



Formula: C₃₉H₆₁NO₅

MW: 623.92

MDL: MFCD01459821

TNP: TNP00020



LogP: -3.33

LogS: -2.43

Acceptors: 5

Donors: 0

Rotation Bonds: 6

Chiral Centers: 9

N+O: 6

LIPINSKY: 3

IUPAC: 1,2,6,6,10,17,20-heptamethyl-12-oxo-17-[(2-piperidylethyl)oxycarbonyl]pentacyc lo[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)OCCN2CC CCC2)(C)CC1)C)C