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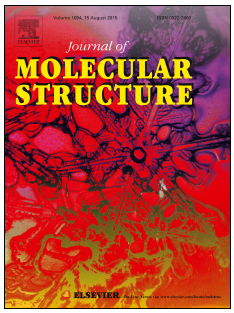

Crystallographic, spectroscopic (FTIR and NMR) and quantum computational calculation studies on bis(2-methoxy-4-((*E*)-prop-1-enyl)phenyl)oxalate

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