



Formula: C₂₀H₂₂O₆

MW: 358.39

MDL: MFCD03412365

TNP:

3,4,2',4',6'-PENTAMETHOXYCHALCONE



LogP: 9.31

LogS: -7.35

Acceptors: 6

Donors: 0

Rotation Bonds: 6

Chiral Centers: 0

N+O: 6

LIPINSKY: 3

IUPAC: (2E)-3-(3,4-dimethoxyphenyl)-1-(2,4,6-trimethoxyphenyl)prop-2-en-1-one

Smiles: c1(c(cc(cc1OC)OC)OC)C(/C=Cc1cc(OC)c(cc1)OC)=O

Specification: 3,4,2',4',6'-PENTAMETHOXYCHALCONE

3,4,2',4',6'-PENTAMETHOXYCHALCONE Usage And Synthesis

3,4,2',4',6'-PENTAMETHOXYCHALCONE