TCI.7

Goal: unfold given tensor to TT via repeated CI factorizations (either CI or prrLU) into 'TCI form':

Defining characteristics of this decomposition: it is built only from <u>one-dimensional slices</u> of F<sub>\$\sigma\$</sub>

TCI algorithms use only <u>local</u> updates of these slices. (on which all tensor indices \$\sigma\$ but one are fixed)

For each CI factorization, say along bond  $\ell$ , the tensor is viewed as a matrix, with row index  $i = (G_1, \dots, G_L)$  with dimension(i) = dcolumn index  $j = (G_{\ell_1}, \dots, G_L)$  with dimension(j) = dCI factorization  $F_{ij} \simeq \widehat{F}_{ij} = [C_L]_{ij} [P_L]_{ji} [R_\ell]_{ij}$  external indices i, j are fixed,  $F_{ij} = \frac{i}{\mathbb{L}_\ell} \int_{\ell_{\ell_1}} \frac{C_{\ell_2}}{C_{\ell_1}} \frac{C_{\ell_2}}{C_{\ell_2}} \frac{C_{\ell_2}}{C$ 

external indices i, j are fixed, internal bonds represent sums over pivot lists:  $j \in \mathcal{I}_{\ell+1}$ 

t lists:  $\int_{-\infty}^{\infty} \int_{\mathbb{R}^{+}} \int_{-\infty}^{\infty} \int_{\mathbb{R}^{+}} \int_{\mathbb{R}^{+}}$ 

Note: all ingredients carry bond index  $\ell$ .

Dimensions of  $\iota$ , j are generically large, 'top-down' CI factorization is impractical. Instead, use 'bottom-up' approach: start 'TCI Ansatz' (1) with small pivot matrices, increase their size via 'sweeps' until desired tolerance is reached.

To be systematic, we introduce some bookkeeping conventions:

External index 
$$\mathcal{L}$$
 ( $\mathcal{L} \in \{1, 2, ..., L\}$ ) takes  $\mathcal{L}$  different values from set  $\mathcal{L}$ 

$$\mathbb{I}_{\ell} = \mathbb{S}_{1} \times \dots \times \mathbb{S}_{\ell} = \text{set of row multi-indices up to site } \mathcal{L}; \qquad \qquad i \in \mathbb{I}_{\ell} \text{ has the form } i = (6_{i_1}, \dots, 6_{\ell}) \tag{SA}$$

$$\mathbb{J}_{\ell} = \mathbb{S}_{\ell} \times \dots \times \mathbb{S}_{\ell} = \text{set of column multi-indices from site } \mathcal{L} \text{ upwards; } \mathcal{J} \in \mathbb{J}_{\ell} \text{ has the form } \mathcal{J} = (\mathcal{S}_{\ell}, \dots, \mathcal{S}_{\ell}) \quad (sb)$$

$$\underline{\mathbb{T}}_{\underline{L}} = \underline{\mathbb{T}}_{\underline{I}} = \text{full configuration space.}$$
 A full configuration  $\overrightarrow{\sigma} \in \underline{\mathbb{T}}_{\underline{L}}$  takes the form  $\overrightarrow{\sigma} = (6, \dots, 6)$ 

$$i_{\ell} \oplus j_{\ell+1} = \left( \underbrace{\sigma_{1}, \dots, \sigma_{\ell}}_{i_{\ell}}, \underbrace{\sigma_{\ell+1}, \dots, \sigma_{\ell}}_{j_{\ell+1}} \right) = \text{concatenation of complementary multi-indices.}$$
 (7)

For each  $\ell$  , we define a list of 'pivot rows'  $\underline{\mathcal{I}}_{\ell} \subseteq \underline{\mathcal{I}}_{\ell}$  and list of 'pivot columns'  $\underline{\mathcal{I}}_{\ell+1} \subseteq \underline{\mathcal{I}}_{\ell+1}$ Also:  $\underline{\mathcal{I}}_{\circ} = \underline{\mathcal{I}}_{\ell+1} = \{(\ )\}$  , with () = empty tuple.  $\underline{\mathcal{I}}_{\ell}$  and  $\underline{\mathcal{I}}_{\ell}$  are lists of lists: e.g. , for  $\underline{\mathcal{L}} = 5$  ,  $\underline{\mathcal{S}} = \{\ \circ, 1\}$ 

Now define zero-, one-, and two-dimensional slices of input tensor  $\digamma$ : (k-dimensional slice has k free indices)

A(I,1)

'pivot matrix' 🦒 (zero-dimensional slice):

3-leg tensor  $\mathsf{T}_{\ell}$  (one-dimensional slice, with free index  $\mathsf{G}_{\ell}$ ):

$$T_{\ell} = F\left(T_{\ell-1}, S_{\ell}, T_{\ell+1}\right) \quad \text{with elements} \qquad [T_{\ell}]_{i\sigma j} \equiv F_{i \oplus (\sigma) \oplus j} = \frac{1}{\prod_{i = \sigma} \prod_{j =$$

For fixed 6 , we define the matrix  $[T_\ell^\sigma]_{ij} \equiv [T_\ell]_{i\sigma j}$ 

4-leg tensor 
$$\Pi_{\ell}$$
 (two-dimensional slice, with free indices  $\sigma_{\ell}$ ,  $\sigma_{\ell+1}$ ): 
$$\Pi_{\ell} = F(\mathcal{I}_{\ell-1}, \mathcal{S}_{\ell}, \mathcal{S}_{\ell+1}, \mathcal{J}_{\ell+2}) \text{ elements } [\Pi_{\ell}]_{i\sigma\sigma'j} \equiv F_{i\oplus(\sigma,\sigma')\oplus j} = \frac{1}{|\Pi_{\ell}|} \frac{1}{|\sigma_{\ell}|} \frac{1}{|$$

With these definitions, the 'TCI approximation'  $\stackrel{\longleftarrow}{\mathsf{F}}$  of  $\stackrel{\longleftarrow}{\mathsf{F}}$  is defined as

$$F_{\sigma} \approx \widetilde{F}_{\sigma} = T_1^{\sigma_1} P_1^{-1} \cdots T_{\ell}^{\sigma_{\ell}} P_{\ell}^{-1} T_{\ell+1}^{\sigma_{\ell+1}} \cdots P_{\mathcal{L}-1}^{-1} T_{\mathcal{L}}^{\sigma_{\mathcal{L}}}, \tag{12a}$$

with independent sums over all row multi-indices  $i_{\ell} \in \mathcal{I}_{\ell}$  and column multi-indices  $j_{\ell+1} \in \mathcal{J}_{\ell+1}$ , for  $\ell=1,...$ ,

- (12) defines the 'TCI form'. It is fully defined by  $\sqrt{\phantom{a}}$  and  $\sqrt{\phantom{a}}$  tensors, i.e. by slices of  $\sqrt{\phantom{a}}$ . These can be constructed if
- (i) one knows the pivot lists  $\{ \mathcal{I}_{\ell}, \mathcal{I}_{\ell+1} \mid \ell=1, ..., \ell-1 \}$  and
- (ii) can read out / evaluate / compute the input tensor 🗧 for any configuration ₹

Any tensor train can be converted exactly into a TCI form (see TCI. ). (Ferrally 2025, Sec. 4.5.1)

TCI.8

The interpolation properties of TCI Ansatz rely on nesting conditions satisfied by its pivot lists. Below, we define these nesting conditions. Their relevance will become clear in subsequent sections.

For any bond &

•  $\mathcal{I}_{\ell}$  is nested w.r.t. to  $\mathcal{I}_{\ell-1}$ , denoted by  $\mathcal{I}_{\ell-1} \subset \mathcal{I}_{\ell}$  if  $\mathcal{I}_{\ell} \subseteq \mathcal{I}_{\ell-1} \times \mathcal{S}_{\ell}$  (1)  $\frac{\mathcal{I}_{\ell} = ((1))}{\mathcal{I}_{1} = ((1))}$ 

Then, removing last index of any element of  $\mathcal{I}_{\ell}$  an yields element of  $\mathcal{I}_{\ell-1}$  , i.e.  $\mathcal{I}_{\ell}$  'descends from'  $\mathcal{I}_{\ell}$  's  $\mathcal{S}_{\ell}$  .

 $\underline{\mathcal{I}_\ell}$  left-nested row-pivot lists  $\mathcal{I}_2 = ((1,0),(1,1))$   $\mathcal{I}_3 = ((1,1,0),(1,0,1))$  $\bar{\mathcal{I}}_4 = ((1, 1, 0, 0))$ 

$$[P_{\ell}]_{ij} = F_{i \oplus j} =$$

Moreover, then  $P_{\ell}$  is a slice of  $T_{\ell}$ :  $[P_{\ell}]_{ij} = F_{i \oplus j} = \frac{\log \ell}{\prod_{i = 1}^{l} \prod_{j = 1}^{l} \prod_{j \in I} \prod_{j \in$ 

In the property of the second strong from the second strong of the second strong from the i.e.  $\mathcal{I}_{\ell}$  'descends from'  $\mathcal{S}_{\ell} \times \mathcal{I}_{\ell+1}$ 

$$[P_{\ell-1}]_{ij} = F_{i \oplus j} = \prod_{i \in \mathcal{I}_{\ell-1}} \prod_{j \in \mathcal{I}_{\ell}} \prod_{j \in$$

Pivots are 'left-nested' up to  $\mbox{\it l}$  if  $\mbox{\it I}_0 < \mbox{\it I}_1 < \cdots < \mbox{\it I}_\ell$ 

$$\mathcal{I}_0 < \mathcal{I}_1 < \dots < \mathcal{I}_\ell$$

Pivots are 'right-nested' up to  $\begin{cases} \ell \end{cases}$  if  $\begin{cases} \mathcal{J}_\ell > \mathcal{J}_{\ell+1} > \cdots > \mathcal{J}_{\mathcal{L}+1} \end{cases}$ 

$$\mathcal{J}_{\ell} > \mathcal{J}_{\ell+1} > \dots > \mathcal{J}_{\mathcal{L}+1}$$

(5)

Pivots are 'fully left-nested' if they are left-nested up to  $\ \mathcal{L}$  - I  $\mathcal{I}_0 < \mathcal{I}_1 < \dots < \mathcal{I}_{\mathcal{L}-1} \,,$ 

$$L_0 < L_1 < \cdots < L_{\mathcal{L}-1},$$

Pivots are 'fully right-nested' if they are right-nested up to 2

$$\mathcal{J}_2 > \mathcal{J}_{\ell+2} > \dots > \mathcal{J}_{\mathcal{L}+1} \tag{6}$$

Pivots are 'fully nested' if they are fully left- and right-nested.

(7)

Pivots are nested w.r.t.  $\sqrt{\phantom{a}}$  if they are left-nested up to  $\ell_{-1}$  and right-nested up to  $\ell_{+1}$ Then, the TCI form is exact on the one-dimensional slice  $T_{\mu}$ 

 $\widetilde{F}_{i\oplus(\sigma)\oplus i} = [T_{\ell}]_{i\sigma i} = F_{i\oplus(\sigma)\oplus i}, \quad \forall i \in \mathcal{I}_{\ell-1}, \ \sigma \in \mathbb{S}_{\ell}, \ j \in \mathcal{J}_{\ell+1}$ 



• Pivots are nested w.r.t.  $\Pi_{\ell}$  if they are left-nested up to  $\ell$ -1 and right-nested up to  $\ell$ +2 (9)



If pivots are <u>fully</u> nested, TCI form is exact on every  $\mathcal{T}_{\ell}$  and  $\mathcal{P}_{\ell}$  , i.e. on all slices used to construct it, thus it is an interpolation. (10)

Properties (8) and (10) are very important. For a proof, see TCI.10.

TCI.9

Goal: obtain TCI approximation  $\hat{F}$  of given tensor F at specified tolerance,  $\|F - \hat{F}\| < \mathcal{I}$ , the finding a minimal set of suitable pivots. (1) by finding a minimal set of suitable pivots.

Basic 2-site TCI algorithm [Fernandez2025, Sec. 4.3.1]

(1) Initialization: start with any configuration  $\hat{\sigma}$ 

$$\mathcal{I}_{\ell} = \left\{ \left( \hat{\sigma}_{\ell}, \dots, \hat{\sigma}_{\ell} \right) \right\}, \quad \mathcal{I}_{\ell+1} = \left\{ \left( \hat{\sigma}_{\ell+1}, \dots, \hat{\sigma}_{\ell} \right) \right\} \quad (2)$$

(2) Sweeping back and forth over  $\ell = 1, ..., L-1$ , perform the following update at each  $\ell$ :

Construct  $\Pi_{\ell} \equiv F(\mathcal{I}_{\ell-1}, \mathbb{S}_{\ell}, \mathbb{S}_{\ell+1}, \mathcal{J}_{\ell+2})$ , view it as a matrix  $F(\underbrace{\mathcal{I}_{\ell-1} \times \mathbb{S}_{\ell}}, \underbrace{\mathbb{S}_{\ell+1} \times \mathcal{J}_{\ell+2}})$ , and prrLU-factorize it:  $\mathbf{I} \qquad \qquad \mathbf{I} \qquad \mathbf$ 

$$[\Pi_{\ell}]_{i_{\ell-1}\sigma_{\ell}\sigma_{\ell+1}j_{\ell+2}} \approx [T_{\ell}^{'\sigma_{\ell}}]_{i_{\ell-1}j_{\ell+1}'} (P_{\ell}')_{j_{\ell+1}'i_{\ell}'}^{-1} [T_{\ell+1}^{'\sigma_{\ell+1}}]_{i_{\ell}'j_{\ell+2}} = \widehat{\Pi}_{\ell}$$

$$\stackrel{\mathcal{I}}{=} \mathbb{I} \quad \mathcal{J} \leq \mathbb{J}$$

$$\text{recall (TCI.3.18)}$$

$$\frac{\Pi_{\ell}}{i_{\ell-1}} \underbrace{\frac{T'_{\ell}}{\sigma_{\ell}} \underbrace{\frac{P'_{\ell-1}}{j_{\ell+1}} \frac{T'_{\ell+1}}{j_{\ell+2}}}_{i_{\ell-1}} \approx \underbrace{\frac{T'_{\ell}}{j_{\ell+1}} \underbrace{\frac{P'_{\ell-1}}{j_{\ell+1}} \frac{T'_{\ell+1}}{j_{\ell+2}}}_{j_{\ell+2}}, \qquad \underbrace{\frac{i}{\mathbb{I}} \underbrace{\frac{C}{J}}_{J} \underbrace{\frac{j}{P^{-i}}}_{I} \underbrace{\frac{k}{J}}_{J}}_{I} \qquad (4)$$

Use new pivot lists to update  $P_{\ell}$  ,  $T_{\ell}$  ,  $T_{\ell+1}$  .

(3) Iterate step (2) until specified tolerance or specified maximum bond dimension is reached.

Eq. (5) shows: new pivot lists  $\mathcal{I}'_{\ell}$  .  $\mathcal{I}'_{\ell+1}$ satisfy nesting conditions (by construction).

However, full nesting (if previously present), may be broken:

one may have  $\mathcal{I}_{\ell} < \mathcal{I}_{\ell+1}$  but not  $\mathcal{I}'_{\ell} < \mathcal{I}_{\ell+1}$  , and one may have  $\mathcal{J}_{\ell} > \mathcal{J}_{\ell+1}$  but not  $\mathcal{J}_{\ell} > \mathcal{J}'_{\ell+1}$  . Reason: since  $\mathbf{I}_{\ell}$ ,  $\mathbf{J}'_{\ell+1}$  play no role in the construction of  $\mathbf{I}'_{\ell}$ ,  $\mathbf{J}'_{\ell+1}$ , some pivots from the former may be missing in the latter; and removal of pivot generally can cause breaking of nesting conditions.

Remedy: if full nesting is desired, it can be restored at the end using 1-site TCI algorithm (see TCI.11).

Important fact: if pivots are nested w.r.t.  $\mathbb{T}_{\ell}$  (left-nested up to  $\ell$ -1 and right-nested up to  $\ell$ +2), then

$$\left[\Pi_{\ell} - \widetilde{\Pi}_{\ell}\right]_{i_{\ell-1}\sigma_{\ell}\sigma_{\ell+1}j_{\ell+2}} = \left[F - \widetilde{F}\right]_{i_{\ell-1}\sigma_{\ell}\sigma_{\ell+1}j_{\ell+2}} \quad \text{for all} \quad \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} \in \mathbf{S}_{\ell} \times \mathbf{S}_{\ell+1} \quad \text{(proof: see TCI.10)} \quad \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} = \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} + \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} + \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} + \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} = \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} + \mathbf{\sigma}_{\ell+1} + \mathbf{\sigma}_{\ell} \cdot \mathbf{\sigma}_{\ell+1} + \mathbf{\sigma$$

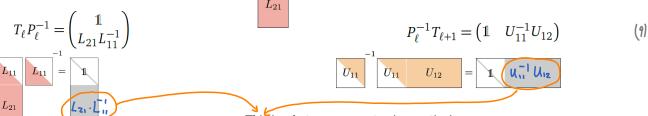
Thus, error of approximating  $\prod_{\ell}$  by  $\widehat{\prod}_{\ell}$  is also error, on this 2-dimensional slice, of approximating  $F_{\widehat{\epsilon}}$  by  $\widehat{F}_{\widehat{\epsilon}}$ . The above algorithm chooses pivots in order to minimize this error.

When factorizing  $\prod_{\ell} \simeq \prod_{\ell} \prod_{\ell \neq 1} V_{\ell}$  via prrLU, the ingredients on the right are constructed as follows: cf. (TCI.6.14-16)

$$\overline{\Pi}_{\ell} = \begin{bmatrix} A_{11} & A_{12} \\ A_{21} & A_{22} \end{bmatrix} \simeq \widetilde{\Pi}_{\ell} = T_{\ell}(P_{\ell})^{-1} T_{\ell+1} = LDU = \begin{pmatrix} P_{\ell} & L_{11}DU_{12} \\ L_{21}DU_{11} & L_{21}DU_{12} \end{pmatrix} = \begin{bmatrix} L_{11} & D \\ L_{21} & D \end{bmatrix} = \begin{bmatrix} L_{11} & D \\ L_{21} & D \end{bmatrix}$$

$$T_{\ell} = \begin{pmatrix} L_{11}DU_{11} \\ L_{21} & D \end{pmatrix} \qquad T_{\ell+1} = \begin{pmatrix} L_{11}DU_{11} & L_{11}DU_{12} \end{pmatrix} \qquad (3)$$

$$P_{\ell} = L_{11}DU_{11}$$
  $T_{\ell} = \begin{pmatrix} L_{11}DU_{11} \\ L_{21}DU_{11} \end{pmatrix}$   $T_{\ell+1} = \begin{pmatrix} L_{11}DU_{11} & L_{11}DU_{12} \end{pmatrix}$   $T_{\ell+1} = \begin{pmatrix} L_{11}DU_{11} & L_{11}DU_{12} \end{pmatrix}$   $T_{\ell+1} = \begin{pmatrix} L_{11}DU_{11} & L_{11}DU_{12} \end{pmatrix}$ 



This is what one computes in practice! Since  $\angle_{II}$  and  $\bigvee_{II}$  are lower- or upper triangular, their inverses can be computed in stable manner using forward/backward substitution

Pivot update method 'reset' vs. 'accumulative'

'reset' mode: replace all pivots in  $\mathcal{I}_{\ell}$ ,  $\mathcal{I}_{\ell+1}$  by all pivots in  $\mathcal{I}_{\ell}$ ,  $\mathcal{I}_{\ell+1}$  (as described above) pro: can discard bad pivots; con: can break nesting conditions

'accumulative' mode: don't discard pivots from  $\mathcal{I}_{\ell_{\ell}}\mathcal{I}_{\ell_{\ell_{\ell}}}$ , just add new ones from  $\mathcal{I}_{\ell_{\ell}}\mathcal{I}_{\ell_{\ell_{\ell}}}$ typically one at a time.

pro: can break nesting conditions; con: cannot discard bad pivots.

Both modes have runtimes  $\mathcal{O}$  ( ) . Accumulative:  $\mathcal{O}$  ( ) per update, needs updates to reach rank  $\chi$  .

Reset: O() per update, but typically converges within a few updates, independent of  $\chi$ 

Pivot <u>search</u> One can use full search, rook search, or block rook search. 0( ), 0( ), 0( Scaling: (works well in combination with reset mode)

## Adding global pivots

In addition to using pivots found via 2-site TCI algorithm, it may be useful to add 'global' pivots based on 'outside' information, such as:

- knowledge of configurations where 😽 is very large;
- doing TCI on a tensor  $\mathcal{F}_1$  that is very similar to a tensor  $\mathcal{F}_1$  whose TCI unfolding  $\hat{\mathcal{F}}_i$  is already known Strategy for adding global pivots:
- split each  $\vec{\sigma}$  as  $\vec{\sigma} = i_{\ell} \oplus j_{\ell+1}$  for all  $\ell = i_{\ell+1}, \ell-1$ ,
- add these  $i_{\ell}$  ,  $j_{\ell+1}$  to the pivot lists  $\mathcal{I}_{\ell}$  ,  $\mathcal{I}_{\ell+1}$
- do prrLU on all pivot matrices ? to discard any spurious pivots
- perform a few sweeps using 2-site TCI in reset mode to stabilize the pivot lists.

## **Ergodicity**

TCI is based on exploration of configuration space, so can encounter ergodicity issues -remaining stuck in subpart of configuration space and not visiting other relevant parts.

If one notices such issues, initialize pivot search with suitably chosen global pivots: Examples:

- Very sparse tensors, where TCI might miss some nonzero entries; remedy: add global pivots for a list of nonzero entries.
- Tensors with discrete symmetries, where exploration may get stuck in one symmetry sector; remedy: add one global pivot per symmetry sector.
- Multivariate functions with very narrow peaks;
   remedy: add global pivots corresponding to peak maxima.

## <u>1-site TCI algorithm</u> [Fernandez2025, Sec. 4.4.1]

Useful for (i) compressing a given TCI to smaller rank; (ii) restoring full nesting; (iii) improve pivots at lower computational cost than 2-site TCI. Limitation: cannot increase bond dimension.

Input: a TT in TCI form. General strategy: Sweep back and forth, compressing  $T_{\ell}$  using prrLU.

Forward sweep: view  $T_\ell \equiv F(\mathcal{I}_{\ell-1}, \mathbb{S}_\ell, \mathcal{J}_{\ell+1})$  as a matrix  $F(\mathcal{I}_{\ell-1} \times \mathbb{S}_\ell, \mathcal{J}_{\ell+1})$ , and prrLU-factorize it:

After complete forward sweep, pivots are fully left-nested:  $\mathcal{I}_0 < \cdots < \mathcal{I}_{\mathcal{L}-1}$ 

Backward sweep: view  $T_{\ell} \equiv F(\mathcal{I}_{\ell-1}, \mathbb{S}_{\ell}, \mathcal{J}_{\ell+1})$  as a matrix  $F(\mathcal{I}_{\ell-1}, \mathbb{S}_{\ell} \times \mathcal{J}_{\ell+1})$ , and prrLU-factorize it:

$$T_{\ell} = \frac{T_{\ell}}{i_{\ell-1}} \int_{\sigma_{\ell}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell}} \int_{i_{\ell-1}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell+1}^{\prime}} \int_{j_{\ell+1}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell+1}^{\prime}} \int_{j_{\ell}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell+1}^{\prime}} \int_{j_{\ell}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell+1}^{\prime}} \int_{j_{\ell}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell+1}^{\prime}} \int_{j_{\ell}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell+1}^{\prime}} \int_{j_{\ell}}^{\tau_{\ell}} \frac{P_{\ell-1}^{\prime-1}}{j_{\ell}^{\prime}} \int_{j_{\ell}}^{\tau_{\ell}} \frac{P_{\ell-1}^$$

$$\begin{aligned} \text{indices of } \overline{\mathsf{T}_{\pmb{\ell}}} : & i_{\ell-l} \in \mathcal{I}_{\ell-l} \\ & \leq j_{\ell+1} \in \mathbb{S}_{\ell} \times \mathcal{J}_{\ell+1} \end{aligned} \qquad \begin{aligned} \text{pivot indices: } i_{\pmb{\ell}-l}' \in \mathcal{I}_{\ell-1}' \subseteq \mathcal{I}_{\ell-1} \\ & j_{\ell}' \in \mathcal{I}_{\ell}' \subset \mathbb{S}_{\ell} \times \mathcal{J}_{\ell+1} \end{aligned} \qquad \Rightarrow \qquad \mathcal{J}_{\ell}' > \mathcal{J}_{\ell+1} \\ & \text{nested!} \end{aligned}$$

After complete backward sweep, pivots are fully right-nested:  $\mathcal{J}_2 > \dots > \mathcal{J}_{\mathcal{L}+1}$ 

Backward sweeping preserves left-nesting only if taking the subset  $\mathcal{I}'_{\ell-1}\subseteq\mathcal{I}_{\ell-1}$  does not remove any pivots. To achieve full nesting, do one more forward sweep at same tolerance. This preserves right-nesting, since all bond dimensions already meet the tolerance, so last forward sweep removes no pivots from  $\mathcal{I}_{\ell+1}$  for  $\ell=1,\ldots,\ell-1$ 

## 0-site TCI algorithm [Fernandez2025, Sec. 4.4.2]

Input: given TT in TCI format. Sweep through pivot matrices, prrLU decomposing each to yield updated pivot lists  $\mathcal{L}'_{\ell}$ ,  $\mathcal{T}'_{\ell+\ell}$  that replace  $\mathcal{L}_{\ell}$ ,  $\mathcal{T}'_{\ell+\ell}$ . Main usage: improve conditioning of  $\mathcal{L}_{\ell}$  by removing 'spurious' pivots. Breaks nesting conditions.