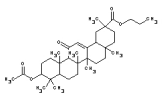

Formula: C35H54O5

MW: 554.81

MDL: MFCD01459828

TNP: TNP00019

LogP: -2.26

LogS: -1.64

Acceptors: 5

Donors: 0

Rotation Bonds: 6

Chiral Centers: 9

N+O: 5

LIPINSKY: 3

IUPAC: 1,2,6,6,10,17,20-heptamethyl-12-oxo-17-(propoxycarbonyl)pentacyclo[12.8.0.0<2, 11>.0<5,10>.0<15,20>]docos-13-en-7-yl acetate

Smiles:
C=12C(C3(CCC4C(C3C(C1)=O)(CCC(OC(=O)C)C4(C)C)C)C)(CCC1(C2CC(C(=O)OCCC)(C)CC1)C)C