Formula: C 20 H 17 NO 6

MW: 367.36

MDL:

TNP: TNP00219

LogP: -2.13

LogS: -3.79

Acceptors: 6

Donors: 0

Rotation Bonds: 1

Chiral Centers: 2
$\mathrm{N}+\mathrm{O}: 7$

LIPINSKY: 4

IUPAC: (6S)-6-((5S)-6-methyl(5,6,7,8-tetrahydro-2H-1,3-dioxolano[4,5-g]isoquinolin-5-yl))-6-hydro-2H-1,3-dioxoleno[4,5-e]isobenzofuran-8-one

Smiles: c12C(O[C@@H](c1ccc1c2OCO1)([C@@H]1(c2c(cc3c(c2)OCO3)CCN1C)))=O

