



Formula: C₂₆H₃₄O₈

MW: 474.55

MDL: MFCD08741676

TNP: TNP00375



LogP: 3.55

LogS: -3.77

Acceptors: 8

Donors: 6

Rotation Bonds: 9

Chiral Centers: 8

N+O: 8

LIPINSKY: 3

IUPAC: 2,3,7-trihydroxy-5-methyl-4-(methylethyl)-6-[1,2,4-trihydroxy-6-methyl-7-(methylethyl)-3-oxo(2,4,7-trihydroinden-5-yl)]-2,3,4,7-tetrahydroinden-1-one

Smiles:

C=C1(C(C2=C(C(C1C=1C(C3=C(C(C1C)C(C)C)C(C(C3=O)O)O)O)O)C(=O)C(O)C2O)C(C)C)C