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Spectroscopic characterization of 1-[3-(1*H*-imidazol-1-yl)propyl]-3-phenylthiourea and assessment of reactive and optoelectronic properties employing DFT calculations and molecular dynamics simulations

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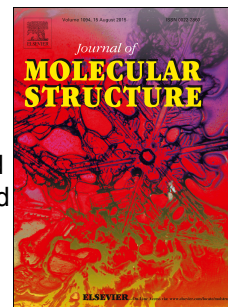
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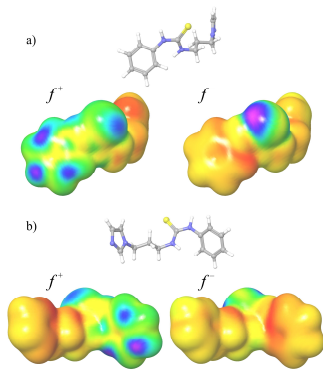
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