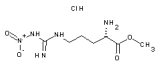

Formula: C7H16ClN5O4

MW: 269.69

MDL NUMBER: MFCD01325054

IUPAC: methyl (2S)-2-amino-5-{[imino(nitroamino)methyl]amino}pentanoate, chloride

Smiles: [N+](NC(=N)NCCCC(C(=O)OC)N)([O-])=O.Cl

ACCEPTORS: 4

DONORS: 5

ROTATION BONDS: 6

N+O: 9

Chiral Centers: 1

LogP: -1.47

LogS: -2.18

LIPINSKI: 4