



Formula: C₂₁H₃₂O₅

MW: 364.48

MDL: MFCD02253413

TNP: TNP00230



LogP: -1.65

LogS: -1.88

Acceptors: 5

Donors: 1

Rotation Bonds: 4

Chiral Centers: 7

N+O: 5

LIPINSKY: 4

IUPAC: 4-hydroxy-11,11-dimethyl-5-(methylethyl)-3,6-dioxotricyclo[8.4.0.0<2,7>]tetrad ec-8-yl acetate

Smiles: C1(C2C(C(=O)C(C1C(C)C)O)C1CCCC(C1CC2OC(=O)C)(C)C)=O