

Structure-reactivity relationships in bithiophenic precursors based on the 3-phenoxythiophene building block

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Résumé en anglais	3-Toluoxythiophene and bithiophenes diversely substituted by toluxo groups have been synthesized. Theoretical, spectroscopic and electrochemical studies show that the number and position of the phenoxy groups exert a strong influence on the geometry of the ground state and cation radical and determine the reactivity of the latter and hence its aptitude for electropolymerization.
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- [1] <http://okina.univ-angers.fr/philippe.leriche/publications>
- [2] <http://okina.univ-angers.fr/pierre.frere/publications>
- [3] <http://okina.univ-angers.fr/jean.roncali/publications>
- [4] <http://okina.univ-angers.fr/publications/ua10294>
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- [6] <http://xlink.rsc.org/?DOI=b502164b>