

Compound	Chemical Structure (SMILES)	EC50 (μM)
1	<chem>CCN(CC)C1=NC(NC2=CC=CC=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	0.24
6a	<chem>CCN(CC)C1=NC(NC2=CC=CC(OC=O)=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	0.16
6b	<chem>CCN(CC)C1=NC(NC2=CC=CC(F)=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	0.14
6c	<chem>CCN(CC)C1=NC(NC2=CC=CC(Cl)=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	2.8
6d	<chem>CCN(CC)C1=NC(NC2=CC=CC([N+][O-])=O)=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	0.2
6e	<chem>CCN(CC)C1=NC(NC2=CC=C(C(O)=O)C=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	0.9
6f	<chem>CCN(CC)C1=NC(NC2=CC=C(F)C=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	0.7
6g	<chem>CCN(CC)C1=NC(NC2=CC=C([N+][O-])=O)C=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	18
6h	<chem>CCN(CC)C1=NC(NC2=CC=C3C(NN=C3)=C2)=NC(OC(C(F)(F)F)[C](F)F)=N1.F</chem>	0.9
6i	<chem>CCN(CC)C1=NC(NC2=CC=C3C(C=NN3)=C2)=NC(OC(C(F)(F)F)[C](F)F)=N1.F</chem>	2.8
6j	<chem>CCN(CC)C1=NC(NC2=CC=C3C(OC=N3)=C2)=NC(OC(C(F)(F)F)[C](F)F)=N1.F</chem>	0.15
6k	<chem>CCN(CC)C1=NC(NC2=CC=C3C(SC=N3)=C2)=NC(OC(C(F)(F)F)[C](F)F)=N1.F</chem>	0.05
7a	<chem>CCN(CC)C1=NC(NC2=CC=CC=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	2
7b	<chem>CCN(CC)C1=NC(NC2=CC=CC(OC=O)=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	0.5
7c	<chem>CCN(CC)C1=NC(NC2=CC=CC(F)=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	1.6
7d	<chem>CCN(CC)C1=NC(NC2=CC=CC(Cl)=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	5.4
7e	<chem>CCN(CC)C1=NC(NC2=CC=CC([N+][O-])=O)=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	1.2
7f	<chem>CCN(CC)C1=NC(NC2=CC=C(C(O)=O)C=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	2.5
7g	<chem>CCN(CC)C1=NC(NC2=CC=C(F)C=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	4.8
11	<chem>CNC1=NC(NC2=CC=CC=C2)=NC(OC(C(F)(F)F)C(F)(F)F)=N1</chem>	0.12
12	<chem>CNC1=NC(NC2=CC=CC=C2)=NC(OCC(F)(F)C(F)F)=N1</chem>	0.04