

Interference by Nickel(II) Salts and Their 5-Methylimidazole-4-carboxylate Coordination Compounds on the Chloroplast Redox Chain

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Nickel(II) salts and their coordination compounds with ethyl 5-methylimidazole-4-carboxylate (emizco), $[\text{Ni}(\text{emizco})_2\text{Cl}_2]$, $[\text{Ni}(\text{emizco})_2\text{Br}_2]$, $[\text{Ni}(\text{emizco})_2(\text{H}_2\text{O})_2](\text{NO}_3)_2 \cdot \text{H}_2\text{O}$, $\text{Ni}(\text{NO}_3)_2$, inhibit photosynthetic electron flow (basal, phosphorylating and uncoupled) and ATP-synthesis, therefore behave as Hill reaction inhibitors. Coordination compounds are more potent inhibitors than the salts. It was found that the target for NiCl_2 ; NiBr_2 and $\text{Ni}(\text{NO}_3)_2$ is at the b_6f level. On the other hand, the complexes $[\text{Ni}(\text{Emizco})_2\text{Cl}_2]$, $[\text{Ni}(\text{Emizco})_2\text{Br}_2]$ and $[\text{Ni}(\text{emizco})_2(\text{H}_2\text{O})_2](\text{NO}_3)_2 \cdot \text{H}_2\text{O}$ binding sites are located at $Q_B(\text{D1})$ -protein and b_6f level. Therefore, they have a common inhibition site located at b_6f avoiding the PQH_2 oxidation. The Q_B inhibition site was corroborated by variable chlorophyll *a* fluorescence yield $[V(j)]$. The emizco ligand has no activity on photosynthetic electron flow.

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