

Compound No.	SMILES	pIC50 (IC50 in microM)
1	<chem>FC1=CC(Cl)=C(CN2C(=O)C(=O)C3=CC(I)=CC=C23)C=C1</chem>	-0.973127854
2	<chem>CC1=NC(SCC(=O)NC2=CC=C(Cl)C=C2F)=NC(=C1)C(F)(F)F</chem>	-2.698970004
3	<chem>ClC1=CC(OC(=O)C2=CC3=C(OC2=O)C=CC=C3)=CN=C1</chem>	0.966576245
4	<chem>COC(=O)C1(C)CCCC2=C3C(=O)C(=O)C4=C(OC=C4C)C3=CC=C12</chem>	-1.324282455
5	<chem>FC(F)(F)C1=CC=C(NC(=O)CSC2=NC(=CC=N2)C2=CC(=NO2)C2=CC=C(Cl)C=C2Cl)C=C1</chem>	-1.176091259
6	<chem>ClC1=CC(OC(=O)C2=C3C=CC=CC3=CC=C2)=CN=C1</chem>	BLINDED
7	<chem>COC1=CC=CC(=C1)C1=C(C#N)C(=O)NC(SCC(=O)NC2=CC=C(C=C2)C(C)=O)=N1</chem>	-1.77815125
8	<chem>CC1CC(C)CN(C1)S(=O)(=O)C1=CC2=C(NC(=O)C2=O)C=C1</chem>	-0.633468456
9	<chem>ClC1=C2C(=O)C(=O)N(CC3=CC4=CC=CC=C4S3)C2=CC=C1</chem>	-1.049218023
10	<chem>CN1C(=O)C(=O)C2=C1C=CC(=C2)S(=O)(=O)N1CCOCC1</chem>	-0.996073654
11	<chem>O=C1N(CC2=CC=CC=C2)C2=C(C=C(C=C2)S(=O)(=O)N2CCOCC2)C1=O</chem>	-1.14176323
12	<chem>CC1=CC(O)=NC(SCC(=O)NC2=CC=C(OC3=CC=C(Cl)C=C3)C=C2)=N1</chem>	BLINDED
13	<chem>COC1=CC(CN2CCN(CC2)S(=O)(=O)C2=CC3=C(NC(=O)C3=O)C=C2)=CC(OC)=C1OC</chem>	-1.50623436
14	<chem>CC1CCN(CC1)S(=O)(=O)C1=CC2=C(C=C1)N(CC1=CC=CC=C1)C(=O)C2=O</chem>	-0.017033339
15	<chem>CSC1=NNC(NC(=O)C2=CC=CS2)=C1S(=O)(=O)C1=CC=C(Cl)C=C1</chem>	-1.176091259
16	<chem>ClC1=CC=C(C=C1)C1=CC=C(O1)C(=O)OC1=CN=CC(Cl)=C1</chem>	1.200659451
17	<chem>[O-][N+](=O)C1=CC(=CC=C1S(=O)(=O)C1=CC=C(Cl)C=C1)C(F)(F)F</chem>	-1.079181246
18	<chem>CCC(SC1=NC(C2=CC=CC(OC)=C2)=C(C#N)C(=O)N1)C(=O)NC1=CC=CC(=C1)C(C)=O</chem>	-1.77815125
19	<chem>NC1=NC=C(C(N)=N1)S(=O)(=O)C1=CC=C(Cl)C=C1</chem>	-0.77815125
20	<chem>CSC1=C(C(C)=C(S1)C1=NC(C)=CS1)C1=CC=NC(SCC(=O)NC2=CC=C(Cl)C=C2)=N1</chem>	-1.041392685
21	<chem>COC1=CC=C(C=C1)C1=NC(SCC(=O)NC2=CC=C(C=C2)C(C)=O)=NC=C1</chem>	-2.602059991
22	<chem>FC(F)(F)C1=CC=CC(NC(=O)CSC2=NC(C3=CC=CC=C3)=C(C#N)C(=O)N2)=C1</chem>	-2.301029996
23	<chem>CC1CC(C)CN(C1)S(=O)(=O)C1=CC=C2N(CC3=CC=C4C=CC=CC4=C3)C(=O)C(=O)C2=C1</chem>	-0.672097858
24	<chem>CC1=C(C=C(O1)C(C)(C)C1=NNC(NS(=O)(=O)C2=CC=CS2)=C1</chem>	BLINDED
25	<chem>ClC(Cl)=C(Cl)C(=O)OC1=CC=C(C=C1)S(=O)(=O)C1=CC=C(OC(=O)C(Cl)=C(Cl)Cl)C=C1</chem>	BLINDED
26	<chem>CC1=CC(=O)NC(SCC(=O)NC2=C(OC3=CC=CC=C3)C=CC(Cl)=C2)=N1</chem>	-2
27	<chem>O=C1N(CC2=CC=C3C=CC=CC3=C2)C2=CC=C(C=C2C1=O)S(=O)(=O)N1CCN(CC1)C1=CC=CC=N1</chem>	-0.741939078
28	<chem>ClC1=CC(OC(=O)C2=CC=NC=C2)=CN=C1</chem>	0.785156152
29	<chem>CC1=C(Cl)C=C(Cl)C(=C1)S(=O)(=O)C1=C(C=C(C=C1[N+](=[O-])=O)C(F)(F)F)[N+](=[O-])=O</chem>	0.522878745
30	<chem>CC1CCN(CC1)S(=O)(=O)C1=CC=C2N(CC3=CC=C4C=CC=CC4=C3)C(=O)C(=O)C2=C1</chem>	-0.227886705
31	<chem>O=C1N(CC2=CC=C3C=CC=CC3=C2)C2=CC=C(C=C2C1=O)S(=O)(=O)N1CCCCC1</chem>	BLINDED
32	<chem>CC1=C(C(C)=C(C#N)C(=N1)S(=O)(=O)C1=CC=CC=C1)[N+](=[O-])=O</chem>	-1.113943352
33	<chem>ClC1=CC(CN2CCN(CC2)S(=O)(=O)C2=CC3=C(NC(=O)C3=O)C=C2)=CC=C1</chem>	-1.501196242
34	<chem>COC1=C(OC)C=C(NC(=O)CSC2=NC(O)=CC(=N2)C2=CC=CC=C2)C=C1</chem>	-1.653212514
35	<chem>CN1CCN(CC1)S(=O)(=O)C1=CC=C2N(CC3=CC=C4C=CC=CC4=C3)C(=O)C(=O)C2=C1</chem>	-1.918606915
36	<chem>COC(OC)C1=CC(O)=NC(SCC(=O)NC2=CC=C(C=C2)C(F)(F)F)=N1</chem>	-2.301029996
37	<chem>IC1=CC=C2N(C\C=C\C3=CC4=CC=CC=C4S3)C(=O)C(=O)C2=C1</chem>	BLINDED
38	<chem>[O-][N+](=O)C1=CC=C(C=C1)S(=O)(=O)C1=CC=C(C=C1)[N+](=[O-])=O</chem>	-1.397940009
39	<chem>CC(=O)C1=CC=CC=C1S(=O)(=O)C1=CC=CC=C1C(O)=O</chem>	-1.204119983
40	<chem>CC1=NC(=CS1)C1=NC(=CS1)C1=NC(SCC(=O)NC2=CC=C(Cl)C=C2)=NC=C1</chem>	-1.146128036
41	<chem>CC1(C)CC2=C(SC(NCC3=CC=CO3)=C2C(=O)C1)C#N</chem>	-1.204119983
42	<chem>CC1=CC=C(C=C1)S(=O)(=O)C1=NC(C)=C(C(C)=C1C#N)[N+](=[O-])=O</chem>	-1.113943352
43	<chem>CC1CCCCN1S(=O)(=O)C1=CC2=C(NC(=O)C2=O)C=C1</chem>	BLINDED
44	<chem>ClC1=CC=C(S1)C1=CSC(CS(=O)(=O)C2=CC=CS2)=N1</chem>	-1.255272505
45	<chem>OC(=O)C1=CC=C(C=C1)S(=O)(=O)C1=CC(Br)=C(O)C(Br)=C1</chem>	-1.204119983

46	FC1=CC=C2N(CC3=CC4=CC=CC=C4S3)C(=O)C(=O)C2=C1	-0.683047038
47	ClC1=CC=C(NC(=O)C2=CC=C(CN3C(=O)C(=O)C4=CC(I)=CC=C34)S2)C=C1	-1.099335278
48	CN1CCN(CC1)S(=O)(=O)C1=CC2=C(C=C1)N(CC1=CC=CC=C1)C(=O)C2=O	-1.827369273
49	CCCC1=CC(Cl)=NC(SCC(=O)NC2=CC=C(Cl)C=C2)=N1	-1.477121255
50	[O-][N+](=O)C1=C(C=CC=C1)C1=CC=C(O1)C(=O)OC1=CN=CC(Cl)=C1	0.681936665
51	O=C1N(CC2=CC3=CC=CC=C3S2)C2=CC=CC=C2C1=O	-1.117602692
52	CC1=COC2=C1C(=O)C(=O)C1=C2C=CC2=C(C)C=CC=C12	-1.587710965
53	CC(C)C1=CC=C(NC(=O)CSC2=NC=CC(=N2)C2=CSC(COC3=C(Cl)C=CC=C3)=N2)C=C1	-2.301029996
54	CC1CCN(CC1)S(=O)(=O)C1=CC2=C(C=C1)N(C)C(=O)C2=O	-1.2509077
55	[O-][N+](=O)C1=C[N+](=[O-])=C(C=C1)S(=O)(=O)C1=CC=C(Cl)C=C1	-1.176091259
56	[O-][N+](=O)C1=CC=CC(=C1)C(=O)OC1=CN=CC(Cl)=C1	0.164943898
57	[O-][N+](=O)C1=C2N(CC3=CC4=CC=CC=C4S3)C(=O)C(=O)C2=CC=C1	-0.301029996
58	[O-][N+](=O)C1=C(C=CC=C1)C(=O)OC1=CN=CC(Cl)=C1	0.477555766
59	[O-][N+](=O)C1=CC=C(C=C1)C1=CC=C(O1)C(=O)OC1=CN=CC(Cl)=C1	BLINDED
60	CCOC(=O)C1=CN=C(SCC(=O)NC2=CC=C(C=C2)[N+](=[O-])=O)N=C1N	-2.397940009
61	COC1=CC(=CC=C1)C1=C(C#N)C(=O)NC(SCC(=O)NC2=CC=C(C=C2)S(N)(=O)=O)=N1	-1.602059991
62	ClC1=CC(NC(=O)CSC2=NC=CC(=N2)C2=CSC(=N2)C2=CC=CC=C2)=CC(Cl)=C1	-0.477121255
63	O=C(CC1=NCCS1)C1=NC=CS1	-1.602059991
64	CN1C(=O)C(=O)C2=C1C=CC(=C2)S(=O)(=O)N1CCN(C)CC1	-1.072984745
65	O=C1NC2=C(C=C(C=C2)S(=O)(=O)N2CCN(CC2)C2=CC=CC=N2)C1=O	-1.710371264
66	O=C1N(CC2=CC=C3C=CC=CC3=C2)C2=CC=C(C=C2C1=O)S(=O)(=O)N1CCOCC1	-1.600646236
67	CCCCC1=CC=C(NC(=O)CSC2=NC(C3CCCC3)=C(C#N)C(=O)N2)C(C)=C1	-2.301029996
68	CC1CC(C)CN(C1)S(=O)(=O)C1=CC2=C(C=C1)N(CC1=CC=CC=C1)C(=O)C2=O	-0.450249108
69	CC(C)(C)C1=CC=C(NC(=O)CSC2=NC(C3CCCC3)=C(C#N)C(=O)N2)C=C1	-2.544068044
70	CC(SC1=NC(C2=CC=CC=C2)=C(C#N)C(=O)N1)C(=O)NC1=CC=C(Cl)C=C1	BLINDED
71	FC(F)(F)C1=NN=C(N1)SC(=O)C1=CC=C(O1)C#CC1=CC=CC=C1	-0.477121255
72	O=C1NC2=C(C=C(C=C2)S(=O)(=O)N2CCOCC2)C1=O	-1.102433706
73	CC(C)C1=CC2=CC=C3C(CCCC3(C)C)=C2C(=O)C1=O	-1.324282455
74	CN1N=C(C)C(=C1NCC1=CC=C(S1)C1=CC=CS1)[N+](=[O-])=O	-1.301029996
75	IC1=CC=C2N(CC3=CC4=CC=CC=C4S3)C(=O)C(=O)C2=C1	0.022276395
76	O=C1N(CC2=CC=C3C=CC=CC3=C2)C2=CC=C(C=C2C1=O)S(=O)(=O)N1CCN(CCC2=CC=CC=C2)CC1	-1.14176323
77	[O-][N+](=O)C1=CC=CC(=C1)C1=CC=C(O1)C(=O)OC1=CN=CC(Cl)=C1	0.301029996
78	ClC1=CC(OC(=O)C2=CN=CC=C2)=CN=C1	0.156767222
79	CC1COC2=C1C(=O)C(=O)C1=C2C=CC2=C(C)C=CC=C12	-1.158362492
80	CC(C)C1=CC=C(NC(=O)CSC2=NC(=CC=N2)C2=CC=CS2)C=C1	BLINDED
81	CCOC1=CC=C(C=C1)N1C(=O)CC(SC2=NC(C)=CC(C)=N2)C1=O	-2.477121255
82	BrC1=C2N(CC3=CC4=CC=CC=C4S3)C(=O)C(=O)C2=CC=C1	0.008773924
83	CSC1=NN(C2=C(C(C)=NN2C)[N+](=[O-])=O)C(C2=CC=CS2)=C1C#N	-1.255272505
84	CC1COC2=C1C(=O)C(=O)C1=C3CCCC(C)(C)C3=CC=C21	-2.35545152
85	IC1=CC=C2N(CC3=CC=C(S3)C(=O)N3CCCC3)C(=O)C(=O)C2=C1	-1.243038049
86	CC1=C(CN2C(=O)C(=O)C3=CC(=CC=C23)C#N)C(C)=NO1	-0.857332496
87	CN1N=C(C=C1C1=CC=C(S1)C(=O)NC1=C(C(C)=NO1)[N+](=[O-])=O)C(F)(F)F	-0.698970004
88	CN1N=C(C=C1C(F)(F)F)C1=CC=C(S1)C1=CC=NC(SCC(=O)NC2=CC=C(Cl)C=C2)=N1	-1
89	CC1=COC2=C1C(=O)C(=O)C1=C3CCCC(C)(CO)C3=CC=C21	-1.394451681
90	O=C(N1CCN(CC1)S(=O)(=O)C1=CC2=C(NC(=O)C2=O)C=C1)C1=CC=CO1	-1.003029471
91	ClC1=CC=C(NC(=O)CSC2=NC=CC(=N2)C2=CC(=NO2)C2=CC=CC=C2)C=C1	-1.176091259

92	<chem>CC1CCN(CC1)S(=O)(=O)C1=CC2=C(NC(=O)C2=O)C=C1</chem>	-0.071882007
93	<chem>O=C1NC2=C(C=C(C=C2)S(=O)(=O)N2CCN(CCC3=CC=CC=C3)CC2)C1=O</chem>	-1.542949849
94	<chem>ClC1=CC=C(NC(=O)CSC2=NC=CC(=N2)C2=CC(=NO2)C2=C(Cl)C=CC=C2)C=C1</chem>	-1.176091259