

PDAF Tutorial

**Implementation of the analysis step
in online mode with a parallel model**



<http://pdaf.awi.de>

PDAF Parallel
Data Assimilation
Framework

Implementation Tutorial for PDAF online with serial model

We demonstrate the implementation
of an online analysis step with PDAF
with a model that is parallelized
using the template routines provided by PDAF

The example code is part of the PDAF source code package
downloadable at <http://pdaf.awi.de>

(This tutorial is compatible with PDAF V1.16 and later)

Implementation Tutorial for PDAF online / parallel model

This is just an example!

For the complete documentation of PDAF's interface
see the documentation
at <http://pdaf.awi.de>

Overview

Focus on Error Subspace Transform Kalman Filter
(ESTKF, Nerger et al., Mon. Wea. Rev. 2012)

4 Parts

a) Global filter

b) Localized filter

We recommend to first implement the global filter. The localized filter re-uses routines of the global filter.

In this tutorial we only cover the case of a parallel model. The implementation using a model without parallelization is described in a separate tutorial.

0a) Files for the Tutorial

Tutorial implementation

Files are in the PDAF package

Directory:

```
/tutorial/classical/online_2D_parallelmodel
```

- Fully working implementations of user codes
- PDAF core files are in `/src`
Makefile refers to it and compiles the PDAF library
- Only need to specify the compile settings (compiler, etc.) by environment variable `PDAF_ARCH`. Then compile with 'make'.

Template files for online mode

Directory: `/templates/online`

- Contains all required files
- Contains also command line parser (convenient but not required)

To generate your own implementation:

1. Copy content of directory
e.g. into sub-directory of model source code
2. Add calls to interface routines to model code
3. Complete user-routines for your model
4. Adapt compilation (e.g. Makefile) and compile
5. Run with assimilation options

PDAF library

Directory: `/src`

- The PDAF library is not part of the template
- PDAF is compiled separately as a library and linked when the assimilation program is compiled
- Makefile includes a compile step for the PDAF library
- One can also `cd` to `/src` and run 'make' there (requires setting of `PDAF_ARCH`)

`$PDAF_ARCH`

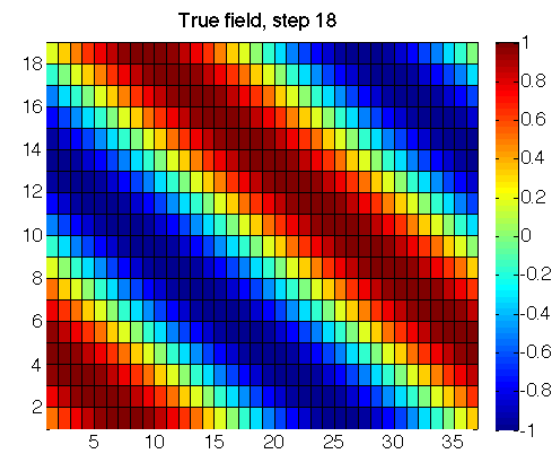
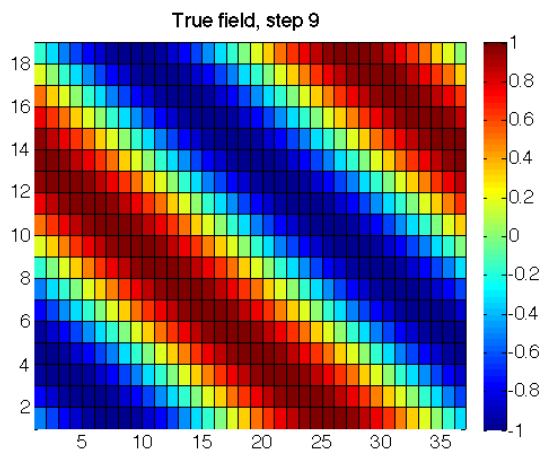
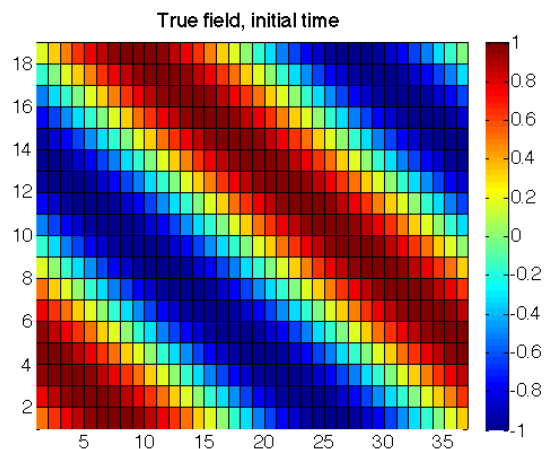
- Environment variable to specify the compile specifications
- Definition files in `/make.arch`
- Define by, e.g.

```
setenv PDAF_ARCH linux_gfortran (tcsh/csh)
export PDAF_ARCH=linux_gfortran (bash)
```


0b) The parallelized model

2D „Model“

- See the separate tutorial slides about the model
- Simple 2-dimensional grid domain
- 36 x 18 grid points (longitude x latitude)
- True state: sine wave in diagonal direction (periodic for consistent time stepping)



Parallel Mode1: Files

The source code of the parallel model consists of the following files:

- mod_model.F90
- mod_parallel_model.F90
- main.F90
- initialize.F90
- integrate.F90

Note: One can nicely compare the source codes of the model without and with parallelization

For clarity, the implementation with PDAF is found in

- main_pdaf.F90
- integrate_pdaf.F90

It allows for easy comparison of the implementations

0c) state vector and observation vector

State vector – some terminology used later

- PDAF performs computations on state vectors
- **State vector**
 - Stores model fields in a single vector
 - Tutorial shows this for one 2-dimensional field
 - Multiple fields are just concatenated into the vector
 - All fields that should be modified by the assimilation have to be in the state vector
- **State dimension**
 - Is the length of the state vector
(the sum of the sizes of the model fields in the vector)
- **Ensemble array**
 - Rank-2 array which stores state vectors in its columns

Observation vector

- **Observation vector**
 - Stores all observations in a single vector
 - Tutorial shows this for one 2-dimensional field
 - Multiple observed fields are just concatenated into the vector
- **Observation dimension**
 - Is the length of the observation vector
(sum of the observations over all observed fields in the vector)
- **Observation operator**
 - Operation that computes the observed part of a state vector
 - Tutorial only selects observed grid points
 - The operation can involve interpolation or integration depending on type of observation

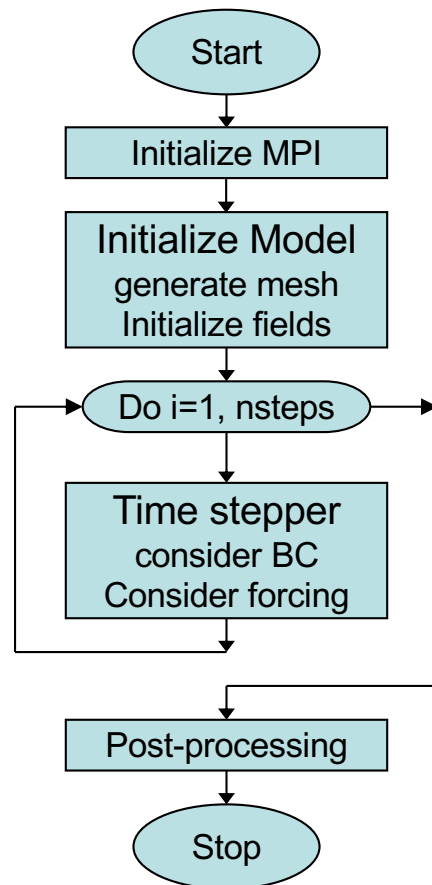
0c) PDAF online mode

Online mode

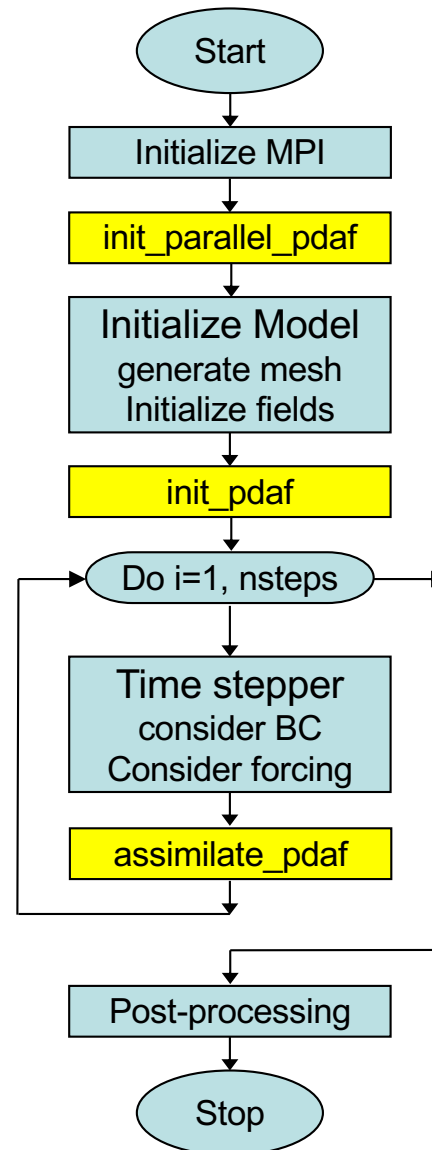
- Combine model with PDAF into single program
 - “model_pdaf”
- Add 3 subroutine calls:
 - `init_parallel_pdaf` - revise parallelization
 - `init_pdaf` - initialize assimilation
 - `assimilate_pdaf` - perform assimilation
- Implement user-supplied routines, e.g. for
 - observation operator
 - initialization of observation vector
 - transfer between state vector and model fields

Program flow with model extended for data assimilation

Simulation Model



Assimilation System



Legend

Model

Extension for data assimilation

Fully parallel configuration

- Tutorial shows implementation for a fully parallel case
 - Number of processes equals ensemble size times number of processes used for a single model task!
- For a more flexible (and complicated) configuration see PDAF's online guide

model_pdaf: General program structure

```
program main_pdaf
```

```
    init_parallel_pdaf
```

- initialize parallelization

```
    initialize
```

- initialize model information

```
    init_pdaf
```

- initialize parameters for PDAF
and read ensemble

```
    integrate
```

- time stepping loop

```
        assimilate_pdaf
```

- compute analysis step
(called inside stepping loop)

```
end program
```

Note:

In the example code, we use different files main.F90 and main_pdaf.F90 to allow for easy comparison

mod_assimilation.F90

Fortran module

- Declares the parameters used to configure PDAF
- Will be included (with 'use') in the user-written routines
- Additions to template necessary for observation handling

0d) Inserting subroutine calls

Where to insert subroutine calls?

`init_parallel_pdaf`

- at the start of the program, but after the `MPI_Init` performed in the code of the parallel model

`init_pdaf`

- after the initialization of the model
i.e. directly before the time stepping loop

`assimilate_pdaf`

- Last operation in the time stepping loop
i.e. just before the 'END DO'

Note: One can add the routines one after the other:

First insert `init_parallel_pdaf` and test the program, then add `init_pdaf`, etc.

init_parallel_pdaf.F90

- It is fully implemented template usable with small adaptations
- Required adaptations
 - Include MPI variables from module of the model:
MPI_COMM_WORLD, COMM_model, mype_model, npes_model
(the latter three variables might be named differently in a model)
 - init_parallel_pdaf defines a model communicator
`comm_model`
(actually it's a set for communicators, one for each model task)
 - Set communicator of the parallel model to `comm_model` at the end
if `init_parallel_pdaf`:
`"my_models_communicator" = comm_model`
(include `my_models_communicator` from module of model)
 - Set variables for number of processes in model and rank of a
process (`npes_model`, `mype_model`) at end of routine

init_parallel_pdaf.F90 (2)

- Parallelization variables for PDAF are declared in Fortran module

```
mod_parallel_pdaf
```

- Important variable:

```
n_modeltasks
```

- Defines number of concurrent model integrations.
 - Has to be equal to ensemble size
 - In the example: Read as 'dim_ens' from command line (using subroutine 'parse')
- **Important:** If the parallel model uses MPI_COMM_WORLD, this has to be replaced! (MPI_COMM_WORLD denotes always all processes in the program)

init_parallel_pdaf.F90 (3) - Example

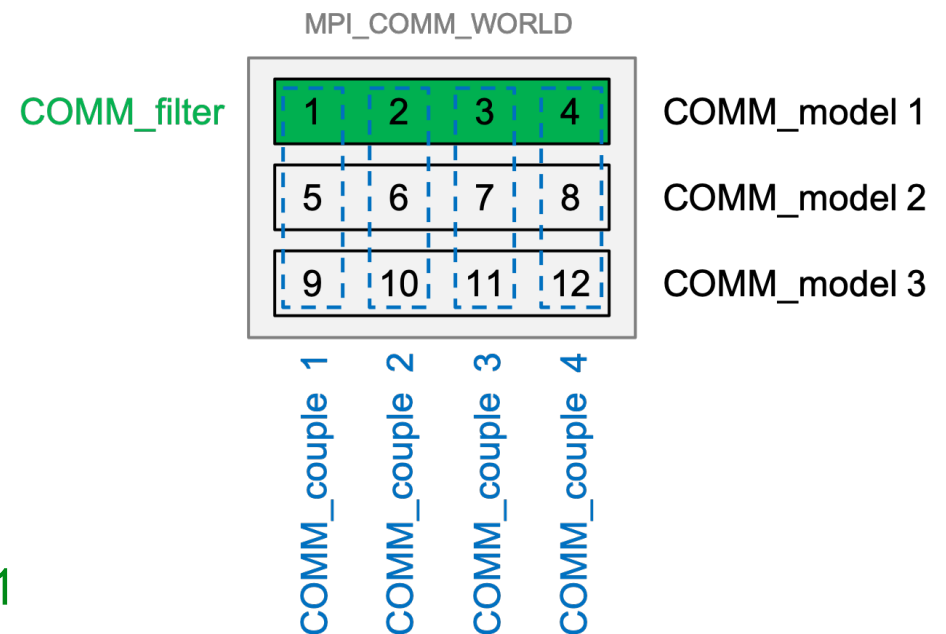
The routine initializes 3 groups of communicators

- COMM_model: Used to run the parallel model forecasts
- COMM_filter: Used to compute the filter
- COMM_couple: Coupling between model and filter processes

These are provided to PDAF when calling PDAF_init

The figure shows an example

- 12 processes in total
- 3 model tasks in parallel
- Each model task uses 4 processes in its COMM_model
- Each COMM_couple links groups of 3 processes to distribute and collect ensemble states
- The filter processes use model task 1



init_parallel_pdaf is coded to provide this configuration when running with 12 processes and setting dim_ens=3

init_pdaf.F90

Routine sets parameters for PDAF, calls `PDAF_init` to initialize the data assimilation, and calls `PDAF_get_state` to prepare the ensemble integrations:

Template contains list of available parameters (declared in and used from `mod_assimilation`)

Independent of the filter algorithm:

- Include information on size of model fields from model
- Define dimension of decomposed state vector

```
dim_state_p = nx_p * ny
```

In call to `PDAF_init`, the name of the user-supplied routine for ensemble initialization routine is specified:

```
init_ens_pdaf
```

init_pdaf.F90 (II)

In call to PDAF_get_state, the names of 3 user-supplied routines are specified:

`next_observation_pdaf`

- Set number of time steps in forecast phase

`distribute_state_pdaf`

- Initialize model fields from state vector

`prepoststep_ens_pdaf`

- poststep routine (compute estimated errors, write state estimate, etc.)

Initially, one can just copy the template routines. One can adapt them later to the particular application.

assimilate_pdaf.F90

Routine just calls a filter-specific routine like

```
PDAF_assimilate_estkf
```

We don't insert `PDAF_assimilate_estkf` directly into the model code

→ because, we need to declare all user-supplied routines as 'EXTERNAL'. This could clutter the model code.

Filter-specific user routines are described next. Initially, one can just copy the template routines.

Note: Template contains calls for `PDAF_assimilate_estkf` and `PDAF_assimilate_lestkf`. Need to adapt for other filters

Differences online and offline

- If you've studied the tutorial for offline mode

Offline

- Separate programs for model and assimilation
- Needed to implement routine `initialize`
- Grid dimensions declared in `mod_assimilation`
- Ensemble information read from files
- `mod_assimilation` contains all field and assimilation variables

Online

- Extend model program for assimilation
- Operations in `initialize` given by model; no changes for assimilation!
- Grid dimensions defined in model code (`mod_model`)
- Ensemble information provided by model fields
- `mod_assimilation` only contains variables for assimilation

Optional routine: `finalize_pdaf.F90`

Call to `finalize_pdaf` can be inserted at the end of the model

Routine contains two calls to `PDAF_print_info`:

```
CALL PDAF_print_info(2)
```

– display information on allocated memory inside PDAF

```
CALL PDAF_print_info(1)
```

– display timing information

(values 3 and 4 also possible for more detailed timers)

Note: `finalize_pdaf` only prints the information for `mype_world==0`

In addition there is

```
CALL PDAF_deallocate()
```

which deallocates internal arrays in PDAF

0e) Forecast phase

Files for PDAF

Template contains all required files

- just need to be filled with functionality

`init_pdaf.F90`

`init_ens_pdaf.F90`

`next_observation_pdaf.F90`

`distribute_state_pdaf.F90`

`collect_state_pdaf.F90`

`init_dim_obs_pdaf.F90`

`obs_op_pdaf.F90`

`init_obs_pdaf.F90`

`prodrinva_pdaf.F90`

`prepoststep_ens_pdaf.F90`

} initialization

} ensemble
forecast

} analysis step

init_pdaf.F90

Routine sets parameters for PDAF and calls `PDAF_init` to initialize the data assimilation:

Template contains list of available parameters (declared in and used from `mod_assimilation`)

For the example set :

1. `dim_ens = 9`
2. `rms_obs = sqrt(0.5)`
3. `filtertype = 6` (for ESTKF)
4. `delt_obs = 2` (assimilate after each 2nd time step)

In call to `PDAF_init`, the name of the ensemble initialization routine is specified:

```
init_ens_pdaf
```

init_ens_pdaf.F90

A *call-back* routine called by PDAF_init:

- Implemented by the user
- Its name is specified in the call to PDAF_init
- It is called by PDAF through a defined interface:

```
SUBROUTINE init_ens_pdaf(filtertype, dim_p,  
                        dim_ens, state_p, Uinv, ens_p, flag)
```

Declarations in header of the routine shows “intent” (input, output):

```
REAL, INTENT(out)      :: ens_p(dim_p, dim_ens)
```

Note:

All call-back routines have a defined interface and show the intent of the variables. Their header comment explains what is to be done in the routine.

init_ens_pdaf.F90 (2)

Initialize ensemble matrix `ens_p` for the start time of the assimilation

1. Include `nx`, `ny`, `nx_p` with `use mod_model`
2. Declare and allocate `real :: field(ny, nx)`
3. Loop over ensemble files (`i=1,dim_ens`)

for each file:

- read ensemble state into `field`
- store **local part** of `field` in column `i` of `ens_p`
(columns `nx_p*mype_model+1 : nx_p*mype_model+nx_p`)

4. Deallocate `field`

Note:

Columns of `ens_p` are state vectors. Store following storage of field in memory (column-wise in Fortran)

The forecast phase

At this point the initialization of PDAF is complete:

- Initial Ensemble of model states is initialized
- Filter algorithm and its parameters are chosen

Next:

- Implement user-routines for forecast phase
- All are call-back routines:
 - User-written, but called by PDAF

Note:

Some variables end with `_p`.

It means that the variable is specific for a process
(its values are different for each process)

next_observation_pdaf.F90

Routine to

- Set number of time steps in next forecast phase
- Set flag to control exit from forecasts (`doexit`)

Most simple setting:

```
include delt_obs from mod_assimilation  
nsteps = delt_obs  
doexit = 0
```

Note: The assimilation program stops when the maximum number of time steps of the model is reached, even if `doexit=0`

next_observation_pdaf.F90 (II)

More sophisticated setting:

- Utilize `stepnow` (current time step) and `total_steps` (total number of time steps given by model).

```
IF (stepnow + nsteps <= total_steps) THEN
    nsteps = delt_obs           ! Forecast length
    doexit = 0                 ! Continue assimilation
ELSE
    nsteps = 0                 ! No more steps
    doexit = 1                 ! Exit assimilation
END IF
```

Note: In the example `doexit=1` is used only inside PDAF and avoids some screen output.

distribute_state_pdaf.F90

Routine to

- Initialize model fields from a state vector
- Routine is provided with the state vector `vector_p`

For the example:

1. Access `nx_p`, `ny` and `field_p` with use `mod_model`

2. Initialize model field from state vector:

```
DO j = 1, nx_p
    field_p(1:ny, j) = state_p(1+(j-1)*ny : j*ny)
END DO
```

prepoststep_ens_pdaf.F90

Post-step routine for the online mode:

Already there in the template:

1. Compute ensemble mean state `state_p`
2. Compute estimated variance vector `variance`
3. Compute estimated root mean square error `rmerror_est`

Possible extensions:

4. Write analysis state (ensemble mean, `state_step*_ana.txt`)
5. Write analysis ensemble into files
(Analogous to reading in `init_ens_pdaf`)
6. Analogously one can write the forecast fields

Completion of forecast phase

At this point the implementation of the forecast phase is practically complete:

- Initial ensemble and PDAF's parameters are set
- The ensemble forecast can be computed

One can now compile the program `model_pdaf` (make `model_pdaf`) to check whether it runs.

Note: It is recommended to compile PDAF with `-DPDAF_NO_UPDATE` at this point as the routine for the analysis step are not yet implemented.

Note: For now, `prepoststep_ens_pdaf` only lets you test the initial ensemble. Testing the forecast fields need implementation of routine `collect_state_pdaf`

1a) Global filter

Running the tutorial program

- `cd` to `/tutorial/classical/online_2D_serialmodel`
- Set environment variable `PDAF_ARCH` or set it in Makefile (e.g. `linux_gfortran_openmpi`)
- Compile by running `'make model_pdaf'` (next slide will discuss possible compile issues)
- Run the program with

```
mpirun -np 18 ./model_pdaf -dim_ens 9
```
- Inputs are read in from `/tutorial/inputs_online`
- Outputs are written in
`/tutorial/classical/online_2D_parallelmodel`
- Plot result, e.g. with 'octave':

```
load state_step10_ana.txt  
pcolor(state_step10_ana)
```

Requirements for compiling PDAF

PDAF requires libraries for BLAS and LAPACK

- Libraries to be linked are specified in the include file for make in `/make.arch` (file according to `PDAF_ARCH`)
- For `$PDAF_ARCH=linux_gfortran_openmpi` the specification is
`LINK_LIBS =-L/usr/lib -llapack -lblas -lm`
- If the libraries are at another non-default location, one has to change the directory name (`/usr/lib`)
- Some systems or compilers have special libraries (e.g. MKL for ifort compiler, or ESSL on IBM/AIX)

PDAF needs to be compiled for double precision

- Needs to be set at compiler time in the include file for make:
- For gfortran: `OPT = -O3 -fdefault-real-8`

Files in the tutorial implementation

/tutorial/inputs_online

- true_stepY.txt true state
- state_ini.txt initial estimate (ensemble mean)
- obs_stepY.txt observations
- ens_X.txt initial ensemble members

/tutorial/classical/online_2D_parallelmodel

(after running model_pdaf)

- state_stepY_ana.txt analysis state estimate
- ens_X_stepY_ana.txt analysis ensemble members

X=1,...,9: ensemble member index

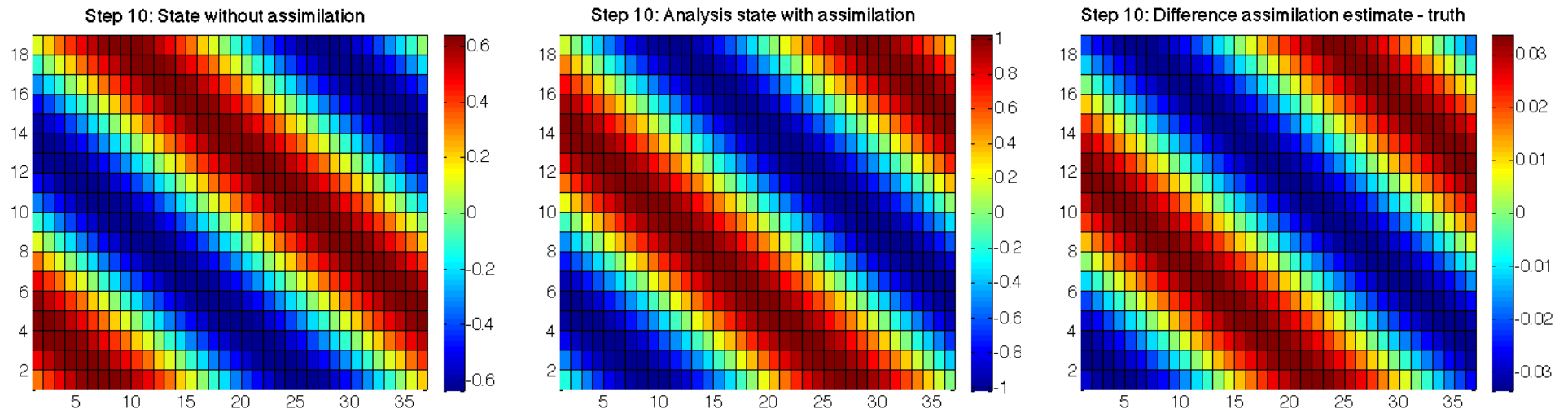
Y=1,...,18: time step index

Note: Files *_for.txt contain forecast fields

Result of the global assimilation

For example, at step 10

- The analysis state (center) is closer to the true field than without assimilation (left)
- Truth and analysis are nearly identical (right)



The analysis step

Next: Implement user-routines for the analysis step

The analysis step needs several user-supplied routines for operations like

- write forecast model fields into state vector
- determine number of available observations
- observation operator acting on a state vector
- initialization of the vector of observations

collect_state_pdaf.F90

Routine to

- Fill state vector with forecasted model fields
- Routine is provided with the state vector `vector_p`

For the example:

1. Access `nx`, `ny` and `field` with use `mod_model`
2. Initialize state vector from model field:

```
DO j = 1, nx_p
  state_p(1+(j-1)*ny : j*ny) = field_p(1:ny, j)
END DO
```

Note: The routine is independent of the filter!

init_dim_obs_pdaf.F90

Routine to

- read observation file
- Count number of observations for **process-local** part of state vector (direct output to PDAF: **dim_obs_p**)

Optional, also

- Initialize array holding **process-local** available observations
- Initialize index array telling index of observation point in **process-local** state vector

The most complicated routine in the example!
(but only about 123 lines)

init_dim_obs_pdaf.F90 (2)

Preparations and reading of observation file:

1. Include `nx`, `ny`, `nx_p` with `use mod_model`
2. declare and allocate real array `obs_field(ny, nx)`
3. Get offset of local part in global state vector
`off_p = Sum over nx_p*ny up to i=mype_filter`
4. read observation file for current time step:

Initialize string 'stepstr' for time step

```
OPEN (12, &  
      file='inputs_online/obs'//stepstr//'.txt', &  
      status='old')  
DO i = 1, ny  
  READ (12, *) obs_field(i, :)  
END DO  
CLOSE (12)
```

init_dim_obs_pdaf.F90 (3)

Count available process-local observations (**dim_obs_p**):

1. Declare integer :: cnt0, cnt_p
2. Now count

```
cnt0 = 0
cnt_p = 0
DO j = 1, nx
  DO i= 1, ny
    cnt0 = cnt0 + 1
    IF (cnt0>off_p .AND. cnt0<=off_p+nx_p*ny) THEN
      IF (obs_field(i,j) > -999.0) cnt_p = cnt_p + 1
    END IF; END DO; END DO
dim_obs_p = cnt_p
```

init_dim_obs_pdaf.F90 (4)

Initialize observation vector (`obs_p`)
and index array (`obs_index_p`):

1. Include `obs_p` and `obs_index_p`
with `use mod_assimilation`
2. Allocate
`obs_p(dim_obs_p), obs_index_p(dim_obs_p)`
(If already allocated, deallocate first)
3. Now initialize ...

Note:

The arrays only contain information about valid observations;
one could store observations already in files in this way.

init_dim_obs_pdaf.F90 (5)

Initialize obs and obs_index

```
cnt0 = cnt_p = cnt0_p = 0    ! Count grid points
DO j = 1, nx
  DO i= 1, ny
    cnt0 = cnt0 + 1
    IF (cnt0>off_p .AND. &
        cnt0<=off_p+nx_p*ny) THEN
      cnt0_p = cnt0_p + 1
      IF (obs_field(i,j) > -999.0) THEN
        cnt_p = cnt_p + 1
        obs_index_p(cnt_p) = cnt0_p    ! Index
        obs_p(cnt_p) = obs_field(i, j) ! observations
      END IF; END IF
    END DO
  END DO
END DO
```

obs_op_pdaf.F90

Implementation of observation operator
acting on some state vector

Input: state vector `state_p`

Output: observed state vector `m_state_p`

1. Include `obs_index_p` by use `mod_assimilation`
2. Select observed grid points from state vector:

```
DO i = 1, dim_obs_p
    m_state_p(i) = state_p(obs_index_p(i))
END DO
```

Note:

`dim_obs_p` is an input argument of the routine

init_obs_pdaf.F90

Fill PDAF's observation vector

Output: vector of observations `observation_p`

1. Include `obs_p` with `use mod_assimilation`
2. Initialize `observation_p`:

```
observation_p = obs_p
```

Note:

This is trivial, because of the preparations in `init_dim_obs_pdaf`!

(However, the operations needed to be separate, because PDAF allocates `observations_p` after the call to `init_dim_obs_pdaf`)

prodrinva_pdaf.F90

Compute the product of the inverse observation error covariance matrix with some other matrix

- Input: Matrix $A_p(\text{dim_obs_p}, \text{rank})$
- Output: Product matrix $C_p(\text{dim_obs_p}, \text{rank})$
(rank is typically $\text{dim_ens}-1$)

1. Declare and initialize inverse observation error variance
`ivariance_obs = 1.0 / rms_obs**2`

2. Compute product:

```
DO j = 1, rank
  DO i = 1, dim_obs_p
    C_p(i, j) = ivariance_obs * A_p(i, j)
  END DO
END DO
```


Done!

The analysis step in online mode with the parallelized model is fully implemented now

The implementation allows you now to use the global filter ESTKF (ETKF and SEIK are usable by adding a call to the corresponding routines PDAF_assimilate_X in assimilate_pdaf)

Not usable are EnKF and SEEK (The EnKF needs some other user files und SEEK a different ensemble initialization)

A complete analysis step

We now have a fully functional analysis step
- if no localization is required!

Possible extensions for a real application:

Adapt routines for

- Multiple model fields
 - Store full fields consecutively in state vector
- Third dimension
 - Extend state vector
- Different observation types
 - Store different types consecutively in observation vector
- Other file type (e.g. binary or NetCDF)
 - Adapt reading/writing routines

Differences between online and offline modes

For the analysis step in online mode:

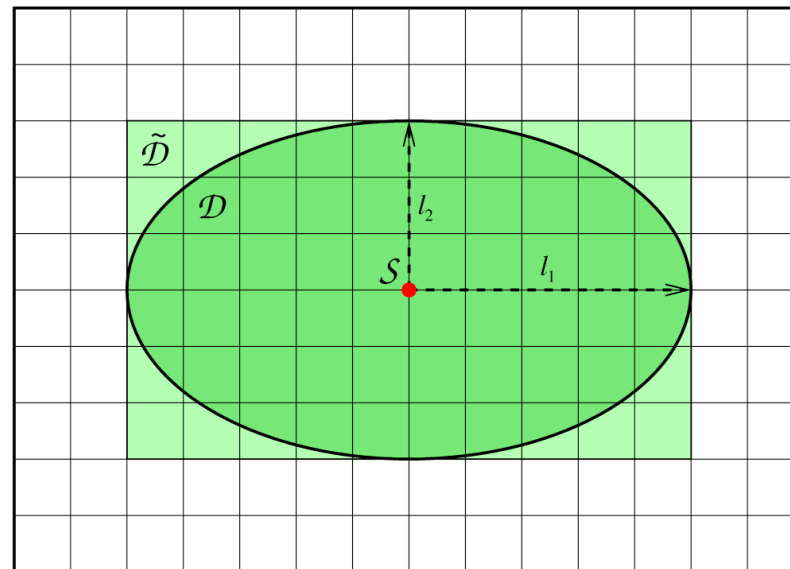
- | | |
|---------------------------------|--|
| <code>collect_state_pdaf</code> | - additional routine for online mode |
| <code>init_dim_obs_pdaf</code> | - read from file for current time step;
include <code>nx</code> , <code>ny</code> from <code>mod_model</code>
instead of <code>mod_assimilate</code> |
| <code>obs_op_pdaf</code> | - identical in online and offline modes |
| <code>init_obs_pdaf</code> | - identical in online and offline modes |
| <code>prodrinva_pdaf</code> | - identical in online and offline modes |

1b) Local filter with parallelized model

Localization

Localization is usually required for high-dimensional systems

- Update small regions (S)
(e.g. single grid points, single vertical columns)
- Consider only observations within cut-off distance (D)
- Weight observations according to distance from S



The FULL observation vector

- A single local analysis at S (single grid point) need observations from domain D
- A loop of local analyses over all S needs all observations
 - This defines the *full* observation vector
- Why distinguish *full* and *all* observations?
 - They can be different in case of parallelization!
- Example:
 - Split domain in left and right halves
 - Some of the analyses in left half need observations from the right side.
 - Depending on localization radius not all observations from the right side might be needed for the left side analyses

Running the tutorial program

- Compile as for the global filter
- Run the program with

```
mpirun -np 18 ./model_pdaf -dim_ens 9 OPTIONS
```

- OPTIONS are always of type `-KEYWORD VALUE`
- Possible OPTIONS are

`-filtertype 7` (select LESTKF if not set in `init_pdaf`)

`-local_range 5.0` (set localization radius, 0.0 by default, any positive value should work)

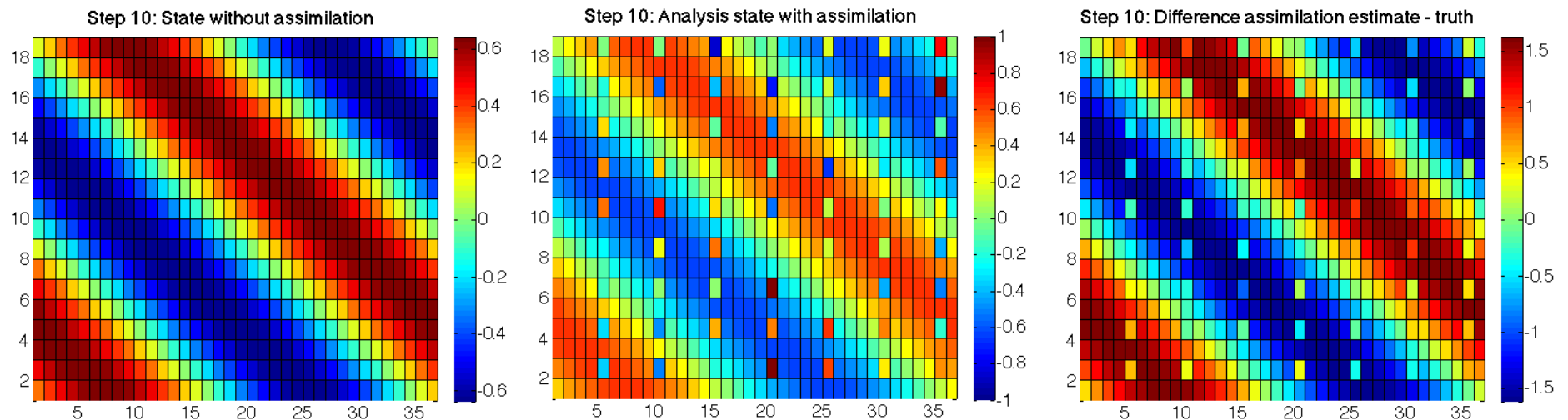
`-locweight 2` (set weight function for localization, default=0 for constant weight of 1; possible are integer values 0 to 4; see `init_pdaf`)

Note: You can run the model e.g. using 18 MPI-processes even on most computers with only 2 processor cores. However, to see a speedup in computing time, you need more physical processors

Result of the local assimilation

```
mpirun -np 9 ./model_pdaf -dim_ens 9 -filtertype 7
```

- Default: zero localization radius (local_range=0.0)
- Change only at observation locations

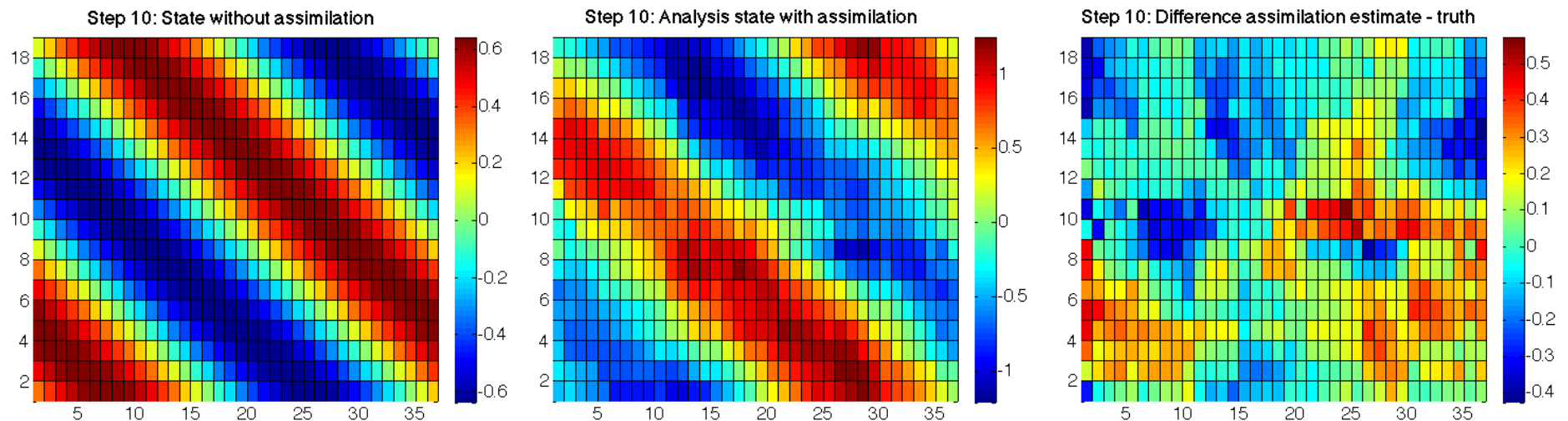


Result of the local assimilation (2)

```
... -filtertype 7 -local_range 10.0
```

- All local analysis domains are influenced (all see observations)
- Up to 16 observations in a single local analysis (average 9.6)

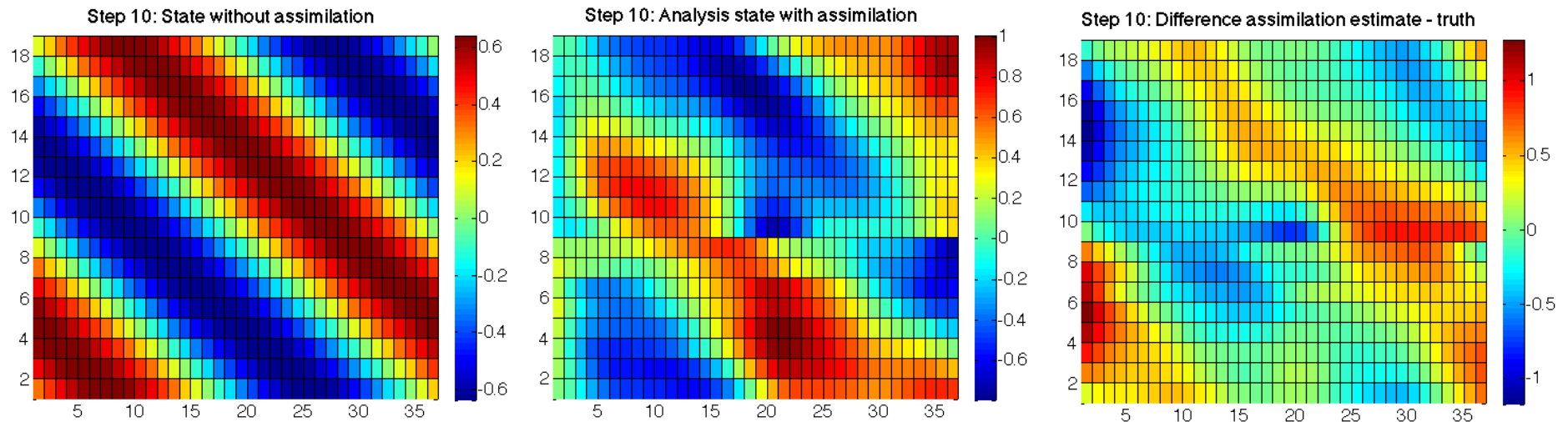
Note: The set up of the experiment favors the global filter because of the shape of the ensemble members



Result of the local assimilation (2)

```
... -filtertype 7 -local_range 10.0 -locweight 2
```

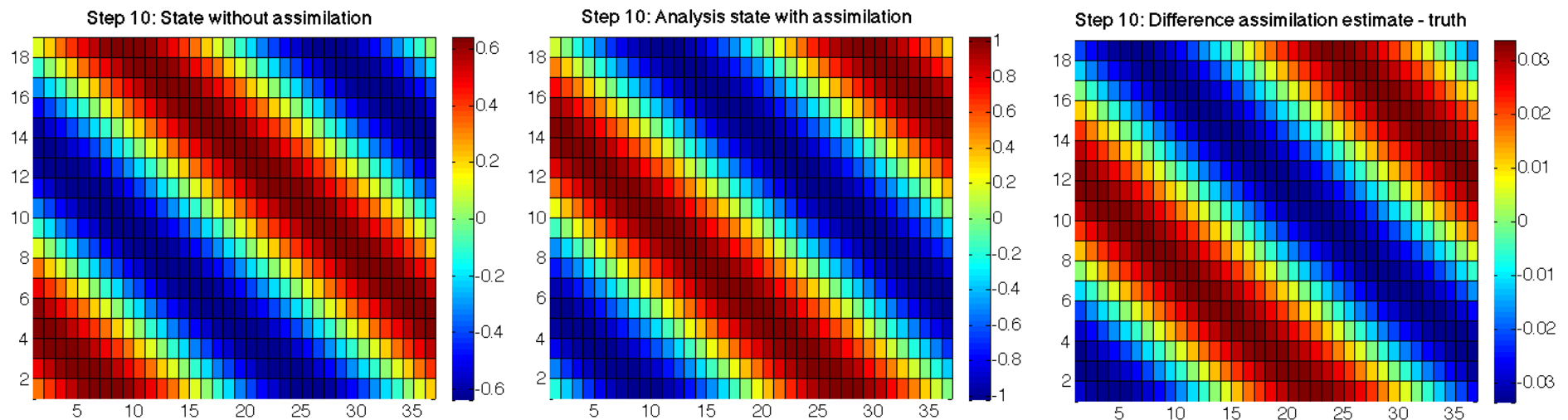
- Observation weighting by 5th-order polynomial
- Analysis field is smoother than before (because of weighting)



Result of the local assimilation (3)

```
... -filtertype 7 -local_range 40.0
```

- Large radius: All local analysis domains see all observations
- Result identical to global filter



Local filter LESTKF

- Localized filters are a variant of the global filters
- User written files for global filter can be widely re-used
- Additional user-written files to handle local part
- No changes to:

```
initialize.F90
```

```
init_ens_pdaf.F90
```

```
prepoststep_ens_pdaf.F90
```

- Change in `init_pdaf.F90`:

```
Set filtertype = 7
```

(You can also set it later on command line)

Local filter LESTKF (2)

Adapt files from global analysis

<code>init_dim_obs_pdaf.F90</code>	→	<code>init_dim_obs_f_pdaf.F90</code>
<code>obs_op_pdaf.F90</code>	→	<code>obs_op_f_pdaf.F90</code>
<code>init_obs_pdaf.F90</code>	→	<code>init_obs_f_pdaf.F90</code>
<code>prodrinva_pdaf.F90</code>	→	<code>prodrinva_l_pdaf</code>

Naming scheme:

`_f_` “full”: operate on all required observations
(without parallelization these are all observations)

`_l_` “local”: operation in local analysis domain or corresponding
local observation domain

Local filter LESTKF (3)

Additional files for local analysis step

`init_n_domains_pdaf.F90`

`init_dim_l_pdaf.F90`

`g2l_state_pdaf.F90`

`l2g_state_pdaf.F90`

`init_dim_obs_l_pdaf.F90`

`g2l_obs_pdaf.F90`

`init_obs_l_pdaf.F90`

localize
state vector

localize
observations

Discuss now the files in the order they are called

init_n_domains_pdaf.F90

Routine to set the number of local analysis domains

Output: `n_domains_p`

For the example: number of process-local grid points (`nx_p * ny`)

To do:

1. Include `nx_p`, `ny` with `use mod_model`
2. Set

```
n_domains_p = nx_p * ny
```

init_dim_obs_f_pdaf.F90

Operations in case of parallelization:

- Read observation file
- Count number of observations for **process-local** part of state vector (`dim_obs_p`)
- Initialize arrays holding **process-local** available observations (`obs_p`) and their coordinates (`coords_obs_p`)
- Initialize index array (`obs_index_p`) telling index of a **process-local** observation in **process-local** state vector
- Initialize **full** number of observations (`dim_obs_f`), vector of observations (`obs_f`), and coordinates (`coords_obs_f`)

“FULL” observation vector:

All observations required for all local analyses in process-local part of state vector (Here: *Full=All* observations for simplicity)

Adapt `init_dim_obs_pdaf` from global filter ...

init_dim_obs_f_pdaf.F90 (2)

Initialize coordinates of process-local observations

For the local filter:

1. Copy functionality from `init_dim_obs_pdaf.F90`
(In the subroutine definition take care that `dim_obs_f` is used instead of `dim_obs_p`)
2. Include `dim_obs_p` with `use mod_assimilation`
3. Initialization of observation coordinates
 - a) Add allocatable REAL arrays `coords_obs_p` and `obs_p` to the routine (don't include `obs_p` from module)
 - b) In the loop where `obs_p` is set add

```
coords_obs_p(1, cnt_p) = REAL (j)
coords_obs_p(2, cnt_p) = REAL (i)
```

init_dim_obs_f_pdaf.F90 (3)

Initialize full quantities (`dim_obs_f`, `obs_f`, `coords_obs_f`)

1. Include allocatable arrays `coords_obs_f` and `obs_f` with `mod_assimilation`
2. Obtain `dim_obs_f` by calling `PDAF_gather_dim_obs_f`
3. Allocate `obs_f` and `coords_obs_f` (deallocate first if already allocated)
4. Obtain `obs_f` by calling `PDAF_gather_obs_f`
5. Obtain `coords_obs_f` by calling `PDAF_gather_obs_f`
6. Add `DEALLOCATE` for `obs_p` and `coords_obs_p`

Note: It is mandatory to call `PDAF_gather_dim_obs_f` once before using the two other functions because it stores dimension information.

Note: The three PDAF functions have been added with PDAF Version 1.13 to avoid that the user implementation needs calls to MPI functions.

Note: `coords_obs_f` has to be a REAL array

obs_op_f_pdaf.F90

Implementation of observation operator
for full observation domain

Difficulty:

- The state vector `state_p` is local to each process
- Full observed vector goes beyond process boundary

Implement two steps:

1. Initialize process-local observed state
2. Gather full observed state vector using MPI

obs_op_f_pdaf.F90 (2)

1. Initialize process-local observed state `m_state_p`

a) Include `dim_obs_p` and `obs_index_p`
with `use mod_assimilation`

b) Declare real allocatable array `m_state_p(:)`

c) Allocate
`m_state_p(dim_obs_p)`

d) Fill the array

```
DO i = 1, dim_obs_p
  m_state_p(i) = state_p(obs_index_p(i))
END DO
```

obs_op_f_pdaf.F90 (3)

2. Get full observed state vector

a) Add variable `INTEGER :: status`

b) Add call to `PDAF_gather_obs_f`:

```
CALL PDAF_gather_obs_f(m_state_p, m_state_f, status)
```

c) Deallocate `m_state_p`

Note: It is mandatory to call `PDAF_gather_dim_obs_f` once before using the two other functions because it stores dimension information. Usually this was already done in `init_dim_obs_f_pdaf`

init_dim_1_pdaf.F90

Set the vector size `dim_1` of the local analysis domain

Further set the coordinates of the local analysis domain and the indices of the elements of the local state vector in the global state vector

Each single grid point is a local analysis domain in the example

1. Set `dim_1 = 1`

init_dim_l_pdaf.F90 (2)

2. Determine coordinates of local analysis domain

1. Compute offset:

`off_p` = Sum over `nx_p*ny` up to `i=mype_filter`

2. Include `coords_l` with use `mod_assimilation`

3. Include `nx`, `ny`, `nx_p` with use `mod_model`

4. Compute `coords_l` from `nx`, `ny`:

```
coords_l(1)
  = real(ceiling(real(domain_p + off_p)/real(ny)))
coords_l(2)
  = real(domain_p + off_p) - (coords_l(1)-1)*ny
```

Note: With parallelization the domain numbering begins with 1 for each process. For the coordinates we also need to count the domains from processes with lower process rank using `off_p`

init_dim_1_pdaf.F90 (3)

3. Set indices of the elements of the local state vector in the global decomposed state vector

a) Include `id_lstate_in_pstate`
with `use mod_assimilation`

b) Allocate `id_lstate_in_pstate(dim_1)`
(Deallocate first if already allocated)

c) Specify the index: It's identical to `domain_p` here
(because we only have a single model variable)

```
id_lstate_in_pstate(1) = domain_p
```


init_dim_obs_l_pdaf.F90

Set the size of the observation vector for the local analysis domain

As for the global filter, this is the longest routine (~102 lines)

Only direct output: **dim_obs_l**

Operations:

1. Include coordinates `coords_l` with `use mod_assimilation`
2. Determine coordinate range for observations
3. Count observations within prescribed localization radius
4. Set index array for local observations (`id_lobs_in_fobs`) and array of distances of local observations (`distance_l`)

Note: The index array in step 4 is re-used for an efficient implementation of `g2l_obs_pdaf`. The local distance array initialized in step 4 is re-used in `prodrinva_l_pdaf` avoiding to recompute distances.

init_dim_obs_l_pdaf.F90 (2)

2. Determine coordinate range for local observations
 1. Declare `real :: limits_x(2), limits_y(2)`
 2. Include `local_range` with `use mod_assimilation`
 3. Set lower and upper limits. E.g. for x-direction

```
limits_x(1) = coords_l(1) - local_range
if (limits_x(1) < 1.0) limits_x(1) = 1.0
limits_x(2) = coords_l(1) + local_range
if (limits_x(2) > real(nx)) limits_x(2) = real(nx)
```

(analogous for y-direction)

Note: Using `limits_x`, `limits_y` is not strictly required, but it makes the search for local observations more efficient.

If the localization is only based on grid point indices, the coordinates could be handled as integer values

init_dim_obs_l_pdaf.F90 (3)

3. Count local observations (within distance `local_range`)

```
dim_obs_l = 0
```

```
DO i = 1, dim_obs_f
```

```
  IF ("coords_obs_f(:,i) within coordinate limits") THEN
```

```
    Compute distance between coords_obs and coords_l
```

```
    IF (distance <= local_range) &
```

```
      dim_obs_l = dim_obs_l + 1
```

```
  END IF
```

```
END DO
```

Note:

For efficiency, we only compute `distance` for observations within coordinate limits `limits_x`, `limits_y`. Valid local observations reside within circle of radius `local_range`.

init_dim_obs_l_pdaf.F90 (4)

4. Set index array for local observations

➤ Index of a local observation in the full observation vector

1. Include `id_obs_in_fobs` and `distance_l`
with `use mod_assimilation`

2. Allocate `id_obs_in_fobs(dim_obs_l)`

3. Fill index array:

```
cnt = 0
```

```
DO i = 1, dim_obs_f
```

```
  IF ("coords_obs(:,i) within coordinate limits") THEN
```

```
    Compute distance between coords_obs and coords_l
```

```
    IF (distance <= local_range) THEN
```

```
      cnt = cnt + 1
```

```
      id_obs_in_fobs(cnt) = i
```

```
      distance_l(cnt) = distance
```

```
  END ...
```

g2l_state_pdaf.F90 & l2g_state_pdaf.F90

g2l_state_pdaf: Initialize state vector for local analysis domain
from global state vector

l2g_state_pdaf: Initialize global state vector
from state vector for local analysis domain

- The templates provide a generic implementation
using the array `id_lstate_in_fstate`

→ We use the templates without any changes!

g2l_obs_pdaf.F90 & init_obs_l_pdaf.F90

g2l_obs_pdaf: Initialize local observed state vector from full observed vector

init_obs_l_pdaf: Initialize local vector of observations

- The templates provide a generic implementation using the array `id_lobs_in_fobs`

→ We use the templates with out any changes!

Note:

`init_obs_l_pdaf` requires that the full observation vector is stored in the array `obs_f`

prodrinva_l_pdaf.F90

Compute the product of the inverse observation error covariance matrix with some other matrix

+ apply observation localization (weighting)

➤ The weighting and the product are fully implemented for a diagonal observation error covariance matrix with constant variance

→ When we re-use the array `distance_1` initialized in `init_dim_obs_1_pdaf`, the template can be used without changes.

Done!

Now, the analysis step for local ESKTF in offline mode is fully implemented.

The implementation allows you now to use the local filter LESTKF (LETKF, LSEIK can be used after adding calls to PDAF_assimilate_X)

Not usable are EnKF and SEEK (PDAF does not have localization for these filters)

For testing one can vary localization parameters:

`local_range` – the localization radius

`locweight` – the weighting method

Default are `local_range=0.0` (observation at single grid point) and `locweight=1` (uniform weight)

2) Hints for adaptations for real models

Implementations for real models

- Tutorial demonstrates implementation for simple model
- You can base your own implementation on the tutorial implementation or the templates provided with PDAF
- Need to adapt most routines, e.g.
 - Specify model-specific state vector and its dimension
 - Adapt `distribute_state` and `collect_state`
 - Adapt routines handling observations
- Further required changes
 - Adapt file output (usually only want to write ensemble mean state in `prepoststep_pdaf`; sometimes possible to use output routines from model)

Multiple fields in state vector

- Tutorial uses a single 2-dimensional field
- All fields that should be updated by the assimilation have to be part of the state vector
- For more fields:
 - concatenate them in the state vector
 - adapt state dimension in `init_pdaf`
 - adapt `init_ens_pdaf`, `collect_state_pdaf`, `distribute_state_pdaf`, `prepoststep_pdaf`
 - For local filters: Adapt full (`_f_`) and local (`_l_`) routines and `g2l_state_pdaf`, `l2g_state_pdaf`, `g2l_obs_pdaf`
- **Note**
 - It can be useful to define a vector storing the offset (position) of each field in the state vector

Multiple observed fields

- In tutorial: observed one field at some grid points
- For several observed fields adapt observation routines:
 - concatenate observed fields in observation vector
 - adapt all observation-handling routines
- **Note**
 - The observation errors can be set differently for each observed field (e.g. using an array `rms_obs`)
 - The localization radius can be set specific for each observed field (observation search in `init_dim_obs_1_pdaf` would use different `local_range` for different fields)
 - One can use spatially varying observation errors using an array `rms_obs` in `prodrinva(_1)_pdaf`

The End!

Tutorial described example implementations

- Online mode of PDAF
- Simple 2D model with parallelization
- Parallelization over ensemble members at the model itself
- Square root filter ESTKF
 - global and with localization
- Extension to more realistic cases possible with limited coding
- Applicable also for large-scale problems

For full documentation of PDAF
and the user-implemented routines
see <http://pdaf.awi.de>