

# On Symmetry and Degeneracy in the Construction of the Adiabatic Connection Based on the Lieb Variational Principle



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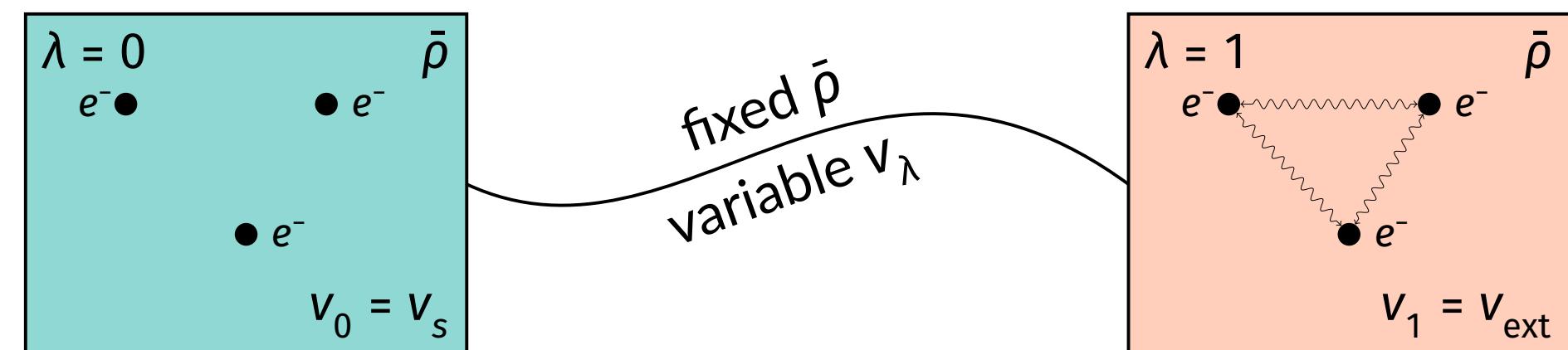


## 1. The Exact Adiabatic Connection

For an  $N$ -electron system, consider a  $\lambda$ -parametrised electronic Hamiltonian

$$\hat{\mathcal{H}}_\lambda(v_\lambda) = \hat{T} + \lambda \hat{W} + \sum_{i=1}^N v_\lambda(\mathbf{r}_i), \quad \text{where } \hat{W} = \sum_{i=1}^N \sum_{j>i}^N \frac{1}{|\mathbf{r}_i - \mathbf{r}_j|},$$

that continuously links the **physical system** to the **non-interacting system**.



The **potential**  $v_\lambda$  is chosen for  $\hat{\mathcal{H}}_\lambda$  to admit the same **ground density**  $\rho$  as  $\hat{\mathcal{H}}_1$ , via the **ground density matrix**  $\hat{\gamma}_\lambda(\rho)$ .

The **adiabatic connection** is defined by  $\mathcal{W}_\lambda(\rho) = \text{tr } \hat{\gamma}_\lambda(\rho) \hat{W}$ , which we calculate accurately for small systems to form approximate models for larger systems.

## 2. The Lieb Variational Principle

Consider the space of densities  $\mathcal{X}$  and its dual, the space of potentials  $\mathcal{X}^*$ :

$$\mathcal{X} = L^3(\mathbb{R}^3) \cap L^1(\mathbb{R}^3), \quad \mathcal{X}^* = L^{3/2}(\mathbb{R}^3) + L^\infty(\mathbb{R}^3).$$

For every  $\lambda$  along the adiabatic connection, given a **wavefunction method M** to compute the **M-ground energy**  $E_\lambda^M(v)$  of  $\hat{\mathcal{H}}_\lambda(v)$  and the M-reference density  $\bar{\rho}_1^M \in \mathcal{X}$ , by finding the **Lieb universal functional**

$$F_\lambda^M(\bar{\rho}_1^M) = \sup_{v \in \mathcal{X}^*} \left[ E_\lambda^M(v) - \int v(\mathbf{r}) \bar{\rho}_1^M(\mathbf{r}) d\mathbf{r} \right] \equiv \sup_{v \in \mathcal{X}^*} G_\lambda^M(v; \bar{\rho}_1^M),$$

one obtains the **M-optimal potential**  $v_\lambda^M$  that supports  $\bar{\rho}_1^M$  as its ground density.

The higher the quality of the wavefunction method M:

- the better  $\bar{\rho}_1^M$  approximates the true reference density  $\rho$ ;
- the lower the Lieb-variational errors  $\|\bar{\rho}_\lambda^M - \bar{\rho}_1^M\|_p$  where  $1 \leq p \leq 3$ ; and
- the better  $v_\lambda^M$  approximates the true potential  $v_\lambda$ .

## 3. Challenges of Degenerate Systems

Consider a Hamiltonian  $\hat{\mathcal{H}}_\lambda(v)$  having a **symmetry group**  $\mathcal{G}_\lambda(v)$  and admitting a **degenerate ground wavefunction**  $\tilde{\Psi}_\lambda^M(v)$  with energy  $E_\lambda^M(v)$  and density  $\bar{\rho}_\lambda^M(\mathbf{r}; v)$ .

### Well-defined gradient?

The Lieb optimisation procedure to obtain the M-optimal potential  $v_\lambda^M$  requires the functional derivative

$$\frac{\delta G_\lambda^M(v; \bar{\rho}_1^M)}{\delta v(\mathbf{r})} = \frac{\delta E_\lambda^M(v)}{\delta v(\mathbf{r})} - \bar{\rho}_1^M(\mathbf{r}).$$

As  $\bar{\rho}_\lambda^M(\mathbf{r}; v)$  is invariant only under a proper subgroup of  $\mathcal{G}_\lambda(v)$ , the functional derivative  $\delta E_\lambda^M(v)/\delta v(\mathbf{r})$  is **not unique**:

$$\frac{\delta E_\lambda^M(v)}{\delta v(\mathbf{r})} = \sum_{i=1}^{|\mathcal{G}_\lambda(v)|} c_i \hat{g}_i \bar{\rho}_\lambda^M(\mathbf{r}; v), \quad g_i \in \mathcal{G}_\lambda(v), \quad c_i \geq 0, \quad \sum_{i=1}^{|\mathcal{G}_\lambda(v)|} c_i = 1.$$

### Symmetry of potential?

At all  $\lambda$ , the M-optimal potential  $v_\lambda^M$  determines the symmetry group of the Hamiltonian:

$$\begin{aligned} \lambda = 1 : \quad v_{\text{ext}} &\mapsto \mathcal{G}_1 && \text{physical system,} \\ \lambda \neq 1 : \quad v_\lambda^M &\mapsto \mathcal{G}_\lambda(v_\lambda^M) && \text{auxiliary systems.} \end{aligned}$$

Is **equality** in the following condition

$$\mathcal{G}_1 \geq \mathcal{G}_\lambda(v_\lambda^M), \quad \lambda \neq 1$$

necessary for  $v_\lambda^M \in \mathcal{X}^*$ ? In other words, should the **auxiliary systems respect** the symmetry of the **physical system**?

### Guarantee of v-representability?

The existence of the adiabatic connection  $\mathcal{W}_\lambda(\rho)$  requires the **reference density**  $\rho$  to be **v-representable**  $\forall \lambda \in [0, 1]$ . However, general v-representability conditions for  $\rho \in \mathcal{X}$  are not well-established. In particular:

- Is  $\bar{\rho}_1^M$  pure-state v-representable  $\forall \lambda \in [0, 1]$ ?
- Is the ensemble density

$$\bar{\rho}_{1,\text{en}}^M = \frac{1}{|\mathcal{G}_1|} \sum_{i=1}^{|\mathcal{G}_1|} \hat{g}_i \bar{\rho}_1^M$$

ensemble-state v-representable  $\forall \lambda \in [0, 1]$ ?