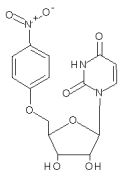

Formula: C15H15N3O8

MW: 365.3

MDL: MFCD00598743

TNP:

LogP: 0.28

LogS: -2.97

Acceptors: 8

Donors: 3

Rotation Bonds: 5

Chiral Centers: 4

N+O: 11

LIPINSKY: 4

IUPAC: 1-{3,4-dihydroxy-5-[(4-nitrophenoxy)methyl]oxolan-2-yl}-1,3-dihydropyrimidine-2,4-dione

Smiles: n1(C2OC(COc3ccc([N+](=O)[O-])cc3)C(C2O)O)c([nH]c(=O)cc1)=O