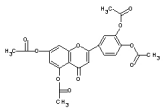

Formula: C23H18O10

MW: 454.39

MDL: MFCD01083639

TNP: TNP00171

LogP: 4.82

LogS: -4.96

Acceptors: 10

Donors: 0

Rotation Bonds: 9

Chiral Centers: 0

N+O: 10

LIPINSKY: 4

IUPAC: 7-acetyloxy-2-(3,4-diacetyloxyphenyl)-4-oxochromen-5-yl acetate

Smiles: c1(c(cc(c2oc3c(c(cc(OC(=O)C)c3)OC(=O)C)c(c2)=O)cc1)OC(=O)C)OC(=O)C