

Quantum Computation

ZX calculus — reference sheet

(December 2022)

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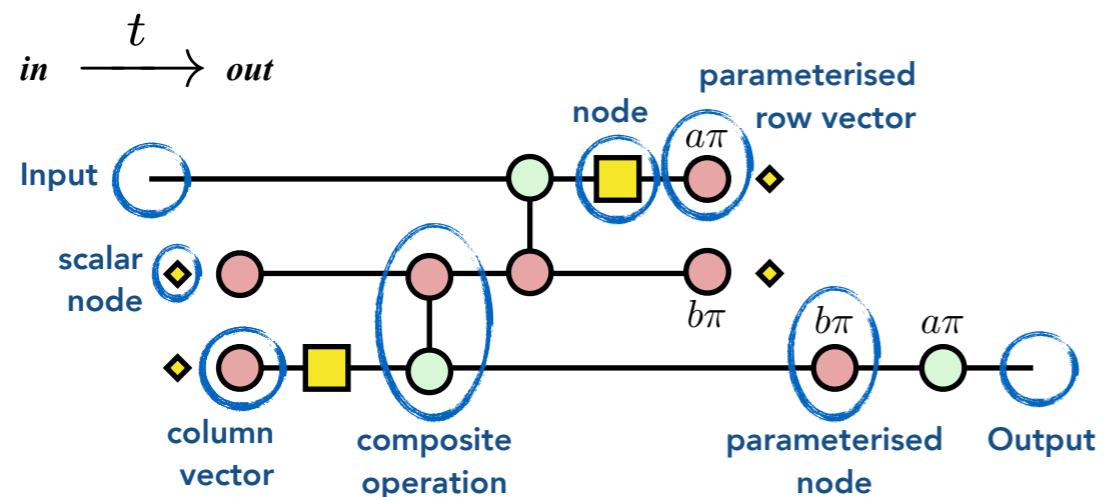
This reference sheet is intended to present the most important (i.e., the most often used) rewrites for the ZX calculus, in the 'well-tempered' normalisation [[arXiv:2006.02557](https://arxiv.org/abs/2006.02557)].

This set of rewrites is incomplete: there are more rewrites, which you can use (and which are necessary to use) to be able to show when any two diagrams represent the same state or operator. The rules presented here are intended to cover essentially any 'simple' calculation in the ZX calculus.

Using diagrammatic notation

The meaning of a diagram

The ZX calculus uses diagrams to represent sequences of transformations on column vectors (and on tensor products of column vectors)



Each diagram represents a matrix, described through a "semantic map" $\llbracket \quad \rrbracket : (\text{diagrams}) \rightarrow (\text{matrices})$

$$\llbracket m \left\{ \begin{array}{c} \text{---} \\ \text{---} \end{array} \right\} n \rrbracket = \begin{bmatrix} \text{some} \\ 2^n \times 2^m \\ \text{matrix} \end{bmatrix} \quad \text{(function application order — right to left)}$$

Composing diagrams

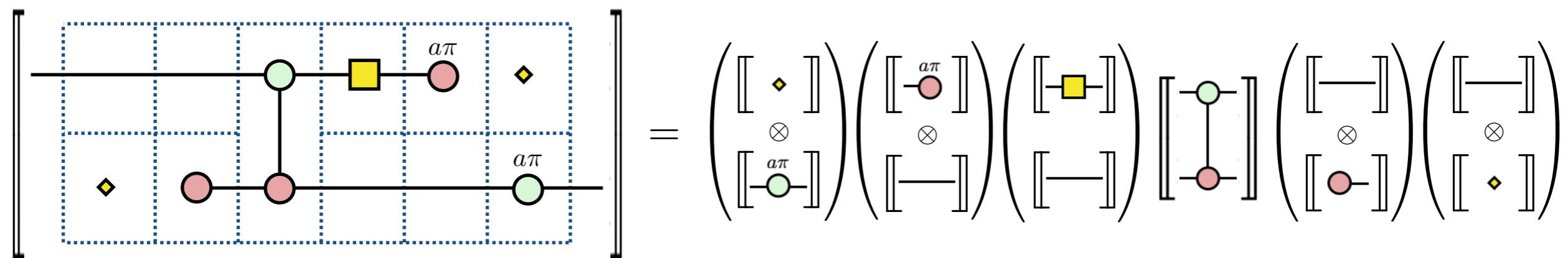
- In parallel ...

$$\llbracket \begin{array}{c} \text{---} \\ D \\ \text{---} \\ \text{---} \\ D' \\ \text{---} \end{array} \rrbracket = \llbracket \begin{array}{c} \text{---} \\ D \\ \text{---} \end{array} \rrbracket \otimes \llbracket \begin{array}{c} \text{---} \\ D' \\ \text{---} \end{array} \rrbracket$$

- In sequence ...

$$\begin{aligned} \llbracket \begin{array}{c} \text{---} \\ D \\ \text{---} \\ \text{---} \\ D' \\ \text{---} \end{array} \rrbracket &= \llbracket \begin{array}{c} \text{---} \\ D' \\ \text{---} \end{array} \rrbracket \circ \llbracket \begin{array}{c} \text{---} \\ D \\ \text{---} \end{array} \rrbracket \\ &\quad \text{function application order (right to left)} \end{aligned}$$

We multiply the matrices / vectors in the same **sequence** as in the diagram — **but** from right to left
(as this is how we transform column vectors)



Essential ZX for circuits

Regarding normalisation: to allow for simple rewrite rules, we define the parametrised nodes " $\theta \text{---}$ ", " $\theta \text{---}$ ", etc. to have normalisation > 1 . We use the " \diamond " scalar node as a normalising factor, to allow us to easily represent unit vectors.

$$\begin{aligned} \left[\begin{array}{c} \diamond \\ \bullet \end{array} \right] &= \langle 0 | 0 \rangle = 1 = \langle + | + \rangle = \left[\begin{array}{c} \diamond \\ \bullet \end{array} \right] \\ \left[\begin{array}{c} \pi \\ \bullet \end{array} \right] &= \langle 1 | 1 \rangle = 1 = \langle - | - \rangle = \left[\begin{array}{c} \pi \\ \bullet \end{array} \right] \\ \left[\begin{array}{c} \pi \\ \bullet \end{array} \right] &= \langle 0 | 1 \rangle = 0 = \langle + | - \rangle = \left[\begin{array}{c} \pi \\ \bullet \end{array} \right] \end{aligned}$$

Single-qubit states and tests

$$\begin{aligned} \left[\begin{array}{c} \diamond \\ \bullet \end{array} \right] &= |0\rangle \\ \left[\begin{array}{c} \diamond \\ \pi \bullet \end{array} \right] &= |1\rangle \\ \left[\begin{array}{c} \diamond \\ a\pi \bullet \end{array} \right] &= |a\rangle \text{ for } a \in \{0, 1\} \end{aligned} \quad \begin{aligned} \left[\begin{array}{c} \bullet \\ \diamond \end{array} \right] &= \langle 0| \\ \left[\begin{array}{c} \pi \bullet \\ \diamond \end{array} \right] &= \langle 1| \\ \left[\begin{array}{c} a\pi \bullet \\ \diamond \end{array} \right] &= \langle a| \text{ for } a \in \{0, 1\} \end{aligned}$$

$$\begin{aligned} \left[\begin{array}{c} \diamond \\ \bullet \end{array} \right] &= |+\rangle \\ \left[\begin{array}{c} \diamond \\ \pi \bullet \end{array} \right] &= |-\rangle \\ \left[\begin{array}{c} \diamond \\ a\pi \bullet \end{array} \right] &= \begin{cases} |+\rangle, & \text{if } a = 0 \\ |-\rangle, & \text{if } a = 1 \end{cases} \end{aligned} \quad \begin{aligned} \left[\begin{array}{c} \bullet \\ \diamond \end{array} \right] &= \langle +| \\ \left[\begin{array}{c} \pi \bullet \\ \diamond \end{array} \right] &= \langle -| \\ \left[\begin{array}{c} a\pi \bullet \\ \diamond \end{array} \right] &= \begin{cases} \langle +|, & \text{if } a = 0 \\ \langle -|, & \text{if } a = 1 \end{cases} \end{aligned}$$

Single-qubit gates

$$\begin{aligned} \left[\begin{array}{c} \text{---} \end{array} \right] &= \left[\begin{array}{c} \text{---} \bullet \end{array} \right] = \left[\begin{array}{c} \bullet \text{---} \end{array} \right] = \mathbb{1} = |0\rangle\langle 0| + |1\rangle\langle 1| \\ \left[\begin{array}{c} \text{---} \square \end{array} \right] &= H = \frac{1}{\sqrt{2}} \begin{bmatrix} 1 & 1 \\ 1 & -1 \end{bmatrix} = |+\rangle\langle 0| + |-\rangle\langle 1| \\ \left[\begin{array}{c} \text{---} \bullet \\ \pi \end{array} \right] &= Z = \begin{bmatrix} 1 & 0 \\ 0 & -1 \end{bmatrix} = |0\rangle\langle 0| - |1\rangle\langle 1| \\ \left[\begin{array}{c} \text{---} \bullet \\ \pi \end{array} \right] &= X = \begin{bmatrix} 0 & 1 \\ 1 & 0 \end{bmatrix} = |+\rangle\langle +| - |-\rangle\langle -| \end{aligned}$$

Simple two-qubit entangled state

$$\left[\begin{array}{c} \diamond \\ \bullet \end{array} \right] = \left[\begin{array}{c} \diamond \\ \bullet \end{array} \right] = |\Phi^+\rangle = \frac{1}{\sqrt{2}}(|0,0\rangle + |1,1\rangle)$$

Two-qubit controlled gates

$$\begin{aligned} \left[\begin{array}{cc} \text{---} & \text{---} \\ \text{---} & \bullet \end{array} \right] &= \text{CNOT} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 0 & 1 \\ 0 & 0 & 1 & 0 \end{bmatrix} = |0\rangle\langle 0| \otimes \mathbb{1} \\ &\quad + |1\rangle\langle 1| \otimes X \\ \left[\begin{array}{cc} \text{---} & \text{---} \\ \text{---} & \square \end{array} \right] &= \text{CZ} = \begin{bmatrix} 1 & 0 & 0 & 0 \\ 0 & 1 & 0 & 0 \\ 0 & 0 & 1 & 0 \\ 0 & 0 & 0 & -1 \end{bmatrix} = |0\rangle\langle 0| \otimes \mathbb{1} \\ &\quad + |1\rangle\langle 1| \otimes Z \end{aligned}$$

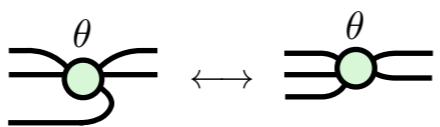
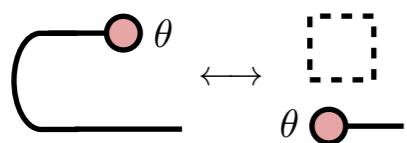
Wire bending

$$\begin{aligned} \left[\begin{array}{c} \bullet \\ \text{---} \end{array} \right] &= \left[\begin{array}{c} \bullet \\ \text{---} \end{array} \right] = \left[\begin{array}{c} \text{---} \\ \bullet \end{array} \right] = |0,0\rangle + |1,1\rangle \\ \left[\begin{array}{c} \text{---} \\ \bullet \end{array} \right] &= \left[\begin{array}{c} \text{---} \\ \bullet \end{array} \right] = \left[\begin{array}{c} \text{---} \end{array} \right] = \langle 0,0| + \langle 1,1| \end{aligned}$$

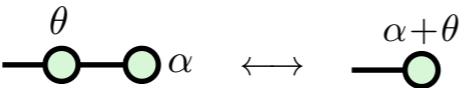
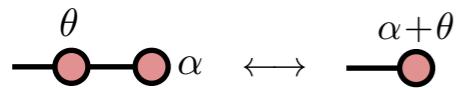
Essential ZX simplifications

Only the connectivity matters

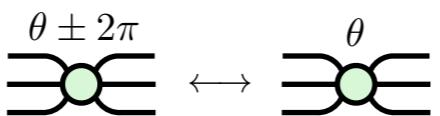
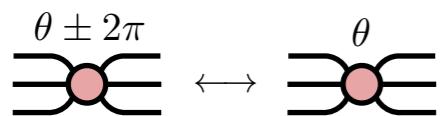
Because the vectors $|0\rangle$, $|+\rangle$, $|1\rangle$, and $|-\rangle$ (the vectors which we use to define the **red** and **green** nodes) all have real coefficients, we may change input wires to output wires and vice-versa when these are composed with a 'cup' or a 'cap' (a wire which bends back on itself).



As "only the connectivity matters", all of the rewrites below also hold true if they are flipped horizontally or vertically. For example, following the 'fusion' rewrites on the next page, we have:

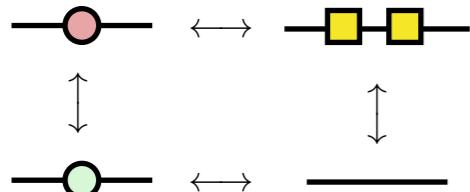


Angles



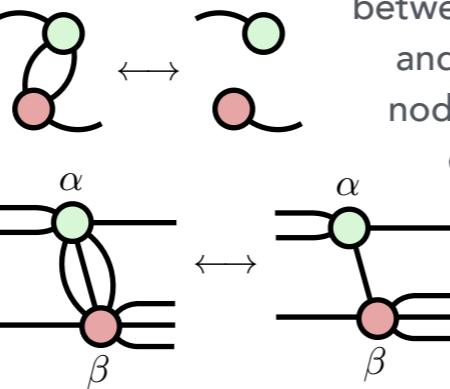
Both **red** and **green** nodes (with any # inputs or outputs): angles are evaluated mod 2π . If no angle is noted, the default angle is 0.

Identity



Phase-free degree-2 **red** and **green** nodes are identity operations; and Hadamards are self-inverse.

Hopf Law



Parallel edges between a **red** and a **green** node can be canceled in pairs

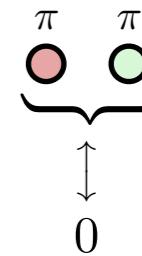
More scalar factors

$$\diamond \leftrightarrow 2^{-1/4}$$

This represents a scalar factor < 1 , which we use for renormalising factors

$$\{ \} \leftrightarrow \sqrt{2}$$

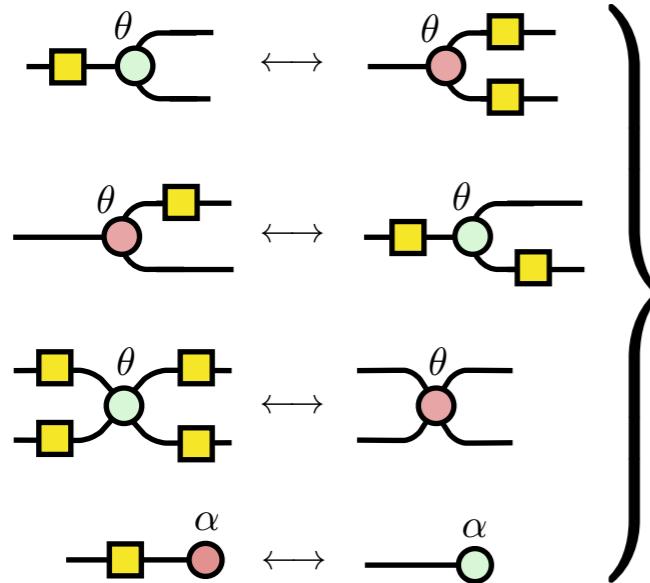
Each of these nodes represent a scalar factor > 1 , and so affect the normalisation



each of these nodes represent a scalar of 0: the larger diagram represents a zero vector or operator

Essential ZX transformations

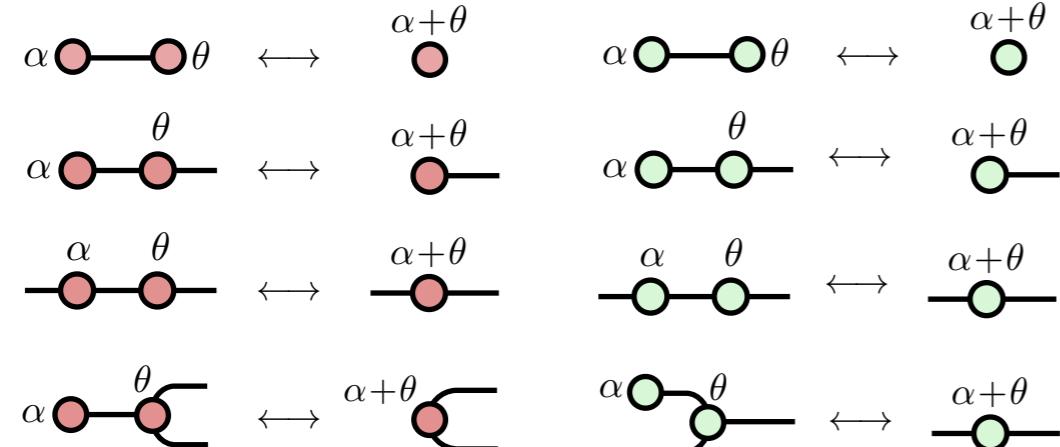
Colour change rules



A **red** or **green** node with Hadamards on some/all of its wires, can be replaced with a colour-switched node with Hadamards on the subset of its wires which did not have Hadamards on them

Every single ZX rewrite remains correct if we change every **red** node to a **green** node, and vice versa. This is precisely because they are related through Hadamad operations in the way described above.

Fusion rules



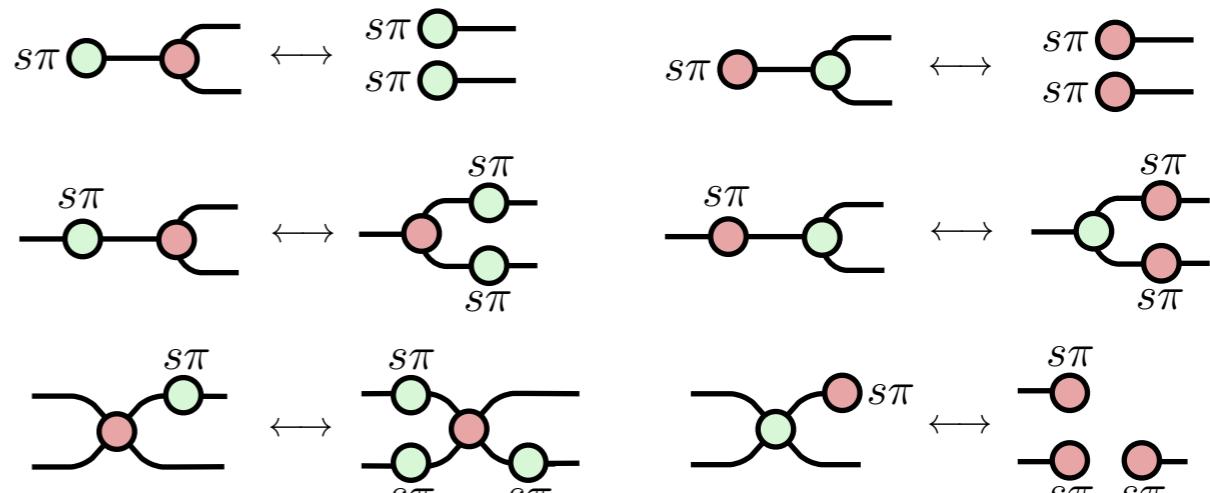
Red nodes
fuse (adding
their angles)

etc.

Green nodes
fuse (adding
their angles)

etc.

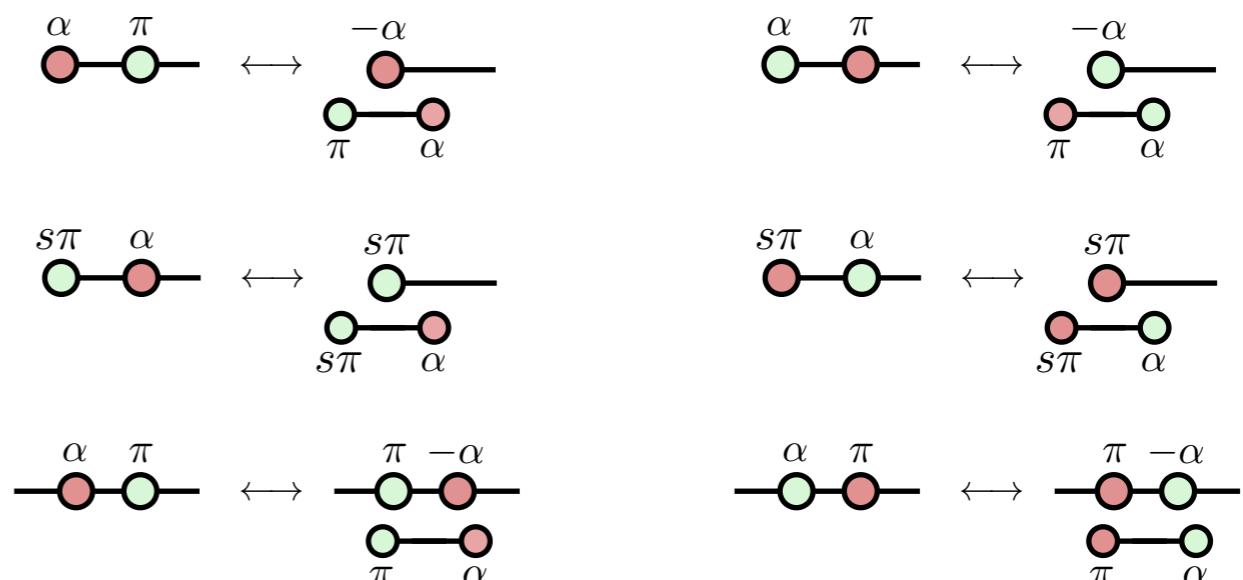
Copying rules



'**Green data**' — states, tests, and rotations with angle in $\pi\mathbb{Z}$ — are copied by phase-free **red nodes**

'**Red data**' — states, tests, and rotations with angle in $\pi\mathbb{Z}$ — are copied by phase-free **green nodes**

π -phase commutation & absorbtion



Up to a scalar factor, '**green π phases**' can commute past, absorb, or be absorbed by **a red a phase**, with a change in sign in **a** if it isn't absorbed

Up to a scalar factor, '**red π phases**' can commute past, absorb, or be absorbed by **a green a phase**, with a change in sign in **a** if it isn't absorbed

More ZX rewrites

Adjoints

To take the adjoint (also known as the Hermitian conjugate) of any diagram — e.g., to represent the inverse of a unitary matrix U , or to represent a measurement outcome $\langle \alpha |$ which corresponds to a quantum state $|\alpha\rangle$ — we flip the diagram horizontally (which yields the transpose of the original diagram) and then negate all of the angles of the nodes (which takes the complex conjugate), e.g.:

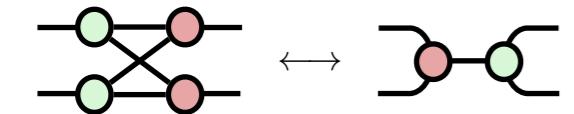
$$\begin{aligned} \left[\begin{array}{c} \text{yellow diamond} \\ \text{green circle} \end{array} \right] &= |\alpha\rangle = \frac{1}{\sqrt{2}}(|0\rangle + e^{i\theta}|1\rangle) \\ \Rightarrow \left[\begin{array}{c} -\theta \\ \text{green circle} \\ \text{yellow diamond} \end{array} \right] &= \langle \alpha | = \frac{1}{\sqrt{2}}(\langle 0| + e^{-i\theta}\langle 1|) \end{aligned}$$

Loops

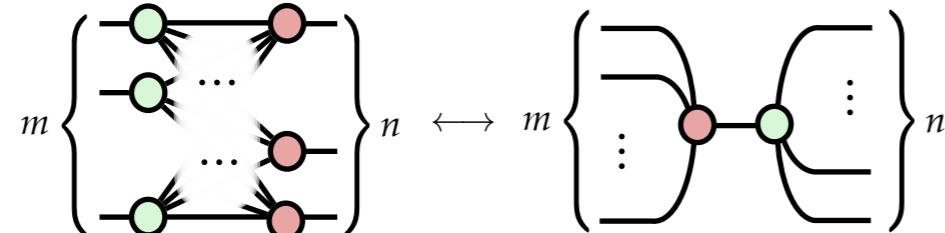


Each closed loop on any **red** or **green** node can be replaced by an isolated phase-free node (denoting a scalar factor of $\sqrt{2}$)

Bialgebra rule



In the general form shown here, each **green** node is connected to each **red** node

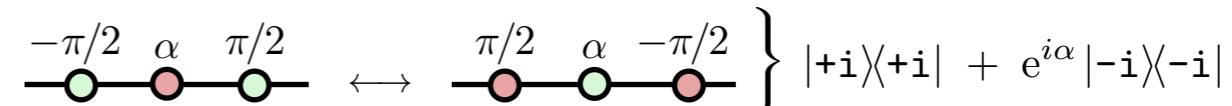


The bialgebra rule applies to sub-diagrams of phase-free **red** and **green** nodes, and describes the complementary relationship between the two bases $|0\rangle/|1\rangle$ and $|+\rangle/|-\rangle$

Y-axis states and rotations



The states on the Y-axis, are the only states that can be represented by either red nodes or green nodes, but require different angles to do so (and they are also related by a global scalar factor)



Rotations about the Y-axis involve interplay between X-axis rotations and Z-axis rotations, which can take multiple equivalent forms

$\pm\pi/2$ rotations and Hadamards

