

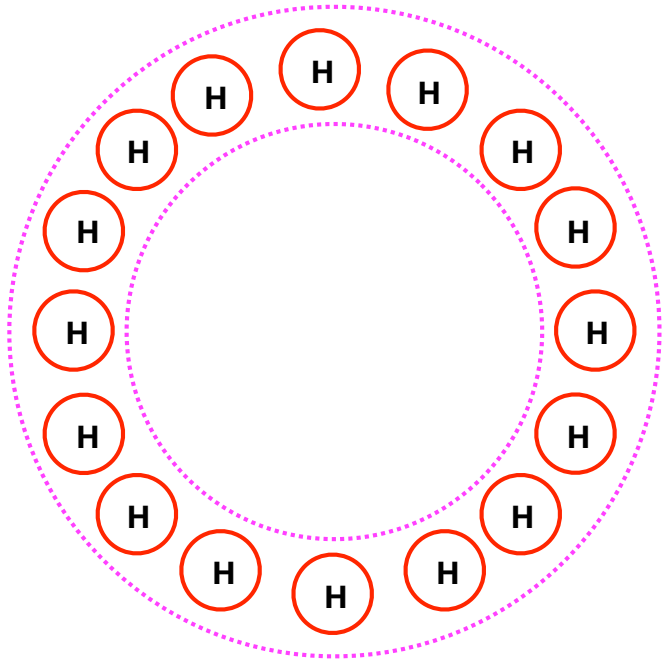
## *Quantum embedding in electronic structure theory*

### *Part 3: Exact embedding of localised orbitals for non-interacting electrons*

***Emmanuel Fromager***

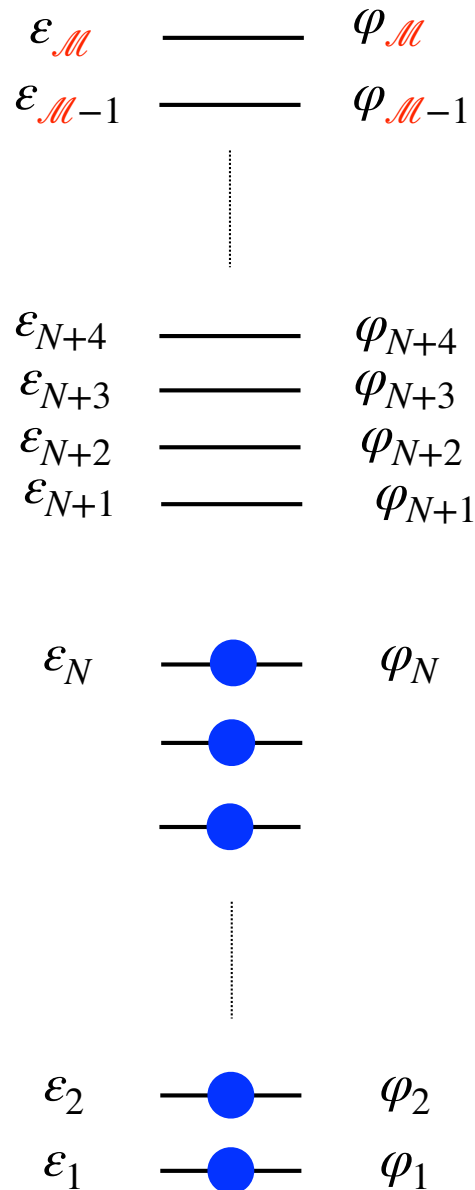
*Laboratoire de Chimie Quantique, Institut de Chimie de Strasbourg,  
Université de Strasbourg, Strasbourg, France.*

## Non-interacting delocalised representation



$$\hat{H} \equiv \sum_{PQ} \langle \varphi_P | \hat{h} | \varphi_Q \rangle \hat{a}_P^\dagger \hat{a}_Q$$

# Non-interacting (delocalised) molecular orbital representation

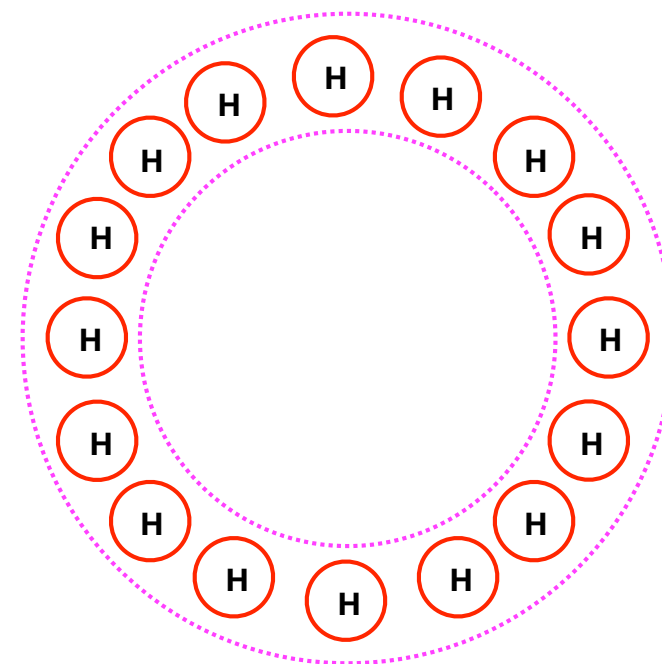
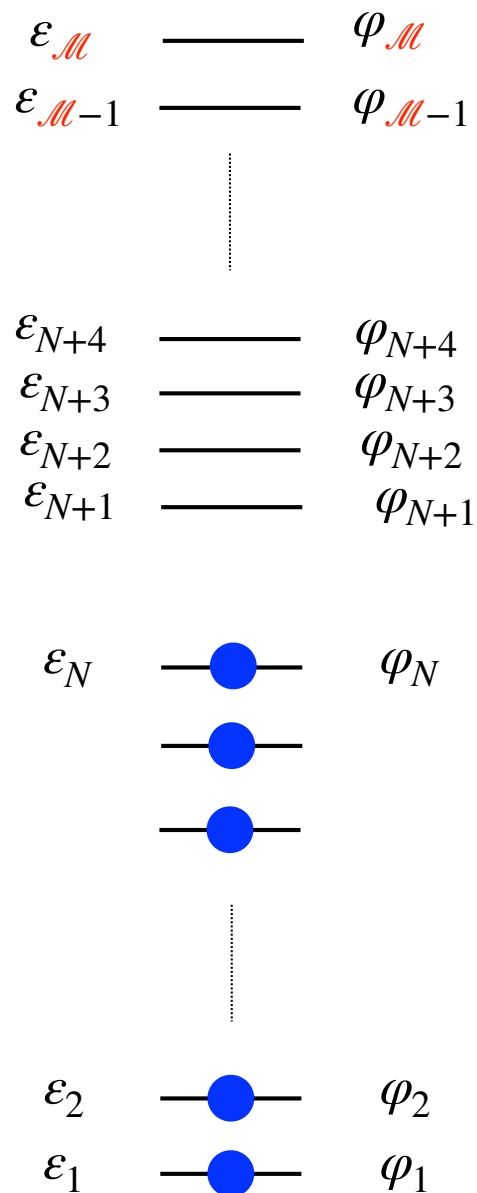


$$\hat{H} \equiv \sum_{PQ} \langle \varphi_P | \hat{h} | \varphi_Q \rangle \hat{a}_P^\dagger \hat{a}_Q$$

The molecular spin-orbitals are simply obtained by solving the **one-electron Schrödinger equation**

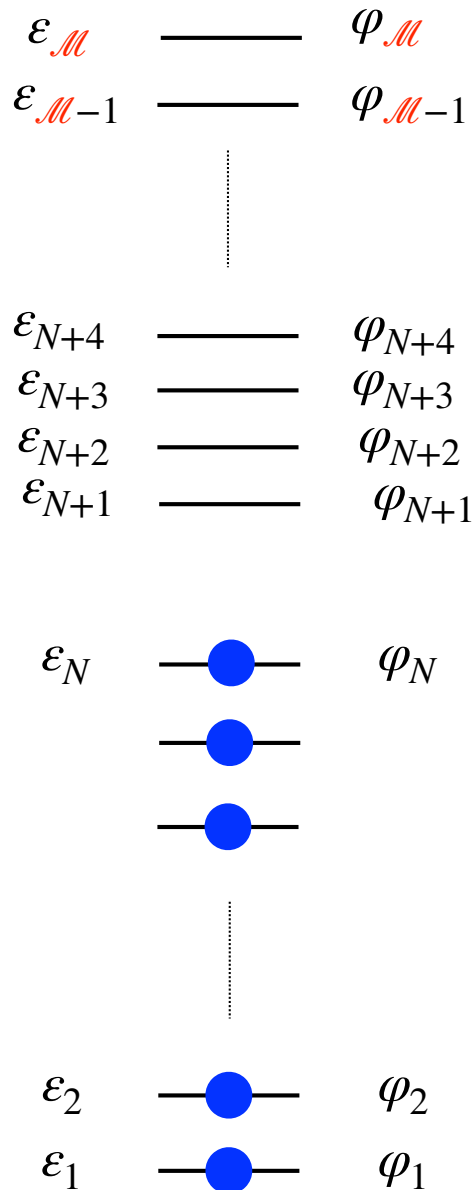
$$\hat{h}\varphi_Q(\mathbf{x}) = \varepsilon_Q \varphi_Q(\mathbf{x})$$

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$$\varphi_P(\mathbf{x}) = \sum_{\nu} C_{\nu P} \chi_{\nu}(\mathbf{x})$$

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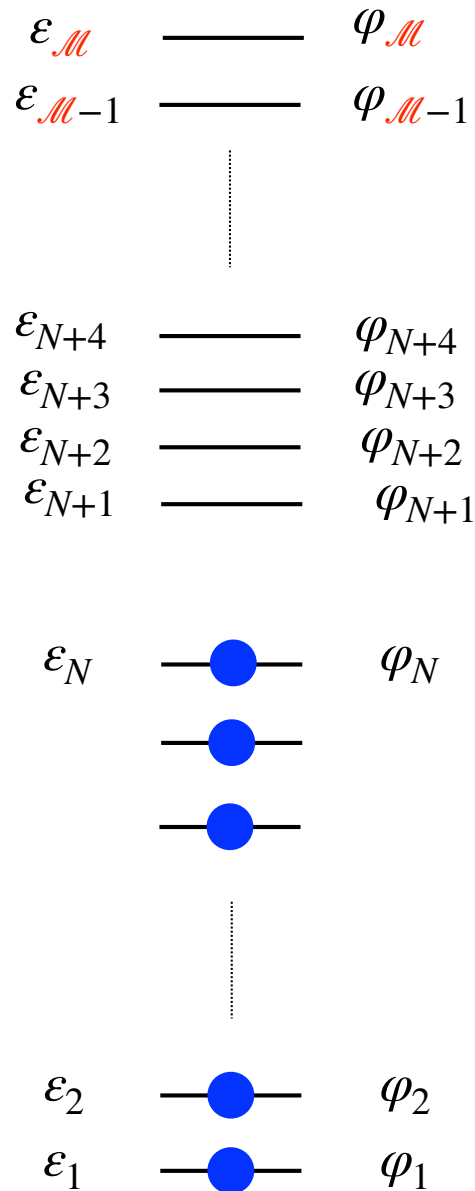
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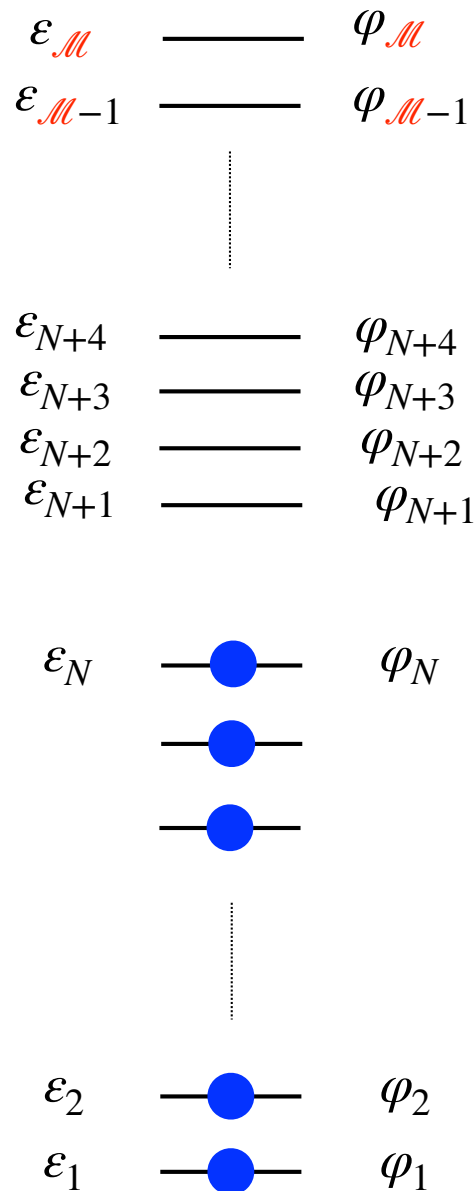
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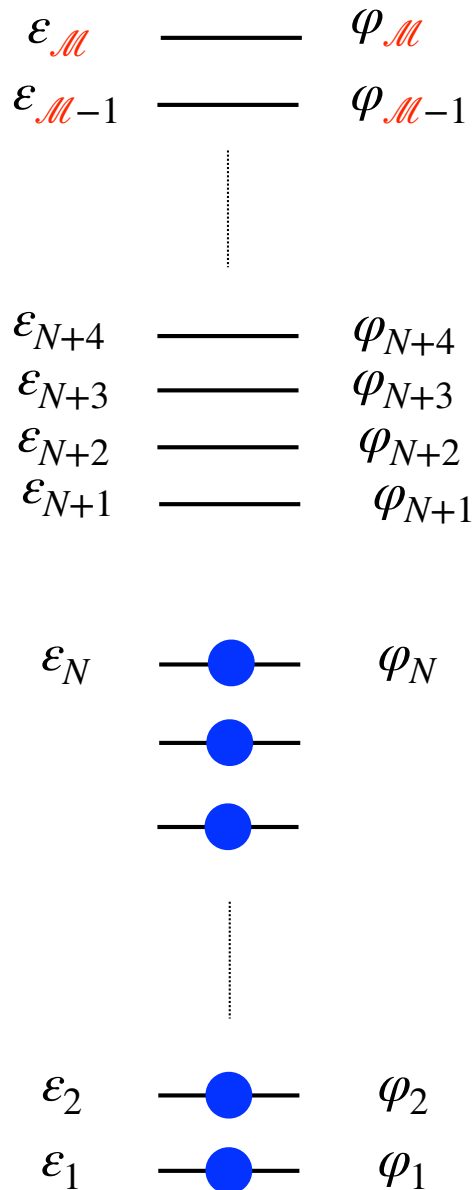


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The exact solutions to the non-interacting Schrödinger equation are Slater determinants  $\hat{a}_{P_1}^\dagger \hat{a}_{P_2}^\dagger \dots \hat{a}_{P_{N-1}}^\dagger \hat{a}_{P_N}^\dagger | \text{vac} \rangle$

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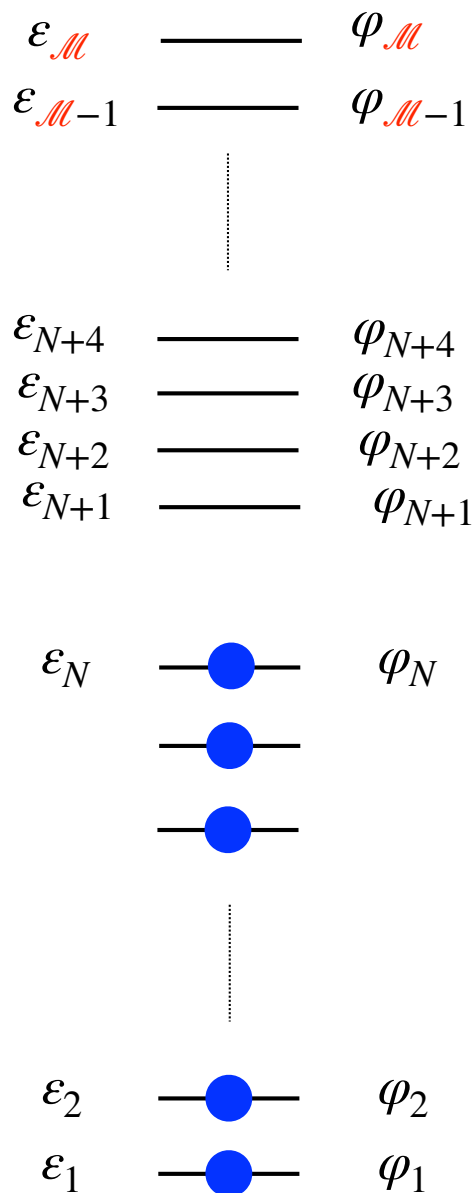
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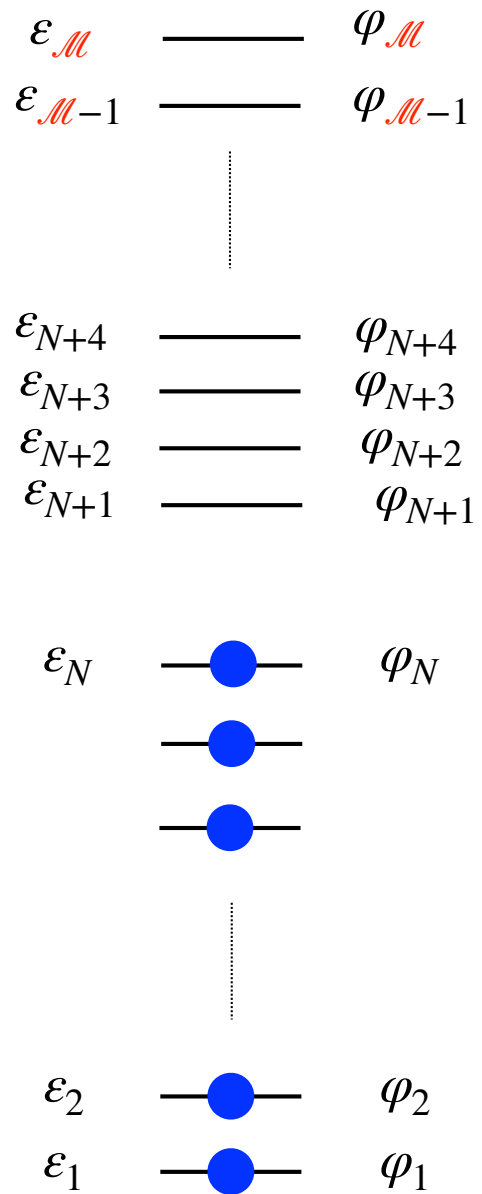


$$\hat{H} = \sum_P \varepsilon_P \hat{a}_P^\dagger \hat{a}_P$$

Spin-orbital **occupation** operator

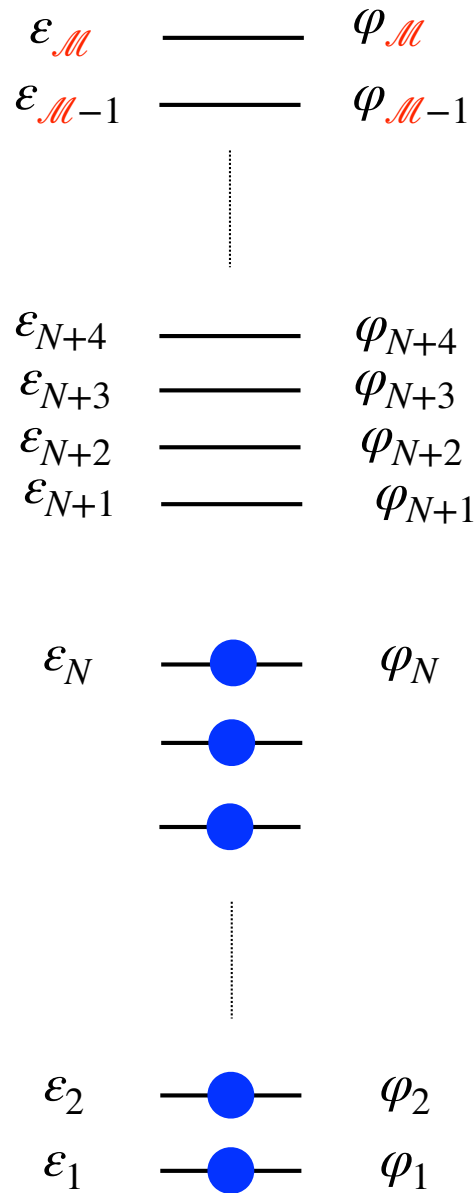
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# Non-interacting (delocalised) molecular orbital representation



$$|\Psi_0\rangle = \hat{a}_1^\dagger \hat{a}_2^\dagger \dots \hat{a}_{N-1}^\dagger \hat{a}_N^\dagger |\text{vac}\rangle$$

# 1RDM in the molecular orbital (mo) representation



$$\gamma_{PQ}^{mo} = \langle \Psi_0 | \hat{a}_P^\dagger \hat{a}_Q | \Psi_0 \rangle$$

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*Is  $\varphi_Q$  occupied in  $\Psi_0$ ?*


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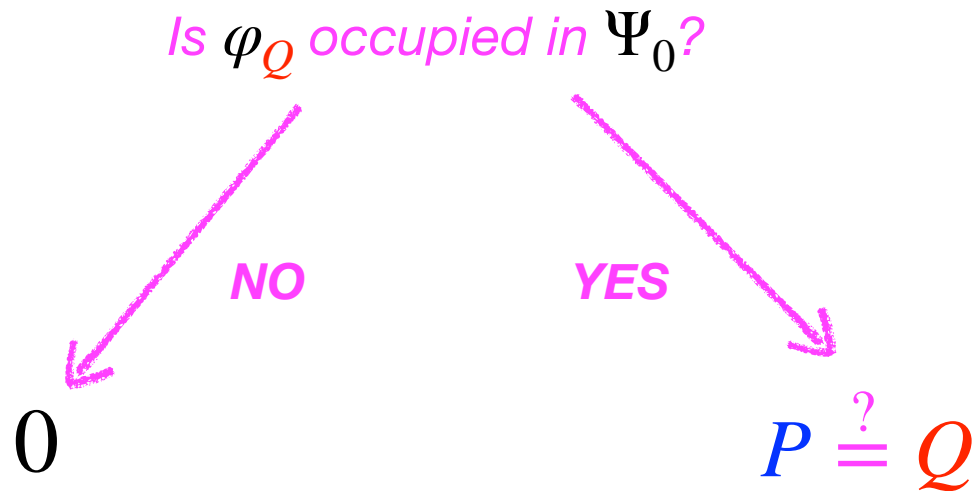
NO

0




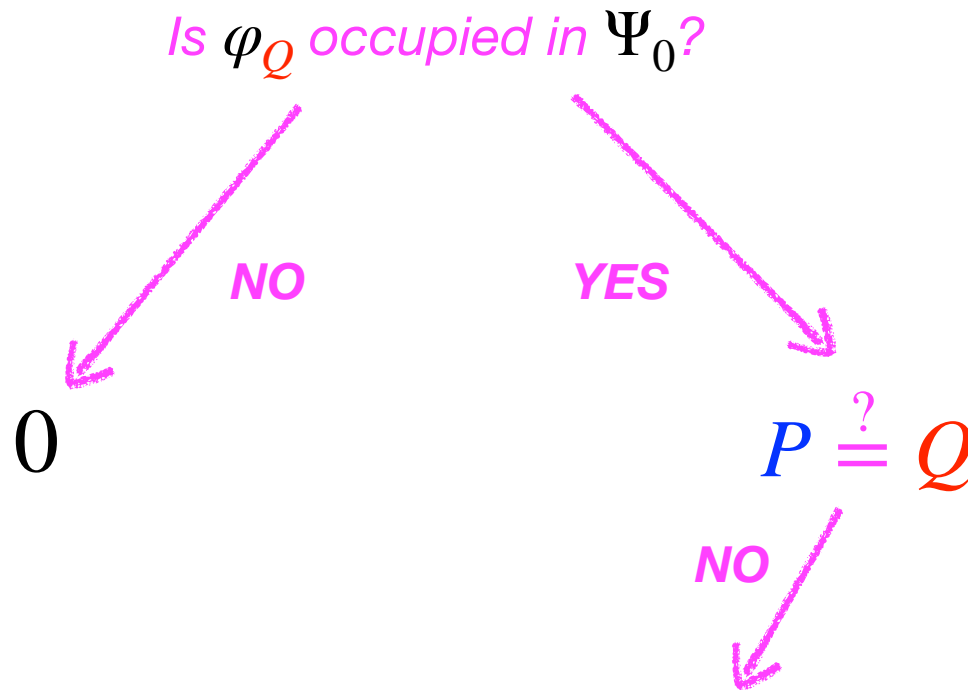
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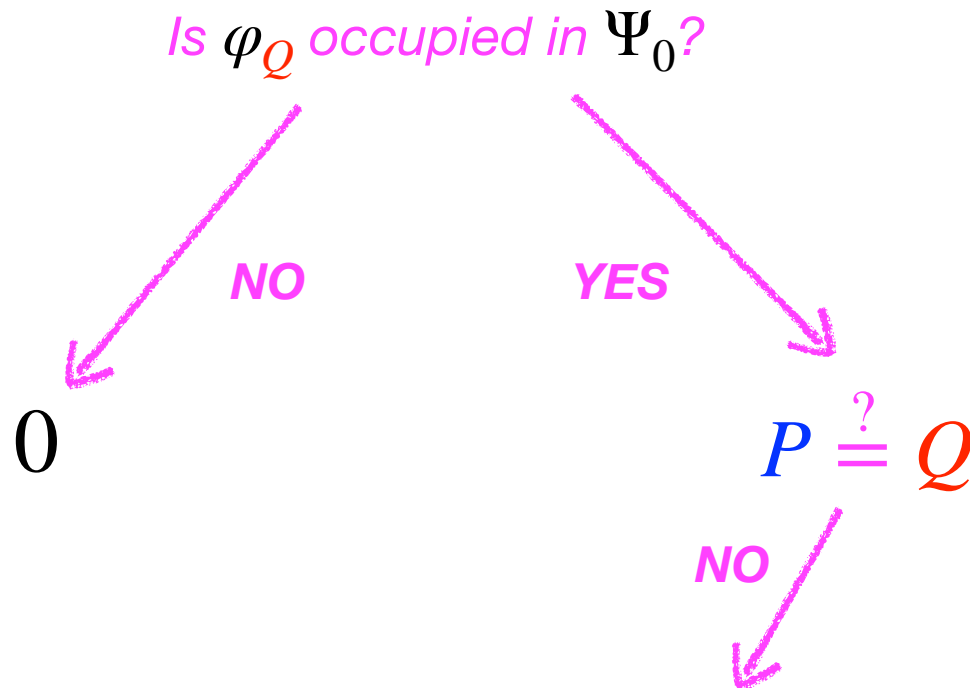


If  $\varphi_P$  is occupied in  $\Psi_0$  then  $\hat{a}_P^\dagger \hat{a}_Q | \Psi_0 \rangle = 0$  **Pauli principle**




## 1RDM in the molecular orbital (mo) representation

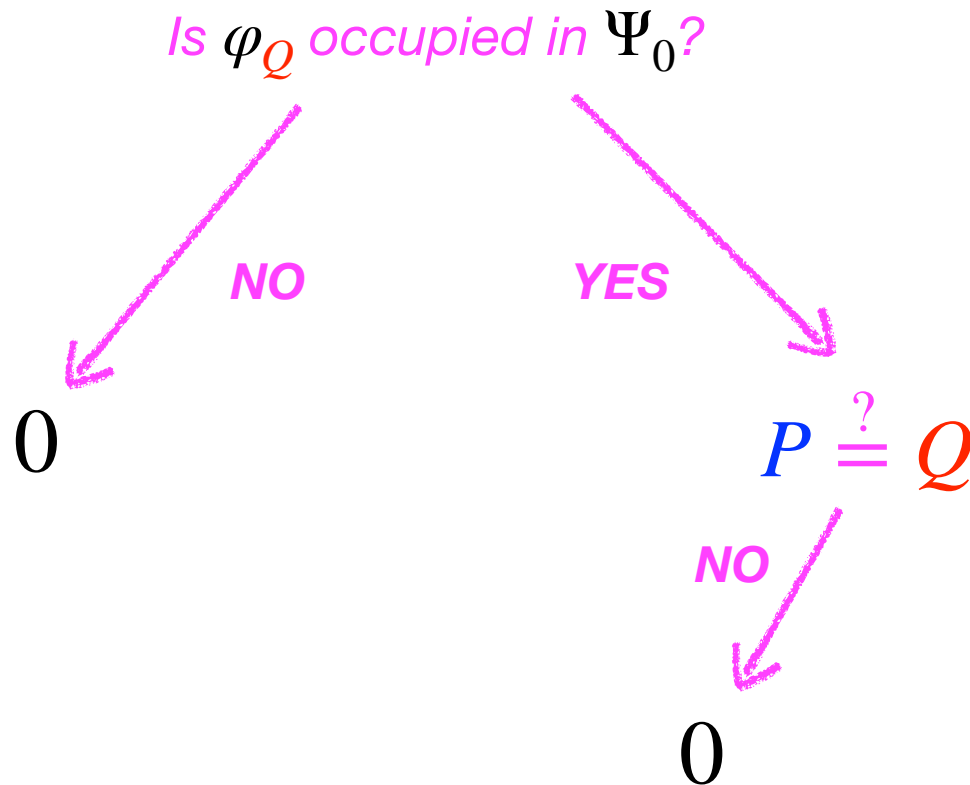
$$\gamma_{PQ}^{mo} = \langle \Psi_0 | \hat{a}_P^\dagger \hat{a}_Q | \Psi_0 \rangle \quad \text{where} \quad |\Psi_0\rangle = \hat{a}_1^\dagger \hat{a}_2^\dagger \dots \hat{a}_{N-1}^\dagger \hat{a}_N^\dagger | \text{vac} \rangle$$





If  $\varphi_P$  is unoccupied in  $\Psi_0$  then  $\hat{a}_P^\dagger \hat{a}_Q | \Psi_0 \rangle \perp | \Psi_0 \rangle \Rightarrow \langle \Psi_0 | \hat{a}_P^\dagger \hat{a}_Q | \Psi_0 \rangle = 0$

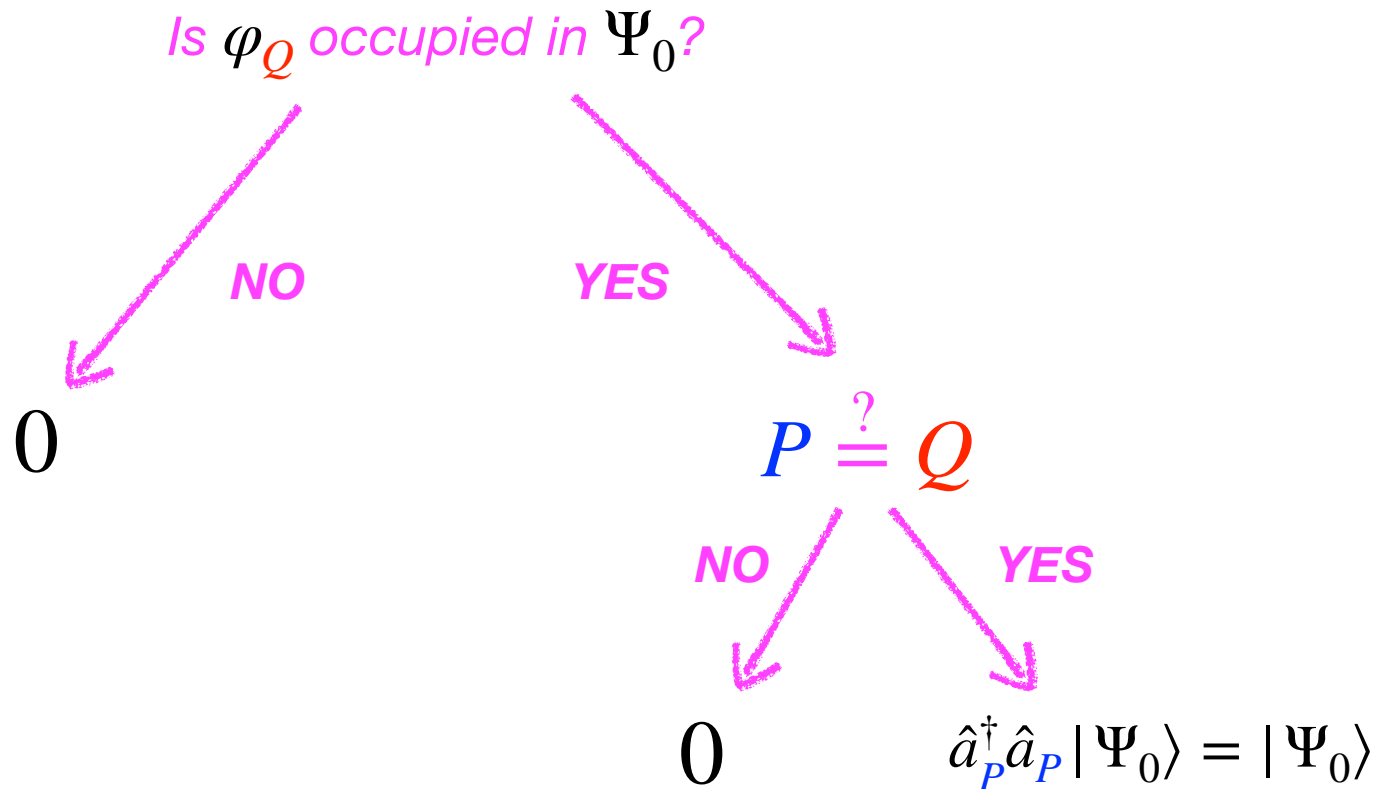
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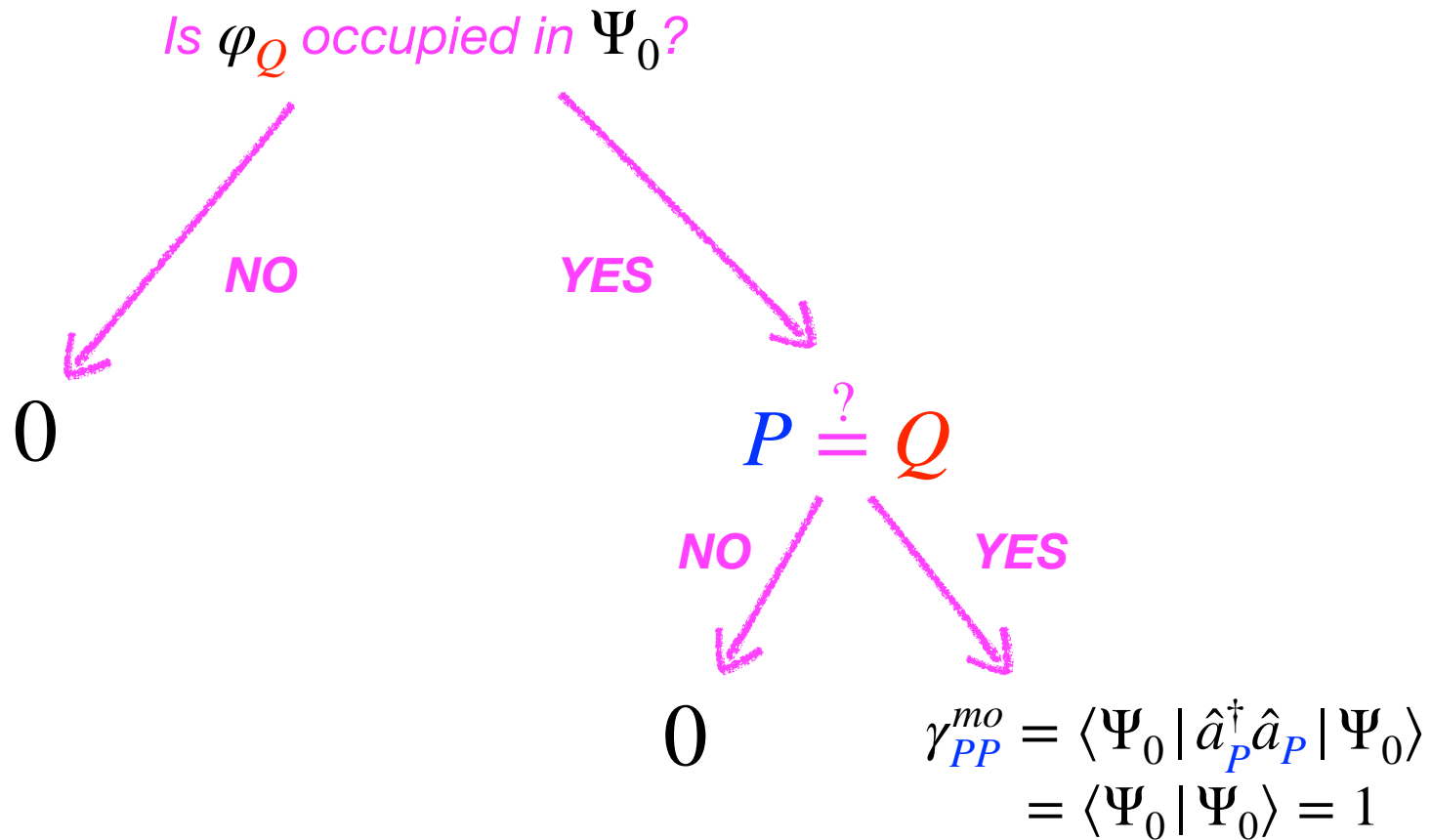
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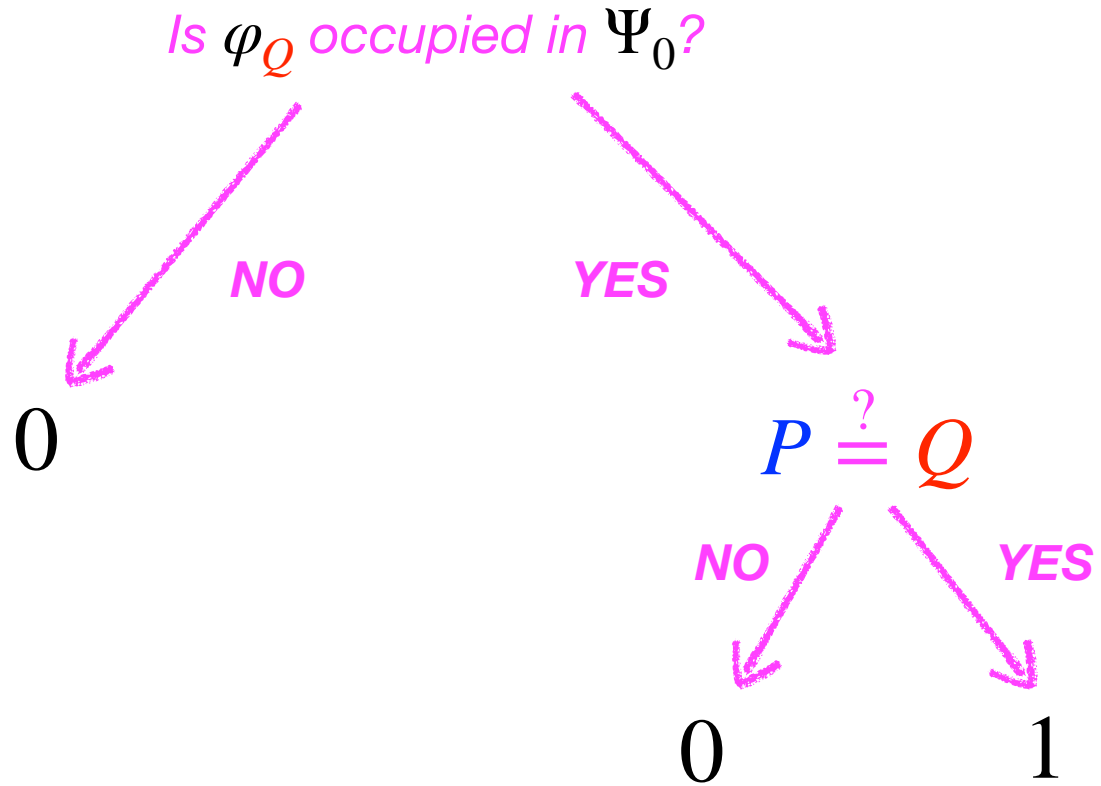
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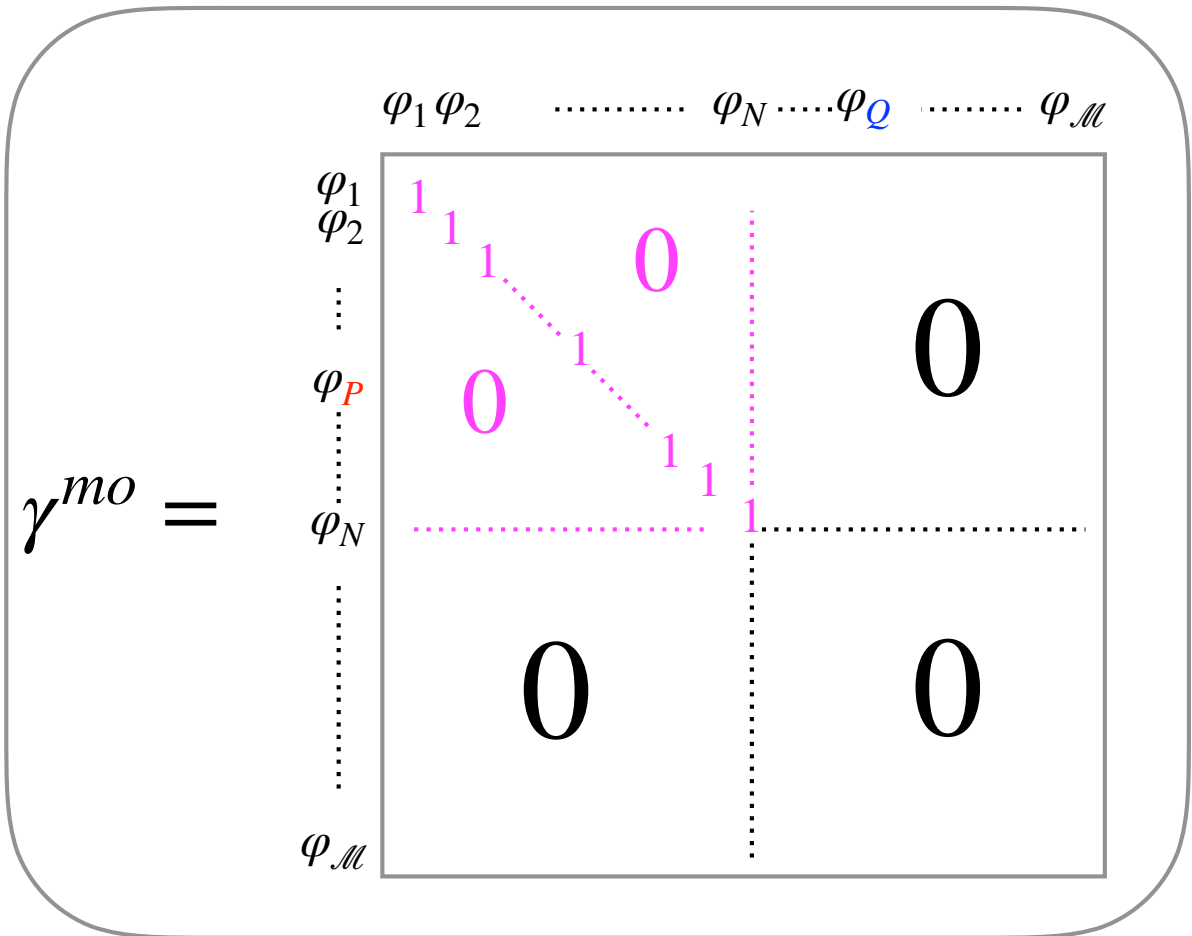
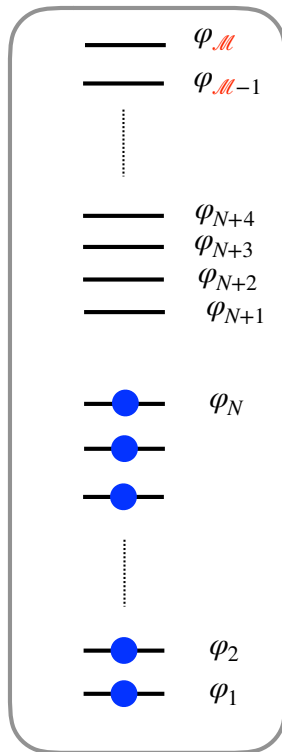
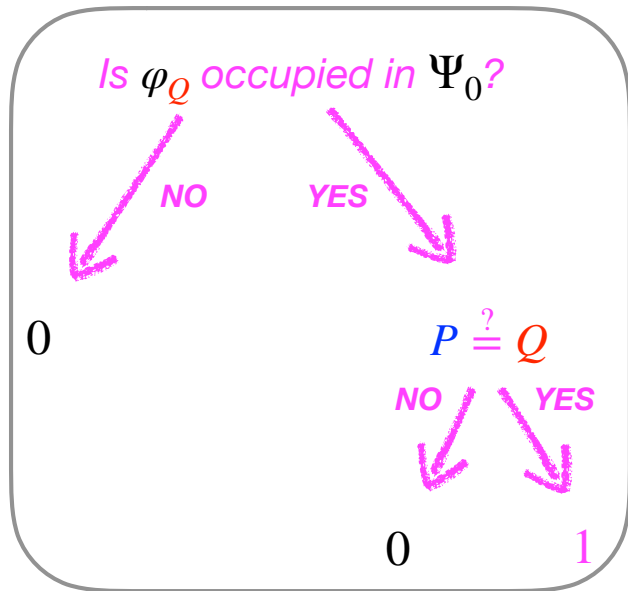
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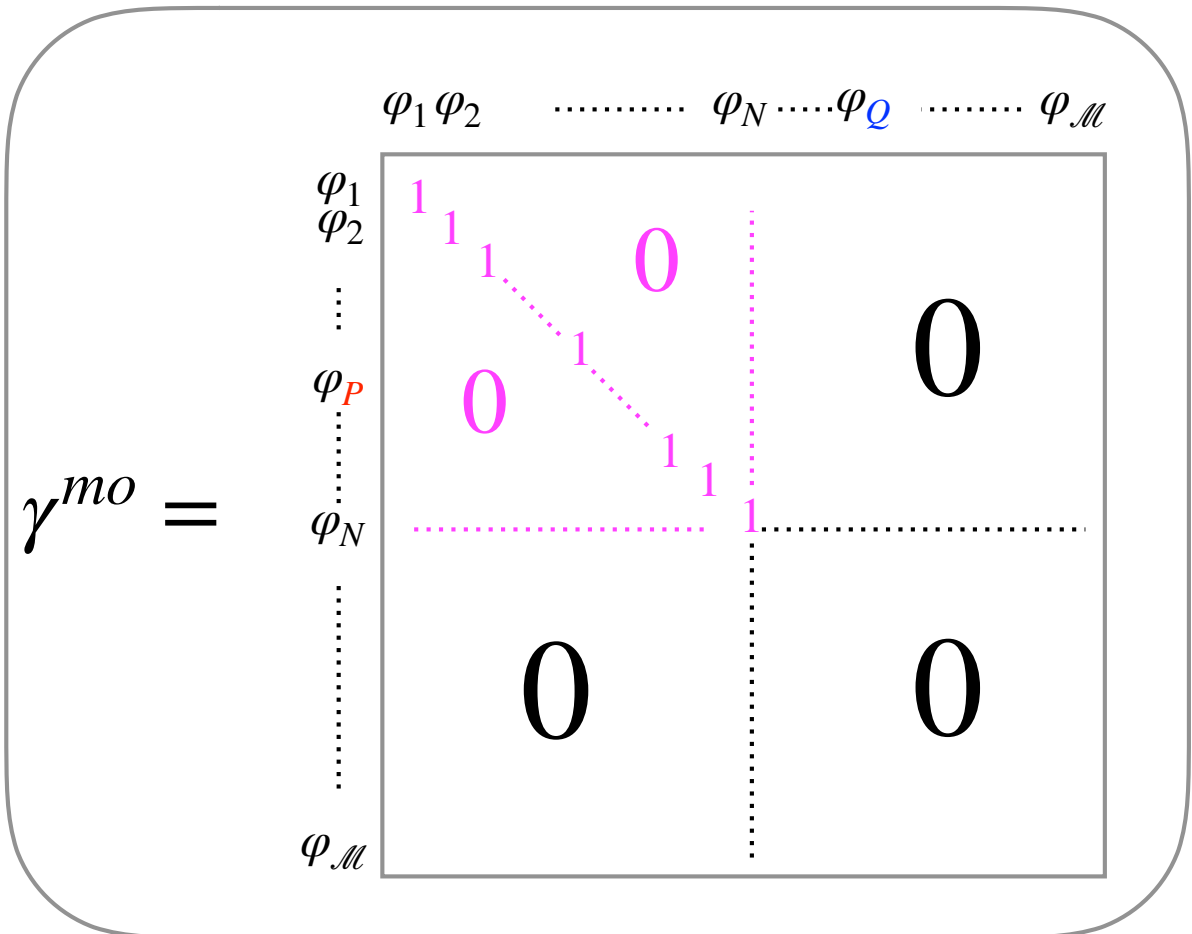
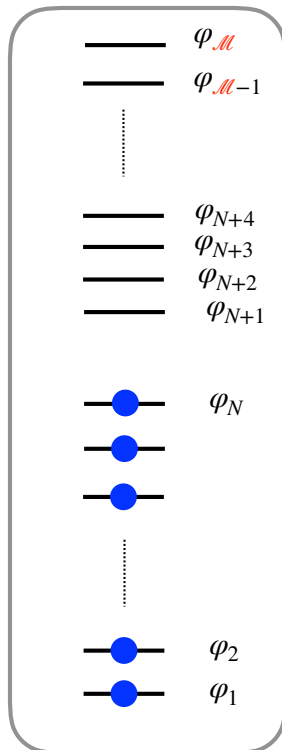
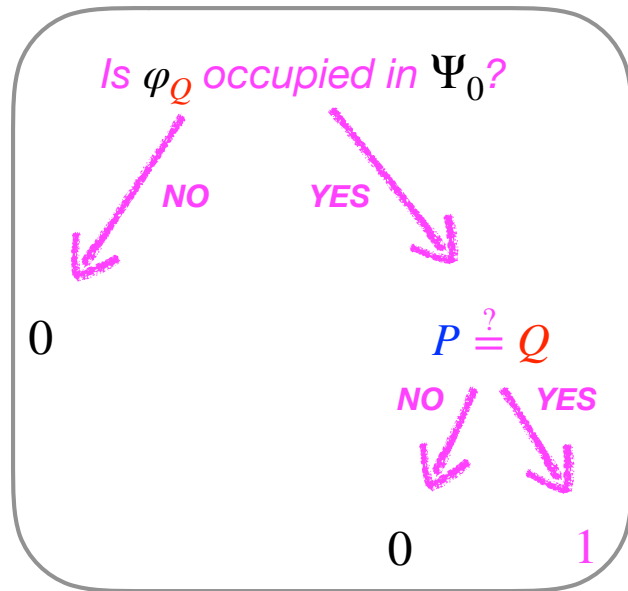
# 1RDM in the molecular orbital (mo) representation



Non-interacting problem solved!



# 1RDM in the molecular orbital (mo) representation



*No entanglement between the molecular orbitals  
in the non-interacting case*

# Idempotency property

$$\gamma^{mo} = \begin{array}{c} \varphi_1 \varphi_2 \quad \dots \quad \varphi_N \dots \varphi_Q \quad \dots \quad \varphi_M \\ \begin{array}{|c|} \hline \begin{array}{c} \varphi_1 \\ \varphi_2 \\ \vdots \\ \varphi_P \\ \varphi_N \\ \vdots \\ \varphi_M \end{array} \\ \hline \end{array} \end{array} = \begin{array}{c} \begin{array}{|c|} \hline \begin{array}{c} 1 \\ 1 \\ \vdots \\ 1 \\ 1 \end{array} \\ \hline \end{array} \begin{array}{|c|} \hline \begin{array}{c} 0 \\ 0 \\ \vdots \\ 0 \\ 0 \end{array} \\ \hline \end{array} \\ \hline \\ \hline \end{array} = [\gamma^{mo}]^2$$

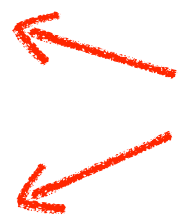
The diagram illustrates the idempotency property of the matrix  $\gamma^{mo}$ . The matrix is partitioned into four quadrants by a vertical dashed line at column  $\varphi_N$  and a horizontal dashed line at row  $\varphi_N$ . The top-left quadrant is an upper triangular matrix with 1s on the diagonal and 0s elsewhere. The top-right, bottom-left, and bottom-right quadrants are all zero matrices. The equation shows that  $\gamma^{mo}$  is equal to its square,  $[\gamma^{mo}]^2$ .



***Turning to the localised picture (useless here although interesting)***

$$|\varphi_P\rangle = \sum_I C_{IP} |\chi_I\rangle$$
$$|\varphi_Q\rangle = \sum_J C_{JQ} |\chi_J\rangle$$

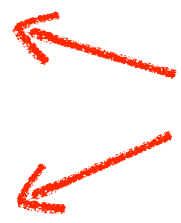
*Localised spin-orbitals*



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$$\langle\varphi_P|\varphi_Q\rangle = \delta_{PQ} = \sum_{IJ} C_{IP}C_{JQ} \langle\chi_I|\chi_J\rangle$$

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*Localised spin-orbitals*

$$\langle \varphi_P | \varphi_Q \rangle = \delta_{PQ} = \sum_{IJ} C_{IP} C_{JQ} \langle \chi_I | \chi_J \rangle$$

*Orthonormalisation procedure*

$\delta_{IJ}$

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*Localised spin-orbitals*

$$\langle\varphi_P|\varphi_Q\rangle = \delta_{PQ} = \sum_{IJ} C_{IP} C_{JQ} \delta_{IJ}$$

*Molecular orbital coefficients matrix*

$$= \sum_I C_{IP} C_{IQ} = \sum_I [\mathbf{C}^T]_{PI} [\mathbf{C}]_{IQ}$$
$$= [\mathbf{C}^T \mathbf{C}]_{PQ}$$

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*Localised spin-orbitals*

*Unitary transformation  
from the delocalised to localised pictures*

$$\mathbf{C}^{-1} = \mathbf{C}^T$$

$$\delta_{PQ} = [\mathbf{C}^T \mathbf{C}]_{PQ}$$

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*1RDM in the delocalised representation*

$$\gamma_{PQ}^{mo} = \sum_{IJ} C_{IP} C_{JQ} \langle \Psi_0 | \hat{c}_I^\dagger \hat{c}_J | \Psi_0 \rangle$$

$$\gamma_{IJ}^{loc} \equiv \gamma_{IJ}$$

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$$\hat{a}_P^\dagger = \sum_I C_{IP} \hat{c}_I^\dagger$$
$$\hat{a}_Q = \sum_J C_{JQ} \hat{c}_J$$

$$\gamma_{PQ}^{mo} = \sum_{IJ} C_{IP} C_{JQ} \gamma_{IJ} = [\mathbf{C}^T]_{PI} \gamma_{IJ}^{loc} [\mathbf{C}]_{JQ}$$



**Turning to the localised picture (useless here although interesting)**

$$\gamma_{PQ}^{mo} = \langle \Psi_0 | \hat{a}_P^\dagger \hat{a}_Q | \Psi_0 \rangle$$



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$$\begin{aligned} \gamma_{PQ}^{mo} &= \sum_{IJ} C_{IP} C_{JQ} \gamma_{IJ} = [\mathbf{C}^T]_{PI} \gamma_{IJ}^{loc} [\mathbf{C}]_{JQ} \\ &= [\mathbf{C}^T \gamma^{loc} \mathbf{C}]_{PQ} \end{aligned}$$

*Turning to the localised picture (useless here although interesting)*

$$\gamma^{loc} = \mathbf{C}\gamma^{mo}\mathbf{C}^T$$

$$\gamma^{mo} = \mathbf{C}^T\gamma^{loc}\mathbf{C}$$

$$\gamma_{PQ}^{mo} = [\mathbf{C}^T\gamma^{loc}\mathbf{C}]_{PQ}$$

**Turning to the localised picture (useless here although interesting)**

$$\gamma^{loc} = \mathbf{C}\gamma^{mo}\mathbf{C}^T$$

$$\gamma_{IJ}^{loc} = \sum_{PQ} C_{IP} \gamma_{PQ}^{mo} C_{JQ} = \sum_{\substack{\text{occupied} \\ \text{spin-MOs}}} C_{IP} C_{JP}$$

**Turning to the localised picture (useless here although interesting)**

$$\gamma^{loc} = \mathbf{C}\gamma^{mo}\mathbf{C}^T \quad \leftarrow \text{Not diagonal!}$$

$$\gamma_{IJ}^{loc} = \langle \hat{c}_I^\dagger \hat{c}_J \rangle_{\Psi_0} = \sum_P^{\text{occupied spin-MOs}} C_{IP}C_{JP} \neq \delta_{IJ}$$

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Any localised spin-orbital  $\chi_I$  is **entangled** with the other spin-orbitals  $\chi_J$

**Turning to the localised picture (useless here although interesting)**

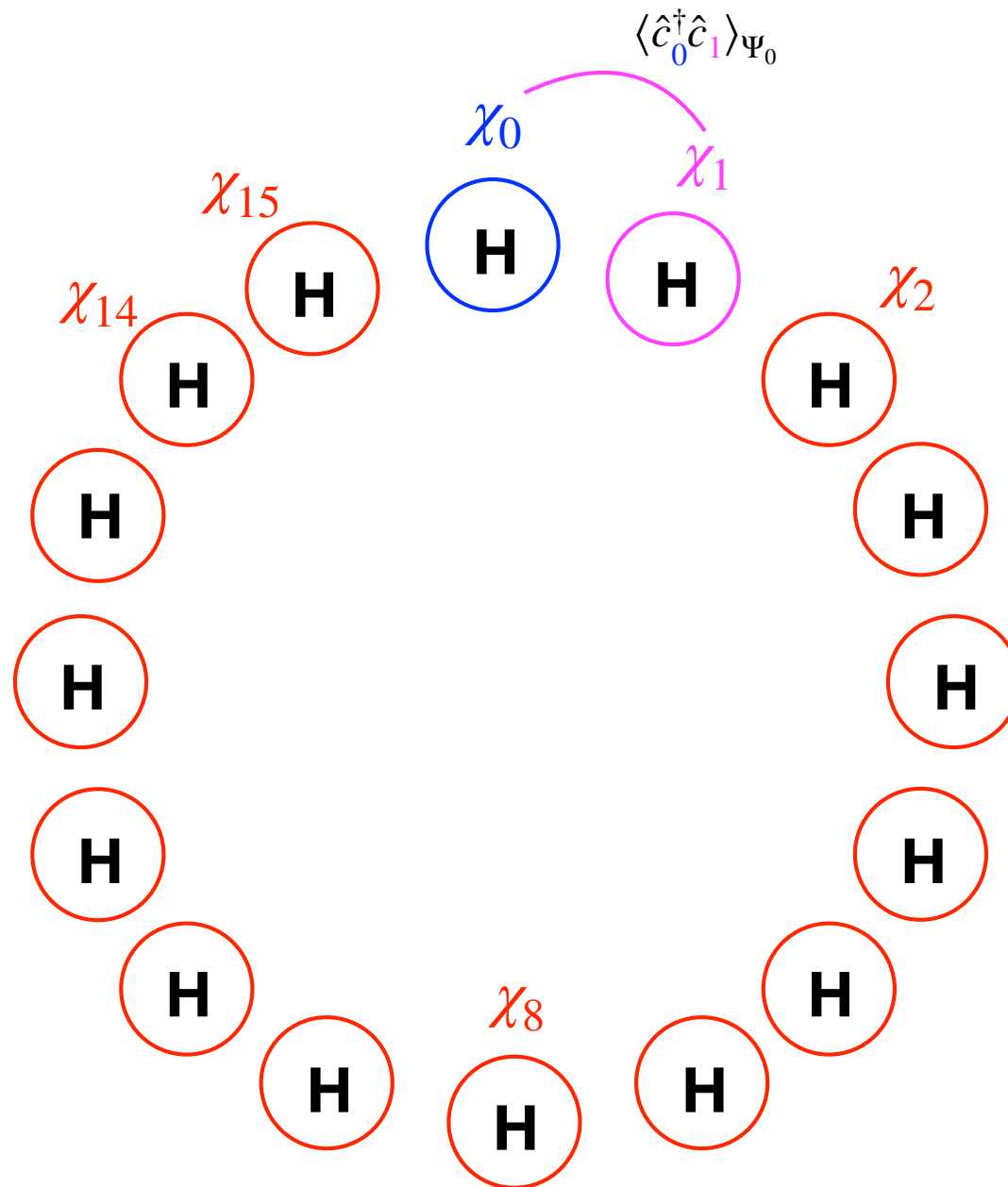
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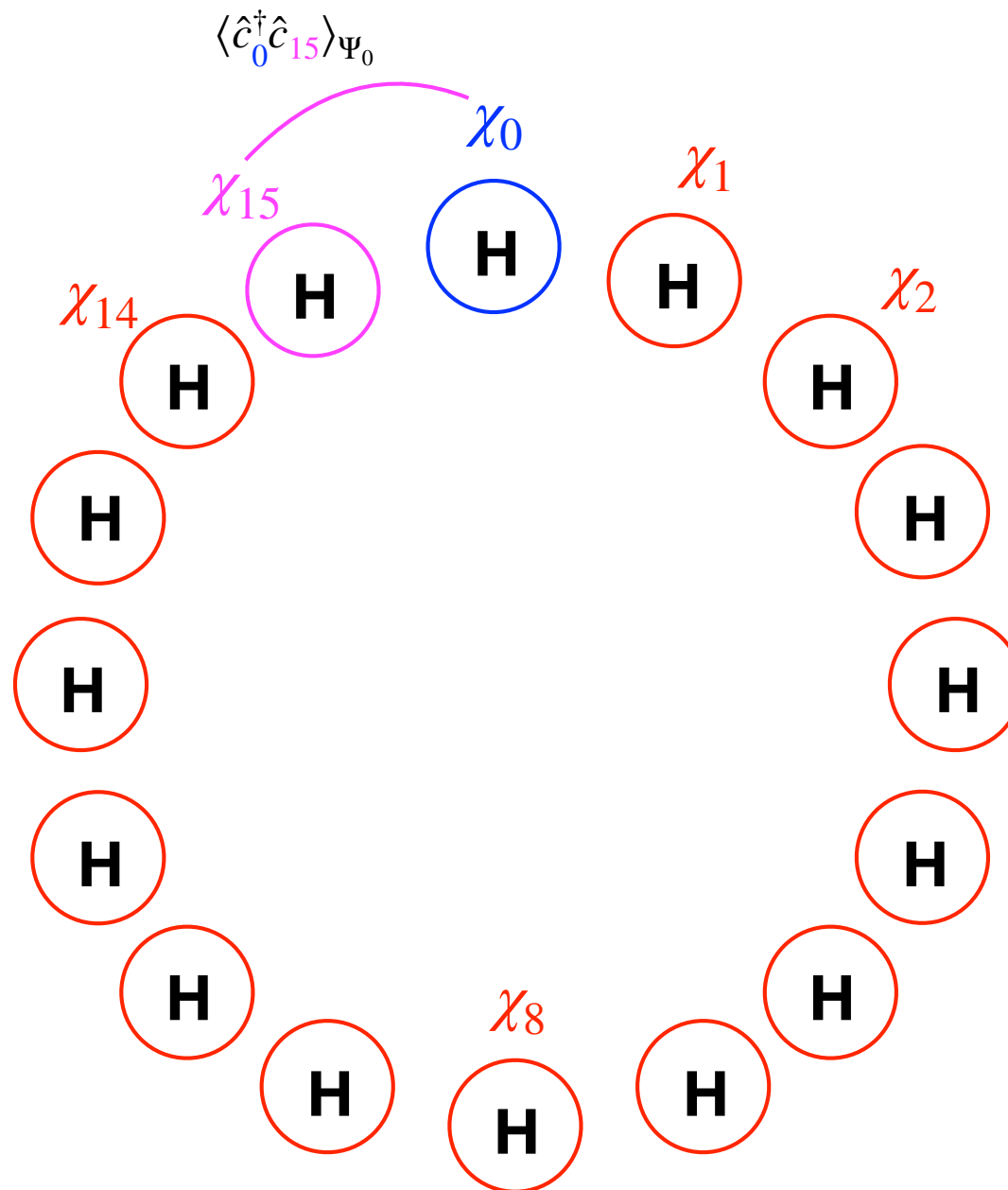
Any localised spin-orbital  $\chi_I$  is **entangled** with the other spin-orbitals  $\chi_J$

**unlike in the delocalised molecular orbital space!**

# Prototypical ring of $L = 16$ hydrogen atoms

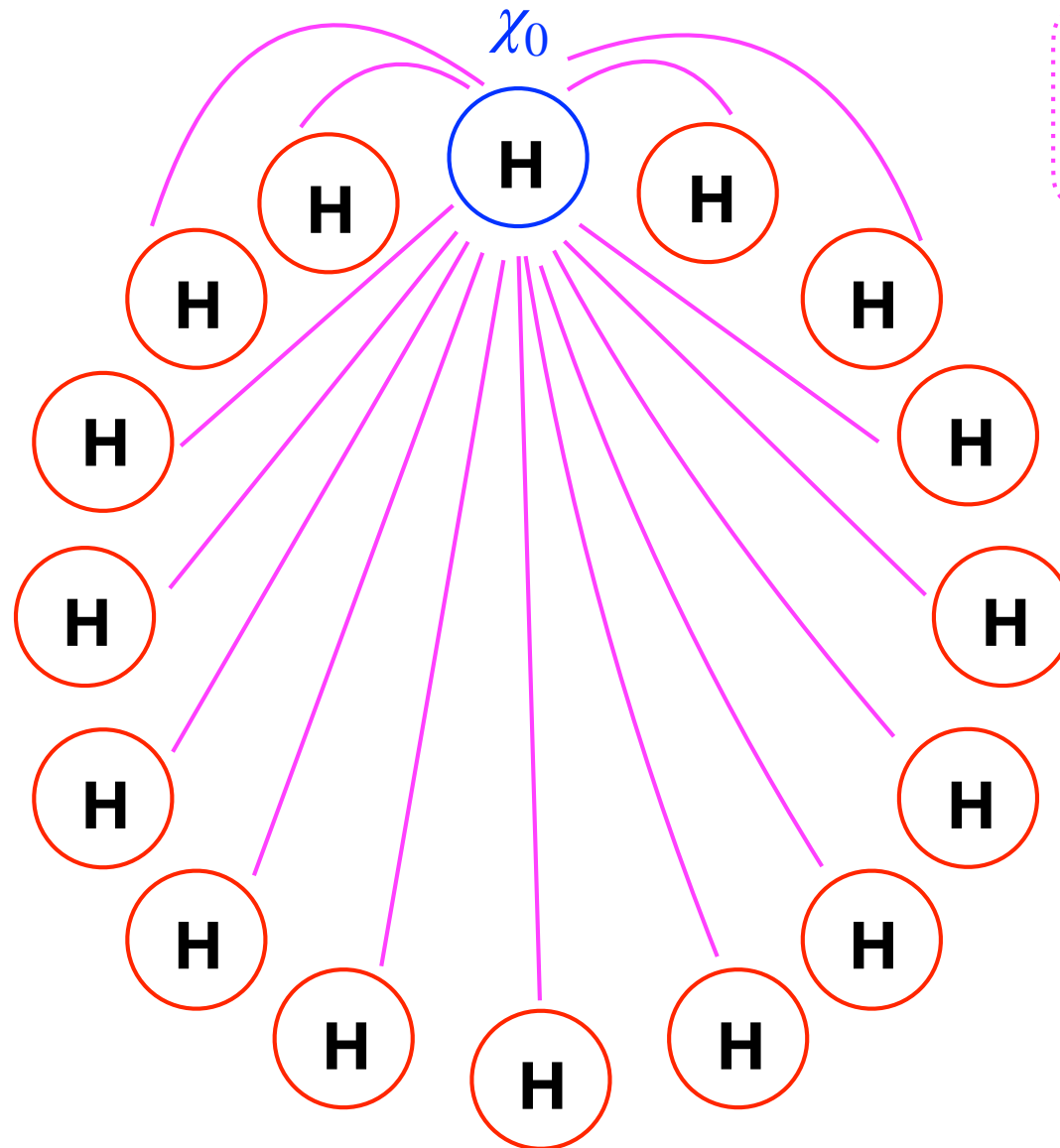


# Prototypical ring of $L = 16$ hydrogen atoms



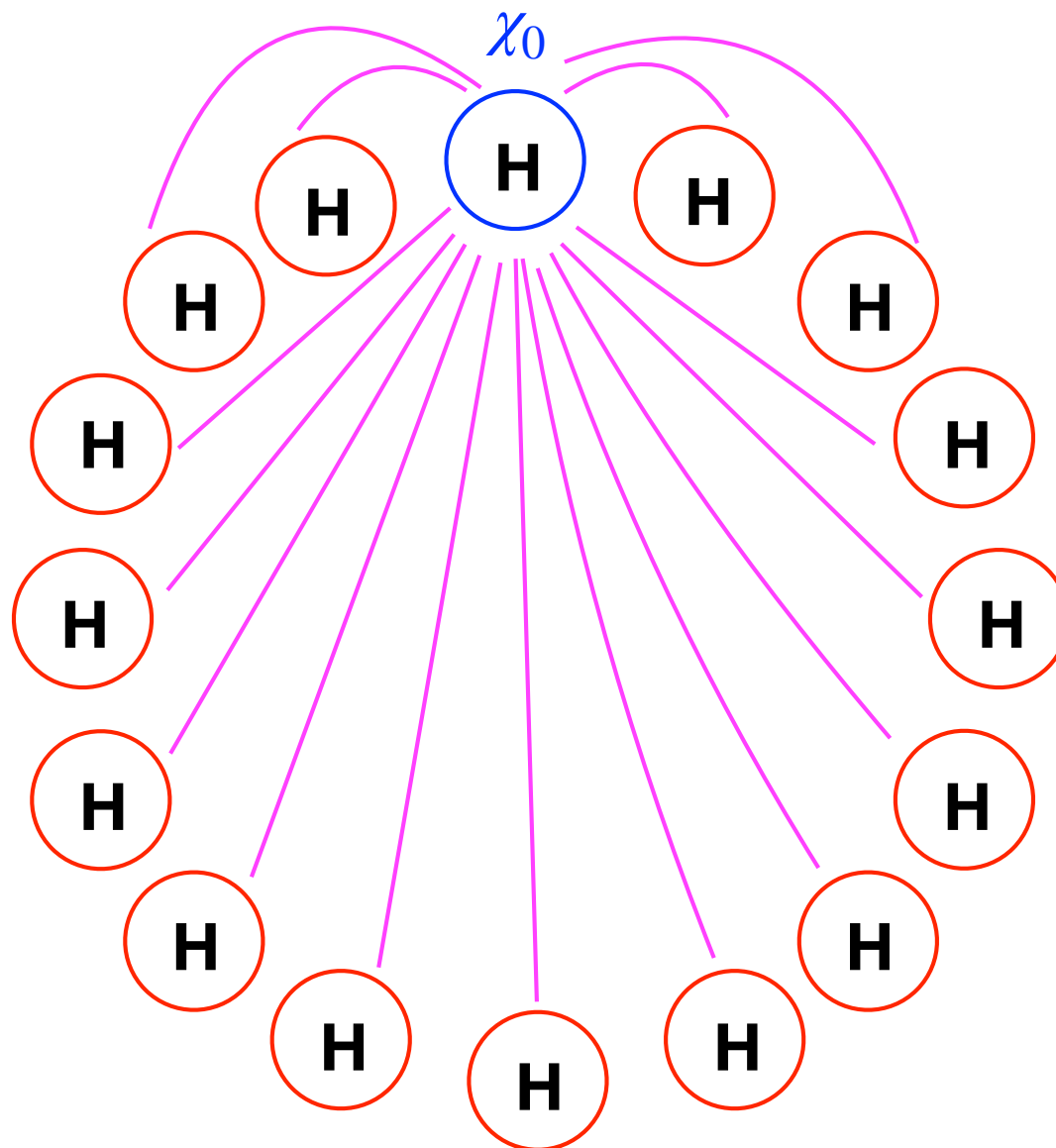


# Prototypical ring of $L = 16$ hydrogen atoms

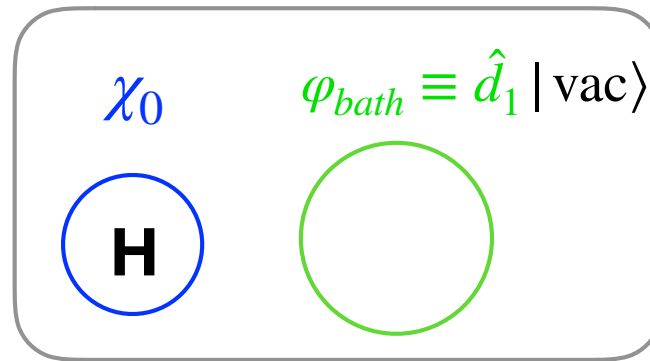


$$\left\{ \langle \hat{c}_0^\dagger \hat{c}_J \rangle_{\Psi_0} \right\}_{1 \leq J \leq 15}$$

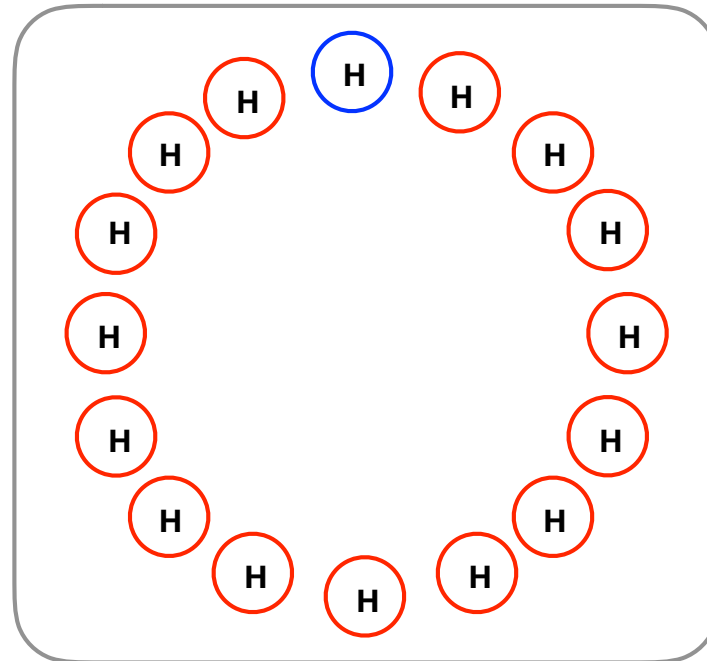
To-be-embedded (so-called *impurity*)  
localised orbital



# Exact density matrix functional embedding

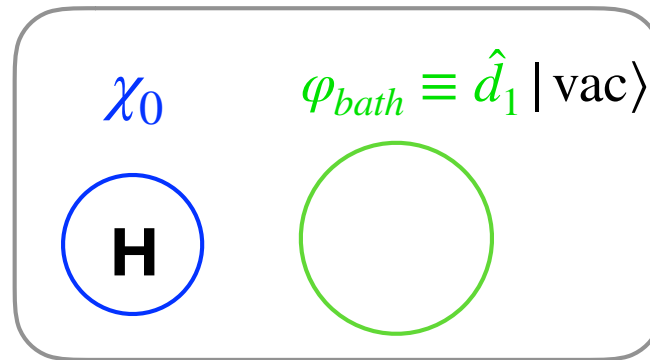


Two-electron system



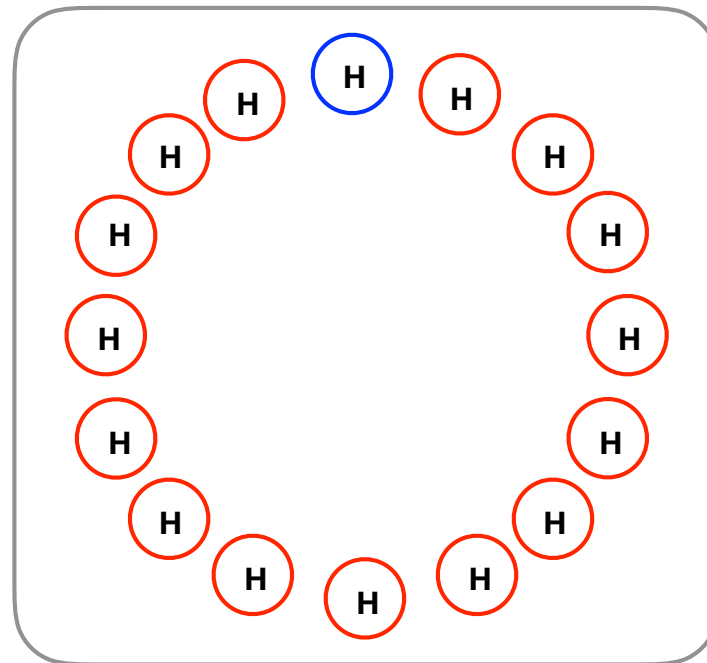
$N$ -electron system

# Exact density matrix functional embedding



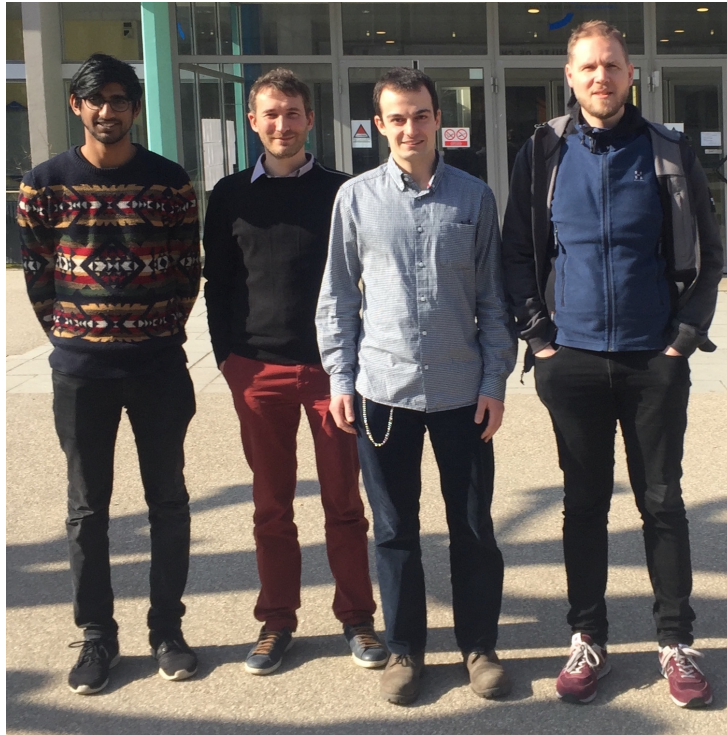
Two-electron system

Let's prove that the embedding is **exact**  
for **non-interacting** electrons

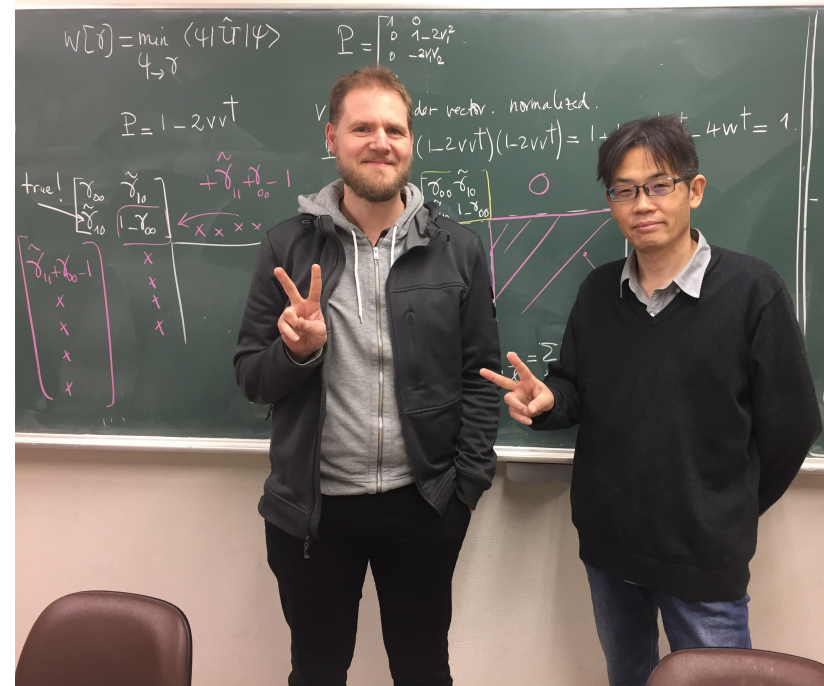


$N$ -electron system

# The “Householder embedding” project

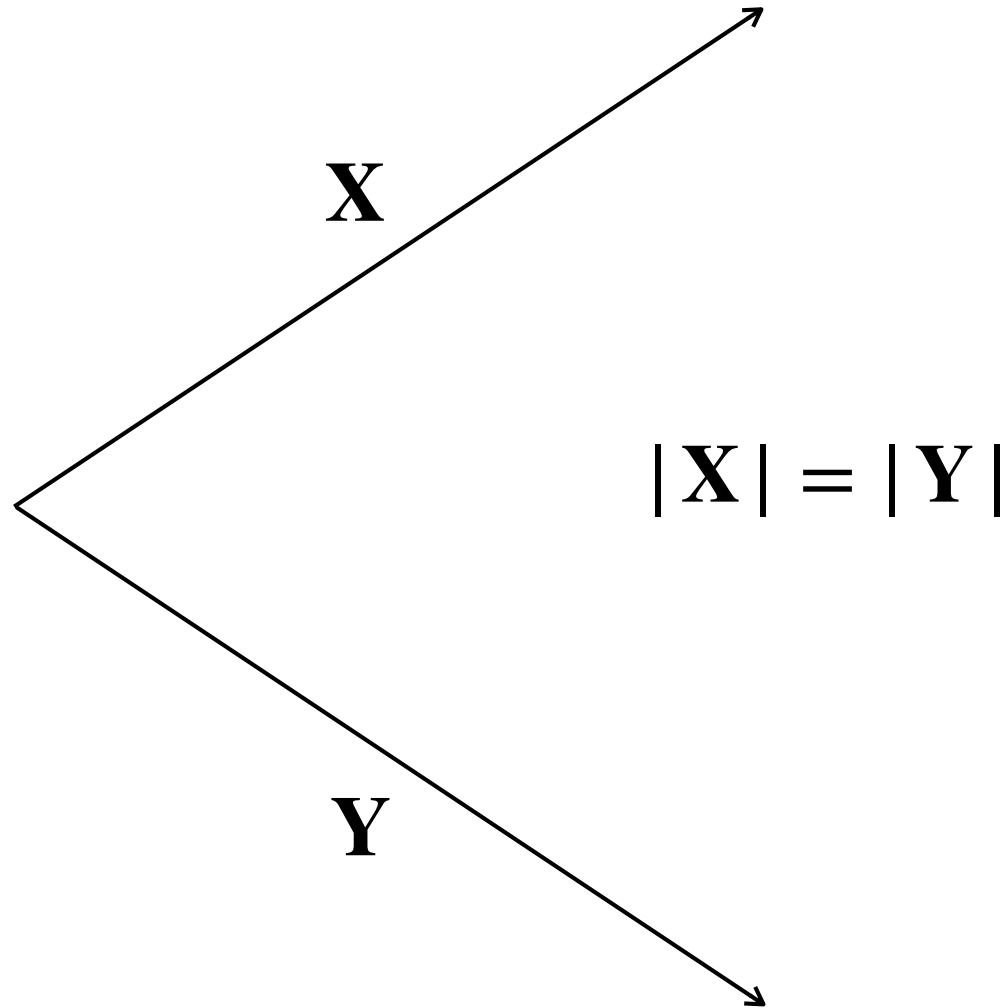


From left to right: **S. Sekaran** (Strasbourg, France),  
**M. Saubanère** (Montpellier, France),  
**L. Mazouin** (Strasbourg, France), and **E.F.**

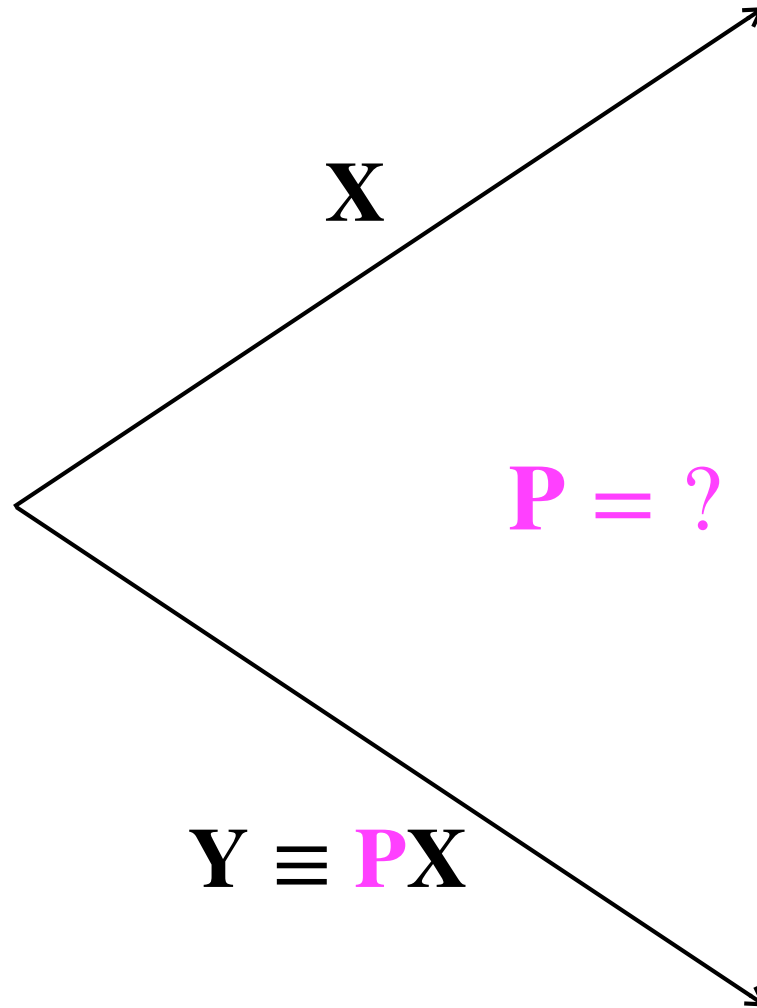


**E.F.** and **M. Tsuchiizu** (Nara, Japan).

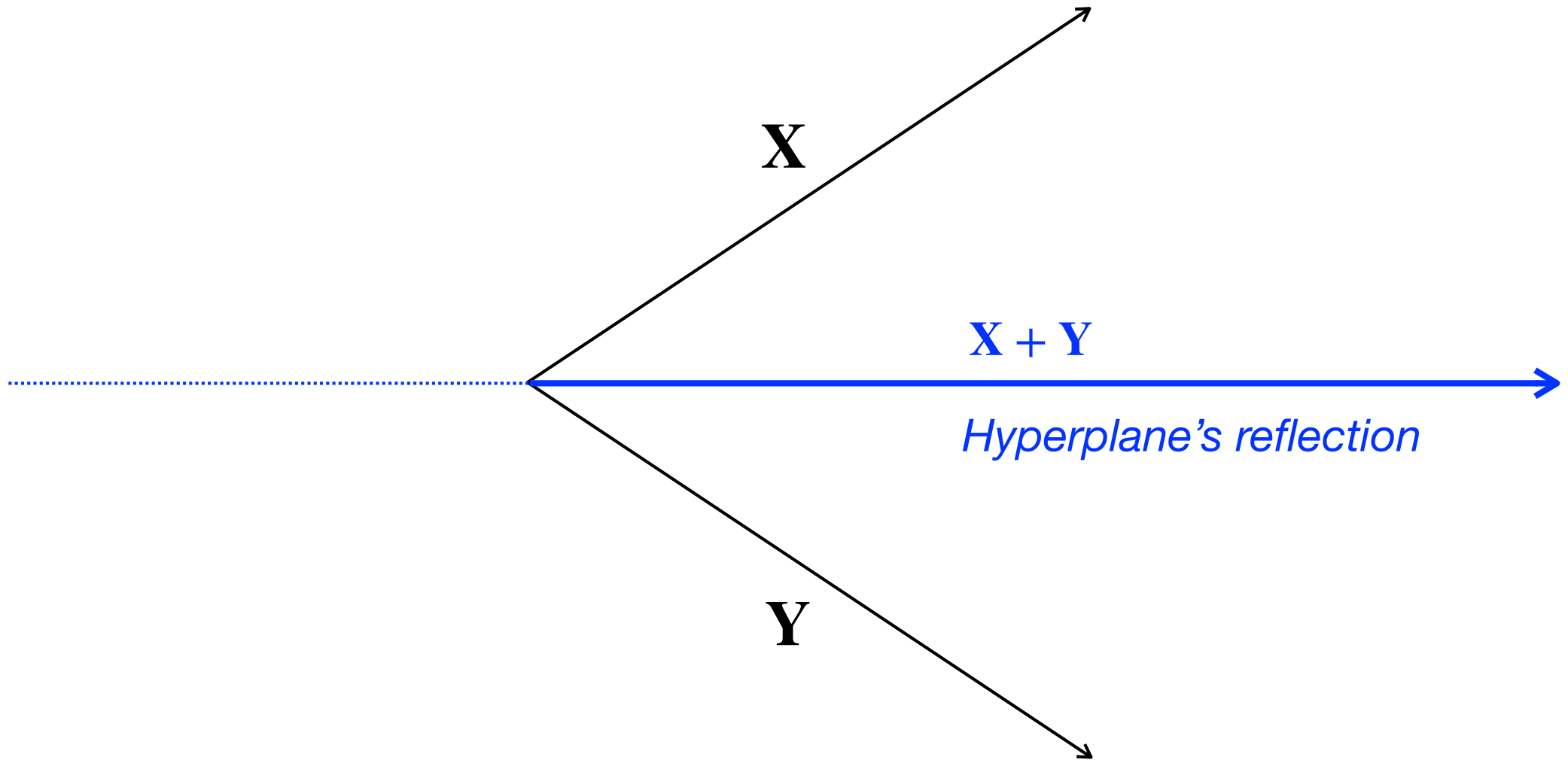
# The Householder transformation



# The Householder transformation

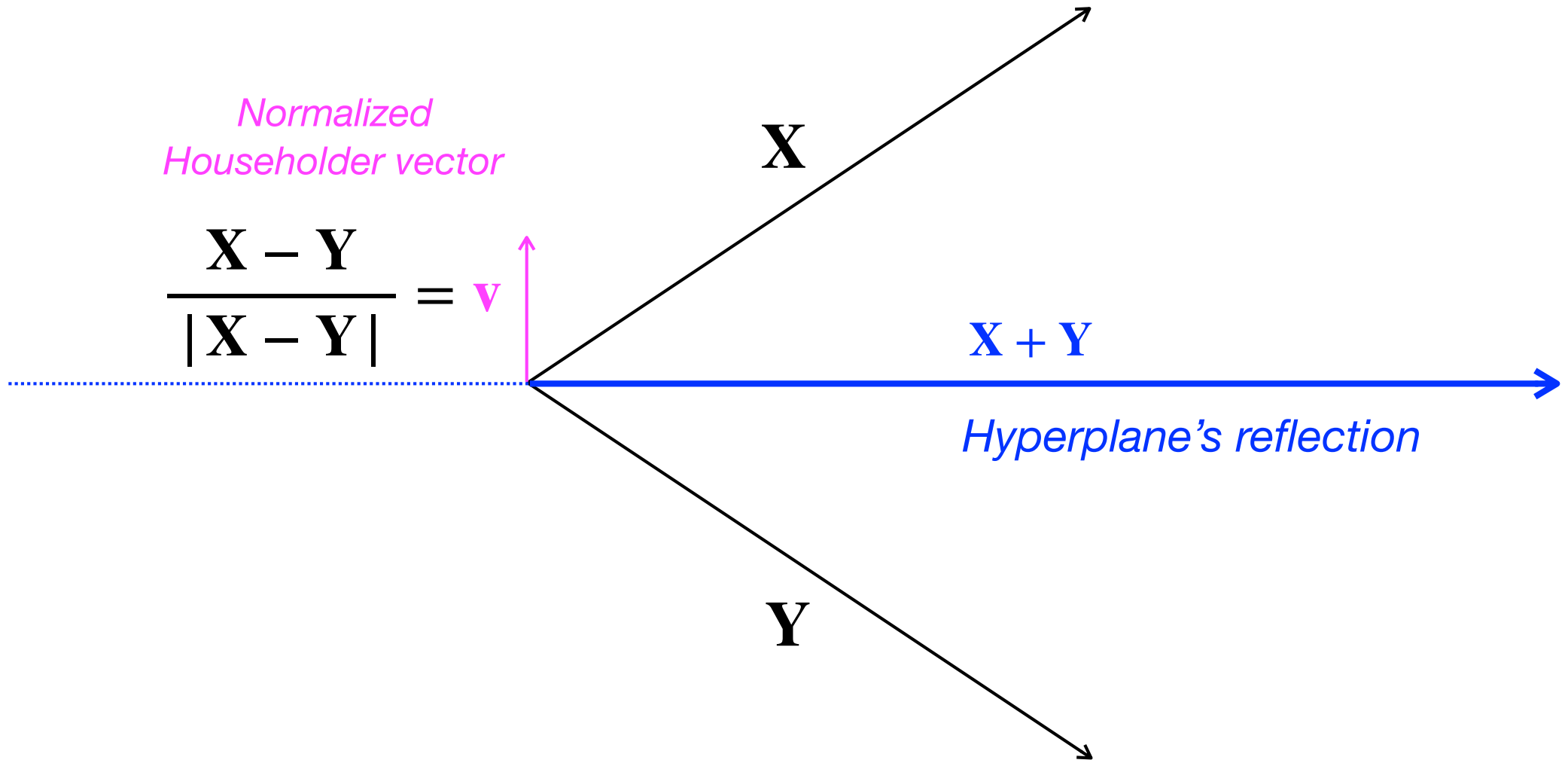


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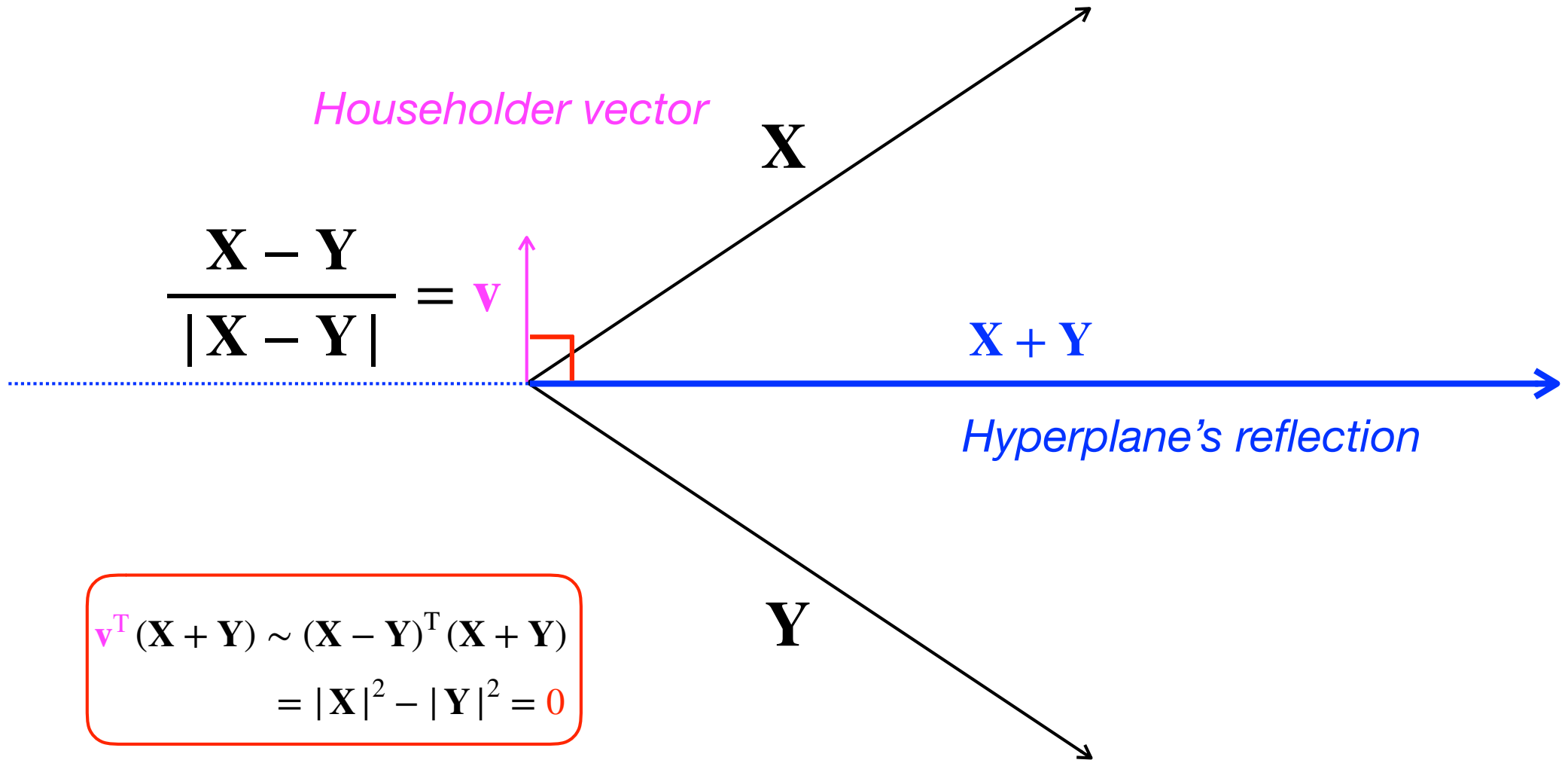




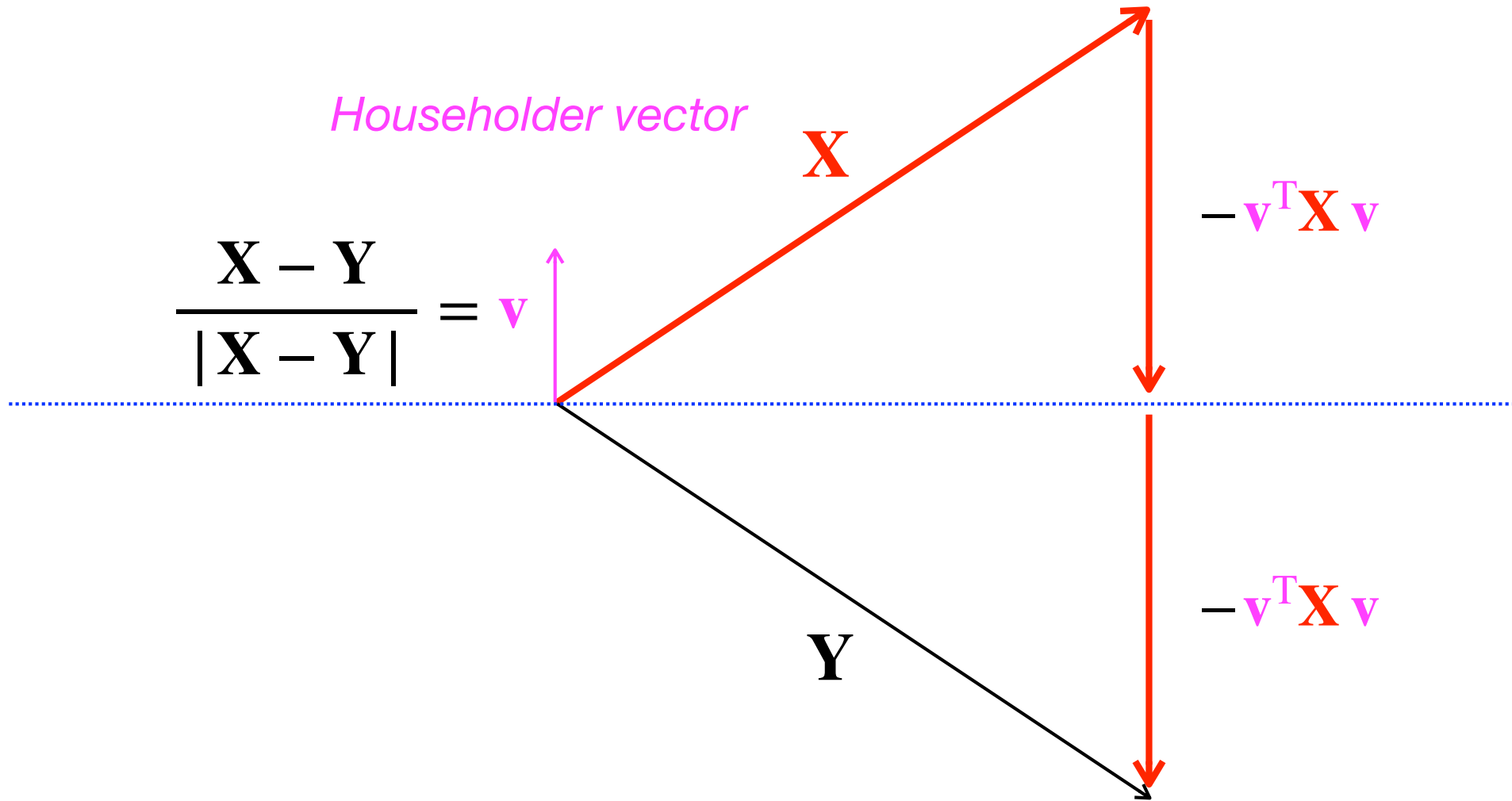
# The Householder transformation



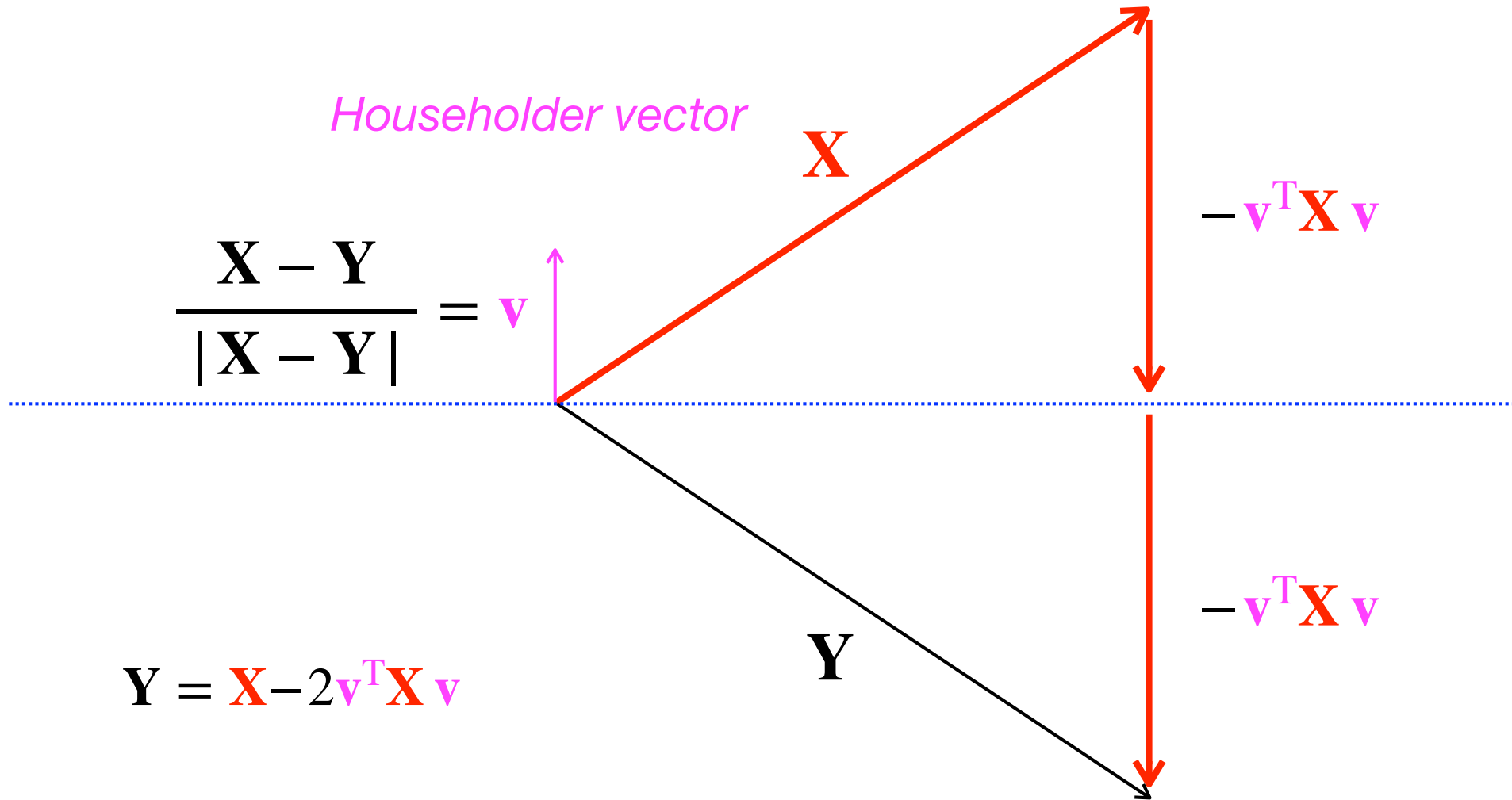
# The Householder transformation



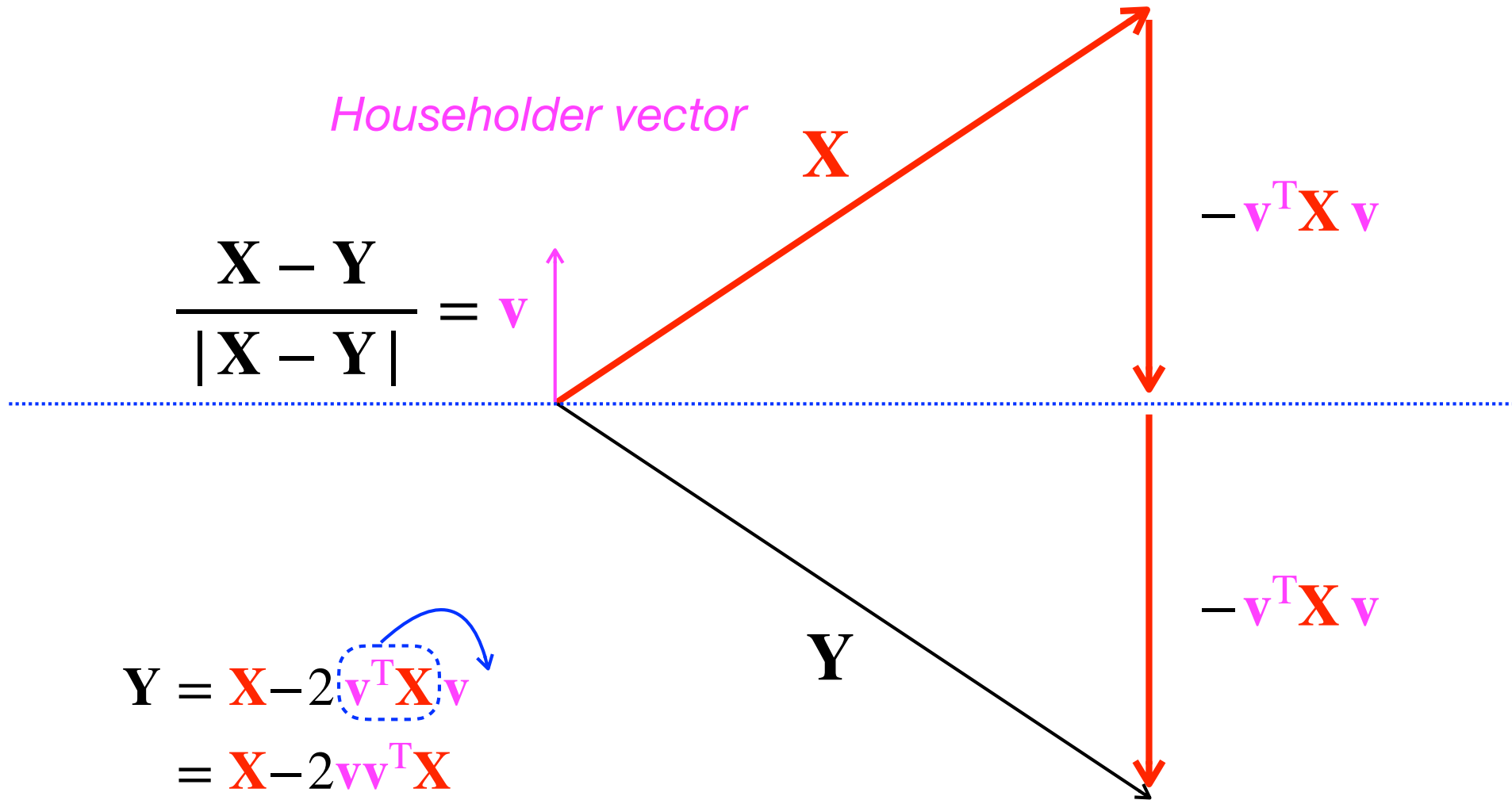
# The Householder transformation



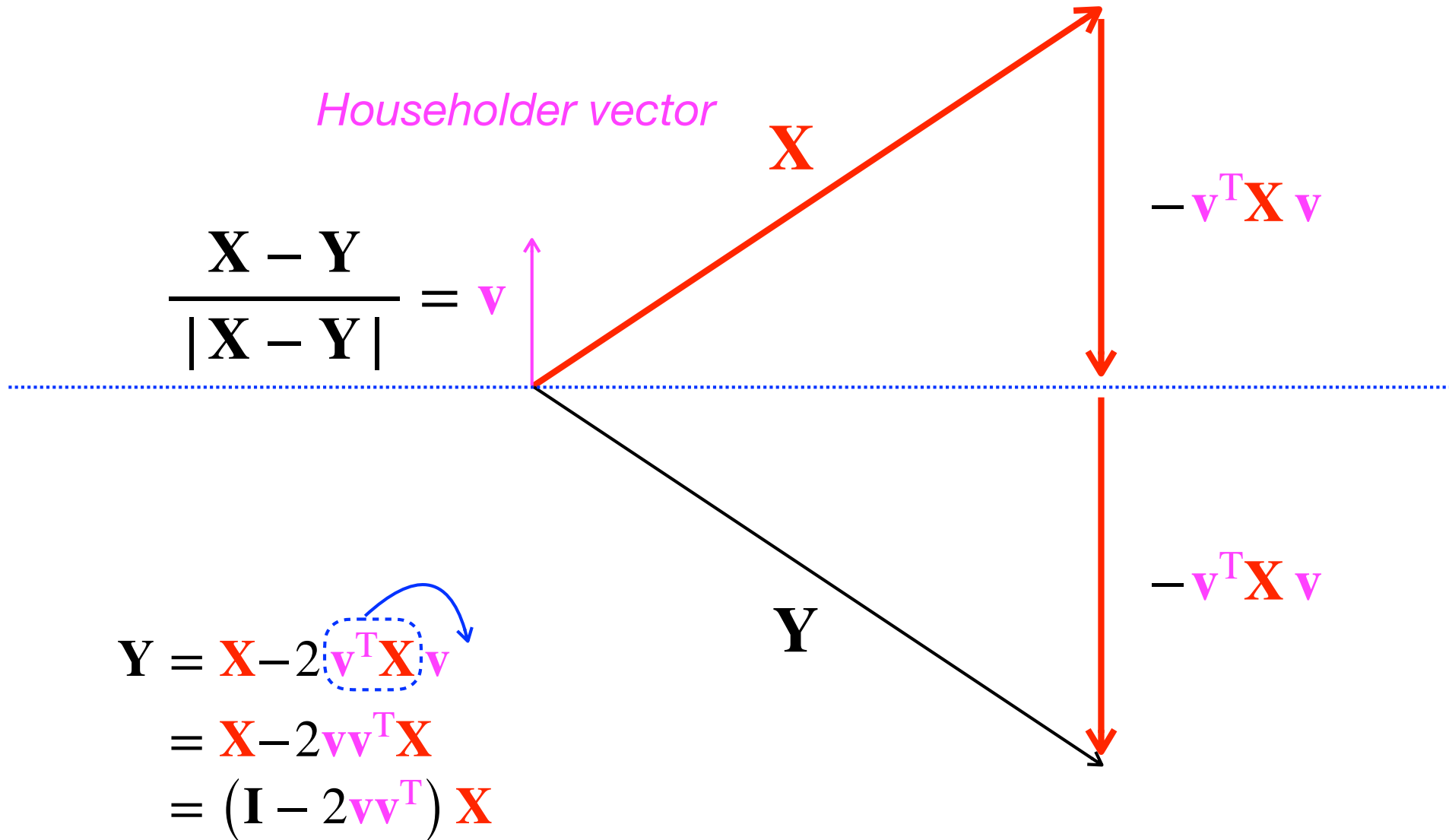
# The Householder transformation



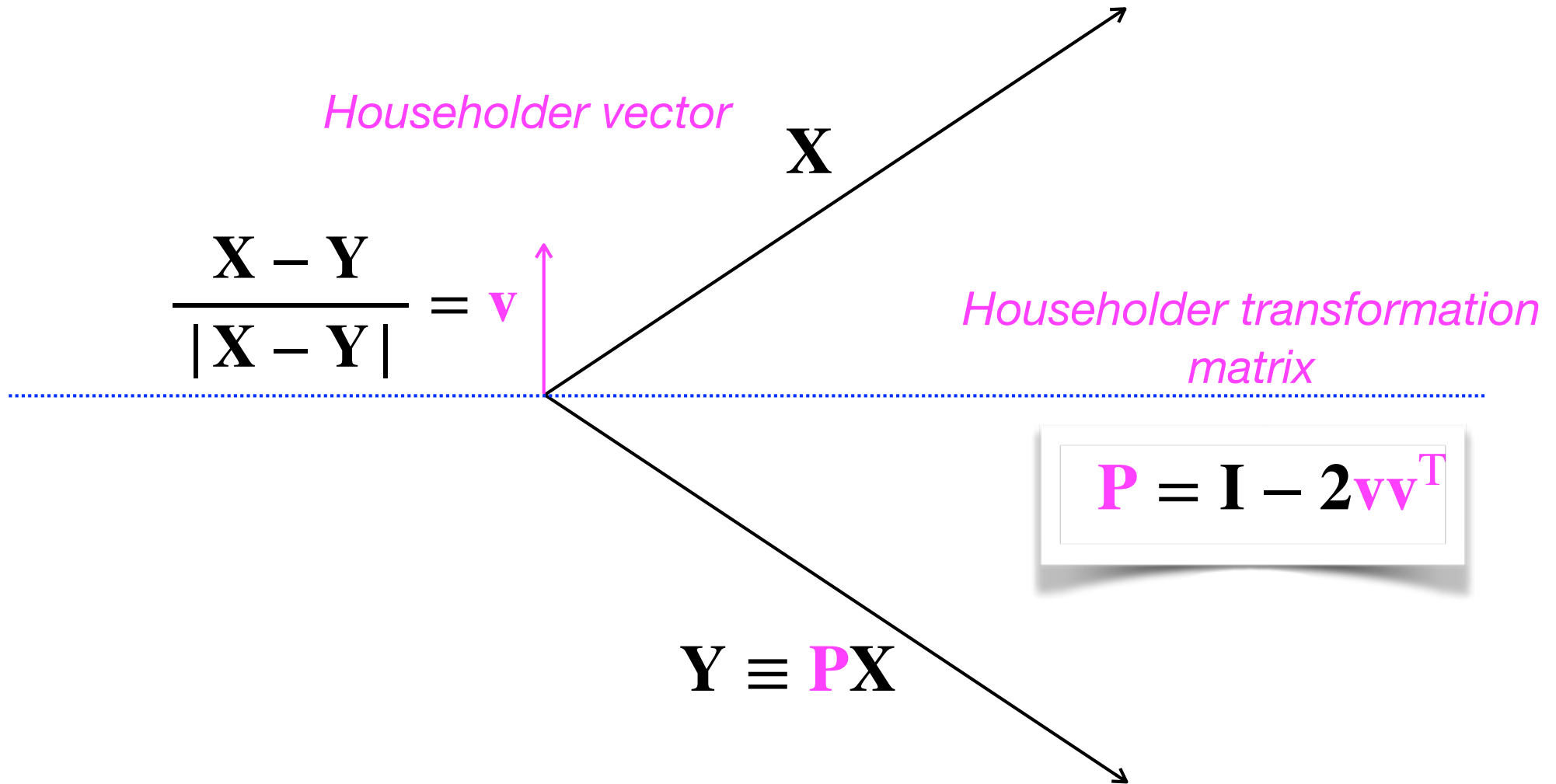
# The Householder transformation



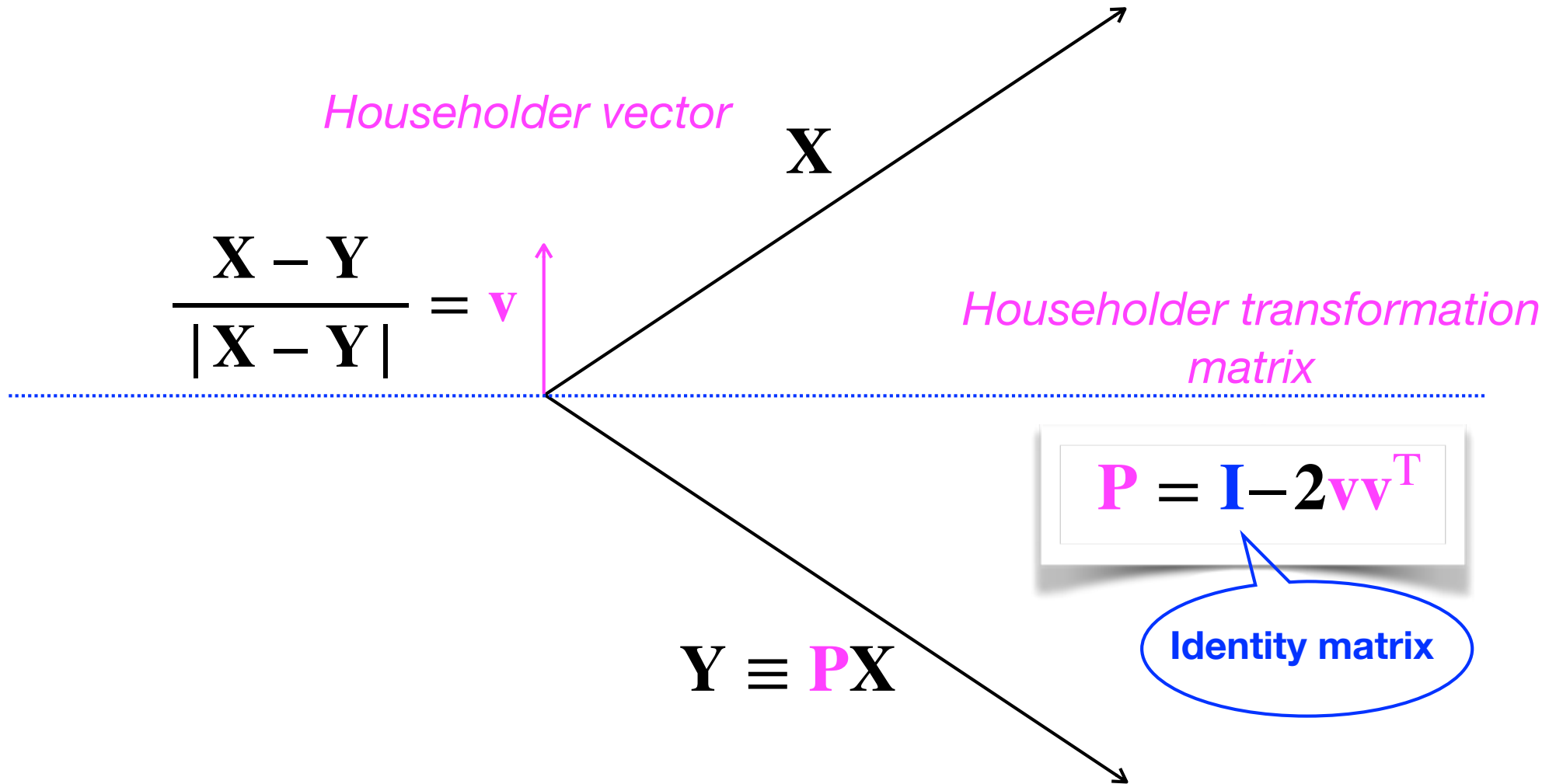
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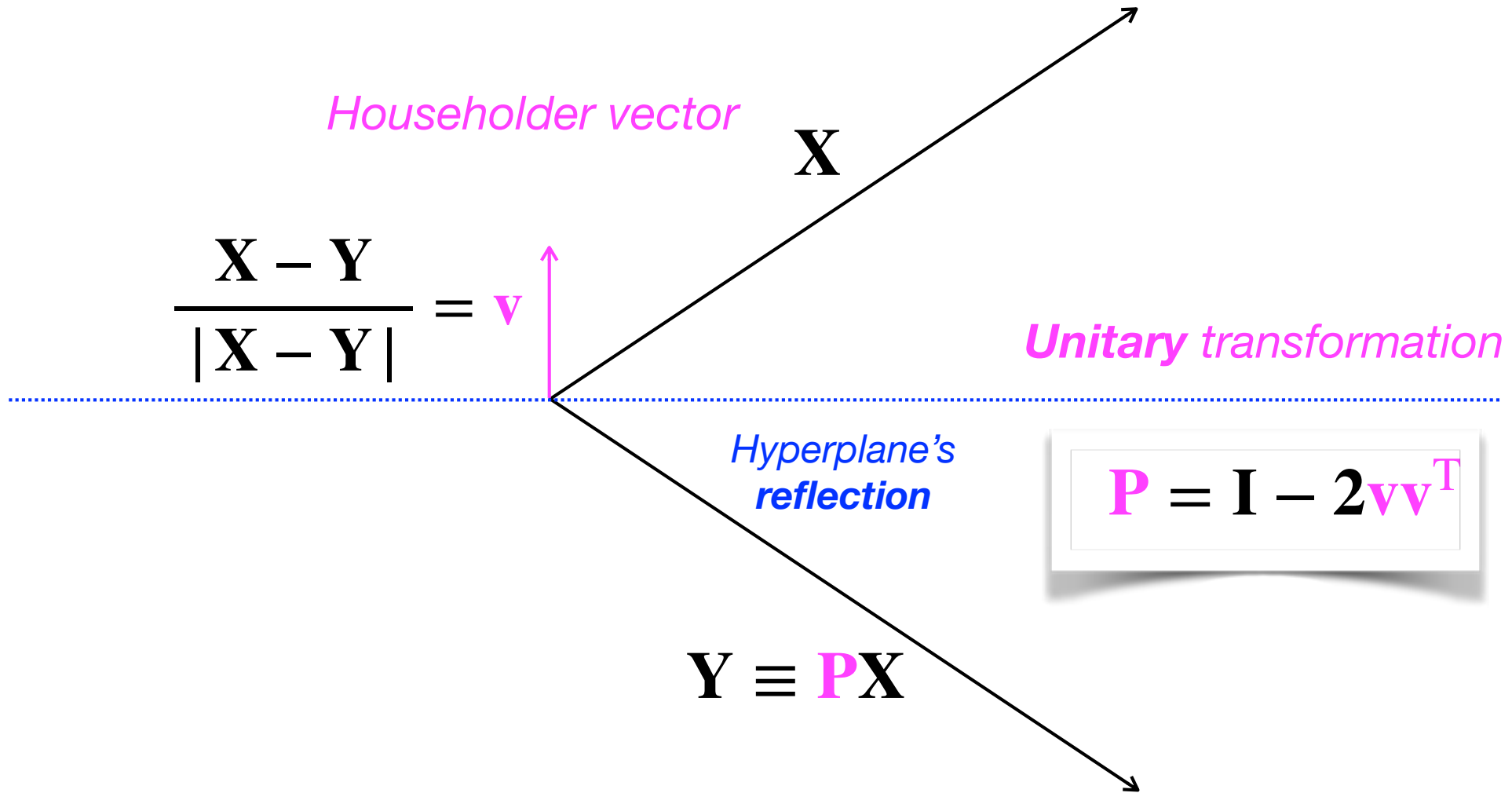


# The Householder transformation





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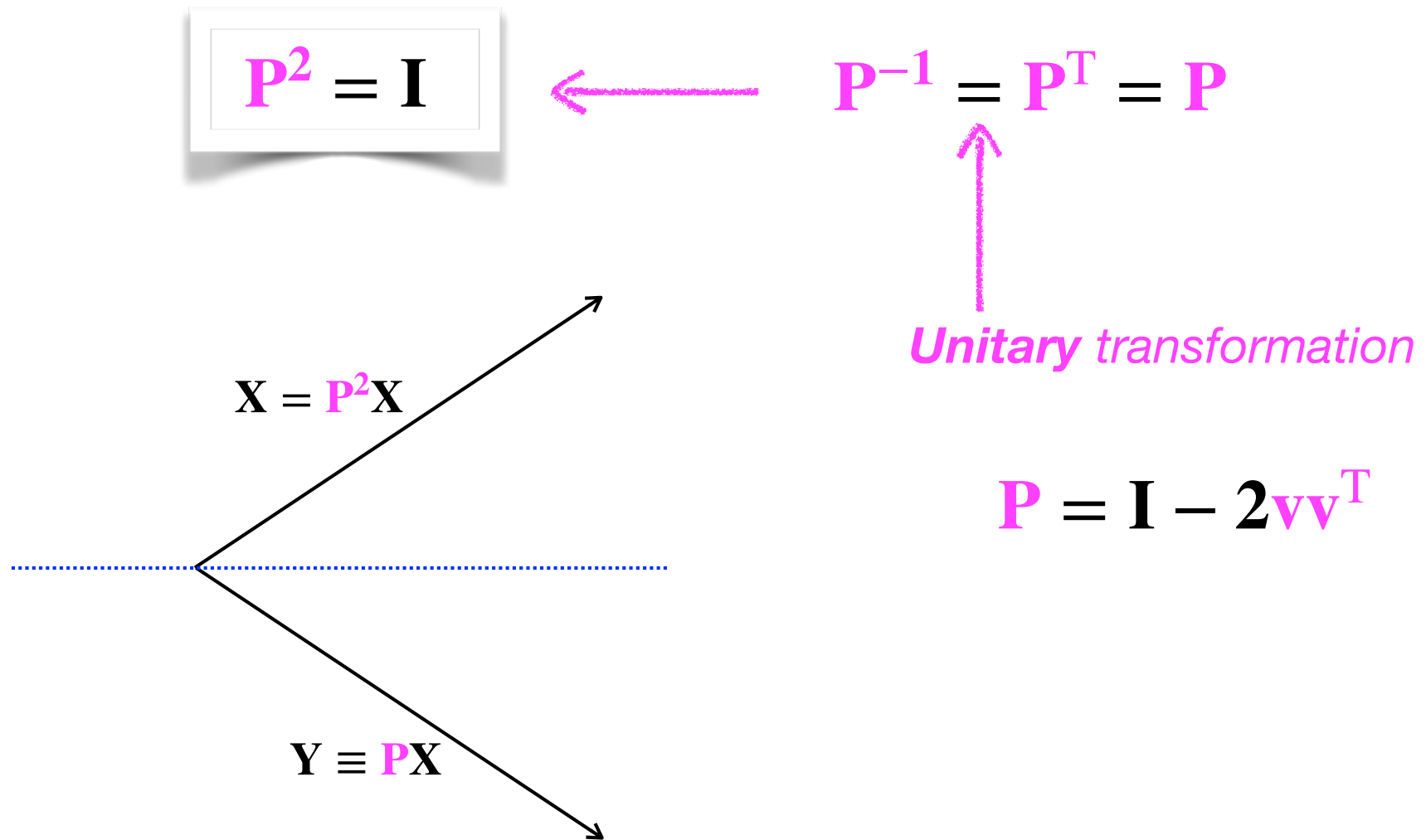
# The Householder transformation

$$\mathbf{P}^{-1} = \mathbf{P}^T = \mathbf{P}$$

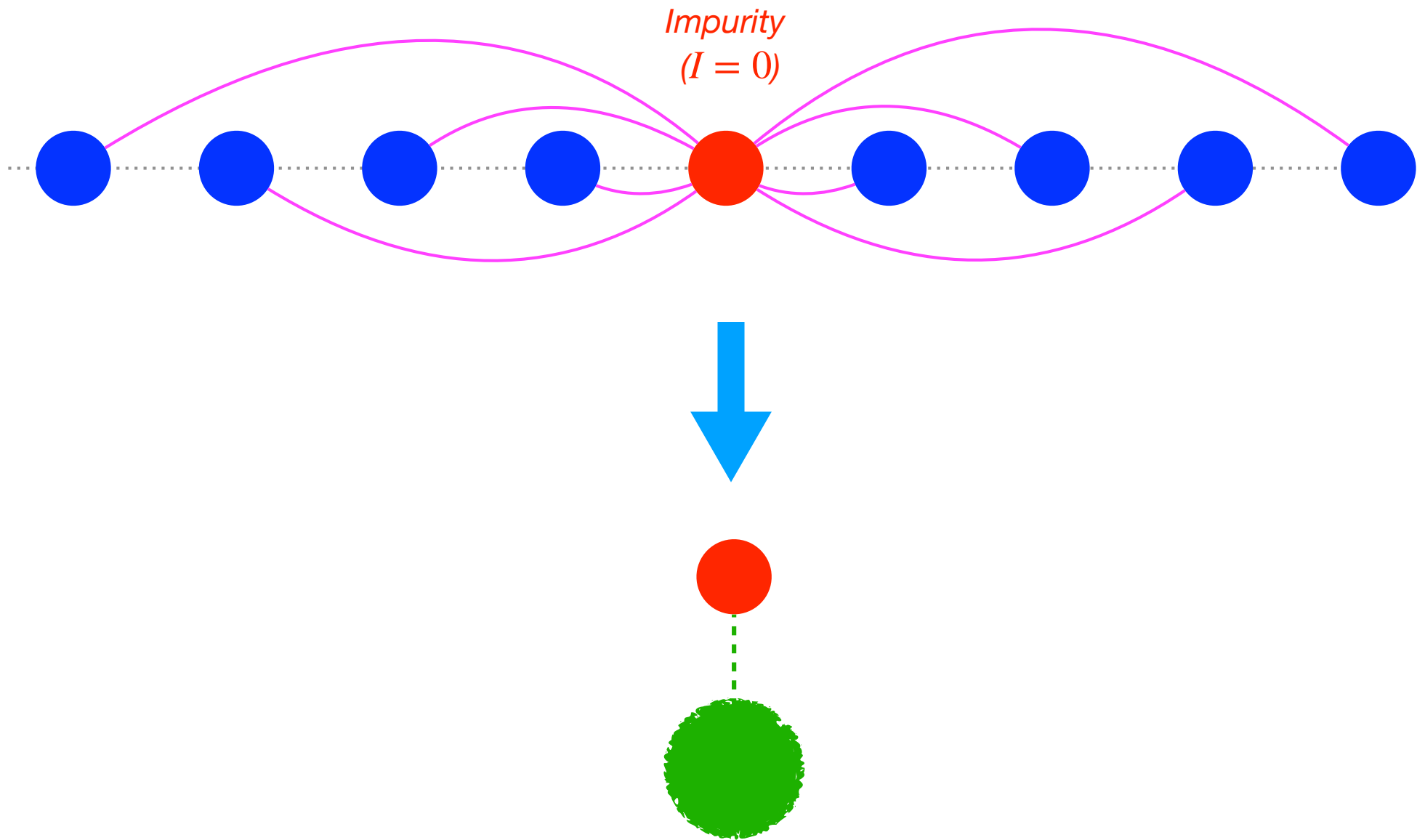
*Unitary transformation*

$$\mathbf{P} = \mathbf{I} - 2\mathbf{v}\mathbf{v}^T$$

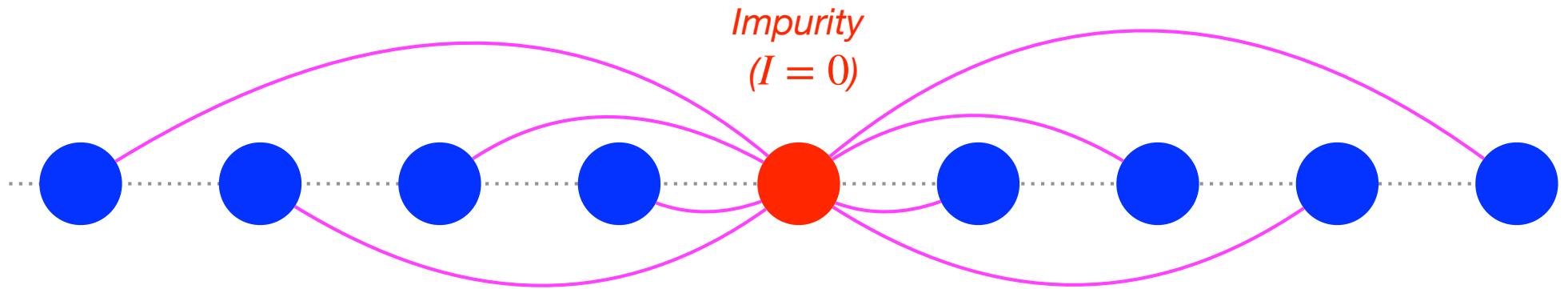
# The Householder transformation



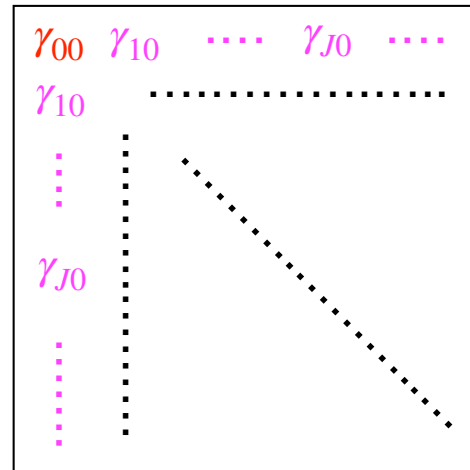
# Householder transformed density matrix embedding



# Householder transformed density matrix embedding

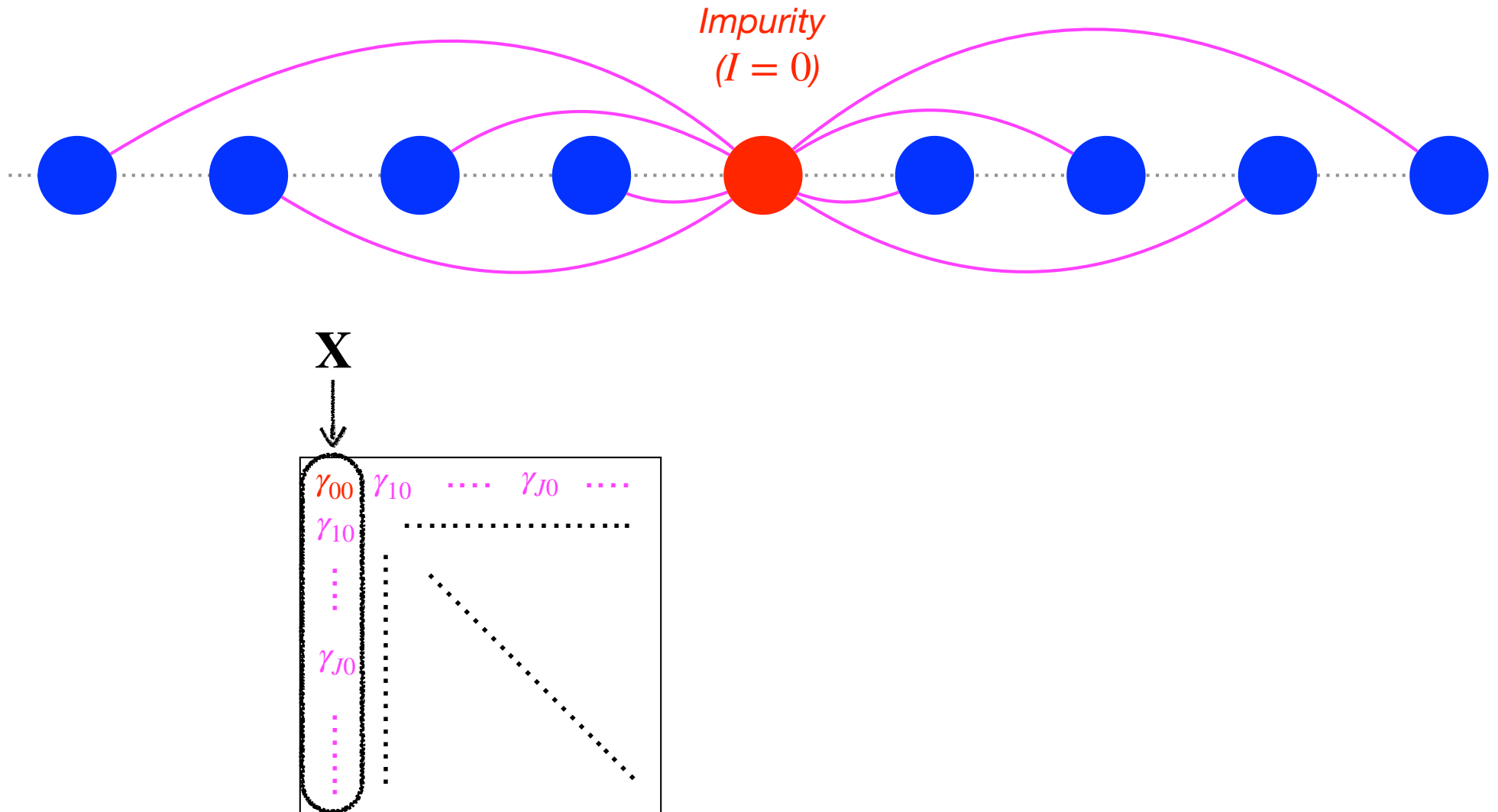


$$\gamma_{IJ} = \langle \hat{c}_I^\dagger \hat{c}_J \rangle \equiv$$

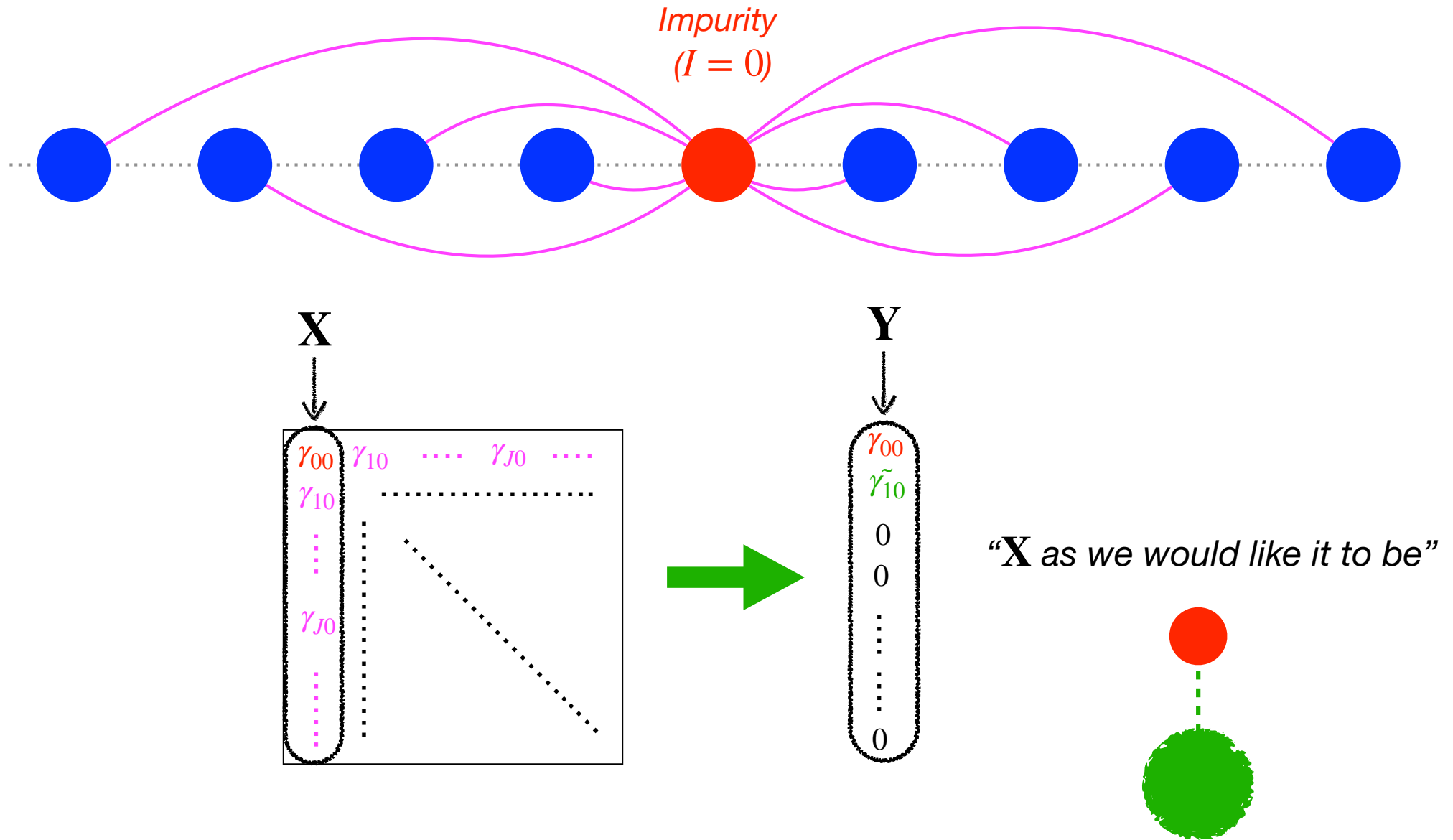


$$\gamma^{loc} \stackrel{\text{notation}}{=} \gamma$$

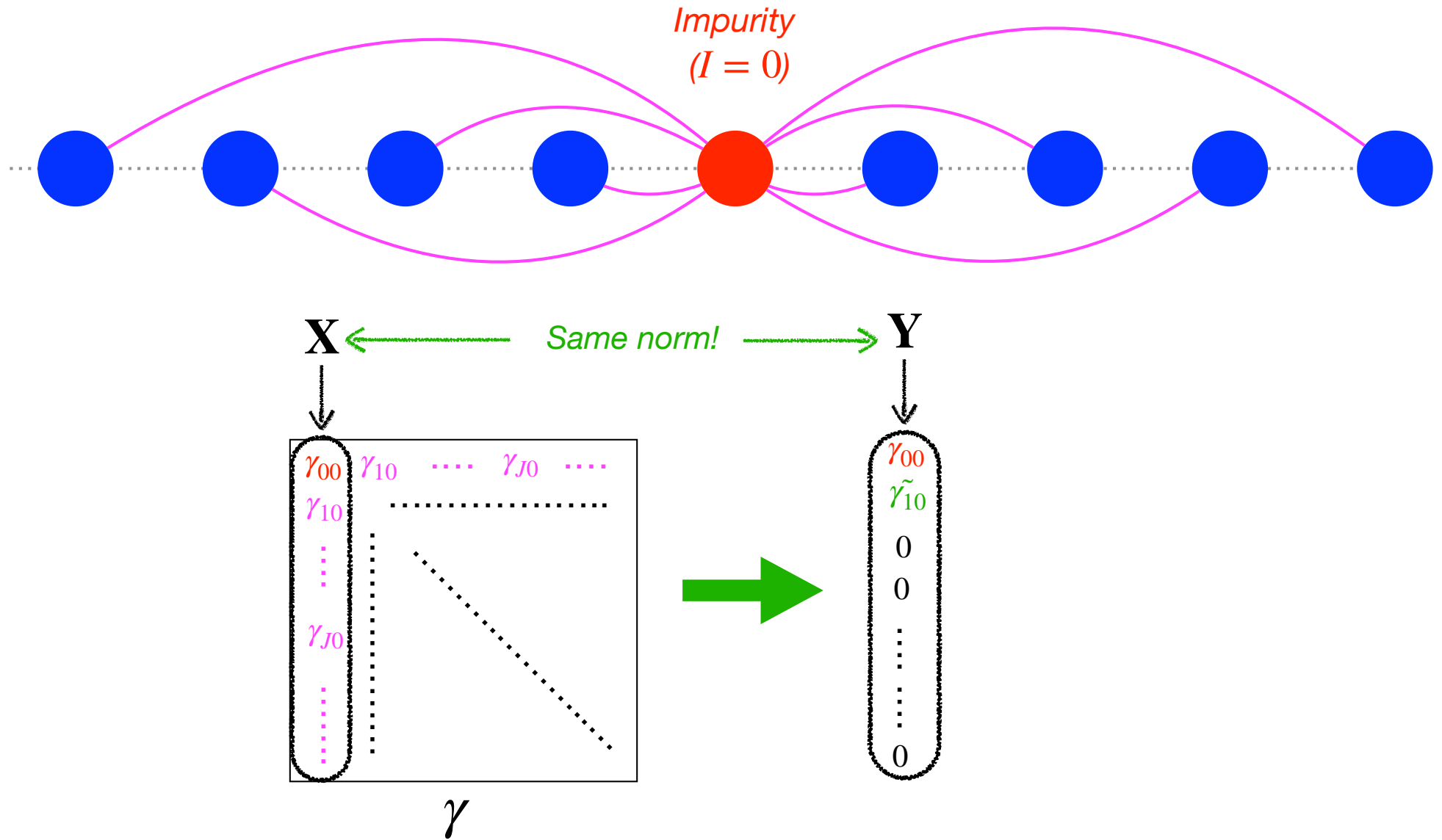
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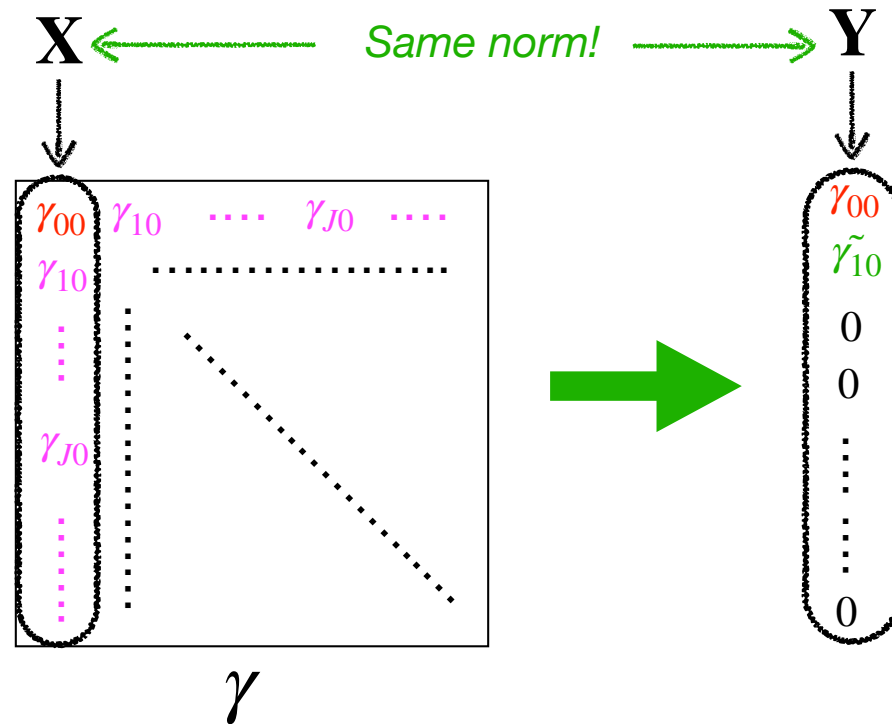
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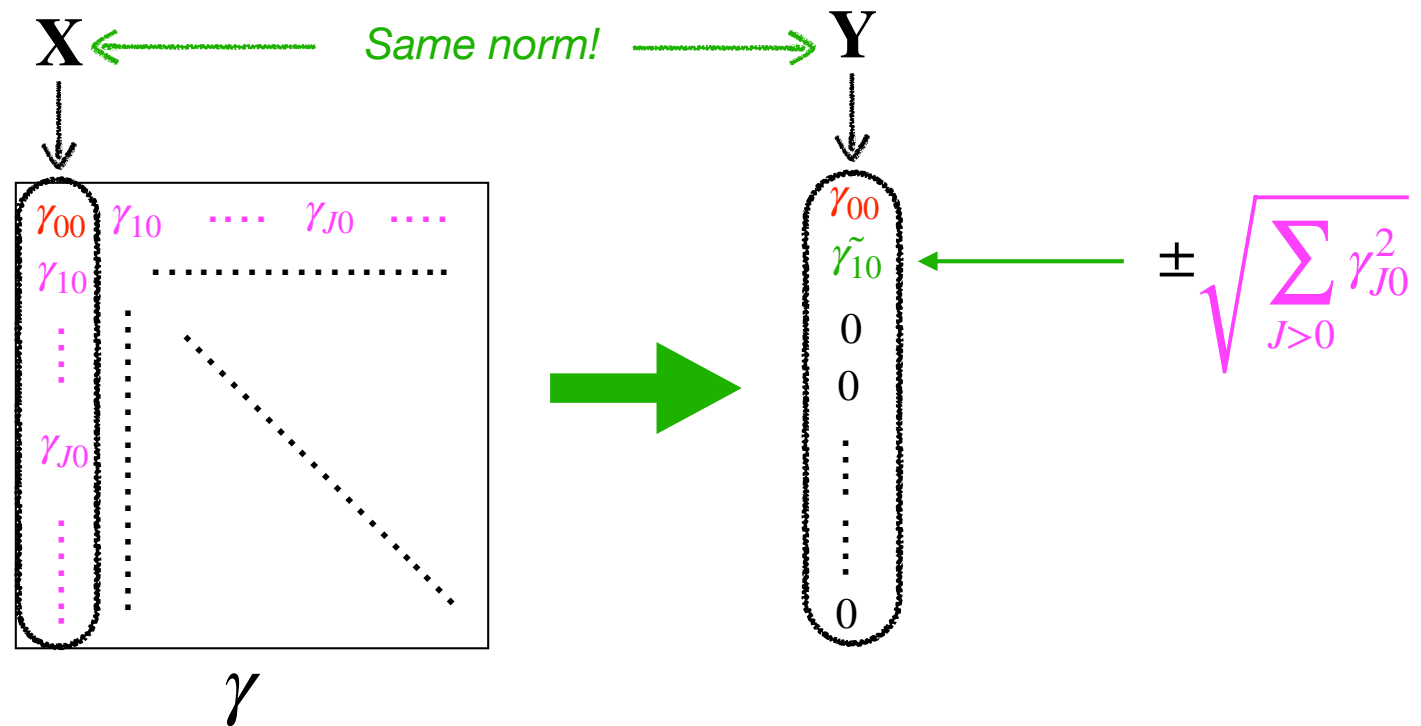


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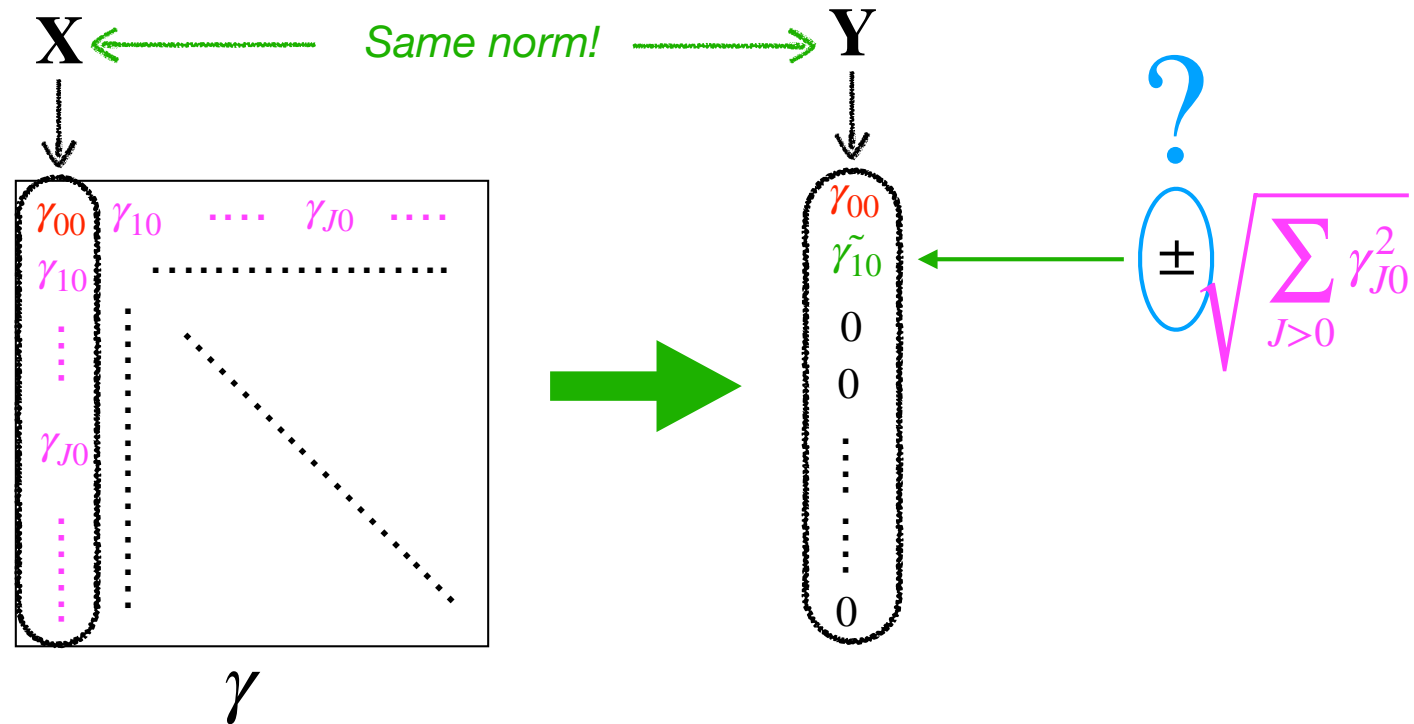
$$|\mathbf{X}|^2 - \gamma_{00}^2 = |\mathbf{Y}|^2 - \gamma_{00}^2 = \tilde{\gamma}_{10}^2 = \sum_{J>0} \gamma_{J0}^2$$



# Householder transformed density matrix embedding



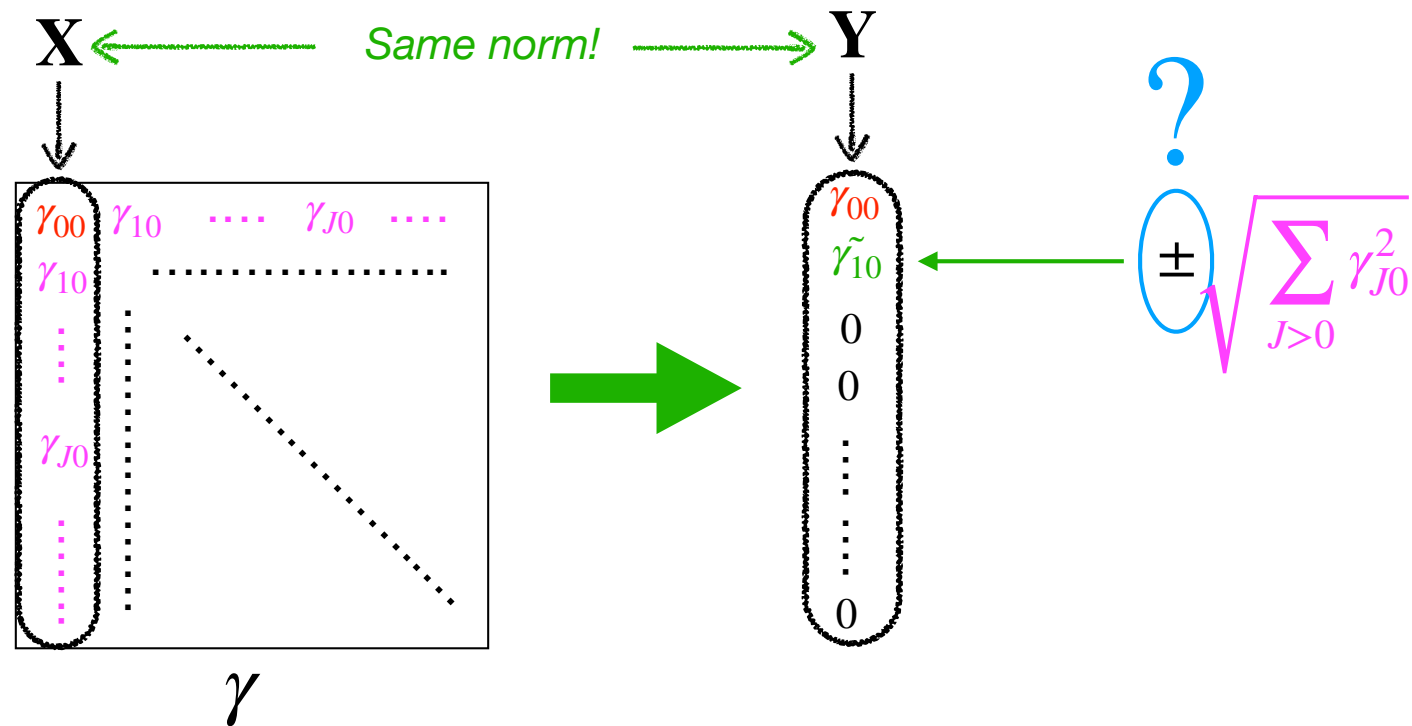
# Householder transformed density matrix embedding



# Householder transformed density matrix embedding

$$\mathbf{v} = \frac{\mathbf{X} - \mathbf{Y}}{|\mathbf{X} - \mathbf{Y}|} \quad \leftarrow \text{Householder vector}$$

where  $|\mathbf{X} - \mathbf{Y}|^2 = 2 \left( |\mathbf{Y}|^2 - \mathbf{X}^T \mathbf{Y} \right) = 2 \left( \tilde{\gamma}_{10}^2 - \gamma_{10} \tilde{\gamma}_{10} \right) = 2 \tilde{\gamma}_{10} \left( \tilde{\gamma}_{10} - \gamma_{10} \right)$

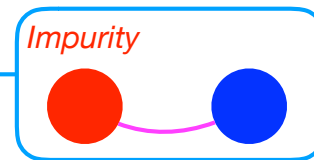


# Householder transformed density matrix embedding

$$\mathbf{v} = \frac{\mathbf{X} - \mathbf{Y}}{|\mathbf{X} - \mathbf{Y}|} \quad \text{Householder vector}$$

where

$$|\mathbf{X} - \mathbf{Y}|^2 = \pm 2 \sqrt{\sum_{J>0} \gamma_{J0}^2} \left( \pm \sqrt{\sum_{J>0} \gamma_{J0}^2} - \gamma_{10} \right)$$



If one single neighbour...

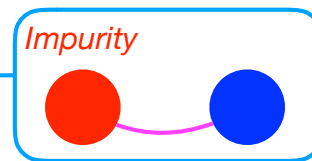
$$|\mathbf{X} - \mathbf{Y}|^2 = \pm 2 |\gamma_{10}| (\pm |\gamma_{10}| - \gamma_{10})$$

# Householder transformed density matrix embedding

$$\mathbf{v} = \frac{\mathbf{X} - \mathbf{Y}}{|\mathbf{X} - \mathbf{Y}|} \quad \text{Householder vector}$$

where

$$|\mathbf{X} - \mathbf{Y}|^2 = (\pm) 2 \sqrt{\sum_{J>0} \gamma_{J0}^2} \left( (\pm) \sqrt{\sum_{J>0} \gamma_{J0}^2} - \gamma_{10} \right)$$

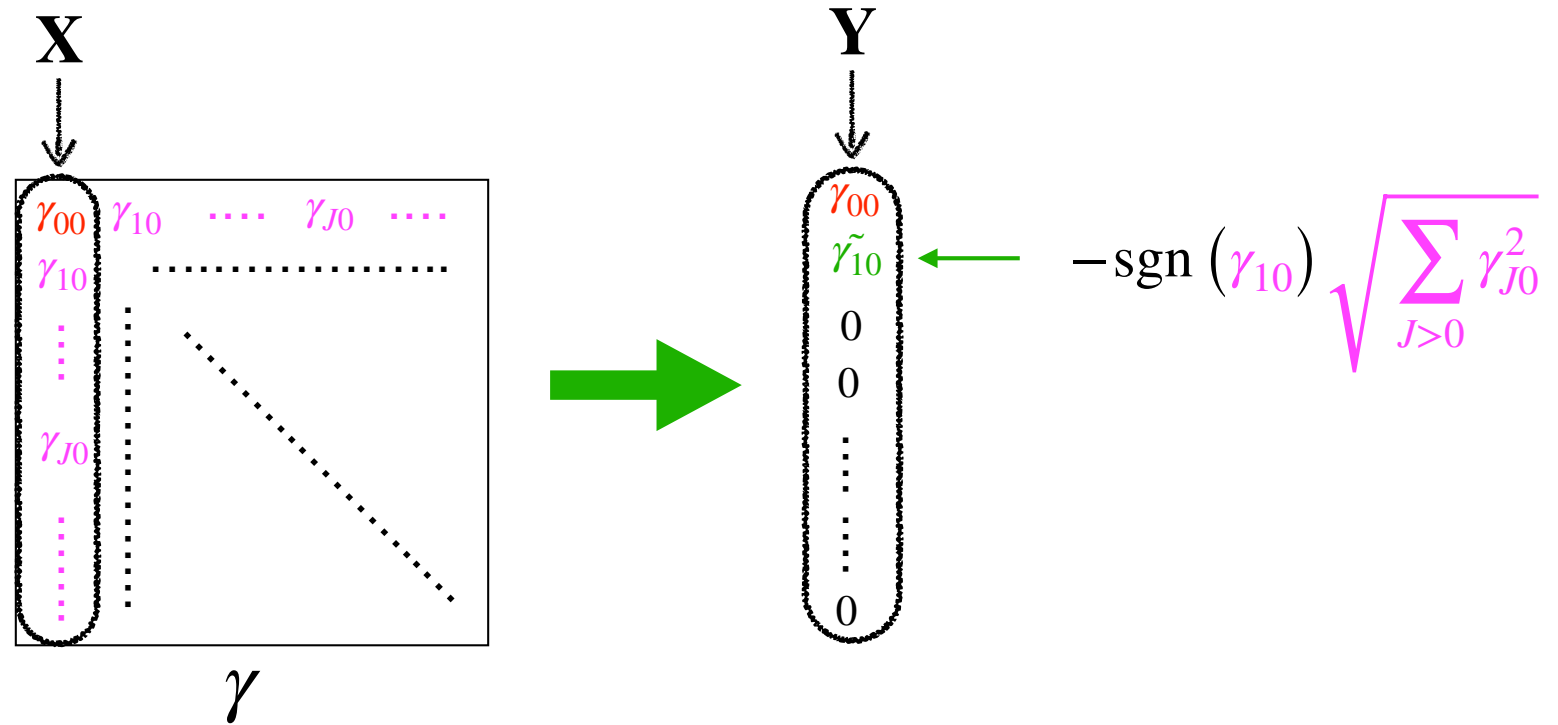


If one single neighbour...

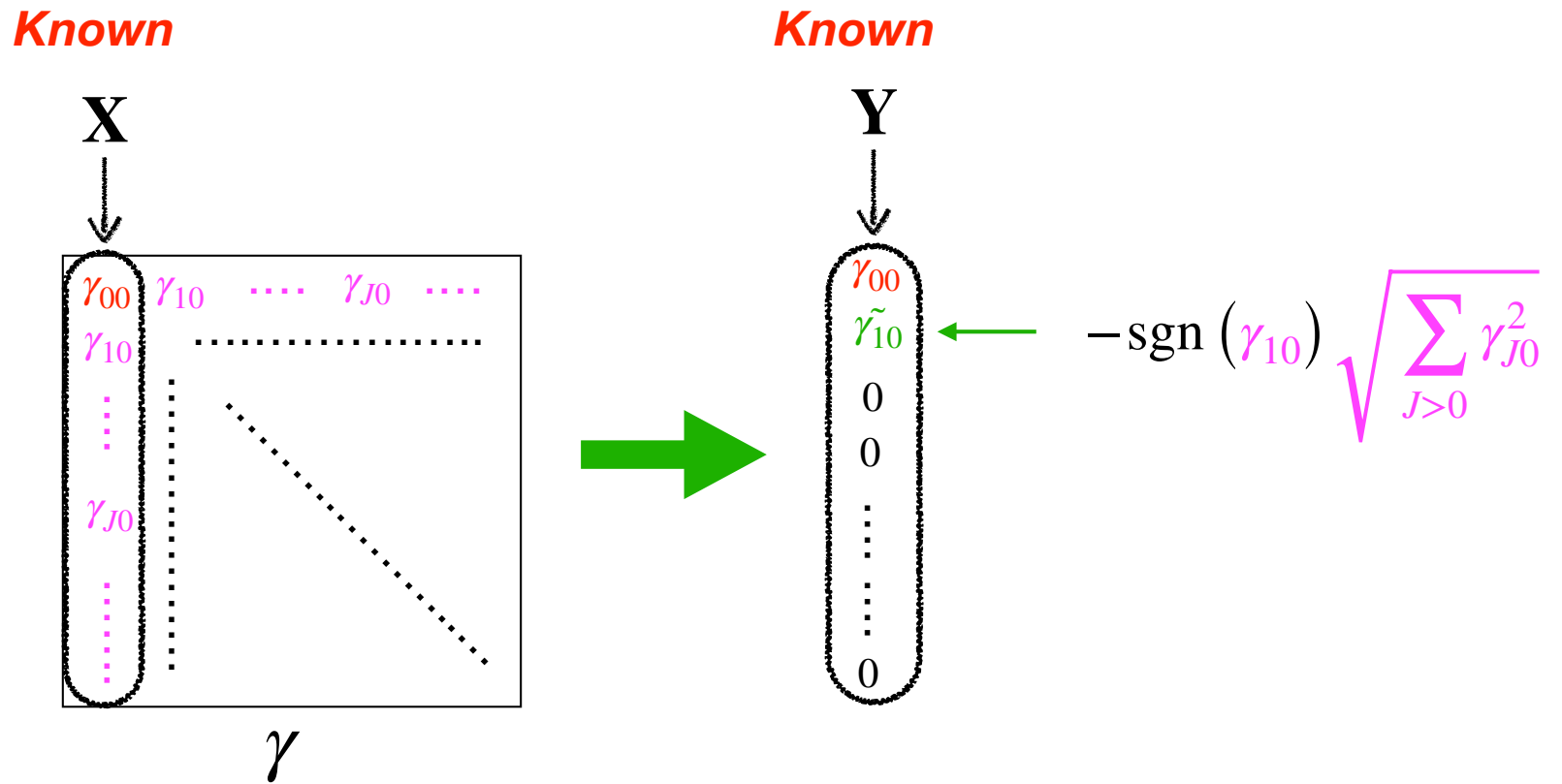
$$|\mathbf{X} - \mathbf{Y}|^2 = (\pm) 2 |\gamma_{10}| \left( (\pm) |\gamma_{10}| - \gamma_{10} \right)$$

choose  $-\text{sgn}(\gamma_{10}) \leftarrow |\mathbf{X} - \mathbf{Y}| \neq 0$

# Householder transformed density matrix embedding

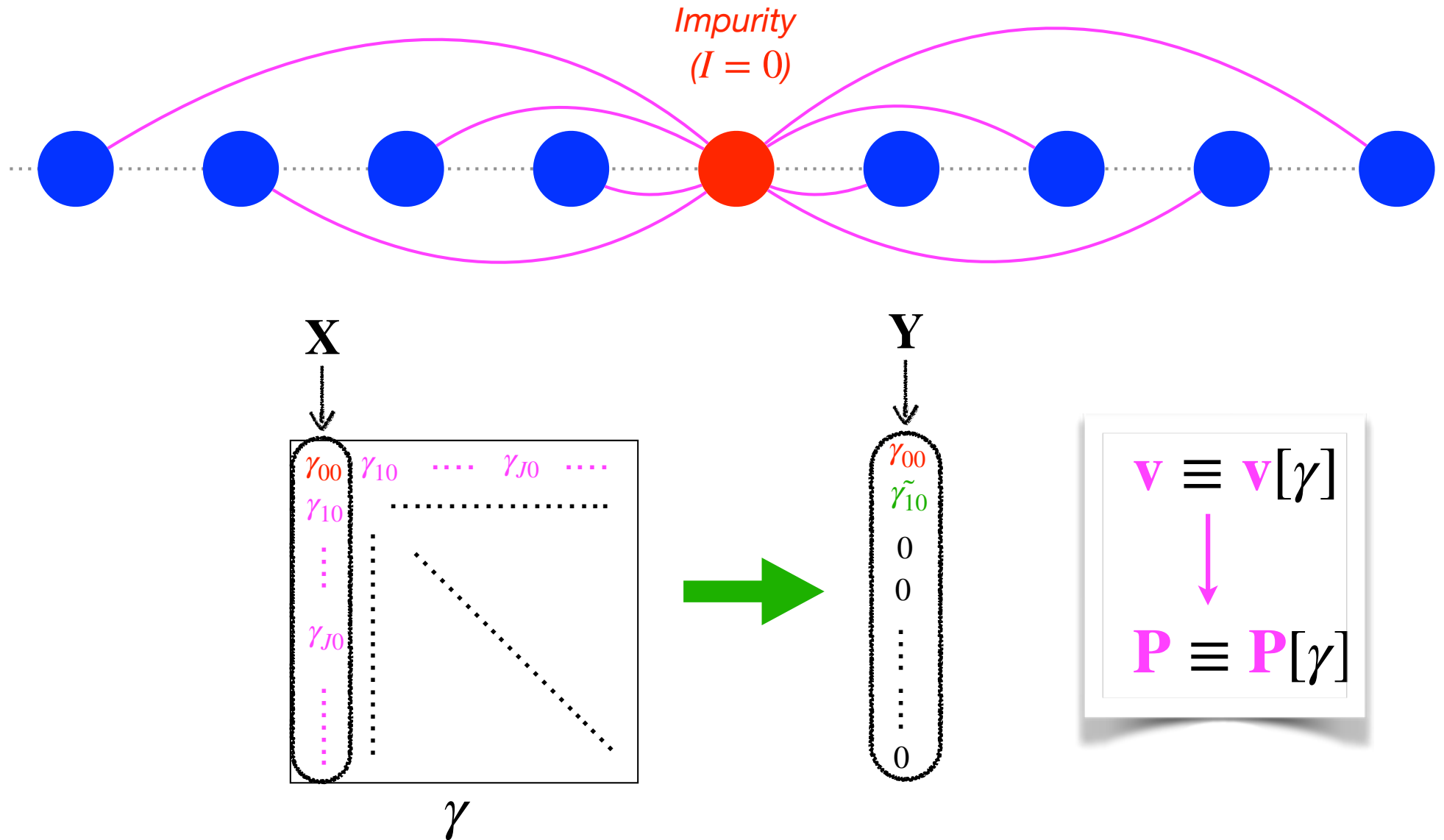


# Householder transformed density matrix embedding





# Householder transformed density matrix embedding



The Householder transformation is an **explicit functional** of the density matrix!

## *Householder representation in second quantization*

$$\mathbf{P} \equiv \mathbf{P}[\gamma] = \mathbf{I} - 2\mathbf{v}\mathbf{v}^T$$

*Unitary Householder transformation matrix*

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From the *localised* to the **Householder** representation

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$$\hat{d}_I^\dagger = \sum_J P_{IJ} \hat{c}_J^\dagger$$

From the **localised** to the **Householder** representation

$$\sum_I P_{KI} \hat{d}_I^\dagger = \sum_J \sum_I P_{KI} P_{IJ} \hat{c}_J^\dagger$$

## Householder representation in second quantization

$$\hat{d}_I^\dagger = \sum_J P_{IJ} \hat{c}_J^\dagger$$

From the **localised** to the **Householder** representation

$$\sum_I P_{KI} \hat{d}_I^\dagger = \sum_J \sum_I P_{KI} P_{IJ} \hat{c}_J^\dagger = \sum_J [\mathbf{P}^2]_{KJ} \hat{c}_J^\dagger = \sum_J \delta_{KJ} \hat{c}_J^\dagger = \hat{c}_K^\dagger$$

## Householder representation in second quantization

$$\hat{d}_I^\dagger = \sum_J P_{IJ} \hat{c}_J^\dagger$$

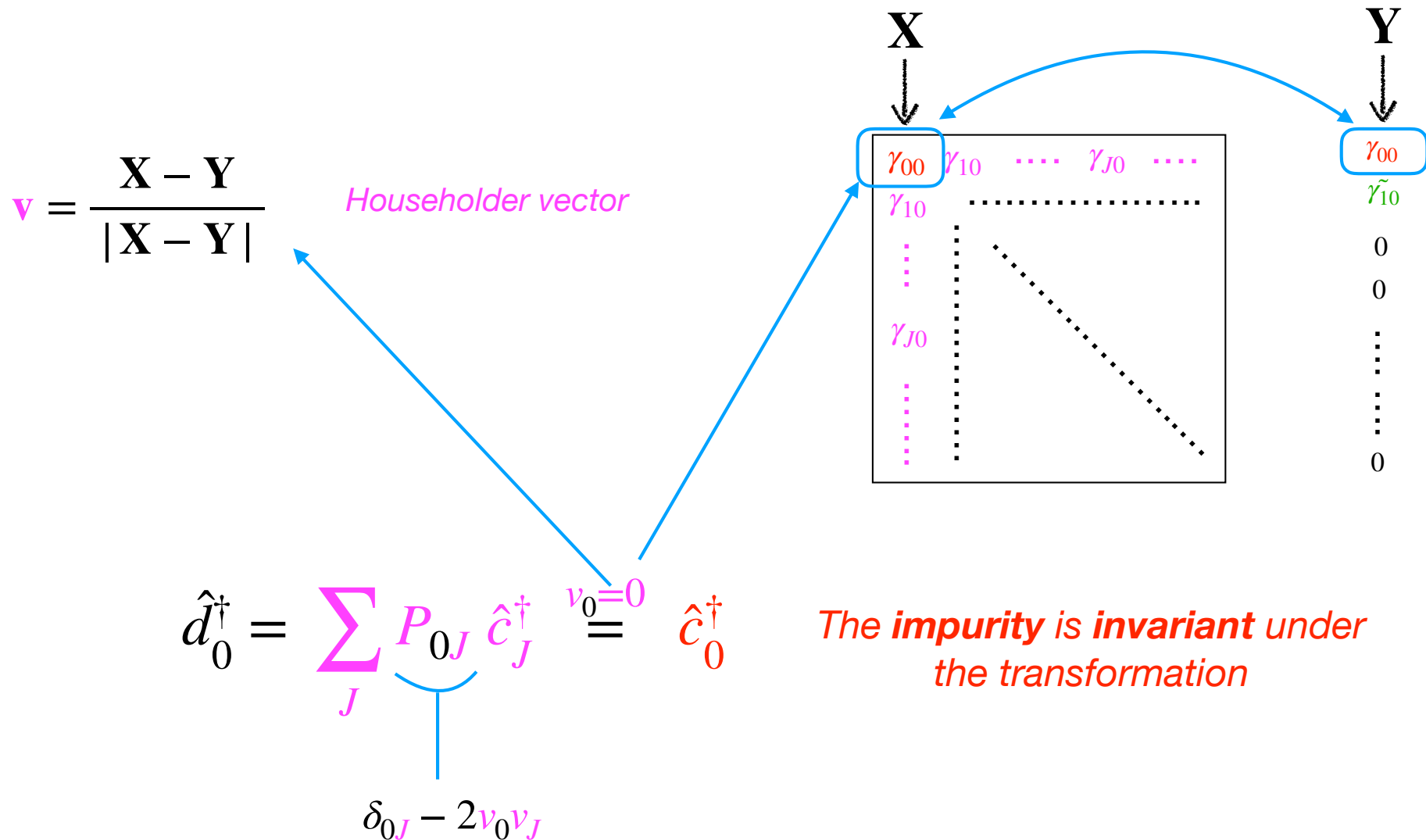
From the **localised** to the **Householder** representation

$$\hat{c}_K^\dagger = \sum_I P_{KI} \hat{d}_I^\dagger$$

From the **Householder** to the **localised** representation



# Householder representation in second quantization



## Householder representation in second quantization

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Unitary Householder transformation matrix

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Creates delocalised **Householder orbitals**

$$\hat{d}_0^\dagger = \hat{c}_0^\dagger$$

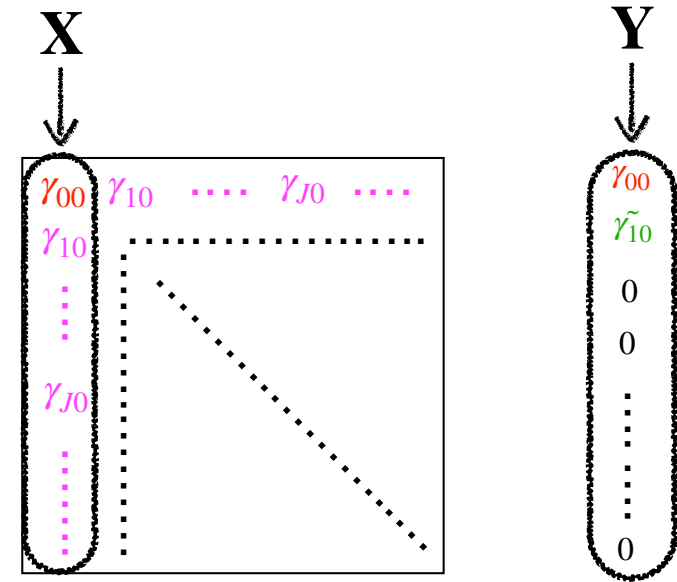
The **impurity** is **invariant** under the transformation

$$\hat{d}_1^\dagger |\text{vac}\rangle = \sum_J P_{1J} |\chi_J\rangle$$

Will play the role of the **bath orbital**

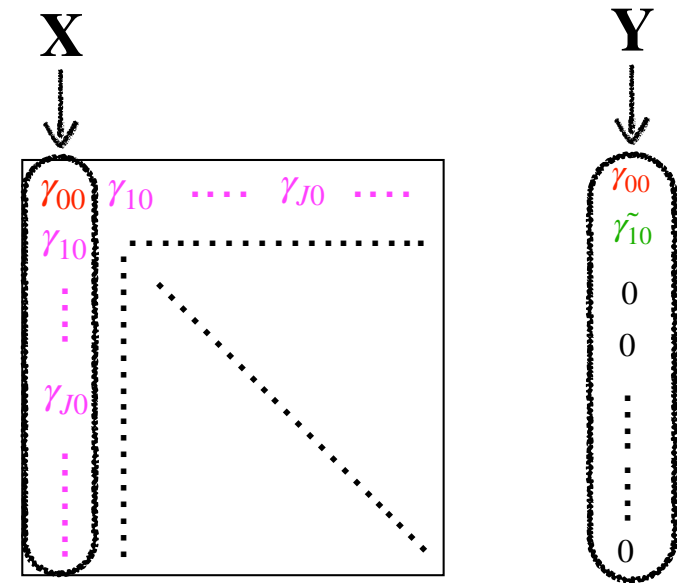
# 1RDM in the Householder representation

$$\begin{aligned}
 \tilde{\gamma}_{J0} &= \langle \Psi_0 | \hat{d}_J^\dagger \hat{d}_0 | \Psi_0 \rangle \\
 &= \sum_I P_{JI} \langle \Psi_0 | \hat{c}_I^\dagger \hat{c}_0 | \Psi_0 \rangle \\
 &= \sum_I P_{JI} \gamma_{I0} \\
 &= [\mathbf{P}\mathbf{X}]_J \\
 &= [\mathbf{Y}]_J \\
 &\stackrel{J>1}{=} 0
 \end{aligned}$$



# 1RDM in the Householder representation

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 \end{aligned}$$



By construction, the *impurity* is *entangled only with the bath*

## *1RDM in the Householder representation*

**Theorem:** As the electrons are **non-interacting**, the **bath** turns out to be **entangled only with the impurity**

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Proof

$$\tilde{\gamma} = \tilde{\gamma}^2$$

**Idempotency property**

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Proof

$$\tilde{\gamma} = \tilde{\gamma}^2 \quad \text{Idempotency property}$$

$$\begin{aligned} \tilde{\gamma}_{J0} &= \langle \Psi_0 | \hat{d}_J^\dagger \hat{d}_0 | \Psi_0 \rangle = [\tilde{\gamma}^2]_{J0} = \sum_K \tilde{\gamma}_{JK} \tilde{\gamma}_{K0} \\ &= \tilde{\gamma}_{J0} \tilde{\gamma}_{00} + \tilde{\gamma}_{J1} \tilde{\gamma}_{10} + \sum_{K>1} \tilde{\gamma}_{JK} \times 0 \end{aligned}$$

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Proof

$$\tilde{\gamma} = \tilde{\gamma}^2 \quad \text{Idempotency property}$$

$$\tilde{\gamma}_{J0} = \tilde{\gamma}_{J0}\tilde{\gamma}_{00} + \tilde{\gamma}_{J1}\tilde{\gamma}_{10}$$



$$\tilde{\gamma}_{J1} = \frac{\tilde{\gamma}_{J0} (1 - \tilde{\gamma}_{00})}{\tilde{\gamma}_{10}}$$



# 1RDM in the Householder representation

**Theorem:** As the electrons are **non-interacting**, the **bath** turns out to be **entangled only with the impurity**

Proof

$$\tilde{\gamma}_{J1} = \frac{\tilde{\gamma}_{J0} (1 - \tilde{\gamma}_{00})}{\tilde{\gamma}_{10}}$$

$$J > 1 \quad \tilde{\gamma}_{J0} = 0$$

No entanglement between the impurity and the orbitals other than the bath

$$\tilde{\gamma}_{J1} = 0$$

# 1RDM in the Householder representation

**Theorem:** As the electrons are **non-interacting**, the **bath** turns out to be **entangled only with the impurity**

Proof

**No entanglement** between the **bath** and the orbitals other than the **impurity!**



$$\tilde{\gamma}_{J1} = \langle \Psi_0 | \hat{d}_J^\dagger \hat{d}_1 | \Psi_0 \rangle = 0$$


# 1RDM in the Householder representation

**Theorem:** As the electrons are **non-interacting**, the **bath** turns out to be **entangled only with the impurity**

Proof

$$\tilde{\gamma}_{J1} = \frac{\tilde{\gamma}_{J0} (1 - \tilde{\gamma}_{00})}{\tilde{\gamma}_{10}}$$

$J = 1$


$$\tilde{\gamma}_{11} + \tilde{\gamma}_{00} = 1$$

## 1RDM in the Householder representation

**Theorem:** As the electrons are **non-interacting**, the **bath** turns out to be **entangled only with the impurity**

Proof

$$\tilde{\gamma}_{11} + \tilde{\gamma}_{00} = \langle \Psi_0 | \hat{d}_1^\dagger \hat{d}_1 | \Psi_0 \rangle + \langle \Psi_0 | \hat{d}_0^\dagger \hat{d}_0 | \Psi_0 \rangle = 1$$

The “**impurity+bath**” cluster contains exactly **one electron (per spin)**

## 1RDM in the Householder representation

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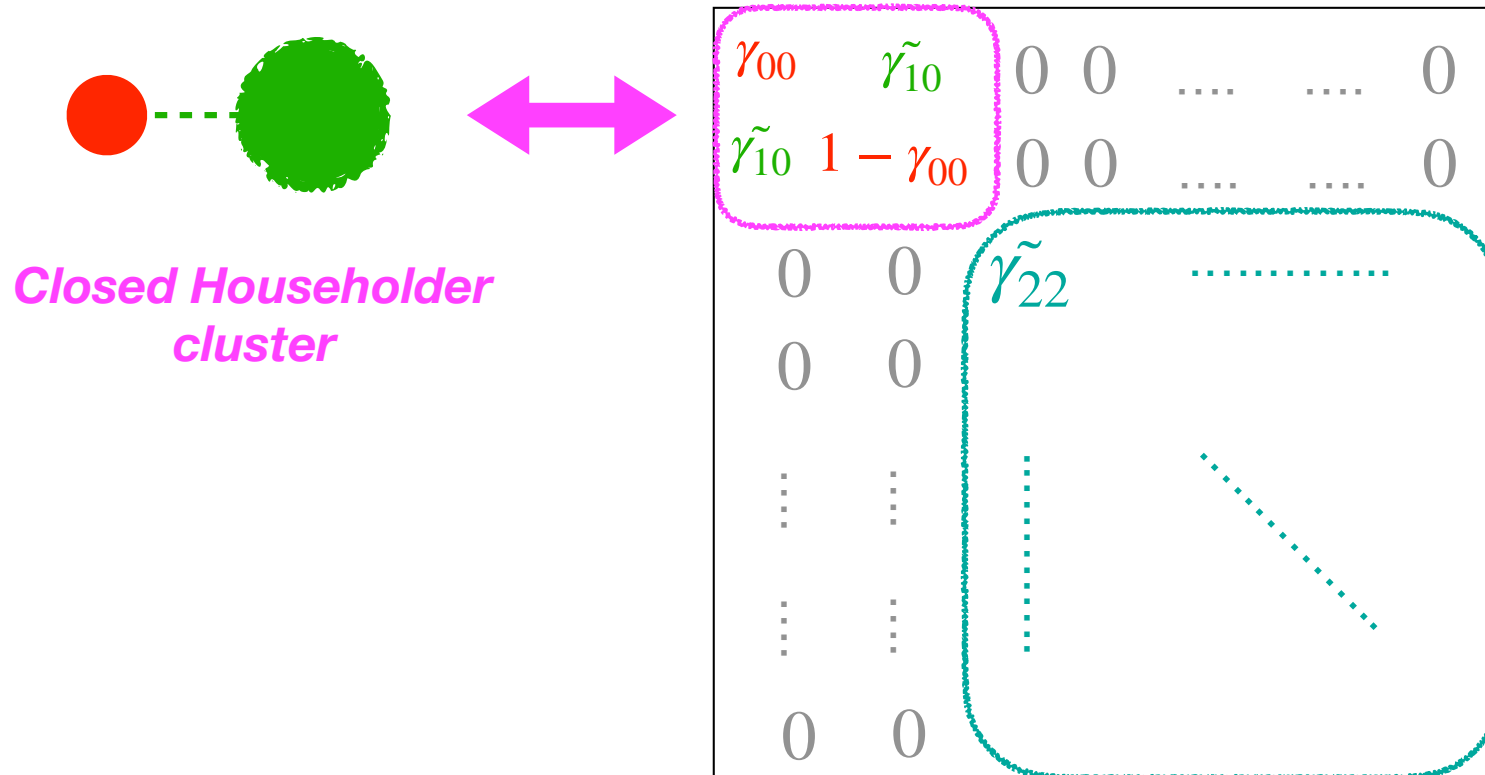
Proof

$$\tilde{\gamma}_{11} + \tilde{\gamma}_{00} = \langle \Psi_0 | \hat{d}_1^\dagger \hat{d}_1 | \Psi_0 \rangle + \langle \Psi_0 | \hat{d}_0^\dagger \hat{d}_0 | \Psi_0 \rangle = 1$$

The “**impurity+bath**” cluster contains exactly **one electron (per spin)**

The cluster is a **closed quantum system** that can be described with a two-electron wave function  $\Psi^{\mathcal{C}}$

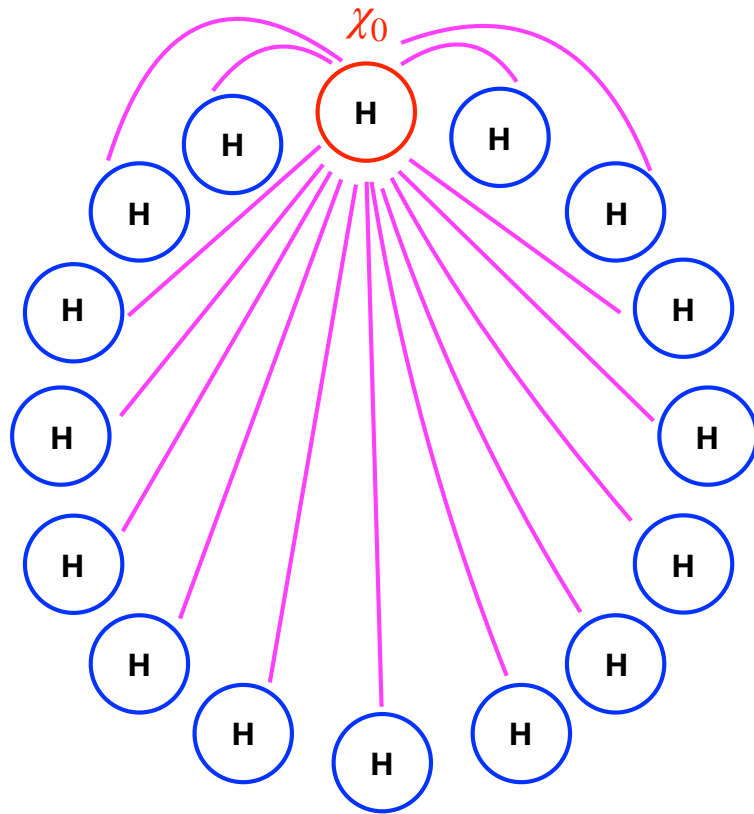
# 1RDM in the Householder representation



## Energy evaluation by fragmentation

$$\hat{H} = \sum_{IJ} \bar{h}_{IJ} \hat{c}_I^\dagger \hat{c}_J \quad \text{Localised representation}$$

## Energy evaluation by fragmentation

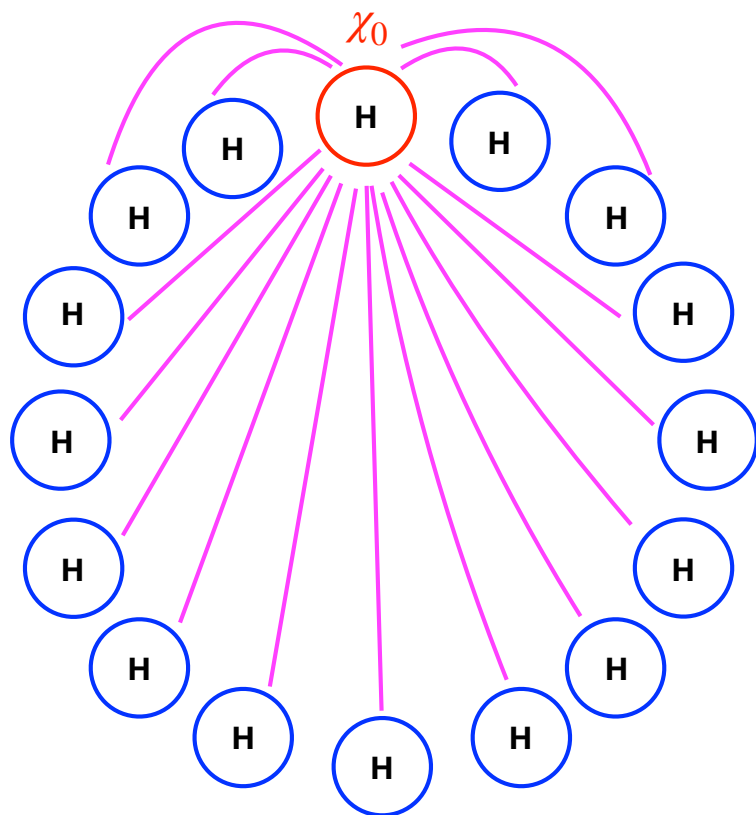


$$\hat{H} = \sum_{IJ} \bar{h}_{IJ} \hat{c}_I^\dagger \hat{c}_J$$

*Localised representation*



## Energy evaluation by fragmentation



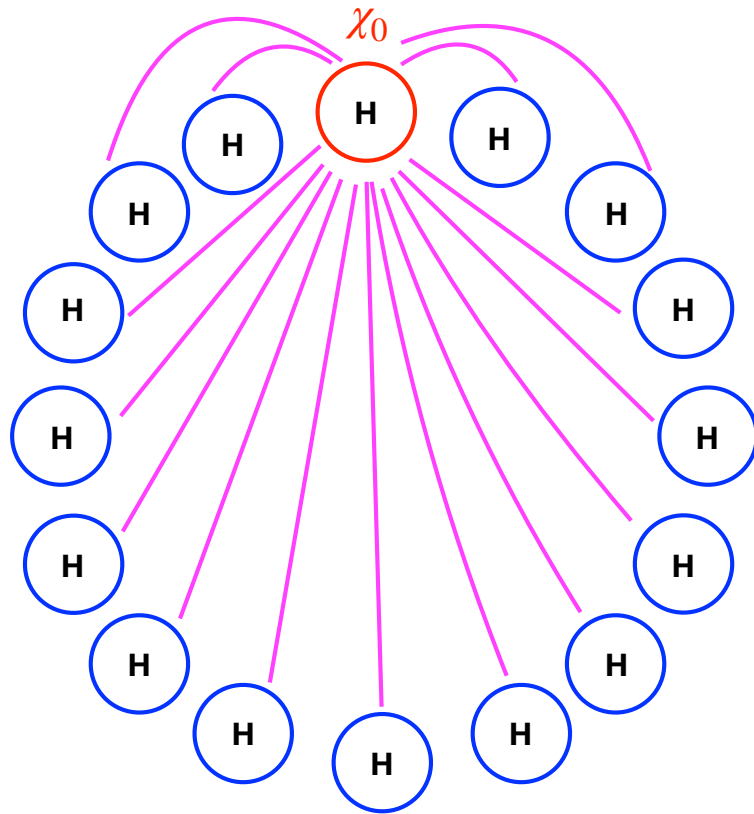
$$\hat{H} = \sum_{IJ} \bar{h}_{IJ} \hat{c}_I^\dagger \hat{c}_J$$

*Localised representation*

*Energy contributions involving the impurity*

$$2 \sum_J \bar{h}_{0J} \langle \Psi_0 | \hat{c}_0^\dagger \hat{c}_J | \Psi_0 \rangle$$

# Energy evaluation by fragmentation



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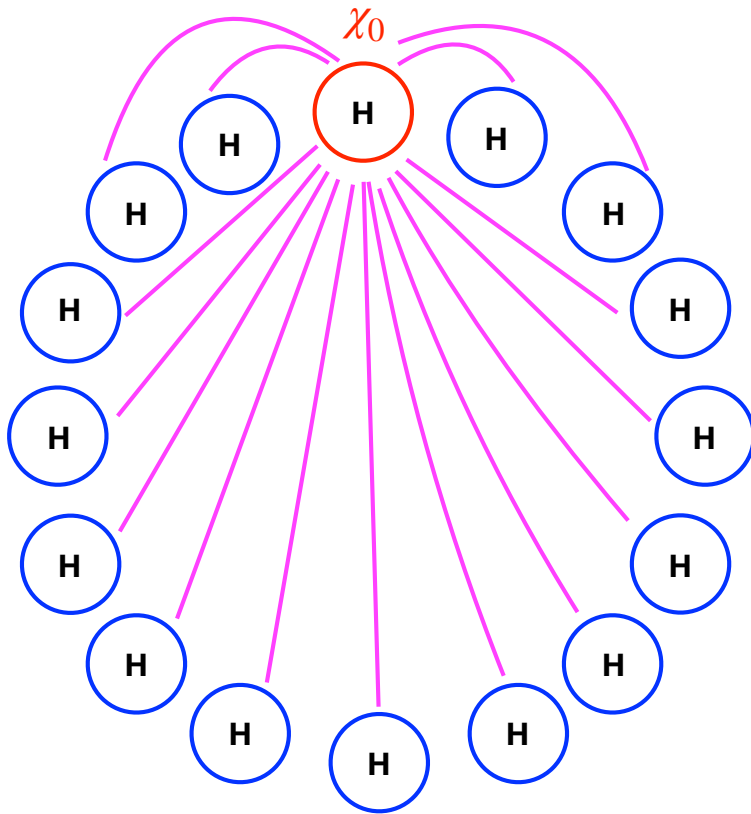
Householder representation



$$= 2 \sum_K \left( \sum_J \bar{h}_{0J} P_{JK} \right) \langle \Psi_0 | \hat{d}_0^\dagger \hat{d}_K | \Psi_0 \rangle$$

$\tilde{h}_{0K}$

## Energy evaluation by fragmentation



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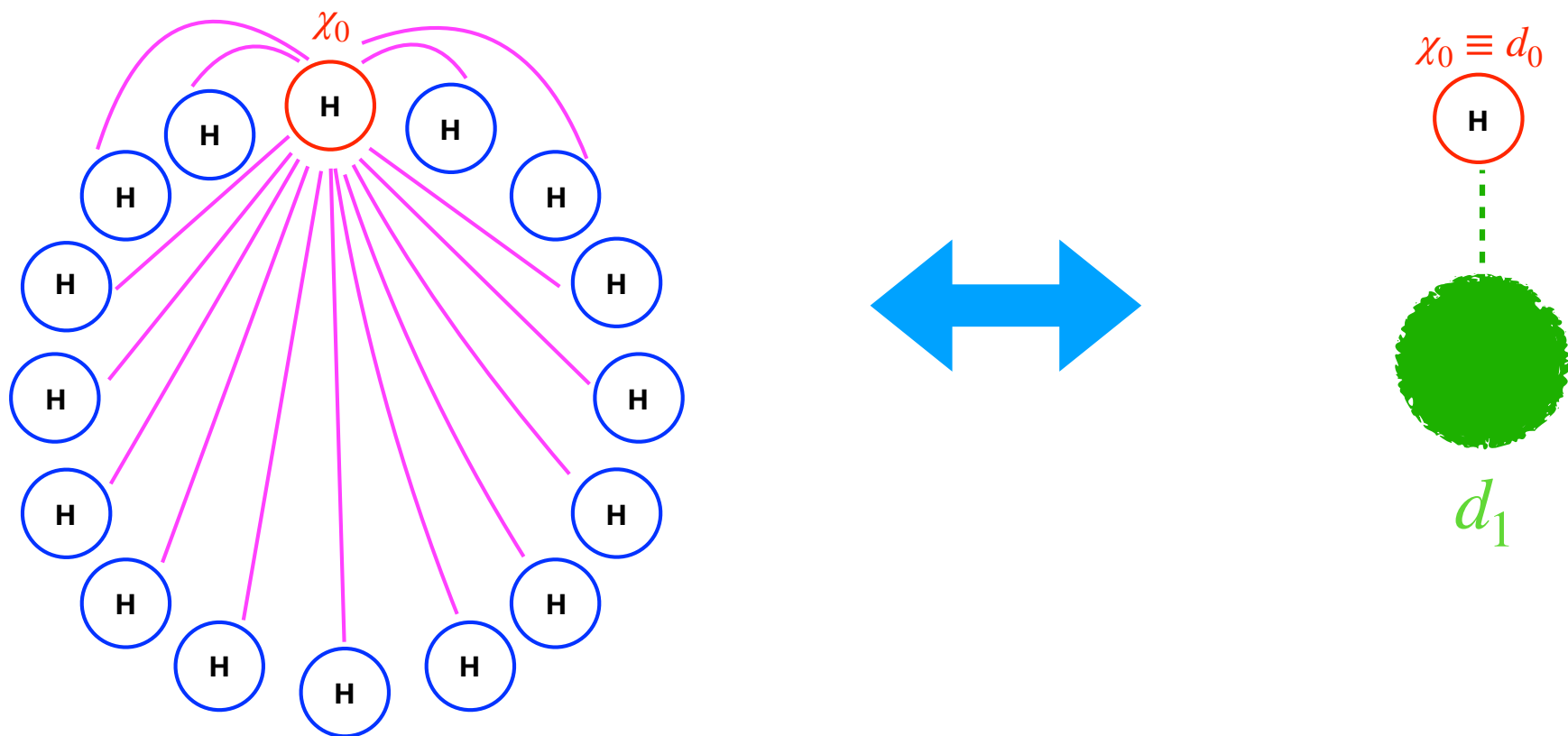
Energy contributions involving the impurity

$$2 \sum_J \bar{h}_{0J} \langle \Psi_0 | \hat{c}_0^\dagger \hat{c}_J | \Psi_0 \rangle$$

$$= 2 \sum_K \left( \sum_J \bar{h}_{0J} P_{JK} \right) \langle \Psi_0 | \hat{d}_0^\dagger \hat{d}_K | \Psi_0 \rangle$$

$$= 2 \sum_K \tilde{h}_{0K} \tilde{\gamma}_{0K} = 2 \left( \tilde{h}_{00} \tilde{\gamma}_{00} + \tilde{h}_{01} \tilde{\gamma}_{01} \right)$$

## Energy evaluation by fragmentation



*Determined from the cluster*

$$2 \sum_J \bar{h}_{0J} \langle \Psi_0 | \hat{c}_0^\dagger \hat{c}_J | \Psi_0 \rangle = 2 \left( \tilde{h}_{00} \tilde{\gamma}_{00} + \tilde{h}_{01} \tilde{\gamma}_{01} \right)$$

## *Approximate embedding for interacting electrons*

*The present embedding approach is **useless for non-interacting electrons** (!)*

## Approximate embedding for interacting electrons

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**We need**  $\gamma_{IJ} = \gamma_{IJ}^{loc} = \sum_P^{occupied\ spin-MOs} C_{IP} C_{JP}$

## Approximate embedding for interacting electrons

We have to solve the Schrödinger equation for the **full system!**

**We need**  $\gamma_{IJ} = \gamma_{IJ}^{loc} = \sum_P^{occupied\ spin-MOs} C_{IP} C_{JP}$

## *Approximate embedding for interacting electrons*

*The present embedding approach is **useless for non-interacting electrons** (!)*

*However, it can be used for describing electron **repulsions** (within the cluster)*



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However, it can be used for describing electron *repulsions (within the cluster)*

$$\hat{H} \equiv \sum_{PQ} \tilde{h}_{PQ} \hat{d}_P^\dagger \hat{d}_Q + \frac{1}{2} \sum_{PQRS} \tilde{g}_{PQRS} \hat{d}_P^\dagger \hat{d}_Q^\dagger \hat{d}_S \hat{d}_R$$

Full Hamiltonian in the **Householder representation**

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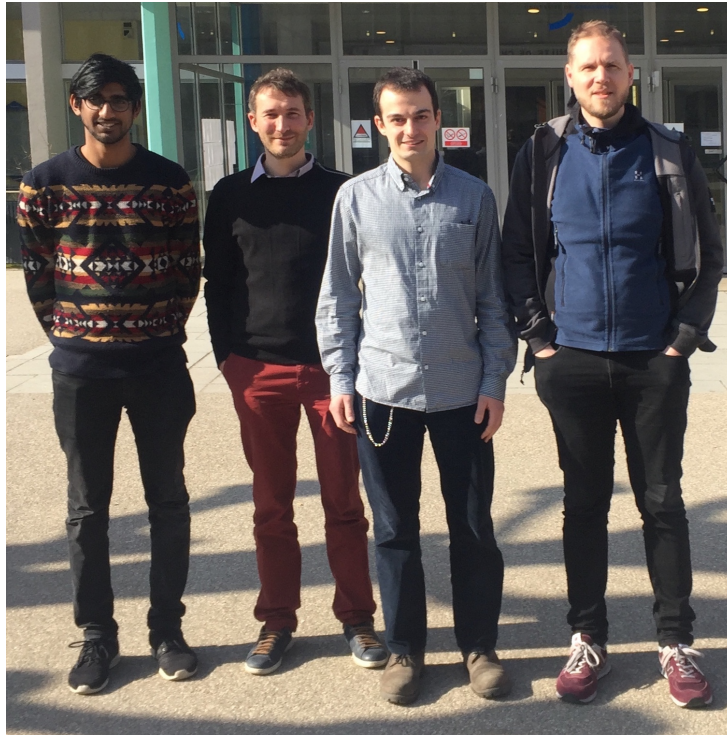


*Projection onto the cluster*

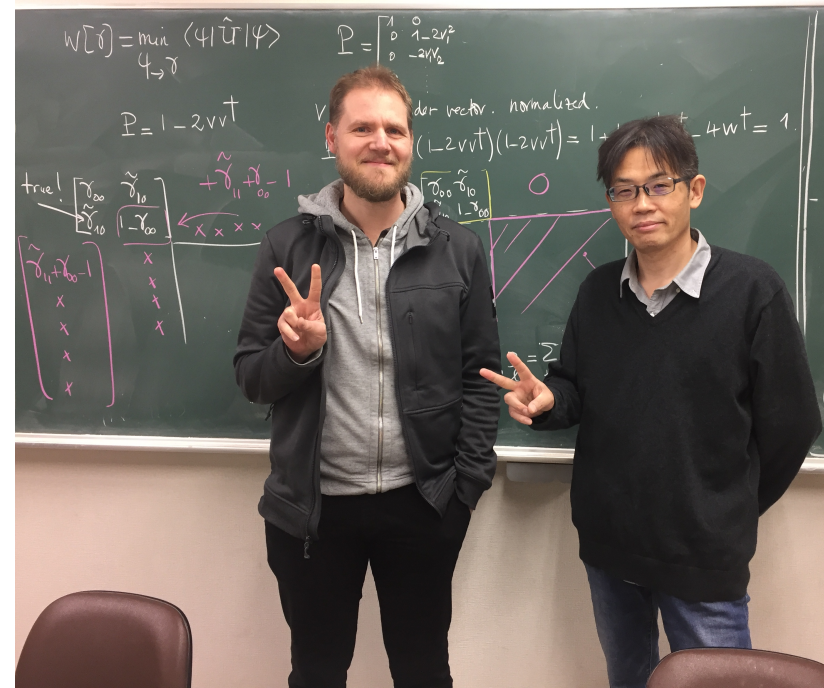
$$\hat{H}^{\mathcal{C}} \equiv \sum_{P,Q \in \mathcal{C}} \tilde{h}_{PQ} \hat{d}_P^\dagger \hat{d}_Q + \frac{1}{2} \sum_{P,Q,R,S \in \mathcal{C}} \tilde{g}_{PQRS} \hat{d}_P^\dagger \hat{d}_Q^\dagger \hat{d}_S \hat{d}_R$$

## ***Application to the 1D Hubbard model***

# The “Householder embedding” project

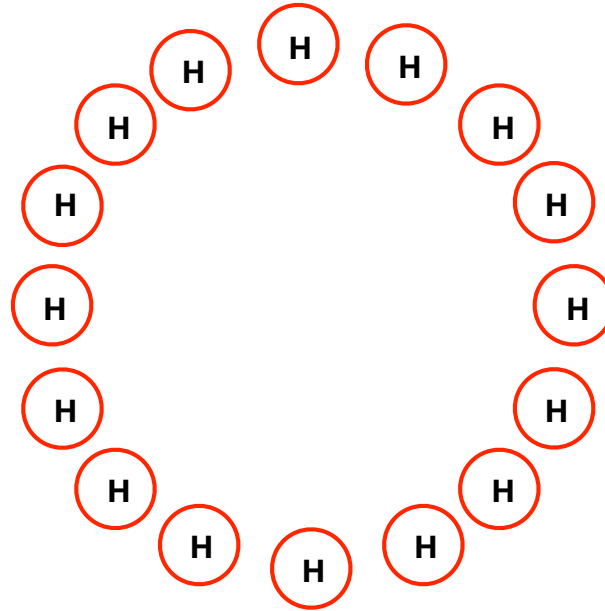


From left to right: **S. Sekaran** (Strasbourg, France),  
**M. Saubanère** (Montpellier, France),  
**L. Mazouin** (Strasbourg, France), and **E.F.**

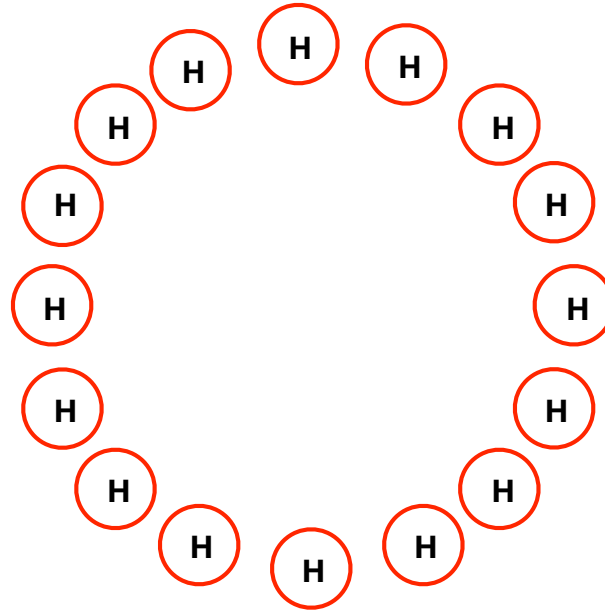


**E.F.** and **M. Tsuchiizu** (Nara, Japan).

## *Prototypical ring of $L$ hydrogen atoms*



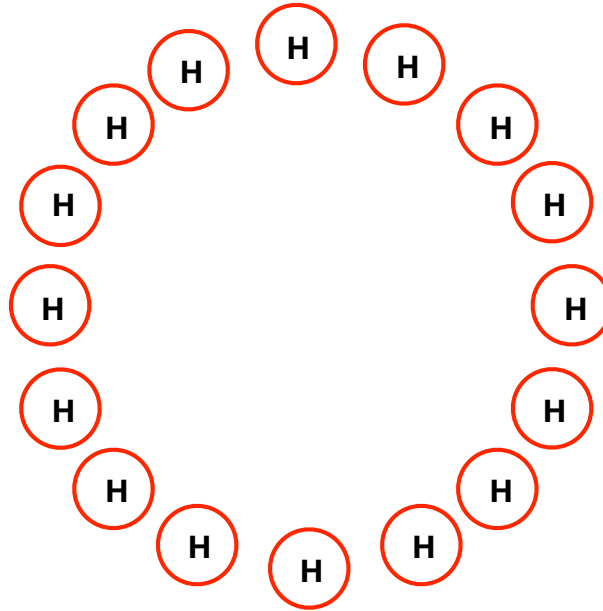
## Prototypical ring of $L$ hydrogen atoms



### Hubbard model

$$\hat{H} \approx -t \sum_{\sigma=\uparrow,\downarrow} \sum_{i=0}^{L-1} \left( \hat{c}_{i\sigma}^\dagger \hat{c}_{(i+1)\sigma} + \hat{c}_{(i+1)\sigma}^\dagger \hat{c}_{i\sigma} \right) + U \sum_{i=0}^{L-1} \hat{c}_{i\uparrow}^\dagger \hat{c}_{i\downarrow}^\dagger \hat{c}_{i\downarrow} \hat{c}_{i\uparrow}$$

## Prototypical ring of $L$ hydrogen atoms

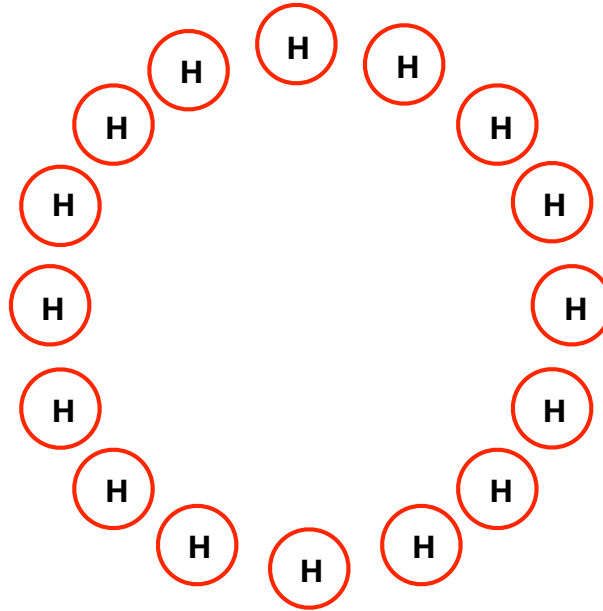


### Hubbard model

$$\hat{H} \approx -t \sum_{\sigma=\uparrow,\downarrow} \sum_{i=0}^{L-1} \left( \hat{c}_{i\sigma}^\dagger \hat{c}_{(i+1)\sigma} + \hat{c}_{(i+1)\sigma}^\dagger \hat{c}_{i\sigma} \right) + U \sum_{i=0}^{L-1} \hat{c}_{i\uparrow}^\dagger \hat{c}_{i\downarrow}^\dagger \hat{c}_{i\downarrow} \hat{c}_{i\uparrow}$$

*Hückel model*      $-t \equiv \beta$

## Prototypical ring of $L$ hydrogen atoms



### Hubbard model

$$\hat{H} \approx -t \sum_{\sigma=\uparrow,\downarrow} \sum_{i=0}^{L-1} \left( \hat{c}_{i\sigma}^\dagger \hat{c}_{(i+1)\sigma} + \hat{c}_{(i+1)\sigma}^\dagger \hat{c}_{i\sigma} \right) + U \sum_{i=0}^{L-1} \hat{c}_{i\uparrow}^\dagger \hat{c}_{i\downarrow}^\dagger \hat{c}_{i\downarrow} \hat{c}_{i\uparrow}$$

Two-electron repulsion  
on each atom only



## Prototypical ring of $L$ hydrogen atoms

$$U/t \ll 1$$

*Weakly correlated regime*

$$U/t \gg 1$$

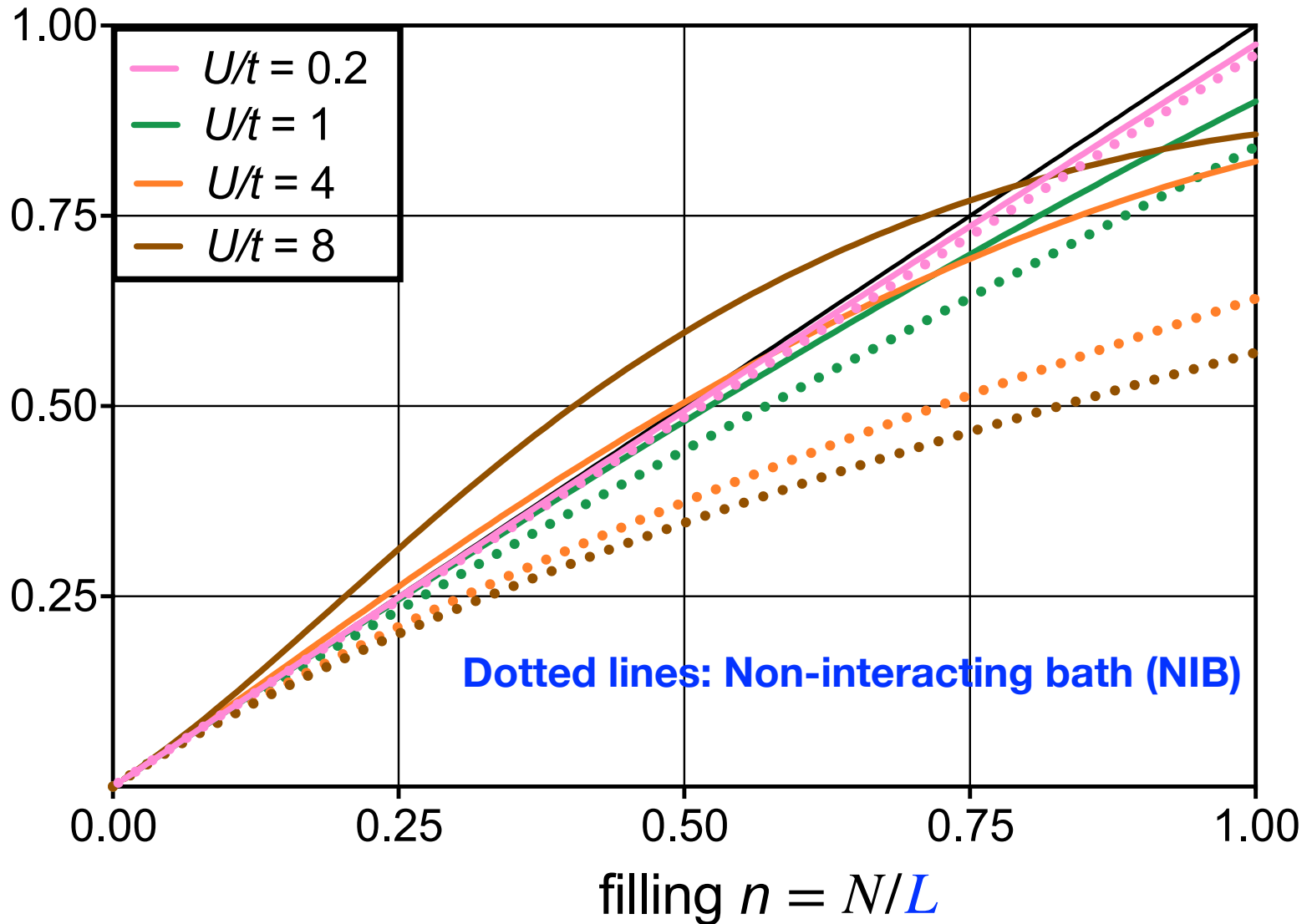
**Strongly** correlated regime

### **Hubbard model**

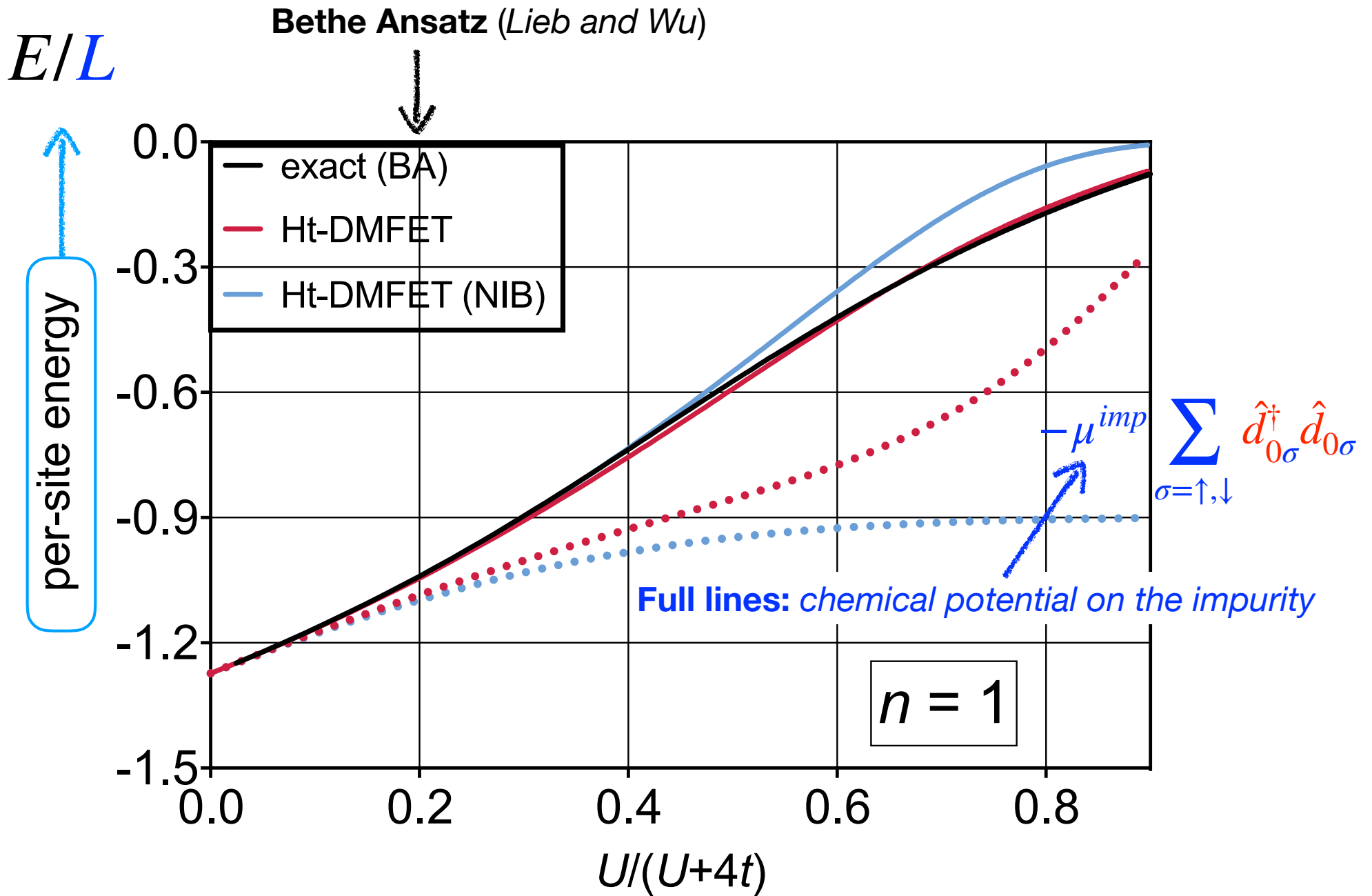
$$\hat{H} \approx -t \sum_{\sigma=\uparrow,\downarrow} \sum_{i=0}^{L-1} \left( \hat{c}_{i\sigma}^\dagger \hat{c}_{(i+1)\sigma} + \hat{c}_{(i+1)\sigma}^\dagger \hat{c}_{i\sigma} \right) + U \sum_{i=0}^{L-1} \hat{c}_{i\uparrow}^\dagger \hat{c}_{i\downarrow}^\dagger \hat{c}_{i\downarrow} \hat{c}_{i\uparrow}$$

$$\sum_{\sigma=\uparrow,\downarrow} \langle \Psi^{\mathcal{E}} | \hat{c}_{0\sigma}^\dagger \hat{c}_{0\sigma} | \Psi^{\mathcal{E}} \rangle$$

impurity site occupation

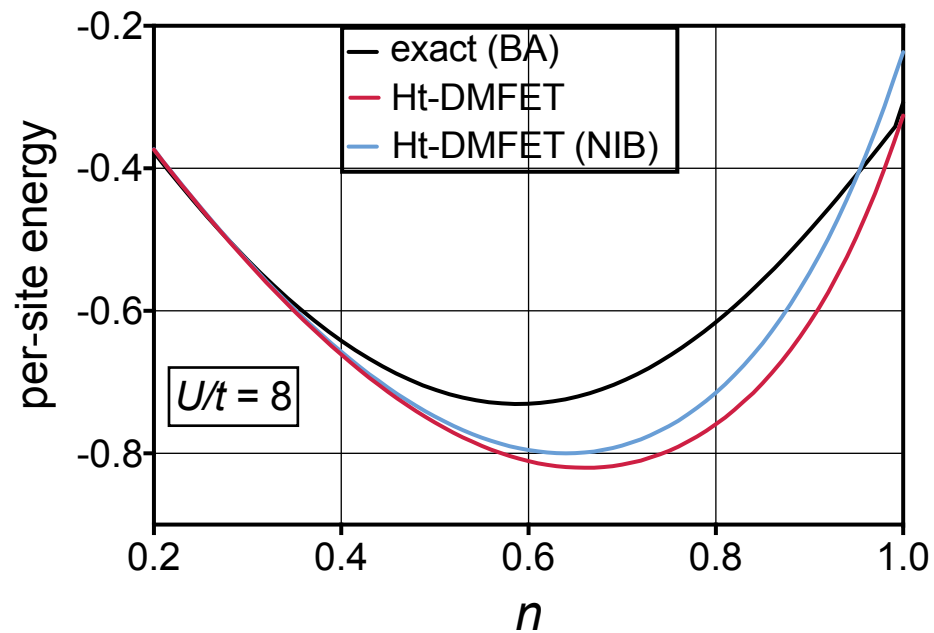
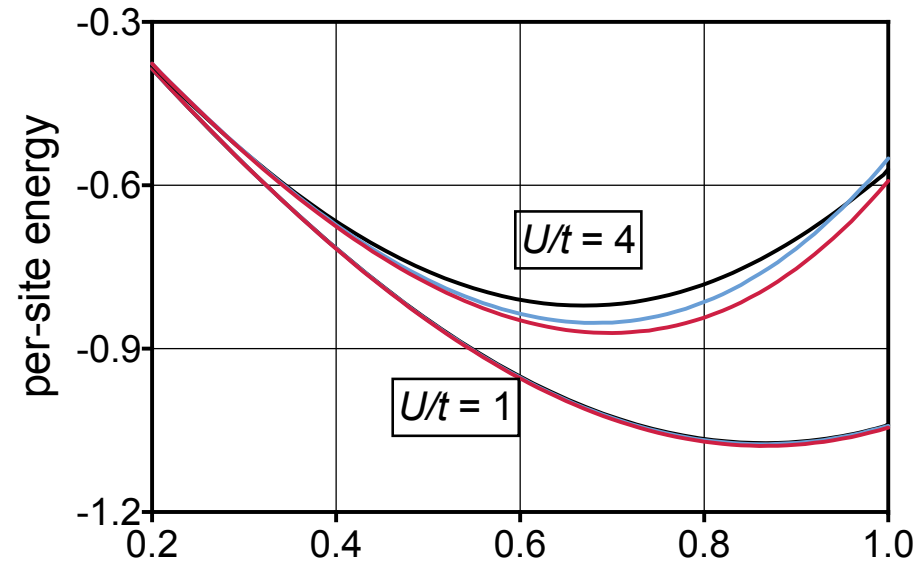


$L = 400$  atoms



$L = 400$  atoms

# Ht-DMFET per-site energies away from half-filling ( $n < 1$ )



$L = 400$  atoms