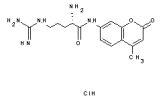

Formula: C16H22ClN5O3

MW: 367.84

MDL: MFCD04221072

TNP:

LogP: 1.21

LogS: -3.6

Acceptors: 3

Donors: 7

Rotation Bonds: 5

Chiral Centers: 1

N+O: 8

LIPINSKY: 3

IUPAC: (2S)-5-(amidinoamino)-2-amino-N-(4-methyl-2-oxochromen-7-yl)pentanamide, chloride

Smiles: Cc1c2c(cc(cc2)NC([C@H](CCCNC(=N)N)N)=O)oc(c1)=O.Cl