



Formula: C₃₀H₄₈O₄

MW: 472.71

CAS: 559-64-8

MDL: MFCD05618209

TNP: TNP00239



LogP: 6.83

LogS:

Acceptors: 4

Donors: 3

Rotation Bonds: 1

Chiral Centers: 9

N+O: 4

LIPINSKY: 3

IUPAC: (2S,5S,18S,1R,15R,21R)-18,21-dihydroxy-1,2,8,8,15,19,19-heptamethylpentacyclo[12.8.0.0.0.0]docos-11-ene-5-carboxylic acid

Smiles:

C1=2[C]([C]3(C([C]4(CC[C@@H]([C@](C4[C@@H](C3)O)(C)C)O)C)CC2C)(CC[C@@]2(C(=O)O)CC[C@](CC12)(C)C)C