



Formula: C<sub>21</sub>H<sub>18</sub>O<sub>11</sub>

MW: 446.37

MDL: MFCD02939904

TNP: TNP00076



LogP: 1.68

LogS: -3.33

Acceptors: 11

Donors: 6

Rotation Bonds: 9

Chiral Centers: 5

N+O: 11

LIPINSKY: 2

IUPAC: 6-(5,8-dihydroxy-4-oxo-2-phenylchromen-7-yloxy)-3,4,5-trihydroxy-2H-3,4,5,6-tetrahydropyran-2-carboxylic acid

Smiles: c1ccc(cc1)c1oc2c(c(OC3C(C(C(C(O3)C(O)=O)O)O)O)cc(c2c(c1)=O)O)O