Intensive Course on Data Assimilation Buenos Aires, Argentina 27 October - 7 November 2008

SPEEDY DA (LETKF)

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Local Ensemble Transform Kalman Filter (LETKF, Ott et al., 2004, Hunt et al., 2007)



✓ Yellow: analysis ensemble and its uncertainty; Blue: background ensemble and its uncertainty; Red: observation and its uncertainty

✓ LETKF, like any other EnKF, provides background and analysis uncertainty estimation in every analysis cycle.

Liu (2007)

ETKF equations: Hunt et. al. 2007

$$\mathbf{\bar{w}}^{a} = \mathbf{\tilde{P}}^{a} (\mathbf{Y}^{b})^{T} \mathbf{R}^{-1} (\mathbf{y}^{o} - \mathbf{\bar{y}}^{b}),$$
$$\mathbf{\tilde{P}}^{a} = [(k-1)\mathbf{I} + (\mathbf{Y}^{b})^{T} \mathbf{R}^{-1} \mathbf{Y}^{b}]^{-1}.$$
$$\mathbf{X}^{b} = [\mathbf{x}_{1}^{b} - \mathbf{\bar{x}}^{b} | \dots | \mathbf{x}_{K}^{b} - \mathbf{\bar{x}}^{b}];$$
$$\mathbf{y}_{i}^{b} = H(\mathbf{x}_{i}^{b}); \mathbf{Y}_{n}^{b} = [\mathbf{y}_{1}^{b} - \mathbf{\bar{y}}^{b} | \dots | \mathbf{y}_{K}^{b} - \mathbf{\bar{y}}^{b}]$$

k is the number of ensemble members.

$$\mathbf{\bar{x}}^{a} = \mathbf{\bar{x}}^{b} + \mathbf{X}^{b} \mathbf{\bar{w}}^{a}$$
 The analysis is a weighed average of the background ensemble members.
 $\mathbf{W}^{a} = [(k-1)\tilde{\mathbf{P}}^{a}]^{1/2}$ $\mathbf{X}^{a} = \mathbf{X}^{b}\mathbf{W}^{a}$ $\overset{X^{a}, X^{b}}{\text{where n is the number of model variables.}}$

Local Ensemble Transform Kalman Filter (LETKF, Ott et al., 2004, Hunt et al., 2007)

Schematic of 2-dimension local patch



- Different local volumes have a great overlap.
- Each observation is used more than once in the data assimilation.
- The analysis in each grid point is highly parallel.

Liu (2007)

Analysis localization in the SPEEDY model.

The shape of the localization boxes is square.



nlocal=4 means a total horizontal size of 9 grid points.

Error covariance inflation:

To avoid filter divergence the background error covariance or the analysis error covariance can be "inflated" by different methods.

In this case a multiplicative scheme with fixed parameter is used to inflate the background error covariance.

$$X^{i}_{b} = (x^{i}_{b} - \overline{x}_{b}) \sqrt{(1 + parameter)}$$

The paremeter inflation factor has to be tuned for each particular system. The value for the SPEEDY model is set in the file letkf_tools.f90

IMPLICIT NONE

PRIVATE

PUBLIC :: set_letkf_tools,das_letkf



Localization by R

Inside the local domain the observations that are closer to the domain center will have greater impact than that which are near the domain border.

To achieve this the observation error of every observation within the local domain is increased by a factor which depends on the inverse of the distance to the domain center.



This helps to avoid the undesirable consequences of sampling errors, reducing the influence of observations that are far away from the patch center. This is especially important in the situation where we have sparce observations.

LETKF Response to a single observation



Response experiment:





Once the location is set we must tell the LETKF that the observational increment is going to be fixed. (edit the letkf_tools.f90 file)



Set the "compiler" variable in the scripts letkf_response.sh (for the PC in the lab use the Intel option).

Run the letkf_response.sh:

./letkf_response.sh response



The results will be available in DAS_result/letkf/response

The files analysis.grd and gues.grd contains the analysis ensemble mean and first gues ensemble mean respectively and they can be opened with GrADS.

In this case you will also have a gues_spread.grd file which contains the first gues ensemble spread data and also can be opened with GrADS.

In the folder DAS_result/letkf/response you will also find a grads script (DA2008_response.gs) to plot the results.

Results...

Response in PS for a single PS observation.



Results...

Response in V (z=4) for a single q(z=4) observation.



Run LETKF of rawinsonde observation network



Edit ex_obs.f90 in the letkf folder



You can run the letkf_response.sh to do only one assimilation or run the letkf.sh script to start and assimilation cycle.

To run an assimilation cycle type:

./letkf.sh exp_name

Where exp_name is the experiment name. The results will be stored in a folder with the name of the experiment under DAS_results/letkf.

In the letkf.sh script there is a variable "STORE" that controls the amount of output generated by the assimilation cycle. If it set to 0, then only the first gues mean and the analysis mean in pressure levels will be stored, else first gues ensemble members, analysis and filtered analysis ensemble members will be stored in sigma and pressure levels. (do not set store = 1 in the lab computers since we don't have enough storage capacity)

Directory description:



Zonally averaged RMSE computation:

After the completion of a DA experiment you can use the rmse_zm.sh and rmse_energy.sh scripts to compute the zonally averaged rmse for each variable and for each time.

Under "letkf" type:

>rmse_zm.sh "EXP_NAME"

Where "EXP_NAME" is the name of the experiment. This will create a folder inside the experiment results folder with the name "rmse", inside this folder a GrADS ctl file and a data file will be created.

As in the previous example the script rmse_energy.sh will compute the RMSE for the total energy.

The time control of the RMSE computation (to change the start and end date for the RMSE computation you will have to edit the rmse_zm.f90 and rmse_energy.f90 files.

The rmse_zm.sh and rmse_energy.sh scripts under "tdvar" computes the RMSE for the 3DVAR experiments.

Data Assimilation Cycle Intercomparisson

To compare the results obtained with different configurations you can use two grads scripts located in /home/user/SPEEDY_DA/DAS_results/letkf/gs and in

/home/user/SPEEDY_DA/DAS_results/tdvar/gs

DA2008_ave.gs Computes the time average of the zonally averaged RMSE for a particular variable at a particular level. Can be used to compare multiple experiments.

>DA2008_timec.gs Shows the temporal evolution of the zonally averaged RMSE at a particular level, latitude and for a particular variable.

2002	×I		
	*WHITE exp1='EXP_FULL' *RED exp2='EXP_NOT' *GREEN exp3='EXP_NOUV' *BLUE exp4='EXP_NOQ' *YELLOW exp5='EXP_SUPUPPER' *BROWN exp6='EXP_NOERRSTAT'		Select the experimets that you to compare
	<mark>W</mark> ar='t' level=4 range1=0 range2=8		Select the variable (U, V, T, F and the level (1 to 7), and the start and end values.
	<pre>'open/'exp1'/rmse/rmse_zm.ctl' 'open/'exp2'/rmse/rmse_zm.ctl' 'open/'exp3'/rmse/rmse_zm.ctl' 'open/'exp4'/rmse/rmse_zm.ctl' 'open/'exp5'/rmse/rmse_zm.ctl' 'open/'exp6'/rmse/rmse_zm.ctl'</pre>	The rest of the script opens the selected experiments, compute and plots the time average.	
	'set z ' <mark>level</mark> 'set grads off' 'set t 1 230' 'set lat -90 90'		

experimets that you'd like re

variable (U, V, T, PS, Q) evel (1 to 7), and the Y axis end values.

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* Draw a line plot comparing the RMSE evolution for different experiments at a fixed lattitude.

*WHITE exp1='EXP_FULL' *RED exp2='EXP_NOT' *GREEN exp3='EXP_NOUV' *BLUE exp4='EXP_NOQ' *YELLOW exp5='EXP_SUPUPPER' *BROWN exp6='EXP_NOERRSTAT' var='u' level=4 lat=20 range1=0 range2=15 open .../'exp1'/rmse/rmse_zm.ctl' open .../'exp2'/rmse/rmse_zm.ctl' 'open ___/'exp3'/rmse/rmse_zm.ctl' open .../'exp4'/rmse/rmse_zm.ctl' 'open .../'exp5'/rmse/rmse_zm.ctl' 'open .../'exp6'/rmse/rmse_zm.ctl' 'set z 'level 'set grads off' 'set t 1 230' 'set lat 'lat

Chose the experiments.

Chose the desidered variable, level, and latitude.

Set the Y axis range.

Example 3DVAR: U at level 4 and latitude 20 N.

EXP_FULL (black), EXP_NOT (red), EXP_NOUV (green), EXP_NOQ (blue), EXP_SUPUPPER (yellow), EXP_NOERRSTAT (brown)



Example 3DVAR: T at level 4.

EXP_FULL (black), EXP_NOT (red), EXP_NOUV (green), EXP_NOQ (blue), EXP_SUPUPPER (yellow), EXP_NOERRSTAT (brown)

