



Formula: C₁₆H₂₁NO₅

MW: 307.35

MDL: MFCD00582419

TNP: TNP00204



LogP: 2.44

LogS: -5.34

Acceptors: 5

Donors: 2

Rotation Bonds: 7

Chiral Centers: 1

N+O: 6

LIPINSKY: 4

IUPAC: 8-(2,3-dihydroxy-3-methylbutoxy)-4-methoxy-1-methylhydroquinolin-2-one

Smiles: c1(n(c2c(OCC(C(O)(C)C)O)ccc2c(c1)OC)C)=O