

References

- Decker, R. B. Formation of Shock-Spike Events at Quasi-Perpendicular Shocks. Journal of Geophysical Research **88** (1983), 9959-9973.
- De Hoffmann, F. and Teller, E. Magneto-Hydrodynamic Shocks. Physical Review **80** (1950), 692-703.
- Krasnov, N. F. Aerodynamics 1. Mir Publishers Moscow 1985, 13-16.
- Lee, M. A. Coupled Hydrodynamic Wave Excitation and Ion Acceleration at Interplanetary Traveling Shocks. Journal of Geophysical Research **88** (1983), 6109-6119.
- Lüst, R. Magneto-Fluid Dynamic Shock Waves. Reviews of Modern Physics **32** (1960), 706-709.
- Morrison, P. The Origin of Cosmic Rays. Encyclopedia of Physics **XLVI/1** 1961, 10-38.
- Parker, E. N. Dynamics of the interplanetary gas and magnetic Fields. Astrophysical Journal **128** (1958), 664-676.
- Parks, G. K. Physics of Space Plasma Addison-Wesley Publishing Company 1991, 2-3.
- Press, W. H., Teukolsky, S. A., Vetterling, W. T., and Flannery, B. P. Numerical Recipes in C Second Edition Cambridge University Press 1992, 856-857.
- Rajamäki, M. and Saarinen, M. Accurate One-Dimensional Computation of Frontal Phenomena by PLIM. Journal of Computational Physics **111** (1994), 62-73.
- Richardson, I. G., Cane, H. V. and von Rosenvinge, T. T. MeV Ion Anisotropies at Interplanetary Shocks: Observations from the ISEE-3/ICE Medium Energy Cosmic Ray Experiment. Proceedings of the 21th International Cosmic Ray Conference 1990, 333-336.
- Ruffolo, D. Interplanetary Transport of Decay Protons from Solar Flare Neutrons. Astro-

physical Journal **382** (1991), 688-698.

Ruffolo, D. Effect of Adiabatic Deceleration on the Focused Transport of Solar Cosmic Rays. Astrophysical Journal **442** (1995), 861-874.

Ruffolo, D. Comparative Analysis of Cosmic Rays from Three Types of Solar Flares. Final Report 1996, 1-3.

Ruffolo, D. and Khumlumlert, T. Formation, Propagation, and Decay of Coherent Pulses of Solar Cosmic Rays. Geophysical Research Letters **22** (1995), 2073-2076.

Ruffolo, D. and Khumlumlert, T. Propagation of Coherent Pulses of Solar Cosmic Rays. Proceedings of the 24th International Cosmic Ray Conference 1995.

Zauderer, E. Partial Differential Equations of Applied Mathematics, Second Edition, John Wiley & Sons 1989, 48-58.



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Appendix A

Operator Splitting Method

This appendix is derived from Numerical Recipes in C of Press et al(1992). The basic idea of operator splitting or time splitting is that: Suppose we have an initial value problem in form

$$\frac{\partial F}{\partial t} = \mathcal{L}F \quad (1)$$

where \mathcal{L} is some operator not necessarily linear. Suppose that it can be written as a linear sum of m pieces as

$$\mathcal{L}F = \mathcal{L}_1F + \mathcal{L}_2F + \dots + \mathcal{L}_mF \quad (2)$$

Suppose that for each of the pieces, we already know how to update F from timestep n to timestep $n + 1$, valid as if that only operator were on the right-hand side:

$$\begin{aligned} \mathcal{L}F &= \mathcal{L}_1F \\ \mathcal{L}F &= \mathcal{L}_2F \\ &\dots \\ \mathcal{L}F &= \mathcal{L}_mF, \end{aligned} \quad (3)$$

we will write these updatings symbolically as

$$\begin{aligned} F^{n+1} &= \mathbf{F}_1(F^n, \Delta t) \\ F^{n+1} &= \mathbf{F}_2(F^n, \Delta t) \\ &\dots \\ F^{n+1} &= \mathbf{F}_m(F^n, \Delta t), \end{aligned} \quad (4)$$

respectively.

Now, one form of operator splitting would be to get from n to $n + 1$ by the following sequence of updatings:

$$\begin{aligned} F^{n+1/m} &= \mathbf{F}_1(F^n, \Delta t) \\ F^{n+2/m} &= \mathbf{F}_2(F^{n+1/m}, \Delta t) \\ &\dots \\ F^{n+1} &= \mathbf{F}_m(F^{n+(m-1)/m}, \Delta t) \end{aligned} \quad (5)$$

For example, a combined advective-diffusion equation, such as

$$\frac{\partial F}{\partial t} = -v \frac{\partial F}{\partial x} + D \frac{\partial^2 F}{\partial x^2} \quad (6)$$

might use an explicit scheme for the advective term, obtaining

$$F_i^{n+1} = F_i^{n-1} - \frac{v\Delta t}{\Delta z}(F_{i+1}^n - F_{i-1}^n) \quad (7)$$

combined with a Crank-Nicholson or other implicit scheme for the diffusion term, obtaining

$$2(1 + \varepsilon)F_i^{n+1} - \varepsilon(F_{i-1}^{n+1} + F_{i+1}^{n+1}) = 2(1 - \varepsilon)F_i^n + \varepsilon(F_{i-1}^n + F_{i+1}^n) \quad (8)$$

where $\varepsilon = D\Delta t/(\Delta z)^2$.

then F^{n+1} can be obtained easily from eqs. (7) and (8) by following the sequence of updatings in eq. (5).

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Appendix B

Program Usage

This work needs 9 files:

wind.c contains main routine for simulating the transport of cosmic rays. It gets input data from the user then calls routines in other files to do their jobs.

decel.c is used to update $F(t, \mu, z, p)$ for the effects of deceleration.

field.c contain routines that depend on the configuration of the interplanetary magnetic field. In this work magnetic field is constants but different both upstream and downstream regions. It takes into account the change of magnetic field at the shock.

initial.c is used to determine the initial values of the distribution function then send these values to the main routine in **wind.c**.

inject.c calculates the injection of the particles after the start of simulation.

nrutil.c contains routines to allocate and unallocate memory for variables in running program.

printout.c prints the data that are computed in the program.

stream.c updates F for the streaming and convection of cosmic ray particles. In the thesis **stream.c** plays important roles in transforming particles across the shock.

These files will be compiled and linked altogether, then the program will be run and it needs input data. The user must key in the following input data:

- initial value of t
- final value of t

- step size
- range for printing
- number of μ points
- length of the simulation region
- number of momentum points
- momentum (in MeV/c)
- particle mass
- solar wind speed divided by c
- scattering mean free path
- scattering power law index
- whether to print extra diagnostic information(0=no, 1=yes)

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Appendix C

Source Code for Simulation

/*

s_k_refl.c --- Febuary 24th, 1997

reflecting at boundary

s_k_noinf.c -- February 10th, 1997

Boundary conditions: no flux inflow

Given $f = 1$ at one point, we observe its behavior at
any points before and after crossing the shock.

s_k_inf.c (v) -- February 5th, 1997

Boundary conditions: inflow of 1.

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s_kink.c (v) -- November 18th, 1996

"Stationary" shock at $z_{\text{length}}/2$. BETASWD and BETASWU are defined as the (constant) solar wind speeds over C *parallel* to the magnetic field.

s_abs.c (t) -- July 6th, 1995

Change s to time.

Change s_{step} to timestep.

Change $\text{Beta}[w]$ to $1/C$ at r_1 and r_2 .

Add constant C to program by define.

s_abs.c -- March 14th, 1994

Removed μ transformations, occasionally skip extra steps instead.

stream.abs.c -- February 27th, 1994

Reworked the μ transformation to account for differential convection.

stream.abs.c -- July 15th, 1993

Inserted a missing factor of μ_{step} . Reorganized the μ transformation so that the first transformation is explicit, the second is implicit. If the latter generates negative numbers,

explicit, upwind differencing is used instead.

stream.abs.c -- February 8th, 1993

Eliminated slide() and corr[w]. Now working with z defined in a corotating frame.

Now before streaming, need to make a transformation (in mu) from the local solar wind frame to the corotating frame.

Need to transform back after streaming...

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stream.abs.c -- November 2nd, 1992

Brought back tosun[w] to keep track of the total density that streams past the inner boundary for a given w.

stream.w.abs.c -- October 29th, 1992

Added a subroutine, slide, to occasionally move f(1) to f(1+1) and to calculate corr[w], so that z in the lab frame is $(1-1+corr[w])*zstep...$

stream.w.abs.c -- September 29th, 1992

Modified for the program, wind.

stream.abs.c -- October 29th, 1990

A SUBROUTINE FOR transport.

MOVES $f(i,1)$ TO $f(i,1+i)$.

ASSUMES ABSORBING BOUNDARY CONDITIONS AT $z = 0$ AND $z = \text{length}$.

*/

#include <math.h>

#include <stdio.h>

#define C 0.1202 /* speed of light in AU/min. */

#define BDOVERBU 1.5 /* B downstream / B upstream */

#define BETASWD 0.0005 /* solar wind speed / C along B downstream
(z < length/2.0)

#define BETASWU 0.002 /* solar wind speed / C along B upstream
(z > length/2.0).

Should have BETASWU > BETASWD.

*/

#define GAMMAL 2.0 /* Power-law exponent below p[1]. */

#define TINY 1e-6

```

extern double   ***f, **fs, *tosun;

void   stream(time,timestep,beta,m,np,p,lprint,nz,zstep,nmu,printextra)
double  time, timestep, *beta, m, *p, *zstep;
int     nmu, np, printextra;
long    lprint, *nz;
{
    double  calc_pd(),calc_pu(),mustari(),refl_pr(),mu_refl();
    double  corr, dpudpd, dpudpr, fpd, fpu, frac, ftemp;
    double  lnp0, lnp1, lower, mu, mucent, mulower, muminus;
    double  munew, muplus=1.0, mustard, mustarr, mustaru, mustep;
    double  muupper, pd, pu, r1, r2, z, expon, pr;
    int     flagm, flagp=1, i, imin, parttrans, u, w, unew;
    long    k, k1, k2, l, l1, lnew;

    double  betaswc(), invert_to_pu(), invert_to_pd(), mu_d(), mu_star();
    double  mu_u();
    double  **temp;

    mustep = 2.0/(double)nmu;

    for (w=1;w<=np;w++) {
        if (w > 1) {
            lnp0 = log(p[w-1]);

```

```

    lnpi = log(p[w]);
}

for(u=1;u<=nmu;u++) for(l=1;l<=nz[w];l++) fs[l][u] = 0.0;

imin = -(nmu-1)/2;

/* constant inflow = 0 */
/*
for (u=1,i=imin;u<=nmu;u++,i++){
    if(i < 0)
        for (l=nz[w]+1+i;l<=nz[w];l++) fs[l][u] = 0.0;
    if(i > 0)
        for (l=1;l<=i;l++) fs[l][u] = 0.0;
}
*/

for (u=1,i=imin;u<=nmu;u++,i++) {
    mu = 2.0*i/(double)nmu;
    corr = 1 - mu*mu*beta[w]*beta[w];
    for (l=1;l<=nz[w];l++) {
if (f[w][l][u] > 0.0) {
        z = (1-nz[w]/2.0-0.5) * zstep[w];
        r1 = (betaswc(z)*C)*((time+ timestep)/zstep[w]);
        r2 = (betaswc(z)*C)*(time/zstep[w]);
        k1 = corr * r1 + 0.000001;
        k2 = corr * r2 + 0.000001;

```

```

k      = k1 - k2;

lnew = l + i + k;

/* reflection at left boundary */
if (lnew < 1) {
fs[1-lnew][nmu+1-u] += f[w][l][u];
}

/* reflection at right boundary */
else if ( lnew > nz[w] ) {
fs[2*nz[w]+1-lnew][nmu+1-u] += f[w][l][u];
}

/* If particles pass the inner boundary ("crash into the
Sun") then flux is lost and added to tosun[w].
*/

/* if (lnew < 1) {
tosun[w] += f[w][l][u];*/

/* If particles cross from upstream to downstream,
take special account of changes in mu and p.
*/

/* } */ else if (l > nz[w]/2 && lnew <= nz[w]/2) {

mustaru = mustar1(m,p[w],&mustard,&mustarr);

```

```

printf("mustaru=%lf\n",mustaru);

/* Case where all particles in [mu-mustep/2.0,
mu+mustep/2.0] are transmitted, or some
are transmitted and some are reflected.

This loop will handle the transmitted flux.
*/

if (mu - mustep/2.0 < mustaru) {
    parttrans = mu + mustep/2.0 > mustaru;
    if (parttrans) {
        mucent = (mustaru + mu - mustep/2.0) / 2.0;
    } else {
        mucent = mu;
    }
    pd = calc_pd(m,p[w],mucent,&dpudpd);
    printf("\n pd = %lf",pd);

    if (pd < p[w] - TINY)
        nrerror("stream: anti-acceleration");

/* Find fpu, the z-interpolated value of f at (t,z,pu) */
/* printf("\n pu = %lf p[w] = %lf",pu,p[w]); */

    if (w > 1) {
printf("\n I am in loop w>1 for transmission");

```

```

l1 = z/zstep[w-1] + nz[w-1]/2.0 + 0.5;
frac = z/zstep[w-1] + nz[w-1]/2.0 + 0.5 - (double)l1;

if (l1 < 1 || l1+1 > nz[w+1]) { /* cannot interpolate */
    lower = 0.0;
} else { /* can interpolate */
    if (f[w-1][l1][u] == 0.0 || f[w-1][l1+1][u] == 0.0) {
        lower = 0.0;
    } else {
        lower = (1.0-frac)*f[w-1][l1][u]
            + frac*f[w-1][l1+1][u];
    }
}
}

```

/* fpu is $F(p[w], \mu_{cent}, zold) * \text{pow}(2.0 - pd/p[w], \text{expon})$,

where pd is the momentum downstream

corresponding to p[w] upstream.

ftemp is $fpu * d(\mu) / d(pd)$. This will be split into

different mu-cells according to the fraction

of the interval $[\mu_u - \text{mstep}/2, \mu_u + \text{mstep}/2]$

which maps into different cells.

*/

```

if (lower <= 0.0) {

```

```

    if (printextra && l%lprint==1 && w==np)

```

```

        printf("\nNote: f[%d][%d][%d] = %le, lower = 0",

```

```

            w, l, u, f[w][l][u]);

```

```

    fpu = f[w][l][u];
} else {
    expon = (log(f[w][l][u])-log(lower))
           / (lnp1-lnp0);
    fpu = f[w][l][u] * pow(2.0-pd/p[w],expon);
}

/* Otherwise, w=1 -> extrapolate from lower p assuming
   f \propto p^-GAMMAL.
*/

} else {

/* fpu is F(p_u,mucent,zold), where p_u is the momentum
   which will change to p_d on the downstream side.
   ftemp is fpu*d(p_u)/d(p_d). This will be split into
   different mu-cells according to the fraction
   of the interval [mu_u - mustep/2,mu_u + mustep/2]
   which maps into different cells.
*/

fpu = f[w][l][u] * pow(2.0-pd/p[w],-GAMMAL);
printf("\n I am out loop w>1 for transmission");
printf("\n total transmitted fpu = %lf",fpu);

}

printf("dpudpd=%lf\n",dpudpd);
ftemp = fpu * dpudpd;

```



```

printf("\n fpu for transmitt= %lf",fpu);
printf("\n ftemp for transmitt = %lf",ftemp);

/* Will loop over all possible unew and munew
   grid points on the downstream side, and
   see how much of ftemp falls in each cell.

   muminus is munew - mustep/2.0 (lower cell boundary),
   transformed back upstream
   muplus is munew + mustep/2.0 (upper cell boundary)
   transformed back upstream

   flagm -> 1 if muminus within original cell
   flagp -> 1 if muplus within original cell

   original cell means from mu-mustep/2.0 to muupper
   in the case of partial transmission
   */

   muupper = mu+mustep/2.0;
   if (parttrans) {
/*       ftemp *= (mustaru-mu+mustep/2.0) / mustep; */
       muupper = mustaru;
printf("\n fraction = %lf", (mustaru-mu+mustep/2.0) / mustep);
printf("\n fraction of ftemp transmitted = %lf", ftemp);
   }

```

```

muminus = -1;

flagm = muminus >= mu-mustep/2.0
      && muminus < muupper;

for (unew=1;unew<=nmu &&
     (munew=(unew-0.5)*mustep-1.0)-mustep/2.0<=mustard;
     unew++,muminus=muplus,flagm=flagp) {

    if (munew+mustep/2.0 >= mustard) {
        flagp = (muplus = mustaru)
                >= mu-mustep/2.0 && muplus < muupper;
    } else {
        flagp = (muplus = mu_u(m,pd,munew+mustep/2.0))
                >= mu-mustep/2.0 && muplus < muupper;
    }

    printf("unew=%d, muplus=%lf\n",unew,muplus);

    if (muminus < mu-mustep/2.0
        && muplus >= muupper) {
        fs[lnew][unew] += ftemp;

if (lnew>=12 && lnew<=14) printf("UD1 fs[%ld] [%d] = %lf\n",lnew,unew,fs[lnew][unew]);
        break;
    } else if (!flagm && flagp) {
        fs[lnew][unew] += ftemp * (muplus-(mu-mustep/2.0))
                        / mustep;

if (lnew>=12 && lnew<=14) printf("UD2 fs[%ld] [%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    } else if (flagm && flagp) {
        fs[lnew][unew] += ftemp * (muplus-muminus)
                        / mustep;

```

```

if (lnew>=12 && lnew<=14) printf("UD3 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    } else if (flagm && !flagp) {
        fs[lnew][unew] += ftemp * (muupper-muminus)
            / mustep;
if (lnew>=12 && lnew<=14) printf("UD4 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    break;
    }
}
}

/* Case where all particles in [mu-mustep/2.0,
mu+mustep/2.0] are reflected, or some
are transmitted and some are reflected.

This loop will handle the reflected flux.
*/

if( mu+mustep/2.0 > mustaru ) {

    /* Reflection: lnew = nz[w] + 1 - lnew. */

    lnew = nz[w] + 1 - lnew;
    parttrans = mu-mustep/2.0 < mustaru;

    /* partially, not completely transmitted */

    if( parttrans ) {

```

```

mucent = (mu+mustep/2.0+mustaru)/2.0;
printf("\n partial reflection");
        } else {
mucent = mu;
printf("\n total reflection");
        }

        pr = refl_pr(m,p[w],mucent,&dpudpr);
        printf("\n pr = %lf",pr);
        /* Find lower. */

        if (w > 1) {
printf("\n I am in loop w>1 for reflection");

        l1 = z/zstep[w-1] + nz[w-1]/2.0 + 0.5;
        frac = z/zstep[w-1] + nz[w-1]/2.0 + 0.5 - (double)l1;

        if (l1 < 1 || l1+1 > nz[w+1]) { /* cannot interpolate */
            lower = 0.0;
        } else { /* can interpolate */
            if (f[w-1][l1][u] == 0.0 || f[w-1][l1+1][u] == 0.0) {
                lower = 0.0;
            } else {
                lower = (1.0-frac)*f[w-1][l1][u]
                    + frac*f[w-1][l1+1][u];
            }
        }
    }
}

```

```

/* fpu is F(p[w],mucent,zold) * pow(2.0-pd/p[w],expon),
   where pd is the momentum downstream
   corresponding to p[w] upstream.
   ftemp is fpu*d(pu)/d(pd). This will be split into
   different mu-cells according to the fraction
   of the interval [mu_u - mustep/2,mu_u + mustep/2]
   which maps into different cells.
*/

if (lower <= 0.0) {
    if (printextra && !lprint==1 && w==np)
        printf("\nNote: f[%d][%d][%d] = %le, lower = 0",
            w,l,u,f[w][l][u]);
    fpu = f[w][l][u];
} else {
    expon = (log(f[w][l][u])-log(lower))
            / (lnp1-lnp0);
    fpu = f[w][l][u] * pow(2.0-pr/p[w],expon);
}

/* Otherwise, w=1 -> extrapolate from lower p assuming
   f \propto p^-GAMMAL.
*/

} else {

/* fpu is F(p_u,mucent,zold), where p_u is the momentum

```

which will change to p_d on the downstream side.
 f_{temp} is $f_{pu} * d(p_u) / d(p_d)$. This will be split into
different μ -cells according to the fraction
of the interval $[\mu_u - \text{mustep}/2, \mu_u + \text{mustep}/2]$
which maps into different cells.

```

*/
printf("\n I am out loop w>1 for reflection");
    fpu = f[w][1][u] * pow(2.0-pr/p[w],-GAMMAL);
printf("\n total reflection fpu = %lf",fpu);
printf("\n partial reflection fpu before = %lf",pow(2.0-pr/p[w],-GAMMAL));
printf("\n partial reflection fpu = %lf",fpu);
    }

    printf("dpudpr=%lf\n",dpudpr);
    ftemp = fpu * dpudpr;
    printf("\n ftemp for reflection = %lf",ftemp);

/* Will loop over all possible unew and munew
   grid points reflected upstream, and
   see how much of ftemp falls in each cell.

   muminus is munew - mustep/2.0 (lower cell boundary),
   transformed back upstream

   muplus is munew + mustep/2.0 (upper cell boundary)
   transformed back upstream

   flagm -> 1 if muminus within original cell

```

```

        flagp -> 1 if muplus within original cell

        original cell means from mulower to mu + mustep/2.0
        in the case of partial transmission
    */

    mulower = mu-mustep/2.0;
    if (parttrans) {
/*          ftemp *= (mu+mustep/2.0-mustaru) / mustep; */
printf("\n fraction = %lf", (mu+mustep/2.0-mustaru) / mustep);
printf("\n fraction of ftemp for reflection = %lf", ftemp);
        mulower = mustaru;
    }

    muminus = mu_refl(m,pr,-mustep/2.0);
    flagm = muminus >= mulower
            && muminus < mu+mustep/2.0;
    for (unew=(nmu+1)/2; unew<=nmu &&
        (munew=(unew-0.5)*mustep-1.0)-mustep/2.0<=mustarr;
        unew++,muminus=muplus,flagm=flag) {
        if (munew+mustep/2.0 >= mustarr) {
            flagp = (muplus = mustaru)
                    >= mulower && muplus < mu+mustep/2.0;
        } else {
            flagp = (muplus = mu_refl(m,pr,munew+mustep/2.0))
                    >= mulower && muplus < mu+mustep/2.0;
        }
    }
}

```

```

}

/* Note that muminus > muplus ! */

printf("unew=%d, muplus=%lf\n",unew,muplus);
if (muplus < mulower
    && muminus >= mu+mustep/2.0) {
    fs[lnew][unew] += ftemp;
if (lnew>=12 && lnew<=14) printf("UR1 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    break;
} else if (!flagp && flagm) {
    fs[lnew][unew] += ftemp * (muminus-(mu-mustep/2.0))
        / mustep;
if (lnew>=12 && lnew<=14) printf("UR2 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    break;
} else if (flagm && flagp) {
    fs[lnew][unew] += ftemp * (muminus-muplus)
        / mustep;
if (lnew>=12 && lnew<=14) printf("UR3 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
} else if (flagp && !flagm) {
    fs[lnew][unew] += ftemp * (mu+mustep/2.0-muplus)
        / mustep;
if (lnew>=12 && lnew<=14) printf("UR4 fs[%ld][%d] = %lf",lnew,unew,fs[lnew][unew]);
}
}
}
}

```



```

/* begin */

/* If particles cross from downstream to upstream,
   take special account of changes in mu and p.

   All particles are transmitted if they cross
   from downstream to upstream.

   If mustard is the value of mu in the downstream
   solar wind frame for which mu is 0 in the de Hoffmann-
   Teller (shock) frame, then mustard < 0.

   Is there a chance that flux for mu < mustard
   will be transmitted? NO, because grid points
   with mu <= 0 are never transmitted (convection
   moves them farther from the shock).

   */

} else if (1 <= nz[w]/2 && lnew > nz[w]/2) {
    mustaru = mustar1(m,p[w],&mustard,&mustarr); /* return mu_u */

    pu = calc_pu(m,p[w],mu,&dpudpd);

    /* pu = p[w];

    dpudpd = 1.0; */

printf("pu/p[%d]=%lf, dpudpd=%lf\n",w,pu/p[w],dpudpd);

```

```

/* printf("\n pu = %lf p[w] = %lf",pu,p[w]); */

/* Find fpd, the z-interpolated value of f at (t,z,pd) */

if (pu < p[w] - TINY) nrerror("stream: anti-acceleration");
if (w > 1) {

    l1 = z/zstep[w-1] + nz[w-1]/2.0 + 0.5;
    frac = z/zstep[w-1] + nz[w-1]/2.0 + 0.5 - (double)l1;

    if (l1 < 1 || l1+1 > nz[w+1]) { /* cannot interpolate */
        lower = 0.0;
    } else { /* can interpolate */
        if (f[w-1][l1][u] == 0.0 || f[w-1][l1+1][u] == 0.0) {
            lower = 0.0;
        } else {
            lower = (1.0-frac)*f[w-1][l1][u]
                + frac*f[w-1][l1+1][u];
        }
    }
}

/* fpd is F(p[w],mu,zold) * pow(2.0-pu/p[w],expon),
   where pu is the momentum upstream
   corresponding to p[w] downstream.
   ftemp is fpd*d(p_d)/d(p_u). This will be split into
   different mu-cells according to the fraction

```

```

of the interval  $[\mu_d - \text{mustep}/2, \mu_d + \text{mustep}/2]$ 
which maps into different cells.

*/

if (lower <= 0.0) {
  if (printextra && 1%lprint==1 && w==np)
    printf("\nNote: f[%d][%d][%d] = %le, lower = 0",
           w,l,u,f[w][l][u]);
  fpd = f[w][l][u];
} else {
  expon = (log(f[w][l][u])-log(lower)) / (lnp1-lnp0);
  fpd = f[w][l][u] * pow(2.0-pu/p[w],expon);
}

/* Otherwise, w=1 -> extrapolate from lower p assuming
   f \propto p-GAMMAL.
*/

} else {
  /* fpd is F(p[w],mu,zold) * pow(2.0-pu/p[w],expon),
     where pu is the momentum upstream
     corresponding to p[w] downstream.
     ftemp is fpd*d(p_d)/d(p_u). This will be split into
     different mu-cells according to the fraction
     of the interval  $[\mu_d - \text{mustep}/2, \mu_d + \text{mustep}/2]$ 
     which maps into different cells.

```

```

*/

    fpd    = f[w][l][u] * pow(2.0-pu/p[w],-GAMMAL);
}

ftemp = fpd / dpudpd;

muplus = 1;

flagp = muplus > mu-mustep/2.0
      && muplus <= mu+mustep/2.0;
for (unew=nmu;unew>=1 &&
     (munew=(unew-0.5)*mustep-1.0)+mustep/2.0 >= mustarr;
     unew--,muplus=muminus,flagp=flagm) {

    if (munew-mustep/2.0 <= mustarr) {
        muminus = -1;
        flagm = 0;
    } else {
        flagm = (muminus = mu_d(m,pu,munew-mustep/2.0))
                > mu-mustep/2.0 && muminus <= mu+mustep/2.0;
    }
    printf("unew=%d, muminus=%lf\n",unew,muminus);
    if (muminus <= mu-mustep/2.0
        && muplus > mu+mustep/2.0) {
        fs[lnew][unew] += ftemp;
    }
    if (lnew>=12 && lnew<=14) printf("DU1 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    break;
} else if (!flagm && flagp) {

```

```

        fs[lnew][unew] += ftemp * (muplus-(mu-mustep/2.0))
                                / mustep;

if (lnew>=12 && lnew<=14) {
    printf("DU2 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    printf("\t muminus=%lf, muplus=%lf\n",muminus,muplus);
    printf("\t ftemp=%lf, fpd=%lf, dpudpd=%lf\n",ftemp,fpd,dpudpd);
}

        break;
    } else if (flagm && flagp) {
        fs[lnew][unew] += ftemp * (muplus-muminus)
                                / mustep;
if (lnew>=12 && lnew<=14) printf("DU3 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    } else if (flagm && !flagp) {
        fs[lnew][unew] += ftemp * (mu+mustep/2.0-muminus)
                                / mustep;

if (lnew>=12 && lnew<=14) {
    printf("DU4 fs[%ld][%d] = %lf\n",lnew,unew,fs[lnew][unew]);
    printf("\t muminus=%lf, muplus=%lf\n",muminus,muplus);
}
    }
}

/* end */

/* Finally, if particles stay on the same side of the shock,
and stay within the simulation region, move to new location.
*/

```

```

    } else if (lnew <= nz[w]) {
        fs[lnew][u] += f[w][l][u];
    }
}
}
}

for (u=1;u<=nmu;u++) {
printf("\t\tu=%d\n",u);
for (l=1;l<=nz[w];l++) {
    f[w][l][u] = fs[l][u];
if (l >= 12 && l <= 14) printf("\t\tf[%d][%d][%d]=%lf\n",w,l,u,f[w][l][u]);
}
}
}
}

/*
compute pd and dpd/dpu when given
mu_u : cos(psi) in upstream local frame and
pu : momentum in upstream local frame
*/

double calc_pd(m,pu,mu_u,dpddpu)

```

```

double m, pu, mu_u, *dpddpu;
{
    double Eu, difEu, puper, pupar, pusper, puspar, difpusper, difpuspar;
    double gswu, pus, difpus, mu_us, difmu_us, pds, difpds, mu_ds, difmu_ds;
    double Eds, difEds, pd, difpd, pdpar, pdper, difpdpar, difpdper, gswd;

    /* now we are in upstream local frame
    */

    Eu = sqrt( pu*pu+m*m );
    difEu = pu/Eu;
    puper = pu*sqrt( 1.0-mu_u*mu_u );
    pupar = pu*mu_u;
    gswu = 1.0/sqrt( 1.0-BETASWU*BETASWU );

    /* Transforming from upstream local frame
    to upstream dHT frame ( observer is in
    upstream dHT frame.)
    */

    pusper = puper;
    difpusper = sqrt( 1.0-mu_u*mu_u );
    puspar = gswu*( pupar-BETASWU*Eu );
    difpuspar = gswu*( mu_u-BETASWU*difEu );
    pus = sqrt( puspar*puspar+pusper*pusper );
    difpus = (pusper*difpusper+puspar*difpuspar)/pus;
    mu_us = puspar/pus;

```

```

difmu_us = (pus*difpuspar-puspar*difpus)/(pus*pus);

/* Transforming from upstream dHT frame to downstream
   dHT frame
*/

pds = pus;
difpds = difpus;
mu_ds = -sqrt( 1.0 - ( 1.0-mu_us*mu_us )*BDOVERBU );
difmu_ds = ( mu_us/mu_ds )*BDOVERBU*difmu_us;
Eds = sqrt( pds*pds + m*m );
difEds = (pds/Eds)*difpds;

/* Transforming from downstream dHT frame to downstream
   local frame
*/

gswd = 1.0/sqrt( 1.0-BETASWD*BETASWD );
pdpar = gswd*( mu_ds*pds+BETASWD*Eds );
difpdpar = gswd*( mu_ds*difpds + pds*difmu_ds + BETASWD*difEds );
pdper = pds*sqrt( 1.0-mu_ds*mu_ds );
difpdper = (-mu_ds*pds*pds*difmu_ds + (1.0-mu_ds*mu_ds)*pds*difpds)/pdper;
pd = sqrt( pdpar*pdpar + pdper*pdper);
difpd = (pdpar*difpdpar + pdper*difpdper)/pd;
*dpddpu = difpd;

return pd;
}

```



```

/*
compute pu: momentum in upstream local frame when given
pd: momentum in downstream local frame
mu_d: mu in downstream local frame
*/

double calc_pu( m, pd, mu_d, dpddpu )
double m, pd, mu_d, *dpddpu;
{
double Ed, difEd, pdpar, difpdpar, pdper, difpdper, pds, difpds, pdspar;
double difpdspar, pdsper, difpdsper, gswd, mu_ds, difmu_ds, mu_us, difmu_us;
double pus, difpus, pusper, difpusper, Eus, difEus, gswu, pu, difpu;
double pupar, difpupar, puper, difpuper;

printf("\n BETASWD should be less than %lf",mu_d*pd/sqrt(pd*pd+m*m));

/*****/

/* in downstream local frame
*/

Ed = sqrt( pd*pd + m*m );

```

```

difEd = pd/Ed;
pdpar = mu_d*pd;
difpdpar = mu_d;
pdper = pd*sqrt( 1.0-mu_d*mu_d );
difpdper = sqrt( 1.0-mu_d*mu_d );

/* Transforming from downstream local frame to
   downstream dHT frame
*/

gswd = 1.0/sqrt( 1.0-BETASWD*BETASWD );
pdspar = gswd*( pdpar-BETASWD*Ed );
difpdspar = gswd*( difpdpar-BETASWD*difEd );
pdsper = pdper;
difpdsper = difpdper;
pds = sqrt( pdspar*pdspar + pdsper*pdsper );
difpds = (pdsper*difpdsper + pdspar*difpdspar)/pds;
mu_ds = pdspar/pds;
difmu_ds = (pds*difpdspar-pdspar*difpds)/(pds*pds);

/* It is possible for a particle to recede
   from the shock.
*/
if(mu_ds <= 0.0)
    printf("\n mu_d = %lf a particle recede from the shock\n",mu_d);

/* Transforming from downstream dHT frame to

```

```

    upstream dHT frame
*/

mu_us = sqrt( 1.0 - (1.0-mu_ds*mu_ds)/BDOVERBU );
difmu_us = (mu_ds/mu_us)*(1.0/BDOVERBU)*difmu_ds;

pus = pds;
difpus = difpds;

Eus = sqrt( pus*pus + m*m );
difEus = (pus/Eus)*difpus;

pusper = pus*sqrt( 1.0 - mu_us*mu_us );
difpusper = (-mu_us*pus*pus*difmu_us+(1.0-mu_us*mu_us)*pus*difpus)/pusper;

/* Transforming from upstream dHT frame to upstream
   local frame
*/

gswu = 1.0/sqrt( 1.0-BETASWU*BETASWU );
pupar = gswu*( mu_us*pus + BETASWU*Eus );
difpupar = gswu*( mu_us*difpus + pus*difmu_us + BETASWU*difEus );
puper = pusper;
difpuper = difpusper;

pu = sqrt( pupar*pupar + puper*puper );
difpu = (pupar*difpupar + puper*difpuper)/pu;
*dpddpu = 1.0/difpu;

return pu;
}

```

```

/*
    check whether a particle in downstream recedes from
    the shock?
*/

int recede( m, pd, mu_d )
double m, pd, mu_d;
{
    double Ed, pdper, pdpar, pdspar, pdsper, pds, mu_ds, gswd;

    /* in downstream local frame
       */

    Ed = sqrt( pd*pd + m*m );
    pdpar = mu_d*pd;
    pdper = pd*sqrt( 1.0 - mu_d*mu_d );

    /* Transforming from downstream local frame to
       downstream dHT frame
       */
    gswd = 1.0/sqrt( 1.0-BETASWD*BETASWD );
    pdspar = gswd*( pdpar - BETASWD*Ed );
    pdsper = pdper;
    pds = sqrt( pdspar*pdspar + pdsper*pdsper );
    mu_ds = pdspar/pds;

```

```

    if ( mu_ds <= 0.0 ) return 1; /* it recedes from the shock */
    if ( mu_ds > 0.0 ) return 0;
}

```

```

/* determine mu_d from pu, mu_u and when

```

- (a) $|\mu_u| > 1 + \text{TINY}$ report error
- (b) $1 - \text{TINY} < \mu_u < 1 + \text{TINY}$ return +1
- (c) $-1 - \text{TINY} < \mu_u < -1 + \text{TINY}$ return -1.

```

*/

```

```

double mu_d( m, pu, mu_u )

```

```

double m, pu, mu_u;

```

```

{

```

```

    double m2, pd, mu_d;

```

```

    double Eu, p_us_par, p_us_per, p_us, mu_us, gamma_us;

```

```

    double mu_ds, p_ds, p_ds_par, p_ds_per, Eds, gamma_ds;

```

```

    m2 = m*m;

```

```

    if( fabs(mu_u) > 1.0+TINY )

```

```

        nrerror("error in routine mu_d(): mu_u out of bounds");

```

```

    else if( 1.0-TINY < mu_u && mu_u < 1.0+TINY )

```

```

        return 1.0;

```

```

else if( -1.0-TINY < mu_u && mu_u < -1.0+TINY )
    return -1.0;

else {

/* energy in upstream local frame
*/

    printf("#19 = %lf\n", pu*pu+m2);
    Eu = sqrt( pu*pu + m2 );

/* transforming from upstream local frame to
upstream dHT frame
*/

    printf("#20 = %lf\n", 1.0-BETASWU*BETASWU);
    gamma_us = 1.0 / sqrt( 1.0 - BETASWU*BETASWU );
    printf("#21 = %lf\n", pu*pu*(1.0-mu_u*mu_u));
    p_us_per = sqrt( pu*pu*( 1.0-mu_u*mu_u ) );
    p_us_par = gamma_us*( pu*mu_u-BETASWU*Eu );
    printf("#22 = %lf\n", p_us_per*p_us_per+p_us_par*p_us_par);
    p_us      = sqrt( p_us_per*p_us_per + p_us_par*p_us_par );
    mu_us     = p_us_par / p_us;

/* transforming from upstream dHT frame
to downstream dHT frame

```

```

*/

printf("#23 = %lf\n",1.0-BETASWD*BETASWD);
gamma_ds = 1.0/sqrt(1.0-BETASWD*BETASWD);
printf("#24 = %lf\n",1.0-(1.0-mu_us*mu_us)*BDOVERBU);
printf("mu_us = %lf\n",mu_us);
mu_ds    = sqrt( 1.0 - ( 1.0 - mu_us*mu_us ) * BDOVERBU );
p_ds     = p_us;
printf("#25 = %lf\n",p_ds*p_ds+m2);
Eds      = sqrt( p_ds*p_ds+m2 );
printf("#26 = %lf\n",p_ds*p_ds*(1.0-mu_ds*mu_ds));
p_ds_per = sqrt( p_ds*p_ds*(1.0-mu_ds*mu_ds));
p_ds_par = gamma_ds*( p_ds*mu_ds+BETASWD*Eds );

/* transforming from downstream dHT frame to
   downstream local frame
*/
printf("#27 = %lf\n",p_ds_per*p_ds_per+p_ds_par*p_ds_par);
pd      = sqrt( p_ds_per*p_ds_per + p_ds_par*p_ds_par );
mu_d    = p_ds_par / pd;

return mu_d;
}
}

/* determine mu_u from pd, mu_d and when

```

- (a) $|\mu_d| > 1+TINY$ report error
- (b) $1-TINY < \mu_d < 1+TINY$ return +1
- (c) $-1-TINY < \mu_d < -1+TINY$ return -1.

*/

```
double mu_u( m, pd, mu_d )
double m, pd, mu_d;
{
    double m2, pu, mu_u;
    double Ed, p_ds_par, p_ds_per, p_ds, mu_ds, gamma_ds;
    double mu_us, p_us, p_us_par, p_us_per, Eus, gamma_us;

    m2 = m*m;

    if( fabs(mu_d) > 1.0+TINY )
        nrerror("error in routine mu_u():mu_d out of bounds");

    else if( 1.0-TINY < mu_d && mu_d < 1.0+TINY )
        return 1.0;

    else if( -1.0-TINY < mu_d && mu_d < -1.0+TINY )
        return -1.0;

    else {
```



```

/* energy in downstream local frame
*/

printf("#28 = %lf\n",pd*pd+m2);

Ed = sqrt( pd*pd + m2 );

/* transforming from downstream local frame to
downstream dHT frame
*/

printf("#29 = %lf\n",1.0-BETASWD*BETASWD);
gamma_ds = 1.0 / sqrt( 1.0 - BETASWD*BETASWD );
printf("#30 = %lf\n",pd*pd*(1.0-mu_d*mu_d));
p_ds_per = sqrt( pd*pd*( 1.0-mu_d*mu_d ));
p_ds_par = gamma_ds*( pd*mu_d-BETASWD*Ed );
printf("#31 = %lf\n",p_ds_per*p_ds_per+p_ds_par*p_ds_par);
p_ds      = sqrt( p_ds_per*p_ds_per + p_ds_par*p_ds_par );
mu_ds     = p_ds_par / p_ds;

/* transforming from downstream dHT frame
to upstream dHT frame
*/

printf("#32 = %lf\n",1.0-BETASWU*BETASWU);
gamma_us = 1.0 / sqrt(1.0-BETASWU*BETASWU);
printf("#33 = %lf\n",1.0-(1.0-mu_ds*mu_ds)/BDOVERBU);
mu_us     = -sqrt( 1.0 - ( 1.0 - mu_ds*mu_ds ) / BDOVERBU );
p_us      = p_ds;

```

```

printf("#34 = %lf\n",p_us*p_us+m2);

Eus      = sqrt( p_us*p_us+m2 );

printf("#35 = %lf\n",p_us*p_us*(1.0-mu_us*mu_us));

p_us_per = sqrt( p_us*p_us*(1.0-mu_us*mu_us));

p_us_par = gamma_us*( p_us*mu_us+BETASWU*Eus );

/* transforming from upstream dHT frame to
   upstream local frame
*/

printf("#36 = %lf\n",p_us_per*p_us_per+p_us_par*p_us_par);

pu      = sqrt( p_us_per*p_us_per + p_us_par*p_us_par );

mu_u    = p_us_par / pu;

return mu_u;
}
}

double mustar1(m,pu,mustard,mustarr)
double m, pu, *mustard, *mustarr;
{
double newsqrt();

double mu_us,mu_u;

double Eu, gswu,gswd,pus,pdpar,pupar;

double a,b,c;

```

```

/* This routine considers mu_ds = 0. */

gswu = 1.0 / newsqrt( 1.0 - BETASWU*BETASWU );
Eu = newsqrt( pu*pu + m*m );
mu_us = -newsqrt(1.0-1.0/BDOVERBU);

a= ( 1.0-mu_us*mu_us ) * gswu*gswu * pu*pu + mu_us*mu_us*pu*pu;
b= -2.0 * pu * BETASWU * Eu * gswu*gswu * ( 1.0-mu_us*mu_us );
c= BETASWU*BETASWU * Eu*Eu * gswu*gswu * ( 1.0-mu_us*mu_us )
  - mu_us*mu_us * pu*pu;
mu_u = ( -b - newsqrt( b*b - 4.0*a*c )) / ( 2.0*a );
*mustarr = ( -b + newsqrt( b*b - 4.0*a*c)) / ( 2.0*a );

/* compute mustard
*/

gswd = 1.0 / newsqrt( 1.0 - BETASWD*BETASWD );
pus = newsqrt(gswu*( mu_u*pu-BETASWU*Eu )*gswu*( mu_u*pu-BETASWU*Eu )
  + pu*pu*(1.0-mu_u*mu_u));
pdpar = gswd*BETASWD*newsqrt( m*m + pus*pus );
*mustard = pdpar / newsqrt( pus*pus+pdpar*pdpar );
return mu_u;
}

/*

```

```

compute pr: momentum reflecting at the shock front back to upstream
        local frame when given
        pu: momentum incident at the shock front in upstream local
            frame
        mu_u: cos(phi) in upstream local frame
*/

double refl_pr( m, pu, mu_u, dpudpr )
double m, pu, mu_u, *dpudpr;
{
    double Eu, difEu, pus, difpus, puspar, difpuspar, pusper, difpusper;
    double gswu, prs, difprs, prspar, difprspar, prsper, difprsper;
    double mu_rs, difmu_rs, Ers, difErs, pr, difpr, prpar, difprpar;
    double prper, difprper, mu_us;

/* in upstream local frame
*/
    Eu = sqrt( pu*pu+m*m );
    difEu = pu/Eu;

/* Transforming from upstream local frame to
upstream dHT frame.
*/

    gswu = 1.0/sqrt( 1.0 - BETASWU*BETASWU );
    puspar = gswu*( mu_u*pu - BETASWU*Eu );
    difpuspar = gswu*( mu_u - BETASWU*difEu );

```

```

pusper = pu*sqrt( 1.0 - mu_u*mu_u );
difpusper = sqrt( 1.0 - mu_u*mu_u );
mu_us = puspar/sqrt( puspar*puspar + pusper*pusper );
if( mu_us >= 0.0 )
    printf("\n in dHT frame mu_us = %lf",mu_us);

/* Reflecting at the shock front in dHT frame back to
upstream
*/

prspar = -puspar;
difprspar = -difpuspar;
prasper = pusper;
difprasper = difpusper;
prs = sqrt( prspar*prspar + prasper*prasper );
difprs = ( prspar*difprspar + prasper*difprasper ) / prs;
mu_rs = prspar/prs;
difmu_rs = ( prs*difprspar - prspar*difprs ) / ( prs*prs );
Ers = sqrt( prs*prs + m*m );
difErs = ( prs/Ers )*difprs;

/* Transforming back from dHT frame to upstream local frame
*/

prpar = gswu*( prspar + BETASWU*Ers );
difprpar = gswu*( mu_rs*difprs + prs*difmu_rs + BETASWU*difErs );
prper = prasper;

```

```

    difprper = difprsper;
    pr = sqrt( prpar*prpar + prper*prper );
    difpr = ( prpar*difprpar + prper*difprper )/pr;
    *dpudpr = 1.0/difpr;
    return pr;
}

/* compute mu after reflecting at the shock front from
upstream local frame back to upstream local frame

NOTE: momentum magnitude conservs in dHT frame
pu -> momentum in upstream local frame
mu -> mu in upstream local frame
*/

double mu_refl( m,pu,mu )
double m,pu,mu;
{
    double pus,pus_par,pus_per,Eu,Eus,gswu,pr;
    double pus_refl_par,pus_refl_per,pu_refl_par,pu_refl_per;

/* in upstream local frame we compute energy
and gamma of particle
*/

```

```

Eu = sqrt( m*m + pu*pu );
gswu = 1.0 / sqrt( 1.0-BETASWU*BETASWU );

/* transforming from upstream local frame to
   dHT frame
*/
pus_par = gswu*( mu*pu - BETASWU*Eu );
pus_per = pu*sqrt( 1.0-mu*mu );

/* reflection in dHT frame conservs energy, and
   momentum change its direction but its magnitude
   does not.
*/

pus_refl_par = -pus_par;
pus_refl_per = pus_per;
Eus = sqrt( pus_refl_par*pus_refl_par+pus_refl_per*pus_refl_per+m*m );

/* transforming back from dHT frame to upstream
   local frame
*/
pu_refl_par = gswu*( pus_refl_par + BETASWU*Eus );
pu_refl_per = pus_refl_per;
pr = sqrt( pu_refl_par*pu_refl_par + pu_refl_per*pu_refl_per );

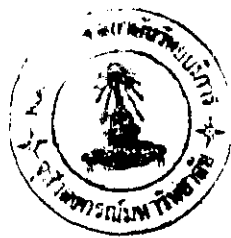
return pu_refl_par/pr;

```

```
}  
  
/* compute square root of x  
   if |x| < TINY  
   it return 0  
*/  
  
double newsqrt( double x )  
{  
    if( x < -TINY ) nrerror("newsqrt error");  
    if( fabs(x) < TINY ) return 0.0;  
    else return sqrt(x);  
}
```



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Curriculum Vitae

Mr. Songklod Riyavong was born on January 2, 1971 in Udonthani. He received his B.Sc. degree in physics from Khon Kaen University in 1992. He received support from the Institute for the Promotion of Teaching Science and Technology for both his undergraduate and graduate studies.



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