

INVERSE METHOD FOR THE RETRIEVAL OF OCEAN VERTICAL PROFILES USING SELF ORGANIZING MAPS AND HIDDEN MARKOV MODELS

Application on Ocean Colour Satellite Image Inversion

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Keywords: Self organising maps, Hidden markov models, Inversion, Geophysical, Chlorophyll-A, Satellite imaging, Inversion.

Abstract: This paper presents a statistical inversion method used to infer 3D data from 2D imaging. The methodology is based on a combination of the Self Organising Maps and the Hidden Markov Models. The method has been validated by inferring the oceanic vertical profiles of Chlorophyll-A based on sea-surface data.

1 INTRODUCTION

The density of satellite observations allowed a semi-continuous observation of the global ocean surface. The two-dimensional images provided by this coverage often contain information on integrated quantities whose vertical distribution is unknown. Depending on the field of study there exist different dynamic approaches for inverting this type of data. However, these approaches are often faced with the problem of non-linearity, and can also be hampered by a lack of knowledge of the complete mechanisms that govern the distributions.

The present paper deals with the inversion of observed sea-surface satellite images, noted \tilde{x}_{obs}^t , $t \in [1 \dots T]$, for retrieving of the vertical distribution of Chlorophyll-A, noted \tilde{x}_{dis}^t , $t \in [1 \dots T]$, using a statistical, non-linear approach.

The methodology we have developed is a mixture of the neuronal algorithms known as Self Organizing Topological Maps (SOM) and the Hidden Markov Models (HMM). The SOM are unsupervised classification algorithms, that allow us to cluster our available data into classes. The classes are arranged on a topological map and connected to each other by a topological similarity distance. In the

present study the SOM classification is applied twice, once on the sea-surface data, and once upon the vertical profiles connected to these images. The resulting topological maps allow us to discretize both data spaces into set amounts of classes.

The second statistical algorithm, the HMM allows us to infer the most likely sequence of some discrete, unobservable states, given a series of discrete, observable states. To do so a set of probability matrices are calculated, corresponding to the dynamic processes of the unobserved states and the links existing between the observed and unobserved states. We use the classes created through the SOMs to discretize the available data and therefore represent both the observable and the unobservable states.

In this paper, we present the results obtained with the methodology developed on a case study at the site of the Bermuda Atlantic Time Series (BATS) (32 N -64 W) of the JGOFS campaign.

2 SELF-ORGANIZING TOPOLOGICAL MAPS

Self-Organising Topological Maps (SOM) are

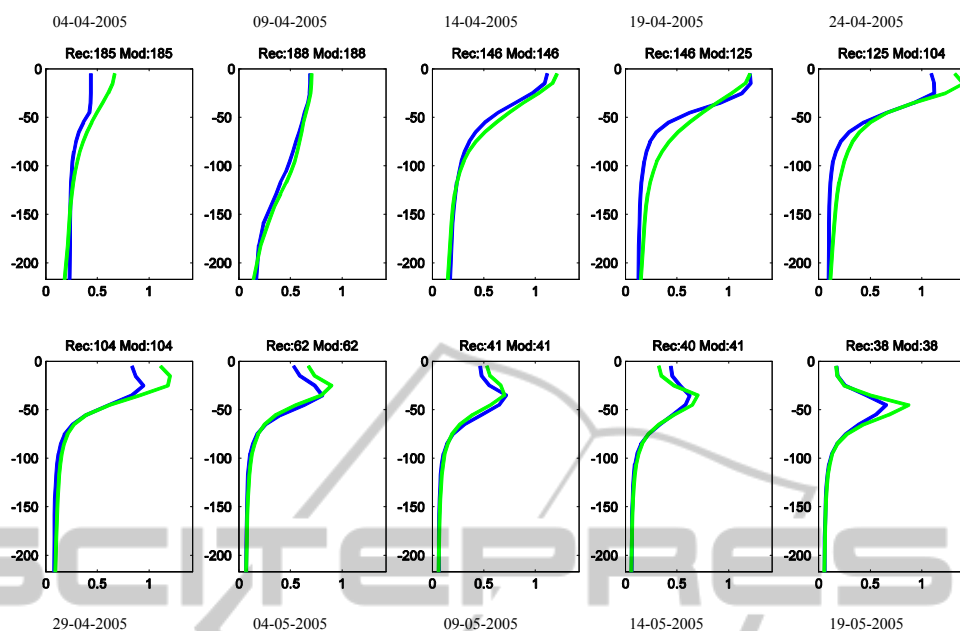


Figure 1: Inversion of ten 5-days steps for the period from 04-04-2005 to 05-19-2005, at BATS. In green, the states provided by the inverse method and in blue the vertical distribution of Chlorophyll-A according to the NEMO-PISCES model. The horizontal axes are in $10^6 \mu\text{mol/L}$ of Chlorophyll-A, while the vertical ones are in meters from the sea surface. The numbers on top correspond, after **Rec**, to the indexes of the classes on M_{dis} attributed to that 5-day step by the inverse method, and, after **Mo** to the indexes attributed by projection of the total profile on M_{dis} . These are not show in the figure.

clustering methods based on neural networks (S, Haykin 1999). They provide a discretization of a learning dataset $A = \{ \tilde{x}_k \in R^p, k = 1 \dots N \}$ into a reduced number of subsets, called classes, $P_i, \{ i = 1 \dots M \}$ that share some common statistical characteristics. Each subset is represented by its referent vector r_i which approaches the mean value of the elements in the class P_i . In our case, we trained two SOMs, one containing the observations, called M_{obs} and one containing the distributions of the unobservable states, called M_{dis} . The number of classes in M_{obs} and M_{dis} are respectively noted N_{obs} and N_{dis} .

The topological aspect of the maps can be justified if we consider the Map as an undirected graph on a two-dimensional lattice whose vertices are the m classes. This graph structure therefore allows the definition of a discrete distance $d(i,j)$ between two classes i and j , defined as the length of the shortest path between i and j on the map. The nature of the SOM training algorithm forces a topological ordering upon the map, and therefore any neighbouring classes c_i and c_j on the map have referent vectors r_i and r_j that are close in the Euclidian sense in the data space R^p . The topological ordering constitutes a major element of our inverse

method, since it allows us to make, latter on, the ergodic assumption for our Markov states.

We define a series of observable events by taking the data from observations related to a given period of time and we label each observation by the index of the class to which it is assigned by using M_{obs} . This classification is done by allocating to each observation $\tilde{x}_{\text{obs}}^t, t \in [1 \dots T]$, in the sequence the index of the class of M_{obs} whose referent is the closest to it in the Euclidian sense. Therefore we obtain the series

$$S_{\text{obs}} = \{ s_{\text{obs}}^t = \underset{i = 1 \dots N_{\text{obs}}}{\text{argmin}_i} (| \tilde{x}_{\text{obs}}^t - r_i^{\text{obs}} |), \quad (1)$$

In the same way, we obtain the time-series of distributions

$$S_{\text{dis}} = \{ s_{\text{dis}}^t = \underset{i = 1 \dots N_{\text{dis}}}{\text{argmin}_i} (| \tilde{x}_{\text{dis}}^t - r_i^{\text{dis}} |), \quad (2)$$

These series S_{obs} and S_{dis} are used by the HMM in order to estimate the probabilistic links that exist between the observable states and the unobservable ones.

The referents r_{dis} of M_{dis} are also used as the vertical profiles in our sequence reconstructions.

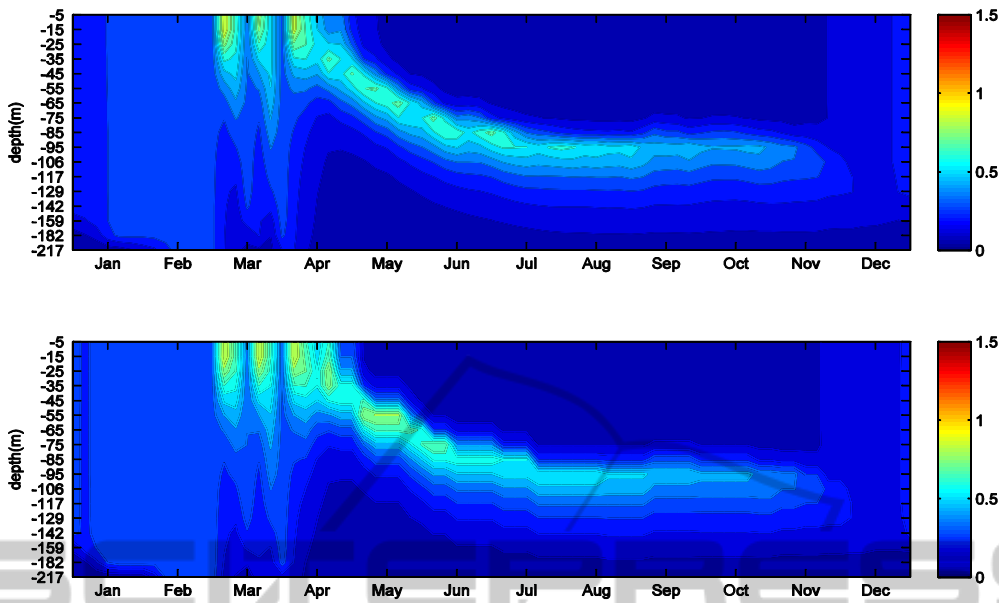


Figure 2: The reconstruction at BATS of the validation year 2005, according to, the NEMO-PISCES MODEL (top graph) and the inverse method result (bottom graph). The colorbar indicates the Chlorophyll-A concentration in $10^6 \mu\text{mol/L}$.

3 HIDDEN MARKOV MODELS

A Markov model is a stochastic model that assumes the first order Markovian property, meaning that each consecutive state of the model depends solely on its previous stat of the model such as

$$P(X_t | X_1 X_2 \dots X_{t-1}) = P(X_t | X_{t-1}) \quad (3)$$

Expanding this principle, a Hidden Markov Model (HMM) is a stochastic model with two sequences. One sequence of unobservable states that follow the first order Markovian property, (represented in our method by S_{dis}), and one sequence of observable states, (represented by S_{obs}), that have a statistical link with the unobservable states (O. Cappé et al., 2005).

We consider two phases, a training one, and a retrieval one. During the training, the Transitions matrix Tr and the Emissions matrix Em are estimated. Tr contains the transition probabilities of the unobserved states

$$tr_{i,j} = P(C_{dis(i,t)} | C_{dis(j,t-1)}) \quad (4)$$

where

$$\sum_{i=1}^{N_{dis}} t_{i,j} = 1 \quad (5)$$

Tr corresponds, in a physical sense, to the underlying dynamics that govern the unobserved states.

Em contains the *à posteriori* probabilities of each observed state to have been emitted by an unobserved state,

$$e_{i,j} = P(C_{dis(i,t)} | C_{obs(j,t)}) \quad (6)$$

where

$$\sum_{j=1}^{N_{dis}} e_{i,j} = 1 \quad (7)$$

Em corresponds, in a physical sense, to the link existing between the observed quantities and the dynamics of the unobserved quantities. Another probability matrix that needs to be calculated is the initial probability matrix Π , with components π_i which represent the average revisit rate of each unobserved state given an infinite sequence. All mentioned probabilities are estimated by using the Baum-Welch algorithm (L. E. Baum et al., 1970), which is a maximum likelihood optimization algorithm, that takes as input the sequences S_{obs} and S_{dis} and outputs the most likely matrices to have generated them through a hidden Markov process.

During the recognition phase, we used the Viterbi algorithm, which is a well-know dynamic programming algorithm (Viterbi AJ, 1967), for inferring the most likely sequence of indexes $S_{dis-est}$ representing the unobserved states, given the previously estimated parameters Tr , Em and Π of the HMM and a sequence of observations $S_{obs-new}$. It is well documented (M.S. Ryan and G.R. Nudd. 1993) that the Viterbi algorithm can face problems due to

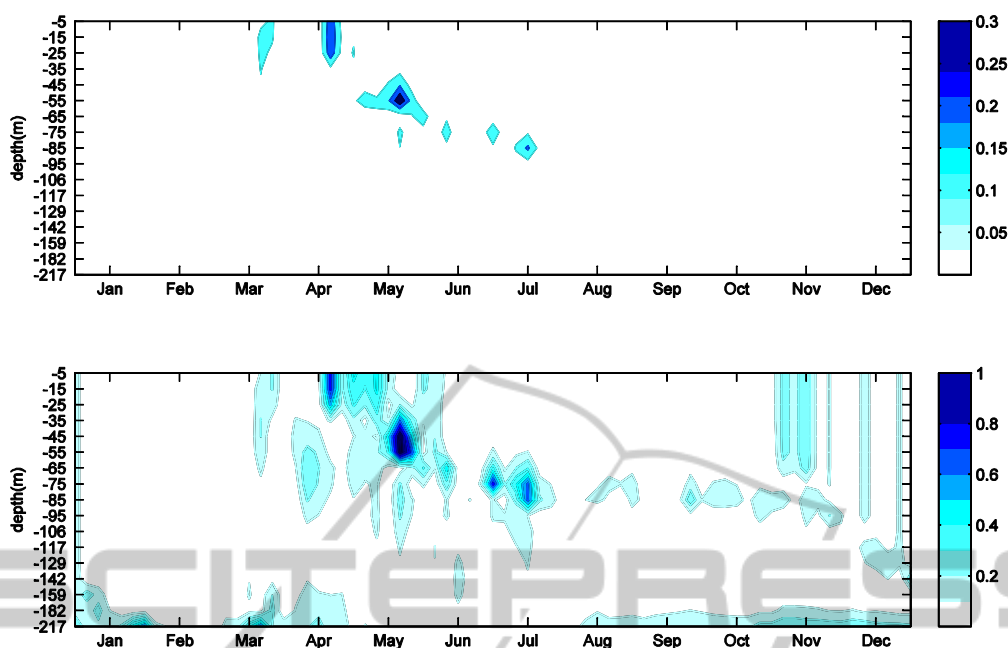


Figure 3: The top image contains the absolute error between the NEMO-PISCES model and the result of the inverse method, at BATS for the validation year 2005. The bottom image contains the absolute relative error between the NEMO-PISCES model and the result of the inverse method.

transitions that were not observed in the training data set. A balance needs to be found between the sizes of the SOM maps that will determine the amount of discretization provided by the method, and the correctness of the allocation of indices. The dimensions of each map are therefore optimized using a validation set. Yet, even with an optimization there will be some situations and transitions that are seldomly encountered in the training data and result in null probabilities in the probability matrices Em_{B-W} and Tr_{B-W} that we estimated in the first pass of the Baum-Welch algorithm.

Due to this usual lack of sufficient data in the concerned domains, Em_{B-W} and Tr_{B-W} need to be adjusted. This is done by taking into account the properties of the topological maps. A major characteristic of the present method is to use the topological order in order to improve the accuracy of the estimated probabilities matrices. The topological maps allow us to modify the probabilities by allowing each state to communicate via a diminutive probability with each of its neighbouring states.

This is done by considering the neighbourhood matrices NM_{obs} and NM_{dis} , of dimensions (N_{obs}, N_{obs}) and (N_{dis}, N_{dis}) , where

$$NM_{SOM}(i,j) = \begin{cases} 1, & \text{if } d(c_i^{SOM}, c_j^{SOM}) < 2 \\ 0, & \text{else} \end{cases} \quad (8)$$

with $d(i,j)$ being the discrete distance of the map. Taking into account the neighbourhood consists in increasing the probability of reaching a class j from a class i , by an amount proportional to the sum of the previously calculated probabilities of reaching the neighbour classes of class j on SOM. In order to favour the data observed during training, we add a weighting term, noted w_c , to the initial probabilities, and we further multiply it by the total length of the training sequences used in the initial Baum-Welch algorithm's pass, noted $T_{training}$, since this length is a measure of confidence in the correctness of the estimated parameters. The matrices obtained are then normalized. The final Em and Tr matrices we use, noted Em_{final} and Tr_{final} , are computed by applying:

$$Em_{final}(i,j) = \frac{w_c * T_{training} * Em_{B-W}(i,j) + \sum_{k=1}^{N_{obs}} (NM_{obs}(i,k) * Em(i,k))}{Em(i,k) + 1} \quad (9)$$

Which is normalized to fit the constraint (7), and

$$Tr_{final}(i,j) = \frac{w_c * T_{training} * Tr_{B-W}(i,j) + \sum_{k=1}^{N_{dis}} (NM_{dis}(i,k) * Tr(i,k))}{Em(i,k)} \quad (10)$$

Which is normalized to fit the constraint (5). For this application w_c is set to 9.

These modifications permit the Viterbi algorithm to circumvent the problems of impossible transitions, or emissions due to insufficient data in the training sequences that resulted in null probabilities in the estimated parameters.

4 APPLICATION FOR THE RESTITUTION OF THE VERTICAL CHLOROPHYLL-A CONCENTRATION THROUGH SEA SURFACE DATA

The bio-geochemical activity of the oceans and the carbon cycle are two parts of a complex feedback system. A change in climate and an increase of the amount of available carbon can affect the primary oceanic production, and in return a change in the bio-geochemical activity affects, by modifying the albedo and carbon fixation rates, the climate and carbon concentration. It is therefore important to be able to determine the oceanic primary production. In recent years, many algorithms have been developed that infer the Chlorophyll-A concentration in ocean surface layers through satellite imaging (Brajard et al., 2008). It has also been proved that the vertical Chlorophyll-A distribution, is correlated with sea surface data (Uitz et al., 2006). Therefore, the determination of the vertical distribution of Chlorophyll-A from sea surface data is a problem that can be solved by the methodology we propose.

One cannot determine the vertical distribution of Chlorophyll-A without first understanding the parameters that influence the development of phytoplankton. It is generally accepted that phytoplankton growth depends on 5 parameters: available radiation, available nutrients, predators and biology, water temperature, water turbidity.

These parameters cannot easily be monitored through a direct approach. Satellite imaging, however, can give us proxy information, which can be used in an empirical approach for determining the vertical distribution of Chlorophyll-A. Specifically, in this study we used: Sea Surface Chlorophyll-A concentration (SCHL), Sea Surface Temperature (SST), Sea Surface Elevation (SSH), Shortwave Radiation (SR) and Wind-speed Intensity (WS).

Since our objective is to validate the theoretical methodology, we used simulated data in order to test the validity of our approach. We therefore approximated the satellite values of the previous

parameters by using the input and output values provided by the NEMO oceanic circulation model coupled to the PISCES bio-geochemical model (C. Moulin, 2008). In order to better simulate the noise and errors inherent to satellite images we added a white noise $z \sim N(0, \hat{\epsilon})$, $\hat{\epsilon} = 1/2 * (\sigma_{schl}, \sigma_{sst}, \sigma_{ssh}, \sigma_{ws}, \sigma_{sr})$ to the parameters that could be gathered from satellite imaging. σ represents the standard derivation of each corresponding surface parameter, as computed on the training data. The application was set at the site of BATS.

The unobserved states that were classified, were the output data vectors containing the average vertical Chlorophyll-A distribution at 17 depth levels (from 5 meters to 217) and temperature distribution at 9 depth levels. These vertical distribution profiles were 5-day averages spanning the period from 1991 to 2007 located in a $2^\circ \times 2^\circ$ square centred on BATS. Therefore M_{dis} belongs to R^{26} (17 levels of Chlorophyll-A + 9 levels of Temperature). M_{obs} belongs to R^5 .

We trained M_{obs} and M_{dis} by taking into account all available profiles at BATS, as well as any adjacent points included in the model. This gave us $9*73*17=11169$ profiles for the construction of the maps. The optimum map sizes, N_{obs} and N_{dis} were determined to be $21*14=294$ classes.

For the estimation of the HHM parameters on the other hand, we take a total of 14 years (1991-2004) for the training, each including seventy-three 5-day steps. Therefore $T_{training}=1022$. We maintained 3 years (2005-2007), or 219 5-day steps, to validate our approach.

The results shown in Figure 1 present the temporary evolution of Chlorophyll-A profiles in ten 5-day steps sequence, from 04-04-2005 to 19-05-2005, at BATS. In green we see the Chlorophyll-A distribution profiles, taken from the referents r_{dis} of M_{dis} , corresponding to the indexes of the reconstructed time series $S_{dis-rec}$. In blue we can see the vertical distribution of Chlorophyll-A according to the NEMO-PISCES model at the same 5-day steps.

In Figure 1 we also have, preceded by the acronym **Rec**, the indexes that constitute time series $S_{dis-rec}$ and, preceded by **Mod**, the indexes we obtain by projecting the corresponding profiles of the NEMO-PISCES model on M_{dis} . In order to avoid confusion, the profiles corresponding to the indexes after **Mod** are not displayed. When the vertical distribution of Chlorophyll-A is known, these indexes would correspond to the optimum reconstruction we could get with M_{dis} . We note this optimum time series as $S_{dis-opt}$. It is interesting to notice that even

when the indexes are not equal, the classes are neighbours on M_{dis} , and the estimated profiles are quite similar to the observed ones.

If we define $S_{dis-2005to2007}$ as the reconstructed time series of indexes of the validation years from 2005 to 2007 and as $S_{dis-opt-2005to2007}$ the corresponding optimum reconstruction we observe that they are in agreement 84,59% of the time. This performance reaches 88,58% when applied on the reconstruction $S_{dis-2005}$ of the year 2005 alone, as compared to its optimum reconstruction. This was probably due to the validation year 2005 having a small variation from the mean year, and presenting often-observed transitions. In the training date we had on average an agreement of 86,46%

In Figures 2 and 3, we applied the inverse method to the full 73 5-day steps serie of the validation year 2005. We can observe that the reconstruction closely fits the results provided by the NEMO-PISCES model. The correlation index of the two images in Figure 2 is 97,30%.

We can notice that the discretization induced by the SOM is apparent, yet the general form and intensity are correctly represented, as it becomes clear in Figure 3, where the error graphs tend to have small values.

5 CONCLUSIONS

In the present paper we have introduced an inversion method based on SOM and HMM, that is able to reconstruct the vertical profiles of Chlorophyll-A based on satellite images. One of its main advantages in the inversion of Chlorophyll-A is that it does not use Gaussian approxiamtions used in other methods (Morel 1988, Uitz et al., 2006), allowing the reconstruction of situations where the distribution is not conforming to a single gaussian curve. An additional benefit of the inversion method presented, is its efficiency in terms of calculations. The method is open ended enough to be applicable for the inversion of the profiles of different biogeophysical parameters based on satellite imaging.

We plan to further validate this method by testing its robustness with satellite imaging and in-situ data, as well as to apply it on different types of profiles, such as oceanic salinity or temperature profiles. A latter goal is to expand the method to take spatial constraints into consideration, and reconstruct 3D profiles.

ACKNOWLEDGEMENTS

We would like to thank the "Délégation Generale de l'Armement" for financing this work.

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