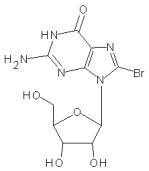

Formula: C10H12BrN5O5

MW: 362.14

MDL: MFCD00382172

TNP:

LogP: -0.57

LogS: -2.67

Acceptors: 5

Donors: 6

Rotation Bonds: 5

Chiral Centers: 4

N+O: 10

LIPINSKY: 3

IUPAC: 2-amino-9-[3,4-dihydroxy-5-(hydroxymethyl)oxolan-2-yl]-8-bromohydropurin-6-one

Smiles: n1(c2c(c([nH]c(n2)N)=O)nc1Br)C1OC(CO)C(C1O)O