Markov state models

Theory, properties, estimation and validation

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Motivation





 ω_e / 10³ rad \cdot s⁻¹

 ω_e / 10³ rad \cdot s⁻¹

Example: BPTI

Simulation of BPTI



Markov state models



Metastability of states allow us to significantly simplify the dynamics of our system of interest

Markov state models

Final state

| | | | Ser - | | |
|--------------|--|-----|-------|-----|-----|
| | Start Start | 96% | 1% | 2% | 1% |
| nitial state | Sec. | 5% | 95% | 0% | 0% |
| | Sec. | 1% | 0% | 97% | 2% |
| | Ser and a ser a se | 1% | 0% | 2% | 97% |

A Markov state model describes the dynamics of a system as conditional transition probabilities

What is meta-stability?





is important.

What is meta-stability?





is important.

Molecular simulations

- Molecular simulations are realizations of stochastic process on $\,\Omega\,$ and are Markovian w.r.t. this space.

$$p(\mathbf{x}, \mathbf{y}; \tau) \, d\mathbf{y} = \mathbb{P}[\mathbf{x}(t + \tau) \in \mathbf{y} + d\mathbf{y} \mid \mathbf{x}(t) = \mathbf{x}]$$
$$\mathbf{x}, \mathbf{y} \in \Omega, \ \tau \in \mathbb{R}_{0+},$$

Transition probabilities are well defined

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$$\mathbf{x}, \mathbf{y} \in \Omega, \ \tau \in \mathbb{R}_{0+},$$

Transition probabilities are well defined

$$p(\mathbf{x}, A; \tau) = \mathbb{P}[\mathbf{x}(t + \tau) \in A | \mathbf{x}(t) = \mathbf{x}]$$
$$= \int_{\mathbf{y} \in A} d\mathbf{y} \ p(\mathbf{x}, \mathbf{y}; \tau).$$

Also applies for regions

Molecular simulations (2)

Ergodicity

No two or more segments of the space Ω are dynamically disconnected from each other.

and

For an infinitely long simulation we will have visited every state $\mathbf{x} \in \Omega$ infinitely many times.

Molecular simulations (3)

Reversibility

Simulations fulfill the detailed-balance condition:

$$\mu(\mathbf{x}) \ p(\mathbf{x}, \mathbf{y}; \tau) = \mu(\mathbf{y}) \ p(\mathbf{y}, \mathbf{x}; \tau)$$

$$\mu(\mathbf{x}) = Z(\beta)^{-1} \exp\left(-\beta H(\mathbf{x})\right)$$

At equilibrium the probability of jumping from any x to any y is the same as jumping from y to x.

An illustration of the transition density



Instead of single realizations we now focus on the evolution of an ensemble

$$p: x \in \Omega \mapsto p(x) \in \mathbb{R}_0^+, \quad \int_{\Omega} dx p(x) = 1$$

Figure courtesy of JH Prinz

Assumptions about the full dynamics

Markovian

$$\mathbb{P}(x_{t+\tau} \in A \mid x_{t_1}, \dots, x_t = x) = \mathbb{P}(x_{t+\tau} \in A \mid x_t = x)$$

Factorization of the dynamics into conditional probabilities

Chapman-Kolmogorov property

$$p_{\tau_1 + \tau_2}(x, A) = \int_{\Omega} p_{\tau_1}(x, y) p_{\tau_2}(y, A) \, \mathrm{d}y$$

Direct combination of conditional probabilities with different lag-times

96% 1% 2% 1% Initial state 5% 95% 0% 0% 1% 0% 97% 2% h C 1% 0% 2% 97%

Final state

Assumptions about the full dynamics

Irreducibility

All states of the state space can be reached from any other state in a finite time. **Ensures unique stationary distribution.**

Ergodicity

No states are disconnected No cyclic dynamics. Ensures time and ensemble average properties are equal.

Reversibility

No net-probability flux at equilibrium. => no energy production/absorption => mass conservation. Not strictly necessary for Markov models

Ensemble view of dynamics



TIME τ

TIME $t+\tau$

A propagator is an operator which transports probability densities in time

$$p_{t+\tau}(x) = [P_{\tau} p_t](x) = \int_{\Omega} dy p_{\tau}(y, x) p_t(y)$$
$$p_{t+\tau}(x) = [P_{\tau} p_t](x) = \int_{\Omega} dy p_{\tau}(y, x) p_t(y)$$

Example dynamics





Propagator depends on lag time



Propagator depends on lag time



Propagator depends on lag time



So why is this?

Implied time-scales

Eigenvalues of the propagator

 $\Pr_{\tau} \phi_i^{P_{\underline{\tau}}} \phi_i = \lambda_i \phi_i$ Chapman-Kolmogorov Implies exponential lag-time dependence





Meta-stability

- We can approximate the propagator by a finite number of processes with non-zero Eigenvalues
- If we have a gap in the Eigenvalue spectrum, we can choose the lag-time in a manner such that we fulfill this assumption
- When we do this, processes faster than the lagtime 'have decayed' or 'are not resolved'.



Prinz et al. (2011) JCP 134, 174105

Estimation

Discretization of Ω



Count matrix

| C _{ij} (1) | A | B | C | D |
|----------------------------|------|------|------|------|
| A | 9963 | 37 | 0 | 0 |
| B | 22 | 9974 | 4 | 0 |
| C | 0 | 2 | 9919 | 79 |
| D | 0 | 0 | 115 | 9885 |

$$C_{ij}(\tau) = \delta(x_{n-\tau} = i, x_n = j)$$

Maximum likelihood estimator

We can express the probability of the observed data - discrete trajectory - given a transition probability matrix of an MSM

$$\mathbb{P}(x_1, \dots, x_t \mid P) = \prod_{k=1}^L p_{x_{k-1}, x_k}$$
$$= p_{x_0, x_1} \cdot \dots \cdot p_{x_L - 1, L}$$
$$= \prod_{ij} p_{ij}^{c_{ij}}$$
$$= p_{11}^{c_{11}} \cdot \dots$$

The aim is then to find the *P* which maximizes this expression -That is, the *Maximum likelihood estimator*.

Analytical solution for Nonreversible case $p_{ij} = 1, \forall i$ We enforce the constraint that the transition probability matrix is row $\mathfrak{stochastic:}_{\forall i}$ $P^{\text{MLE}} \stackrel{j}{=} \operatorname{argmax} \prod^{L} p_{x_{k-1}, x_k}$ $\sum p_{ij} = 1, \stackrel{P}{\forall} i \stackrel{\overline{k=1}}{\underset{L}{\overset{P}{i}}}$ $P^{\text{MLE}} = \underset{P}{\operatorname{argmax}} \prod_{k=1}^{k} p_{x_{k-1}, x_k}$ k=1

• One can show the estimator is simply: $P^{\text{MLE}} = \operatorname{argmax} \prod_{k=1,x_k}^{L} p_{x_{k-1},x_k}$

$$\hat{p}_{ij} = \frac{\hat{C}_{ij}}{\sum_{j} \hat{C}_{ij}} = 1$$

$$\hat{p}_{ij} = \frac{\sum_{n=\tau}^{L} \delta(x_{n-\tau} = i, x_n = j)}{\sum_{k=\tau}^{L} \delta(x_{n-\tau} = i, x_n = j)}$$

Prinz et al. (2011) JCP 134, 174105

Reversible estimator

- Enforces the detailed balance condition.
- No exact analytical solution:
 - Fixed-point iteration algorithm available.
 - Approximate solutions.
- Implemented in PyEMMA

- The less simulation data we have, the more ambiguous the solution of the likelihood problem will be.
- Consequently, if we limit ourselves to the MLE, we are *ignorant* as to how **robust** our inferred MSM is.
- One way to quantify the uncertainty of MSMs is through **Bayesian inference**

Likelihood from before

$$\mathbb{P}(x_i, \dots, x_t \mid P) = p(C \mid P) \propto \prod_{i,j=1}^n p_{ij}^{c_{ij}}$$

Likelihood from before

$$\mathbb{P}(x_i, \dots, x_t \mid P) = p(C \mid P) \propto \prod_{i,j=1}^n p_{ij}^{c_{ij}}$$

Introduction of prior information

$$p(P \mid C) \propto p(C \mid P)p(P)$$

The prior can encode useful constraints: row-stochasticity, reversibility, fixed stationary distribution, sparsity etc

Inference is done by MCMC sampling



Noé (2008) JCP 128, 244103 Trendelkamp-Schroer & Noé (2013) JCP 138, 164113

Alternative estimators

Transition(-based) Reweighting Analysis Method

- Allows taking into account simulation data from multiple thermodynamic ensembles.
- That means, we can use data from enhanced sampling simulations together with unbiased simulation data to generate models more efficiently.
- More about this tomorrow.



Wu et al. PNAS 2016, 113(23), E3221–E3230

Implemented in PyEMMA

Augmented Markov models

- Enables integration of external information into the estimation of Markov state models.
- Fx use of experimental constraints from biophysical experiments such as NMR.
- A notebook tutorial distributed with PyEMMA 2.5 and up.



Olsson et al. PNAS 2017, 114(31), pp. 8265-8270. doi: 10.1073/pnas.1704803114

Implemented in PyEMMA

Analysis of our estimate

| P _{ij} (1) | A | B | C | D | projected timescales | original timescales |
|----------------------------|--------|--------|--------|--------|-------------------------|------------------------|
| A | 0,9963 | 0,0037 | | | ∞ | ∞ |
| B | 0,0022 | 0,9974 | 0,0004 | | 2,746 | 17,671 |
| C | | 0,0002 | 0,9919 | 0,0079 | 165 | 1,610 |
| D | | | 0,0115 | 0,9885 | 51 | 538 |

Time-scales are always under-estimated

Increasing the lag-time

| COU | IN | T |
|-----|----|---|
| MAT | R | X |

| C _{ij} (100) | A | B | C | D |
|------------------------------|------|------|------|------|
| A | 9533 | 477 | 40 | 0 |
| B | 1644 | 8014 | 262 | 80 |
| C | 0 | 40 | 9025 | 935 |
| D | 0 | 0 | 1366 | 8634 |

| original timescales | projected timescales |
|------------------------|-------------------------|
| ∞ | ∞ |
| 17,671 | 15,397 |
| 1,610 | 1211 |
| 538 | 379 |

May improve $e^{\hat{C}_{ij}} = \frac{L/\tau}{\text{stim}} \sum_{a=1}^{L} \delta(x_{a} - p_{a} - p_{a}) = \delta(x_{a} - p_{a} - p_{a})$

Figure courtesy of JH Prinz

Projection/discretization error



Projection/discretization error



Figure courtesy of JH Prinz

Known problems

- Observations (projections, discretizations) are in many cases <u>not Markovian</u>
- However, we are often interested in understanding the full system not just the observation.
- Since we often have a lot of freedom to choose the projections and discretization, it is important to chose one which is as Markovian as possible.

Validation

Chapman-Kolmogorov test

Compare the evolution of the data with the model



General scheme for Markov state model generation

- Discretize a suitable projection of your data.
- Construct a transition matrix.
- Estimate the number of meta-stable states (timescale gap)
- Perform Chapman-Kolmogorov test.

Analysis

Useful predictions from a MSM

Common properties

- Relaxation time-scales
- Dominant processes
- Stationary distribution (thermodynamics)
- Meta-stable sets (more about this later)
- Correlation functions (spectroscopic observables)
- Mean first passage times
- Path probabilities

Spec roscopic observables



Noé et al. Dynamical fingerprints for probing individual relaxation processes in biomolecular dynamics with simulations and kinetic experiments. Proc. Nat. Acad. Sci. USA 108, 4822–4827 (2011).

Olsson & Noé Mechanistic Models of Chemical Exchange Induced Relaxation in Protein NMR. 139, 200–210 JACS (2017)

Summary

- Markov state models are derived coarse-grained models of the full original (Markovian) dynamics.
- MSMs may be parameterized (estimated/learned) from simulation data to compute properties of interest.
- MSMs are particularly useful if the projection/ discretization error can be minimized: then the predicted quantities match the original.

Questions?

Otherwise it's time for the practical.