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typedef std::vector<double> VD;

void model(double t, double PS, double PR, double STARCH, double NIT, double RESPIRATION, VD&
params, VD& states) {

    /*
    PS is  $\mu\text{mol C6} / (\text{g}\cdot\text{h})$ 
    PR is  $\mu\text{mol C2} / (\text{g}\cdot\text{h})$ 
    RESPIRATION is  $\mu\text{mol C6} / (\text{g}\cdot\text{h})$ 
    */

    // Extract parameters

    // HPR
    double vmax_hpr = params[0];
    double km_hpr = params[1];

    // NR
    double vmax_nr = params[2];
    double km_nr = params[3];

    // GS
    double vmax_gs = params[4];
    double km_gs_GLU = params[5];
    double km_gs_AMMONIUM = params[6];

    // GOGAT
    double vmax_gogat = params[7];
    double km_gogat_GLN = params[8];
    double km_gogat_KG = params[9];

    // remaining parameters
    double alpha = params[10];
    double hp_to_bmexp = params[11];
    double hp_to_mf = params[12];
    double mf_to_cit = params[13];
    double cit_to_mf = params[14];
    double cit_to_akg = params[15];
    double aa_to_bmexp = params[16];
    double ser_to_aa = params[17];
    double correct = params[18];
    double gdc_vmax = params[19];
    double shmt_vmax = params[20];
    double km_gly_gdc = params[21];
    double km_gly_shmt = params[22];
    double gs_inactivation = params[23];

    // Extracting states
    double GLY = states[0];
    double SER = states[1];
    double AMMONIUM = states[2];
    double AA = states[3];
    double GLU = states[4];
    double GLN = states[5];
    double HP = states[6];
    double MF = states[7];
    double CIT = states[8];
    double KG = states[9];

    // only 5.0425% Cytosol Koffler2013
    NIT = NIT*0.050425;

    // Enzyme kinetics
    double HPR = (vmax_hpr*SER) / (km_hpr + SER);
    double NR = (vmax_nr*NIT) / (km_nr + NIT);

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double GOGAT = (vmax_gogat*GLN*KG) / (km_gogat_KG*km_gogat_GLN + km_gogat_KG*GLN +
km_gogat_GLN*KG + GLN*KG);
double GS = (vmax_gs*GLU*AMMONIUM) / (km_gs_GLU*km_gs_AMMONIUM + km_gs_GLU*AMMONIUM +
km_gs_AMMONIUM*GLU + GLU*AMMONIUM);

double GDC = (gdc_vmax*GLY) / (km_gly_gdc + GLY);

//shmt_vmax = gdc_vmax;
//km_gly_shmt = km_gly_gdc;
double SHMT = (shmt_vmax*GLY) / (km_gly_shmt + GLY);

if(SHMT > GDC) {
    SHMT = GDC;
}

// factors:
// c2 flux to c6 flux
double factor0 = 2./6.;
// AA C6 N1.23
// ca. 14.3197% of proteom consists of Ser
double factor1 = 0.143197;
// C6 to C5
double factor2 = 6./5.;

// nightly flux in gly & maximal HPR rate
double PR_new = PR;
if(t > 8.) {
    // PR_new = PR + alpha;
}
PR_new = PR + alpha;
if(HPR > (PR_new/2.)) {
    HPR = PR_new/2.;
}

/*
respiration is reduced by photorespiration
But only if GDC*GLY is smaller than respiration
*/
double RESPIRATION_new = RESPIRATION + (GDC)*factor0;
if(RESPIRATION_new > 0.) {
    RESPIRATION_new = 0.;
}

if( (t < 8.) && (RESPIRATION_new < 0.) ) {
    PS = PS - RESPIRATION_new;
}

/*
regulation of gs and nr during night
flux from ser to aa only during night
*/
double ser_to_aa_new = 0.;
if(t > 8.) {
    ser_to_aa_new = ser_to_aa;
    GS = GS*gs_inactivation;
}

double GLU_TO_AA = NR*correct;

/*
N balance in order to calculate flux from Glu to aKG
*/
double GLU_GLYOXYLATE_AT = PR_new - HPR;
if(GLU_GLYOXYLATE_AT < 0.) {
    GLU_GLYOXYLATE_AT = 0.;
}

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}

// ODE System
double ddtGLY = states[0] = PR_new - GDC - SHMT;
double ddtSER = states[1] = (1./3.)*GDC + (2./3.)*SHMT - HPR - SER*ser_to_aa_new -
AA*aa_to_bmexp*factor1;
double ddtAMMONIUM = states[2] = GDC + NR - GS;
double ddtAA = states[3] = GLU_TO_AA - AA*aa_to_bmexp + (SER*ser_to_aa_new)/2.;
double ddtGLU = states[4] = 2.*GOGAT - GS - GLU_TO_AA - GLU_GLYOXYLATE_AT;
double ddtGLN = states[5] = GS - GOGAT;
double ddtHP = states[6] = PS + HPR/2. - HP*hp_to_bmexp - HP*hp_to_mf - STARCH;
double ddtMF = states[7] = HP*hp_to_mf + RESPIRATION_new - MF*mf_to_cit + CIT*cit_to_mf;
double ddtCIT = states[8] = MF*mf_to_cit - CIT*cit_to_mf - CIT*cit_to_akg;
double ddtKG = states[9] = factor2*CIT*cit_to_akg + GLU_GLYOXYLATE_AT - GOGAT;
double ddtbilanz = states[10] = PR_new/3 + PS + RESPIRATION_new -HP*hp_to_bmexp -
AA*aa_to_bmexp*(6./5.);
}

```