



Formula: C₁₅H₁₀O₅

MW: 270.24

MDL: MFCD03412391

TNP: TNP00581



LogP: 1.53

LogS: -4.6

Acceptors: 5

Donors: 3

Rotation Bonds: 4

Chiral Centers: 0

N+O: 5

LIPINSKY: 4

IUPAC: 3,6-dihydroxy-2-(3-hydroxyphenyl)chromen-4-one

Smiles: c1(c(c(=O)c2c(o1)ccc(O)c2)O)c1cc(O)ccc1

Specification: NA

NA Usage And Synthesis NA