



Formula: C₂₆H₃₆O₈

MW: 476.57

MDL:

TNP: TNP00008



LogP: 6.09

LogS: -4.92

Acceptors: 8

Donors: 1

Rotation Bonds: 9

Chiral Centers: 7

N+O: 8

LIPINSKY: 3

IUPAC: 7-(3,4-dimethoxyphenylcarbonyloxy)-9-hydroxy-1,5-dimethyl-9-(methylethyl)-4-ox
atricyclo[6.3.0.0<3,5>]undec-11-yl acetate

Smiles: C1C(OC(=O)C)C2(C)C(C(OC(c3ccc(c(c3)OC)OC)=O)CC3(OC3C2)C)C1(C(C)C)O