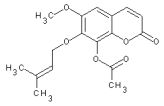

Formula: C17H18O6

MW: 318.33

MDL: MFCD01881962

TNP: TNP00185

LogP: 4.44

LogS: -4.68

Acceptors: 6

Donors: 0

Rotation Bonds: 6

Chiral Centers: 0

N+O: 6

LIPINSKY: 4

IUPAC: 6-methoxy-7-(3-methylbut-2-enyloxy)-2-oxochromen-8-yl acetate

Smiles: c1(c(c2c(cc1OC)ccc(o2)=O)OC(C)=O)OCC=C(C)C