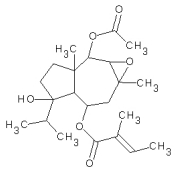

Formula: C22H34O6

MW: 394.51

MDL: MFCD02940454

TNP: TNP00012

LogP: 6.76

LogS: -5.54

Acceptors: 6

Donors: 1

Rotation Bonds: 7

Chiral Centers: 7

N+O: 6

LIPINSKY: 3

IUPAC: 2-acetyloxy-9-hydroxy-1,5-dimethyl-9-(methylethyl)-4-oxatricyclo[6.3.0.0<3,5>] undec-7-yl (2E)-2-methylbut-2-enoate

Smiles: C1C2(C(C3C(CC(OC(/C(=CC)C)=O)C2C(O)(C(C)C)C1)(C)O3)OC(=O)C)C