

```

/*****
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*
* This file contains C code for an MCMC algorithm constructed
* to fit a hierarchical model that incorporates the idea of
* temporally dependent partitions.
*
* I will include model details at a later date
*
*****/

#include "Rutil.h"

#include <R_ext/Lapack.h>
#include <R.h>
#include <Rmath.h>

#include <stdio.h>
#include <stdlib.h>
#include <time.h>

/*****
* The following are the inputs of the function that are read from R
*
* draws = total number of MCMC draws
* burn = number of MCMC draws discarded as burn-in
* thin = indicates how much MCMC chain should be thinned
*
* nsubject = integer for the number of subjects/units in data set
* ntime = integer for the number of time points
* y = double nsubject x ntime matrix containing response for each subject at time t
* s1 = nsubject x 1 vector containing spatial coordinate one
* s2 = nsubject x 1 vector containing spatial coordinate two
* slp = nsubject x 1 vector containing spatial coordinate one for prediction (currently not used)
* s2p = nsubject x 1 vector containing spatial coordinate two for prediction (currently not used)
*
* M = double - indicating value of M associated with cohesion (scale parameter of DP).
* alpha = double - prior probability of being pegged, starting value only if is TRUE
* modelPrior = vector - containing values for prior distributions as follows
*     m0 - mean phi0, s20 - variance of phi0
*     A - upper bound of sigma*_{jt}
*     A0 - upper bound of tau
*     A1 - upper bound of lam
*     a - alpha_t shape 1 parameter, b - alpha_t shape 2 parameter
*     be - scale parameter of eta.
*
* global_alpha = integer - logical indicating whether to make alpha time-specific or one global alpha.
* alpha_0 = integer - logical indicating whether alpha = 0 or not.
* etal_0 = integer - logical indicating whether etal = 0 or not.
* phil_0 = integer - logical indicating whether phil = 0 or not.
* sPPM = integer - logical indicating whether to use spatial information or not
*
* SpatialCohesion = integer indication which cohesion to use
*     1 -Auxiliary
*     2- Double dipper
*
* cParms - vector holding values employed in the cohesion
* mh -
* verbose - logical indicating if information should be printed to screen
*
* OUTPUT
* Si -
* mu -
* sig2 -
* etal
* theta -
* tau2 -
* phi0 -
* phil -
* lam2 -
* gamma -
* alpha.out -
* like
* lpml -
* waic -
*****/

void mcmc_drpm_ar1(int *draws, int *burn, int *thin, int *nsubject, int *ntime,
                  double *y, double *s1, double *s2, double *M,
                  double *modelPriors,
                  int *global_alpha, int *alpha_0, int *etal_0, int *phil_0,
                  int *sPPM, int *SpatialCohesion, double *cParms, double *mh, int *verbose,
                  int *Si, double *mu, double *sig2, double *etal, double *theta, double *tau2,

```

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double *phi0, double *phil, double *lam2, int *gamma, double *alpha_out,
double *fitted, double *llike, double *lpml, double *waic){

// i - MCMC iterate
// ii - MCMC iterate that is saved
// j - subject iterate
// jj - second subject iterate
// t - time iterate
// k - cluster iterate
// p - prediction iterate

int i, ii, j, jj, t, k;
ii = 0;

int nout = (*draws - *burn)/(*thin);

if(*verbose){
    Rprintf("nsubject = %d\n", *nsubject);
    Rprintf("ntime = %d\n", *ntime);
    Rprintf("nout = %d\n", nout);
}

// =====
//
// Memory vectors to hold MCMC iterates for non cluster specific parameters
//
// =====

// This variable is used to create a "buffer" zone of memory so that updating
// things on time boundary do not need special attention in the algorithm since
// I have to look at time period before and after when updating partitions
int ntime1 = *ntime + 1;

// I am adding one more year as an empty vector
// so that the C program does not crash.
int gamma_iter[(*nsubject)*(ntime1)];
int Si_iter[(*nsubject)*(ntime1)];
int nclus_iter[ntime1];

double *etal_iter = R_VectorInit(*nsubject, 0.0);
double *theta_iter = R_VectorInit(ntime1, 0.0);
double *tau2_iter = R_VectorInit(ntime1, 1.0);

double phi0_iter = 0.0;
double phil_iter = 0.0;
double lam2_iter = 1.0;

double *alpha_iter = R_VectorInit(ntime1, 0.0);

// =====
//
// Memory vectors to hold MCMC iterates for cluster specific parameters
//
// =====

double *muh = R_VectorInit((*nsubject)*(ntime1), 0.0);
double *sig2h = R_VectorInit((*nsubject)*(ntime1), 0.5);

int nh[(*nsubject)*(ntime1)];

// =====
//
// Initialize a few parameter vectors
//
// =====

// Initialize Si according to covariates
// I am adding one time period here to have
// scratch memory (never filled in) so that
// I can avoid dealing with boundaries in algorithm
for(j = 0; j < *nsubject; j++){
    for(t = 0; t < ntime1; t++){ // Note I am not initializing the "added time memory"
        Si_iter[j*(ntime1) + t] = 1;
        gamma_iter[j*(ntime1) + t] = 0;
        nh[j*(ntime1) + t] = 0;
        if(t==1) Si_iter[j*ntime1 + t] = 1;
        if(t==*ntime) Si_iter[j*(ntime1) + t] = 0;
    }
}
}

```

```

1;

// Initial enumeration of number of subjects per cluster;
for(j = 0; j < *nsubject; j++){
    for(t = 0; t < *ntime; t++){
        nh[(Si_iter[j*(ntime1)+t]-1)*(ntime1) + t] = nh[(Si_iter[j*(ntime1)+t]-1)*(ntime1) + t] +
    }
}

// Initialize the number of clusters
for(t = 0; t < *ntime; t++){
    nclus_iter[t] = 0;
    for(j = 0; j < *nsubject; j++){
        if(nh[j*(ntime1) + t] > 0) nclus_iter[t] = nclus_iter[t] + 1;
    }
}
nclus_iter[*ntime] = 0;

// =====
//
// scratch vectors of memory needed to update parameters
//
// =====

// stuff needed to update gamma vectors
int nclus_red=0, nh_red[*nsubject], n_red=0, gt;
int nclus_redtmp=0, nh_redtmp[*nsubject], n_redtmp=0, nh_tmp[*nsubject];
int nh_redtmp_no_zero[*nsubject], nh_red_no_zero[*nsubject],nh_tmp_no_zero[*nsubject];
double lpp_full=0.0, lpp_red=0.0;

int nh_red_1[*nsubject];
int nclus_redtmp_1=0, nh_redtmp_1[*nsubject], n_redtmp_1=0, nh_tmp_1[*nsubject];
int nh_redtmp_no_zero_1[*nsubject], nh_red_no_zero_1[*nsubject],nh_tmp_no_zero_1[*nsubject];
double lpp_full_1=0.0, lpp_red_1=0.0;

for(j=0; j<*nsubject; j++){
    nh_tmp[j] = 0; nh_red[j] = 0; nh_redtmp[j] = 0;
    nh_redtmp_no_zero[j] = 0; nh_red_no_zero[j] = 0; nh_tmp_no_zero[j] = 0;

    nh_tmp_1[j] = 0; nh_red_1[j] = 0; nh_redtmp_1[j] = 0;
    nh_redtmp_no_zero_1[j] = 0; nh_red_no_zero_1[j] = 0; nh_tmp_no_zero_1[j] = 0;
}

// stuff that I need to update Si (the partition);
int compt1[(*nsubject)],comptml[(*nsubject)],comp2t[(*nsubject)],comptpl[(*nsubject)];
int rho_tmp[*nsubject], Si_tmp[*nsubject], Si_tmp2[*nsubject];
int oldLab[*nsubject], reorder[*nsubject];
intiaux, Rindx1, Rindx2, n_tmp, nclus_tmp, n_tmp_1, nclus_tmp_1, rho_comp, indx;
doubleauxm, auxs, mudraw, sigdraw, maxph, denph, cprobh, uu, lCo, lCn, lCn_1;
double *ph = R_VectorInit(*nsubject, 0.0);
double *phtmp = R_VectorInit(*nsubject, 0.0);
double *probh = R_VectorInit(*nsubject, 0.0);
double *slo = R_Vector(*nsubject);
double *s2o = R_Vector(*nsubject);
double *sln = R_Vector(*nsubject);
double *s2n = R_Vector(*nsubject);

for(j=0; j<(*nsubject); j++){
    compt1[j] = 0; comptml[j] = 0, comp2t[j]=0, comptpl[j]=0;
}

// stuff I need to update etal
double elo, eln, logito, logitn, one_phisq;

// stuff I need to update muh and sig2h
double mstar, s2star, sumy, sume2;
double nsig, osig, llo, llm, llr;
double *mu_tmp = R_VectorInit(*nsubject, 0.0);
double *sig2_tmp = R_VectorInit(*nsubject, 1.0);

// stuff that I need for theta and lam2
double summu, nt, ot, lam2tmp, phlsq, sumt, opl, npl, ol, nl;

// stuff that I need to update alpha
int sumg;
double astar, bstar,alpha_tmp;

// Stuff to compute lpml, likelihood, and WAIC
int like0, nout_0=0;
double lpml_iter, elppdWAIC;
double *CPO = R_VectorInit((*nsubject)*(ntime1), 0.0);
double *like_iter = R_VectorInit((*nsubject)*(ntime1), 0.0);
double *fitted_iter = R_VectorInit((*nsubject)*(ntime1), 0.0);

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double *mnlike = R_VectorInit((*nsubject)* (ntime1), 0.0);
double *mnllike = R_VectorInit((*nsubject)* (ntime1), 0.0);

// stuff to predict
int gpred[*nsubject], nh_pred[*nsubject], predSi_iter[*nsubject];

// =====
//
// Prior parameter values
//
// =====

// upper bound for sig, tau, lam
double A=modelPriors[2];
double A0=modelPriors[3];
double A1=modelPriors[4];

// priors for phi0
double m0 = modelPriors[0], s20 = modelPriors[1];

// priors for alpha
double a = modelPriors[5], b = modelPriors[6];

//priors for etal
double b_etal = modelPriors[7];

if(*verbose){
    Rprintf("Prior values: m0 = %.2f, s20 = %.2f\n \v Asig = %.2f, Atau = %.2f, Alam = %.2f,\n \t a = %.2f,
b = %.2f, b_etal = %.2f\n",
        m0, s20, A, A0, A1, a, b, b_etal);
}

// DP weight parameter
double Mdp = *M;

// Rprintf("Mdp = %f\n", Mdp);

// Cohesion auxiliary model paramaters for Cohesions 3 and 4
double k0=cParms[1], v0=cParms[2];
double *mu0 = R_VectorInit(2,cParms[0]);
double *L0 = R_VectorInit(2*2,0.0);
L0[0] = cParms[3]; L0[3] = cParms[3];
// Rprintf("k0 = %f\n", k0);
// RprintVecAsMat("L0", L0, 2, 2);

// RprintVecAsMat("mh", mh, 1, 5);
// M-H step tunning parameter
double csigSIG=mh[0], csigTAU=mh[1], csigLAM=mh[2], csigETA1=mh[3], csigPHI1=mh[4];

// Rprintf("csigETA1 = %f\n", csigETA1);

GetRNGstate();

// =====
//
// start of the mcmc algorithm;
//
// =====

for(i = 0; i < *draws; i++){
    if(*verbose){
        if((i+1) % 5000 == 0){
            time_t now;
            time(&now);

            Rprintf("mcmc iter = %d ===== \n", i+1);
            Rprintf("%s", ctime(&now));
        }
    }

    // Start updating gamma and partition for each time period
    for(t = 0; t < *ntime; t++){

        Rprintf("t = %d\n", t);

        //////////////////////////////////////
        //

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```

// begin by updating gamma (pegged) parameters
//
// The complete partition does not change as gammas change
for(j=0; j<*nsubject; j++){
    nh_red[j]=0; nh_redtmp[j]=0; nh_tmp[j] = 0;
}

for(k = 0; k < nclus_iter[t]; k++){
    nh_tmp[k] = nh[k*(ntime1) + t];
}

// Note this value does not change even as gamma changes. The only
// concern is that rho_t-1 and rho_t are compatible.
lpp_full = partition_prob_crp(nh_tmp, nclus_iter[t], Mdp, *nsubject, 1);

//
Rprintf("lpp_full = %f\n", lpp_full);

// find the reduced partition information
// i.e., number of units clustered, number of clusters, size of clusters;
n_red = 0; n_redtmp=0;
for(j = 0; j < *nsubject; j++){
    if(gamma_iter[j*(ntime1) + t] == 1){
        nh_red[Si_iter[j*(ntime1) + t]-1] = nh_red[Si_iter[j*(ntime1) + t]-1]+1;
        n_red=n_red+1;

        nh_redtmp[Si_iter[j*(ntime1) + t]-1] = nh_redtmp[Si_iter[j*(ntime1) + t]-1]-1;
        n_redtmp=n_redtmp+1;
    }
}

nclus_red = 0, nclus_redtmp=0;
for(j = 0; j < *nsubject; j++){
    if(nh_red[j] > 0) nclus_red = nclus_red + 1;
    if(nh_redtmp[j] > 0) nclus_redtmp = nclus_redtmp + 1;
}

//
Rprintf("nclus_red = %d\n", nclus_red);
Rprintf("n_red = %d\n \n \n \n", n_red);

remove_zero(nh_red, *nsubject, nh_red_no_zero);

//
RprintIVecAsMat("nh_red_no_zero", nh_red_no_zero, 1, *nsubject);

//
RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);
RprintIVecAsMat("Si_iter", Si_iter, *nsubject, ntime1);

for(j = 0; j < *nsubject; j++){
    Rprintf("t = %d\n", t);
    Rprintf("j = %d\n", j);

    RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);
    Rprintf("gamma_iter[j*(ntime1) + t] = %d\n", gamma_iter[j*(ntime1) + t]);

    // at time period one, all gammas are zero (none are 'pegged')
    if(t == 0){
        gamma_iter[j*(ntime1) + t] = 0;
    } else {
        // this may need to be updated depending on if the value of gamma changes

        RprintIVecAsMat("nh_red_no_zero", nh_red_no_zero, 1, *nsubject);

        lpp_red = partition_prob_crp(nh_red_no_zero, nclus_red, Mdp, n_red, 1);

        Rprintf("lpp_full = %f\n", lpp_full);
        Rprintf("lpp_red = %f\n", lpp_red);

        // If gamma is 1 at current MCMC iterate, then there are no
        // concerns about partitions being incompatible as gamma changes

```

```
// from 1 to 0.
```

```
if(gamma_iter[j*(ntime1) + t] == 1){
```

```
//      Rprintf("Starting at gamma=1 and staying at gamma=1 \n");
//      // if gamma remains 1, then no changes
//      ph[1] = lpp_full - lpp_red + log(alpha_iter[t]);
//      Rprintf("ph[1] = %f\n", ph[1]);

//      // if gamma moves from 1 to 0 need to remove one unit from
//      // rho_t.R

1      nh_redtmp[Si_iter[j*ntime1+t]-1] = nh_red[Si_iter[j*ntime1+t]-1] -
      n_redtmp = n_redtmp - 1;

      remove_zero(nh_redtmp, *nsubject, nh_redtmp_no_zero);

//      RprintfIVecAsMat("nh_redtmp", nh_redtmp, 1, *nsubject);
//      RprintfIVecAsMat("nh_redtmp_no_zero", nh_redtmp, 1, *nsubject);

      nclus_redtmp=0;
      for(jj = 0; jj < *nsubject; jj++){
          if(nh_redtmp[jj] > 0) nclus_redtmp = nclus_redtmp + 1;
      }

//      Rprintf("Starting at gamma=1 and moving to gamma=0 \n");
//      RprintfIVecAsMat("nh_red", nh_red, 1, *nsubject);
//      RprintfIVecAsMat("nh_redtmp", nh_redtmp, 1, *nsubject);

//      Rprintf("lpp_full = %f\n", lpp_full);
      lpp_red = partition_prob_crp(nh_redtmp_no_zero, nclus_redtmp, Mdp,
      n_redtmp, 1);

//      Rprintf("lpp_red = %f\n", lpp_red);

      ph[0] = lpp_full - lpp_red + log(1-alpha_iter[t]);

}
```

```
// if gamma's current value is 0, then care must be taken when
// trying to change from gamma=0 to gamma=1 as the partitions may
// no longer be compatible
```

```
if(gamma_iter[j*(ntime1) + t] == 0){
```

```
//      Rprintf("lpp_red = %f\n", lpp_red);
//      // if gamma remains zero. nothing changes to evaluate
//      Rprintf("Starting at zero and staying at zero \n");

//      Rprintf("lpp_full = %f\n", lpp_full);
//      Rprintf("lpp_red = %f\n", lpp_red);
//      Rprintf("log(1-alpha_iter[t]) = %f\n", log(1-alpha_iter[t]));

      ph[0] = lpp_full - lpp_red + log(1-alpha_iter[t]);

//      Rprintf("ph[0] = %f\n", ph[0]);
//      // to move from gamma=0 to gamma=1 need to make sure that partitio
ns      // remain compatible, if compatible .

//      RprintfIVecAsMat("Si_iter", Si_iter, *nsubject, ntime1);

//      // to move from gamma=0 to gamma=1 need to add unit to rho_t.R

//      Rprintf("Starting at zero and moving to one \n");

      nh_redtmp[Si_iter[j*ntime1 + t] - 1] = nh_red[Si_iter[j*ntime1+t]-
1] + 1;

      n_redtmp = n_red + 1;
      nclus_redtmp=0;
      for(jj = 0; jj < *nsubject; jj++){
          if(nh_redtmp[jj] > 0) nclus_redtmp = nclus_redtmp + 1;
      }

//      RprintfIVecAsMat("nh_redtmp", nh_redtmp, 1, *nsubject);
//      Rprintf("n_redtmp = %d\n", n_redtmp);
//      Rprintf("nclus_redtmp = %d\n", nclus_redtmp);
```

```

// To determine compatibility, I need to make sure that
// comparison of the reduced partitions is being made with
// correct cluster labeling. I try to do this by identifying
// the sets of units and sequentially assigning "cluster labels"
// starting with set that contains the first unit. I wonder if
// there is a way to do this in C with out using loops? Who
// can I ask about this?

// Get rho_t | gamma_t = 1 and rho_{t-1} | gamma_t = 1
// when gamma_{it} = 1;
Rindx1 = 0;
for(jj = 0; jj < *nsubject; jj++){
    if(gamma_iter[jj*ntime1 + (t)] == 1){
        comptml[Rindx1] = Si_iter[jj*ntime1 + (t-1)];
        complt[Rindx1] = Si_iter[jj*ntime1 + (t)];
        Rindx1 = Rindx1 + 1;
    }
    // I need to include this because determine what happens w

    // gamma goes from 0 to 1;
    if(jj == j){
        comptml[Rindx1] = Si_iter[jj*ntime1 + (t-1)];
        complt[Rindx1] = Si_iter[jj*ntime1 + (t)];
        Rindx1 = Rindx1 + 1;
    }
}

// Rprintf("Rindx1 = %d\n", Rindx1);
// RprintIVecAsMat("comptml", comptml, 1, *nsubject);
// RprintIVecAsMat("complt", complt, 1, *nsubject);

rho_comp = compatibility(comptml, complt, Rindx1);

// Rprintf("rho_comp = %d\n", rho_comp);
if(rho_comp==1){

    remove_zero(nh_redtmp, *nsubject, nh_redtmp_no_zero);

    RprintIVecAsMat("nh_redtmp_no_zero", nh_redtmp_no_zero, 1,
*nsubject);

    lpp_red = partition_prob_crp(nh_redtmp_no_zero, nclus_redt

    Rprintf("lpp_full = %f\n", lpp_full);
    Rprintf("lpp_red = %f\n", lpp_red);
    Rprintf("log(alpha_iter[t]) = %f\n", log(alpha_iter[t]));

    ph[1] = lpp_full - lpp_red + log(alpha_iter[t]);

} else {

    ph[1] = log(0); // partitions are not compatible

}

// Rprintf("ph[0] = %f\n", ph[0]);

// RprintVecAsMat("ph", ph, 1, 2);

}

// RprintVecAsMat("ph = ", ph, 1, 2);

maxph = ph[0]; if(maxph < ph[1]) maxph=ph[1];
Rprintf("maxph = %f\n", maxph);

ph[0] = exp(ph[0] - maxph); ph[1] = exp(ph[1] - maxph);

// RprintVecAsMat("ph = ", ph, 1, 2);

probh[1] = ph[1]/(ph[0] + ph[1]);

// RprintVecAsMat("probh = ", probh, 1, 2);
// Rprintf("probh[1] = %f\n", probh[1]);

gt = rbinom(1,probh[1]);

// Rprintf("gt = %d\n", gt);

```

```

        if(gt != gamma_iter[j*(ntime1) + t]){
            gamma_iter[j*(ntime1) + t] = gt;
            n_red = n_redtmp;
            nclus_red = nclus_redtmp;
            nh_red[Si_iter[j*(ntime1)+t]-1] = nh_redtmp[Si_iter[j*(ntime1)+t]-
1];

        } else {

            nh_redtmp[Si_iter[j*(ntime1)+t]-1] = nh_red[Si_iter[j*(ntime1)+t]-
1];

            nclus_redtmp = nclus_red;
            n_redtmp = n_red;

        }

    }

    remove_zero(nh_red, *nsubject, nh_red_no_zero);

//      Rprintf("gamma_iter[j*(ntime1) + t] = %d\n", gamma_iter[j*(ntime1) + t]);
//      Rprintf("nclus_red = %d\n", nclus_red);
//      Rprintf("n_red = %d\n", n_red);
//      RprintIVecAsMat("nh_red = ", nh_red, 1, *nsubject);
//      RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);
//      RprintIVecAsMat("Si_iter", Si_iter, *nsubject, ntime1);

    }

//      RprintIVecAsMat("Si_iter", Si_iter, *nsubject, ntime1);
//      RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);

//      Rprintf("Begin updating partition for time %d\n", t+1);

//      //////////////////////////////////////
//      // update partition
//      //////////////////////////////////////
//      // The cluster probabilities depend on four partition probabilities
//      //
//      // rho_t
//      // rho_t.R
//      // rho_t+1
//      // rho_t+1.R
//      //
//      // I have switched a number of times on which of these needs to be computed
//      // and which one can be absorbed in the normalizing constant. Right now I am
//      // leaning towards Pr(rho_t+1) and Pr(rho_t+1.R) can be absorbed. But I need
//      // to use rho_t.R and rho_t+1.R to check compatibility as I update rho_t.
//      //
//      //////////////////////////////////////

    for(jj = 0; jj < *nsubject; jj++){

        rho_tmp[jj] = Si_iter[jj*(ntime1) + t];

    }

//      RprintIVecAsMat("rho_tmp", rho_tmp, 1, *nsubject);

//      // It seems to me that I can use some of the structure used to carry
//      // out Algorithm 8 from previous code to keep track of empty clusters
//      // etc.
    for(j = 0; j < *nsubject; j++){
        Rprintf("t ===== %d\n", t);
        Rprintf("j ===== %d\n", j);

        // Only need to update partition relative to units that are not pegged
        if(gamma_iter[j*(ntime1) + t] == 0){

            if(nh[(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t] > 1){

                // Observation belongs to a non-singleton ...
                nh[(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t] = nh[(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t] - 1;

            }else{

                // Observation is a member of a singleton cluster ...

```



```

//
iaux = Si_iter[j*(ntime1) + t];
Rprintf("iaux = %d\n", iaux);
if(iaux < nclus_iter[t]){

// Need to relabel clusters. I will do this by swapping cl
// Si_iter[j] and nclus_iter along with cluster specific p
arameters;

// All members of last cluster will be assigned subject j'
for(jj = 0; jj < *nsubject; jj++){
    if(Si_iter[jj*(ntime1) + t] == nclus_iter[t]){
        Si_iter[jj*(ntime1) + t] = iaux;
    }
}

Si_iter[j*(ntime1) + t] = nclus_iter[t];

// The following steps swaps order of cluster specific par
// so that the newly labeled subjects from previous step r
// their correct cluster specific parameters
auxs = sig2h[(iaux-1)*ntime1 + t];
sig2h[(iaux-1)*ntime1 + t] = sig2h[(nclus_iter[t]-1)*(ntime1)+t];
sig2h[(nclus_iter[t]-1)*(ntime1)+t] = auxs;

auxm = muh[(iaux-1)*ntime1 + t];
muh[(iaux-1)*ntime1 + t] = muh[(nclus_iter[t]-1)*(ntime1)+t];
muh[(nclus_iter[t]-1)*(ntime1)+t] = auxm;

// the number of members in cluster is also swapped with t
nh[(iaux-1)*(ntime1)+t] = nh[(nclus_iter[t]-1)*(ntime1)+t];
nh[(nclus_iter[t]-1)*(ntime1)+t] = 1;

}

// Now remove the ith obs and last cluster;
nh[(nclus_iter[t]-1)*(ntime1)+t] = nh[(nclus_iter[t]-1)*(ntime1)+t] - 1;
nclus_iter[t] = nclus_iter[t] - 1;

}

for(jj = 0; jj < *nsubject; jj++){
    rho_tmp[jj] = Si_iter[jj*(ntime1) + t];
}

//
RprintIVecAsMat("Si_iter", Si_iter, *nsubject, ntime1);
RprintIVecAsMat("nh ", nh, *nsubject, ntime1);
RprintIVecAsMat("rho_tmp", rho_tmp, 1, *nsubject);

//
Rprintf("nclus_iter[t] = %d\n", nclus_iter[t]);

for(k = 0; k < nclus_iter[t]; k++){
    Rprintf("k == %d\n\n", k);

    // Beginning of spatial part
    lCo = lCn = 0.0;
    if(*sPPM==1){
        indx = 0;
        for(jj = 0; jj < *nsubject; jj++){
            if(Si_iter[jj*(ntime1) + t] == k+1 & j != jj){

//
Rprintf("indx = %d\n", indx);

```

```

slo[indx] = s1[jj];
s2o[indx] = s2[jj];

sln[indx] = s1[jj];
s2n[indx] = s2[jj];

indx = indx+1;
}
if(j == jj){
sln[nh[k*(ntime1) + t]] = s1[jj];
s2n[nh[k*(ntime1) + t]] = s2[jj];

}

}

RprintVecAsMat("slo", slo, 1, nh[k*(ntime1) + t]);
RprintVecAsMat("s2o", s2o, 1, nh[k*(ntime1) + t]);
RprintVecAsMat("sln", sln, 1, nh[k*(ntime1) + t]+1);
RprintVecAsMat("s2n", s2n, 1, nh[k*(ntime1) + t]+1);
Rprintf("Cohesion = %d\n", *SpatialCohesion);

lCo = Cohesion3_4(slo, s2o, mu0, k0, v0, L0, nh[k*(ntime1)
lCn = Cohesion3_4(sln, s2n, mu0, k0, v0, L0, nh[k*(ntime1)

}
// End of spatial part

Rprintf("lCo = %f\n", lCo);
Rprintf("lCn = %f\n", lCn);
rho_tmp[j] = k+1;

RprintIVecAsMat("rho_tmp", rho_tmp, 1, *nsubject);
RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);
RprintIVecAsMat("Si_iter", Si_iter, *nsubject, ntime1);

// First need to check compatability
Rindx2=0;
for(jj = 0; jj < *nsubject; jj++){
if(gamma_iter[jj*ntime1 + (t+1)] == 1){
comp2t[Rindx2] = rho_tmp[jj];
comptpl[Rindx2] = Si_iter[jj*ntime1 + (t+1)];
Rindx2 = Rindx2 + 1;
}
}

Rprintf("Rindx2 = %d\n", Rindx2);
RprintIVecAsMat("comp2t", comp2t, 1, *nsubject);
RprintIVecAsMat("comptpl", comptpl, 1, *nsubject);

// check for compatibility
rho_comp = compatibility(comp2t, comptpl, Rindx2);

Rprintf("rho_comp = %d\n", rho_comp);

if(rho_comp != 1){

ph[k] = log(0); // Not compatible

} else {
// Need to compute Pr(rhot), Pr(rhot.R), Pr(rhot+1), Pr(rh
ot+1.R)

for(jj = 0; jj < *nsubject; jj++){
nh_tmp[jj] = 0;
nh_redtmp[jj] = 0;
nh_tmp_1[jj] = 0;
nh_redtmp_1[jj] = 0;
}
n_tmp = 0;
n_redtmp = 0;
n_tmp_1 = 0;
n_redtmp_1 = 0;

for(jj = 0; jj < *nsubject; jj++){
nh_tmp[rho_tmp[jj]-1] = nh_tmp[rho_tmp[jj]-1]+1;
n_tmp=n_tmp+1;
if(gamma_iter[jj*ntime1 + t] == 1){
nh_redtmp[rho_tmp[jj]-1] = nh_redtmp[rho_t
mp[jj]-1]+1;

```

```

n_redtmp = n_redtmp+1;
}

nh_tmp_1[Si_iter[jj*ntime1 + (t+1)]-1] = nh_tmp_1[
n_tmp_1=n_tmp_1+1;
if(gamma_iter[jj*ntime1 + t+1] == 1){
nh_redtmp_1[Si_iter[jj*ntime1 + (t+1)]-1]
n_redtmp_1 = n_redtmp_1+1;
}
}

RprintIVecAsMat("nh_tmp", nh_tmp, 1, *nsubject);
RprintIVecAsMat("nh_redtmp", nh_redtmp, 1, *nsubject);

Rprintf("nsubject = %d\n", *nsubject);
Rprintf("n_redtmp = %d\n", n_redtmp);

remove_zero(nh_tmp, *nsubject, nh_tmp_no_zero);
remove_zero(nh_redtmp, *nsubject, nh_redtmp_no_zero);

remove_zero(nh_tmp_1, *nsubject, nh_tmp_no_zero_1);
remove_zero(nh_redtmp_1, *nsubject, nh_redtmp_no_zero_1);

RprintIVecAsMat("nh_tmp_no_zero", nh_tmp_no_zero, 1, *nsubj
ect);
RprintIVecAsMat("nh_redtmp_no_zero", nh_redtmp_no_zero, 1,
*nsubject);

nclus_redtmp=0;
nclus_tmp=0;
nclus_redtmp_1=0;
nclus_tmp_1=0;
for(jj = 0; jj < *nsubject; jj++){
if(nh_redtmp[jj] > 0) nclus_redtmp = nclus_redtmp

if(nh_tmp[jj] > 0) nclus_tmp = nclus_tmp + 1;
if(nh_redtmp_1[jj] > 0) nclus_redtmp_1 = nclus_red

if(nh_tmp_1[jj] > 0) nclus_tmp_1 = nclus_tmp_1 + 1

}

Rprintf("nclus_tmp = %d\n", nclus_tmp);
Rprintf("nclus_redtmp = %d\n", nclus_redtmp);

lpp_full = partition_prob_crp(nh_tmp_no_zero, nclus_tmp, M
lpp_red = partition_prob_crp(nh_redtmp_no_zero, nclus_redt

lpp_full_1 = partition_prob_crp(nh_tmp_no_zero_1, nclus_tm
lpp_red_1 = partition_prob_crp(nh_redtmp_no_zero_1, nclus_

Rprintf("lpp_full = %f\n", lpp_full);
Rprintf("lpp_red = %f\n", lpp_red);

Rprintf("lpp_full_1 = %f\n", lpp_full_1);
Rprintf("lpp_red_1 = %f\n", lpp_red_1);

Rprintf("lpp_full = %f\n", lpp_full);
Rprintf("lpp_red = %f\n", lpp_red);

Rprintf("muh[k*(ntime1) + t] = %f\n", muh[k*(ntime1) + t])
Rprintf("sigh[k*(ntime1) + t] = %f\n", sqrt(sig2h[k*(ntime
1) + t]));

Rprintf("y[j*(*ntime) + t] = %f\n", y[j*(*ntime) + t]);
Rprintf("nh[k] = %d\n", nh_tmp[k]);
Rprintf("dnorm(y[j*(*ntime) + t], muh[k*(ntime1) + t], sqr
t(sig2h[k*(ntime1) + t]), 1) = %f\n", dnorm(y[j*(*ntime) + t], muh[k*(ntime1) + t], sqrt(sig2h[k*(ntime1) + t]), 1
));

if(t==0){

ph[k] = dnorm(y[j*(*ntime) + t],

```

```

        muh[k*(ntime1) + t],
        sqrt(sig2h[k*(ntime1) + t]), 1) +
lpp_full - lpp_red +
lpp_full_1 - lpp_red_1 +
lCn - lCo; // Spatial part of cohesion fun
ction;

    }
    if(t > 0){
        ph[k] = dnorm(y[j*(ntime) + t],
            muh[k*(ntime1) + t] +
            etal_iter[j]*y[j*(ntime)
                sqrt(sig2h[k*(ntime1) + t]*
                    (1-etal_iter[j]*etal_iter[
lpp_full - lpp_red +
lpp_full_1 - lpp_red_1 +
lCn - lCo; // Spatial part of cohesion fun
ction;

    }

    // use this to test if MCMC draws from prior are correct
    ph[k] = lpp_full - lpp_red + lpp_full_1 - lpp_red_1;

    //
    Rprintf("ph[k] = %f\n", ph[k]);

    }

}

//
RprintVecAsMat("ph = ", ph, 1, nclus_iter[t] );

// Determine if E.U. gets allocated to a new cluster
// Need to check compatibility first

//
Rprintf("nclus_iter[t] = %d\n", nclus_iter[k]);

rho_tmp[j] = nclus_iter[t]+1;

//
RprintIVecAsMat("rho_tmp", rho_tmp, 1, *nsubject);
RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);

// First need to check compatability
Rindx1 = 0, Rindx2=0;
for(jj = 0; jj < *nsubject; jj++){
    if(gamma_iter[jj*ntime1 + (t+1)] == 1){
        comp2t[Rindx2] = rho_tmp[jj];
        comptp1[Rindx2] = Si_iter[jj*ntime1 + (t+1)];
        Rindx2 = Rindx2 + 1;
    }
}

//
Rprintf("Rindx2 = %d\n", Rindx2);
//
RprintIVecAsMat("comp2t", comp2t, 1, *nsubject);
//
RprintIVecAsMat("comptp1", comptp1, 1, *nsubject);

// check for compatibility
rho_comp = compatibility(comp2t, comptp1, Rindx2);

//
Rprintf("rho_comp = %d\n", rho_comp);

if(rho_comp != 1){
    ph[nclus_iter[t]] = log(0); // going to own cluster is not compati
ble;
} else {

//
    RprintIVecAsMat("nh ", nh, *nsubject, ntime1);

//
    Rprintf("mu_iter[t] = %f\n", theta_iter[t]);
    Rprintf("sqrt(tau2_iter[t]) = %f\n", sqrt(tau2_iter[t]));

    mudraw = rnorm(theta_iter[t], sqrt(tau2_iter[t]));
    sigdraw = runif(0, A);

//
    Rprintf("mudraw = %f\n", mudraw);
//
    Rprintf("sigdraw = %f\n", sigdraw);
//
    Rprintf("y[j*(ntime) + t] = %f\n", y[j*(ntime) + t]);

//
    RprintIVecAsMat("nh_tmp", nh_tmp, 1, nclus_iter[t]);

```

```

    for(jj = 0; jj < *nsubject; jj++){
        nh_tmp[jj] = 0;
        nh_redtmp[jj] = 0;
        nh_tmp_1[jj] = 0;
        nh_redtmp_1[jj] = 0;
    }
    n_tmp = 0;
    n_redtmp = 0;
    n_tmp_1 = 0;
    n_redtmp_1 = 0;

    for(jj = 0; jj < *nsubject; jj++){
        nh_tmp[rho_tmp[jj]-1] = nh_tmp[rho_tmp[jj]-1]+1;
        n_tmp=n_tmp+1;
        if(gamma_iter[jj*ntime1 + t] == 1){
            nh_redtmp[rho_tmp[jj]-1] = nh_redtmp[rho_tmp[jj]-1
] + 1;

            n_redtmp = n_redtmp+1;
        }

        nh_tmp_1[Si_iter[jj*ntime1 + (t+1)]-1] = nh_tmp_1[Si_iter[
jj*ntime1 + (t+1)]-1]+1;

        n_tmp_1=n_tmp_1+1;
        if(gamma_iter[jj*ntime1 + t+1] == 1){
            nh_redtmp_1[Si_iter[jj*ntime1 + (t+1)]-1] = nh_red
tmp_1[Si_iter[jj*ntime1 + (t+1)]-1]+1;

            n_redtmp_1 = n_redtmp_1+1;
        }
    }

    // RprintIVecAsMat("nh_tmp", nh_tmp, 1, *nsubject);
    // RprintIVecAsMat("nh_redtmp", nh_redtmp, 1, *nsubject);

    remove_zero(nh_tmp, *nsubject, nh_tmp_no_zero);
    remove_zero(nh_redtmp, *nsubject, nh_redtmp_no_zero);

    remove_zero(nh_tmp_1, *nsubject, nh_tmp_no_zero_1);
    remove_zero(nh_redtmp_1, *nsubject, nh_redtmp_no_zero_1);

    // RprintIVecAsMat("nh_tmp_no_zero",nh_tmp_no_zero, 1, *nsubject);
    // RprintIVecAsMat("nh_redtmp_no_zero", nh_redtmp_no_zero, 1, *nsubje
ct);

    nclus_redtmp=0;
    nclus_tmp=0;
    nclus_redtmp_1=0;
    nclus_tmp_1=0;
    for(jj = 0; jj < *nsubject; jj++){
        if(nh_redtmp[jj] > 0) nclus_redtmp = nclus_redtmp + 1;
        if(nh_tmp[jj] > 0) nclus_tmp = nclus_tmp + 1;
        if(nh_redtmp_1[jj] > 0) nclus_redtmp_1 = nclus_redtmp_1 +
1;

        if(nh_tmp_1[jj] > 0) nclus_tmp_1 = nclus_tmp_1 + 1;
    }

    // Rprintf("nclus_tmp = %d\n", nclus_tmp);
    // Rprintf("nclus_redtmp = %d\n", nclus_redtmp);

    lpp_full = partition_prob_crp(nh_tmp_no_zero, nclus_tmp, Mdp, *nsu
bject, 1);

    lpp_red = partition_prob_crp(nh_redtmp_no_zero, nclus_redtmp, Mdp,
n_redtmp, 1);

    lpp_full_1 = partition_prob_crp(nh_tmp_no_zero_1, nclus_tmp_1, Mdp
, *nsubject, 1);

    lpp_red_1 = partition_prob_crp(nh_redtmp_no_zero_1, nclus_redtmp_1
, Mdp, n_redtmp_1, 1);

    // RprintIVecAsMat("nh_tmp", nh_tmp, 1, nclus_iter[t]+1);
    // Rprintf("nclus_iter = %d\n", nclus_iter[t]+1);

    lpp_full = partition_prob_crp(nh_tmp_no_zero, nclus_tmp, Mdp, *nsu
bject, 1);

    lpp_red = partition_prob_crp(nh_redtmp_no_zero, nclus_redtmp, Mdp,
n_redtmp, 1);

    lpp_full_1 = partition_prob_crp(nh_tmp_no_zero_1, nclus_tmp_1, Mdp
, *nsubject, 1);

    lpp_red_1 = partition_prob_crp(nh_redtmp_no_zero_1, nclus_redtmp_1
, Mdp, n_redtmp_1, 1);

    // Rprintf("lpp_full = %f\n", lpp_full);
    // Rprintf("lpp_red = %f\n", lpp_red);

```

```

// Rprintf("lpp_full_1 = %f\n", lpp_full_1);
// Rprintf("lpp_red_1 = %f\n", lpp_red_1);

// Rprintf("dnorm(y[j*(*ntime) + t], mudraw, sigdraw, 1) = %f\n", dno
rm(y[j*(*ntime) + t], mudraw, sigdraw, 1));

// spatial part
lCn_1 = 0.0;
Rprintf("lCn_1 = %f\n", lCn_1);
Rprintf("sPPM = %d\n", *sPPM);
if(*sPPM==1){
    s1o[0] = s1[j];
    s2o[0] = s2[j];
    lCn_1 = Cohesion3_4(s1o, s2o, mu0, k0, v0, l0, 1, *Spatial
Cohesion, 1);
}

// Rprintf("lCn_1 = %f\n", lCn_1);

if(t==0){
    ph[nclus_iter[t]] = dnorm(y[j*(*ntime) + t], mudraw, sigdr
lpp_full -
lpp_red +
- lpp_red_1 +
lCn_1; //t
his is spatial part
}
if(t > 0){
    ph[nclus_iter[t]] = dnorm(y[j*(*ntime) + t],
mudraw + etal_iter[j]*y[j*(*ntime)
sigdraw*sqrt(1-etal_iter[j]*etal_i
lpp_full -
lpp_red +
- lpp_red_1 +
lCn_1; //t
his is spatial part
}

// ph[nclus_iter[t]] = lpp_full - lpp_red + lpp_full_1 - lpp_red_1;

// Rprintf("ph[nclus_iter[t]] = %f\n", ph[nclus_iter[t]]);
}

// RprintVecAsMat("ph = ", ph, 1, nclus_iter[t] + 1);
// RprintIVecAsMat("rhotmp = ", rho_tmp, 1, *nsubject);

// Now compute the probabilities
for(k = 0; k < nclus_iter[t]+1; k++) phtmp[k] = ph[k];

R_rsort(phtmp, nclus_iter[t]+1) ;

// RprintVecAsMat("phtmp ", phtmp, 1, nclus_iter[t]+1);

maxph = phtmp[nclus_iter[t]];

// Rprintf("maxph = %f\n", maxph);

denph = 0.0;
for(k = 0; k < nclus_iter[t]+1; k++){
    ph[k] = exp(ph[k] - maxph);
    ph[k] = pow(exp(ph[k] - maxph), (1 - exp(-0.0001*(i+1))));
    denph = denph + ph[k];
}

// RprintVecAsMat("ph", ph, 1, nclus_iter[t]+1);

for(k = 0; k < nclus_iter[t]+1; k++){
    probh[k] = ph[k]/denph;

```

```

    }
    Rprintf("denph = %f\n", denph);

    RprintVecAsMat("probh", probh, 1, nclus_iter[t]+1);

    uu = runif(0.0,1.0);
    Rprintf("uu = %f\n", uu);

    cprobh= 0.0;;
    for(k = 0; k < nclus_iter[t]+1; k++){

        cprobh = cprobh + probh[k];

        if (uu < cprobh){

           iaux = k+1;
            break;

        }

    }

    Rprintf("iaux = %d\n \n \n", iaux);

    if(iaux <= nclus_iter[t]){

        Si_iter[j*(ntime1) + t] = iaux;
        nh[(Si_iter[j*(ntime1) + t]-1)*(ntime1)+t] = nh[(Si_iter[j*(ntime1
) + t]-1)*(ntime1)+t] + 1;

        rho_tmp[j] = iaux;
    }else{

        nclus_iter[t] = nclus_iter[t] + 1;
        Si_iter[j*(ntime1) + t] = nclus_iter[t];
        nh[(Si_iter[j*(ntime1) + t]-1)*(ntime1)+t] = 1;
        rho_tmp[j] = nclus_iter[t];

        muh[(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t] = mudraw;
        sig2h[(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t] = sigdraw*sigdraw;

    }

    Rprintf("Si_iter[j*(ntime1) + t] = %d\n", Si_iter[j*(ntime1) + t]);
    RprintVecAsMat("muh", muh, *nsubject, ntime1);
    RprintVecAsMat("sig2h", sig2h, *nsubject, ntime1);
    RprintIVecAsMat("Si_iter ", Si_iter, *nsubject, ntime1);
    RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);
    RprintIVecAsMat("nh ", nh, *nsubject, ntime1);
    RprintIVecAsMat("nclus_iter", nclus_iter, 1, ntime1);

}

}

RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);
RprintIVecAsMat("Si_iter ", Si_iter, *nsubject, ntime1);
RprintIVecAsMat("nh ", nh, *nsubject, ntime1);
RprintIVecAsMat("nclus_iter", nclus_iter, 1, ntime1);

for(j = 0; j < *nsubject; j++){
    Si_tmp[j] = Si_iter[j*(ntime1) + t];
    Si_tmp2[j] = 0;
    reorder[j] = 0;
}

// I believe that I have to make sure that groups are order so that
// EU one is always in the group one, and then the smallest index not
// with group 1 anchors group 2 etc.

relabel(Si_tmp, *nsubject, Si_tmp2, reorder, oldLab);

RprintIVecAsMat("gamma_iter", gamma_iter, *nsubject, ntime1);
RprintIVecAsMat("Si_tmp2", Si_tmp2, 1, *nsubject);
RprintIVecAsMat("reorder", reorder, 1, *nsubject);
RprintIVecAsMat("oldLab", oldLab, 1, nclus_iter[t]);

RprintIVecAsMat("Si_iter ", Si_iter, *nsubject, ntime1);
RprintIVecAsMat("nh ", nh, *nsubject, ntime1);
RprintIVecAsMat("nclus_iter", nclus_iter, 1, ntime1);

RprintVecAsMat("muh", muh, *nsubject, ntime1);

```

```

// RprintVecAsMat("sig2h", sig2h, *nsubject, ntime1);

for(j=0; j<*nsubject; j++){
    Si_iter[j*(ntime1) + t] = Si_tmp2[j];
}

for(k = 0; k < nclus_iter[t]; k++){
    mu_tmp[k] = muh[k*(ntime1)+t];
    sig2_tmp[k] = sig2h[k*(ntime1)+t];
}

for(k = 0; k < nclus_iter[t]; k++){
    nh[k*(ntime1)+t] = reorder[k];
    muh[k*(ntime1)+t] = mu_tmp[oldLab[k]-1];
    sig2h[k*(ntime1)+t] = sig2_tmp[oldLab[k]-1];
}

// RprintIVecAsMat("Si_iter ", Si_iter, *nsubject, ntime1);
// RprintIVecAsMat("nh ", nh, *nsubject, ntime1);
// RprintIVecAsMat("nclus_iter", nclus_iter, 1, ntime1);

// RprintVecAsMat("muh", muh, *nsubject, ntime1);
// RprintVecAsMat("sig2h", sig2h, *nsubject, ntime1);

// for(k = 0; k < nclus_iter[t]; k++) sig2h[k*(ntime1)+t] = 1.0;

for(k = 0; k < nclus_iter[t]; k++){

// Rprintf("k = %d\n", k);

// ////////////////////////////////////////
// //
// // udate muh //
// //
// ////////////////////////////////////////

Rprintf("sumy = %f\n", sumy);
Rprintf("nh[k*(ntime1)+t] = %d\n", nh[k*(ntime1)+t]);
Rprintf("sig2h[k*(ntime1) + t] = %f\n", sig2h[k*(ntime1) + t]);
Rprintf("tau2_iter[t] = %f\n", tau2_iter[t]);
Rprintf("theta_iter[t] = %f\n", theta_iter[t]);

if(t==0){
    sumy = 0.0;
    for(j = 0; j < *nsubject; j++){
        if(Si_iter[j*(ntime1) + t] == k+1){
            sumy = sumy + y[j*(ntime1)+t];
        }
    }
    s2star = 1/((double) nh[k*(ntime1)+t]/sig2h[k*(ntime1) + t] + 1/tau2_iter[
t]);
    mstar = s2star*( (1/sig2h[k*(ntime1) + t])*sumy + (1/tau2_iter[t])*theta_i
ter[t]);

}
if(t > 0){
    sumy = 0.0;
    sume2 = 0.0;
    for(j = 0; j < *nsubject; j++){
        if(Si_iter[j*(ntime1) + t] == k+1){
            sume2 = sume2 + 1.0/(1-etal_iter[j]*etal_iter[j]);
            sumy = sumy + (y[j*(ntime1)+t] - etal_iter[j]*y[j*(ntime1)
+t-1])/
            (1-etal_iter[j]*etal_iter[j]);
        }
    }

    s2star = 1/((1.0/sig2h[k*(ntime1) + t])*sume2 + 1/tau2_iter[t]);
    mstar = s2star*( (1.0/sig2h[k*(ntime1) + t])*sumy + (1/tau2_iter[t])*theta
_iter[t]);

}

// Rprintf("sume2 = %f\n", sume2);
// Rprintf("sumy = %f\n", sumy);
// Rprintf("mstar = %f\n", mstar);
// Rprintf("sqrt(s2star) = %f\n", sqrt(s2star));

```



```

        muh[k*(ntime1) + t] = rnorm(mstar, sqrt(s2star));

//      muh[k] = 0.0;
//      Rprintf("muh[k*(ntime1) + t] = %f\n", muh[k*(ntime1) + t]);
//
//      RprintVecAsMat("muh", muh, *nsubject, ntime1);

////////////////////////////////////
//                                                                    //
//      udate sig2h                                                    //
//                                                                    //
////////////////////////////////////
osig = sqrt(sig2h[k*(ntime1) + t]);
nsig = rnorm(osig, csigSIG);

if(nsig > 0.0 & nsig < A){
    lln = 0.0;
    llo = 0.0;
    if(t == 0){
        for(j = 0; j < *nsubject; j++){
            if(Si_iter[j*(ntime1) + t] == k+1){
                llo = llo + dnorm(y[j*(ntime)+t], muh[k*(ntime1)
+ t], osig,1);
                lln = lln + dnorm(y[j*(ntime)+t], muh[k*(ntime1)
+ t], nsig,1);
            }
        }
    }
    if(t > 0){
        for(j = 0; j < *nsubject; j++){
            if(Si_iter[j*(ntime1) + t] == k+1){
                llo = llo + dnorm(y[j*(ntime)+t], muh[k*(ntime1)
+ t] +
                    etal_iter[j]*y[j*(ntime) + t-1],
osig*sqrt(1-etal_iter[j]*etal_iter[j]),1);
                lln = lln + dnorm(y[j*(ntime)+t], muh[k*(ntime1)
+ t] +
                    etal_iter[j]*y[j*(ntime) + t-1],
nsig*sqrt(1-etal_iter[j]*etal_iter[j]),1);
            }
        }
    }
}

//      Rprintf("ms = %f\n", ms);
//      Rprintf("osig = %f\n", osig);
//      Rprintf("nsig = %f\n", nsig);
lln = llo + dunif(osig, 0.0, A, 1);
lln = lln + dunif(nsig, 0.0, A, 1);

//
//      Rprintf("llo = %f\n", llo);
//      Rprintf("lln = %f\n", lln);

llr = lln - llo;
uu = runif(0,1);

//
//      Rprintf("llr = %f\n", llr);
//      Rprintf("log(uu) = %f\n", log(uu));

if(log(uu) < llr){
    sig2h[k*(ntime1) + t] = nsig*nsig;
}

//      sig2h[k*(ntime1) + t] = 1.0;

}

//      Rprintf("sig2h[k*(ntime1) + t] = %f\n", sig2h[k*(ntime1) + t]);
//
//      RprintVecAsMat("sig2h", sig2h, *nsubject, ntime1);

}

//
//      RprintVecAsMat("muh", muh, *nsubject, ntime1);
//      RprintVecAsMat("sig2h", sig2h, *nsubject, ntime1);

```

```

////////////////////////////////////
//
//
// update theta (mean of mh)
//
//
//
////////////////////////////////////
summu = 0.0;
for(k = 0; k < nclus_iter[t]; k++){
    summu = summu + muh[k*(ntime1) + t];
    Rprintf("nh[k*(ntime1)+t] = %d\n", nh[k*(ntime1)+t]);
}
Rprintf("summu = %f\n", summu);
Rprintf("nclus_iter[t] = %d\n", nclus_iter[t]);

philsq = phil_iter*phil_iter;
lam2tmp = lam2_iter*(1.0 - philsq);

if(t==0){
    Rprintf("t = %d\n", t);

    s2star = 1.0/((double) nclus_iter[t]/tau2_iter[t] + 1.0/lam2_iter + philsq/lam2tmp);

    mstar = s2star*( (1.0/tau2_iter[t])*summu +
                    (1.0/lam2_iter)*phi0_iter +
                    (1.0/lam2tmp)*phil_iter*(theta_iter[t+1]-phi0_iter*(1-phil_iter))

);

    Rprintf("mstar = %f\n", mstar);
    Rprintf("sqrt(s2star) = %f\n", sqrt(s2star));

} else if(t==(ntime-1)){

    Rprintf("t = %d\n", t);
    s2star = 1.0/((double) nclus_iter[t]/tau2_iter[t] + 1.0/lam2tmp);
    mstar = s2star*((1.0/tau2_iter[t])*summu +
                    (1.0/lam2tmp)*(phi0_iter*(1-phil_iter) + phil_iter*theta_iter[t-1]

));
    Rprintf("mstar = %f\n", mstar);
    Rprintf("sqrt(s2star) = %f\n", sqrt(s2star));

} else {

    s2star = 1.0/((double) nclus_iter[t]/tau2_iter[t] + (1.0 + philsq)/lam2tmp);
    mstar = s2star*( (1.0/tau2_iter[t])*summu +
                    (1.0/lam2tmp)*(phil_iter*(theta_iter[t-1] + theta_iter[t+1]) +
                    phi0_iter*(1.0 - phil_iter)*(1.0 - phil_iter)));

}

Rprintf("mstar = %f\n", mstar);
Rprintf("sqrt(s2star) = %f\n", sqrt(s2star));

theta_iter[t] = rnorm(mstar, sqrt(s2star));

Rprintf("theta_iter = %f\n", theta_iter[t]);

////////////////////////////////////
//
//
// update tau2 (variance of mh)
//
//
//
////////////////////////////////////
ot = sqrt(tau2_iter[t]);
nt = rnorm(ot,csigTAU);

if(nt > 0){

    llm = 0.0;
    llo = 0.0;
    for(k = 0; k < nclus_iter[t]; k++){

        llm = llm + dnorm(muh[k*(ntime1) + t], theta_iter[t], ot,1);
        llm = llm + dnorm(muh[k*(ntime1) + t], theta_iter[t], nt,1);

    }

    Rprintf("ms = %f\n", ms);
    Rprintf("osig = %f\n", osig);
    Rprintf("nsig = %f\n", nsig);
    llo = llo + dunif(ot, 0.0, A0, 1);

```

```

lln = lln + dunif(nt, 0.0, A0, 1);

//      Rprintf("llo = %f\n", llo);
//      Rprintf("lln = %f\n", lln);

llr = lln - llo;
uu = runif(0,1);

//      Rprintf("llr = %f\n", llr);
//      Rprintf("log(uu) = %f\n", log(uu));

if(log(uu) < llr){
    tau2_iter[t] = nt*nt;

//      tau2_iter[t] = 5*5;
}

}

//      Rprintf("tau2_iter = %f\n", tau2_iter[t]);

}

//      RprintIVecAsMat("Si_iter ", Si_iter, *nsubject, ntime1);
//      //////////////////////////////////////
//      //
//      // update etal (temporal correlation parameter at likelihood)
//      //
//      //////////////////////////////////////
if(*etal_0==0){
    for(j = 0; j < *nsubject; j++){
//      Rprintf("j = %d\n", j);

//      Rprintf("etal_iter = %f\n", etal_iter[j]);

        elo = etal_iter[j];
        eln = rnorm(elo, csigETA1);

//      Rprintf("elo = %f\n", elo);
//      Rprintf("eln = %f\n", eln);

        if(eln < 1 & eln > -1){
            llo=lln=0.0;
            for(t=1; t<*ntime; t++){
                llo = llo + dnorm(y[j*(ntime)+t],
//      muh[(Si_iter[j*(nt
ime1) + t]-1)*(ntime1) + t] +
                elo*y[j*(ntime)+t-1],
                sqrt(sig2h[(Si_iter[j*(ntime1) + t]-1)*(nt
ime1) + t]*
//      (1-elo*elo)), 1);

                lln = lln + dnorm(y[j*(ntime)+t],
//      muh[(Si_iter[j*(nt
ime1) + t]-1)*(ntime1) + t] +
                eln*y[j*(ntime)+t-1],
                sqrt(sig2h[(Si_iter[j*(ntime1) + t]-1)
//      *(ntime1) + t]*
//      (1-eln*eln)), 1);

            }

//      Rprintf("llo = %f\n", llo);
//      Rprintf("lln = %f\n", lln);

            logito = log(0.5*(elo + 1)) - log(1 - 0.5*(elo+1));
            logitn = log(0.5*(eln + 1)) - log(1 - 0.5*(eln+1));

//      Rprintf("logito = %f\n", logito);
//      Rprintf("logitn = %f\n", logitn);

//      Rprintf("(1/b_etal)*fabs(logito - 0.0) = %f\n", (1/b_etal)*fabs(logito - 0
.0));
//      Rprintf("(1/b_etal)*fabs(logitn - 0.0) = %f\n", (1/b_etal)*fabs(logitn - 0

```

```

.0));

//
//
Rprintf("fabs(logito - 0.0) = %f\n", fabs(logito - 0.0));
Rprintf("fabs(logitn - 0.0) = %f\n", fabs(logitn - 0.0));

llo = llo + -log(2*b_etal) - (1/b_etal)*fabs(logito - 0.0);
lln = llm + -log(2*b_etal) - (1/b_etal)*fabs(logitn - 0.0);

//
//
Rprintf("llo = %f\n", llo);
Rprintf("lln = %f\n", llm);

llr = llm - llo;
uu = runif(0,1);

if(llr > log(uu)) etal_iter[j] = eln;

//
Rprintf("etal_iter = %f\n", etal_iter[j]);

}

}

//
RprintVecAsMat("etal", etal_iter, 1, *nsubject);

////////////////////////////////////
//
// update alpha
//
//
//
////////////////////////////////////
if(*alpha_0 == 0){
    if(*global_alpha == 1){
        sumg = 0;
        for(j = 0; j < *nsubject; j++){
            for(t = 1; t < *ntime; t++){

                sumg = sumg + gamma_iter[j*ntime1 + t];

            }

        }

//
Rprintf("sumg = %d\n", sumg);
astar = (double) sumg + a;
bstar = (double) ((*nsubject)*(*ntime-1) - sumg) + b;

//
Rprintf("astar = %f\n", astar);
Rprintf("bstar = %f\n", bstar);

alpha_tmp = rbeta(astar, bstar);

for(t=0;t<*ntime;t++){alpha_iter[t] = alpha_tmp;}
Rprintf("alpha_iter = %f\n", alpha_iter);

} else {

    for(t = 0; t < *ntime; t++){
        sumg = 0;
        for(j = 0; j < *nsubject; j++){
            sumg = sumg + gamma_iter[j*ntime1 + t];
        }

//
Rprintf("sumg = %d\n", sumg);

astar = (double) sumg + a;
bstar = (double) ((*nsubject) - sumg) + b;

//
Rprintf("astar = %f\n", astar);
Rprintf("bstar = %f\n", bstar);

alpha_iter[t] = rbeta(astar, bstar);

    }

}

alpha_iter[0] = 0.0;

}

```

```

// RprintVecAsMat("alpha_iter", alpha_iter, 1, *ntime);

////////////////////////////////////
//
//
// update phi0
//
//
//
////////////////////////////////////
phisq = phil_iter*phil_iter;
one_phisq = (1-phil_iter)*(1-phil_iter);
lam2tmp = lam2_iter*(1.0 - phisq);
// Rprintf("lam2tmp = %f\n", lam2tmp);
sumt = 0.0;
for(t=1; t<*ntime; t++){
// Rprintf("t = %d\n", t);
// Rprintf("theta_iter[t] = %f\n", theta_iter[t]);
// Rprintf("theta_iter[t-1] = %f\n", theta_iter[t-1]);
// Rprintf("phil_iter = %f\n", phil_iter);
// Rprintf("(theta_iter[t] - phil_iter*theta_iter[t-1]) = %f\n", (theta_iter[t] - phil_iter*th
eta_iter[t-1]));

sumt = sumt + (theta_iter[t] - phil_iter*theta_iter[t-1]);
// Rprintf("sumt = %f\n", sumt);
}

s2star = 1.0/(((*ntime-1)*(one_phisq/lam2tmp) + (1/lam2_iter) + (1/s20)));
mstar = s2star*((1.0-phil_iter)/lam2tmp)*sumt + (1/lam2_iter)*theta_iter[0] + (1/s20)*m0;

// Rprintf("sumt = %f\n", sumt);
// Rprintf("mstar = %f\n", mstar);
// Rprintf("s2star = %f\n", s2star);

// Rprintf("m0 = %f\n", m0);
// Rprintf("s20 = %f\n", s20);

phi0_iter = rnorm(mstar, sqrt(s2star));

// Rprintf("phi0_iter = %f\n", phi0_iter);

////////////////////////////////////
//
//
// update phil
//
//
//
////////////////////////////////////

if(*phil_0==0){
op1 = phil_iter;
np1 = rnorm(op1, csigPHI1);

// Rprintf("op1 = %f\n", op1);
// Rprintf("np1 = %f\n", np1);

if(np1 > -1 & np1 < 1){
llo = 0.0, llm = 0.0;
for(t=1; t < *ntime; t++){
// Rprintf("t = %d\n", t);
// Rprintf("theta_iter[t] = %f\n", theta_iter[t]);
// Rprintf("theta_iter[t-1] = %f\n", theta_iter[t-1]);
// Rprintf("lam2_iter = %f\n", lam2_iter);
// Rprintf("lam2_iter*(1.0 - op1*op1) = %f\n", lam2_iter*(1.0 - op1*op1));
// Rprintf("lam2_iter*(1.0 - np1*np1) = %f\n", lam2_iter*(1.0 - np1*np1));
// Rprintf("phi0_iter + op1*theta_iter[t-1] = %f\n", phi0_iter + op1*theta_it
er[t-1]);
// Rprintf("phi0_iter + np1*theta_iter[t-1] = %f\n", phi0_iter + np1*theta_it
er[t-1]);

llo = llo + dnorm(theta_iter[t], phi0_iter*(1-op1) + op1*theta_iter[t-1],
sqrt(lam2_iter*(1.0 - op1*op1)), 1);
llm = llm + dnorm(theta_iter[t], phi0_iter*(1-np1) + np1*theta_iter[t-1],
sqrt(lam2_iter*(1.0 - np1*np1)), 1);

// Rprintf("llo = %f\n", llo);
// Rprintf("llm = %f\n", llm);
}
// Rprintf("llo = %f\n", llo);
// Rprintf("llm = %f\n", llm);
llo = llo + dunif(op1, -1,1, 1);
llm = llm + dunif(np1, -1,1, 1);

```

```

        llr = lln - llo;
        Rprintf("llr = %f\n", llr);
        if(llr > log(runif(0,1))) phil_iter = npl;
    }
}
Rprintf("phil_iter = %f\n", phil_iter);

////////////////////////////////////
//
// update lam2
//
//
//
////////////////////////////////////
philsq = phil_iter*phil_iter;
ssq = 0.0;
for(t=1; t<*ntime; t++){
    ssq = ssq + (theta_iter[t] - (phi0_iter*(1-phil_iter) + phil_iter*theta_iter[t-1]))*
    (theta_iter[t] - (phi0_iter*(1-phil_iter) + phil_iter*theta_iter[t-1]));
}
ssq = 1.0/(1.0 - philsq)*ssq + (theta_iter[0]-phi0_iter)*(theta_iter[0]-phi0_iter);

astar = 0.5*(ntime) + al;
bstar = 0.5*ssq + 1/bl;

lam2_iter = 1.0/rgamma(astar, 1/bstar);

// Update lambda with a MH step
philsq = phil_iter*phil_iter;

ol = sqrt(lam2_iter);
nl = rnorm(ol, csigLAM);
if(nl > 0.0){
    lln = 0.0;
    llo = 0.0;
    for(t=1; t<*ntime; t++){
        llo = llo + dnorm(theta_iter[t],
            phi0_iter*(1-phil_iter) + phil_iter*theta_iter[t-1], ol*sqrt(1-p
hilsq),1);
        lln = lln + dnorm(theta_iter[t],
            phi0_iter*(1-phil_iter) + phil_iter*theta_iter[t-1], nl*sqrt(1-p
hilsq),1);
    }
    llo = llo + dnorm(theta_iter[0], phi0_iter, ol, 1) + dunif(ol, 0.0, Al, 1);
    lln = lln + dnorm(theta_iter[0], phi0_iter, nl, 1) + dunif(nl, 0.0, Al, 1);

    llr = lln - llo;
    uu = runif(0,1);

    if(log(uu) < llr){
        lam2_iter = nl*nl;
    }
}

Rprintf("lam2_iter = %f\n", lam2_iter);

////////////////////////////////////
//
// predict partition for new time period
//
// THIS HAS YET TO BE FINISHED
//
////////////////////////////////////
/*
for(p = 0; p < *npred; p++){
    for(j=0; j<*nsubject; j++){
        nh_pred[j] = 0;
        predSi_iter[j*(npred) + p] = 0;
    }
    RprintIVecAsMat("nh_pred", nh_pred, 1, *nsubject);

    if(*alpha_0 == 1){
        n_red = 0;
        for(j=0; j<*nsubject; j++){
            gpred[j] = rbinom(1,0.0);

            if(gpred[j] == 1){
                nh_pred[Si_iter[j*(ntime1)+(*ntime)-1] - 1] = nh_pred[Si_iter[j*(n
time1)+(*ntime)-1] - 1] + 1;

```

```

        n_red = n_red + 1;

        predSi_iter[j*(npred) + p] = Si_iter[j*(ntime1)+(*ntime)-1];
    }

}

if(*alpha_0 == 0){
    if(*global_alpha == 1){

        n_red = 0;
        for(j=0; j<*nsubject; j++){

            gpred[j] = rbinom(1,alpha_iter[1]);

            if(gpred[j] == 1){
                nh_pred[Si_iter[j*(ntime1)+(*ntime)-1] - 1] = nh_pred[Si_i
ter[j*(ntime1)+(*ntime)-1] - 1] + 1;

                n_red = n_red + 1;

                predSi_iter[j*(npred) + p] = Si_iter[j*(ntime1)+(*ntime)-
1];
            }

        }

    }else {

    }

}

// RprintIVecAsMat("predSi_iter", predSi_iter, npred, nsubject);
// RprintIVecAsMat("gpred", gpred, 1, nsubject);
// RprintIVecAsMat("nh_pred", nh_pred, 1, nsubject);
// Rprintf("n_red = %d\n", n_red);

remove_zero(nh_pred, nsubject, nh_tmp_no_zero);
// RprintIVecAsMat("nh_tmp_no_zero", nh_tmp_no_zero, 1, nsubject);

nclus_tmp = 0;
for(j=0; j<*nsubject; j++){
    if(nh_tmp_no_zero[j] > 0){
        nclus_tmp = nclus_tmp + 1;
    }else{
        break;
    }
}

// Rprintf("nclus_tmp = %d\n", nclus_tmp);

for(j=0; j<*nsubject; j++){
    Rprintf("j = %d\n", j);
    if(gpred[j] == 0){
        for(k = 0; k < nclus_tmp; k++){
            probh[k] = nh_pred[k]/(n_red + Mdp);
        }
        probh[nclus_tmp] = Mdp/(n_red + Mdp);

        RprintVecAsMat("probh = ", probh, 1, nclus_tmp+1);

        uu = runif(0.0,1.0);

        cprobh= 0.0;;
        for(k = 0; k < nclus_tmp+1; k++){

            cprobh = cprobh + probh[k];

            if (uu < cprobh){

               iaux = k+1;
                break;
            }

        }

        Rprintf("iaux = %d\n", iaux);
        if(iaux <= nclus_tmp){

            predSi_iter[j*(npred) + p] = iaux;
            nh_pred[iaux-1] = nh_pred[iaux-1] + 1;
        }else{

            nclus_tmp = nclus_tmp + 1;
            predSi_iter[j*(npred) + p] = nclus_tmp;

```

```

        nh_pred[(predSi_iter[j*(npred) + p]-1)*(npred)+p] = 1;
    }
    n_red = n_red + 1;

// RprintIVecAsMat("predSi_iter", predSi_iter, npred, nsubject);
// RprintIVecAsMat("nh_pred", nh_pred, 1, nsubject);
// Rprintf("nclus_tmp = %d\n", nclus_tmp);
// Rprintf("n_red = %d\n", n_red);
    }
}

*/
//
//
// evaluating likelihood that will be used to calculate LPML and WAIC?
// (see page 81 Christensen Hansen and Johnson)
//
//
//
if((i > (*burn-1)) & ((i+1) % (*thin) == 0)){
    like0=0;
    for(j = 0; j < nsubject; j++){
// Rprintf("j = %d\n", j);
        for(t = 0; t < *ntime; t++){
// Rprintf("t = %d\n", t);

// Rprintf("Si_iter[j*(ntime1) + t] = %d\n", Si_iter[j*(ntime1) + t]);
// Rprintf("(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t = %d\n", (Si_iter[j*(nti
me1) + t]-1)*(ntime1) + t);

            mudraw = muh[(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t];
            sigdraw = sqrt(sig2h[(Si_iter[j*(ntime1) + t]-1)*(ntime1) + t]);

// Rprintf("mudraw = %f\n", mudraw);
// Rprintf("sigdraw = %f\n", sigdraw);
// Rprintf("etal_iter[j] = %f\n", etal_iter[j]);
// Rprintf("y[j*(ntime)+t] = %f\n", y[j*(ntime)+t]);
// Rprintf("y[j*(ntime)+t-1] = %f\n", y[j*(ntime)+t-1]);

            if(t == 0){
// Rprintf("mudraw = %f\n", mudraw);
// Rprintf("sigdraw = %f\n", sigdraw);
// Rprintf("y[j*(ntime)+t] = %f\n", y[j*(ntime)+t]);

                like_iter[j*(ntime)+t] = dnorm(y[j*(ntime)+t], mudraw, sigdraw, 1);
                fitted_iter[j*(ntime)+t] = mudraw;
// Rprintf("like_iter = %f\n", like_iter[j*(ntime)+t]);
            }
            if(t > 0){

// Rprintf("mudraw + etal_iter[j]*y[j*(ntime)+t-1] = %f\n", mudraw +
// etal_iter[j]*y[j*(ntime)+t-1]);

                like_iter[j*(ntime)+t] = dnorm(y[j*(ntime)+t],
                                                    mudraw + e
                                                    sigdraw*sq
rt(1-etal_iter[j]*etal_iter[j]), 1);
                fitted_iter[j*(ntime)+t] = mudraw + etal_iter[j]*y[j*(ntime)+t-1];

            }

// Rprintf("like_iter = %f\n", like_iter[j*(ntime)+t]);

// These are needed for WAIC
mnlike[j*(ntime)+t] = mnlike[j*(ntime)+t] + exp(like_iter[j*(ntime)+t])
mnllike[j*(ntime)+t] = mnllike[j*(ntime)+t] + (like_iter[j*(ntime)+t])/

// Rprintf("mnlike = %f\n", mnlike[j*(ntime)+t]);
// Rprintf("mnllike = %f\n", mnllike[j*(ntime)+t]);

// if(exp(like_iter[j*(ntime)+t]) < 1e-320) like0=1;

CPO[j*(ntime)+t] = CPO[j*(ntime)+t] + (1/(double) nout)*(1/exp(like_iter
[j*(ntime)+t]));

    }
}

// Rprintf("like0 = %d\n", like0);
// if(like0==1) nout_0 = nout_0 + 1;
// if(i == (*draws-1)) Rprintf("xb = %f\n", xb);
// Rprintf("nout_0 = %d\n", nout_0);

```



```

//          if(like0==0){
//              Rprintf("nout - nout_0 = %d\n", nout - nout_0);
//              for(j = 0; j < *nsubject; j++){
//                  for(t = 0; t < *ntime; t++){
//                      Rprintf("like_iter[j*(*ntime)+t] = %f\n", like_iter[j*(*ntime)+t])
//                  }
//                  Rprintf("exp(like_iter[j*(*ntime)+t]) = %f\n", exp(like_iter[j*(*n
time)+t]));
//                  CPO[j*(*ntime)+t] = CPO[j*(*ntime)+t] + (1/exp(like_iter[j*(*ntime
)+t]));
//
//                      Rprintf("CPO = %f\n", CPO[j*(*ntime)+t]);
//
//              }
//          }
//      }
//      RprintVecAsMat("llike", like_iter, *nsubject, *ntime);
//  }

////////////////////////////////////
//
//          //
//          // Save MCMC iterates
//          //
//          //
//          //
//          //
//          if((i > (*burn-1)) & ((i+1) % *thin == 0)){

//          Rprintf("ii = %d\n", ii);
//          RprintVecAsMat("theta", theta_iter, 1, *ntime);

//          for(t = 0; t < *ntime; t++){
//              alpha_out[ii*(*ntime) + t] = alpha_iter[t];
//              theta[ii*(*ntime) + t] = theta_iter[t];
//              tau2[ii*(*ntime) + t] = tau2_iter[t];

//          for(j = 0; j < *nsubject; j++){
//              Rprintf("(Si_iter[j]-1)*(ntime1) + t = %d\n", (Si_iter[j*(ntime1) + t]-1)*
(ntime1) + t);

//              sig2[(ii*(*nsubject) + j)*(*ntime) + t] = sig2h[(Si_iter[j*(ntime1) + t]-1
)*(*ntime1) + t];

//              mu[(ii*(*nsubject) + j)*(*ntime) + t] = muh[(Si_iter[j*(ntime1) + t]-1)*(n
time1) + t];

//              Si[(ii*(*nsubject) + j)*(*ntime) + t] = Si_iter[j*ntime1 + t];
//              gamma[(ii*(*nsubject) + j)*(*ntime) + t] = gamma_iter[j*ntime1 + t];

//              llike[(ii*(*nsubject) + j)*(*ntime) + t] = like_iter[j*(*ntime)+t];
//              fitted[(ii*(*nsubject) + j)*(*ntime) + t] = fitted_iter[j*(*ntime)+t];

//          }

//      }

//          for(j=0; j<*nsubject; j++){

//              etal[ii*(*nsubject) + j] = etal_iter[j];

//          }

//          phil[ii] = phil_iter;
//          phi0[ii] = phi0_iter;
//          lam2[ii] = lam2_iter;

//          ii = ii+1;
//          Rprintf("ii = %d\n", ii);

//      }

}

lpml_iter=0.0;
for(t = 0; t < *ntime; t++){
//      Rprintf("t = %d\n", t);
//      for(j = 0; j < *nsubject; j++){
//          Rprintf("j = %d\n", j);

```

```

//          Rprintf("CPO = %f\n", CPO[j*(ntime)+t]);

//          lpml_iter = lpml_iter - log((1/(double) nout-nout_0)*CPO[j*(ntime)+t]);
//          lpml_iter = lpml_iter - log(CPO[j*(ntime)+t]);

    }

//  Rprintf("nout_0 = %d\n", nout_0);
//  lpml[0] = lpml_iter;
//  Rprintf("lpml_iter = %f\n", lpml_iter);

elppdWAIC = 0.0;
//  RprintVecAsMat("mnlike", mnlike, 1, (*nsubject)*(ntime1));
//  RprintVecAsMat("mnllike", mnllike, 1, (*nsubject)*(ntime1));

for(j = 0; j < *nsubject; j++){
    for(t = 0; t < *ntime; t++){
        elppdWAIC = elppdWAIC + (2*mnllike[j*(ntime)+t] - log(mnlike[j*(ntime)+t]));
    }
}
waic[0] = -2*elppdWAIC;

PutRNGstate();

}

```