# Using low-rank tensor formats to enable computations of cancer progression models in large state spaces

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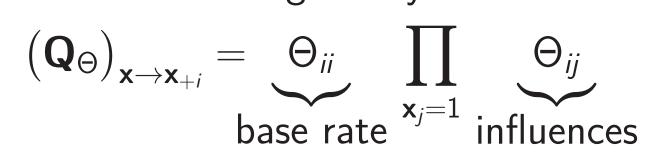
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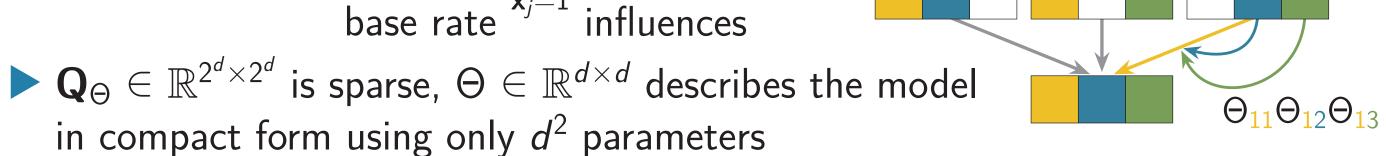
# Summary

- $\triangleright$  Comprehensive cancer progression models should include a large number d of genomic events
- Mutual Hazard Networks model the progression process using only  $d^2$ parameters [1]
  - Computational complexity of a straightforward implementation still scales exponentially in d
  - Calculations using  $\gtrsim 25$  events are computationally infeasible [2]
- Tensor Trains allow for cost-effective storage and calculations for high-dimensional tensors
  - This method reduces the computational complexity from exponential to polynomial in d

## Mutual Hazard Network (MHN) model

- MHN models tumor progression as a continuous-time Markov chain on the  $2^d$ -dimensional state space of possibly active events [1]
- Events can only occur one at a time
- Transition rates are given by





 $\triangleright$  Time marginal probability distribution from  $\Theta$ :



 $\triangleright$  Optimal  $\Theta$  matrices are found by optimizing the time-marginalized Kullback-Leibler divergence from the given data distribution  $\mathbf{p}_{\mathcal{D}}$ :

$$S_{\mathsf{KL}}(\mathbf{p}_{\Theta}) = \sum_{\mathbf{p}} (\mathbf{p}_{\mathcal{D}})_{\mathbf{x}} \log ((\mathbf{p}_{\Theta})_{\mathbf{x}})$$

Gradients can be calculated analytically:

$$\frac{\partial S_{\mathsf{KL}}}{\partial \Theta_{ij}} = \sum_{\mathbf{y}, \mathbf{z}} \sum_{\mathbf{x}} \frac{\partial S_{\mathsf{KL}}}{\partial (\mathbf{p}_{\Theta})_{\mathbf{x}}} (\mathbf{Id} - \mathbf{Q}_{\Theta})_{\mathbf{x}\mathbf{y}}^{-1} \left( \frac{\partial \mathbf{Q}_{\Theta}}{\partial \Theta_{ij}} \right)_{\mathbf{y}\mathbf{z}} (\mathbf{p}_{\Theta})_{\mathbf{z}}$$

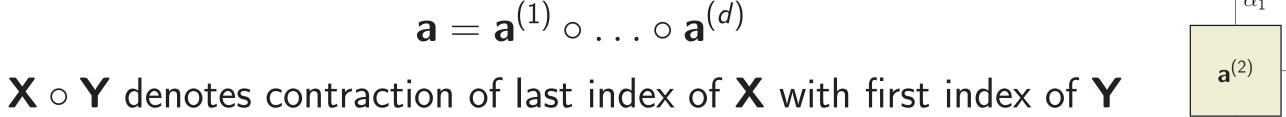
► KL divergence and gradient calculation time is dominated by solution time of two linear equations:

 $(\mathsf{Id} - \mathbf{Q}_{\Theta}) \, \mathbf{p}_{\Theta} = \mathbf{p}_{\varnothing} \qquad (\mathsf{Id} - \mathbf{Q}_{\Theta})^{\mathsf{T}} \, \mathbf{q} = \frac{\partial S_{\mathsf{KL}}}{\partial \, \mathbf{p}_{\Theta}}$ 

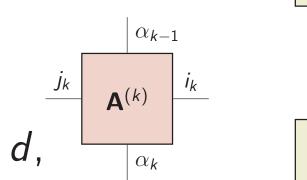
# Tensor Train (TT) representation

A d-dimensional tensors  $\mathbf{a} \in \mathbb{C}^{n_1 \times ... \times n_d}$  can be written as a product of d Tensor Train cores  $\mathbf{a}^{(k)} \in \mathbb{C}^{r_{k-1} \times n_k \times r_k}$ :





Similar for operators  $\mathbf{A} \in \mathbb{C}^{(m_1 \times ... \times m_d) \times (n_1 \times ... \times n_d)}$ : TT cores  $\mathbf{A}^{(k)} \in \mathbb{C}^{r_{k-1} \times m_k \times n_k \times r_k}$ 



 $\triangleright$  Storage cost is reduced from exponential to linear in d, but additional dependency on TT ranks  $r_k$  is introduced

- Many arithmetic operations can be performed directly in the TT format, reducing the computational complexity [3]:

  - Superposition  $\lambda \mathbf{a} + \nu \mathbf{b}$ :  $\mathcal{O}(dn(r_{\mathbf{a}} + r_{\mathbf{b}})^2)$   $n := \max(n_k)$ > Inner product  $\langle \mathbf{a}, \mathbf{b} \rangle$ :  $\mathcal{O}(dnr_{\mathbf{a}}r_{\mathbf{b}}(r_{\mathbf{a}} + r_{\mathbf{b}}))$  $m := \max(m_k)$
  - Operator-by-Tensor product  $\mathbf{Ab}$ :  $\mathcal{O}(dmn(r_{\mathbf{A}}r_{\mathbf{b}})^2)$
- $r_{\mathbf{X}} := \max((r_{\mathbf{X}})_k)$
- $\triangleright$  Linear equations  $\mathbf{A}\mathbf{x} = \mathbf{b}$  can also be solved efficiently directly in this format

# References

- [1] R. Schill, S. Solbrig, T. Wettig, and R. Spang, Modelling cancer progression using Mutual Hazard Networks, Bioinformatics 36 (January, 2020) 241.
- [2] P. Georg, L. Grasedyck, M. Klever, R. Schill, R. Spang, and T. Wettig, Low-rank tensor methods for Markov chains with applications to tumor progression models, Journal of Mathematical Biology 86 (December, 2022).
- [3] P. Georg, Tensor Train Decomposition for solving high-dimensional Mutual Hazard Networks, PhD thesis, Universität Regensburg, October, 2022.

#### **Tensor Trains for MHN**

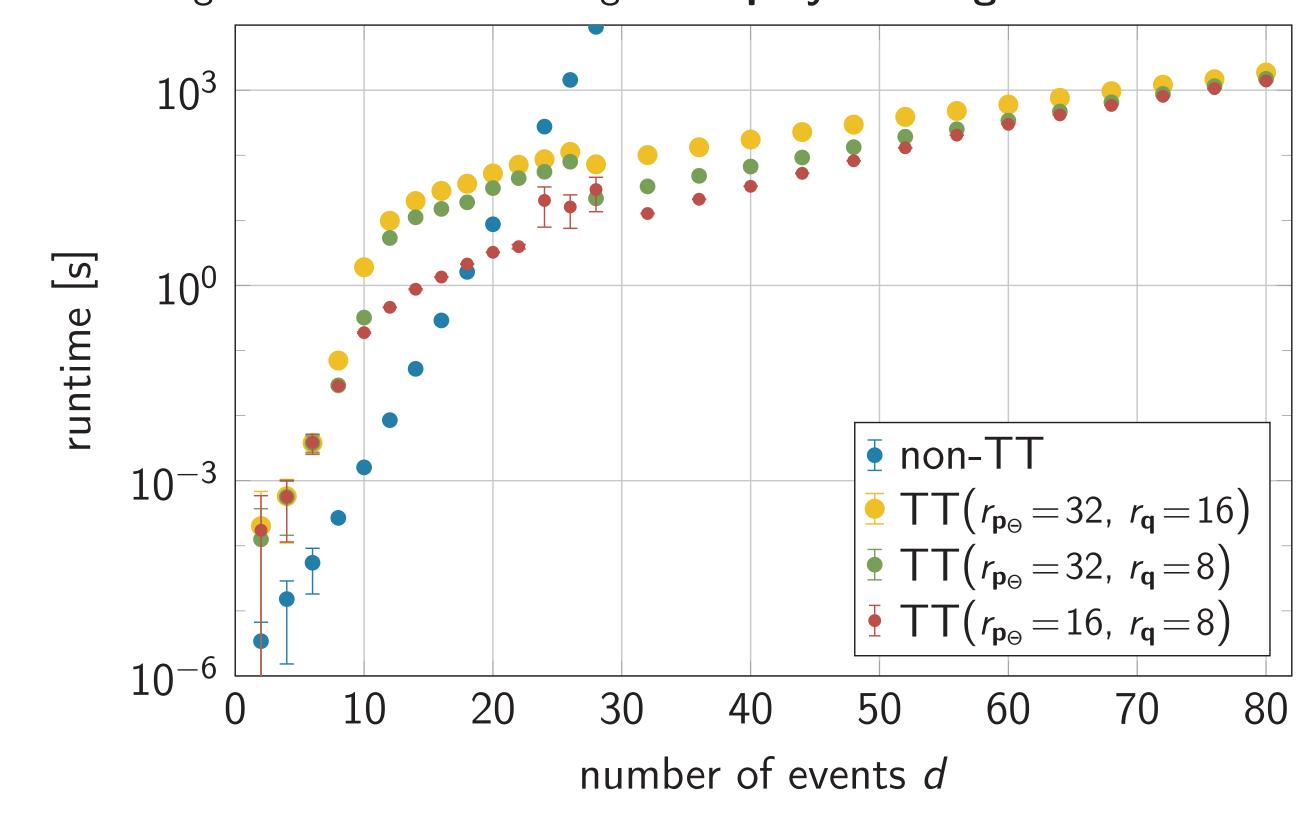
- ightharpoonup Events are binary  $\rightarrow n_k = 2$  for all mode sizes
- $ightharpoonup \mathbf{Q}_{\Theta}$  can naturally be written as a Tensor Train [1]:

$$\mathbf{Q}_{\Theta} = \sum_{i=1}^{d} \left( \bigotimes_{j=1}^{i-1} \underbrace{\begin{pmatrix} 1 & 0 \\ 0 & \Theta_{ij} \end{pmatrix}}_{\in \mathbb{R}^{1 \times 2 \times 2 \times 1}} \otimes \underbrace{\begin{pmatrix} -\Theta_{ii} & 0 \\ \Theta_{ii} & 0 \end{pmatrix}}_{\in \mathbb{R}^{1 \times 2 \times 2 \times 1}} \otimes \underbrace{\bigotimes_{j=i+1}^{d} \underbrace{\begin{pmatrix} 1 & 0 \\ 0 & \Theta_{ij} \end{pmatrix}}_{\in \mathbb{R}^{1 \times 2 \times 2 \times 1}} \right)$$

- $\mathbf{Q}_{\Theta}$  is a sum of d rank-1 Tensor Trains
- All TT-ranks of  $\mathbf{Q}_{\Theta}$  are equal to d
- $\mathbf{p}_{\varnothing}$  is a canonical unit Tensor Train, with all TT-ranks equal to 1
- $ightharpoonup \mathbf{p}_{\Theta}$  and  $\mathbf{q}$  can be calculated in the TT format (max. TT ranks  $r_{\mathbf{p}_{\Theta}}$  and  $r_{\mathbf{q}}$ )
- $\blacktriangleright$  For gradients, each nonzero entry in  $\mathbf{p}_{\mathcal{D}}$  has to be treated individually
  - One linear equation has to be solved for each nonzero entry (usually  $\sim \! 1000)$
  - This can be parallelized trivially

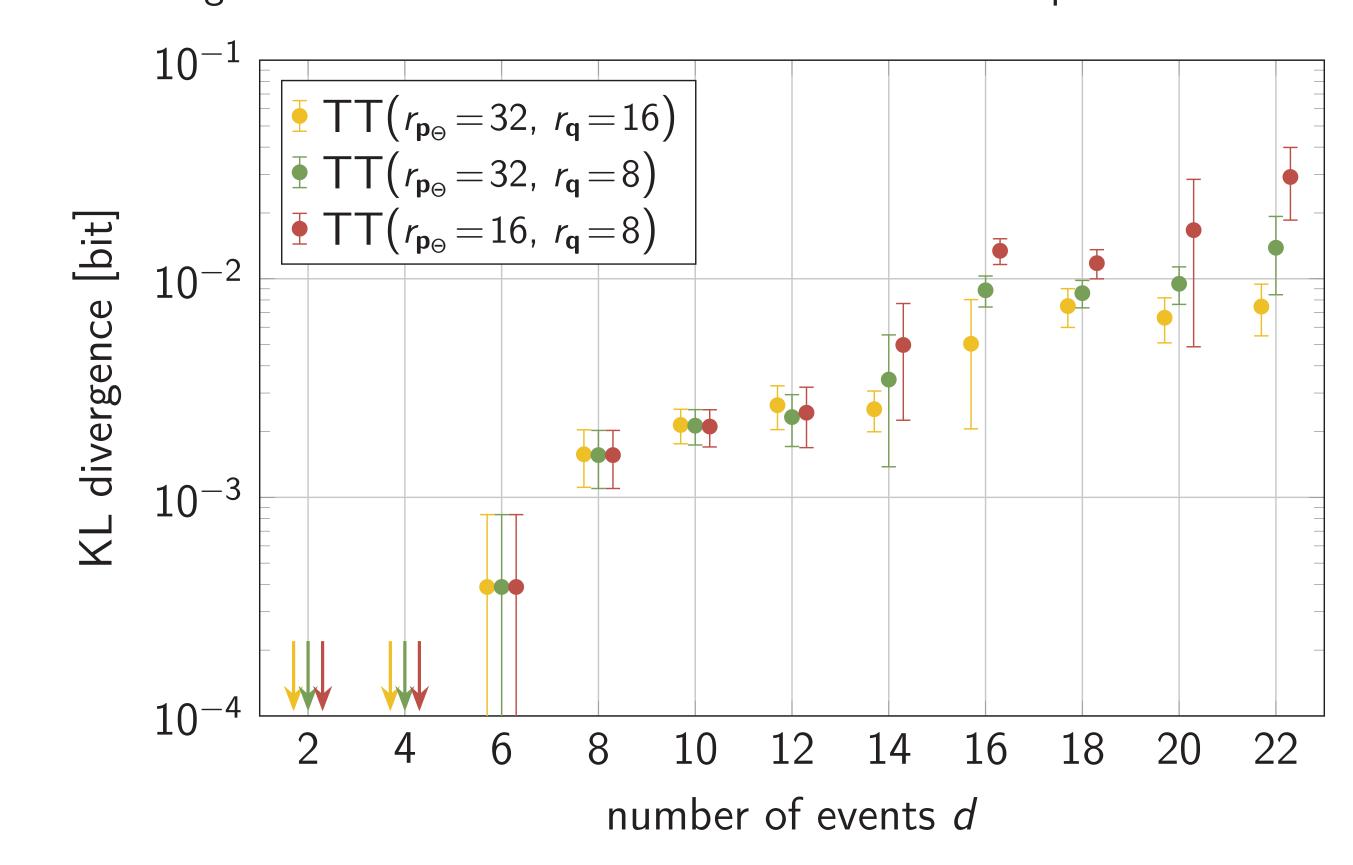
#### Results: Runtime speedup

- ightharpoonup Runtime for one score and gradient evaluation at  $\Theta=$  independence model
- $ightharpoonup p_{\mathcal{D}}$  constructed from 1000 random samples
- Runtime grows with  $\sim d^{5.4}$  for large  $d \Rightarrow$  polynomial growth!



# Results: Accuracy of the TT solution

 $\triangleright$  KL divergence from exact result to TT solution after full optimization of  $\Theta$ 



# **Code availablility**

- C++ library for TT-calculations pRC: gitlab.com/pjgeorg/pRC
- ► Application-specific C++ library cMHN that utilizes pRC for MHN-calculations: soon to be open-source

## **Future improvements**

- ► Reduce runtime by accelerating solution of linear equations in TT format
- Include formation of metastases in the model