



Formula: C₂₁H₂₀O₁₀

MW: 432.38

MDL: MFCD01084295

TNP: TNP00551



LogP: -2.05

LogS: -2.3

Acceptors: 10

Donors: 6

Rotation Bonds: 3

Chiral Centers: 5

N+O: 10

LIPINSKY: 3

IUPAC: 3-{4-[(2S,4S,5S,3R,6R)-3,4,5-trihydroxy-6-(hydroxymethyl)(2H-3,4,5,6-tetrahydropyran-2-yloxy)]phenyl}-5,7-dihydroxychromen-4-one

Smiles:

c1(c2c(cc(cc2O)O)occ1c1ccc(O[C@H]2[C@@H]([C@@H](O)[C@@H]([C@H](O2)CO)O)O)cc1)=O