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//===== print solution profile to text file at
certain times
#include "global.h"
#include <stdio.h>
void output(const int W, double X[], double Y[], double T[][W+2],
double Fx[][W+1], double Fy[][W+1], double E[][W+1], double p[][W+1],
double time, int nsteps, double tend){
//      double flux, energy, phase;
//      FILE *OUT;
//      FILE *TEMP;
//      FILE *PHASE;
//      FILE *ENTH; // initialize file var
if(!nsteps){ // if haven't begun time stepping
    OUT = fopen("outputs/values.o", "w"); // new file
    fprintf(OUT, "#nstep time (ms)  width (cm) depth(cm)
energy (J)\n");
    TEMP = fopen("outputs/temp.o", "w"); // new file
    //fprintf(TEMP, "#Temperatures by timesteps\n");
    PHASE = fopen("outputs/phase.o", "w"); // new file
    //fprintf(PHASE, "#Liquid fraction by timestep\n");
    ENTH = fopen("outputs/enth.o", "w"); // new file
    //fprintf(ENTH, "#Enthalpies by timestep\n");
}
else{
    OUT = fopen("outputs/values.o", "a"); // old file
    TEMP = fopen("outputs/temp.o", "w"); // old file
    PHASE = fopen("outputs/phase.o", "w"); // old file
    ENTH = fopen("outputs/enth.o", "w"); // old file
}
// hidden time, steps, error
// position and profile in columns
//=====
TEMPERATURES
// inside temperatures only, walls not necessary
for(int j = 1; j <= M; j++){
    for(int i = 1; i <= M; i++){
        fprintf(TEMP, "%22.15e ", T[i][j]);
    }
    fprintf(TEMP, "\n");
}
fprintf(TEMP, "\n");
fclose(TEMP);
//===== PHASES
for(int j = 1; j <= M; j++){
    for(int i = 1; i <= M; i++){
        fprintf(PHASE, "%22.15e ", p[i][j]);
    }
    fprintf(PHASE, "\n");
}
fprintf(PHASE, "\n");

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fclose(PHASE);
//===== ENTHALPIES
// just for funsies
for(int j = 1; j <= M; j++){
    for(int i = 1; i <= M; i++){
        fprintf(ENTH, "%22.15e ", E[i][j]);
    }
    fprintf(ENTH, "\n");
}
fprintf(ENTH, "\n");
fclose(ENTH);
//=====
TIMESTAMPS/ETC
double xl, xr, yd, width, depth, energy;
int il = M/2, ir = M/2, id = M;
for(int l = M/2; l <= M; l++){
    for(int k = 1; k <= M; k++){ // bounds for width of
pool
        if(p[k-1][l] < 1 && p[k+1][l] == 1){
            if(k < il){
                il = k;}
        }
        if(p[k-1][l] == 1 && p[k+1][l] < 1){
            if(k > ir){
                ir = k;}
        }
    }
}
for(int k = 1; k <= M; k++){
    if(p[M/2][k-1] < 1 && p[M/2][k+1] == 1){
        id = k;}
}
xl = a + ((double)il - 0.5)*dx;
xr = a + ((double)ir - 0.5)*dx;
yd = a + ((double)id - 0.5)*dx;
width = 100*(xr - xl); // in cm
depth = 100*(b - yd); // in cm
energy = Q0*time; // in Joules
fprintf(OUT, "%i %6.4f %6.4f %6.4f %6.4f\n", nsteps,
100*time, width, depth, energy);
//fprintf(OUT, "# Error up to this time: %.15e\n\n", ERR);
//
// COMPLETION BAR
if(nsteps == 0){
    printf("0%%      20%%      40%%      60%%
80%%      100%%\n");
    printf(" ");
}
if( time > nbar*tend/50 ){
    //printf("%3.0f%%\n", 100*time/tend);

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        printf("=");
        fflush(stdout);
        nbar = nbar + 1;
    }
    if(width > 2){
        maxwidth = 1;
        int barleft = 50 - nbar;
        for(int g = 0; g <= barleft; g++){
            printf("=");
        }

        /*for(int j = M+1; j >= 0; j--){
            for(int i = 0; i <= M+1; i++){
                fprintf(OUT, "%22.15e ", T[i][j]);
            }
            fprintf(OUT, "\n");
        }*/

        /*for(int i = 0; i <= M+1; i++){
            if(i == 0){
                fprintf(OUT, "%18.15e %18.15e %18.15e\n",
X[i], T[i], F[i]);
            }
            if(i > 0 && i <= M){
                fprintf(OUT, "%18.15e %18.15e %18.15e %18.15e
%.3f\n", X[i], T[i], F[i], E[i], p[i]);
            }else if(i > M){
                fprintf(OUT, "%.15e %.15e\n", X[i], T[i]);
            }
        }*/
        //fprintf(OUT, "\n"); // new line to separate times of
printing
        fclose(OUT);
    }

```