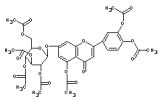

Formula: C35H34O18

MW: 742.64

MDL: MFCD02940831

TNP: TNP00244

LogP: -2.1

LogS: -2.61

Acceptors: 18

Donors: 0

Rotation Bonds: 18

Chiral Centers: 5

N+O: 18

LIPINSKY: 2

IUPAC: 2-(3,4-diacetyloxyphenyl)-4-oxo-7-[3,4,5-triacetyloxy-6-(acetyloxymethyl)(2H-3 ,4,5,6-tetrahydropyran-2-yloxy)]chromen-5-yl acetate

Smiles:
c1(c(cc(c2oc3c(c(cc(OC4C(C(C(C(COC(C)=O)O4)OC(=O)C)OC(=O)C)OC(=O)C)c3)OC(=O)C)c(c2)=O)cc1)OC(=O)C)OC(=O)C