Interpolation of Molecular Dynamics with Bi-Directional Neural Networks

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Solution is a trajectory in phase space

$$x(t) = \begin{bmatrix} r(t) \\ p(t) \end{bmatrix} = \int_{t_0}^t f\left(\begin{bmatrix} r(t) \\ p(t) \end{bmatrix}, t\right) dt$$



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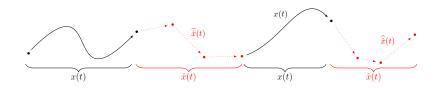
Can we learn the phase space dynamics with a ML algorithm?

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Model Architectures

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HamiltonianNetwork

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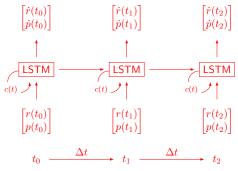
RNN and LSTM

$$\dot{x}(t) = \begin{bmatrix} \dot{r}(t) \\ \dot{p}(t) \end{bmatrix} = f_{\theta} \left(\begin{bmatrix} r(t_0) \\ p(t_0) \end{bmatrix}, \dots, \begin{bmatrix} r(t) \\ p(t) \end{bmatrix} \right], t \right)$$



Learning Dynamical Systems with LSTM

- Best performing due to fewest assumptions and flexible parameterization
- Memory cell c(t) to selectively read and write information
- ullet Outputs $[\dot{r}(t),\dot{p}(t)]^T$ are integrated to obtain solution $\hat{x}(t)$



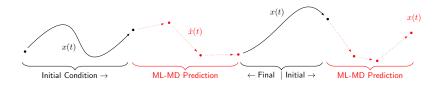
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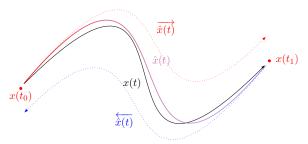


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$$\hat{x}(t) = (1 - \lambda(t)) \ \overrightarrow{\hat{x}(t)} + \lambda(t) \ \overleftarrow{\hat{x}(t)}$$
 (1)

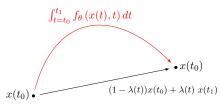


For time-reversible solutions, we obtain

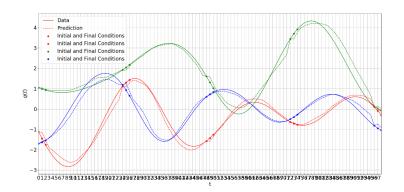
$$\hat{x}(t) = (1 - \lambda(t)) \ \overrightarrow{\hat{x}(t)} + \lambda(t) \ \overleftarrow{\hat{x}(t)}$$

$$= \underbrace{(1 - \lambda(t))x(t_0) + \lambda(t)x(t_1)}_{\text{low frequency components}} + \underbrace{\int_{t=t_0}^{t_1} f_{\theta}\left(x(t), t\right) dt}_{\text{high frequency components}}$$

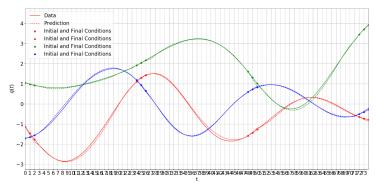
 Adiabatic connection frees the ML model to model high frequency signals



 Unidirectional LSTM architecture for Benzene MD trajectory interpolating over 20 time steps



- Bidirectional LSTM architecture for Benzene MD trajectory interpolating over 20 time steps
- Final condition and additional bidirectional training smooth trajectories significantly



 Single initial and final condition already good for sufficient performance by bidirectional LSTM

