

# Singular values of covariance matrices under localization

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We have two ensembles representing the same system. The ensemble  $X$  is a coarse representation of the state, but has many ensemble members. The ensemble  $Z$  is a fine representation of the sates, but has few ensemble members.

$$\dim X = N_x, N_{ex}$$

$$\dim Z = N_z, N_{ez}$$

In order to find the eigenvalues and eigenvectors of the sample correlation of  $X$ , we take the singular value decomposition of  $\tilde{X} = (N_{ex} - 1)^{-1/2}(X - \bar{X})/\text{sd}(X)$ , where  $\text{sd}(X)$  is the standard deviation of each element of  $X$ .

$$U_x S_x V_x^T = \tilde{X}$$

This means that,

$$U_x S_z^2 U_x^T = \hat{C}_x$$

where  $\hat{C}_x$  is the sample correlation of  $X$ .

We want to then use  $U_x$ , or its leading columns, to estimate the leading  $U_z$ . To do this, we interpolate  $U_x$  to the  $z$  space and then use QR factorization to ensure the interpolated  $u_x$ 's are orthonormal:

$$U_x = \text{interp}(U_x)$$

$$U_{xi}, R = \text{QR}(U_{xi}).$$

We then must choose how many of the columns of  $U_{xi}$  should be used. The best way to make this choice is unclear. A few ways that I am considering are:

1. Keep  $U_{xi}$  based on the cumulative sum of their corresponding eigenvalues.
2. Keep  $U_{xi}$  based on the rate of change of their corresponding eigenvalues.
3. Keep  $U_{xi}$  based on some measure of the length scale produced by the low rank approximation of  $\hat{C}_x$  and the residual.
4. Keep  $U_{xi}$  based on how orthogonal the interpolated vector is in  $z$  space before orthogonalization.

It should be noted that keeping all columns can be detrimental to the assimilation process. If some of the columns of  $U_{xi}$  are representing small scale structures, then in the following steps they will still be used to represent some of these small scale structures.

After choosing which columns of  $U_{xi}$  to keep, we must determine what eigenvalues they should have. We do this by taking  $\lambda = u_{xi}^T \hat{C}_z u_{xi}$  as the eigenvalue of  $uxi$  for our approximation of  $C_z$ . Alternatively, we can take  $\lambda = (\tilde{Z}^T u_{xi})^T (\tilde{Z}^T u_{xi})$  where  $\tilde{Z}$  is defined similarly as  $\tilde{X}$ . This will give us the leading eigenvectors and an approximation of their eigenvalues of  $C_Z$ .

These eigenvectors and eigenvalues represent the large scale structure of the problem. To then find the small scales of the problem, we take

$$C_z^\perp = C_z^\parallel - \hat{C}_z$$

where

$$C_z^\parallel = U_{xi} \Lambda_{xi} U_{xi}^T$$

and  $\Lambda_{xi}$  is the diagonal matrix with the  $\lambda$ s described above on the diagonal. We can then localize  $C_z^\perp$  to get at the small scales that are represented in  $Z$ .

We must choose how to localize  $C_z^\perp$ . One reasonable expectation is that the scales in  $C_z^\perp$  will be shorter or similar to  $N_z/N_x$ . This choice will also be affected by the choice of how many columns of  $U_{xi}$  to keep. Once the localization matrix  $L$  is chosen, we can then generate our localized correlation matrix as:

$$C_z^{loc} = C_z^\parallel + L \circ C_z^\perp$$

and the corresponding covariance matrix as:

$$P_z^{loc} = D_z C_z^{loc} D_z.$$

where  $D_z$  is the diagonal matrix with sample standard deviations of  $Z$  on the diagonal.

We can then find the leading eigenvectors and eigenvalues of  $C_z^{loc}$ :

$$Q \Lambda Q^T = C_z^{loc}$$

and use them to transform the  $z$  variable. First, we must calculate a whitening transformation

$$T_w = \Lambda^{-1/2} Q^T$$

and its right inverse

$$T_w^i = Q \Lambda^{1/2}$$

and the singular value decomposition

$$U \Sigma V^T = R^{-1/2} H T_w^i.$$

We can then define transformations for both  $z$  and  $y$

$$T_z = V^T T_w$$

$$T_y = U^T R^{-1/2}$$

We then know  $z^* = T_z z$  and  $y^* = T_y y$  have identity covariance matrices and

$$y^* = \Sigma z^* + \epsilon^*$$

$$U^T R^{-1/2} y = \Sigma V^T \Lambda^{-1/2} Q^T z + \epsilon^*$$

$$y = R^{1/2} U \Sigma V^T \Lambda^{-1/2} Q^T z + \epsilon$$

$$y = R^{1/2} R^{-1/2} H Q \Lambda^{1/2} \Lambda^{-1/2} Q^T z + \epsilon$$

$$y = H z + \epsilon$$

The question is: can I do this better? Specifically:

1. Can I calculate the eigenvectors and values of  $C^{loc}$  without generating  $C^{\parallel}$  or  $C^{\perp}$  explicitly?
2. Can I calculate the eigenvectors and values of  $P$  from those of  $C$  without generating  $C$ ?
3. Can I calculate the singular vectors and values of  $R^{-1/2} H T_w^i$  from the singular value decomposition of  $R$ ,  $H$ , and  $T_w^i$ ?