b.dat', status='replace')
open (unit=6, file = 'v.dat', status='replace')
open (unit=7, file = 'v_b.dat', status='replace')
open (unit=8, file = 'Theta0.dat', status='replace')
open (unit=9, file = 'Theta1.dat', status='replace')

do i=1,6
    write(0,*) k(i),ks(k(i))
end do

write(*,*) "writing stuff"
do i=0, n_t-1
    write (1,*) x_t(i)
    write (2,'(*(2X, ES14.6E3))') Phi(i,k(1)),Phi(i,k(2)),Phi(i,k(3)),Phi(i,k(4)),Phi(i,k(5)),Phi(i,k(6))
    write (3,'(*(2X, ES14.6E3))') Psi(i,k(1)),Psi(i,k(2)),Psi(i,k(3)),Psi(i,k(4)),Psi(i,k(5)),Psi(i,k(6))
    write (4,'(*(2X, ES14.6E3))')
        delta(i,k(1)),delta(i,k(2)),delta(i,k(3)),delta(i,k(4)),delta(i,k(5)),delta(i,k(6))
    write (5,'(*(2X, ES14.6E3))')
        delta_b(i,k(1)),delta_b(i,k(2)),delta_b(i,k(3)),delta_b(i,k(4)),delta_b(i,k(5)),delta_b(i,k(6))
    write (6,'(*(2X, ES14.6E3))') v(i,k(1)),v(i,k(2)),v(i,k(3)),v(i,k(4)),v(i,k(5)),v(i,k(6))
    write (7,'(*(2X, ES14.6E3))') v_b(i,k(1)),v_b(i,k(2)),v_b(i,k(3)),v_b(i,k(4)),v_b(i,k(5)),v_b(i,k(6))
    write (8,'(*(2X, ES14.6E3))')
        Theta(i,0,k(1)),Theta(i,0,k(2)),Theta(i,0,k(3)),Theta(i,0,k(4)),Theta(i,0,k(5)),Theta(i,0,k(6))
    write (9,'(*(2X, ES14.6E3))')
        Theta(i,1,k(1)),Theta(i,1,k(2)),Theta(i,1,k(3)),Theta(i,1,k(4)),Theta(i,1,k(5)),Theta(i,1,k(6))

end do

write(*,*) "closing files "
do i=0, 9
    close(i)
end do

end subroutine write_to_file_evolution_mod

end module evolution_mod

```
