

MOCCA

a Multi-sourced Ocean Carbonate Chemistry Analysis

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1 Introduction

Deep learning has been heavily used in climate science in recent years (Reichstein et al., 2019), with applications ranging from forecasting, statistical down-scaling, pattern identification, process parameterization, emulation of physical models, to data interpolation. The success of deep learning in such data-driven applications is based on the versatility of neural networks Goodfellow et al. (2016). In theory, these can be used to learn any functional relationship between predictors and variables of interest, as reflected by the universal approximation theorem Hornik et al. (1989).

A particularly important application of deep learning is the interpolation of sparse ship-based measurements of the fugacity of carbon dioxide (fCO₂) in the surface ocean. A globally consistent and complete field of fCO₂ is necessary to estimate the air-sea CO₂ flux and thus necessary to estimate the fraction of anthropogenic carbon emissions that is taken up by the ocean. The global carbon budget (Friedlingstein et al., 2023), an annual assessment of global carbon emissions and sinks, currently estimates the oceanic sink from seven observation-based products. While based on very similar underlying fCO₂ data

and similar mostly satellite-based predictors, these seven products follow different approaches for interpolating the sparse fCO_2 data.

Four of these products are based on feed-forward neural networks: The CMEMS-LSCE-FFNNv2 (Chau et al., 2022) utilizes a 100-member neural network ensemble, bootstrapping from the months before and after a fCO_2 measurements and leaving the months with fCO_2 measurements for independent evaluation. The MPI-SOMFFN (Landschützer et al., 2016) builds on a two step procedure, where first different clusters of similar ocean conditions are determined using a self-organizing map approach and then neural networks are trained to predict fCO_2 in each cluster separately. Similarly, OS-ETHZ-GRaCER (Gregor and Gruber, 2021) provides an ensemble of varying cluster assignments with neural-network-based fCO_2 regression in each cluster. NIES-ML3 (Zeng et al., 2022) is based on three model estimates, from a random forest, a gradient boost machine, and a feed-forward neural network. The remaining three observation-based products build on multiple linear regressions for A_T and C_T (fundamental variables to calculate fCO_2 and other carbonate system variables; JMA-MLR; Iida et al., 2021), extreme gradient boosting to predict the missfit between global ocean biogeochemical models and fCO_2 measurements (LDEO-HPD; Gloege et al., 2022), and a autoregressive multiple linear regression approach (Jena-MLS; Rödenbeck et al., 2022).

The largest uncertainty in these spatially and temporally interpolated fCO_2 fields roots in the sparsity and uneven distribution of the underlying fCO_2 measurements that are collected in the Surface Ocean CO_2 Atlas (Bakker et al., 2016). In particular, measurements are sparse in high-latitude regions and particularly in the Southern Ocean. One approach to soften this issue is applying neural-network based regression separately in clusters with similar ocean-biogeochemical conditions (Landschützer et al., 2016; Gregor and Gruber, 2021), grouping data-sparse regions with others with similar conditions. MOCCA, instead, attempts to tackle these limitations by not only using fCO_2 measurements for training, but also pH measurements from biogeochemical Argo floats (Johnsson et al., 2017), which provide critical additional data in the Southern Ocean, and other biogeochemical data as provided by the Global Ocean Data Analysis Project (GLODAPv2; Olsen et al., 2016). Furthermore, the neural network in MOCCA is pretrained on CMIP6 Earth system models to foster inference of surface fCO_2 in sparsely sampled ocean regions with the prior knowledge about functional relationships from Earth system models. The use of multiple sources of data is enabled by a flexible model structure that allows to train on multiple target variables. Similarly to Iida et al. (2021), this flexible approach also provides coherent estimates for all variables of the oceanic carbonate system.

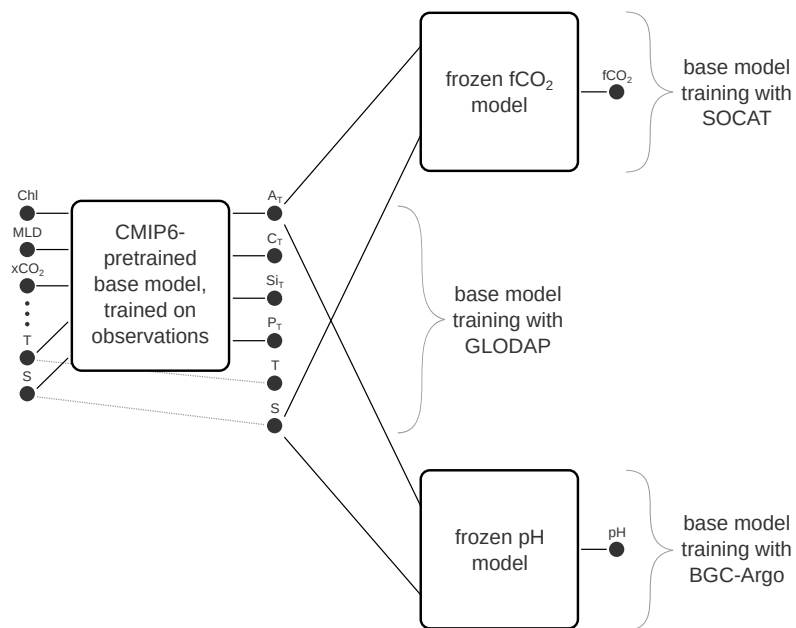


Figure 1: Scheme displaying the overall model architecture of MOCCA.

2 Model architecture

2.1 Overview

MOCCA is based on a three-step procedure (Figure 2). In a first step, surrogate neural network models are trained to fit the functional relationship between a set of variables needed to solve the oceanic carbonate system (*carbonate chemistry drivers*) and two specific carbonate chemistry variables, pH and fugacity of CO_2 (Section 2.2). The functional relationship is learned from the numerical carbonate chemistry package *mocsy 2.0* (Orr and Epitalon, 2015). This step is performed first to identify appropriate hyper parameters and training procedure in a situation where near-perfect learning is possible.

In a second step, a neural network is trained on model output from CMIP6 Earth system models to learn the statistical relationships between a set of predictors and the carbonate chemistry drivers. This model will serve as a prior estimate to improve inference for fCO_2 and pH in undersampled regions where only few fCO_2 and pH measurements are available.

In a third step, the CMIP6-pretrained model from the previous step is trained with observational data for fCO_2 from the Surface Ocean CO_2 Atlas (SOCAT), carbonate chemistry drivers from GLODAP, and pH from BGC-Argo floats. To do so, the loss is either directly calculated from the output of the base model (GLODAP), or the fCO_2 or pH models are appended to the base model (for SOCAT and BGC-Argo respectively). In the latter case, the layers of the fCO_2 or pH models are frozen to prevent further parameter updates.

2.2 Surrogate models for mocsy fCO_2 and pH

In a first step, surrogate multi-layer perceptron models were trained to replace numerical solution of the oceanic carbonate system. To do so, samples for total alkalinity (A_T), dissolved inorganic carbon (C_T), temperature (T), salinity (S), total silicate (Si_T), and total phosphate (P_T) were randomly generated from uniform distributions (Table 1).

The training data size was set to 5 000 000 samples and the validation data size was set to 1 000 000 samples. Mocsy 2.0 (Orr et al., 2015a) was then used to calculate fCO_2 and pH for these samples. After normalizing per feature with the means and standard deviations given in Table 1, multilayer perceptron models with three identical hidden layers were trained with mocsy fCO_2 and pH as labels.

For fCO_2 , model complexity was iteratively increased until a desired maximum deviation of less than $1 \mu\text{atm}$ was reached (the measurement uncertainty for pCO_2 as reported by Orr et al, 2015b)¹. During training, learning rate was decreased from 10^{-3} to 10^{-5} following an exponential learning rate schedule over 10 000 epochs. The decay in learning rate was chosen to shift from an initial identification of an optimal region in the parameter space to finding an

¹For reference, fCO_2 across the generated samples varies between $0.001 \mu\text{atm}$ and $5108 \mu\text{atm}$.

	minimum	maximum
A_T	$1000 \mu\text{mol kg}^{-1}$	$3000 \mu\text{mol kg}^{-1}$
C_T	$1000 \mu\text{mol kg}^{-1}$	A_T
T	-2°C	35°C
S	10 PSU	50 PSU
Si_T	$0 \mu\text{mol kg}^{-1}$	$134 \mu\text{mol kg}^{-1}$
P_T	$0 \mu\text{mol kg}^{-1}$	$4 \mu\text{mol kg}^{-1}$

Table 1: Minima and maxima of the uniform distributions used to generate samples for the fCO_2 and pH surrogate models. The range for A_T was chosen such that it easily encompasses open ocean variations in A_T . That for C_T is limited to values lower A_T since larger C_T do not occur in the ocean. The ranges for T and S are chosen according to Lueker et al., 2000, whose parameterizations for K_1 and K_2 are used in mocsy. Finally, the maximum values for Si_T and P_T were chosen to be the global maxima found in the monthly climatologies for Si_T and P_T from World Ocean Atlas 2023. These uniform distributions have means $(\text{max} + \text{min})/2$ and standard deviations $(\text{max} - \text{min})/\sqrt{12}$, except for C_T where mean and standard deviation are given by $\text{min} + (\text{max} - \text{min})/4$ and $(\text{max} - \text{min}) \cdot \sqrt{7/144}$, respectively. These means and standard deviations are used for feature normalization.

optimal set of parameters for which the mean squared error over the training and validation sets converges to a similar and low value.

Hidden layer size was increased from an initial 64 hidden layer units (8833 trainable parameters), 80 units (13601 parameters), 96 units (19393 parameters), 128 units (34049 parameters), to 160 units (52801 parameters). The largest model² hit a maximum deviation of $1.08 \mu\text{atm}$ over the validation set (Figure 2). The same model architecture was then also used to train the pH model, resulting in a maximum deviation over the validation set of 0.0044.

With root mean squared errors (RMSE) of $0.026 \mu\text{atm}$ (fCO_2 model) and 0.00008 (pH model), these surrogate models provide a precision that is comparable to numerical carbonate chemistry packages: Orr et al., 2015b report a desired numerical uncertainty of $0.1 \mu\text{atm}$ and 0.0003, respectively.

The accuracy of the fCO_2 and pH neural network models is similar for the six million random samples for A_T , C_T , T , S , Si_T , and P_T from the CMIP6 models (section 2.3). Specifically, RMSE is $0.029 \mu\text{atm}$ and 0.00006, respectively, and the maximum deviations are $0.24 \mu\text{atm}$ and 0.0006, respectively.

²To give some context about the model complexity: The number of paramters of this model is comparable to that of a Taylor expansion of a function with six arguments to 15th order. Assuming an efficient use of the MLP parameters, a good fit to the numerical solution from mocsy is thus expected.

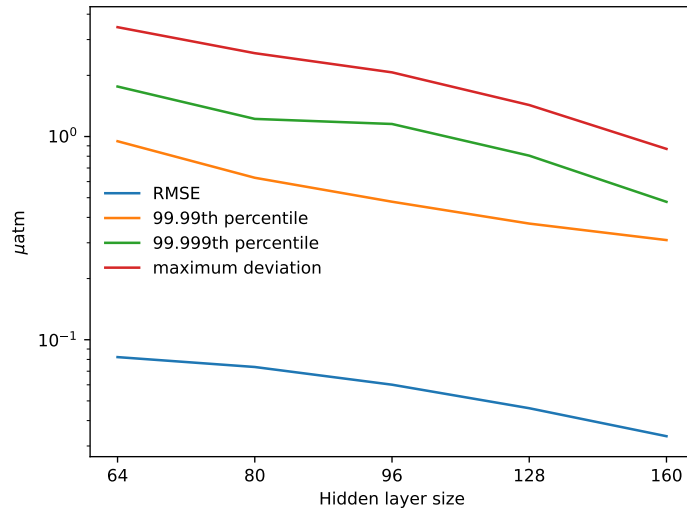


Figure 2: The evolution of root mean squared error (RMSE), the 99.99th and 99.999th percentiles of deviation and the maximum deviation between the surrogate model and mocsy fCO_2 over the validation data set with 1000 000 randomly generated samples.

2.3 CMIP6-pretrained base model

As a first step, we pre-train a neural network on the statistical relationships between predictor variables and sea surface A_T , C_T , Si_T , and P_T . The predictor variables chosen here are sea surface temperature (SST), sea surface salinity (SSS), mixed layer depth (MLD), sea surface height (SSH), chlorophyll-a concentration (chl), sea ice concentration (ice), 10 m easterly wind (u), 10 m northerly wind (v), dry volume molar ratio of CO_2 in the atmosphere (CO_2), sinus and cosinus of the month of year ($\sin(m/12 \cdot 2\pi)$ and $\cos(m/12 \cdot 2\pi)$), as well as latitude and longitude encoded using spherical coordinates following Gade (2010). The representations for month of year and location are used to avoid discontinuities in the months between December and January and in longitude at the dateline ($\pm 180^\circ E$).

The neural network is trained on 6 million samples (5 million for training and 1 million for validation) that were evenly drawn from three CMIP6 Earth system models: the UKESM1-0-LL, the MPI-ESM1-2-LR, and the CMCC-ESM2. Despite a high-resolution version of the MPI model, these were the only ones to provide all predictors on monthly-mean resolution. Data from multiple Earth system models were used to ensure that the model learns general relationships that are not specific to a certain Earth system model with respective biases. The model data is taken from the period 1993-2022 (*historical* simulation until 2014, followed by *SSP2-4.5*). The predictors MLD and chl were log-transformed prior to training to foster learning. The log transformation was applied only where the resulting distribution was more Gaussian (where the test statistic of a Kolmogorov-Smirnov test, i.e. the maximum difference between the empirical distribution function of the data and the distribution function of a standard normal distribution, was smaller after transformation). Finally the predictors for the CMIP6 pre-trained base model are normalized to zero mean and unit variance and the 4 labels are normalized as specified in Table 1 (to match the normalization used for the fCO_2 and pH surrogate models).

For the training, the same architecture that was already used to train the surrogate models (despite not applying a final ELU non-linearity on the output layer, such that the last hidden layer is linearly projected) is used, again with 10 000 epochs and a batch size of 1000. Root mean squared errors of $6.1 \mu\text{mol kg}^{-1}$ (A_T), $6.3 \mu\text{mol kg}^{-1}$ (C_T), $0.8 \mu\text{mol kg}^{-1}$ (Si_T), and $0.03 \mu\text{mol kg}^{-1}$ (P_T) are obtained (on the validation set and after backtransforming to unnormalized labels). As such, RMSEs are more than a magnitude lower than the standard deviations of A_T , C_T , Si_T , and P_T in the CMIP6 model data samples, given by 111.7, 99.1, 26.3, and $0.5 \mu\text{mol kg}^{-1}$, respectively³. Likely due to the large size of the training data set relative to the model complexity, overfitting appears not to be a problem. The validation loss continuously decreases, ending up 11 % larger than the training loss⁴.

³RMSE equals the standard deviation for a model that always predicts the mean of a variable. Hence, a lower RMSE implies skill in predicting spatial and temporal variations in the variable.

⁴Technically speaking, the training data loss is calculated for all batches in an epoch

Predictor	Data product
Sea surface temperature	ESA SST CCI and C3S ¹
Sea ice fraction	ESA SST CCI and C3S ¹
Sea surface salinity	CMEMS ARMOR3D L4 ²
Mixed layer depth	CMEMS ARMOR3D L4 ²
Sea surface height	CMEMS L4 Sea Surface Heights ³
Chlorophyll-a concentration	Copernicus-GlobColour ⁴
CO ₂ mole fraction	Manua Loa Hawaii in-situ data ⁵
Eastward near-surface wind	ECMWF Reanalysis v5 (ERA5) ⁶
Northward near-surface wind	ECMWF Reanalysis v5 (ERA5) ⁶

Table 2: List of predictors used to predict A_T , C_T , Si_T , and P_T and the data product used for each predictor. ¹<https://doi.org/10.48670/moi-00169>; only available until Oct. 2022, ²<https://doi.org/10.48670/moi-00052>, ³<https://doi.org/10.48670/moi-00148>, ⁴<https://doi.org/10.48670/moi-00281>; only available from Sept. 1997 on, ⁵https://scrippsco2.ucsd.edu/data/atmospheric_co2/primary_mlo_co2_record.html, ⁶<https://www.ecmwf.int/en/forecasts/dataset/ecmwf-reanalysis-v5>

A part of this error can be explained by the fact that the neural network mapping is necessarily imperfect since the three models imply different statistical relationships between predictors and labels. Creating 6 million samples only from the UKESM1-0-LL model, the root mean squared errors become significantly smaller, now being $4.0 \mu\text{mol kg}^{-1}$ for A_T , $4.2 \mu\text{mol kg}^{-1}$ for C_T , $0.5 \mu\text{mol kg}^{-1}$ for Si_T , and $0.02 \mu\text{mol kg}^{-1}$ for P_T . The remaining errors should be mainly due to insufficient information in the predictor variables to fully explain the variations in the four concentrations. As such, it can be assumed that no observation-based regression based on these predictors will perform significantly better. Adding further predictors may still enhance the skill of a model. However, an experiment with total cloud cover and precipitation as additional atmospheric predictors resulted only in marginal improvements of errors.

2.4 SOCAT-, GLODAP-, and BGC-Argo-based model tuning

The data products used for each predictor are listed in Table 2.

separately and then averaged. It is thus calculated slightly different than the validation loss.

Label	Data product
fCO ₂	SOCATv2023 ¹
A _T , C _T , Si _T , P _T	GLODAPv2 2023 ²
pH	BGC-Argo ³

Table 3: List of label data used to train the base model. ¹<https://socat.info/index.php/2023/06/20/v2023-release/>, ²<https://glodap.info/index.php/merged-and-adjusted-data-product-v2-2023/>, ³<https://www.seanoe.org/data/00311/42182/>

3 Model performance

4 Discussion and conclusions

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