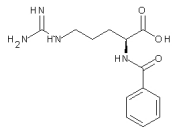

Formula: C13H18N4O3

MW: 278.31

CAS: 154-92-7

MDL: MFCD00063818

TNP:

LogP: -3.5

LogS: -1.81

Acceptors: 3

Donors: 6

Rotation Bonds: 6

Chiral Centers: 1

N+O: 7

LIPINSKY: 3

IUPAC: (2S)-5-(amidinoamino)-2-(phenylcarbonylamino)pentanoic acid

Smiles: C(N[C@H](C(=O)O)CCCNC(=N)N)(=O)c1ccccc1