



Formula: C<sub>18</sub>H<sub>23</sub>NO<sub>6</sub>

MW: 349.38

MDL: MFCD02940837

TNP: TNP00182



LogP: -0.14

LogS: -2.81

Acceptors: 6

Donors: 2

Rotation Bonds: 8

Chiral Centers: 3

N+O: 7

LIPINSKY: 4

IUPAC: 1-(4,8-dimethoxy(4,3a-dihydrofurano[2,3-b]quinolin-7-yloxy))-3-methylbutane-2, 3-diol

Smiles: c1(c(c2c(cc1)C(C1C(=N2)OC=C1)OC)OC)OCC(C(C)(O)C)O