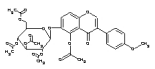

Formula: C32H32O15

MW: 656.6

MDL: MFCD02940843

TNP: TNP00186

LogP: 1.1

LogS: -3.04

Acceptors: 15

Donors: 0

Rotation Bonds: 13

Chiral Centers: 5

N+O: 15

LIPINSKY: 2

IUPAC: {3,4,5-triacetyloxy-6-[5-acetyloxy-3-(4-methoxyphenyl)-4-oxochromen-6-yloxy]-2 H-3,4,5,6-tetrahydropyran-2-yl}methyl acetate

Smiles:
COc1ccc(c2c(c3c(c(OC4OC(C(C(C4OC(C)=O)OC(=O)C)OC(C)=O)COC(C)=O)ccc3oc2)OC(C)=O)=O)cc1