Learning Inverse Dynamics Models in O(n) time with LSTM networks

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Abstract-Inverse dynamics model learning is crucial for modern robots where analytic models cannot capture the complex dynamics of compliant actuators, elasticities, mechanical inaccuracies, frictional effects or sensor noise. However, such models are highly nonlinear and millions of samples are needed to encode a large number of motor skills. Thus, current state of the art model learning approaches like Gaussian Processes or LWPR which scale exponentially with the data cannot be applied. In this work, we developed an inverse dynamics model learning approach based on a long-short-term-memory (LSTM) network with a time complexity of O(n). We evaluated the approach on a KUKA LWR arm that was used in object manipulation skills with various loads. In a comparison to Gaussian Processes we show that LSTM networks achieve better prediction performances and that they can be trained on large datasets with more than 100,000 samples in a few seconds. Moreover, due to the small training batch size of, e.g., 128 samples, the network can be continuously improved in lifelong learning scenarios.

I. INTRODUCTION

In robotics, predictions based on dynamics models are essential for control, object manipulation or planning [13], [20]. However, for modern robots with its dozens of compliant actuators and thousands of noisy tactile or visual sensors, these models have to be learned because of unmodeled effects like manufacturing uncertainties, sensory noise or any form of unmodeled dynamics in presence of contacts [3]. The challenge is that even small model inaccuracies can result in catastrophic behavior [21] and for learning a rich repertoire of versatile skills millions of samples in high dimensional observational spaces need to be processed [23].

Gaussian Processes (GPs) [26], [19] are state of the art regression techniques to learn such model. They are widely used in machine learning because their hyper parameters can be optimized through maximizing the marginal likelihood and thus, it requires little effort to adapt the model to new datasets or problem domains. However, the main limitation of GPs is that the computational complexity scales exponentially in $O(n^3)$ due to the need to invert a $n \times n$ kernel matrix.

To reduce this computational demands local approximations based on sparse GPs [17], mixture of experts approaches [24], [18], drifting GPs [12], or based on local partitions of the data [14] were proposed. However, also these local approximations scale quadratically with the number of samples, i.e., in $O(n^2)$ and are thus limited to few thousand samples [14].

Local approximations like the locally weighted regression approach [25] or even simpler lazy learning approaches like locally weighted regression would scale linear with the number of samples in O(n). However, tuning the often large number of hyper parameters, i.e., the Gaussian bandwidth parameters of many local models can be as challenging as the original model learning task.

We developed in this work an inverse dynamics model learning approach based on a recurrent neural network with a time complexity of O(N). To learn from long time series data we utilized gating mechanism in Long-short-termmemory (LSTM) networks [8], [22]. We demonstrate how LSTMs can learn the dynamics model of a torque controlled robot arm and compare the state of the art approach GPs. The model-based LSTM controller achieves superior prediction performances and scales linearly with the number of data samples. These results can be exploited in neural model-based control approaches that can learn from highdimensional visual and tactile data.



Fig. 1. Illustration of four snap shots of the investigated manipulation task. The robot had to push the flask that was filled with 200, 300, or 400 ml of liquid. The goal of this work is to predict the joint torques which depend on the fill level given joint angles, velocities and accelerations.

II. METHODS

We start with a problem definition for learning inverse dynamics models and subsequently discuss Gaussian Processes

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(GPs) and Long-short-term-memory (LSTM) networks implementations.

A. Problem Definition

We denote the vectors for joint angles, for joint velocities and for joint accelerations by q, \dot{q} , \ddot{q} , respectively. For a robot with d joints these vectors are of the dimension $\mathbb{R}^{d \times 1}$. Using these definitions the general inverse dynamics model of a robot is given by

$$\boldsymbol{\tau} = M(\boldsymbol{q})\boldsymbol{\ddot{q}} + h(\boldsymbol{q},\boldsymbol{\dot{q}}) + \boldsymbol{\epsilon}(\boldsymbol{q},\boldsymbol{\dot{q}},\boldsymbol{\ddot{q}}), \qquad (1)$$

where τ denotes the unknown joint torques which we want to learn. The inertia matrix is denoted by M(q) and the term $h(q, \dot{q})$ combines effects of Coriolis and centripetal forces, friction and of gravitational forces, i.e.,

$$h(oldsymbol{q}, \dot{oldsymbol{q}}) = C(oldsymbol{q}, \dot{oldsymbol{q}}) + f_r(oldsymbol{\dot{q}}) + g(oldsymbol{q})$$
 .

All unmodeled dynamics such as elasticities in the mechanical designs, model parameter inaccuracies in the masses or inertiae, vibrational effects, Stribeck friction, couplings and sensor noise are modeled by the term $\epsilon(q, \dot{q}, \ddot{q})$ in Eq. 1.

The goal of this work is to learn such inverse dynamics models in Eq. 1 including all effects of the unmodeled dynamics. We formulate this learning problem as a standard regression task. Given some input vector x the goal is to learn the function

$$\boldsymbol{y} = f(\boldsymbol{x}) + \boldsymbol{\zeta} : \mathbb{R}^{3d \times 1} \mapsto \mathbb{R}^{d \times 1}, \qquad (2)$$

where $\boldsymbol{y} = \boldsymbol{\tau}$ and $\boldsymbol{x} = [\boldsymbol{q}^T, \boldsymbol{\dot{q}}^T, \boldsymbol{\ddot{q}}^T]^T$. The variable $\boldsymbol{\zeta}$ denotes zero mean Gaussian noise with a standard deviation of σ_y .

Using a dataset $\mathfrak{D} = \langle x_t, y_t \rangle_{t=1,...,n}$ of *n* input-output pairs, we evaluate the trained models by computing the error

$$MSE = \frac{1}{dn} \sum_{j=1}^{d} \sum_{t=1}^{n} \left(\hat{y}_{t}^{[j]} - \tilde{y}_{t}^{[j]} \right)^{2}, \qquad (3)$$

where $\hat{y} = [\hat{y}^{[1]}, ..., \hat{y}^{[d]}]^T$ denotes the true label and \tilde{y} the model prediction.

In the following two subsections we will discuss how Gaussian Processes (GPs) and how Long-short-term-memory (LSTM) networks can be used to learn this mapping.

B. Inverse Dynamics Model Learning with Gaussian Processes

Gaussian Processes (GPs) are state of the art model learning or regression approaches [26], [19] that were successfully used for learning inverse dynamics models in robotic applications [14], [3]. For comprehensive discussions we refer to [21], [13]. Here we briefly discuss them as we will use them for a comparison in our experiments.

GPs represent a distribution over inverse dynamics models in Eq. 2 of the form $f \sim GP(m,k)$. This representation is fully defined by the mean m and the covariance k. We chose as covariance function a *Matérn kernel* [11]. It is a generalization of the *squared-exponential kernel* that has an additional parameter ν which controls the smoothness of the resulting function. The smoothing parameter can be beneficial for learning local models.We used *Matérn kernels* with $\nu = 5/2$ that are defined by

$$k(\boldsymbol{x}_p, \boldsymbol{x}_q) = \sigma^2 \, \frac{1}{2^{\nu-1} \, \Gamma(\nu)} \, A^{\nu} \, \mathbf{H}_{\nu} \, A + \sigma_y^2 \, \delta_{pq} \,,$$

where Γ is the gamma function, $A = (2\sqrt{\nu}||\boldsymbol{x}_p - \boldsymbol{x}_q||)/l$ and H_{ν} is a modified *Bessel* function [2]. The length-scale parameter of the kernel is denoted by σ , the variance of the latent function is denoted by σ and δ_{pq} is the *Kronecker* delta function (which is one if p = q and zero otherwise). Note that for $\nu = 1/2$ the *Matérn kernel* implements the *squaredexponential kernel*. In our experiments we optimized the hyper parameters $\boldsymbol{\theta} = [\sigma, l, \sigma_y]$ by maximizing the marginal likelihood [19].

Computing Predictions: Given a test sample x_* the predictive distribution is defined by

$$p(\tilde{\boldsymbol{y}}|\boldsymbol{\mathfrak{D}}, \boldsymbol{x}_*, \boldsymbol{\theta}) = \mathcal{N}(\mu_{GP}, \sigma_{GP}), \qquad (4)$$

with $\mu_{GP} = \boldsymbol{k}_*^T \boldsymbol{K}^{-1} \boldsymbol{y}$ and $\sigma_{GP} = k_{**} - \boldsymbol{k}_*^T \boldsymbol{K}^{-1} \boldsymbol{k}_*$. The matrix entries in \boldsymbol{K} are $K_{pq} = k(\boldsymbol{x}_p, \boldsymbol{x}_q)$, the scalar $k_{**} = k(\boldsymbol{x}, \boldsymbol{x})$, and $\boldsymbol{k}_* = k(\boldsymbol{X}, \boldsymbol{x}_*)$ with $\boldsymbol{X} = [\boldsymbol{x}_1, ..., \boldsymbol{x}_n]$.

C. Inverse Dynamics Model Learning with LSTM Networks

Long-short-term-memory (LSTM) networks are popular recurrent neural networks for modeling long time series [9]. Special gating mechanisms were added to classical recurrent neural networks to avoid the vanishing gradient problem in back propagation through time [8], [22]. In this subsection, we will give a short introduction to LSTMs and discuss the model with respect to inverse dynamics model learning problems.

In the following, subscripts will be used like in the previous subsections to index samples like in the dataset $\mathfrak{D} = \langle \boldsymbol{x}_t, \boldsymbol{y}_t \rangle$ at time t. However, in addition superscripts are introduced to denote the layer l = 1, ..., L, where the output at the last layer $\boldsymbol{h}_t^L = \tilde{\boldsymbol{y}}_t$ is used to compute the prediction error, e.g. by evaluating Eq. 3. The input to the first layer is denoted by $\boldsymbol{a}_t^{l=1} = \boldsymbol{x}_t$ and the next layer's input is set to the prediction of the previous layer, i.e., $\boldsymbol{a}_t^{l+1} = \boldsymbol{h}_t^l$.

At layer l a LSTM network is fully defined by the equations

$$\begin{split} & \boldsymbol{i}_{t}^{l} \ = \ \sigma(\boldsymbol{W}_{xi}^{l} \, \boldsymbol{a}_{t}^{l} + \boldsymbol{W}_{hi}^{l} \, \boldsymbol{h}_{t-1}^{l} + \boldsymbol{b}_{i}^{l}) \,, \\ & \boldsymbol{f}_{t}^{l} \ = \ \sigma(\boldsymbol{W}_{xf}^{l} \, \boldsymbol{a}_{t}^{l} + \boldsymbol{W}_{hf}^{l} \, \boldsymbol{h}_{t-1}^{l} + \boldsymbol{b}_{f}^{l}) \,, \\ & \boldsymbol{c}_{t}^{l} \ = \ \boldsymbol{f}_{t}^{l} \odot \boldsymbol{c}_{t-1}^{l} + \boldsymbol{i}_{t}^{l} \odot \tanh(\boldsymbol{W}_{xc}^{l} \boldsymbol{a}_{t}^{l} + \boldsymbol{W}_{hc}^{l} \boldsymbol{h}_{t-1}^{l} + \boldsymbol{b}_{c}^{l}) \,, \\ & \boldsymbol{o}_{t}^{l} \ = \ \sigma(\boldsymbol{W}_{xo}^{l} \, \boldsymbol{a}_{t}^{l} + \boldsymbol{W}_{ho}^{l} \, \boldsymbol{h}_{t-1}^{l} + \boldsymbol{b}_{o}^{l}) \,, \\ & \boldsymbol{h}_{t}^{l} \ = \ \boldsymbol{o}_{t}^{l} \odot \sigma(\boldsymbol{c}_{t}^{l}) \,, \end{split}$$

where the vectors i, f and o denote the input, the forget and the output gate, and c represents the memory cell of the network. The symbol $\sigma(.)$ denotes the sigmoid function and \odot is the *Hadamard* or element-wise product.

We used a simple model implementation that neglects recurrent connections from the memory cell to the gates [15] or dropout implementations [7].



Fig. 2. Hyper parameter optimization results using LSTMs. Shown are the training and the test error for variations of the number of hidden neurons $\in \{2, 5, 10, 15, 20\}$, the batch size $\in \{8, 16, 32, 64, 128, 256, 512\}$ and the number of epochs $\in \{1, 10, 20, 50, 100\}$ used during training. Bold numbers mark the used parameter value. We also evaluated different numbers of layers, which is not shown as more than two layers did not lead to any improvement.

Computing Predictions: The hidden value or output of LSTMs for inverse dynamics model learning is given by

$$\tilde{\boldsymbol{y}}_t = \boldsymbol{o}_t^L \odot \sigma(\boldsymbol{c}_t^L)$$

where L denotes the number of layers in the LSTM network.

Implementation Details: All experiments were performed using an efficient *tensorflow* [1] implementation of LSTM networks. For the gradient based updates adaptive learning rates were implemented [5]. Prior to our comparison to Gaussian Processes we optimized the batch size, the number of epochs and the number of layers using our real robot dataset. These results are shown in Figure 2.

III. RESULTS

We evaluated the prediction performance and the computational times for training and for generating predictions on two datasets. A synthetic dataset of one dimensional (d = 1)*sine, triangle* and *sawtooth* functions is used to explore the robustness to noise.



Fig. 3. Training data of a synthetic time series prediction task. The left panel shows the training data for the *low noise* condition with zero mean Gaussian noise with a standard deviation of 0.01. In the right, a standard deviation of 0.1 was used.

In a *dynamics* model learning task using a torque controlled robot arm of KUKA, we evaluated the prediction performances and the required computational time for generating online predictions. The first five joints (d = 5) of the arm were used which results in a 15 dimensional input space, i.e., positions, velocities and accelerations. The five dimensional torque vectors are treated as unknowns. All results show average statistics over 20 cross validation experiments.

Note that we optimized the model parameters of the LSTM network on the real robot dataset. The results for the optimal batch size, the number of epochs and the number of layers are shown in Figure 2.

A. On the effect of noise on the prediction accuracy

TABLE I

Mean squared errors of the predictions on the synthetic dataset for additive zero mean Gaussian noise with a standard deviation of 0.01 (low noise) and 0.1 (large noise condition). The \pm symbol denotes the standard deviation over 20 experiments.

	Naive	GPs	LSTMs
low noise	0.264 ± 0.084	0.030 ± 0.011	$\textbf{0.010}\pm0.004$
large noise	0.310 ± 0.078	0.086 ± 0.016	0.075 ± 0.014

We generated two synthetic datasets with additive zero mean Gaussian noise (with $\sigma = \{0.01, 0.1\}$) to test the robustness of the regression techniques to noise. The data is illustrated in Figure 3 and shows an overlay of three functions, i.e., a *sine*, a *triangle* and *sawtooth* function.

With increasing noise level, the mean squared error of the model's predictions increased. In both conditions, LSTM



Fig. 4. Comparison of regression models on the synthetic time series prediction task. In the box, prediction results for eight consecutive time steps at phase $\pi/2$ are shown. The dark solid line denotes the ground truth target values.

networks outperform GPs. These results are shown in Table I. In Figure 4, we additionally illustrate some example predictions.

We also compared to a naive approach which simply computes the output based on the sum of the current state and the current velocity. This approach fails to compute accurate predictions at the turning points of the functions, which is shown in the inlay in Figure 4.

B. On the benefit of memory models for sequential data



Fig. 5. Comparison of regression models on sequential and non-sequential **real-robot** data, where we eliminated all temporal correlations through random permutations. While there is no significant difference in the GP results, the LSTM network can exploit temporal correlations in the data, i.e, the robot joint angle, the velocity and the acceleration trajectories.

An important question is if models with memory like LSTMs can capture and exploit temporal correlations in the inverse dynamics model learning data. To test that we compared the prediction performance of LSTMs trained on the **real robot** sequential data with models trained on data where we removed all temporal correlations through random permutations. These results and a comparison to GPs are shown in Figure 5. Obviously, GPs showed no significant difference in the prediction error. However, LSTMs could exploit the temporal correlations in the data.

C. Dynamics model learning for torque control

The predicted joint torques are highly nonlinear signals which are shown for two arm reaching motions in Figure 8. Both motions have a time horizon of about 4.1 seconds, which is denoted by the vertical line.

The first panel in Figure 8 illustrates the mean squared error of the predictions of GPs and LSTMs. The remaining panels show predictions of GPs and LSTMs of all five joint torques in the KUKA robot. From these predictions one can see that already a small difference in the mean squared error, i.e., GPs: 0.014 ± 0.0006 and LSTMs: 0.005 ± 0.002 , results in large deviations from the true joint torque signals.

For computing these results the GPs and the LSTM networks were trained with 1000 samples. In Figure 6 we illustrate the effect of the sample size on the prediction performance. On small datasets with less than 1000 samples GPs achieve better results. However, with increasing sample numbers LSTM networks outperform GPs. For 10,000 samples optimizing the length-scale parameters of the GPs was computationally intractable. Even more dramatically is the computational benefit which is discussed in the next subsection.



Fig. 6. Comparison of the mean squared error of GPs and LSTMs for an increasing number of data samples. The solid lines show the mean values and the shaded areas denote the standard deviation over 10 runs. Note that for 1000 samples a more detailed comparison is shown in Fig. 8.

D. On the effect of dataset size on the computational time

TABLE II

Computational times in seconds on the real robot dataset with 1,000, 10,000, 50,000 and 100,000 training samples. The \pm symbol denotes

The standard deviation over 10 experiments.

	1k	10k	50k	100k
GPs	$33.7~\pm~3.9$	$3196~\pm~355$	9 $10^4 (25h)$	_
LSTMs	2.9 ± 0.2	28.1 ± 0.6	141.1 ± 1.2	282.4 ± 2.8

Theoretically LSTM networks scale linearly with the number of samples in O(n), while GPs scale exponentially with $O(n^3)$. To test these theoretical time complexities, we conducted a numerical evaluation using the real robot data. The results of this comparison of GPs and LSTMs is shown in Figure 7. For small datasets with up to 1000 samples both approaches need similar resources, however with more than 1000 samples GPs could not produce predictions in a



Fig. 7. Comparison of the computational time of GPs and LSTMs for an increasing number of data samples. The solid lines show the mean values and the shaded areas denote the standard deviation. The inlay shows results for less than 1148 samples.

reasonable amount of time, e.g., for 10,000 samples, training the GPs took more than 24 hours as shown in Table II. LSTM networks in contrast could be trained in 28.1 seconds and for learning from 112,761 samples, training took 282.4 seconds. Note that for all evaluations a standard PC with 8 cores operating at 3.4 GHz and 16 GB ram was used.

IV. CONCLUSIONS

We demonstrated in this work that complex inverse dynamics models of a real robot arm can be learned in linear time from large datasets with more than 100,000 samples. The prediction error of the used Long-short-term-memory (LSTM) networks decreased exponentially with the number of samples and even for small datasets we outperformed the state of the art approach — Gaussian Processes (GPs).

A notable difference of the neural network predictions compared to the ones of GPs is the *nonexistence* of a variance or prediction uncertainty. In various approaches however, it was shown that such variance estimates can be used to improve a movement representation or to adjust the controller stiffness, see for example [16], [4], [10]. Given the recent investigations aiming at learning a predictive variance in deep networks, e.g., by exploiting the stochasticity of the outputs when using *dropout* [6], the LSTM model could be extended which is part of future work.

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Fig. 8. Joint torque prediction results of GPs and LSTMs on a **real robot test dataset**. The first panel shows the mean squared error of both approaches for 1000 training samples. The remaining panels show the predicted joint torques of all five joints of the KUKA LWR arm. The dark line denotes the ground truth target values and the vertical line marks the end of one trial of robot arm reaching.