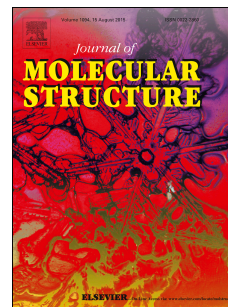


# Accepted Manuscript

New insights into the electrophilic aromatic substitution mechanism of tricyanovinylolation reaction involving tetracyanoethylene and *N,N*-dimethylaniline: An interpretation based on density functional theory calculations

Vitor H. Menezes da Silva, Natália M. Monezi, Rômulo A. Ando, Ataulpa A.C. Braga



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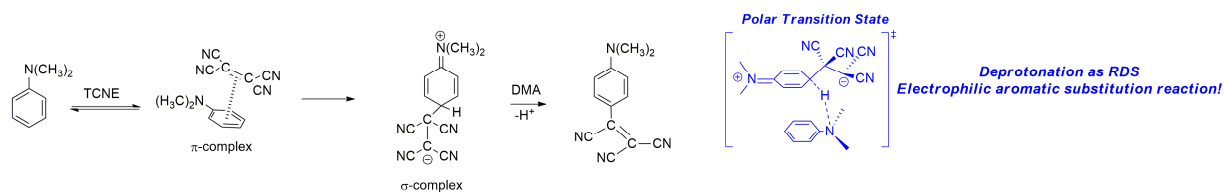
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