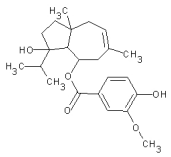

Formula: C23H32O5

MW: 388.5

MDL: MFCD02940467

TNP: TNP00014

LogP: 1.71

LogS: -3.21

Acceptors: 5

Donors: 2

Rotation Bonds: 7

Chiral Centers: 4

N+O: 5

LIPINSKY: 4

IUPAC: 10-hydroxy-4,7-dimethyl-10-(methylethyl)bicyclo[5.3.0]dec-4-en-2-yl 4-hydroxy-3-methoxybenzoate

Smiles: C1C(OC(=O)c2cc(c(cc2)O)OC)C2C(CCC2(C)CC=C1C)(C(C)C)O