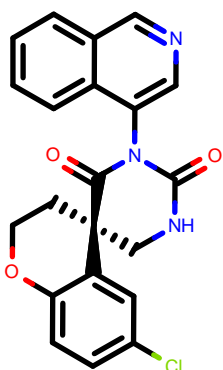
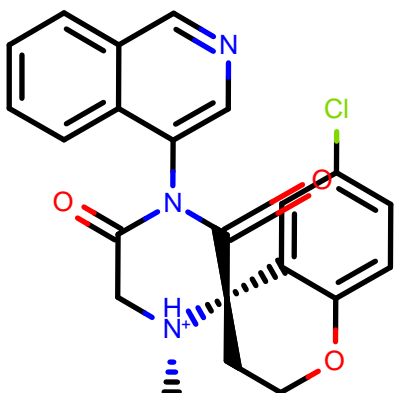


LON-WEI-4d77710c-38_1



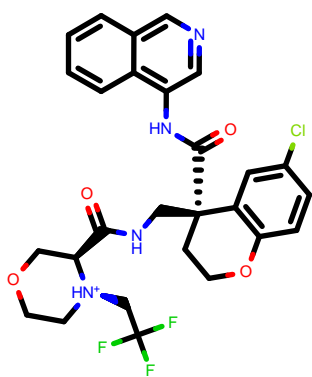
CID:	LON-WEI-4d77710c-38_1
SMILES:	<chem>CCCCN(CCCNC(=O)Nc1cn(c(=O)c2c1cccc2)CC(C)C)c3ccccc3</chem>
RUN:	RUN227
DDG (kcal/mol):	-9.94
dDDG (kcal/mol):	0.22

DAR-DIA-9e4459de-11_16



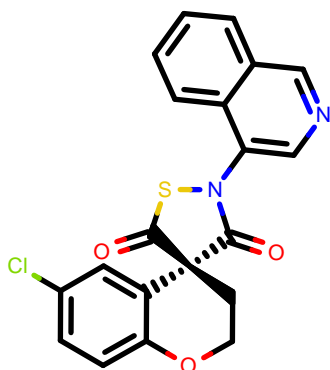
CID:	DAR-DIA-9e4459de-11_16
SMILES:	<chem>c1cc2c(c1)NCOCCOCCOCC3CC4C(C3)CNC4N(C)C1=O</chem>
RUN:	RUN1424
DDG (kcal/mol):	-9.72
dDDG (kcal/mol):	0.31

ALP-POS-347519b5-1_34



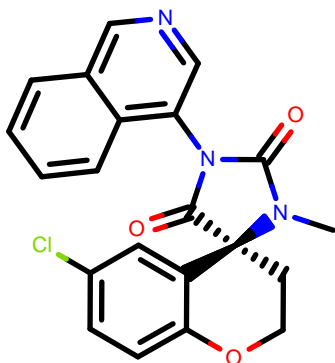
CID:	ALP-POS-347519b5-1_34
SMILES:	<chem>CSi(O)(O)N(B)1C(C@@H)2C@@H3CC(C@@H)3(C)C@@H2(C)H(C)1C(O)Nc4ccc(F)(F)F</chem>
RUN:	RUN4249
DDG (kcal/mol):	-9.49
dDDG (kcal/mol):	0.20

LON-WEI-5e7d1b3e-44_1



CID:	LON-WEI-5e7d1b3e-44_1
SMILES:	<chem>CC(C)Cn1cc(c2cccc2c1=O)NC(=O)NCc3ccccc3</chem>
RUN:	RUN1353
DDG (kcal/mol):	-6.32
dDDG (kcal/mol):	0.22

EDJ-MED-ee07cf00-15_1



CID: EDJ-MED-ee07cf00-15_1

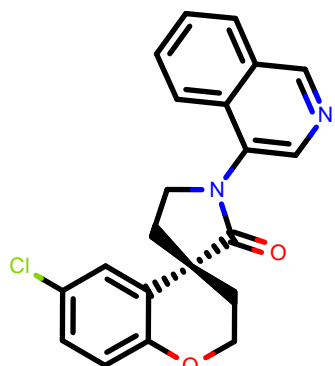
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H](c3cccc(c3)C)NC(=O)C4ccc(cc4)n5cnn5

RUN: RUN2840

DDG (kcal/mol): -6.02

dDDG (kcal/mol): 0.19

MAK-UNK-6ca90168-25_1



CID: MAK-UNK-6ca90168-25_1

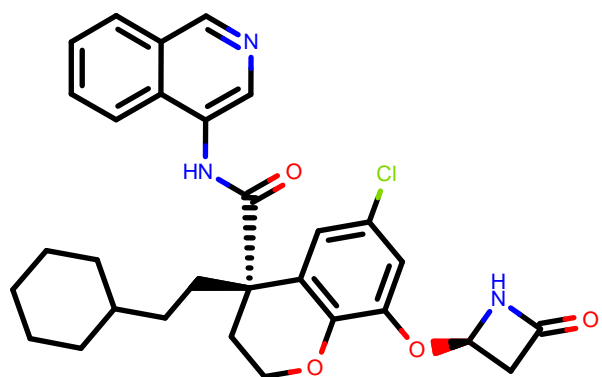
SMILES: c1ccc2c(c1)cnc2C(=O)C(=O)Nc3cccc(c3)C#N

RUN: RUN7

DDG (kcal/mol): -5.78

dDDG (kcal/mol): 0.19

ALP-POS-3b848b35-1_1



CID: ALP-POS-3b848b35-1_1

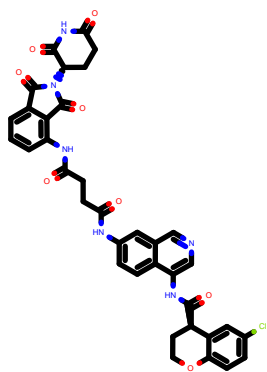
SMILES: COc1cc(cc(c1)Cl)CC(=O)Nc2cnc3c2cccc3

RUN: RUN57

DDG (kcal/mol): -5.38

dDDG (kcal/mol): 0.32

DAR-DIA-6a508060-9_2



CID: DAR-DIA-6a508060-9_2

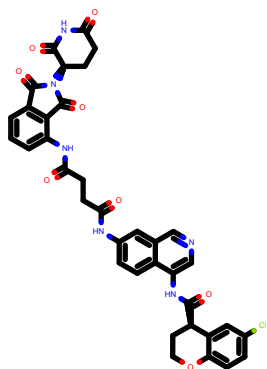
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CCC5CCCC5

RUN: RUN350

DDG (kcal/mol): -4.68

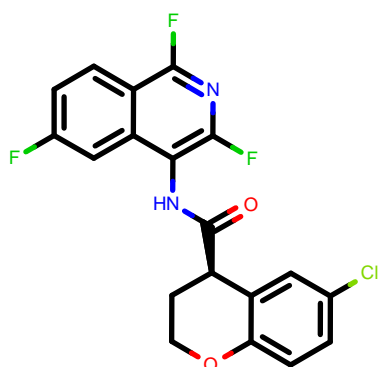
dDDG (kcal/mol): 0.40

DAR-DIA-0d514e7d-31_11



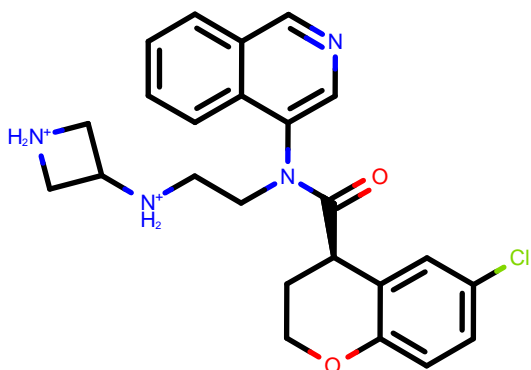
CID:	DAR-DIA-0d514e7d-31_11
SMILES:	<chem>C[C@@H]1CCO[C@H]2C=CC(=C[C@H]2[C@H]1C(=O)Nc3ncoc4c3ccoc4)Cl</chem>
RUN:	RUN844
DDG (kcal/mol):	-4.50
dDDG (kcal/mol):	0.39

VLA-UCB-05e51b3f-7_1



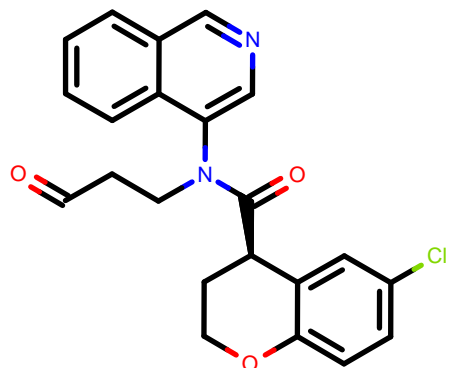
CID:	VLA-UCB-05e51b3f-7_1
SMILES:	<chem>CC(=O)N1CC[NH+](CC1)CC(=O)N(c2ncoc3c2ccoc3)C(=O)[C@H]4CCOc5c4cc(c5)Cl</chem>
RUN:	RUN322
DDG (kcal/mol):	-4.48
dDDG (kcal/mol):	0.18

MAT-POS-f7918075-2_2



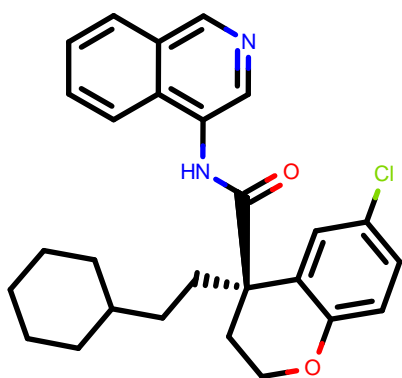
CID:	MAT-POS-f7918075-2_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3COCc4c3cc(cc4)Cl</chem>
RUN:	RUN186
DDG (kcal/mol):	-4.45
dDDG (kcal/mol):	0.47

BEN-DND-a7517465-5_1



CID:	BEN-DND-a7517465-5_1
SMILES:	<chem>c1cc2cncc(c2c(c1)F)NC(=O)[C@@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN1481
DDG (kcal/mol):	-4.43
dDDG (kcal/mol):	0.34

ALF-EVA-0b412456-5_2



CID: ALF-EVA-0b412456-5_2

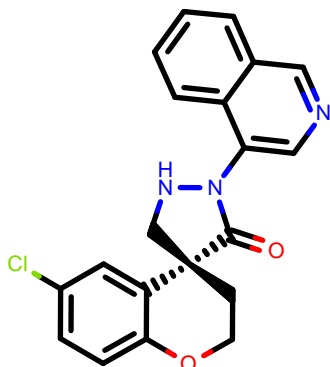
SMILES: c1cc2nccc(e2cc1C[NH3+])NC(=O)Cc3cc(cc(c3)Cl)O[C@H]4CC(=O)N4

RUN: RUN2759

DDG (kcal/mol): -4.40

dDDG (kcal/mol): 0.16

MIC-UNK-d36ab305-4_1



CID: MIC-UNK-d36ab305-4_1

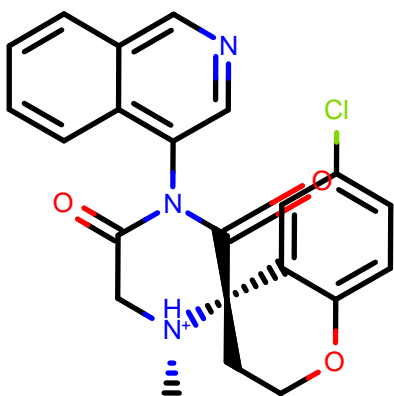
SMILES: CN(C)c1ccc(cc1)[C@@H](c2ccccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN158

DDG (kcal/mol): -4.35

dDDG (kcal/mol): 0.20

MAT-POS-4223bc15-22_3



CID: MAT-POS-4223bc15-22_3

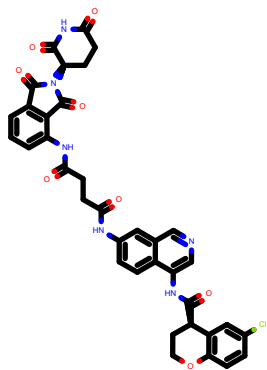
SMILES: COC(=O)C[N@@H+1]Cc2ccc(cc2[C@H](C1)C(=O)Nc3cncc4c3cccc4)Cl

RUN: RUN4103

DDG (kcal/mol): -4.31

dDDG (kcal/mol): 0.26

DAR-DIA-0f2f46c9-3_2



CID: DAR-DIA-0f2f46c9-3_2

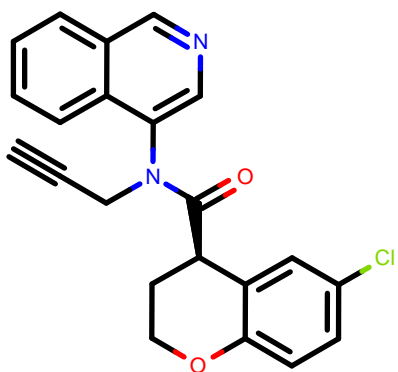
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CC[N@](c4c3cc(cc4)Cl)S(=O)(=O)N

RUN: RUN3227

DDG (kcal/mol): -4.20

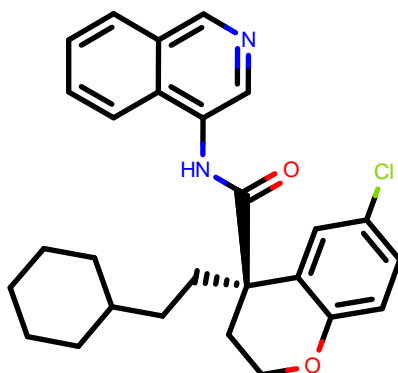
dDDG (kcal/mol): 0.20

LON-WEI-4d77710c-50_1



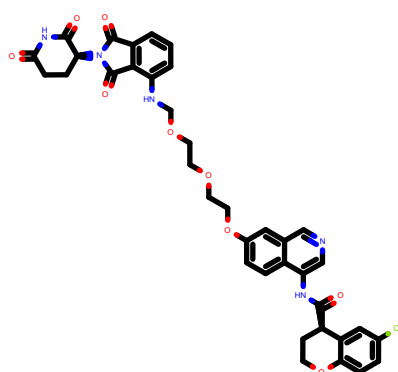
CID:	LON-WEI-4d77710c-50_1
SMILES:	<chem>CCCC[N+](=O)[O-](CCNC(=O)Nc1cn(c(=O)c2c1cccc2)C)Cc3ccccc3</chem>
RUN:	RUN242
DDG (kcal/mol):	-4.17
dDDG (kcal/mol):	0.23

EDG-MED-ba1ac7b9-31_4



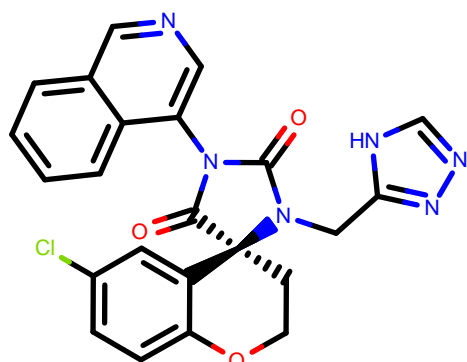
CID:	EDG-MED-ba1ac7b9-31_4
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)C[C@]3(C)CCOC4=CC=C(C=C4)CC(=O)N5CCN(H)CC5[C@@H]6CCCCO6</chem>
RUN:	RUN2743
DDG (kcal/mol):	-4.00
dDDG (kcal/mol):	0.17

MAK-UNK-ffc90da7-9_5



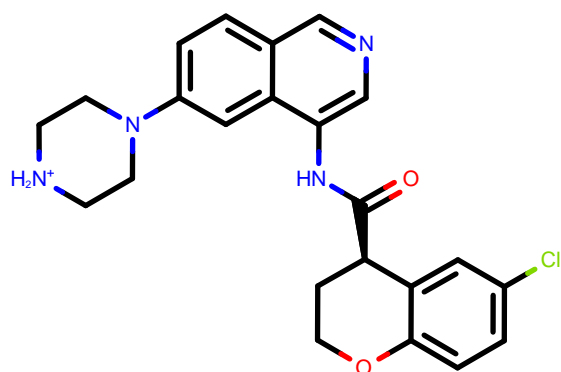
CID:	MAK-UNK-ffc90da7-9_5
SMILES:	<chem>C[C@@H]1[C@@H](CCO1)SC[C@H](C)C2CC3C(C2)CNC3NC(=O)C4CCCC(C4)Cl</chem>
RUN:	RUN716
DDG (kcal/mol):	-3.88
dDDG (kcal/mol):	0.55

DAR-DIA-0d514e7d-10_1



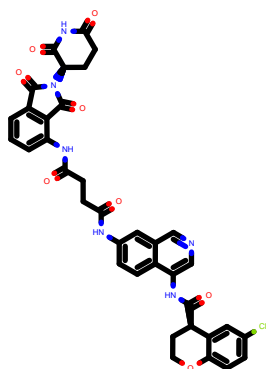
CID:	DAR-DIA-0d514e7d-10_1
SMILES:	<chem>C[C@H]1COc2c(cc(cc2Nc3cnn(c3)C)Cl)[C@@H]1C(=O)Nc4cnc5c4cccc5</chem>
RUN:	RUN814
DDG (kcal/mol):	-3.85
dDDG (kcal/mol):	0.22

MAT-POS-af71705c-2_1



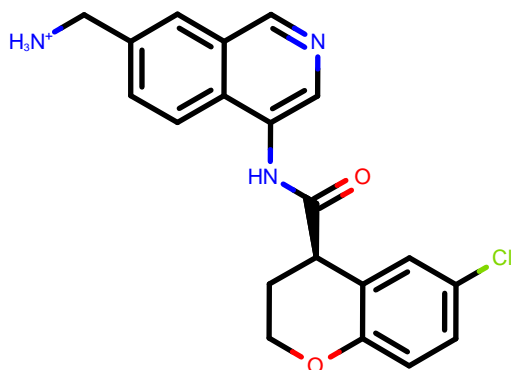
CID:	MAT-POS-af71705c-2_1
SMILES:	<chem>CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)NC3=[NH+]N(CC3)c4ccccc4</chem>
RUN:	RUN56
DDG (kcal/mol):	-3.85
dDDG (kcal/mol):	0.24

MAK-UNK-8be7dca9-8_2



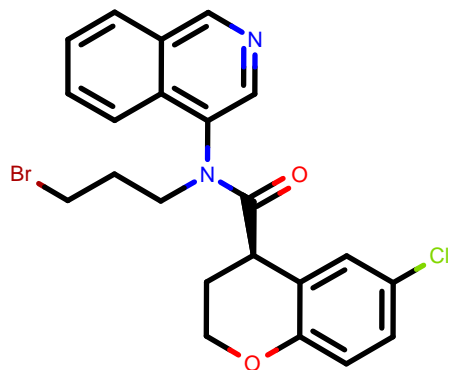
CID:	MAK-UNK-8be7dca9-8_2
SMILES:	<chem>c1cc2nccc(c2cc1CC(=O)[O-])NC(=O)[C@H]3CCOc4c3cc(cc4)Cl</chem>
RUN:	RUN509
DDG (kcal/mol):	-3.77
dDDG (kcal/mol):	0.46

MAT-POS-f7918075-2_1



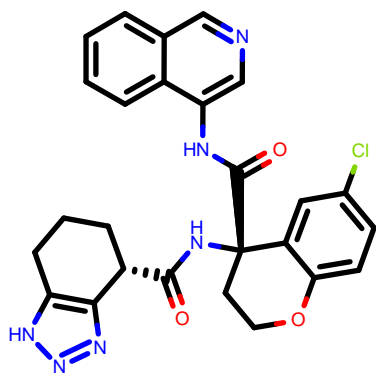
CID:	MAT-POS-f7918075-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOc4c3cc(cc4)Cl</chem>
RUN:	RUN185
DDG (kcal/mol):	-3.65
dDDG (kcal/mol):	0.24

ALP-UNI-8e43a71e-15_31



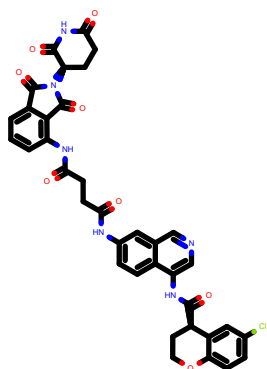
CID:	ALP-UNI-8e43a71e-15_31
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)CC(=O)NCC@H5C[C@@H]6[C@@H]7C[C@@H]8C[C@@H]9C6C(=O)N</chem>
RUN:	RUN3017
DDG (kcal/mol):	-3.44
dDDG (kcal/mol):	0.13

MAK-UNK-c749d764-33_8



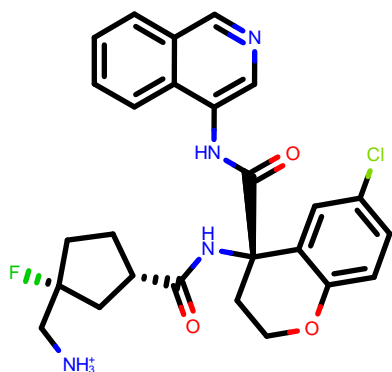
CID:	MAK-UNK-c749d764-33_8
SMILES:	<chem>CS(=O)(=O)N(c1cncc2c1cccc2)C(=O)C[C@H]3CCC[C@H][C@@H]3O)C(F)</chem>
RUN:	RUN1087
DDG (kcal/mol):	-3.41
dDDG (kcal/mol):	0.23

EDJ-MED-e4b030d8-10_1



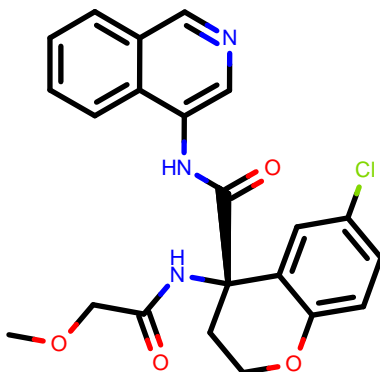
CID:	EDJ-MED-e4b030d8-10_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4c3c(c(cc4)Cl)Cl</chem>
RUN:	RUN293
DDG (kcal/mol):	-3.36
dDDG (kcal/mol):	0.45

ADA-UCB-dc2b944c-5_1



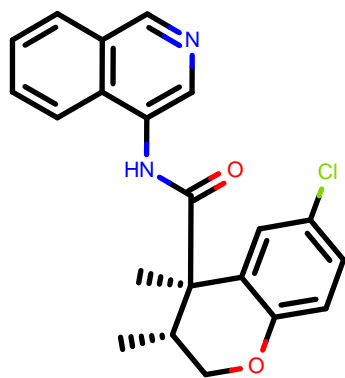
CID:	ADA-UCB-dc2b944c-5_1
SMILES:	<chem>Cc1ccc2c(c1)[C@@H](CCO2)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN601
DDG (kcal/mol):	-3.28
dDDG (kcal/mol):	0.39

MIC-UNK-bcd487e9-10_1



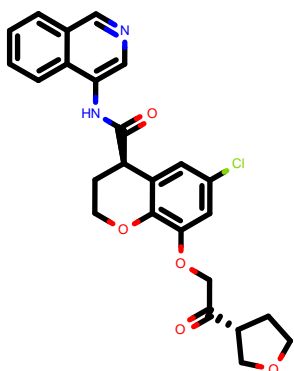
CID:	MIC-UNK-bcd487e9-10_1
SMILES:	<chem>Cn1cnc(n1)CN(c2cccc(c2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN599
DDG (kcal/mol):	-3.28
dDDG (kcal/mol):	0.26

MIC-UNK-d36ab305-6_1



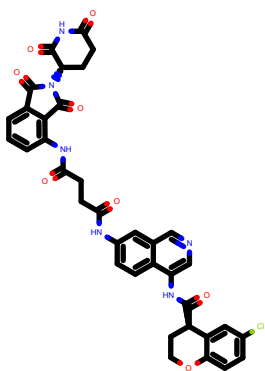
CID:	MIC-UNK-d36ab305-6_1
SMILES:	<chem>CN(C)c1ccc(cc1)N(c2cccc(c2)Cl)C(=O)Cc3cncc4c3cccc4</chem>
RUN:	RUN155
DDG (kcal/mol):	-3.26
dDDG (kcal/mol):	0.18

ALP-UNI-0676e700-18_1



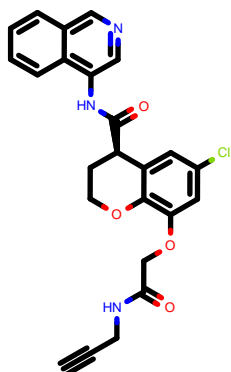
CID:	ALP-UNI-0676e700-18_1
SMILES:	<chem>Cc1nn2n1[C@@H](N)C(=O)Nc2c(Cl)C(=O)Nc3ccccc3</chem>
RUN:	RUN2462
DDG (kcal/mol):	-3.25
dDDG (kcal/mol):	0.36

VLA-UCB-50c39ae8-3_1



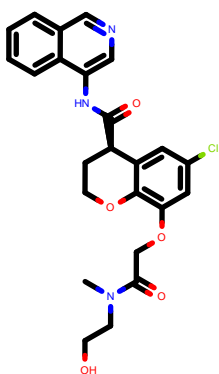
CID:	VLA-UCB-50c39ae8-3_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3C(=O)CC[C@@]4(C3=O)CCOC5c4cc(cc5)Cl</chem>
RUN:	RUN379
DDG (kcal/mol):	-3.25
dDDG (kcal/mol):	0.41

ALP-POS-2da19ca7-5_1



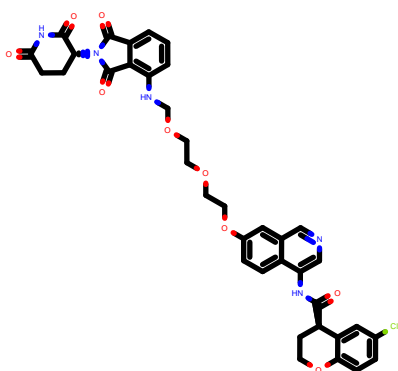
CID:	ALP-POS-2da19ca7-5_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCCC(C5)c6n(n-1)nn6</chem>
RUN:	RUN2379
DDG (kcal/mol):	-3.23
dDDG (kcal/mol):	0.37

ALP-UNI-3496895b-10_1



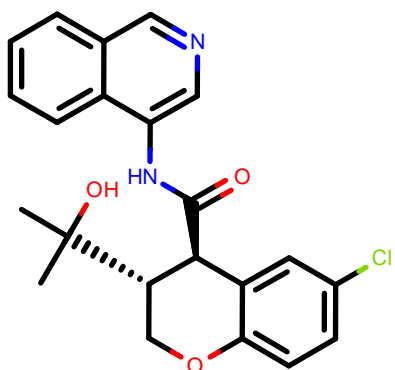
CID:	ALP-UNI-3496895b-10_1
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@@]3(C)CCOc4c3cc(cc4)C)CC(=O)N5CCN(CC5)CC(CO)(F)F</chem>
RUN:	RUN2532
DDG (kcal/mol):	-3.23
dDDG (kcal/mol):	0.38

MIC-UNK-9582b2c5-1_8



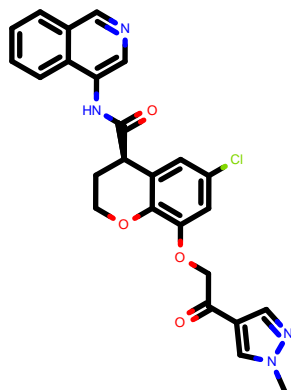
CID:	MIC-UNK-9582b2c5-1_8
SMILES:	<chem>CC(=O)N1CC[C@H]2[C@H](C1)CN(C(=O)[C@H]2c3cccc(c3)Cl)c4nccc5c4cccc5</chem>
RUN:	RUN261
DDG (kcal/mol):	-3.22
dDDG (kcal/mol):	0.53

LON-WEI-4d77710c-56_1



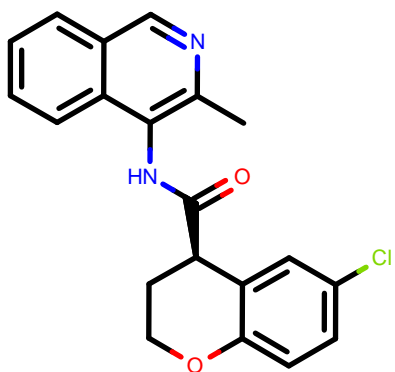
CID:	LON-WEI-4d77710c-56_1
SMILES:	<chem>CC(C)Cn1cc(c2cccc2c1=O)NC(=O)Nc3ccsc3C(=O)OC</chem>
RUN:	RUN248
DDG (kcal/mol):	-3.21
dDDG (kcal/mol):	0.17

MAT-POS-89e65850-1_1



CID:	MAT-POS-89e65850-1_1
SMILES:	<chem>CO[C@@]1(CC(=O)Nc2c1cc(cc2)Cl)C(=O)Nc3nccc4c3cccc4</chem>
RUN:	RUN2184
DDG (kcal/mol):	-3.19
dDDG (kcal/mol):	0.31

MAT-POS-b5746674-108_2



CID: MAT-POS-b5746674-108_2

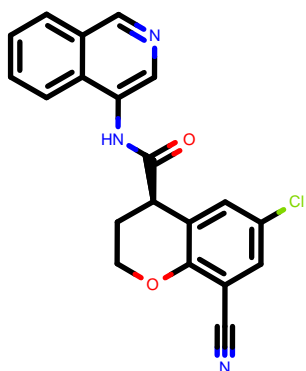
SMILES: Cc1ccc(cc1)[C@H](N)C(=O)C[C@@H](C2)CNC(=O)Nc3cn(c(=O)c4c3cccc4)CC(C)C

RUN: RUN89

DDG (kcal/mol): -3.19

dDDG (kcal/mol): 0.18

MAK-UNK-ffc90da7-4_7



CID: MAK-UNK-ffc90da7-4_7

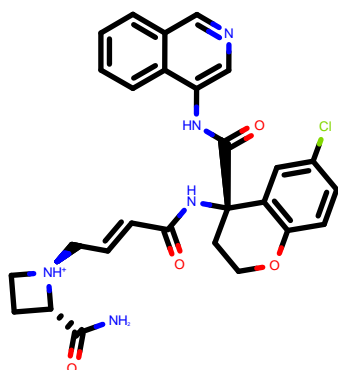
SMILES: C[C@@H](C)[C@H](N)C(=O)C[C@@H](C)CNC(=O)Nc3cn(c(=O)c4c3cccc4)C#N

RUN: RUN704

DDG (kcal/mol): -3.18

dDDG (kcal/mol): 0.19

MAK-UNK-83e0a0b4-4_2



CID: MAK-UNK-83e0a0b4-4_2

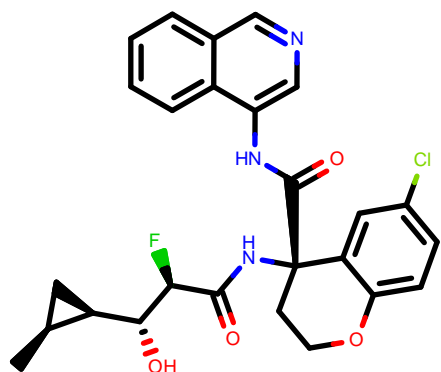
SMILES: CCCC1ccc2c(c1)[C@H](C)C(=O)N(C)CNC3C(N)C(=O)C4c3cccc4

RUN: RUN734

DDG (kcal/mol): -3.18

dDDG (kcal/mol): 0.42

ADA-UCB-dc2b944c-11_1



CID: ADA-UCB-dc2b944c-11_1

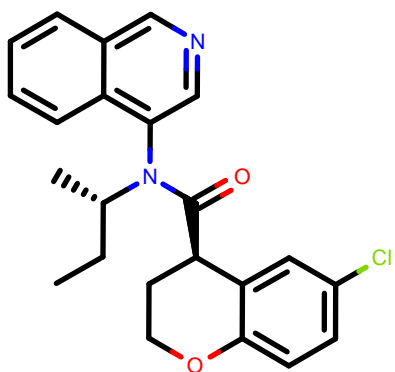
SMILES: c1ccc2c(c1)c(cnc2Br)NC(=O)[C@@H](C)CNC(=O)Nc3cn(c(=O)c4c3cc(cc4)Cl

RUN: RUN610

DDG (kcal/mol): -3.16

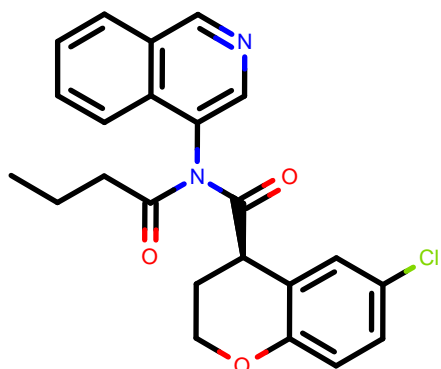
dDDG (kcal/mol): 0.30

DAR-DIA-ecdbc7dd-11_1



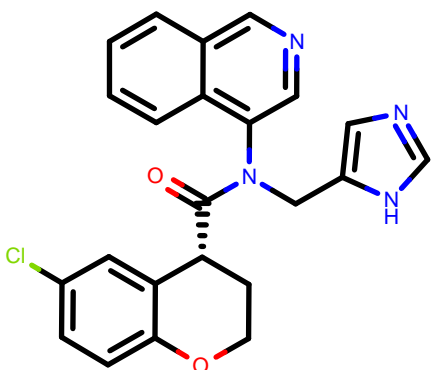
CID:	DAR-DIA-ecdbc7dd-11_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)C@@@3(CCOc4c3cc(cc4)Cl)C[NH+]5CCCCC5</chem>
RUN:	RUN2897
DDG (kcal/mol):	-3.15
dDDG (kcal/mol):	0.11

LON-WEI-4d77710c-28_1



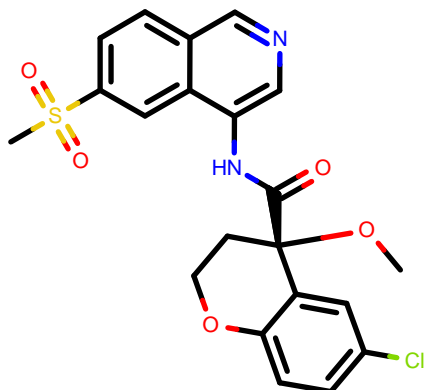
CID:	LON-WEI-4d77710c-28_1
SMILES:	<chem>CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)NCc3cc(ccc3OC)OC</chem>
RUN:	RUN218
DDG (kcal/mol):	-3.15
dDDG (kcal/mol):	0.28

EDG-MED-ba1ac7b9-25_4



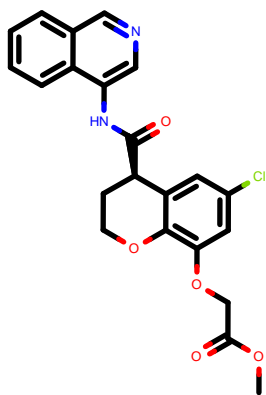
CID:	EDG-MED-ba1ac7b9-25_4
SMILES:	<chem>C[C@H]1CN(CCN@H+1CCO)C(=O)C[C@@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4ncc5c4cccc5</chem>
RUN:	RUN2714
DDG (kcal/mol):	-3.13
dDDG (kcal/mol):	0.14

ALP-POS-fab80cf2-1_1



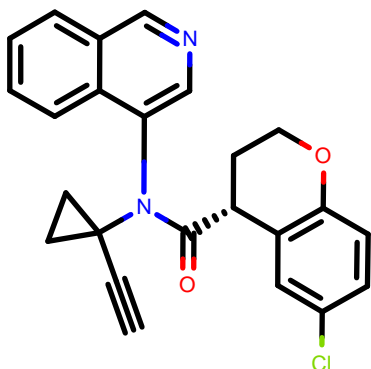
CID:	ALP-POS-fab80cf2-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)C([C@@H]4CC(=O)N4)(F)F</chem>
RUN:	RUN1514
DDG (kcal/mol):	-3.13
dDDG (kcal/mol):	0.21

MAT-POS-f9802937-9_1



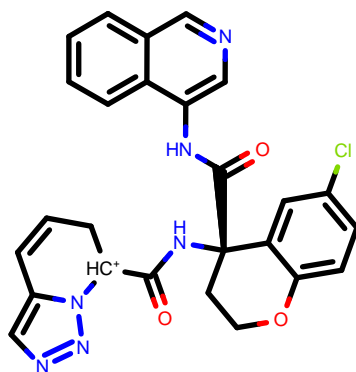
CID:	MAT-POS-f9802937-9_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCNc4c3cc(c(c4)Cl)Cl</chem>
RUN:	RUN2403
DDG (kcal/mol):	-3.11
dDDG (kcal/mol):	0.31

LON-WEI-4d77710c-54_1



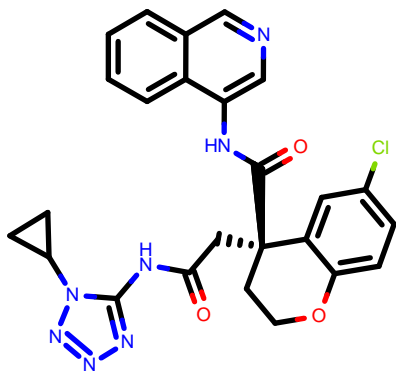
CID:	LON-WEI-4d77710c-54_1
SMILES:	<chem>CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)N3CCC(CC3)Cc4ccccc4</chem>
RUN:	RUN244
DDG (kcal/mol):	-3.09
dDDG (kcal/mol):	0.24

JOH-SUS-a69c159d-6_2



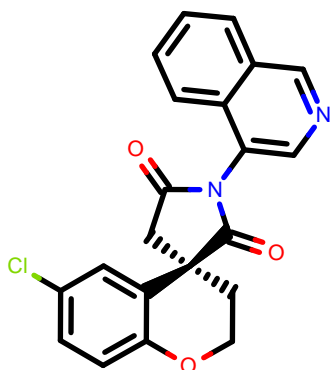
CID:	JOH-SUS-a69c159d-6_2
SMILES:	<chem>c1ccc2c(c1)c(c(nc2F)F)NC(=O)[C@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN1127
DDG (kcal/mol):	-3.07
dDDG (kcal/mol):	0.23

MAK-UNK-c749d764-33_4



CID:	MAK-UNK-c749d764-33_4
SMILES:	<chem>CS(=O)(=O)N(c1cncc2c1ccccc2)C(=O)C[C@H]3CCCC@H(C@H)3O)C(F)F</chem>
RUN:	RUN1082
DDG (kcal/mol):	-3.07
dDDG (kcal/mol):	0.33

LON-WEI-adc59df6-52_2



CID: LON-WEI-adc59df6-52_2

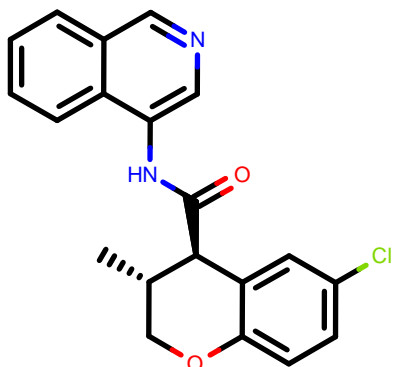
SMILES: CC(C)(C)c1ccc(cc1)N([C@H](c2ccc3c2ccc3)C(=O)NC(C)(C)C(=O)C=C

RUN: RUN3

DDG (kcal/mol): -3.06

dDDG (kcal/mol): 0.25

MIC-UNK-d36ab305-5_1



CID: MIC-UNK-d36ab305-5_1

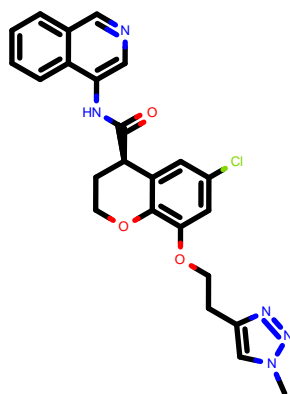
SMILES: CN(C)c1ccc(cc1)N(c2cccc(c2)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN165

DDG (kcal/mol): -3.02

dDDG (kcal/mol): 0.17

MAT-POS-e9e99895-13_2



CID: MAT-POS-e9e99895-13_2

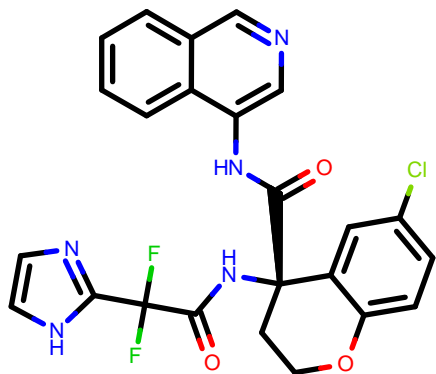
SMILES: C[C@]([c1ccc(c(c1)Cl)Cl)C(=O)Nc2cnc3c2ccc3)NC(=O)[C@@H]4CCN@@H]([C]4)C

RUN: RUN2268

DDG (kcal/mol): -3.00

dDDG (kcal/mol): 0.35

MIC-UNK-50cce87d-11_2



CID: MIC-UNK-50cce87d-11_2

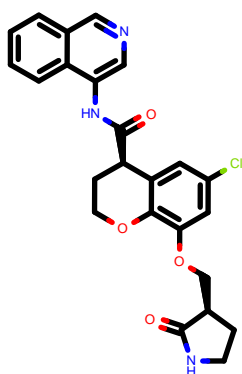
SMILES: Cc1cccc2c1c(cnc2)N3CCC[C@H](C3=O)c4cccc(c4)Cl

RUN: RUN678

DDG (kcal/mol): -2.99

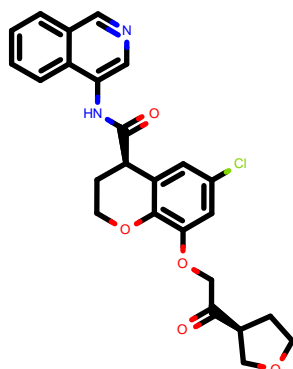
dDDG (kcal/mol): 0.25

EDG-MED-5d232de5-8_1



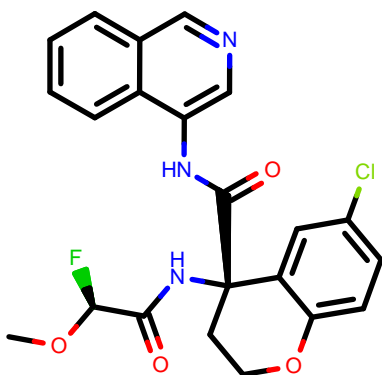
CID:	EDG-MED-5d232de5-8_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3CC[C@H](C3=O)c4cccc(c4)Cl</chem>
RUN:	RUN2370
DDG (kcal/mol):	-2.98
dDDG (kcal/mol):	0.42

EDG-MED-ba1ac7b9-5_1



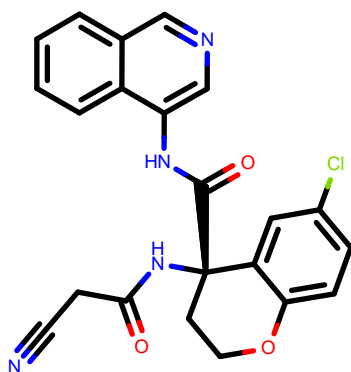
CID:	EDG-MED-ba1ac7b9-5_1
SMILES:	<chem>CC[C@H](CO)[NH+]1CCN(CC1)Cl=O[C]C@ @2(COCc3c2cc(e3)Cl)C(=O)Nc4nc5c4cccc5</chem>
RUN:	RUN2634
DDG (kcal/mol):	-2.98
dDDG (kcal/mol):	0.48

MAK-UNK-8be7dca9-3_2



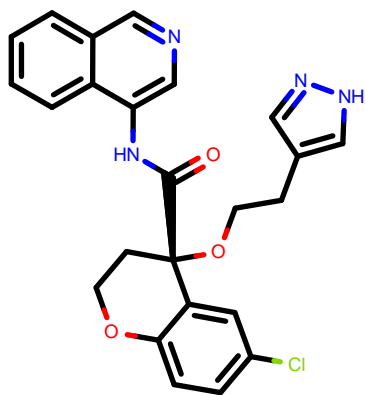
CID:	MAK-UNK-8be7dca9-3_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3c4cc(ccc4OC[C@ @H]3C[NH3+])Cl</chem>
RUN:	RUN496
DDG (kcal/mol):	-2.97
dDDG (kcal/mol):	0.22

MIC-UNK-cdc2493e-8_6



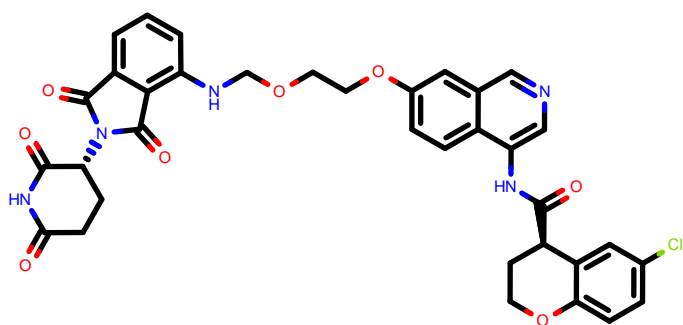
CID:	MIC-UNK-cdc2493e-8_6
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)N(c3cccc(c3)Cl)C@H4CC[C@ @H]5CCCC[C@H]5C4</chem>
RUN:	RUN541
DDG (kcal/mol):	-2.96
dDDG (kcal/mol):	0.23

MAT-POS-3b92565d-3_2



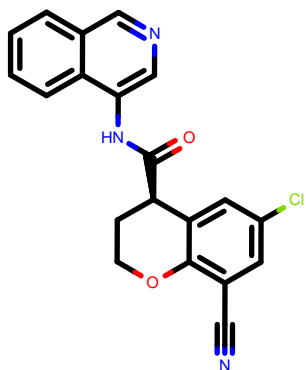
CID:	MAT-POS-3b92565d-3_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)O[C@H]4CC(=O)N4</chem>
RUN:	RUN112
DDG (kcal/mol):	-2.91
dDDG (kcal/mol):	0.32

RAL-THA-1d44ff04-5_1



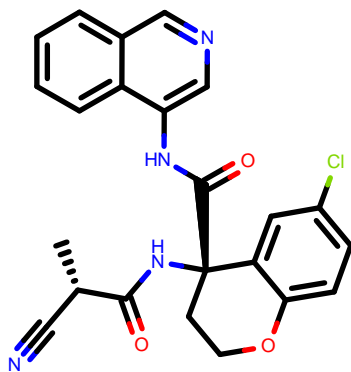
CID:	RAL-THA-1d44ff04-5_1
SMILES:	<chem>CNC(=O)c1cc(cc(c1)Cl)CC(=O)Nc2cncc3c2cccc3</chem>
RUN:	RUN440
DDG (kcal/mol):	-2.90
dDDG (kcal/mol):	0.45

LON-WEI-5e7d1b3e-56_1



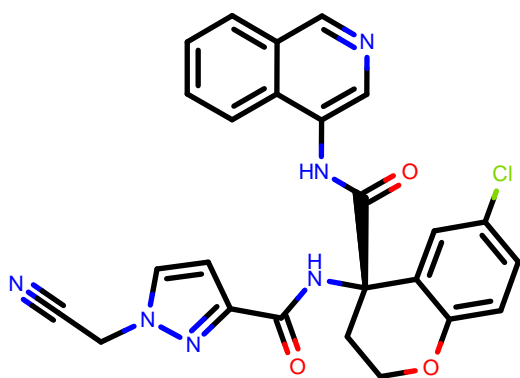
CID:	LON-WEI-5e7d1b3e-56_1
SMILES:	<chem>CC(C)Cn1cc(c2cccc2c1=O)NC(=O)Nc3ccsc3C(=O)OC</chem>
RUN:	RUN1367
DDG (kcal/mol):	-2.88
dDDG (kcal/mol):	0.22

ADA-UCB-dc2b944c-14_1



CID:	ADA-UCB-dc2b944c-14_1
SMILES:	<chem>C[C@H]1COc2ccc(cc2[C@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN612
DDG (kcal/mol):	-2.87
dDDG (kcal/mol):	0.24

MAK-UNK-c749d764-31_6



CID: MAK-UNK-c749d764-31_6

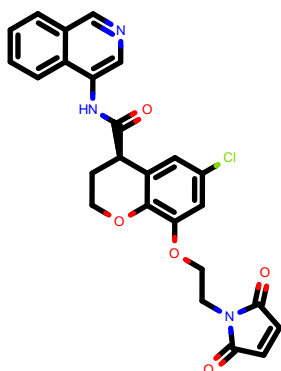
SMILES: CC/C=C/(c1cncc2c1cccc2)C(=O)C[C@H]3CCCC[C@@H]([C@H]3O)C(F)F

RUN: RUN1076

DDG (kcal/mol): -2.83

dDDG (kcal/mol): 0.23

ALP-UNI-0676e700-24_2



CID: ALP-UNI-0676e700-24_2

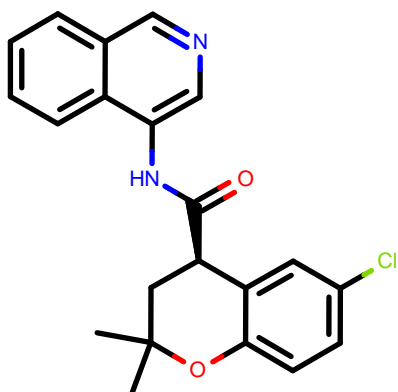
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3[C@@H]4C=CC(=O)C(C)CNC(=O)[C@H]3[C@@H]4O)N(C)C(F)F

RUN: RUN2473

DDG (kcal/mol): -2.83

dDDG (kcal/mol): 0.37

MAT-POS-2492181e-9_1



CID: MAT-POS-2492181e-9_1

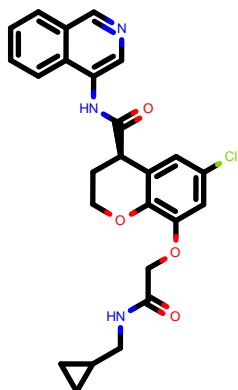
SMILES: CC1CC[NH+](CC1)C2ccc(cc2)NC(=O)Nc3cn(c(=O)c4c3cccc4)CC(C)C

RUN: RUN102

DDG (kcal/mol): -2.82

dDDG (kcal/mol): 0.22

ALP-UNI-0676e700-11_1



CID: ALP-UNI-0676e700-11_1

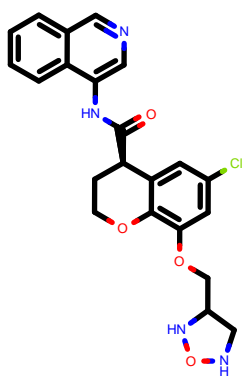
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3[C@@H]4C=CC(=O)C(C)CNC(=O)[C@H]3[C@@H]4O)N(C)C(F)F

RUN: RUN2460

DDG (kcal/mol): -2.82

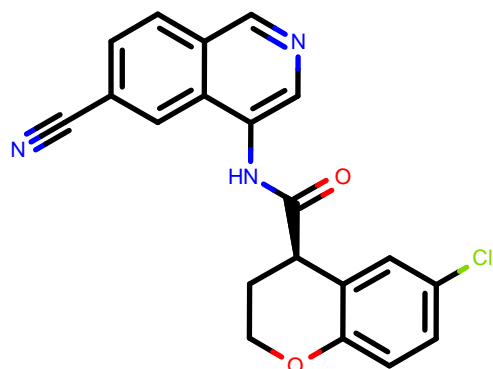
dDDG (kcal/mol): 0.33

ALP-POS-e0fe77e5-7_1



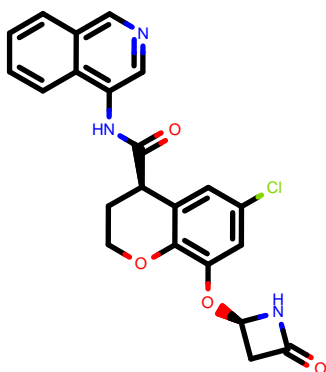
CID:	ALP-POS-e0fe77e5-7_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3CCC[C@]4(C3=O)CCOc5c4cc(cc5)Cl</chem>
RUN:	RUN2325
DDG (kcal/mol):	-2.82
dDDG (kcal/mol):	0.34

ALP-POS-f13221e1-3_1



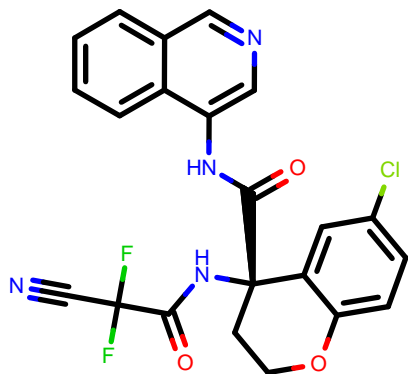
CID:	ALP-POS-f13221e1-3_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)CCc3cccc(c3)Cl</chem>
RUN:	RUN184
DDG (kcal/mol):	-2.81
dDDG (kcal/mol):	0.19

EDG-MED-ba1ac7b9-28_3



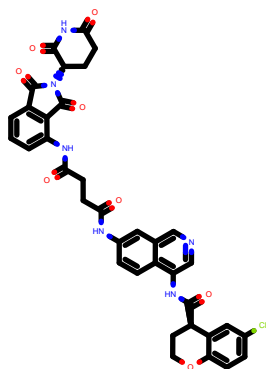
CID:	EDG-MED-ba1ac7b9-28_3
SMILES:	<chem>C[N+]([H])(CCO)C1CCN(CC1)C(=O)C[C@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4nc5c4cccc5</chem>
RUN:	RUN2734
DDG (kcal/mol):	-2.80
dDDG (kcal/mol):	0.17

MIC-UNK-cdc2493e-21_4



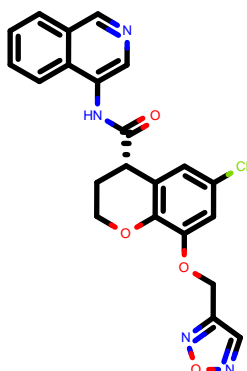
CID:	MIC-UNK-cdc2493e-21_4
SMILES:	<chem>CC(=O)N(C)[C@H]1CC[C@H](C1)N(c2cccc(c2)Cl)C(=O)Nc3nc4c3cccc4</chem>
RUN:	RUN572
DDG (kcal/mol):	-2.78
dDDG (kcal/mol):	0.24

ALP-UNI-8e43a71e-5_14



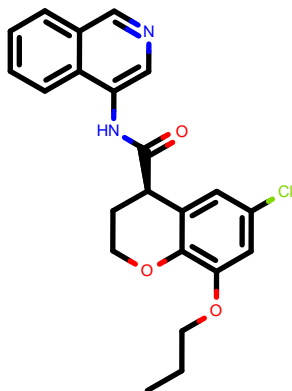
CID:	ALP-UNI-8e43a71e-5_14
SMILES:	<chem>C1C=CN(C1)C(=O)C(C(=O)N2C=CC=C2)C(=O)N3C=CC=C3</chem>
RUN:	RUN2958
DDG (kcal/mol):	-2.78
dDDG (kcal/mol):	0.24

ALP-POS-e0fe77e5-4_1



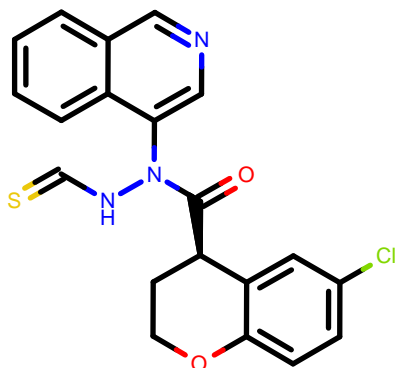
CID:	ALP-POS-e0fe77e5-4_1
SMILES:	<chem>C1=CC=C(C=C1)C(=O)N2C=CC=C2C(=O)N3C=CC=C3</chem>
RUN:	RUN2324
DDG (kcal/mol):	-2.78
dDDG (kcal/mol):	0.24

DAR-DIA-9e4459de-15_1



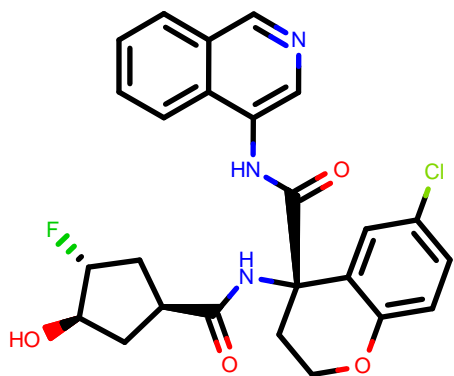
CID:	DAR-DIA-9e4459de-15_1
SMILES:	<chem>C1=CC=C(C=C1)C(=O)N2C=CC=C2C(=O)N3C=CC=C3</chem>
RUN:	RUN1441
DDG (kcal/mol):	-2.77
dDDG (kcal/mol):	0.33

LON-WEI-4d77710c-29_1



CID:	LON-WEI-4d77710c-29_1
SMILES:	<chem>CCOC(=O)C1=CC=C(C=C1)C(=O)N2C=CC=C2</chem>
RUN:	RUN221
DDG (kcal/mol):	-2.76
dDDG (kcal/mol):	0.23

MAK-UNK-ffc90da7-9_1



CID: MAK-UNK-ffc90da7-9_1

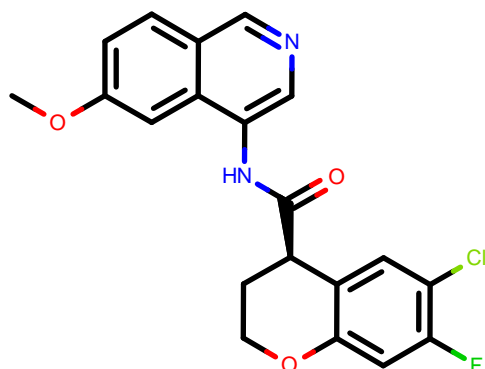
SMILES: C[C@@H]1[C@@H](CCO1)SC[C@@H](C)c2ccc3c(c2)nc3NC(=O)C4CCCC(C4)Cl

RUN: RUN709

DDG (kcal/mol): -2.74

dDDG (kcal/mol): 0.25

LAU-MED-88a3970a-14_1



CID: LAU-MED-88a3970a-14_1

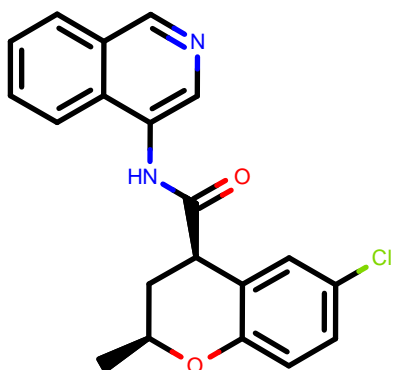
SMILES: COCCc1cc(cc2c1OCC[C@@H]2C(=O)Nc3cnc4c3cccc4)Cl

RUN: RUN1510

DDG (kcal/mol): -2.74

dDDG (kcal/mol): 0.21

DAR-DIA-0cde14eb-50_1



CID: DAR-DIA-0cde14eb-50_1

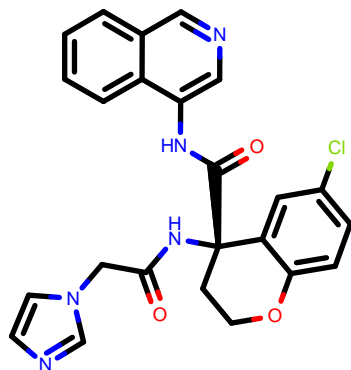
SMILES: CC1(CC1)c2cccc(c2)NC(=O)Nc3cnc4c3cccc4

RUN: RUN15

DDG (kcal/mol): -2.73

dDDG (kcal/mol): 0.21

MAK-UNK-919546f0-4_1



CID: MAK-UNK-919546f0-4_1

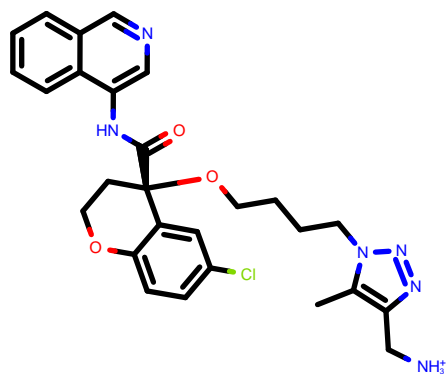
SMILES: c1ccc2c(c1)nc2NC(=O)Cc3cccc3O

RUN: RUN484

DDG (kcal/mol): -2.71

dDDG (kcal/mol): 0.31

JAG-UCB-706446eb-1_1



CID: JAG-UCB-706446eb-1_1

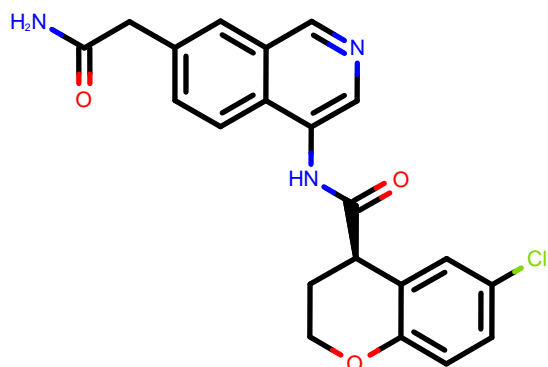
SMILES: c1ccc2c(c1)cnc2NC(=O)C@@3(CCOc4c3cc(cc4)C)CCC(=O)NCC(F)(F)F

RUN: RUN621

DDG (kcal/mol): -2.71

dDDG (kcal/mol): 0.39

ERI-UCB-ce40166b-6_2



CID: ERI-UCB-ce40166b-6_2

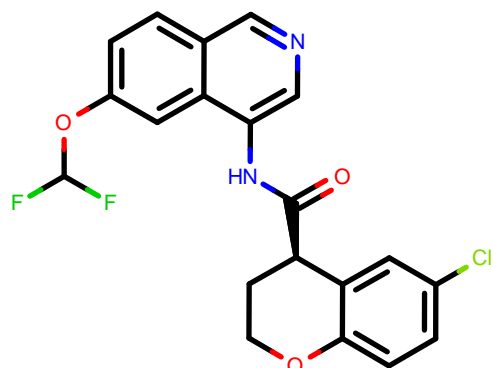
SMILES: c1ccc2c(c1)cnc2CC(=O)Nc3cc(cc(c3)C)O[C@H]4CCC(=O)N4

RUN: RUN43

DDG (kcal/mol): -2.71

dDDG (kcal/mol): 0.20

LON-WEI-4d77710c-7_1



CID: LON-WEI-4d77710c-7_1

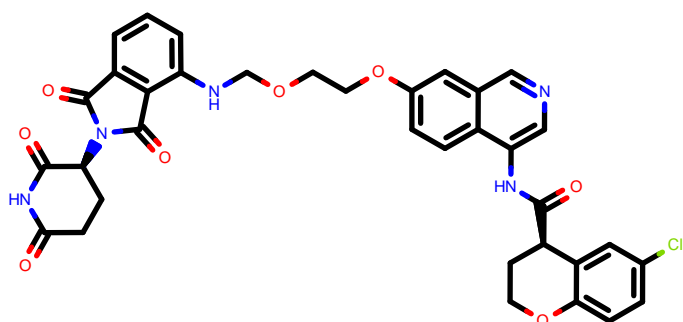
SMILES: Cn1cc(c2ccccc2c1=O)NC(=O)NCCC[N@@H+]3CCc4ccccc4C3

RUN: RUN197

DDG (kcal/mol): -2.71

dDDG (kcal/mol): 0.24

MIC-UNK-9582b2c5-2_1



CID: MIC-UNK-9582b2c5-2_1

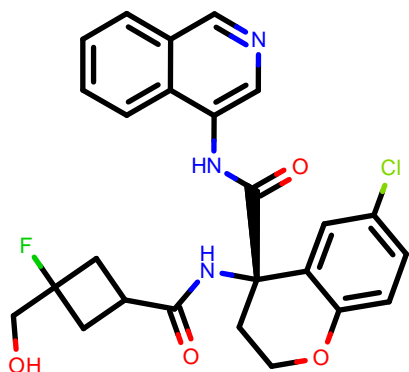
SMILES: CC(=O)N1CC[C@@H]2[C@@H](C1)C[C@@H](C(=O)N2c3cccc(c3)C)C4cnc5c4cccc5

RUN: RUN262

DDG (kcal/mol): -2.70

dDDG (kcal/mol): 0.55

EDJ-MED-28ec730d-3_1



CID: EDJ-MED-28ec730d-3_1

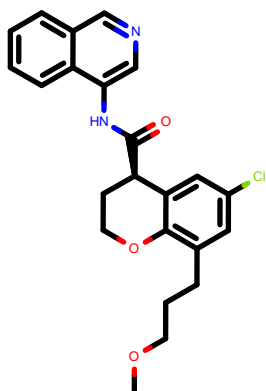
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OC5CCOCC5

RUN: RUN652

DDG (kcal/mol): -2.69

dDDG (kcal/mol): 0.26

VLA-UCB-05e51b3f-1_1



CID: VLA-UCB-05e51b3f-1_1

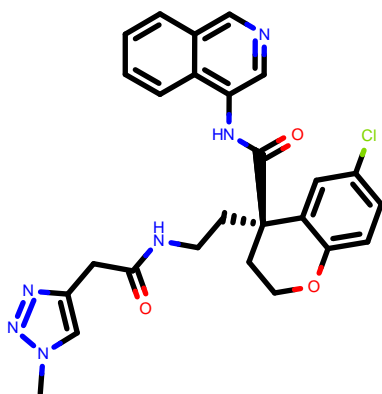
SMILES: CC(=O)N1CC[NH+](CC1)CC(=O)Nc2cnc3c2ccccc3

RUN: RUN315

DDG (kcal/mol): -2.69

dDDG (kcal/mol): 0.27

RAL-THA-e002e396-1_1



CID: RAL-THA-e002e396-1_1

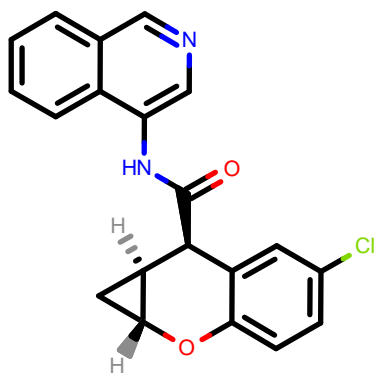
SMILES: CS(=O)(=O)C[C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cnc4c3ccccc4

RUN: RUN3451

DDG (kcal/mol): -2.69

dDDG (kcal/mol): 0.19

NIR-WEI-f9286bb6-4_2



CID: NIR-WEI-f9286bb6-4_2

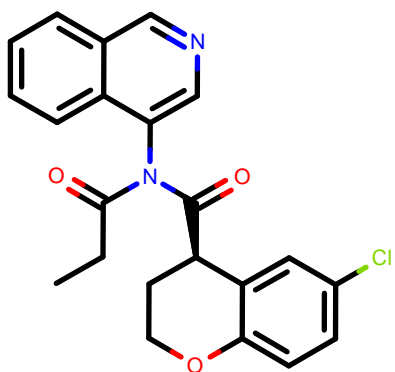
SMILES: CC(C)(C)NC(=O)[C@H](c1cnc2c1cccc2)N(Cc3ccccc3Cl)C(=O)C=C

RUN: RUN142

DDG (kcal/mol): -2.65

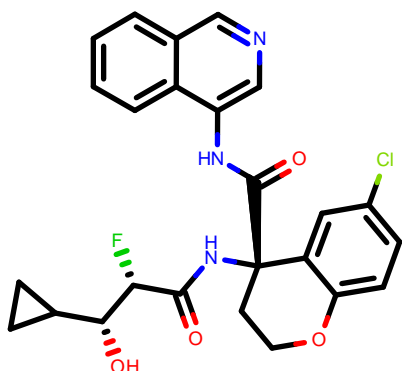
dDDG (kcal/mol): 0.14

LON-WEI-4d77710c-47_1



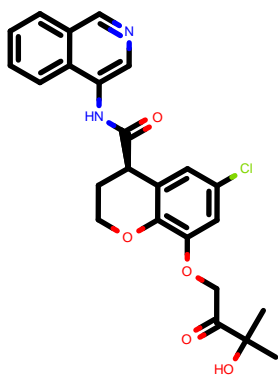
CID:	LON-WEI-4d77710c-47_1
SMILES:	<chem>CCc1cccc1N(CC)C(=O)Nc2cn(c(=O)c3c2cccc3)C</chem>
RUN:	RUN236
DDG (kcal/mol):	-2.64
dDDG (kcal/mol):	0.29

MIC-UNK-cdc2493e-21_3



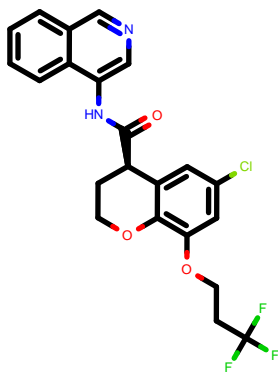
CID:	MIC-UNK-cdc2493e-21_3
SMILES:	<chem>CC(=O)N(C)[C@@H]1CC[C@H](C1)N(c2cccc(c2)Cl)C(=O)Nc3cccc4c3cccc4</chem>
RUN:	RUN571
DDG (kcal/mol):	-2.64
dDDG (kcal/mol):	0.27

ALP-POS-5bb456a5-2_6



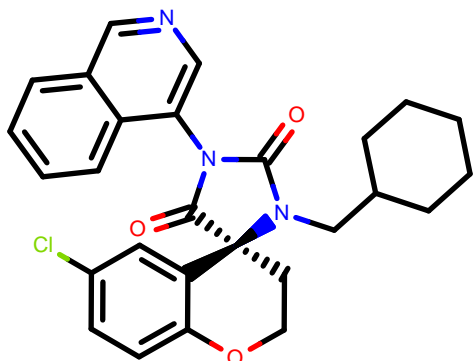
CID:	ALP-POS-5bb456a5-2_6
SMILES:	<chem>CC(C)(C)C(O)COc1ccc(Cl)cc1C(=O)Nc2cn(c(=O)c3c2cccc3)C</chem>
RUN:	RUN2427
DDG (kcal/mol):	-2.63
dDDG (kcal/mol):	0.42

ED_-GRI-5b13fbe2-35_1



CID:	ED_-GRI-5b13fbe2-35_1
SMILES:	<chem>C1CCC2C(C1)CNC(=O)[C@@]3(C)C(OC(=O)C(F)(F)F)OCCN5C(C(=O)N)C6=CC(NH2)=CC=C6N</chem>
RUN:	RUN1566
DDG (kcal/mol):	-2.62
dDDG (kcal/mol):	0.35

DAR-DIA-0d514e7d-32_13



CID: DAR-DIA-0d514e7d-32_13

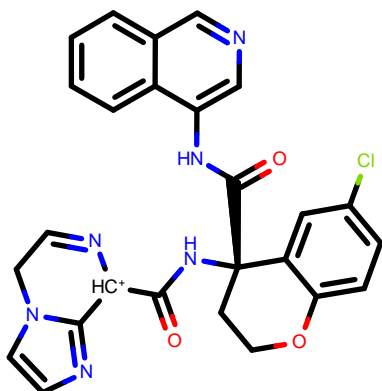
SMILES: c1ccc2c(c1)cncc2NC(=O)C@@H]3[C@@H]4[C@H]4CO[C@H]5[C@@H]3C=C(C=C5)Cl

RUN: RUN862

DDG (kcal/mol): -2.62

dDDG (kcal/mol): 0.23

MAK-UNK-c749d764-31_8



CID: MAK-UNK-c749d764-31_8

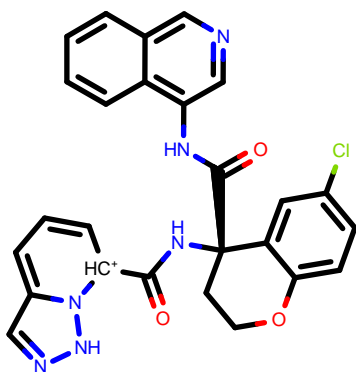
SMILES: CC/C=C(/c1cncc2c1cccc2)C(=O)C[C@H]3CCC[C@H]([C@H]3O)C(F)F

RUN: RUN1077

DDG (kcal/mol): -2.58

dDDG (kcal/mol): 0.27

EDJ-MED-841e0cf0-4_2



CID: EDJ-MED-841e0cf0-4_2

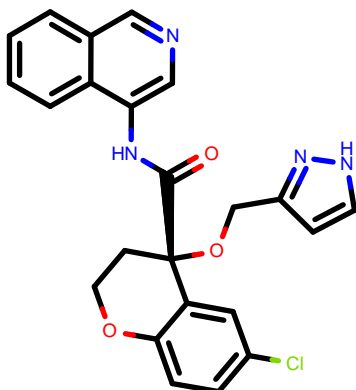
SMILES: C[C@@]1(CN@]1)Cc2c1cc(cc2)Cl)S(=O)(=O)C(=O)Nc3cncc4c3cc(cc4)S(=O)(=O)C

RUN: RUN3834

DDG (kcal/mol): -2.57

dDDG (kcal/mol): 0.17

RAL-THA-1d44ff04-2_1



CID: RAL-THA-1d44ff04-2_1

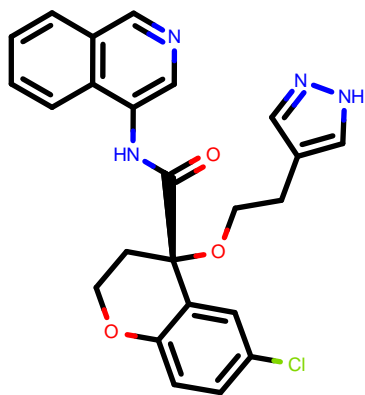
SMILES: c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)OCc4[nH]n4

RUN: RUN436

DDG (kcal/mol): -2.57

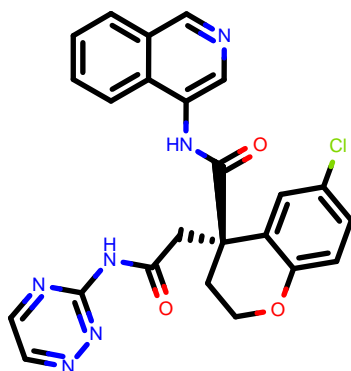
dDDG (kcal/mol): 0.22

DAR-DIA-f6ee7aeb-4_1



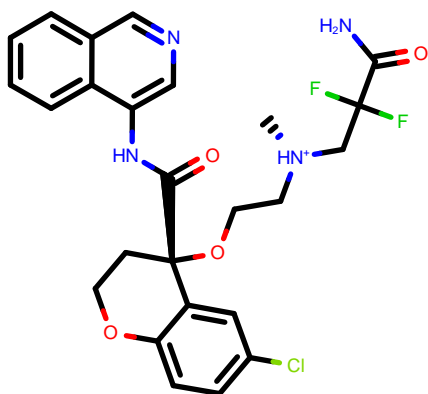
CID:	DAR-DIA-f6ee7aeb-4_1
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(C@H)(C@H)CC3=O)4ccc(cc4)C(=O)CC(F)(F)F)c5c[nH]c5=O</chem>
RUN:	RUN3411
DDG (kcal/mol):	-2.56
dDDG (kcal/mol):	0.19

MAK-UNK-c749d764-26_5



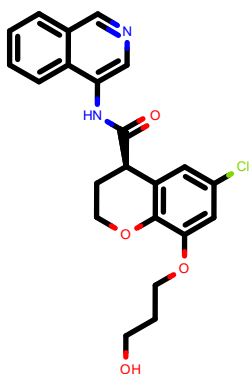
CID:	MAK-UNK-c749d764-26_5
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)C[C@@H]3CCCC[C@@H]1([C@H]3O)Cl</chem>
RUN:	RUN1050
DDG (kcal/mol):	-2.56
dDDG (kcal/mol):	0.30

MIC-UNK-cdc2493e-10_2



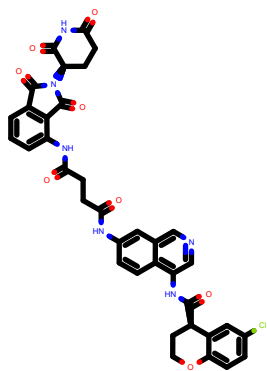
CID:	MIC-UNK-cdc2493e-10_2
SMILES:	<chem>CC(=O)N[C@H]1CC[C@@H](C1)N(c2ccccc2)C(=O)Nc3nccc4c3cccc4</chem>
RUN:	RUN546
DDG (kcal/mol):	-2.55
dDDG (kcal/mol):	0.34

PET-UNK-c5865d42-2_2



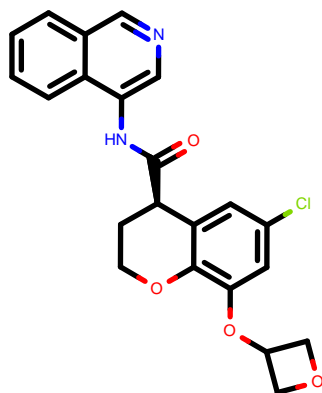
CID:	PET-UNK-c5865d42-2_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)Cc3cc(cc(c3)Cl)C([C@H]4CCN4)(F)F</chem>
RUN:	RUN429
DDG (kcal/mol):	-2.54
dDDG (kcal/mol):	0.35

ALP-UNI-8e43a71e-5_10



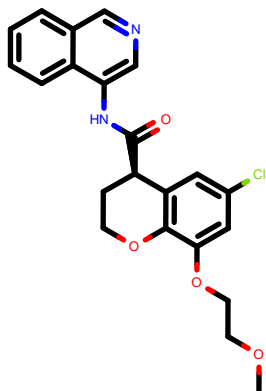
CID:	ALP-UNI-8e43a71e-5_10
SMILES:	<chem>C[C@H]1CN[C@@H]2C[C@@H]1NC(=O)C[C@H]2(COC1C=C2C(=O)C1)C(=O)Nc4nc5c6cccc5S(=O)(=O)C</chem>
RUN:	RUN2956
DDG (kcal/mol):	-2.54
dDDG (kcal/mol):	0.25

PET-UNK-c9c1e0d8-4_2



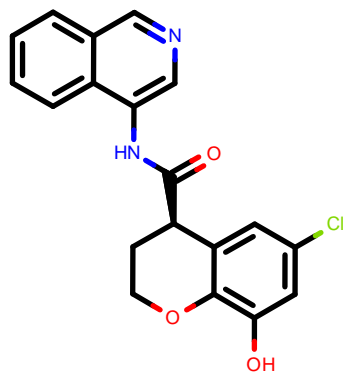
CID:	PET-UNK-c9c1e0d8-4_2
SMILES:	<chem>c1ccc2c(c1)cncc2N3CC[C@H](C3=O)c4cccc(c4)Cl</chem>
RUN:	RUN123
DDG (kcal/mol):	-2.54
dDDG (kcal/mol):	0.30

MAT-POS-1f3f1a6f-1_2



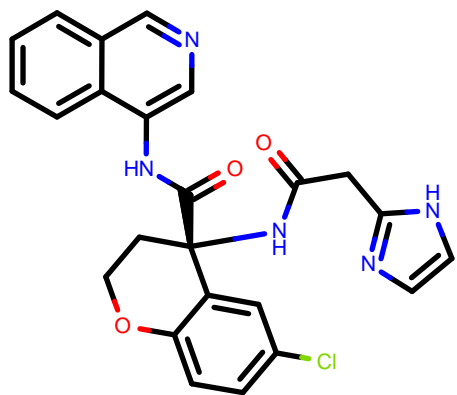
CID:	MAT-POS-1f3f1a6f-1_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H](c3ccc(c(c3)Cl)Cl)[NH3+]</chem>
RUN:	RUN2275
DDG (kcal/mol):	-2.54
dDDG (kcal/mol):	0.33

MIC-UNK-9582b2c5-2_5



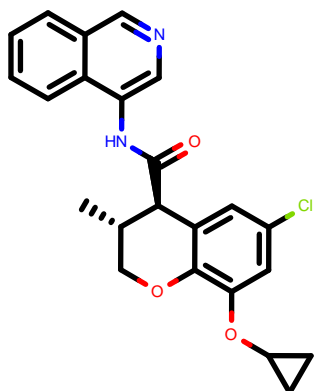
CID:	MIC-UNK-9582b2c5-2_5
SMILES:	<chem>CC(=O)N1CC[C@H]2[C@@H]1C[C@H](C1)C[C@H](C1=O)N2c3cccc(c3)Cl)c4cccc5c4cccc5</chem>
RUN:	RUN266
DDG (kcal/mol):	-2.52
dDDG (kcal/mol):	0.29

MIC-UNK-bcd487e9-9_1



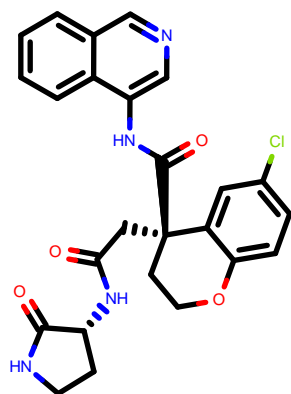
CID:	MIC-UNK-bcd487e9-9_1
SMILES:	<chem>Cn1cc(nn1)CN(c2cccc(c2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN597
DDG (kcal/mol):	-2.51
dDDG (kcal/mol):	0.22

EDJ-MED-c314995a-1_1



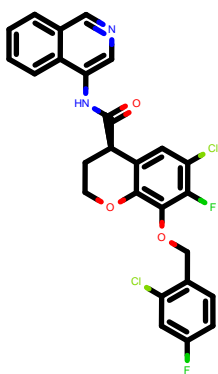
CID:	EDJ-MED-c314995a-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)N(CCC3CCCCC3)c4cccc(c4)Cl</chem>
RUN:	RUN126
DDG (kcal/mol):	-2.51
dDDG (kcal/mol):	0.25

ERI-UCB-d6de1f3c-5_1



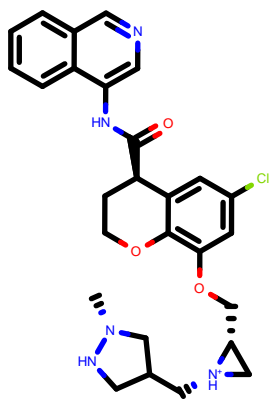
CID:	ERI-UCB-d6de1f3c-5_1
SMILES:	<chem>c1ccc2c(c1)cncc2C(=O)N3C[C@@H](N(C(=O)C3)c4cccc(c4)Cl)CC5CCCC5</chem>
RUN:	RUN1095
DDG (kcal/mol):	-2.51
dDDG (kcal/mol):	0.28

ALP-POS-ce760d3f-6_1



CID:	ALP-POS-ce760d3f-6_1
SMILES:	<chem>COc1ccc2cncc(c2c1)NC(=O)C3=CCOc4c3cc(cc4)Cl</chem>
RUN:	RUN1458
DDG (kcal/mol):	-2.49
dDDG (kcal/mol):	0.33

EDG-MED-ba1ac7b9-13_4



CID: EDG-MED-ba1ac7b9-13_4

SMILES:

C1N=C2C=CC=CC=C2N1C(=O)C3CCOC4=CC=C(C=C4)C3

RUN:

RUN2665

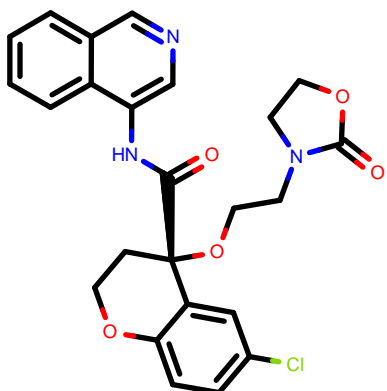
DDG (kcal/mol):

-2.47

dDDG (kcal/mol):

0.50

ALP-POS-02c6a514-32_1



CID: ALP-POS-02c6a514-32_1

SMILES:

CC(C)(C)C1CC(C1)N(C)C(=O)C2CCOC3=CC=C(C=C3)C2

RUN:

RUN644

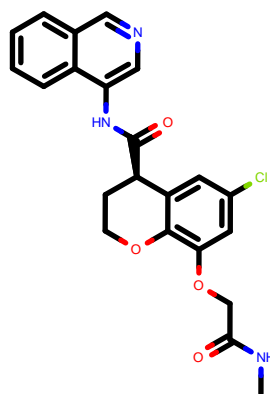
DDG (kcal/mol):

-2.47

dDDG (kcal/mol):

0.32

KAD-UNI-80f122c8-3_5



CID: KAD-UNI-80f122c8-3_5

SMILES:

C1CCOC2=CC=C(C=C2)C1C(=O)N(C)C(=O)C3CCOC4=CC=C(C=C4)C3

RUN:

RUN2293

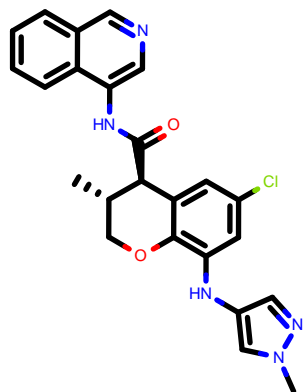
DDG (kcal/mol):

-2.46

dDDG (kcal/mol):

0.28

MIC-UNK-c66144cb-2_2



CID: MIC-UNK-c66144cb-2_2

SMILES:

C1CCOC2=CC=C(C=C2)C1C(=O)N(C)C(=O)C3CCOC4=CC=C(C=C4)C3

RUN:

RUN131

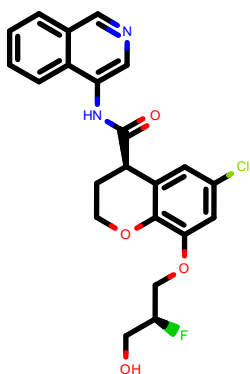
DDG (kcal/mol):

-2.46

dDDG (kcal/mol):

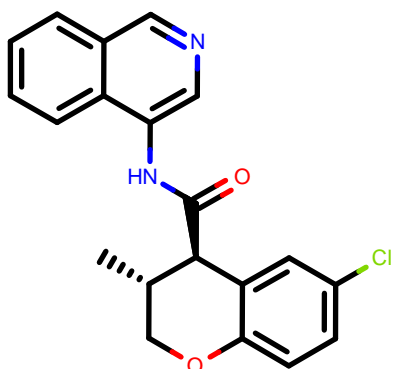
0.30

ALP-UNI-0676e700-10_1



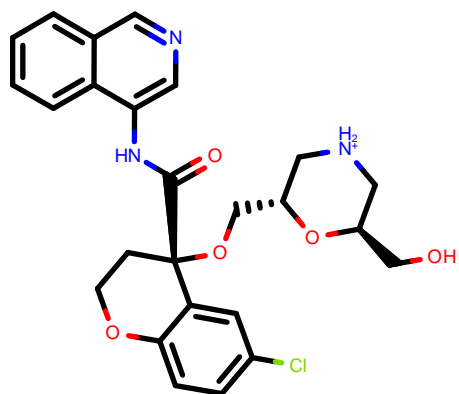
CID:	ALP-UNI-0676e700-10_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)CNC(=O)c5ccc6nnnn6c5</chem>
RUN:	RUN2453
DDG (kcal/mol):	-2.46
dDDG (kcal/mol):	0.40

DAR-DIA-0cde14eb-47_1



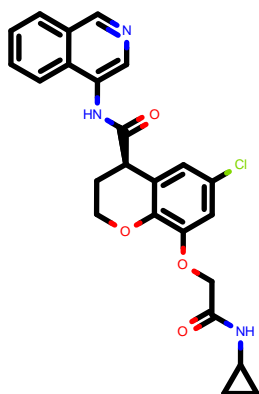
CID:	DAR-DIA-0cde14eb-47_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)Cc3cccc(c3)C4(CC4)Cl</chem>
RUN:	RUN11
DDG (kcal/mol):	-2.45
dDDG (kcal/mol):	0.20

DAR-DIA-6a508060-3_1



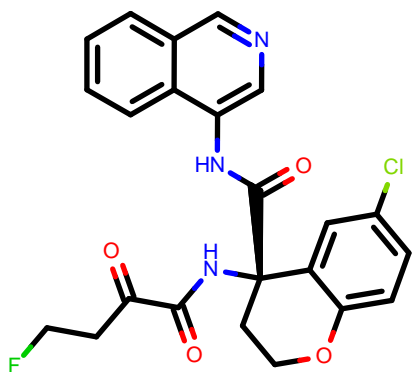
CID:	DAR-DIA-6a508060-3_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4c3cc(cc4)C#N</chem>
RUN:	RUN336
DDG (kcal/mol):	-2.45
dDDG (kcal/mol):	0.32

ALP-POS-2da19ca7-7_8



CID:	ALP-POS-2da19ca7-7_8
SMILES:	<chem>C[C@H]1CN[C@@H]2C[C@@H]1NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)CNC(=O)S(=O)(=O)C</chem>
RUN:	RUN2392
DDG (kcal/mol):	-2.44
dDDG (kcal/mol):	0.34

MIC-UNK-cdc2493e-8_2



CID: MIC-UNK-cdc2493e-8_2

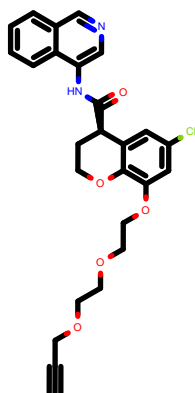
SMILES: c1ccc2c(c1)ncnc2NC(=O)N(c3cccc(c3)Cl)[C@H]4CC[C@@H]5CCCC[C@H]5C4

RUN: RUN538

DDG (kcal/mol): -2.43

dDDG (kcal/mol): 0.27

ALP-UNI-3496895b-1_1



CID: ALP-UNI-3496895b-1_1

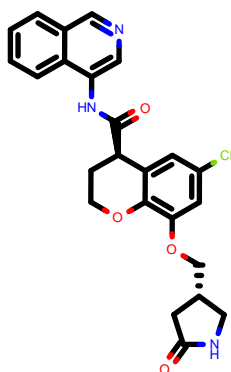
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3(COCc4c3cc(c4)Cl)CC(=O)NCCS(=O)(=O)N@]R[C@@H]5CCCC5

RUN: RUN2499

DDG (kcal/mol): -2.43

dDDG (kcal/mol): 0.49

MAT-POS-fce787c2-8_2



CID: MAT-POS-fce787c2-8_2

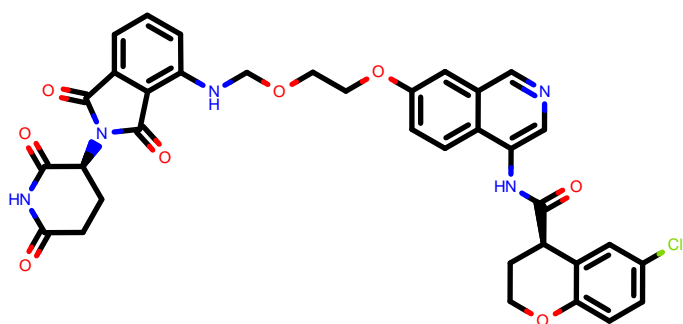
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@H](C#N)c3ccc(c(c3)Cl)Cl

RUN: RUN2155

DDG (kcal/mol): -2.42

dDDG (kcal/mol): 0.35

PET-UNK-1901c25b-1_1



CID: PET-UNK-1901c25b-1_1

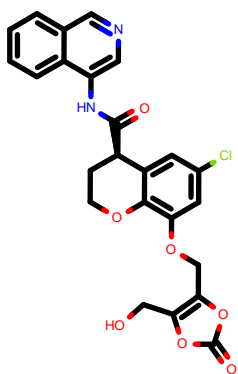
SMILES: CN(C)c1ccc(cc1)N(Cc2ccsc2)C(=O)Cc3cnc4c3cccc4

RUN: RUN302

DDG (kcal/mol): -2.41

dDDG (kcal/mol): 0.46

EDJ-MED-d203f206-16_1



CID: EDJ-MED-d203f206-16_1

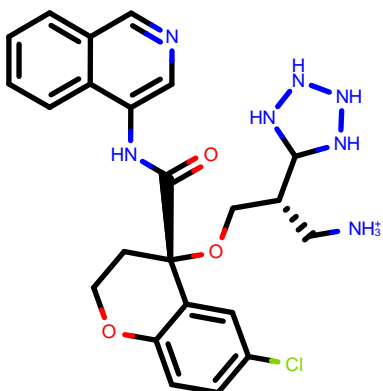
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4cc3cc(cc4)Cl)CC(=O)N5CCCC[C@H]5c6n[n-]nn6

RUN: RUN2578

DDG (kcal/mol): -2.39

dDDG (kcal/mol): 0.31

ALP-UNI-8e43a71e-15_24



CID: ALP-UNI-8e43a71e-15_24

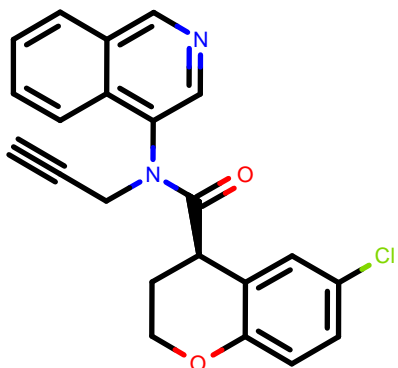
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4cc3cc(cc4)Cl)C[C@H]5C[C@H]6[C@@H]5C[C@@H]6Cl=O

RUN: RUN3005

DDG (kcal/mol): -2.39

dDDG (kcal/mol): 0.19

DAR-DIA-ecdbc7dd-12_1



CID: DAR-DIA-ecdbc7dd-12_1

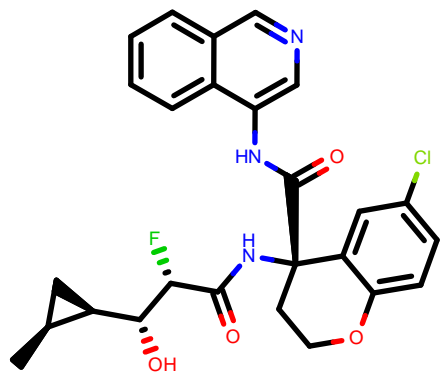
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCNc4c3cc(cc4)Cl)C[NH+]5CCCCC5

RUN: RUN2894

DDG (kcal/mol): -2.38

dDDG (kcal/mol): 0.10

MAK-UNK-ffc90da7-2_2



CID: MAK-UNK-ffc90da7-2_2

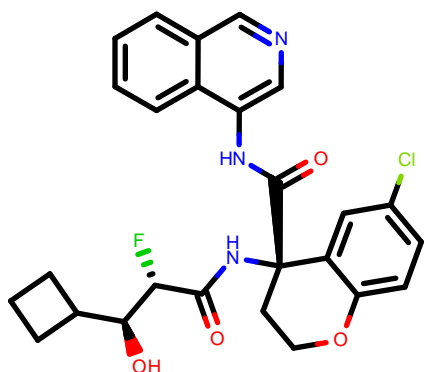
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3CCCC[C@@H]3[C@@H]3O)Cl

RUN: RUN682

DDG (kcal/mol): -2.38

dDDG (kcal/mol): 0.29

VLA-UCB-34f3ed0c-1_1



CID: VLA-UCB-34f3ed0c-1_1

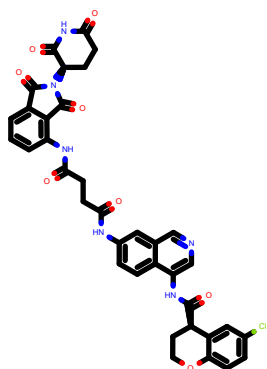
SMILES: C=C(C#N)C(=O)N(c1cncc2c1cccc2)C(=O)[C@@H]3CCOc4c3cc(cc4)Cl

RUN: RUN627

DDG (kcal/mol): -2.37

dDDG (kcal/mol): 0.29

ALP-POS-477dc5b7-1_1



CID: ALP-POS-477dc5b7-1_1

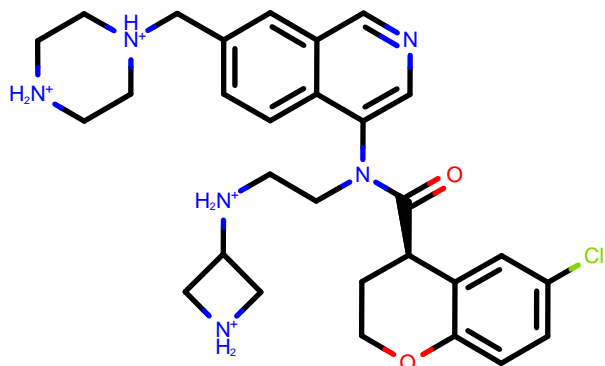
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCCC4c3cc(cc4)Cl

RUN: RUN294

DDG (kcal/mol): -2.36

dDDG (kcal/mol): 0.49

ALP-UNI-b33a865d-1_1



CID: ALP-UNI-b33a865d-1_1

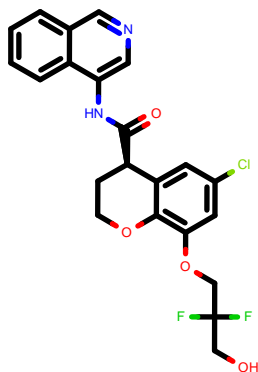
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCN(c4c3cc(cc4)Cl)C5cnc[nH]5

RUN: RUN354

DDG (kcal/mol): -2.36

dDDG (kcal/mol): 0.56

VLA-UNK-83c3754c-1_1



CID: VLA-UNK-83c3754c-1_1

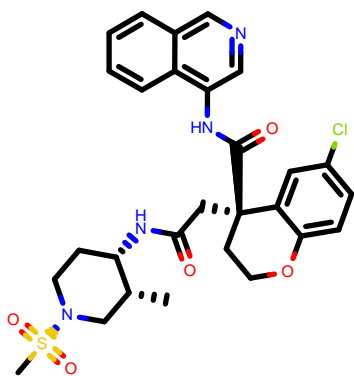
SMILES: c1ccc2c(c1)cncc2N3C(=O)[C@@]4(COc5c4cc(cc5)Cl)NC3=O

RUN: RUN2320

DDG (kcal/mol): -2.36

dDDG (kcal/mol): 0.37

ALP-POS-a0a4abd7-3_2



CID: ALP-POS-a0a4abd7-3_2

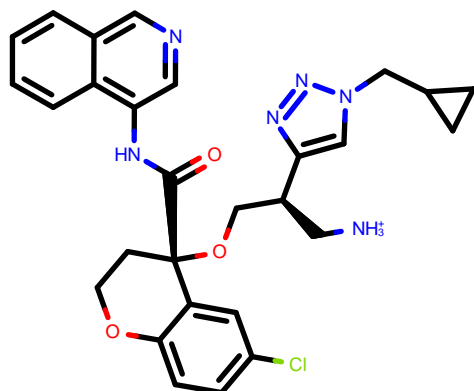
SMILES: Cn1c(=O)c(c[nH]c1=O)NC(C@H)2(CCOc3c2ccc(c3)C)C(=O)Nc4nccc5c4ccc5

RUN: RUN3570

DDG (kcal/mol): -2.34

dDDG (kcal/mol): 0.25

MIC-UNK-cdc2493e-2_1



CID: MIC-UNK-cdc2493e-2_1

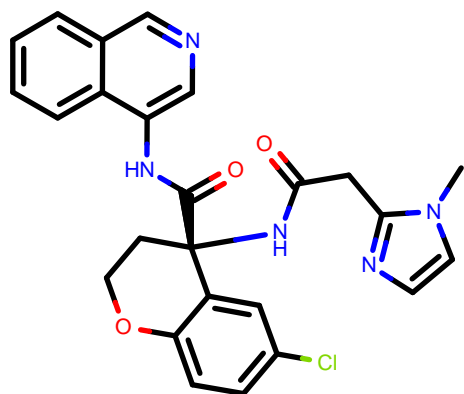
SMILES: c1ccc2c(c1)ncc2NC(=O)N(CC[C@@H]3CCCCO3)c4cccc(c4)Cl

RUN: RUN522

DDG (kcal/mol): -2.33

dDDG (kcal/mol): 0.33

MAK-UNK-8be7dca9-4_1



CID: MAK-UNK-8be7dca9-4_1

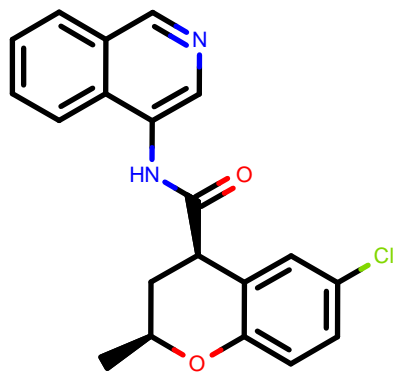
SMILES: c1cc2nccc(c2cc1N3CC[NH2+]CC3)NC(=O)[C@@H]4CCOCc5c4cc(cc5)Cl

RUN: RUN502

DDG (kcal/mol): -2.33

dDDG (kcal/mol): 0.22

MIC-UNK-d36ab305-1_2



CID: MIC-UNK-d36ab305-1_2

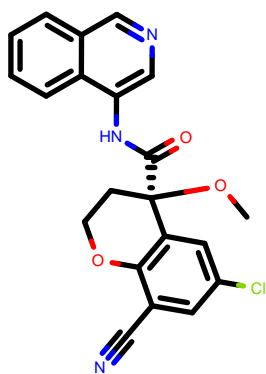
SMILES: CC(=O)Nc1ccc(cc1)[C@H](c2cccc(c2)C)C(=O)Nc3nccc4c3cccc4

RUN: RUN152

DDG (kcal/mol): -2.32

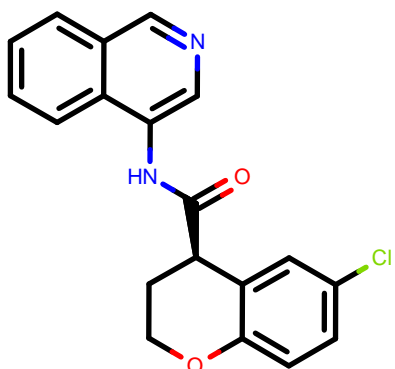
dDDG (kcal/mol): 0.18

VLA-UCB-29506327-1_1



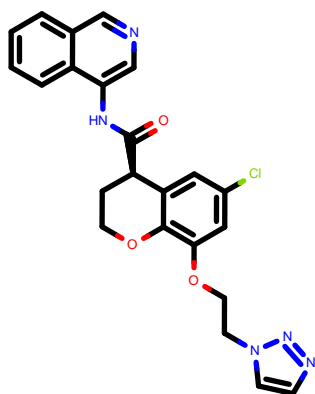
CID:	VLA-UCB-29506327-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3C(=O)[C@]4(CCOC5c4cc(cc5)Cl)NC3=O</chem>
RUN:	RUN1404
DDG (kcal/mol):	-2.32
dDDG (kcal/mol):	0.09

ERI-UCB-ce40166b-9_1



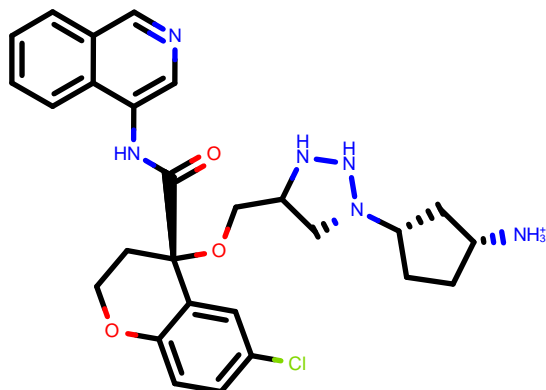
CID:	ERI-UCB-ce40166b-9_1
SMILES:	<chem>c1ccc2c(c1)cncc2CC(=O)Nc3cc(cc(c3)O)[C@@H]4CC(=O)N4C#N</chem>
RUN:	RUN45
DDG (kcal/mol):	-2.31
dDDG (kcal/mol):	0.19

ALF-EVA-5b152d2f-6_1



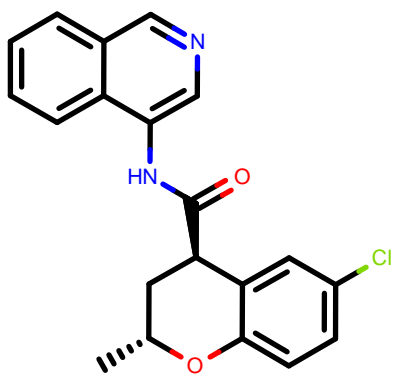
CID:	ALF-EVA-5b152d2f-6_1
SMILES:	<chem>C[C@H]1CCOc2ccc(cc2[C@@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN2359
DDG (kcal/mol):	-2.31
dDDG (kcal/mol):	0.35

MIC-UNK-91acba05-5_2



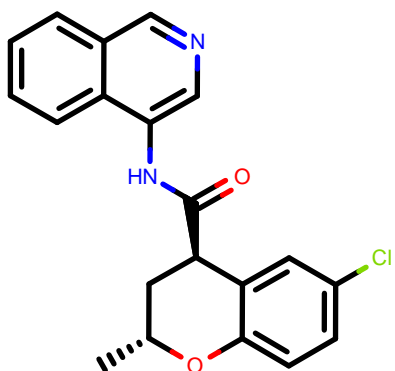
CID:	MIC-UNK-91acba05-5_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCNc4c3cc(cc4)Cl</chem>
RUN:	RUN477
DDG (kcal/mol):	-2.30
dDDG (kcal/mol):	0.35

VLA-UCB-1dbca3b4-3_1



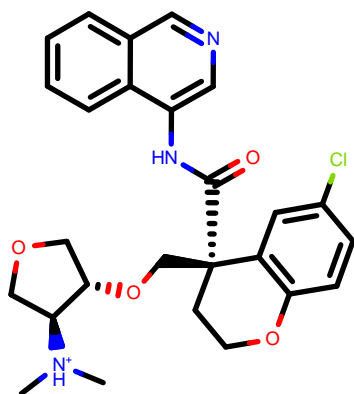
CID:	VLA-UCB-1dbca3b4-3_1
SMILES:	<chem>c1ccc2c(c1)cncc2n3c(cnc3=O)c4cccc(c4)Cl)[O-]</chem>
RUN:	RUN153
DDG (kcal/mol):	-2.29
dDDG (kcal/mol):	0.20

ADA-UCB-6c2cb422-1_1



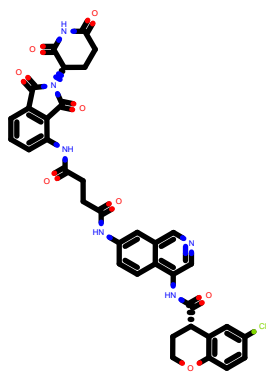
CID:	ADA-UCB-6c2cb422-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)Cc3cccc(c3)Cl</chem>
RUN:	RUN53
DDG (kcal/mol):	-2.28
dDDG (kcal/mol):	0.19

DAR-DIA-0d514e7d-31_10



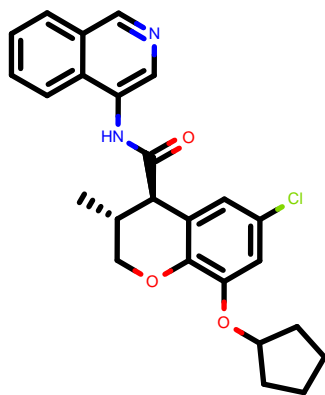
CID:	DAR-DIA-0d514e7d-31_10
SMILES:	<chem>C[C@H]1CC[C@@H]2C=CC(=C[C@H]2[C@H]1C(=O)Nc3ccc4c3ccc4)Cl</chem>
RUN:	RUN842
DDG (kcal/mol):	-2.28
dDDG (kcal/mol):	0.23

ALP-UNI-8e43a71e-7_2



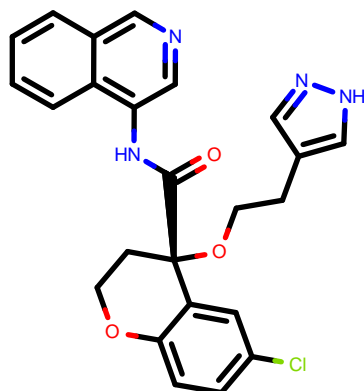
CID:	ALP-UNI-8e43a71e-7_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]3(C)C(CO)C4C3C(CO)C(=O)N5CCN(CC5)CCC#N</chem>
RUN:	RUN2963
DDG (kcal/mol):	-2.27
dDDG (kcal/mol):	0.40

PET-UNK-c9c1e0d8-1_1



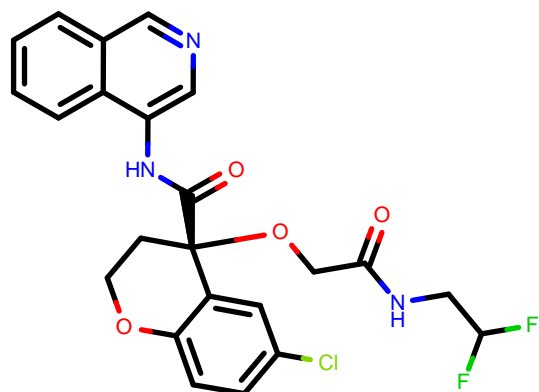
CID:	PET-UNK-c9c1e0d8-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)Cc3ccccc(c3)Br</chem>
RUN:	RUN119
DDG (kcal/mol):	-2.27
dDDG (kcal/mol):	0.29

EDJ-MED-ee07cf00-11_4



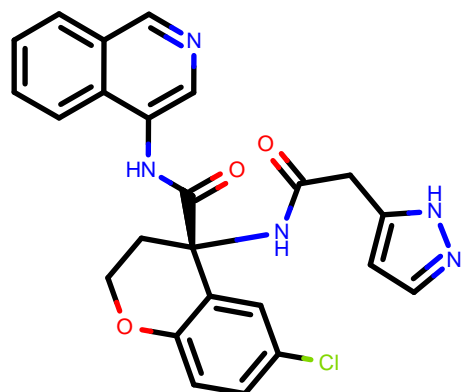
CID:	EDJ-MED-ee07cf00-11_4
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)C[C@H](c3ccccc3)C1NC(=O)C[C@H](C)C1=O</chem>
RUN:	RUN2828
DDG (kcal/mol):	-2.27
dDDG (kcal/mol):	0.18

MAK-UNK-ffc90da7-9_2



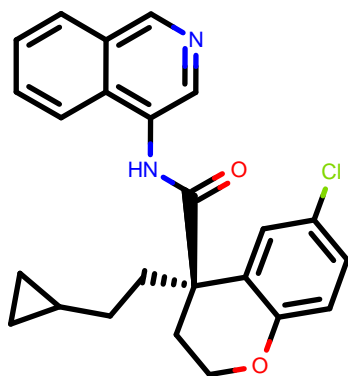
CID:	MAK-UNK-ffc90da7-9_2
SMILES:	<chem>C[C@H]1[C@@H](CCO1)SC[C@@H](C)c2ccc3c(c2)cncc3NC(=O)Cc4cccc(c4)Cl</chem>
RUN:	RUN710
DDG (kcal/mol):	-2.26
dDDG (kcal/mol):	0.27

EDJ-MED-e4b030d8-7_1



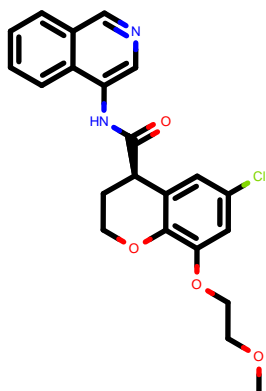
CID:	EDJ-MED-e4b030d8-7_1
SMILES:	<chem>COc1cc(cc2c1OCC[C@H]2C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN289
DDG (kcal/mol):	-2.25
dDDG (kcal/mol):	0.31

EDG-MED-ba1ac7b9-22_2



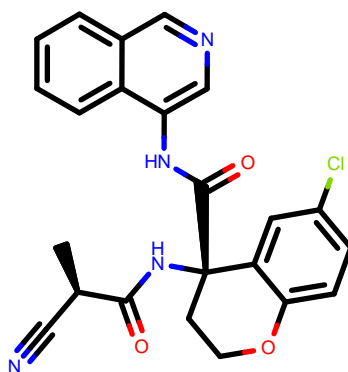
CID:	EDG-MED-ba1ac7b9-22_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@H]3[C@@H](C)CCOC4=CC=C(Cl)C=C4C(=O)N5C6CC6C=C(Cl)C=C5C1=O</chem>
RUN:	RUN2703
DDG (kcal/mol):	-2.25
dDDG (kcal/mol):	0.10

DAR-DIA-9e4459de-15_6



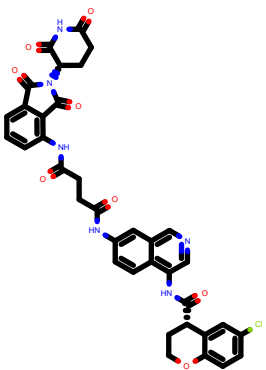
CID:	DAR-DIA-9e4459de-15_6
SMILES:	<chem>c1cc2c(c1)NC(=O)CCOC(=O)N3CCCC3C(=O)N[C@@H]4C(=O)N5C6CC6C=C(Cl)C=C5C1=O</chem>
RUN:	RUN1448
DDG (kcal/mol):	-2.25
dDDG (kcal/mol):	0.37

ADA-UCB-dc2b944c-13_1



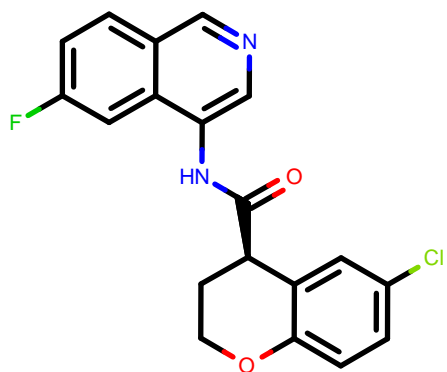
CID:	ADA-UCB-dc2b944c-13_1
SMILES:	<chem>C[C@@H]1COc2ccc(cc2[C@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN611
DDG (kcal/mol):	-2.25
dDDG (kcal/mol):	0.28

ALP-POS-fe871b40-14_1



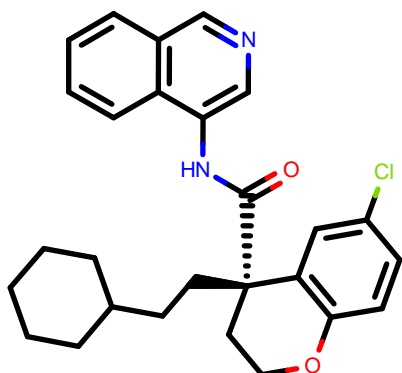
CID:	ALP-POS-fe871b40-14_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CC(=O)Nc4c3cc(ccn4)Cl</chem>
RUN:	RUN3133
DDG (kcal/mol):	-2.25
dDDG (kcal/mol):	0.49

PET-UNK-431b3bfb-1_1



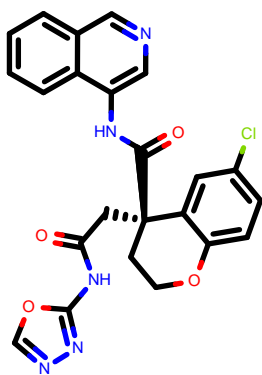
CID:	PET-UNK-431b3bfb-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3CCO[C@H](C3=O)c4cccc(c4)Cl</chem>
RUN:	RUN271
DDG (kcal/mol):	-2.25
dDDG (kcal/mol):	0.19

LON-WEI-4d77710c-42_1



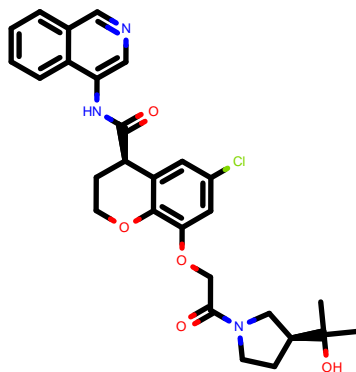
CID:	LON-WEI-4d77710c-42_1
SMILES:	<chem>CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)N[C@@H]3CCCC4c3ccccc4</chem>
RUN:	RUN229
DDG (kcal/mol):	-2.24
dDDG (kcal/mol):	0.22

MAK-UNK-c749d764-22_3



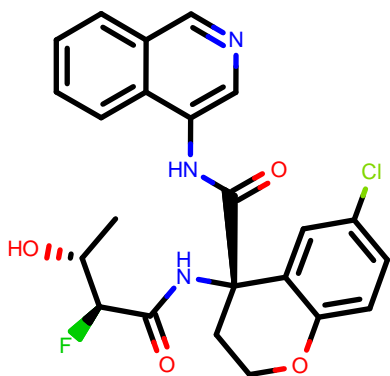
CID:	MAK-UNK-c749d764-22_3
SMILES:	<chem>C[C@@H](C)[NH2+][C][C@H]1CCCCO1)[C@@H](c2ccccc2)C(=O)Nc3ncoc4c3ccccc4</chem>
RUN:	RUN1024
DDG (kcal/mol):	-2.24
dDDG (kcal/mol):	0.30

ALP-POS-5bb456a5-2_5



CID:	ALP-POS-5bb456a5-2_5
SMILES:	<chem>C[C@@H](CN@@)CC[C@@H]1NC(=O)C[C@@]2(CCOc3c2cc(c3)Cl)C(=O)Nc4ncoc5c4ccccc5S(=O)(=O)C</chem>
RUN:	RUN2426
DDG (kcal/mol):	-2.21
dDDG (kcal/mol):	0.40

VLA-UCB-34f3ed0c-14_1



CID: VLA-UCB-34f3ed0c-14_1

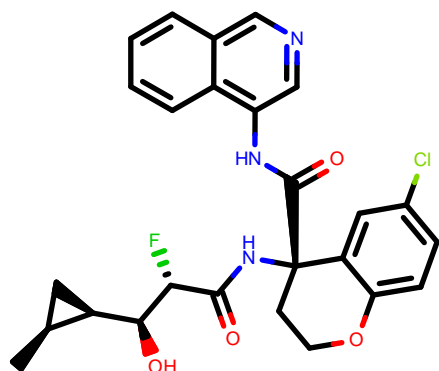
SMILES: c1ccc2c(c1)cnc2N3C(=O)[C@@H](C(=O)O)C(=O)N(C3=O)CC6CCCC6

RUN: RUN636

DDG (kcal/mol): -2.21

dDDG (kcal/mol): 0.27

PET-UNK-824b5c6a-1_1



CID: PET-UNK-824b5c6a-1_1

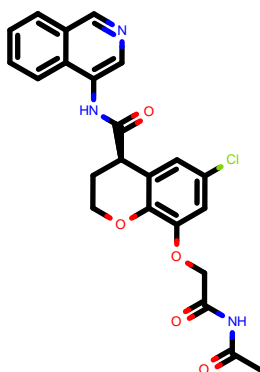
SMILES: CCO[C@]1(CCOC2c1cc(cc2)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN3287

DDG (kcal/mol): -2.20

dDDG (kcal/mol): 0.15

ALP-UNI-0676e700-27_1



CID: ALP-UNI-0676e700-27_1

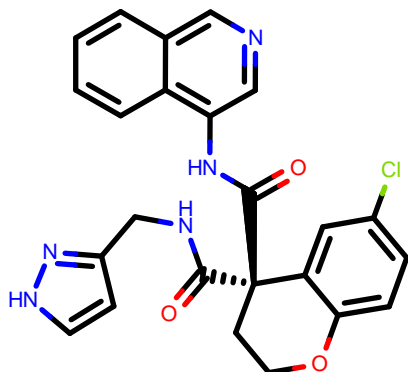
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H](C(=O)O)C(=O)N(C3=O)CC6CCCC6

RUN: RUN2475

DDG (kcal/mol): -2.20

dDDG (kcal/mol): 0.30

ALF-EVA-650655fc-3_2



CID: ALF-EVA-650655fc-3_2

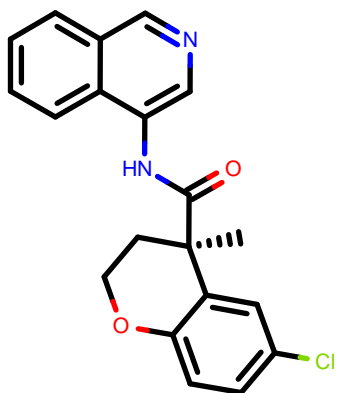
SMILES: c1ccc2c(c1)cnc2NC(=O)N(CCC3CCCC3)C(=O)N(C3=O)CC6CCCC6

RUN: RUN2856

DDG (kcal/mol): -2.19

dDDG (kcal/mol): 0.16

MAT-POS-2492181e-8_1



CID: MAT-POS-2492181e-8_1

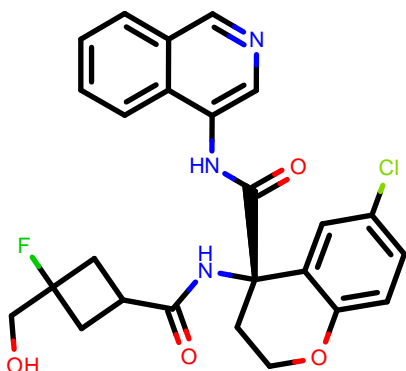
SMILES: CC1CCN(CC1)c2ccc(cc2)NC(=O)Nc3cn(c(=O)c4c3cccc4)CC(C)C

RUN: RUN98

DDG (kcal/mol): -2.18

dDDG (kcal/mol): 0.22

JAG-UCB-706446eb-6_1



CID: JAG-UCB-706446eb-6_1

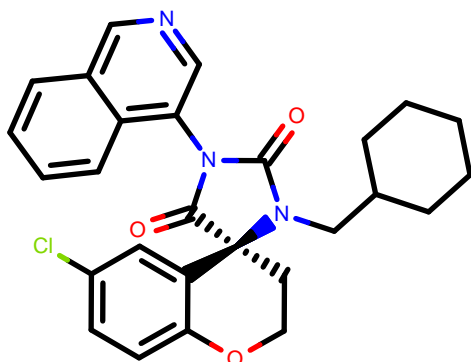
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(C)CCOc4ccc(cc4)C1CC(=O)Nc5cc[nH]5

RUN: RUN618

DDG (kcal/mol): -2.18

dDDG (kcal/mol): 0.26

MIC-UNK-0a05c952-1_3



CID: MIC-UNK-0a05c952-1_3

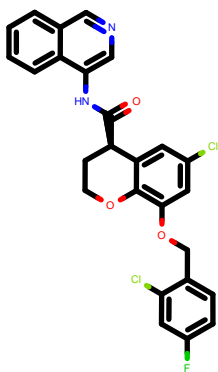
SMILES: c1ccc2c(c1)cncc2N3[C@@H](C[C@H](C3=O)c4cccc(c4)Cl)[C@@H]5CO5

RUN: RUN3498

DDG (kcal/mol): -2.18

dDDG (kcal/mol): 0.20

ALP-POS-ce760d3f-5_1



CID: ALP-POS-ce760d3f-5_1

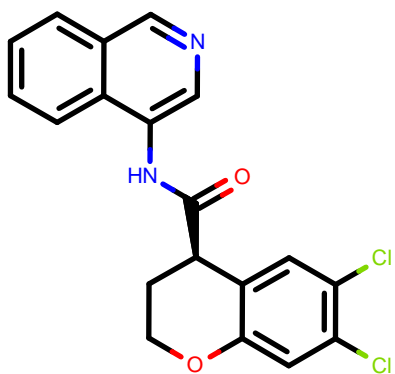
SMILES: c1cc2cncc(c2cc1F)NC(=O)C3=CCOc4ccc(cc4)Cl

RUN: RUN1459

DDG (kcal/mol): -2.17

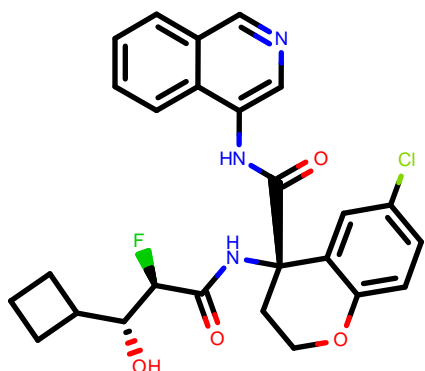
dDDG (kcal/mol): 0.37

MIC-UNK-50cce87d-3_1



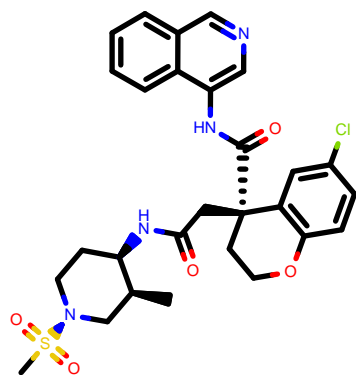
CID:	MIC-UNK-50cce87d-3_1
SMILES:	<chem>Cc1cccc2c1c(cnc2)NC(=O)Cc3cccc(c3)Cl</chem>
RUN:	RUN662
DDG (kcal/mol):	-2.16
dDDG (kcal/mol):	0.26

MIC-UNK-5a93dd5f-9_3



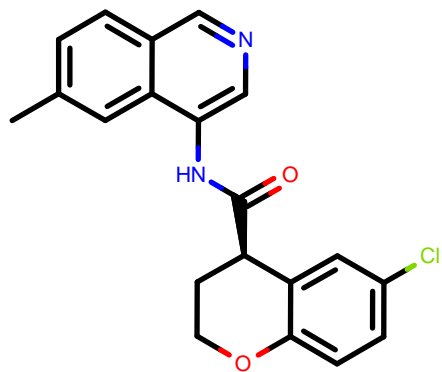
CID:	MIC-UNK-5a93dd5f-9_3
SMILES:	<chem>CN(C)[C@@H]1CC[N@H+](C1)[C@@H](C)[C@@H](c2cccc(c2)Cl)C(=O)Nc3nccc4c3cccc4</chem>
RUN:	RUN780
DDG (kcal/mol):	-2.16
dDDG (kcal/mol):	0.27

MAK-UNK-c749d764-2_5



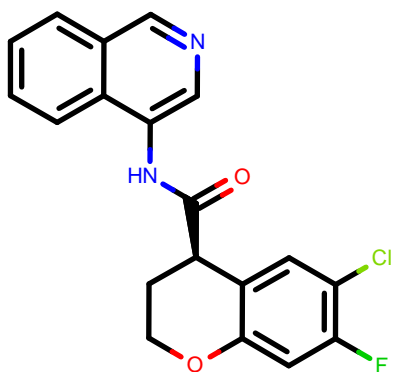
CID:	MAK-UNK-c749d764-2_5
SMILES:	<chem>CC[C@@H]1CCC[C@@H]([C@@H]1O)CC(=O)Nc2nccc3c2cccc3</chem>
RUN:	RUN901
DDG (kcal/mol):	-2.16
dDDG (kcal/mol):	0.27

DAR-DIA-0d514e7d-31_13



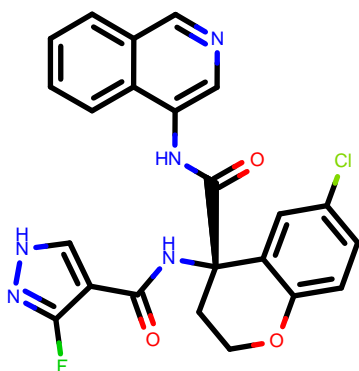
CID:	DAR-DIA-0d514e7d-31_13
SMILES:	<chem>C[C@@H]1CCO[C@@H]2C=CC=C[C@H]2[C@H]1C(=O)Nc3nccc4c3cccc4Cl</chem>
RUN:	RUN846
DDG (kcal/mol):	-2.15
dDDG (kcal/mol):	0.19

ALP-POS-75715966-2_1



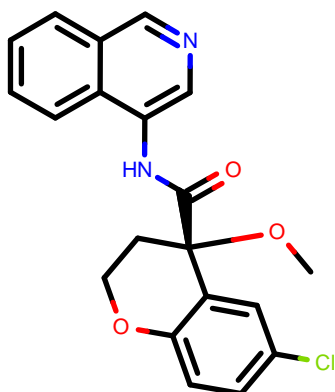
CID:	ALP-POS-75715966-2_1
SMILES:	<chem>CN(C)c1ccc(cc1)N(Cc2cccc(c2)Cl)C(=O)Cc3cncc4c3cccc4</chem>
RUN:	RUN1490
DDG (kcal/mol):	-2.14
dDDG (kcal/mol):	0.20

ADA-UCB-dc2b944c-8_1



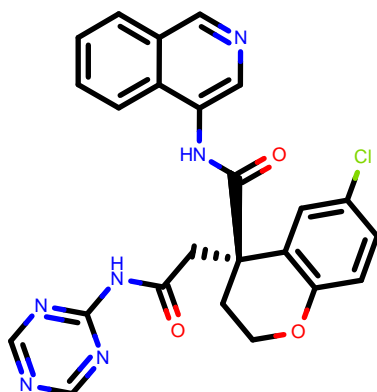
CID:	ADA-UCB-dc2b944c-8_1
SMILES:	<chem>c1cc2c(cncc2NC(=O)[C@@H]3CCOCc4c3cc(cc4)Cl)c(c1)Cl</chem>
RUN:	RUN606
DDG (kcal/mol):	-2.14
dDDG (kcal/mol):	0.25

DAR-DIA-0d514e7d-31_7



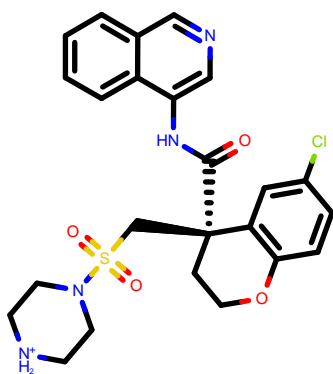
CID:	DAR-DIA-0d514e7d-31_7
SMILES:	<chem>C[C@@H]1CCO[C@H]2C=CC(=C[C@H]2[C@@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN840
DDG (kcal/mol):	-2.13
dDDG (kcal/mol):	0.22

NAU-LAT-2fed8305-4_2



CID:	NAU-LAT-2fed8305-4_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)C[N@H+]3CCC=C(C3)F</chem>
RUN:	RUN1110
DDG (kcal/mol):	-2.12
dDDG (kcal/mol):	0.24

EDG-MED-90036822-13_1



CID: EDG-MED-90036822-13_1

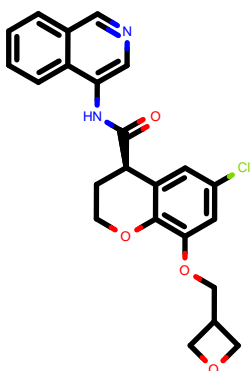
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)NC(=O)CO

RUN: RUN1670

DDG (kcal/mol): -2.12

dDDG (kcal/mol): 0.20

MAT-POS-1f3f1a6f-3_2



CID: MAT-POS-1f3f1a6f-3_2

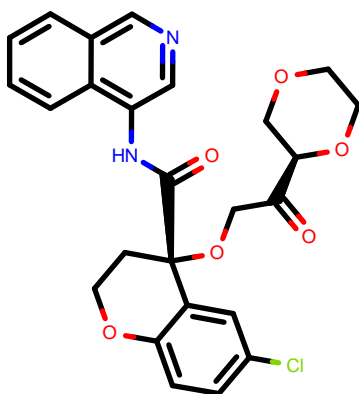
SMILES: CC(=O)N[C@H](c1ccc(c(c1)Cl)Cl)C(=O)Nc2cnc3c2cccc3

RUN: RUN2282

DDG (kcal/mol): -2.11

dDDG (kcal/mol): 0.34

NIR-THE-2069301b-1_2



CID: NIR-THE-2069301b-1_2

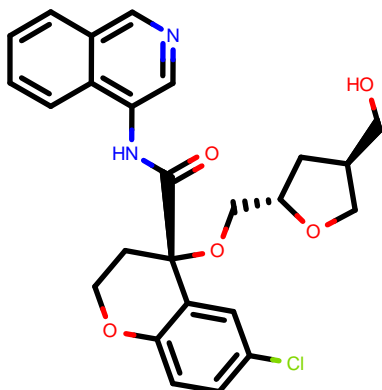
SMILES: C=C(C(=O)N(c1cnc2c1cccc2)C(=O)[C@H]3COc4c3cc(cc4)Cl)F

RUN: RUN3321

DDG (kcal/mol): -2.11

dDDG (kcal/mol): 0.20

MAT-POS-dd3ad2b5-2_2



CID: MAT-POS-dd3ad2b5-2_2

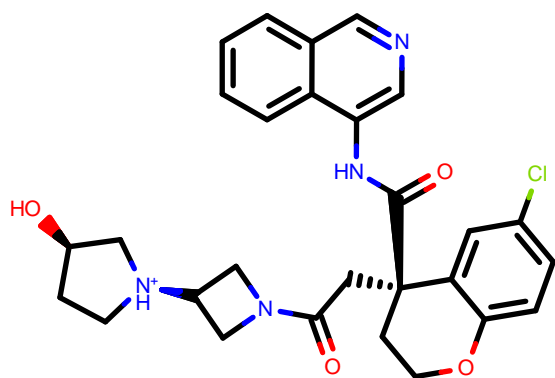
SMILES: CC(=O)N1Cc2ccc(cc2[C@H](C1)C(=O)Nc3cncc4c3cccc4)Cl

RUN: RUN3534

DDG (kcal/mol): -2.11

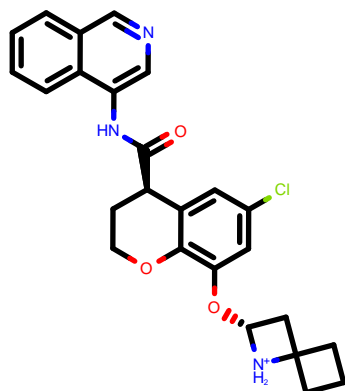
dDDG (kcal/mol): 0.19

GIA-UNK-80c9bc96-1_1



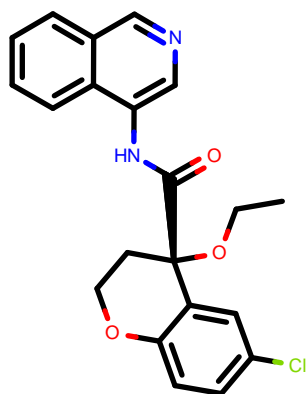
CID:	GIA-UNK-80c9bc96-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3CC[C@@H](C3=O)c4cccc(c4)Cl</chem>
RUN:	RUN1101
DDG (kcal/mol):	-2.11
dDDG (kcal/mol):	0.40

MAT-POS-b5746674-101_1



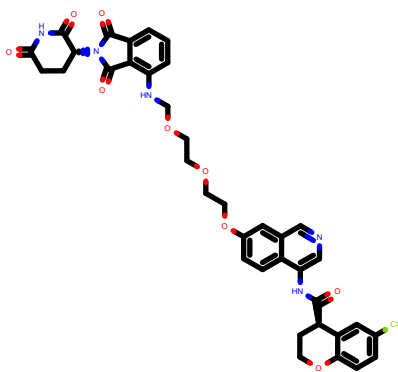
CID:	MAT-POS-b5746674-101_1
SMILES:	<chem>CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)N3CC[C@@H](C3)c4cccc4</chem>
RUN:	RUN77
DDG (kcal/mol):	-2.11
dDDG (kcal/mol):	0.47

DAR-DIA-9e4459de-11_14



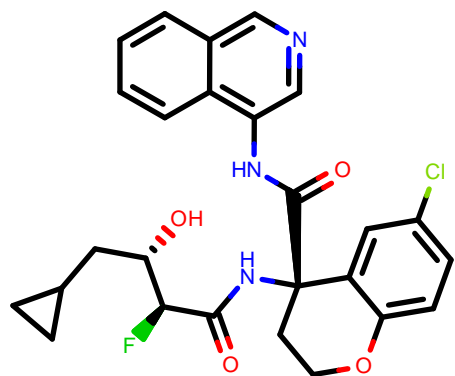
CID:	DAR-DIA-9e4459de-11_14
SMILES:	<chem>c1cc2c(c1)NCOCCOCCOCC3CC4C(C)CNC4N(C)O[C@@H](C3)C5COC6C5C(C6)Cl(C)C(C)O</chem>
RUN:	RUN1422
DDG (kcal/mol):	-2.11
dDDG (kcal/mol):	0.18

ALP-POS-477dc5b7-1_2



CID:	ALP-POS-477dc5b7-1_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCc4c3cc(cc4)Cl</chem>
RUN:	RUN296
DDG (kcal/mol):	-2.10
dDDG (kcal/mol):	0.48

MIC-UNK-0a05c952-1_6



CID: MIC-UNK-0a05c952-1_6

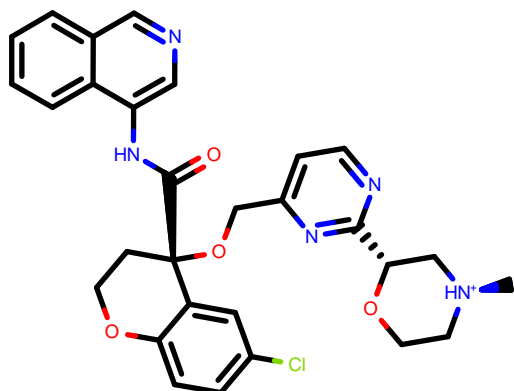
SMILES: c1ccc2c(c1)cncc2N3[C@H](C[C@@H](C3=O)c4cccc(c4)Cl)[C@H]5CO5

RUN: RUN3502

DDG (kcal/mol): -2.09

dDDG (kcal/mol): 0.19

DAR-DIA-23e5a6a0-10_2



CID: DAR-DIA-23e5a6a0-10_2

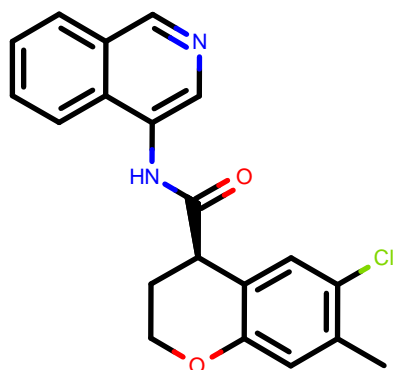
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCO4c3cc(cc4C[C@H]5CC6([NH2+][5]OC6)Cl

RUN: RUN421

DDG (kcal/mol): -2.09

dDDG (kcal/mol): 0.33

DAR-DIA-0d514e7d-13_1



CID: DAR-DIA-0d514e7d-13_1

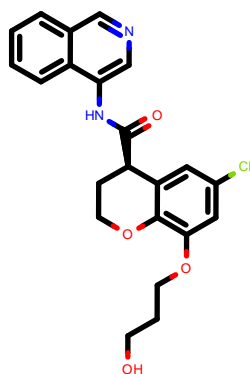
SMILES: C[C@H]1COc2c(cc(cc2[C@@H]1C(=O)Nc3ncc4c3cccc4)Cl)c5ccc(cc5)F

RUN: RUN815

DDG (kcal/mol): -2.09

dDDG (kcal/mol): 0.25

MAT-POS-e9e99895-6_1



CID: MAT-POS-e9e99895-6_1

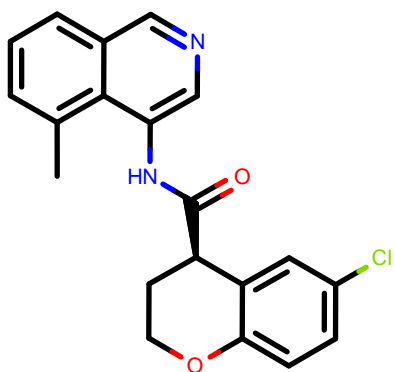
SMILES: C[C@@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2ncc3c2cccc3)NC(=O)CN4CCN(C4=O)C5CC5

RUN: RUN2256

DDG (kcal/mol): -2.09

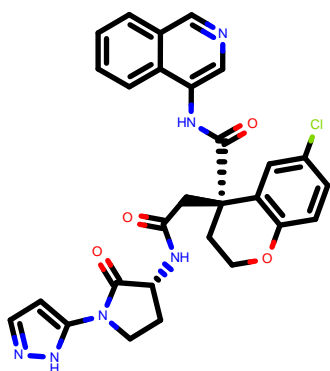
dDDG (kcal/mol): 0.30

ALP-POS-477dc5b7-3_2



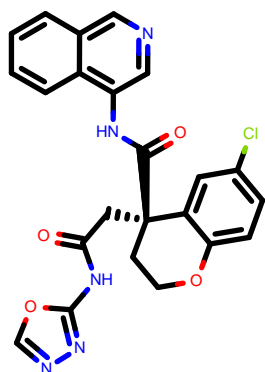
CID:	ALP-POS-477dc5b7-3_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CC5CC5</chem>
RUN:	RUN307
DDG (kcal/mol):	-2.08
dDDG (kcal/mol):	0.18

ALP-POS-696356e4-1_1



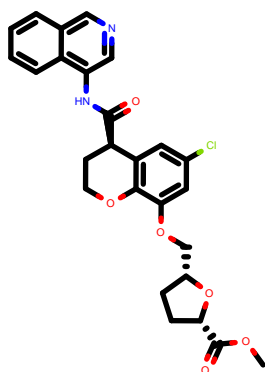
CID:	ALP-POS-696356e4-1_1
SMILES:	<chem>c1cc2cnc(c2cc1F)NC(=O)[C@@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN1314
DDG (kcal/mol):	-2.05
dDDG (kcal/mol):	0.28

MAT-POS-932d1078-1_4



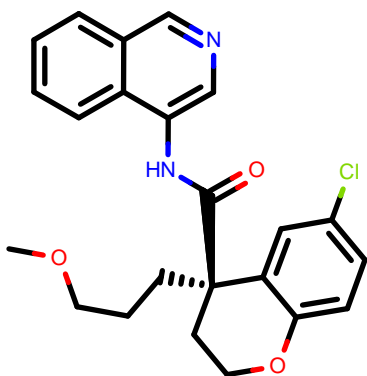
CID:	MAT-POS-932d1078-1_4
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CNC(=O)[C@H]5COC(N)@5(C)F)F</chem>
RUN:	RUN3589
DDG (kcal/mol):	-2.05
dDDG (kcal/mol):	0.10

EDJ-MED-d203f206-6_1



CID:	EDJ-MED-d203f206-6_1
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CNC(=O)[C@H]5COC(N)@5(C)F)F</chem>
RUN:	RUN2568
DDG (kcal/mol):	-2.04
dDDG (kcal/mol):	0.37

MIC-UNK-9582b2c5-1_2



CID: MIC-UNK-9582b2c5-1_2

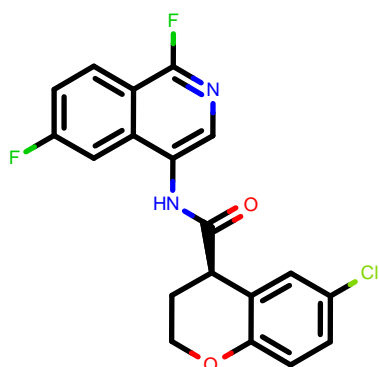
SMILES: CC(=O)N1CC[C@H]2[C@@H](C1)CN(C(=O)[C@@H]2c3ccccc3)Cl)c4ccc5c4ccccc5

RUN: RUN253

DDG (kcal/mol): -2.03

dDDG (kcal/mol): 0.29

LON-WEI-4d77710c-45_1



CID: LON-WEI-4d77710c-45_1

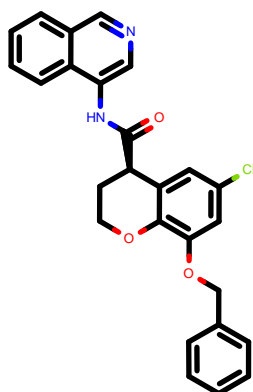
SMILES: Cn1cc(c2ccccc2c1=O)NC(=O)N3CCN(CC3)C(=O)c4ccccc4

RUN: RUN235

DDG (kcal/mol): -2.03

dDDG (kcal/mol): 0.21

DAR-DIA-9e4459de-15_9



CID: DAR-DIA-9e4459de-15_9

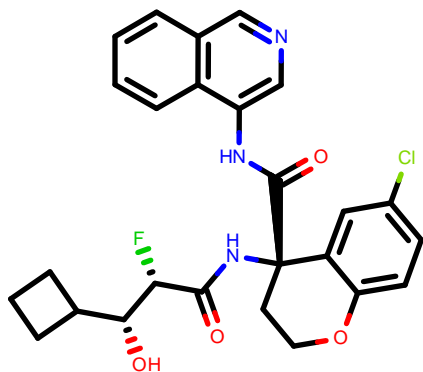
SMILES: c1cc2c(c1)NC(=O)CC(C(=O)N3CCCC3)C(=O)[C@@H]2c3ccccc3)Cl)c4ccc5c4ccccc5

RUN: RUN1449

DDG (kcal/mol): -2.03

dDDG (kcal/mol): 0.32

MIC-UNK-5a93dd5f-12_6



CID: MIC-UNK-5a93dd5f-12_6

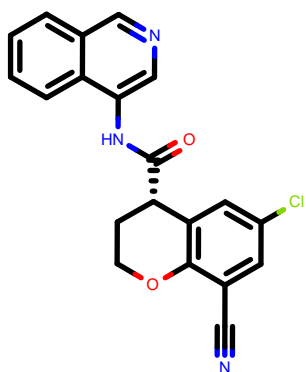
SMILES: c1ccc2c(c1)nc2NC(=O)[C@@H](c3ccccc3)Cl)N@@[C@@H](C4)NHH5CCCC5

RUN: RUN795

DDG (kcal/mol): -2.03

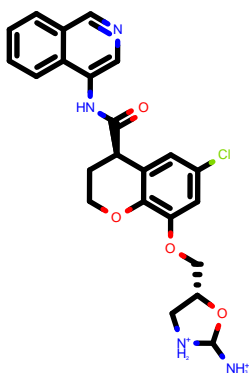
dDDG (kcal/mol): 0.35

DAR-DIA-5ff57136-2_1



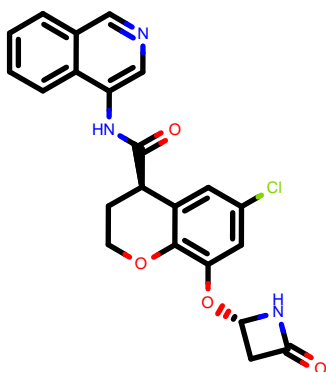
CID:	DAR-DIA-5ff57136-2_1
SMILES:	<chem>C[C@H]1COc2ccc(cc2[C@@H]1C(=O)N(c3cncc4c3cccc4)C(=O)C=C)Cl</chem>
RUN:	RUN1373
DDG (kcal/mol):	-2.02
dDDG (kcal/mol):	0.14

KAD-UNI-8a629cb0-33_1



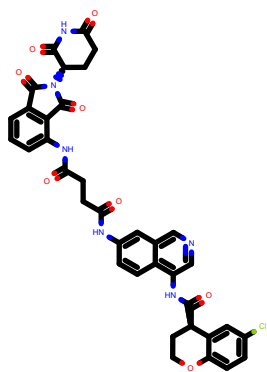
CID:	KAD-UNI-8a629cb0-33_1
SMILES:	<chem>CC[C@H](C[NH+])1CCN(CC1)C(=O)C[C@]2(C)COCc3c2cc(cc3)Cl)C(=O)Nc4cncc5c4cccc5)O</chem>
RUN:	RUN2120
DDG (kcal/mol):	-2.01
dDDG (kcal/mol):	0.42

NIR-WEI-f9286bb6-4_1



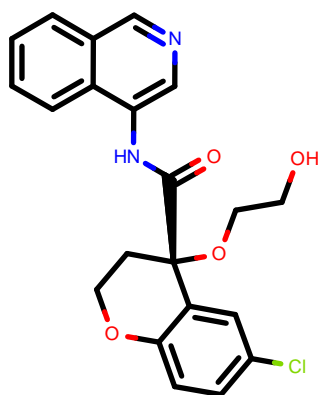
CID:	NIR-WEI-f9286bb6-4_1
SMILES:	<chem>CC(C)(C)NC(=O)[C@@H](c1cncc2c1cccc2)N(Cc3cccc3Cl)C(=O)C=C</chem>
RUN:	RUN138
DDG (kcal/mol):	-2.01
dDDG (kcal/mol):	0.26

MAR-UCB-6ab2ec87-2_1



CID:	MAR-UCB-6ab2ec87-2_1
SMILES:	<chem>C[C@@H]1[C@H](c2cc(ccc2N1)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3022
DDG (kcal/mol):	-2.01
dDDG (kcal/mol):	0.25

DAR-DIA-23e5a6a0-3_2



CID: DAR-DIA-23e5a6a0-3_2

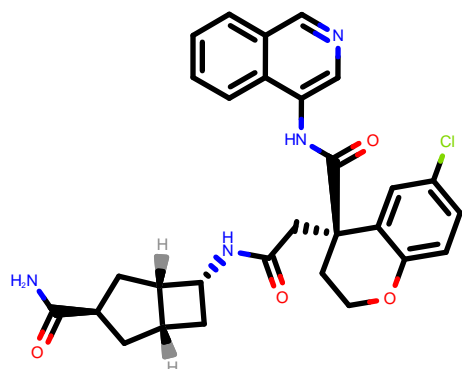
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4c3cc(cc4O)[C@H]5CC6([NH2+][5]COC6)Cl

RUN: RUN406

DDG (kcal/mol): -2.01

dDDG (kcal/mol): 0.23

LON-WEI-5e7d1b3e-29_1



CID: LON-WEI-5e7d1b3e-29_1

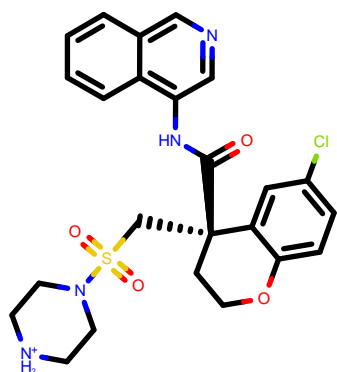
SMILES: CCOC(=O)Cc1csc(n1)NC(=O)Nc2cn(c(=O)c3c2cccc3)CC(C)C

RUN: RUN1340

DDG (kcal/mol): -2.01

dDDG (kcal/mol): 0.35

LAU-MED-88a3970a-13_1



CID: LAU-MED-88a3970a-13_1

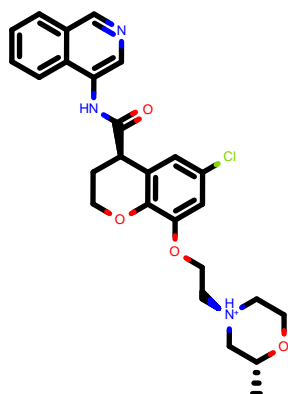
SMILES: COCCc1cc(cc2c1OCC[C@H]2C(=O)Nc3cncc4c3cccc4)Cl

RUN: RUN1509

DDG (kcal/mol): -2.00

dDDG (kcal/mol): 0.40

EDJ-MED-d203f206-2_1



CID: EDJ-MED-d203f206-2_1

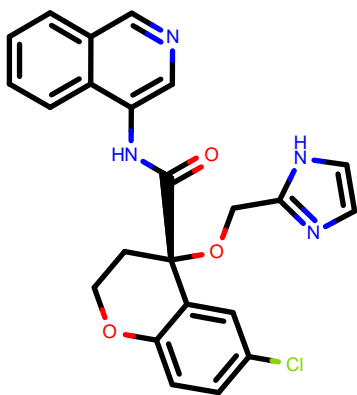
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CC(=O)Nc5ncon5

RUN: RUN2563

DDG (kcal/mol): -2.00

dDDG (kcal/mol): 0.54

MAK-UNK-8be7dca9-3_4



CID: MAK-UNK-8be7dca9-3_4

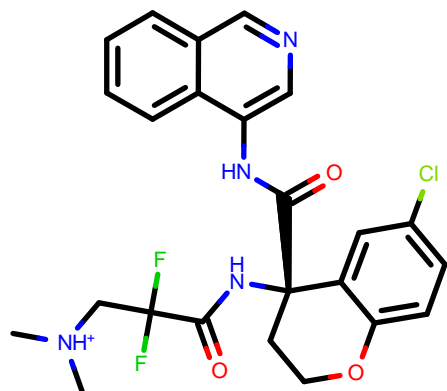
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H]3c4cc(ccc4OC[C@H]3C[NH3+])Cl

RUN: RUN498

DDG (kcal/mol): -1.99

dDDG (kcal/mol): 0.21

JAG-UCB-706446eb-5_1



CID: JAG-UCB-706446eb-5_1

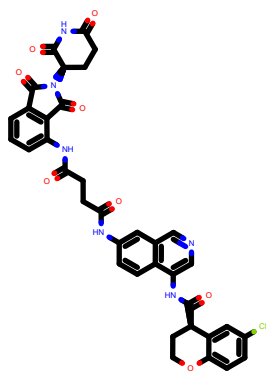
SMILES: COCCN([C@@H]1CCCOC1)C(=O)C[C@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4cnc5c4cccc5

RUN: RUN623

DDG (kcal/mol): -1.99

dDDG (kcal/mol): 0.28

ALP-UNI-8e43a71e-2_11



CID: ALP-UNI-8e43a71e-2_11

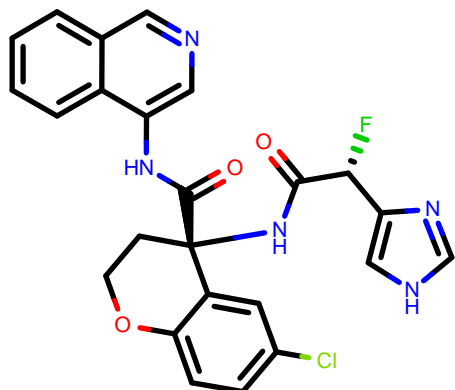
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@]3(COC(=O)C(=O)C)CC(=O)N5CC[C@H]6[C@@H]5OCCN6H+P(=O)(O)O

RUN: RUN2934

DDG (kcal/mol): -1.98

dDDG (kcal/mol): 0.21

MIC-UNK-91acba05-2_1



CID: MIC-UNK-91acba05-2_1

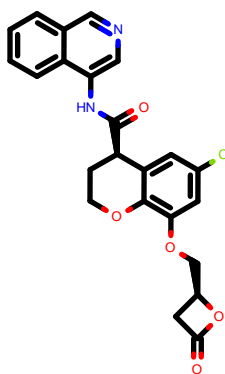
SMILES: CN1C[C@@H](c2cc(ccc2C1=O)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN469

DDG (kcal/mol): -1.97

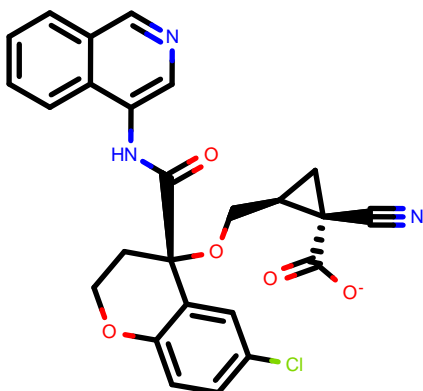
dDDG (kcal/mol): 0.29

MAT-POS-5d65ec79-1_1



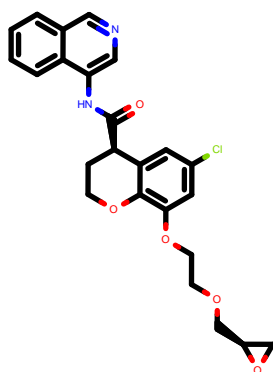
CID:	MAT-POS-5d65ec79-1_1
SMILES:	<chem>CN(C)C(=O)C[C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3ncc4c3cccc4</chem>
RUN:	RUN2080
DDG (kcal/mol):	-1.97
dDDG (kcal/mol):	0.41

VLA-UCB-50c39ae8-6_1



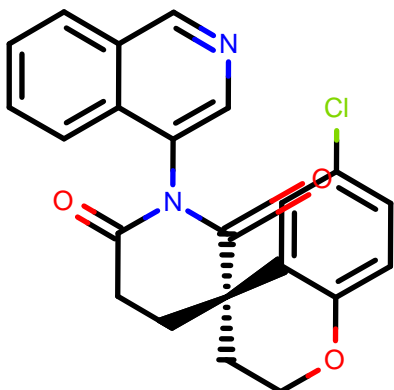
CID:	VLA-UCB-50c39ae8-6_1
SMILES:	<chem>c1ccc2c(c1)cncc2N(C(=O)CC)C(=O)[C@@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN385
DDG (kcal/mol):	-1.97
dDDG (kcal/mol):	0.30

MAT-POS-f9802937-7_1



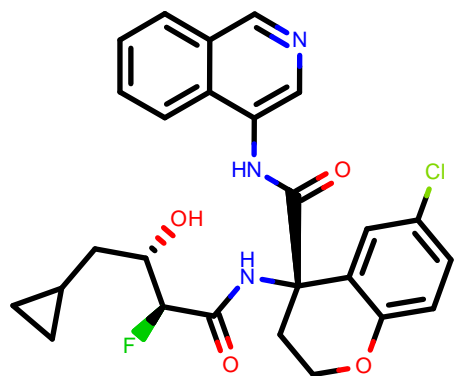
CID:	MAT-POS-f9802937-7_1
SMILES:	<chem>COc1ccc2cncc(c2c1)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN2398
DDG (kcal/mol):	-1.97
dDDG (kcal/mol):	0.42

BAR-COM-0f94fc3d-47_1



CID:	BAR-COM-0f94fc3d-47_1
SMILES:	<chem>Cc1ccc2c(c1)c(cc2)CC(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN8
DDG (kcal/mol):	-1.96
dDDG (kcal/mol):	0.18

MIC-UNK-5a93dd5f-3_15



CID: MIC-UNK-5a93dd5f-3_15

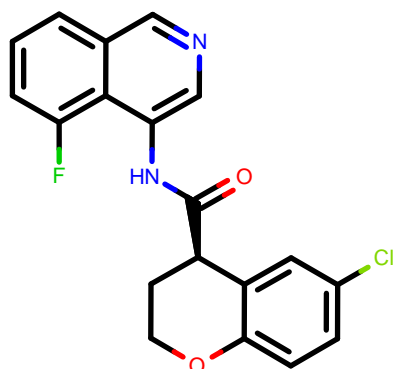
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H](c3cccc(c3)Cl)[N@H]4CC[C@H]5CCCC[C@H]5C4

RUN: RUN755

DDG (kcal/mol): -1.96

dDDG (kcal/mol): 0.33

VLA-UCB-05e51b3f-10_1



CID: VLA-UCB-05e51b3f-10_1

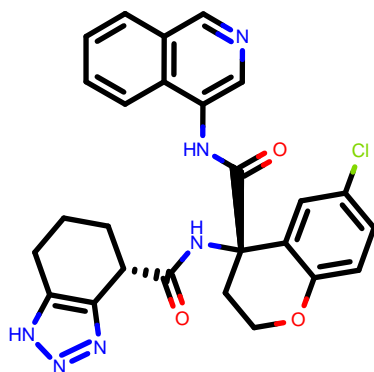
SMILES: C=CC(=O)N(c1cncc2c1cccc2)C(=O)[C@H]3CCOCc4c3cc(cc4)Cl

RUN: RUN318

DDG (kcal/mol): -1.95

dDDG (kcal/mol): 0.21

EDJ-MED-009f762b-5_2



CID: EDJ-MED-009f762b-5_2

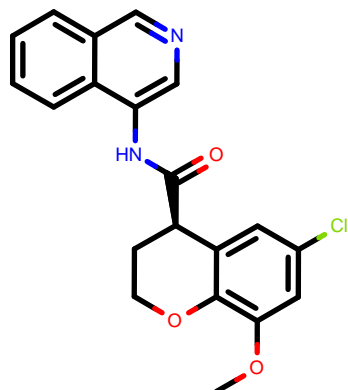
SMILES: CN(C)C(=O)C[N@H]1Cc2cccc(cc2)[C@@H](C1)C(=O)Nc3ncc4c3cc(cc4)F)Cl

RUN: RUN3915

DDG (kcal/mol): -1.95

dDDG (kcal/mol): 0.14

DAR-DIA-0cde14eb-52_1



CID: DAR-DIA-0cde14eb-52_1

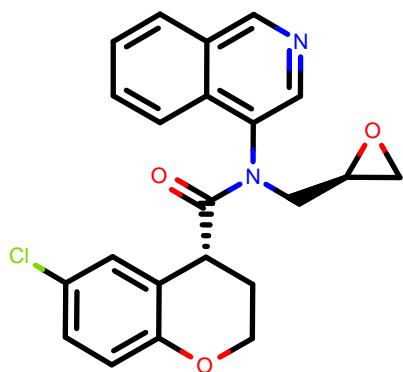
SMILES: c1ccc2c(c1)cncc2NC(=O)Nc3cccc(c3)C4(CC4)F

RUN: RUN18

DDG (kcal/mol): -1.95

dDDG (kcal/mol): 0.23

MAT-POS-4223bc15-11_16



CID: MAT-POS-4223bc15-11_16

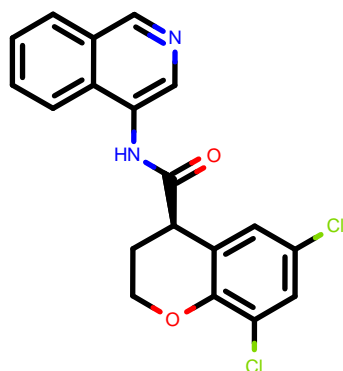
SMILES: C[C@H]1CC[N@](C1S(=O)(=O)[N@]2Cc3ccc(cc3[C@H](C2)C(=O)Nc4nccc5c4ccc5)Cl

RUN: RUN4043

DDG (kcal/mol): -1.95

dDDG (kcal/mol): 0.13

MAK-UNK-ffc90da7-2_3



CID: MAK-UNK-ffc90da7-2_3

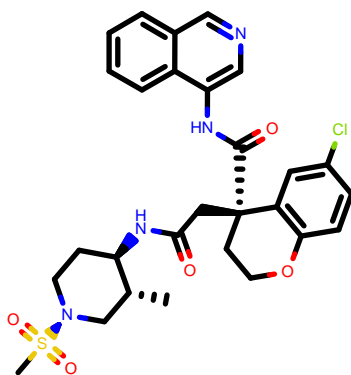
SMILES: c1ccc2c(c1)cncc2NC(=O)C[C@@H]3CCC[C@H]1[C@@H]3O1

RUN: RUN683

DDG (kcal/mol): -1.94

dDDG (kcal/mol): 0.19

LON-WEI-5e7d1b3e-6_1



CID: LON-WEI-5e7d1b3e-6_1

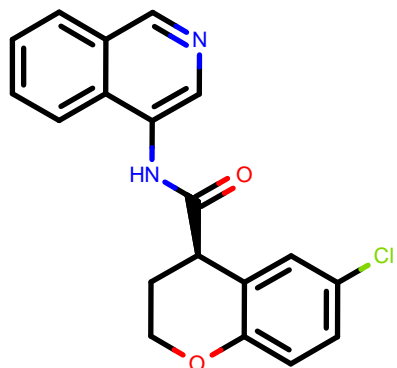
SMILES: CC(=O)c1cccc(c1)NC(=O)Nc2cn(c(=O)c3c2cccc3)C

RUN: RUN1312

DDG (kcal/mol): -1.94

dDDG (kcal/mol): 0.26

DAR-DIA-0cde14eb-53_1



CID: DAR-DIA-0cde14eb-53_1

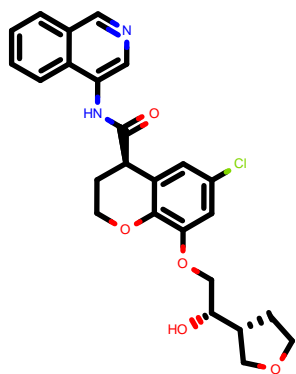
SMILES: c1ccc2c(c1)cncc2NC(=O)Nc3cccc(c3)C4(C4)I

RUN: RUN20

DDG (kcal/mol): -1.94

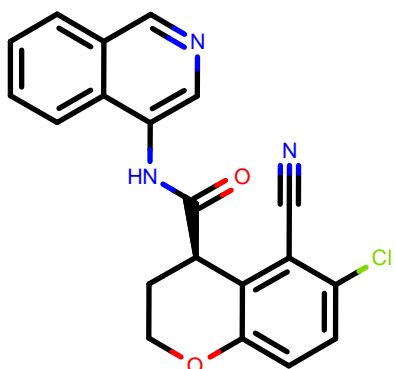
dDDG (kcal/mol): 0.21

ALP-UNI-0676e700-25_1



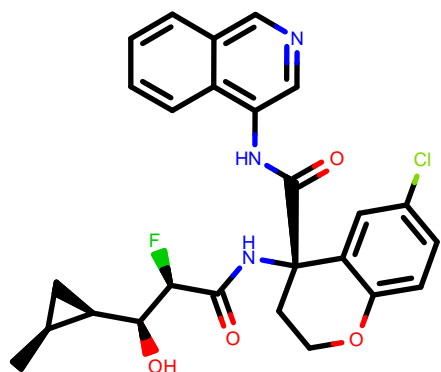
CID:	ALP-UNI-0676e700-25_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CNC(=O)c5cc6nccn6c(=O)[nH]5</chem>
RUN:	RUN2470
DDG (kcal/mol):	-1.93
dDDG (kcal/mol):	0.41

MAK-UNK-ffc90da7-2_7



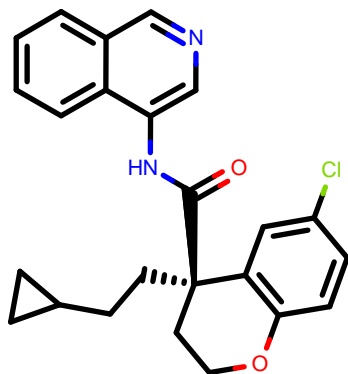
CID:	MAK-UNK-ffc90da7-2_7
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)C[C@@H]3CCCC[C@H]1([C@H]3O)Cl</chem>
RUN:	RUN692
DDG (kcal/mol):	-1.93
dDDG (kcal/mol):	0.20

EDJ-MED-6d9ff7d0-7_1



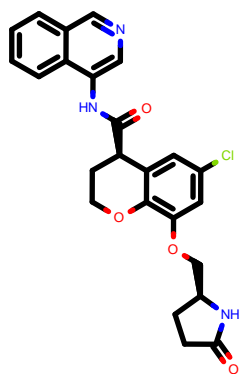
CID:	EDJ-MED-6d9ff7d0-7_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)N[C@@H](C)C(=O)N5</chem>
RUN:	RUN3431
DDG (kcal/mol):	-1.93
dDDG (kcal/mol):	0.14

MAR-TRE-04c86cea-39_1



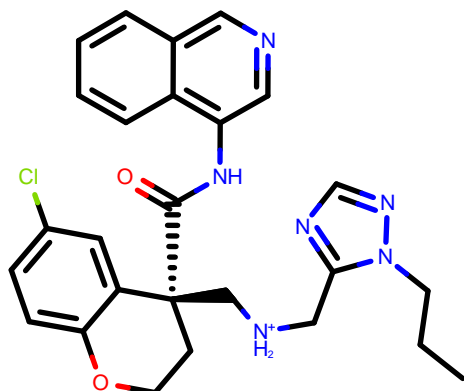
CID:	MAR-TRE-04c86cea-39_1
SMILES:	<chem>COc1ccccc1n2cc(c3ccccc3c2=O)C(=O)Nc4ccc(nc4)OC</chem>
RUN:	RUN52
DDG (kcal/mol):	-1.93
dDDG (kcal/mol):	0.24

EDG-MED-5d232de5-1_1



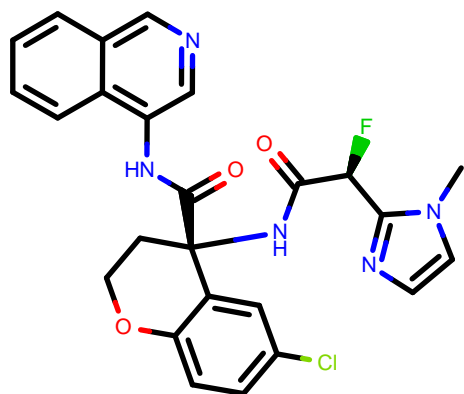
CID:	EDG-MED-5d232de5-1_1
SMILES:	<chem>c1cc2cncc(c2cc1F)NC(=O)[C@@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN2363
DDG (kcal/mol):	-1.92
dDDG (kcal/mol):	0.39

EDJ-MED-1b5395f9-3_1



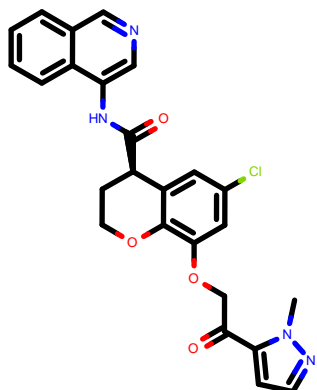
CID:	EDJ-MED-1b5395f9-3_1
SMILES:	<chem>CCc1ccc2cncc(c2c1)NC(=O)[C@H]3C[N@@]([C@@H]3CCOCc4c3cc(cc4)Cl)S(=O)(=O)C</chem>
RUN:	RUN4465
DDG (kcal/mol):	-1.92
dDDG (kcal/mol):	0.35

MIC-UNK-91acba05-1_1



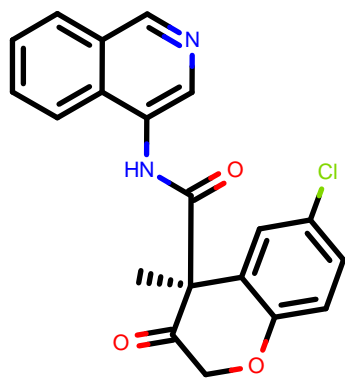
CID:	MIC-UNK-91acba05-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CNC(=O)c4c3cc(cc4)Cl</chem>
RUN:	RUN4466
DDG (kcal/mol):	-1.92
dDDG (kcal/mol):	0.31

EDJ-MED-d203f206-37_1



CID:	EDJ-MED-d203f206-37_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCS(=O)CC5</chem>
RUN:	RUN2599
DDG (kcal/mol):	-1.91
dDDG (kcal/mol):	0.31

DAR-DIA-23aa0b97-19_1



CID: DAR-DIA-23aa0b97-19_1

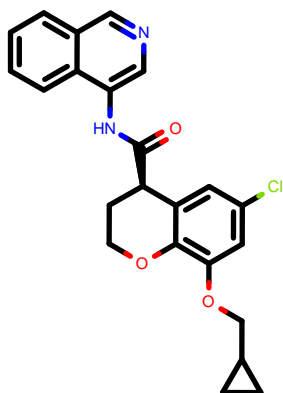
SMILES: c1ccc2c(c1)cncc2CC(=O)Nc3ccccc(c3)C#N

RUN: RUN1

DDG (kcal/mol): -1.91

dDDG (kcal/mol): 0.22

ED_-GRI-5b13fbe2-11_1



CID: ED_-GRI-5b13fbe2-11_1

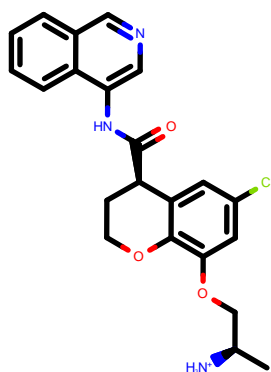
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(C)CCOC4c3cc(cc4)C1OCCN5OC(NH+)[C]C5)C1(=O)N

RUN: RUN1541

DDG (kcal/mol): -1.91

dDDG (kcal/mol): 0.29

MAT-POS-e9e99895-9_2



CID: MAT-POS-e9e99895-9_2

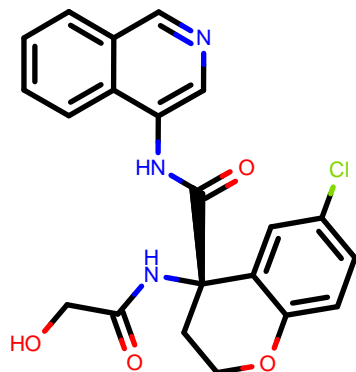
SMILES: C[C@@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2nccc3c2ccc3)NC(=O)COC4ccc(cc4)C1(=O)N

RUN: RUN2262

DDG (kcal/mol): -1.90

dDDG (kcal/mol): 0.41

MIC-UNK-cdc2493e-12_2



CID: MIC-UNK-cdc2493e-12_2

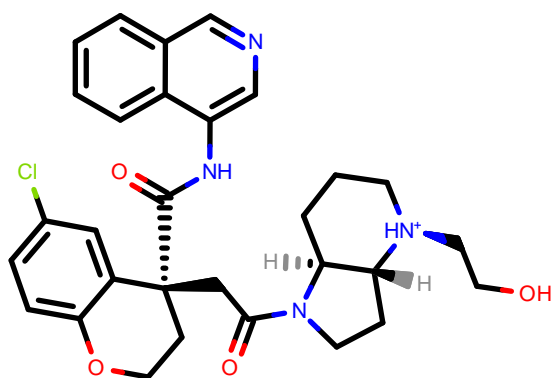
SMILES: C[NH+](C)[C@H]1CC[C@@H](C1)N(c2cccc(c2)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN549

DDG (kcal/mol): -1.90

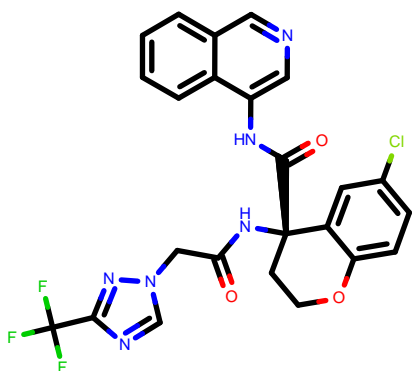
dDDG (kcal/mol): 0.25

RAL-THA-8416115c-11_2



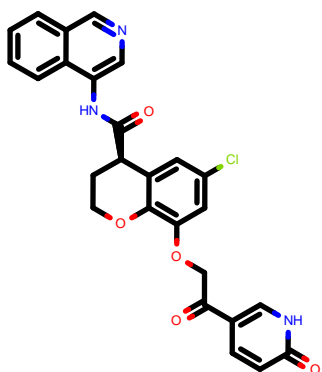
CID:	RAL-THA-8416115c-11_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@H]3CCN(c4c3cc(cc4)Cl)C5cnc[nH]5</chem>
RUN:	RUN1288
DDG (kcal/mol):	-1.90
dDDG (kcal/mol):	0.55

RAL-THA-8416115c-1_2



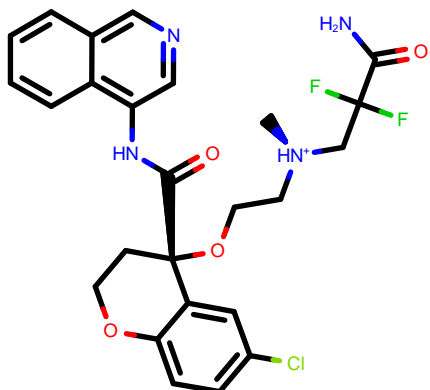
CID:	RAL-THA-8416115c-1_2
SMILES:	<chem>c1ccc(cc1)CN2CC[C@H](c3c2ccc(c3)Cl)C(=O)Nc4ncnc5c4cccc5</chem>
RUN:	RUN1248
DDG (kcal/mol):	-1.90
dDDG (kcal/mol):	0.34

LEE-CAM-7ab9b158-4_1



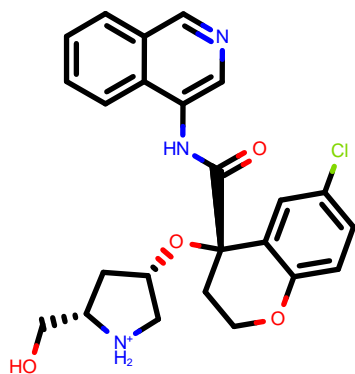
CID:	LEE-CAM-7ab9b158-4_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@H]3[C@@H](COC4=C3C=CC(=O)N4)COC[C@H]5CN6C[C@H](O5)COC6=O</chem>
RUN:	RUN2209
DDG (kcal/mol):	-1.89
dDDG (kcal/mol):	0.35

ERI-UCB-b3e6b0c2-18_1



CID:	ERI-UCB-b3e6b0c2-18_1
SMILES:	<chem>c1cc2c(c1)CN3CC[NH2+]CC3)cnc2N4CC[C@H]5(C4=O)CO6c5cc(cc6)Cl</chem>
RUN:	RUN3052
DDG (kcal/mol):	-1.89
dDDG (kcal/mol):	0.28

MIC-UNK-91acba05-2_2



CID: MIC-UNK-91acba05-2_2

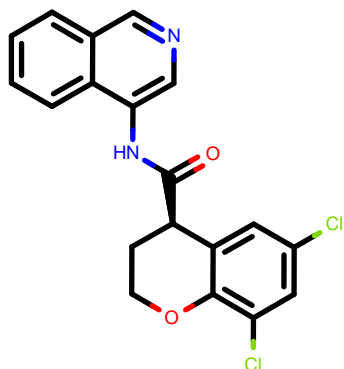
SMILES: CN1C[C@H](c2cc(ccc2C1=O)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN470

DDG (kcal/mol): -1.89

dDDG (kcal/mol): 0.36

MIC-UNK-5a93dd5f-12_5



CID: MIC-UNK-5a93dd5f-12_5

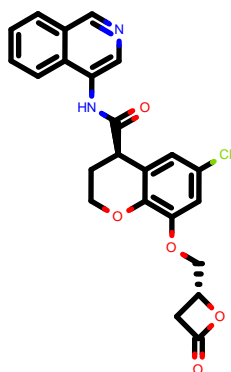
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H](c3cccc(c3)Cl)N@@@4CC[C@H](C4)NH+5CCCC5

RUN: RUN794

DDG (kcal/mol): -1.86

dDDG (kcal/mol): 0.20

LEE-CAM-7ab9b158-2_1



CID: LEE-CAM-7ab9b158-2_1

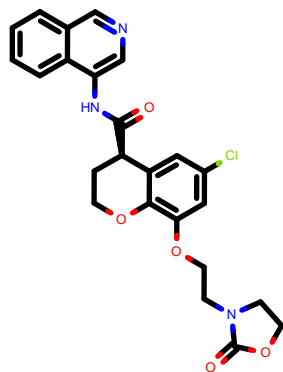
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H](c3(CCOc4c3cc(cc4)Cl)COCc5[nH]c(=O)c6c(n5)CCOC6

RUN: RUN2207

DDG (kcal/mol): -1.86

dDDG (kcal/mol): 0.37

ALP-POS-5bb456a5-2_2



CID: ALP-POS-5bb456a5-2_2

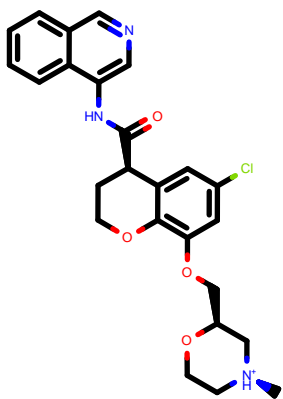
SMILES: C[C@@H]1C[N@@H]CC[C@@H]1NC1=O[C@@H]2(COCc3c2cc(cc3)Cl)C1=O)Nc4nc5c4cccc5)S(=O)(=O)C

RUN: RUN2424

DDG (kcal/mol): -1.86

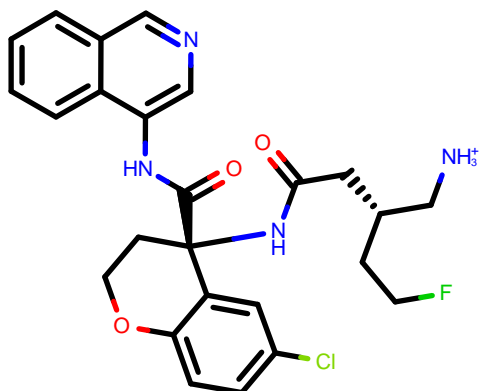
dDDG (kcal/mol): 0.41

MAT-POS-fce787c2-9_1



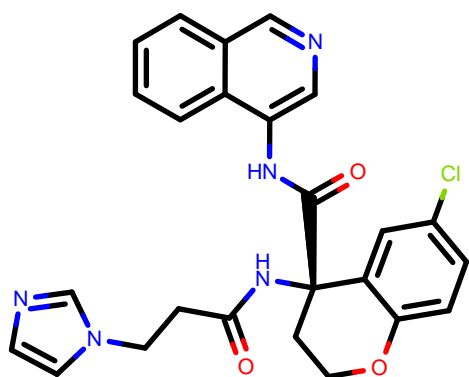
CID:	MAT-POS-fce787c2-9_1
SMILES:	<chem>C[C@@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2cncc3c2cccc3)[NH2+]</chem> C
RUN:	RUN2156
DDG (kcal/mol):	-1.85
dDDG (kcal/mol):	0.44

ADA-UCB-dc2b944c-4_1



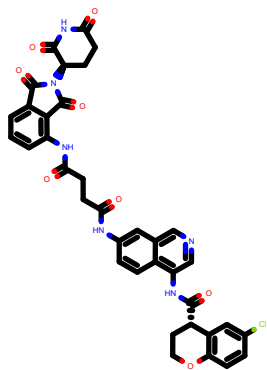
CID:	ADA-UCB-dc2b944c-4_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]34C[C@H]3COc5c4cc(cc5)Cl</chem>
RUN:	RUN600
DDG (kcal/mol):	-1.83
dDDG (kcal/mol):	0.36

NAU-LAT-a5c7d7cb-5_2



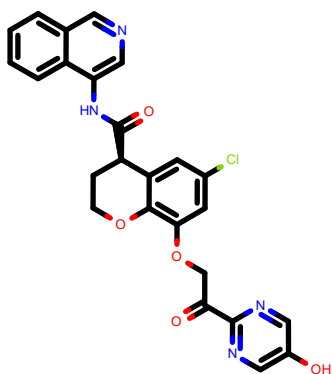
CID:	NAU-LAT-a5c7d7cb-5_2
SMILES:	<chem>CC(=O)N1CC[NH+](CC1)[C@H](c2ccc(cc2)N(C)C)C(=O)Cc3cncc4c3cccc4</chem>
RUN:	RUN580
DDG (kcal/mol):	-1.83
dDDG (kcal/mol):	0.24

VLA-UNK-9a7dc93f-5_1



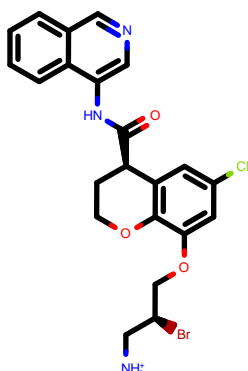
CID:	VLA-UNK-9a7dc93f-5_1
SMILES:	<chem>CO[C@@H](c1cc(c(c(c1)Cl)F)F)C(=O)Nc2cncc3c2cccc3</chem>
RUN:	RUN3085
DDG (kcal/mol):	-1.83
dDDG (kcal/mol):	0.45

EDJ-MED-6864a934-2_1



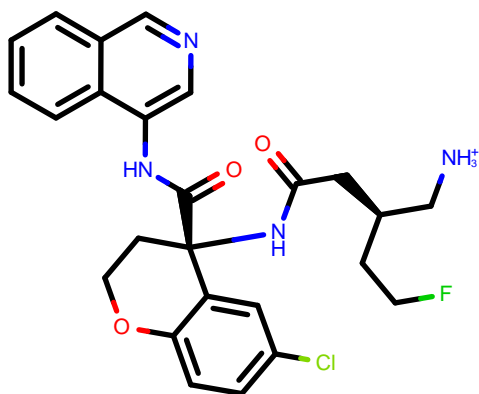
CID:	EDJ-MED-6864a934-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCCOc4c3cc(cc4)C)NC(=O)c5cc6n(n5)CCO6</chem>
RUN:	RUN2606
DDG (kcal/mol):	-1.83
dDDG (kcal/mol):	0.39

MAT-POS-e9e99895-5_2



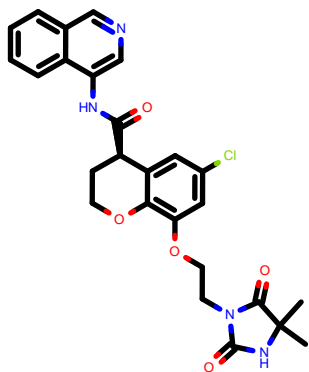
CID:	MAT-POS-e9e99895-5_2
SMILES:	<chem>C[C@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2cncc3c2ccc3)NC(=O)c4cc5n(n4)CCO5</chem>
RUN:	RUN2250
DDG (kcal/mol):	-1.83
dDDG (kcal/mol):	0.44

ADA-UCB-dc2b944c-17_1



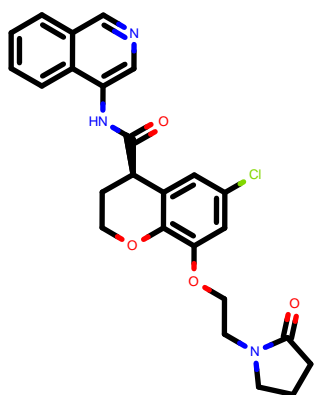
CID:	ADA-UCB-dc2b944c-17_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)N(CCC3CCCCC3)c4cc(ccc4Cl)Cl</chem>
RUN:	RUN615
DDG (kcal/mol):	-1.82
dDDG (kcal/mol):	0.28

MAT-POS-e9e99895-8_2



CID:	MAT-POS-e9e99895-8_2
SMILES:	<chem>Cc1nc2n(n1)C[C@H](CC2)C(=O)N[C@@](C)(C)c3cc(c(c3)Cl)C(=O)Nc4ncc5c4ccc5</chem>
RUN:	RUN2252
DDG (kcal/mol):	-1.82
dDDG (kcal/mol):	0.37

ALP-UNI-0676e700-20_1



CID: ALP-UNI-0676e700-20_1

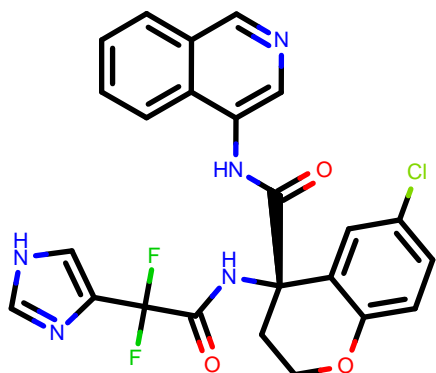
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)C)CNC(=O)c5ccc(cn5)n6cnnn6

RUN: RUN2467

DDG (kcal/mol): -1.82

dDDG (kcal/mol): 0.42

NAU-LAT-a5c7d7cb-2_2



CID: NAU-LAT-a5c7d7cb-2_2

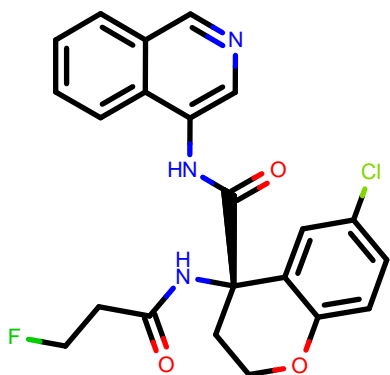
SMILES: CC(=O)N1CC[NH+](CC1)[C@H](c2cccc(c2)Cl)C(=O)Nc3ncc4c3cccc4

RUN: RUN582

DDG (kcal/mol): -1.81

dDDG (kcal/mol): 0.33

MAK-UNK-ffc90da7-7_1



CID: MAK-UNK-ffc90da7-7_1

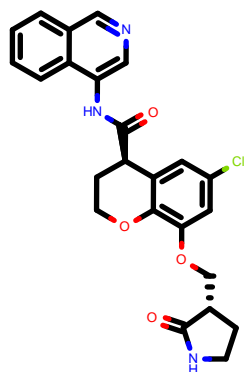
SMILES: CC(C)OC[C@@@H]([C@@@H](c1cccc(c1)Cl)C(=O)Nc2ncc3c2cccc3)[NH2+]C

RUN: RUN707

DDG (kcal/mol): -1.80

dDDG (kcal/mol): 0.27

EDG-MED-5d232de5-5_1



CID: EDG-MED-5d232de5-5_1

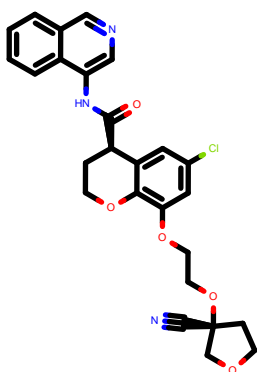
SMILES: CN1CC[C@@@H](c2c1ccc(c2)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN2367

DDG (kcal/mol): -1.80

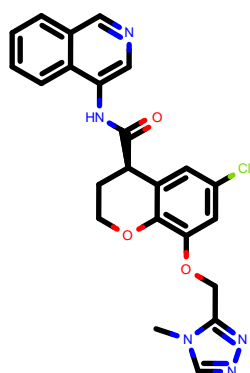
dDDG (kcal/mol): 0.41

EDJ-MED-d203f206-4_1



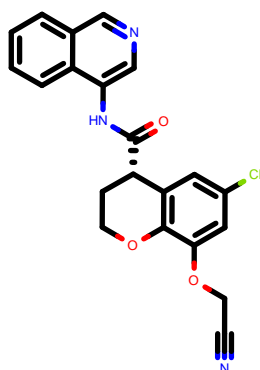
CID:	EDJ-MED-d203f206-4_1
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)C@H3[C@@O]c4ccccc4C1CC(O)N[C@@H]5C[C@@H]6[C@@H]7C[C@@H]8C[C@@H]9</chem>
RUN:	RUN2566
DDG (kcal/mol):	-1.80
dDDG (kcal/mol):	0.58

ALP-POS-5bb456a5-1_3



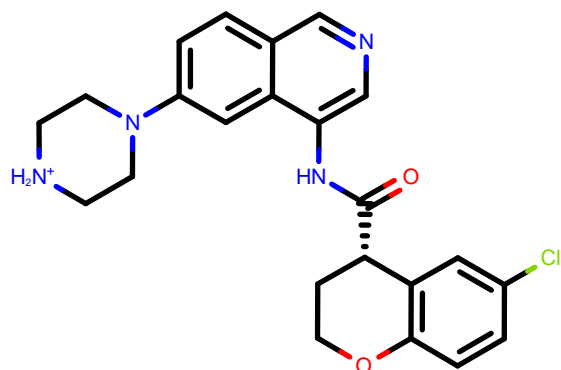
CID:	ALP-POS-5bb456a5-1_3
SMILES:	<chem>C1C@H]2CN@H]3CC[C@@H]1NC(=O)C[C@@H]2[C@@O]c3ccccc3C1C(O)N4c5cccnc5S(=O)(=O)C</chem>
RUN:	RUN2409
DDG (kcal/mol):	-1.79
dDDG (kcal/mol):	0.39

MAT-POS-e9e99895-2_4



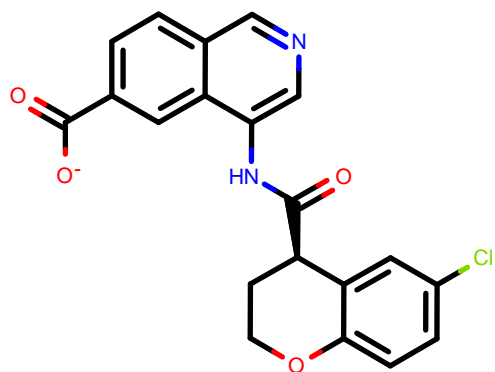
CID:	MAT-POS-e9e99895-2_4
SMILES:	<chem>CC(C)[N@H+]1CCO[C@H](C1)C(=O)N[C@@H]2[C@@O]c3ccccc3C1C(O)Nc3cccnc3</chem>
RUN:	RUN2237
DDG (kcal/mol):	-1.79
dDDG (kcal/mol):	0.19

MAT-POS-2492181e-7_1



CID:	MAT-POS-2492181e-7_1
SMILES:	<chem>CC1CC[NH+](CC1)CCNC(=O)Nc2cn(c(=O)c3c2cccc3)CC(C)C</chem>
RUN:	RUN100
DDG (kcal/mol):	-1.79
dDDG (kcal/mol):	0.13

LON-WEI-4d77710c-58_1



CID: LON-WEI-4d77710c-58_1

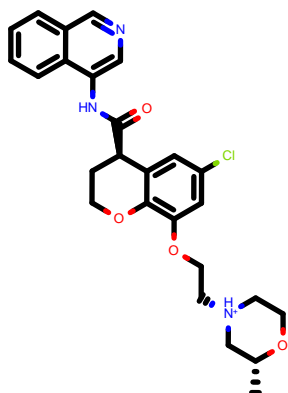
SMILES: CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)Nc3nc4ccc(cc4s3)OC

RUN: RUN250

DDG (kcal/mol): -1.79

dDDG (kcal/mol): 0.19

ALP-UNI-3496895b-15_4



CID: ALP-UNI-3496895b-15_4

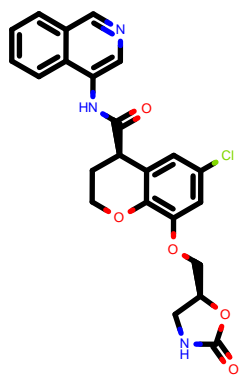
SMILES: c1ccc2c(c1)ncnc2NC(=O)N[C@@H]3CCOC4=CC=C(C=C3)C[C@@H]5C[C@@H]6C[C@@H]7C[C@H]8C[C@H]9C[C@H]8C[C@H]9N

RUN: RUN2534

DDG (kcal/mol): -1.79

dDDG (kcal/mol): 0.40

EDJ-MED-d203f206-39_1



CID: EDJ-MED-d203f206-39_1

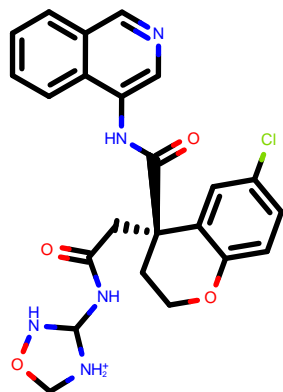
SMILES: c1ccc2c(c1)ncnc2NC(=O)N[C@@H]3CCOC4=CC=C(C=C3)C[C@@H]5C[C@@H]6C[C@@H]7C[C@H]8C[C@H]9C[C@H]8C[C@H]9N

RUN: RUN2601

DDG (kcal/mol): -1.79

dDDG (kcal/mol): 0.44

MAK-UNK-c749d764-22_5



CID: MAK-UNK-c749d764-22_5

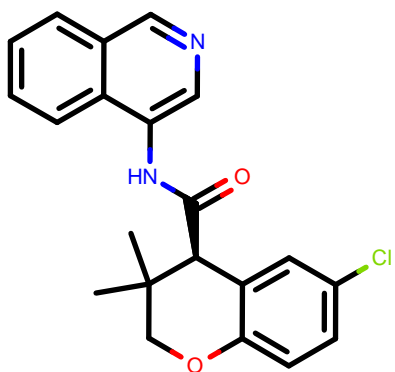
SMILES: C[C@@H](C)[C@H]2NC(=O)N[C@@H]3CCOC1=C[C@@H](C2)C[C@@H]4C[C@@H]5C[C@@H]4C[C@@H]5N

RUN: RUN1026

DDG (kcal/mol): -1.78

dDDG (kcal/mol): 0.27

MAT-POS-3b92565d-1_1



CID: MAT-POS-3b92565d-1_1

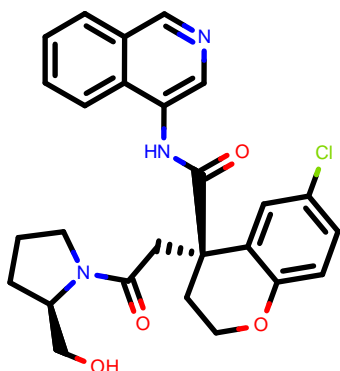
SMILES: COc1ccc(cc1CC(=O)Nc2cncc3c2cccc3)Cl

RUN: RUN106

DDG (kcal/mol): -1.78

dDDG (kcal/mol): 0.19

MAK-UNK-c749d764-33_5



CID: MAK-UNK-c749d764-33_5

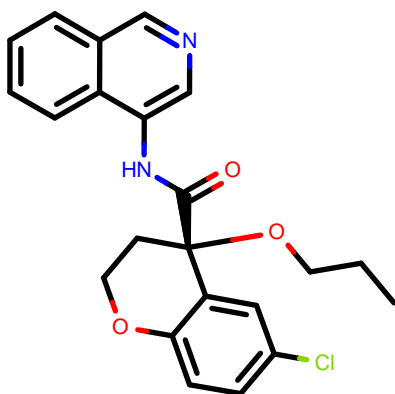
SMILES: CS(=O)(=O)N(c1cncc2c1cccc2)C(=O)C[C@@H]3CCC[C@@H]4[C@H]3O[C@H]4(F)F

RUN: RUN1084

DDG (kcal/mol): -1.78

dDDG (kcal/mol): 0.29

DAR-DIA-9e4459de-13_1



CID: DAR-DIA-9e4459de-13_1

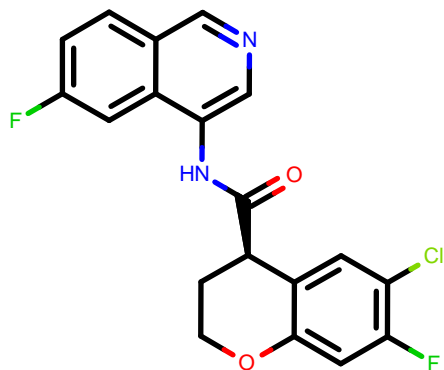
SMILES: c1cc2c(c1)NCCCC3ccc4c(c3)cncc4N(C)C[C@@H]5CCOC6c5c(c6)Clc7c(c2O)[C@@H]7C(C)C1=O)NC7=O)O

RUN: RUN1425

DDG (kcal/mol): -1.78

dDDG (kcal/mol): 0.23

ED_-GRI-5b13fbe2-74_2



CID: ED_-GRI-5b13fbe2-74_2

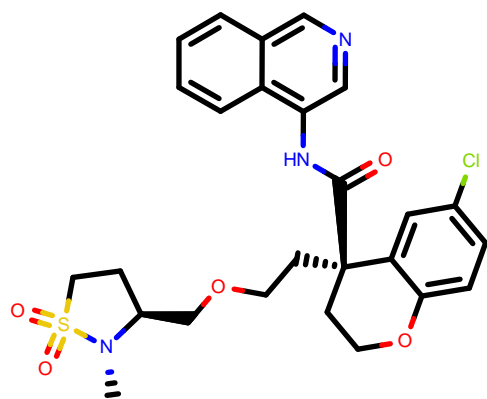
SMILES: C[NH+]C(C)C[C@@H](C1)2nccc3c2c(C)C[C@@H]3COC4c3cc(c4)Cl)C1=O)Nc5cncc6c5cccc6

RUN: RUN1632

DDG (kcal/mol): -1.77

dDDG (kcal/mol): 0.21

EDJ-MED-6d9ff7d0-1_1



CID: EDJ-MED-6d9ff7d0-1_1

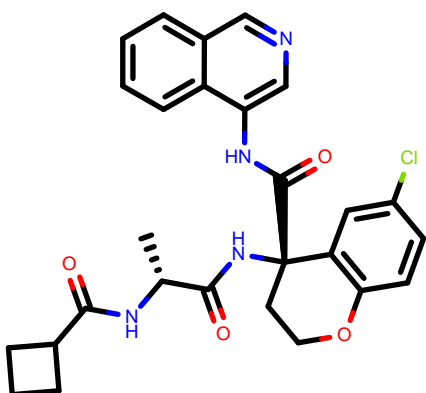
SMILES: COC(C)C[NH2+][C@@]1(COC2C1CC(C)C)C(=O)Nc3ccc4c3cccc4

RUN: RUN3427

DDG (kcal/mol): -1.77

dDDG (kcal/mol): 0.19

RAL-THA-8416115c-12_1



CID: RAL-THA-8416115c-12_1

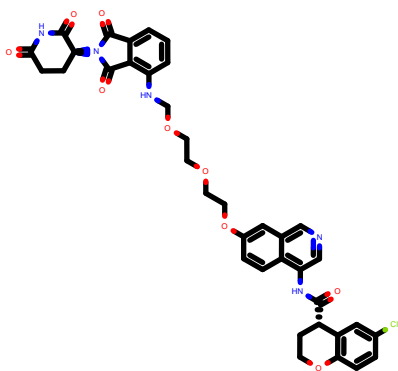
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]([H])3CCN(c4c3cc(cc4)Cl)Cc5[n-]nn5

RUN: RUN1291

DDG (kcal/mol): -1.77

dDDG (kcal/mol): 0.30

DAR-DIA-6a508060-1_2



CID: DAR-DIA-6a508060-1_2

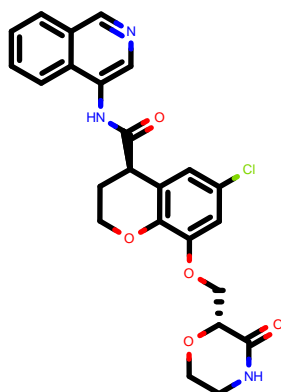
SMILES: c1ccc2c(c1)cncc2CC(=O)[C@@]([H])3CCOCc4c3cc(cc4)Cl

RUN: RUN333

DDG (kcal/mol): -1.77

dDDG (kcal/mol): 0.51

ALP-POS-5bb456a5-2_1



CID: ALP-POS-5bb456a5-2_1

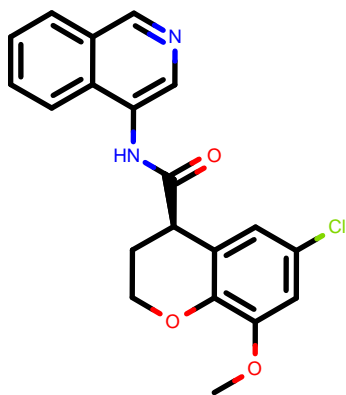
SMILES: C[C@@]([H])1CN[C@@]([H])2C(=O)C[C@@]([H])3CCOCc4c3cc(cc4)Cl

RUN: RUN2422

DDG (kcal/mol): -1.77

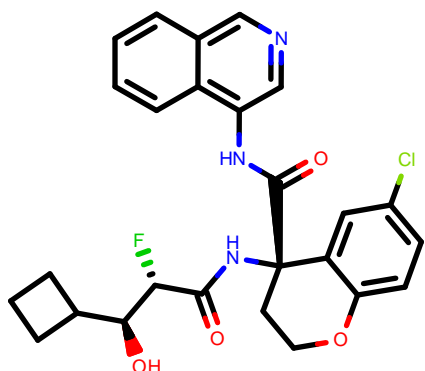
dDDG (kcal/mol): 0.41

MAK-UNK-ffc90da7-7_2



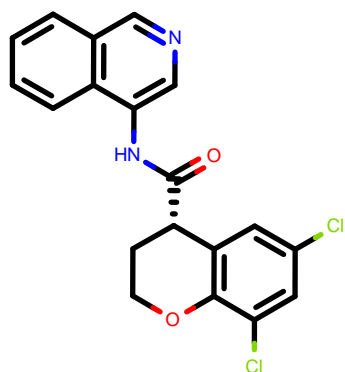
CID:	MAK-UNK-ffc90da7-7_2
SMILES:	<chem>CC(C)OC[C@H]([C@@H](c1cccc(c1)Cl)C(=O)Nc2nccc3c2cccc3)[NH2+]C</chem>
RUN:	RUN708
DDG (kcal/mol):	-1.76
dDDG (kcal/mol):	0.21

ALF-EVA-b701bd13-6_2



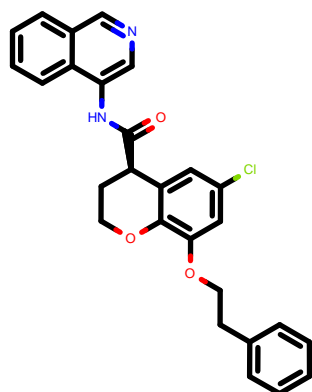
CID:	ALF-EVA-b701bd13-6_2
SMILES:	<chem>c1ccc2c(c1)nccc2NC(=O)[C@H]3CCCC(=O)Nc4c3cc(c(c4)Cl)Cl</chem>
RUN:	RUN3641
DDG (kcal/mol):	-1.76
dDDG (kcal/mol):	0.17

DAR-DIA-0d514e7d-1_1



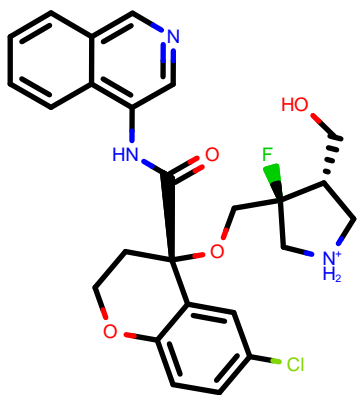
CID:	DAR-DIA-0d514e7d-1_1
SMILES:	<chem>C[C@H]1COc2c(cc2OC)Cl][C@@H]1C(=O)Nc3nccc4c3cccc4</chem>
RUN:	RUN804
DDG (kcal/mol):	-1.76
dDDG (kcal/mol):	0.11

ED_-GRI-5b13fbe2-50_1



CID:	ED_-GRI-5b13fbe2-50_1
SMILES:	<chem>c1ccc2c(c1)nccc2NC(=O)[C@@H]3[C@@H](COc4ccccc4)OC[C@@H]3C(=O)Nc5c3cc(c(c5)O)O</chem>
RUN:	RUN1587
DDG (kcal/mol):	-1.76
dDDG (kcal/mol):	0.50

VLA-UNK-4b5c0188-1_1



CID: VLA-UNK-4b5c0188-1_1

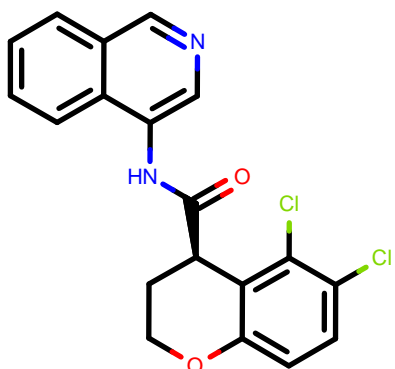
SMILES: c1ccc2c(c1)cncc2N(C[C@@H]3CO3)C(=O)[C@@H]4CCOC5c4cc(cc5)Cl

RUN: RUN3074

DDG (kcal/mol): -1.76

dDDG (kcal/mol): 0.23

MAK-UNK-ffc90da7-2_1



CID: MAK-UNK-ffc90da7-2_1

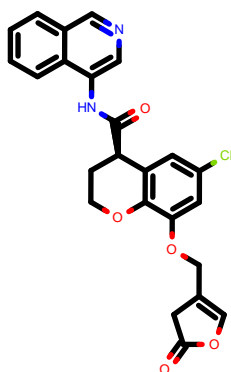
SMILES: c1ccc2c(c1)cncc2NC(=O)C[C@@H]3CCCC[C@@H]4[C@@H]3O4Cl

RUN: RUN681

DDG (kcal/mol): -1.75

dDDG (kcal/mol): 0.18

EDG-MED-5d232de5-6_1



CID: EDG-MED-5d232de5-6_1

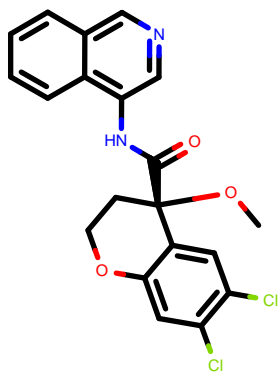
SMILES: CN1CC[C@H]1(c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN2371

DDG (kcal/mol): -1.75

dDDG (kcal/mol): 0.34

MAT-POS-23a8a11a-1_1



CID: MAT-POS-23a8a11a-1_1

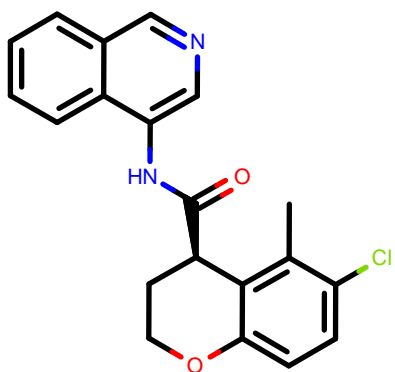
SMILES: c1ccc2c(c1)cncc2NC(=O)Cc3ccc(c(c3)Cl)Cl

RUN: RUN1406

DDG (kcal/mol): -1.75

dDDG (kcal/mol): 0.23

MAK-UNK-ffc90da7-2_4



CID: MAK-UNK-ffc90da7-2_4

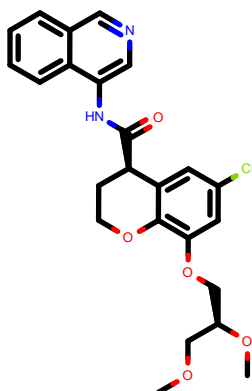
SMILES: c1ccc2c(c1)cncc2NC(=O)C[C@H]3CCCC[C@H]1([C@@H]3O)Cl

RUN: RUN684

DDG (kcal/mol): -1.75

dDDG (kcal/mol): 0.17

KAD-UNI-8a629cb0-8_1



CID: KAD-UNI-8a629cb0-8_1

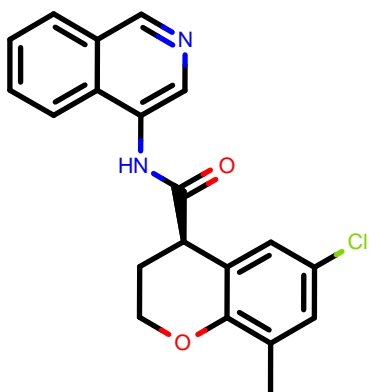
SMILES: C[C@H](Cl)C(O)N[NH+]CCN(C)C1C(O)C[C@H]2(C)CCO3C2C(C)C1(O)N4CCCC4

RUN: RUN2092

DDG (kcal/mol): -1.74

dDDG (kcal/mol): 0.50

MIC-UNK-50cce87d-9_2



CID: MIC-UNK-50cce87d-9_2

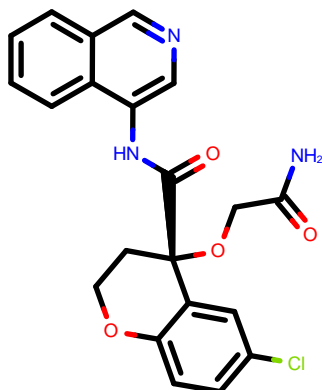
SMILES: c1cc(cc(c1)Cl)[C@H]2CCCN(C2=O)c3cncc4c3c(cc4)F

RUN: RUN675

DDG (kcal/mol): -1.74

dDDG (kcal/mol): 0.18

RAL-THA-1d44ff04-1_1



CID: RAL-THA-1d44ff04-1_1

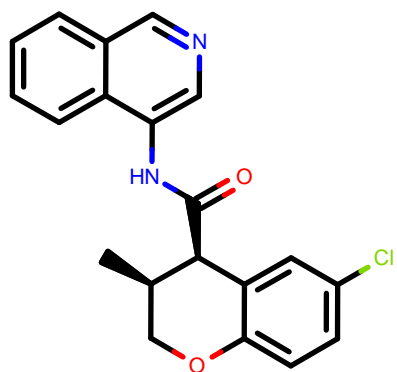
SMILES: c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)OCc4[nH]ccn4

RUN: RUN434

DDG (kcal/mol): -1.74

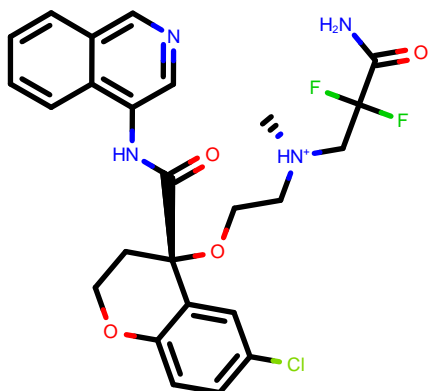
dDDG (kcal/mol): 0.20

ERI-UCB-ce40166b-8_1



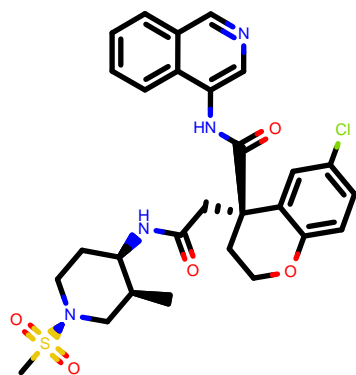
CID:	ERI-UCB-ce40166b-8_1
SMILES:	<chem>c1ccc2c(c1)cncc2CC(=O)Nc3cc(cc(c3)Cl)Oc4ccncc4</chem>
RUN:	RUN46
DDG (kcal/mol):	-1.74
dDDG (kcal/mol):	0.18

ERI-UCB-b3e6b0c2-5_1



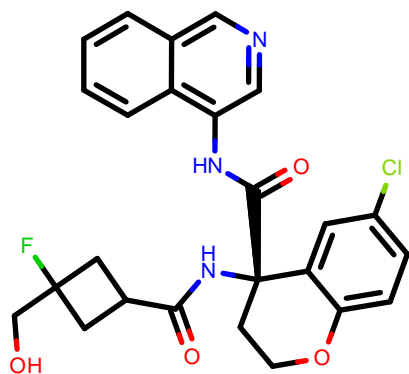
CID:	ERI-UCB-b3e6b0c2-5_1
SMILES:	<chem>c1cc2nccc(c2cc1CN3CC[NH2+][CC3]NC(=O)[C@@H]4CNc5c4cc(cc5)Cl</chem>
RUN:	RUN3037
DDG (kcal/mol):	-1.74
dDDG (kcal/mol):	0.21

MAK-UNK-c749d764-12_6



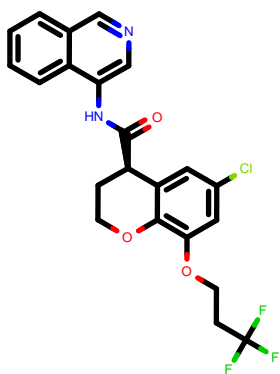
CID:	MAK-UNK-c749d764-12_6
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)C[C@@H]3CCC[C@@H]1([C@H]3O)Cl</chem>
RUN:	RUN957
DDG (kcal/mol):	-1.73
dDDG (kcal/mol):	0.40

MAT-POS-78e1d523-2_2



CID:	MAT-POS-78e1d523-2_2
SMILES:	<chem>CNC(=O)[C@@]1(CCOc2c1cc(cc2)Cl)CC(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3281
DDG (kcal/mol):	-1.73
dDDG (kcal/mol):	0.14

DAR-DIA-9e4459de-15_13



CID: DAR-DIA-9e4459de-15_13

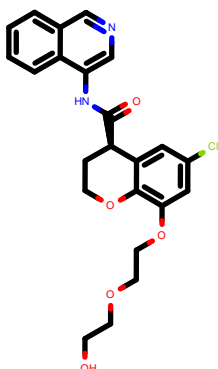
SMILES: c1cc2c(c1)NC(=O)CCC(=O)Nc3ccc4c(c3)ncoc4NC(=O)C@H]3CCOCc5cc(oc6)Clc1c(c2O)C@H]7CC(=O)NC7=O)O

RUN: RUN1452

DDG (kcal/mol): -1.73

dDDG (kcal/mol): 0.36

EDJ-MED-6864a934-10_1



CID: EDJ-MED-6864a934-10_1

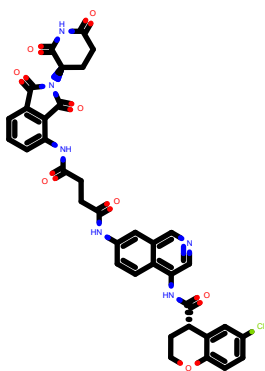
SMILES: c1ccc2c(c1)ncoc2NC(=O)C[C@@]3(CCOc4c3cc(oc4)Cl)NC(=O)C@H]C5(CCCO)O)O

RUN: RUN2614

DDG (kcal/mol): -1.73

dDDG (kcal/mol): 0.31

VLA-UNK-db5e3064-2_2



CID: VLA-UNK-db5e3064-2_2

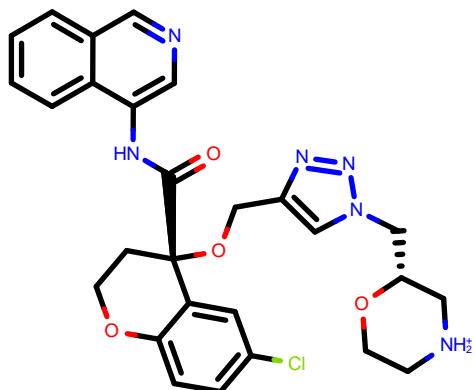
SMILES: c1ccc2c(c1)ncoc2n3c(c(oc3=O)c4cccc(c4)Cl)[O-]

RUN: RUN3096

DDG (kcal/mol): -1.73

dDDG (kcal/mol): 0.38

DAR-DIA-23e5a6a0-2_2



CID: DAR-DIA-23e5a6a0-2_2

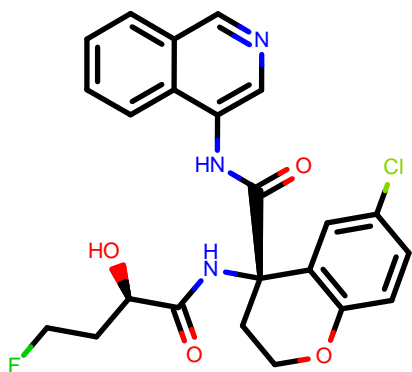
SMILES: c1ccc2c(c1)ncoc2NC(=O)C[C@@]3(CCOc4c3cc(oc4O)C@H]5CC6([NH2+][5]C)CC6)Cl

RUN: RUN404

DDG (kcal/mol): -1.72

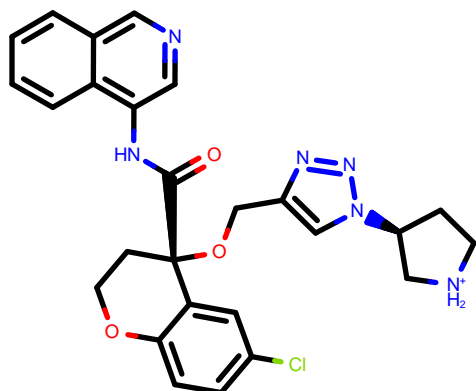
dDDG (kcal/mol): 0.32

MIC-UNK-cdc2493e-23_1



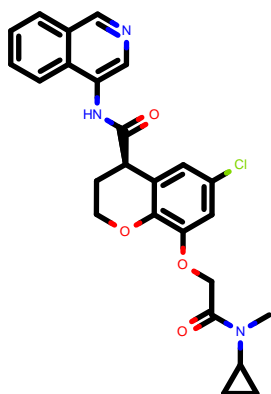
CID:	MIC-UNK-cdc2493e-23_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)N(CCn3cnnc3)c4cccc(c4)Cl</chem>
RUN:	RUN573
DDG (kcal/mol):	-1.72
dDDG (kcal/mol):	0.28

ROB-UNI-322e8f70-3_2



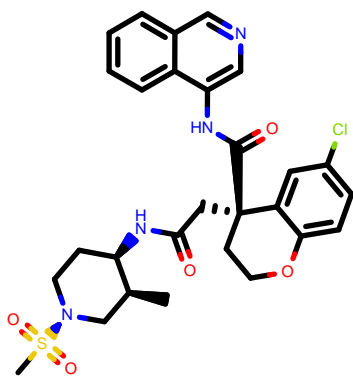
CID:	ROB-UNI-322e8f70-3_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CNC(=O)c4c3cc(cc4)Cl</chem>
RUN:	RUN3155
DDG (kcal/mol):	-1.72
dDDG (kcal/mol):	0.17

EDJ-MED-6864a934-9_1



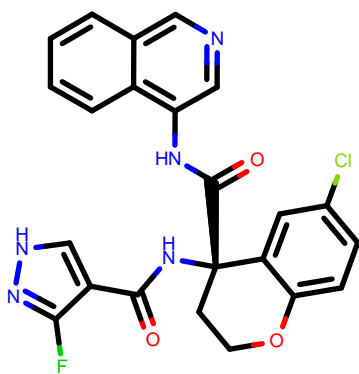
CID:	EDJ-MED-6864a934-9_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(C)COCc4c3cc(cc4)CNC(=O)CSc5cnc(n5)CC#N</chem>
RUN:	RUN2615
DDG (kcal/mol):	-1.71
dDDG (kcal/mol):	0.37

ALP-POS-696356e4-1_2



CID:	ALP-POS-696356e4-1_2
SMILES:	<chem>c1cc2cncc(c2cc1F)NC(=O)[C@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN1316
DDG (kcal/mol):	-1.70
dDDG (kcal/mol):	0.39

DAR-DIA-0f2f46c9-7_2



CID: DAR-DIA-0f2f46c9-7_2

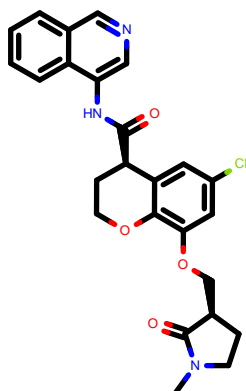
SMILES: CNS(=O)(=O)[N@]1CC[C@@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN3235

DDG (kcal/mol): -1.70

dDDG (kcal/mol): 0.13

MAT-POS-fce787c2-7_2



CID: MAT-POS-fce787c2-7_2

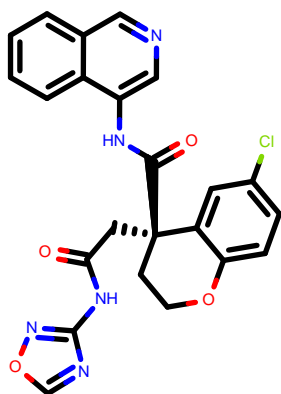
SMILES: C[C@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2cncc3c2cccc3)[NH3+]

RUN: RUN2154

DDG (kcal/mol): -1.70

dDDG (kcal/mol): 0.42

MAK-UNK-3875bbc8-1_2



CID: MAK-UNK-3875bbc8-1_2

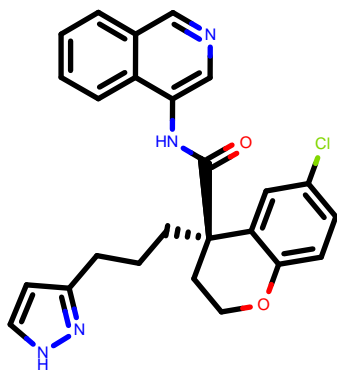
SMILES: Cc1ccc2c(c1)[C@H](CCO2)C(=O)N(C)c3cncc4c3cccc4

RUN: RUN799

DDG (kcal/mol): -1.70

dDDG (kcal/mol): 0.24

VLA-UCB-1dbca3b4-17_2



CID: VLA-UCB-1dbca3b4-17_2

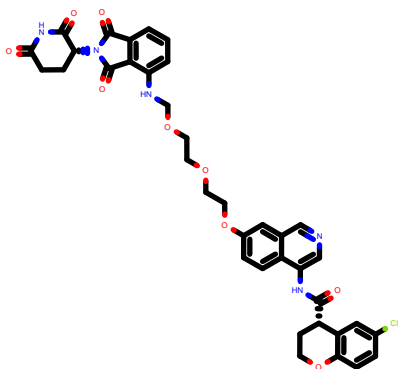
SMILES: c1ccc2c(c1)cncc2n3c(c([nH]c3=O)c4cccc(c4)Cl)[O-]

RUN: RUN167

DDG (kcal/mol): -1.69

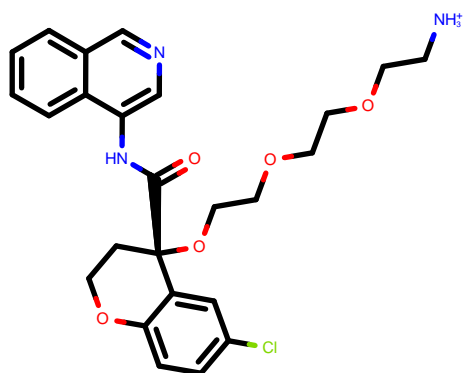
dDDG (kcal/mol): 0.26

EDJ-MED-e4b030d8-3_1



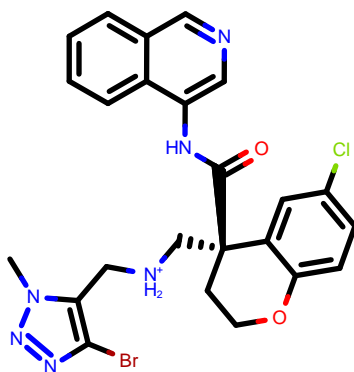
CID:	EDJ-MED-e4b030d8-3_1
SMILES:	<chem>C[C@H]1C[C@H](c2cc(ccc2O1)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN285
DDG (kcal/mol):	-1.69
dDDG (kcal/mol):	0.40

DAR-DIA-6a508060-13_1



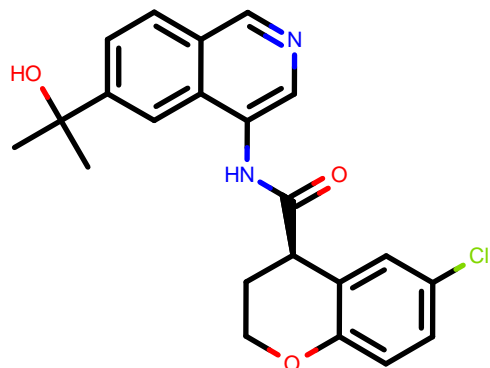
CID:	DAR-DIA-6a508060-13_1
SMILES:	<chem>CS(=O)(=O)NCC[C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN360
DDG (kcal/mol):	-1.69
dDDG (kcal/mol):	0.60

KAD-UNI-8a629cb0-21_1



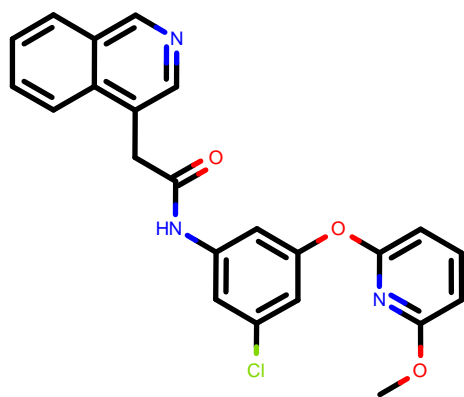
CID:	KAD-UNI-8a629cb0-21_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCN(CC5)C(CO)(F)F</chem>
RUN:	RUN2111
DDG (kcal/mol):	-1.68
dDDG (kcal/mol):	0.42

NIR-THE-d08c3b48-1_1



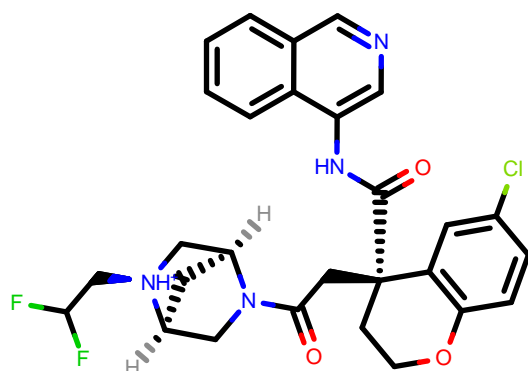
CID:	NIR-THE-d08c3b48-1_1
SMILES:	<chem>CC(=O)N1CC[NH+](CC1)CC(=O)Nc2cncc3c2cccc3</chem>
RUN:	RUN189
DDG (kcal/mol):	-1.68
dDDG (kcal/mol):	0.25

EDG-MED-ba1ac7b9-15_8



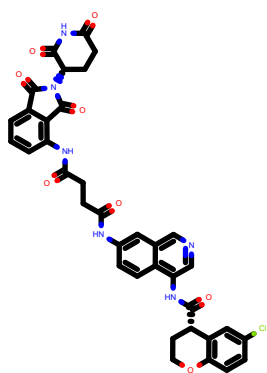
CID:	EDG-MED-ba1ac7b9-15_8
SMILES:	<chem>C[C@H]1CN[H+](CCN1C(=O)C[C@]2(CCOC3C2cc(cc3)Cl)C(=O)Nc4ccc5c4ccc5)C</chem>
RUN:	RUN2681
DDG (kcal/mol):	-1.68
dDDG (kcal/mol):	0.33

WIL-UNI-0732ac76-2_1



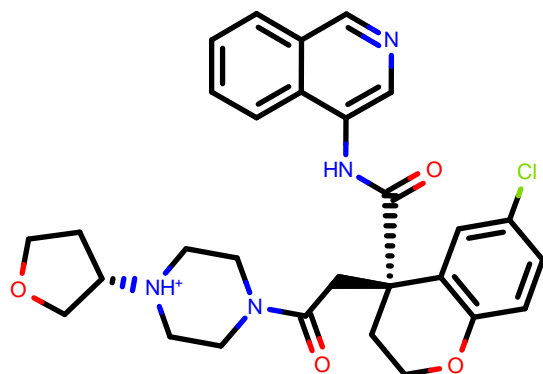
CID:	WIL-UNI-0732ac76-2_1
SMILES:	<chem>C[C@@]1CN[C@@H]1c1ccc2c1c(O)Nc3cccc3Cl(O)c4ccc5c4c5ccccc5F)C(=O)C</chem>
RUN:	RUN1177
DDG (kcal/mol):	-1.67
dDDG (kcal/mol):	0.26

PET-UNK-4880b143-1_1



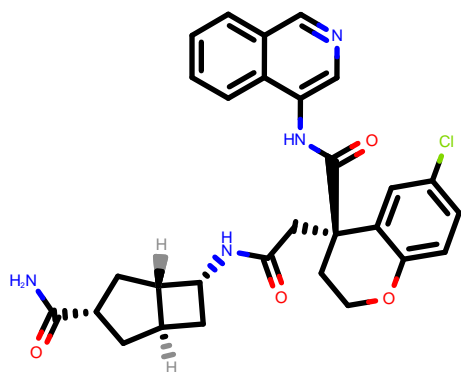
CID:	PET-UNK-4880b143-1_1
SMILES:	<chem>CS(=O)(=O)CCO[C@]11(CCOc2c1cc(cc2)Cl)C(=O)Nc3ccc4c3cccc4</chem>
RUN:	RUN3294
DDG (kcal/mol):	-1.67
dDDG (kcal/mol):	0.43

MAK-UNK-c749d764-16_7



CID:	MAK-UNK-c749d764-16_7
SMILES:	<chem>C[C@@]1H[N](c1ccc2c1cccc2)C(=O)C[C@H]3CCC[C@H]1([C@@H]3O)C(F)F)O</chem>
RUN:	RUN981
DDG (kcal/mol):	-1.67
dDDG (kcal/mol):	0.29

DAR-DIA-0d514e7d-32_26



CID: DAR-DIA-0d514e7d-32_26

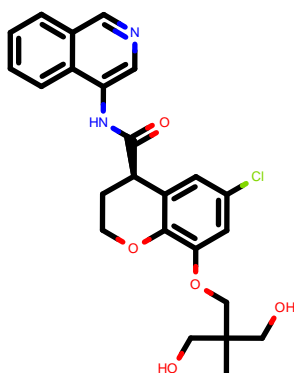
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3[C@@H]4C[C@@H]4C[C@@H]5[C@H]3C=C(C=C5)Cl

RUN: RUN875

DDG (kcal/mol): -1.67

dDDG (kcal/mol): 0.40

MAT-POS-e9e99895-13_8



CID: MAT-POS-e9e99895-13_8

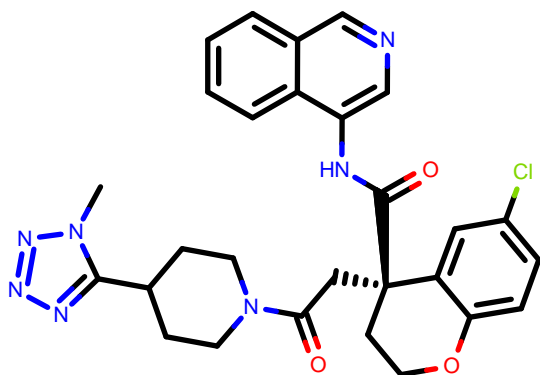
SMILES: C[C@@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2cncc3c2ccc3)NC(=O)[C@H]4CC(N@H+)[C4]C

RUN: RUN2274

DDG (kcal/mol): -1.67

dDDG (kcal/mol): 0.46

MIC-UNK-5a93dd5f-3_3



CID: MIC-UNK-5a93dd5f-3_3

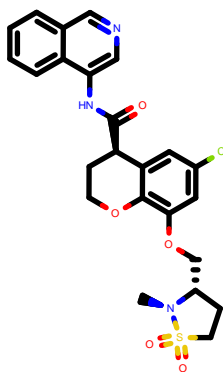
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3[C@@H](c3ccccc3)C[C@@H]4CC[C@@H]5CCCC[C@@H]5C4

RUN: RUN742

DDG (kcal/mol): -1.66

dDDG (kcal/mol): 0.41

ALP-UNI-76695c4f-8_1



CID: ALP-UNI-76695c4f-8_1

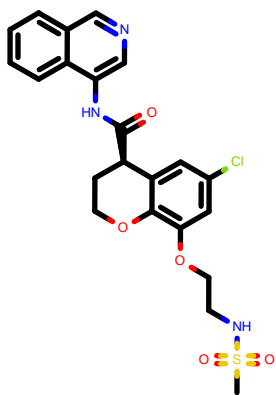
SMILES: C[N@@]1C1CN(C1)Cl(-O)C[C@@]2(CCOC3=CC(=O)C3)C1(-O)N4ncc5c4ccc5(Si(-O)(-O)C

RUN: RUN2177

DDG (kcal/mol): -1.66

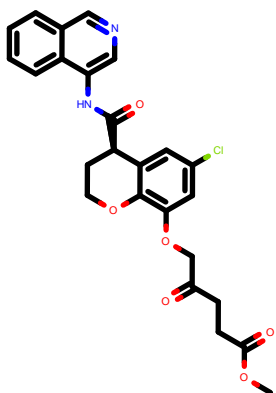
dDDG (kcal/mol): 0.40

MAT-POS-e9e99895-11_1



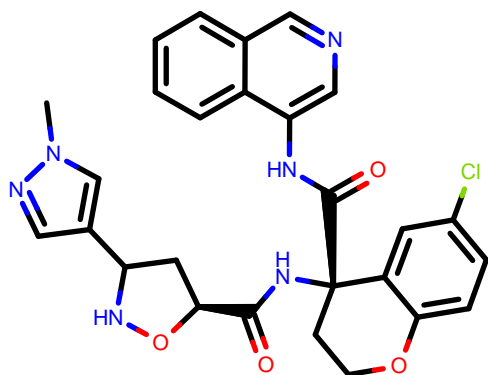
CID:	MAT-POS-e9e99895-11_1
SMILES:	<chem>C[C@@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2ncc3c2cccc3)NC(=O)COC</chem>
RUN:	RUN2263
DDG (kcal/mol):	-1.65
dDDG (kcal/mol):	0.45

JOH-UNI-ea72002d-5_2



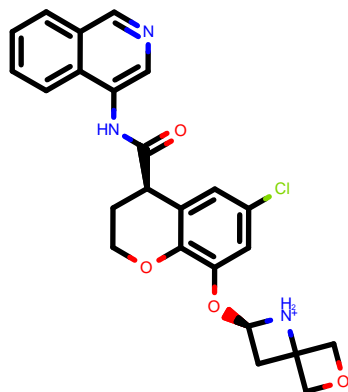
CID:	JOH-UNI-ea72002d-5_2
SMILES:	<chem>C=CS(=O)(=O)NN(c1ncc2c1cccc2)C(=O)[C@]3(CCOc4c3cc(cc4)Cl)F</chem>
RUN:	RUN2493
DDG (kcal/mol):	-1.65
dDDG (kcal/mol):	0.40

ALP-POS-966f8da6-2_2



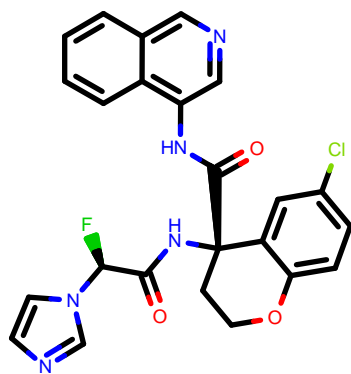
CID:	ALP-POS-966f8da6-2_2
SMILES:	<chem>CC(=O)N1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3ncc4c3cccc4</chem>
RUN:	RUN1226
DDG (kcal/mol):	-1.65
dDDG (kcal/mol):	0.30

MAT-POS-af71705c-1_1



CID:	MAT-POS-af71705c-1_1
SMILES:	<chem>Cn1cc(c2ccccc2c1=O)NC(=O)Nc3ccccn3</chem>
RUN:	RUN54
DDG (kcal/mol):	-1.65
dDDG (kcal/mol):	0.42

DAR-DIA-0f2f46c9-13_2



CID: DAR-DIA-0f2f46c9-13_2

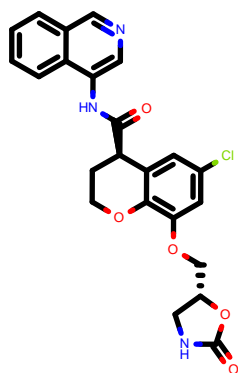
SMILES: CN(C)S(=O)(=O)N@[1]CC[C@@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN3248

DDG (kcal/mol): -1.64

dDDG (kcal/mol): 0.12

MAT-POS-f9802937-2_2



CID: MAT-POS-f9802937-2_2

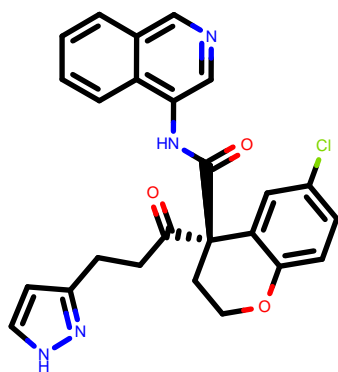
SMILES: CN1c2ccc(cc2[C@](CC1=O)(C(=O)Nc3cncc4c3cccc4)O)Cl

RUN: RUN2393

DDG (kcal/mol): -1.64

dDDG (kcal/mol): 0.36

ERI-UCB-b3e6b0c2-4_1



CID: ERI-UCB-b3e6b0c2-4_1

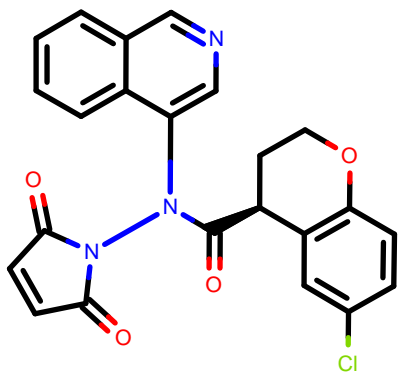
SMILES: C[N@@1]C[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3ccc(cc4)CN5CC[NH2+]CC5

RUN: RUN3036

DDG (kcal/mol): -1.64

dDDG (kcal/mol): 0.16

MAK-UNK-c749d764-12_2



CID: MAK-UNK-c749d764-12_2

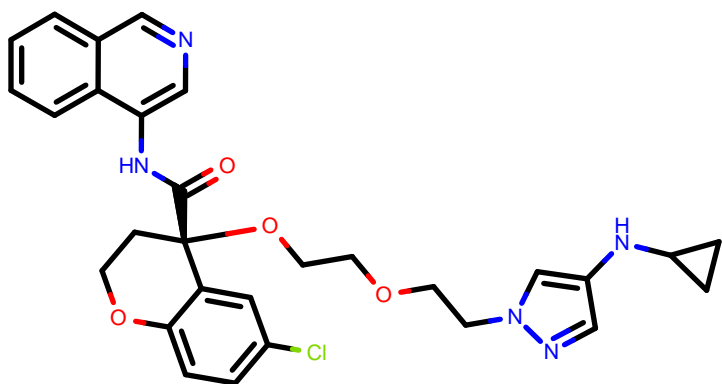
SMILES: c1ccc2c(c1)ncnc2NC(=O)C[C@H]3CCCC[C@@H]3[C@@H](Cl)Cl

RUN: RUN951

DDG (kcal/mol): -1.64

dDDG (kcal/mol): 0.36

MIC-UNK-cdc2493e-10_1



CID: MIC-UNK-cdc2493e-10_1

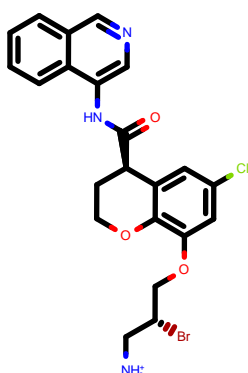
SMILES: CC(=O)N[C@@H]1CC[C@@H](C1)N(c2cccc(c2)C)C(=O)Nc3ncc4c3cccc4

RUN: RUN545

DDG (kcal/mol): -1.64

dDDG (kcal/mol): 0.37

MAT-POS-fce787c2-10_1



CID: MAT-POS-fce787c2-10_1

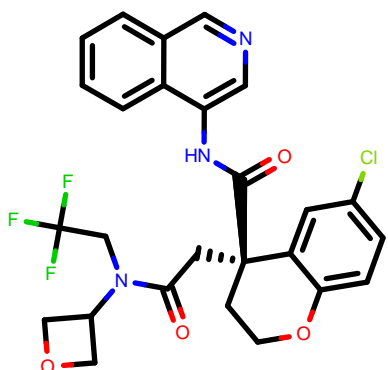
SMILES: CN(C)CC[NH2+][C@@H](c1ccc(c(c1)Cl)Cl)C(=O)Nc2ncc3c2cccc3

RUN: RUN2160

DDG (kcal/mol): -1.63

dDDG (kcal/mol): 0.38

FRA-DIA-a1f3a927-1_1



CID: FRA-DIA-a1f3a927-1_1

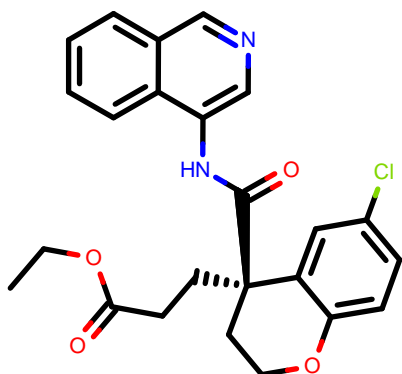
SMILES: C=C(c1cncc2c1cccc2)C(=O)[C@@H](c1ccc(c(c1)Cl)Cl)C(=O)Nc2cnc3c2cccc3

RUN: RUN1176

DDG (kcal/mol): -1.63

dDDG (kcal/mol): 0.26

MAT-POS-2492181e-2_1



CID: MAT-POS-2492181e-2_1

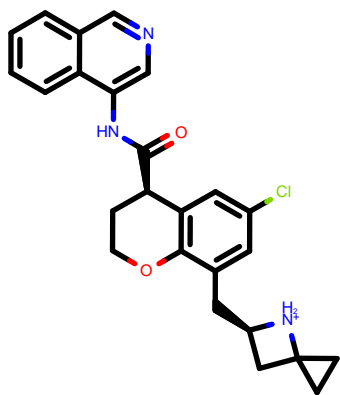
SMILES: Cn1cc(c2cccc2c1=O)NC(=O)NCC[NH+]3CCCC3

RUN: RUN94

DDG (kcal/mol): -1.62

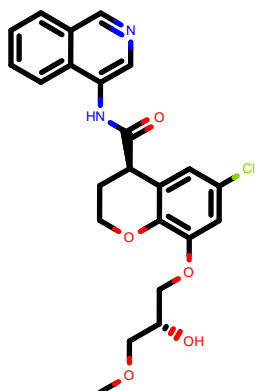
dDDG (kcal/mol): 0.24

MAT-POS-3b92565d-9_2



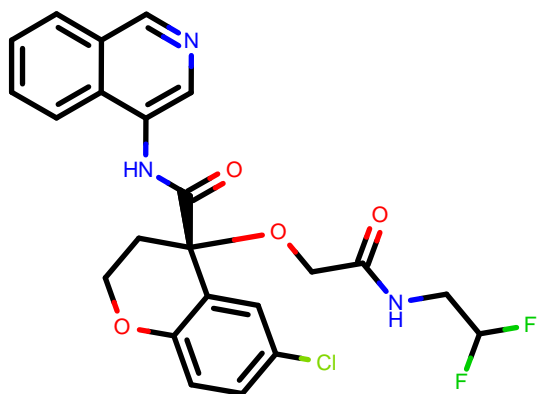
CID:	MAT-POS-3b92565d-9_2
SMILES:	<chem>Cn1cc(e2ccccc2c1=O)NC(=O)Cc3cc(cc(c3)Cl)O[C@H]4CC(=O)N4</chem>
RUN:	RUN120
DDG (kcal/mol):	-1.62
dDDG (kcal/mol):	0.37

MAT-POS-e69ad64a-2_2



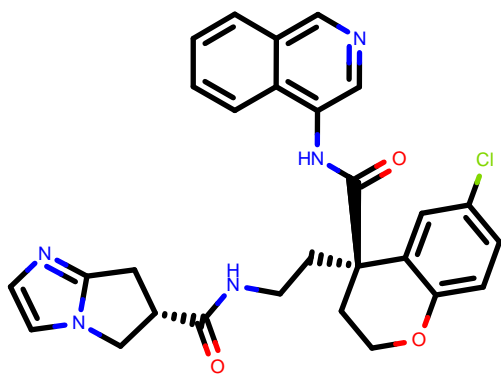
CID:	MAT-POS-e69ad64a-2_2
SMILES:	<chem>C=CC(=O)N(c1cncc2c1cccc2)C(=O)[C@H]3COc4c3cc(cc4)Cl</chem>
RUN:	RUN2333
DDG (kcal/mol):	-1.61
dDDG (kcal/mol):	0.42

MIC-UNK-d854bf4c-7_2



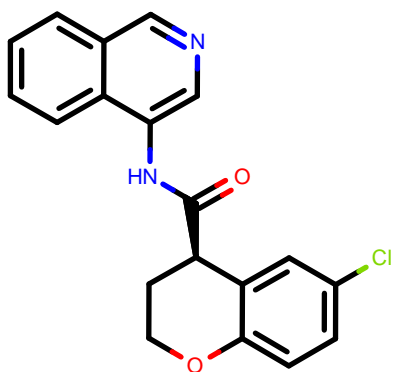
CID:	MIC-UNK-d854bf4c-7_2
SMILES:	<chem>CS(=O)(=O)N1CCC2(CC1)CCN(C(=O)[C@H]2c3cccc(c3)Cl)c4ncc5c4cccc5</chem>
RUN:	RUN3342
DDG (kcal/mol):	-1.60
dDDG (kcal/mol):	0.15

DAR-DIA-0587064e-26_2



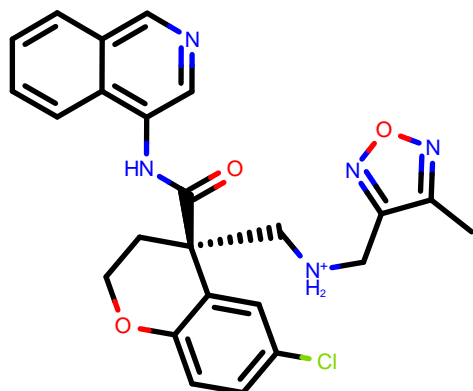
CID:	DAR-DIA-0587064e-26_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCOCc4c3cc(c(c4OCc5ccc(cc5Cl)F)F)Cl</chem>
RUN:	RUN3389
DDG (kcal/mol):	-1.60
dDDG (kcal/mol):	0.22

DAR-DIA-eace69ff-19_1



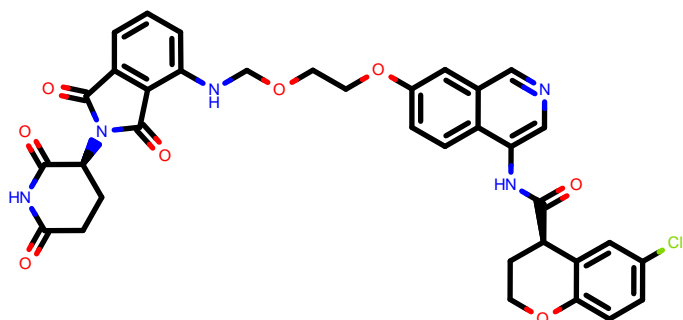
CID:	DAR-DIA-eace69ff-19_1
SMILES:	<chem>CCc1cc2c(cnc2c(c1)OC3CCCCC3)C[NH+](C)C</chem>
RUN:	RUN2
DDG (kcal/mol):	-1.59
dDDG (kcal/mol):	0.20

JOH-UNI-6e27fddc-4_1



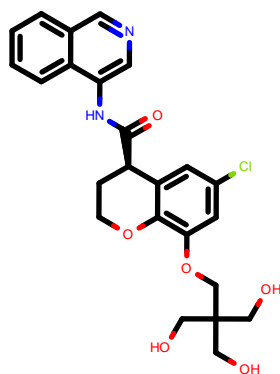
CID:	JOH-UNI-6e27fddc-4_1
SMILES:	<chem>CO[C@]1(CCOC2c1cc(cc2)Cl)/C(=C/c3ncc4c3cccc4)/F</chem>
RUN:	RUN4369
DDG (kcal/mol):	-1.58
dDDG (kcal/mol):	0.26

DAR-DIA-6a508060-9_1



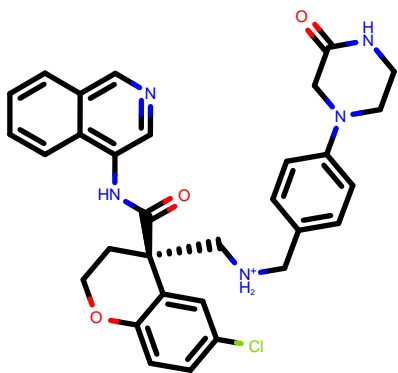
CID:	DAR-DIA-6a508060-9_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CCCC5CCCC5</chem>
RUN:	RUN348
DDG (kcal/mol):	-1.58
dDDG (kcal/mol):	0.48

EDG-MED-ba1ac7b9-9_3



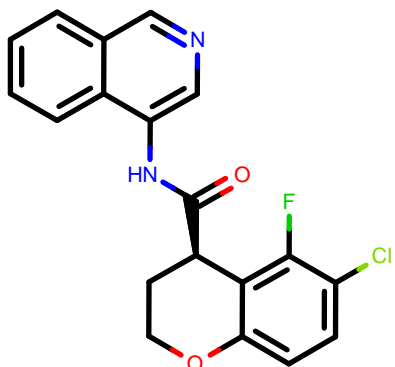
CID:	EDG-MED-ba1ac7b9-9_3
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCCC[C@H]5c6[nH]cn6</chem>
RUN:	RUN2648
DDG (kcal/mol):	-1.58
dDDG (kcal/mol):	0.42

ALP-POS-347519b5-1_50



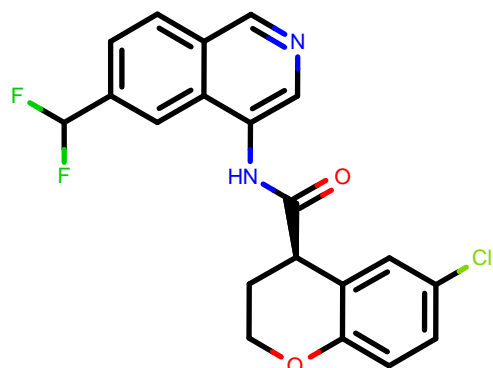
CID:	ALP-POS-347519b5-1_50
SMILES:	<chem>CS(=O)(=O)[N@@]1C[C@@H]2[C@@H]3CC[C@@H]3[C@@H]2[C@@H]1C(=O)Nc4ccc5c6ccccc5</chem>
RUN:	RUN4257
DDG (kcal/mol):	-1.58
dDDG (kcal/mol):	0.27

MAK-UNK-ffc90da7-9_8



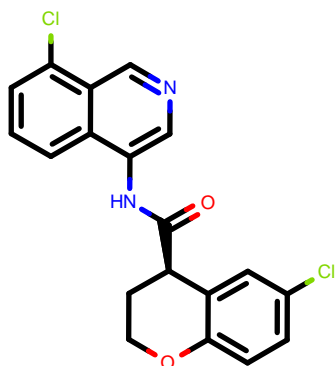
CID:	MAK-UNK-ffc90da7-9_8
SMILES:	<chem>C[C@@H]1[C@@H](CCO1)SC[C@@H](C)C2=CC=C(C(=O)Nc3ccc(F)c(Cl)c3)C2</chem>
RUN:	RUN719
DDG (kcal/mol):	-1.57
dDDG (kcal/mol):	0.20

LON-WEI-4d77710c-18_1



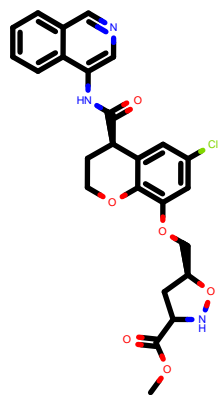
CID:	LON-WEI-4d77710c-18_1
SMILES:	<chem>Cn1cc(c2ccccc2c1=O)NC(=O)Nc3ccc(cc3OC)OC</chem>
RUN:	RUN203
DDG (kcal/mol):	-1.57
dDDG (kcal/mol):	0.23

MAT-POS-b5746674-106_1



CID:	MAT-POS-b5746674-106_1
SMILES:	<chem>C[C@@H]1[C@@H](c1cncc1)N(CCCOC)C(=O)Nc2cn(c(=O)c3c2ccccc3)CC(C)C</chem>
RUN:	RUN86
DDG (kcal/mol):	-1.57
dDDG (kcal/mol):	0.18

EDJ-MED-40433386-2_1



CID: EDJ-MED-40433386-2_1

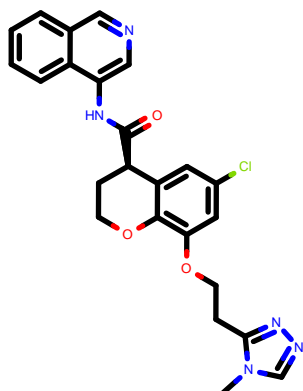
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CNC(=O)c5cn[nH]c5S(=O)(=O)N

RUN: RUN2553

DDG (kcal/mol): -1.57

dDDG (kcal/mol): 0.38

EDJ-MED-fcba3f31-9_1



CID: EDJ-MED-fcba3f31-9_1

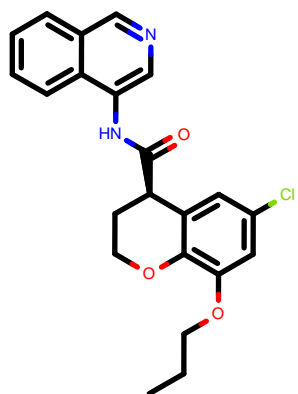
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)COC[C@H]5CC(=O)N5

RUN: RUN2548

DDG (kcal/mol): -1.57

dDDG (kcal/mol): 0.37

ED_-GRI-5b13fbe2-27_1



CID: ED_-GRI-5b13fbe2-27_1

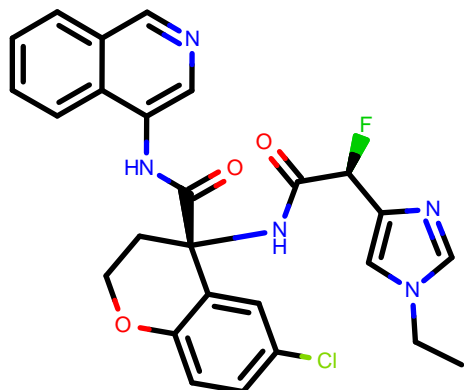
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OCCn5ccn5CN6CC[NH2+]CC6

RUN: RUN1557

DDG (kcal/mol): -1.56

dDDG (kcal/mol): 0.31

PET-UNK-a17c93d1-1_1



CID: PET-UNK-a17c93d1-1_1

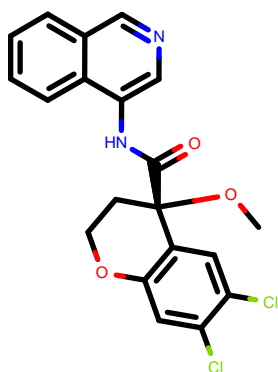
SMILES: c1ccc2c(c1)cncc2NC(=O)N(c3cccc(c3)Cl)OCC4CCCC4

RUN: RUN520

DDG (kcal/mol): -1.56

dDDG (kcal/mol): 0.27

DAR-DIA-0d514e7d-31_8



CID: DAR-DIA-0d514e7d-31_8

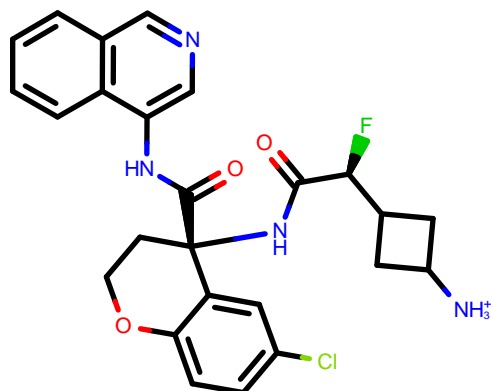
SMILES: C[C@H]1CCO[C@H]2C=CC(=C[C@H]2[C@@H]1C(=O)Nc3ccc4c3ccc4)Cl

RUN: RUN841

DDG (kcal/mol): -1.56

dDDG (kcal/mol): 0.21

MAK-UNK-8be7dca9-1_1



CID: MAK-UNK-8be7dca9-1_1

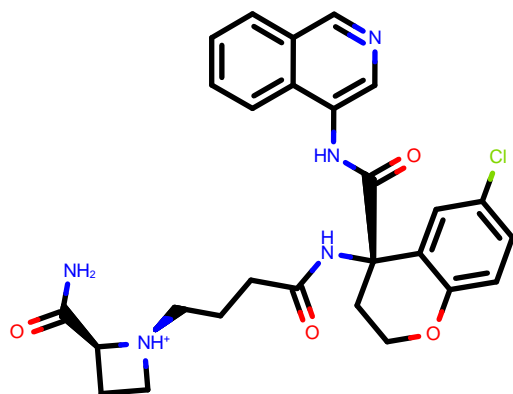
SMILES: c1cc2c(cc1CN3CC[NH2+][CC3]nccc2NC(=O)[C@@H]4CCOc5c4cc(cc5)Cl

RUN: RUN491

DDG (kcal/mol): -1.56

dDDG (kcal/mol): 0.27

MIC-UNK-91acba05-3_2



CID: MIC-UNK-91acba05-3_2

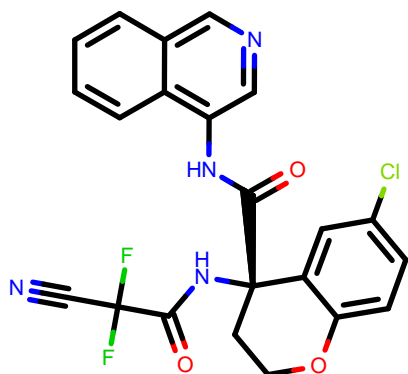
SMILES: c1ccc2c(c1)nccc2NC(=O)[C@@H]3CCS(=O)(=O)c4c3cc(cc4)Cl

RUN: RUN473

DDG (kcal/mol): -1.55

dDDG (kcal/mol): 0.33

NIR-THE-47736cde-2_1



CID: NIR-THE-47736cde-2_1

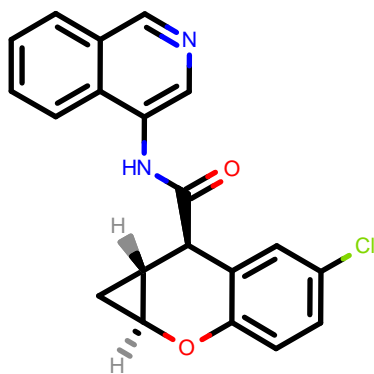
SMILES: C=CC(=O)N(c1cnc2c1cccc2)C(=O)[C@@H]3COc4c3cc(cc4)Cl

RUN: RUN3328

DDG (kcal/mol): -1.55

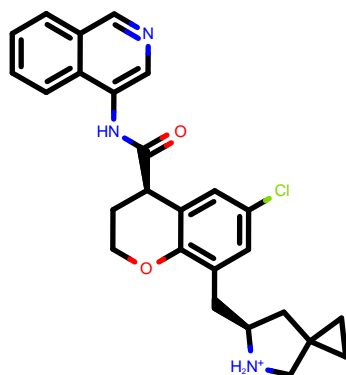
dDDG (kcal/mol): 0.10

LON-WEI-4d77710c-17_1



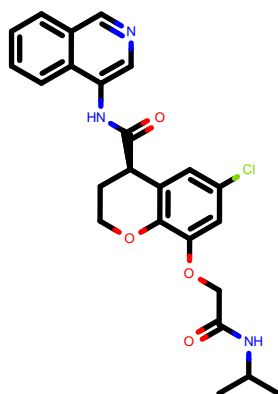
CID:	LON-WEI-4d77710c-17_1
SMILES:	<chem>CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)NCCc3ccc(cc3)Cl</chem>
RUN:	RUN205
DDG (kcal/mol):	-1.55
dDDG (kcal/mol):	0.13

DAR-DIA-0cde14eb-60_1



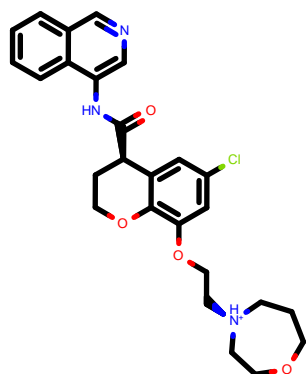
CID:	DAR-DIA-0cde14eb-60_1
SMILES:	<chem>C[C@@H](c1cccc(c1)C2(CC2)C#N)C(=O)Nc3ncc4c3cccc4</chem>
RUN:	RUN30
DDG (kcal/mol):	-1.55
dDDG (kcal/mol):	0.47

EDJ-MED-d08626de-5_1



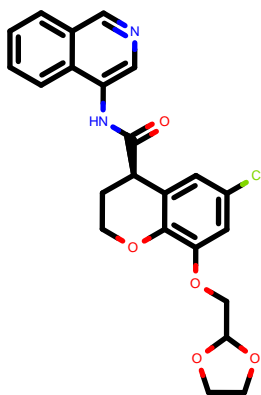
CID:	EDJ-MED-d08626de-5_1
SMILES:	<chem>CO[C@@@]1(CCOC2c1cc(c(c2)Cl)Cl)C(=O)Nc3ncc4c3cccc4</chem>
RUN:	RUN2345
DDG (kcal/mol):	-1.54
dDDG (kcal/mol):	0.29

EDG-MED-ba1ac7b9-5_4



CID:	EDG-MED-ba1ac7b9-5_4
SMILES:	<chem>CC[C@H](CO)[NH+]1CCN(CC1)C(=O)C[C@]2(CCOC3c2cc(cc3)Cl)C(=O)Nc4ncc5c4cccc5</chem>
RUN:	RUN2641
DDG (kcal/mol):	-1.53
dDDG (kcal/mol):	0.60

KAD-UNI-8a629cb0-25_1



CID: KAD-UNI-8a629cb0-25_1

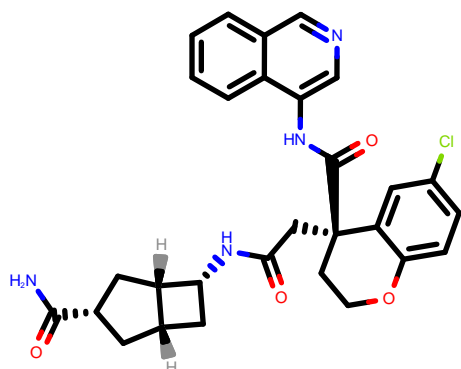
SMILES: c1ccc2c1)nc2NC(=O)C(=O)N3CCN(CC3)C(=O)C4=CC=C(C=C4)OC5=CC=CC=C5

RUN: RUN2107

DDG (kcal/mol): -1.53

dDDG (kcal/mol): 0.42

LON-WEI-5e7d1b3e-45_1



CID: LON-WEI-5e7d1b3e-45_1

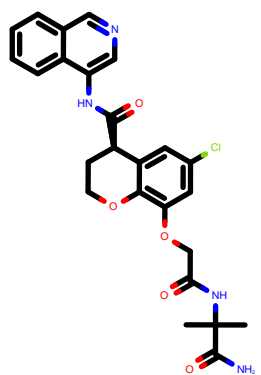
SMILES: Cn1cc(c2ccccc2c1=O)NC(=O)N3CCN(CC3)C(=O)c4cccc4

RUN: RUN1354

DDG (kcal/mol): -1.53

dDDG (kcal/mol): 0.41

ALP-UNI-3496895b-5_4



CID: ALP-UNI-3496895b-5_4

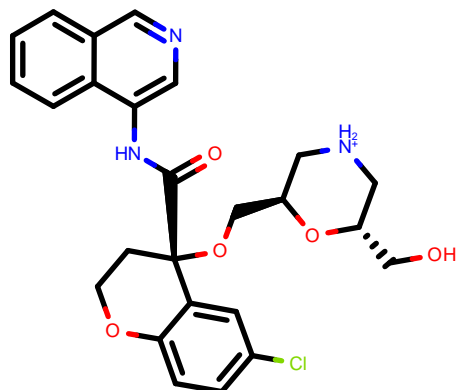
SMILES: C(C)(C)C(=O)NC(=O)C(=O)N3CCN(CC3)C(=O)C4=CC=C(C=C4)OC5=CC=CC=C5

RUN: RUN2515

DDG (kcal/mol): -1.52

dDDG (kcal/mol): 0.44

VLA-UCB-05e51b3f-16_1



CID: VLA-UCB-05e51b3f-16_1

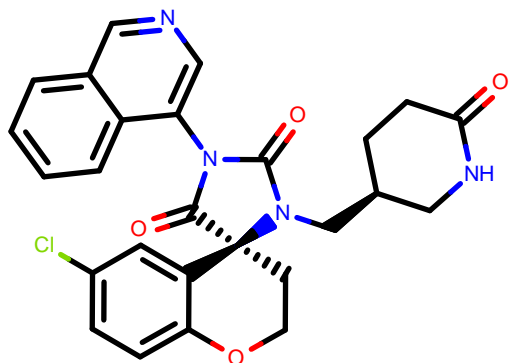
SMILES: C=CC(=O)N(c1cncc2c1cccc2)C(=O)Cc3cccc(c3)Cl

RUN: RUN328

DDG (kcal/mol): -1.52

dDDG (kcal/mol): 0.45

MAK-UNK-c749d764-19_8



CID: MAK-UNK-c749d764-19_8

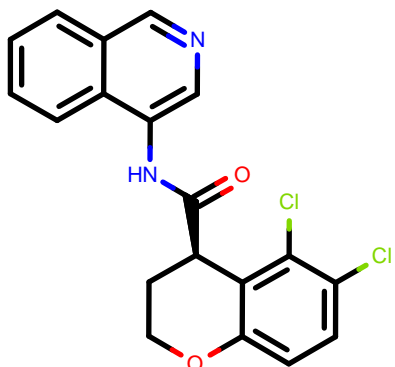
SMILES: c1ccc2c(c1)cncc2N(C[NH3+])C(=O)C[C@H]3CCC[C@H]([C@H]3O)C(F)F

RUN: RUN1005

DDG (kcal/mol): -1.52

dDDG (kcal/mol): 0.25

MAT-POS-b5746674-103_1



CID: MAT-POS-b5746674-103_1

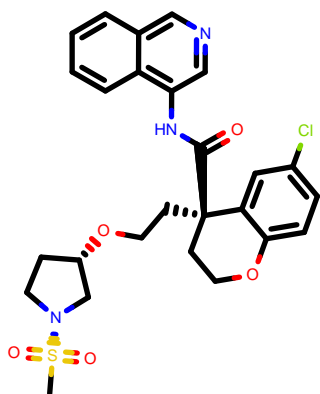
SMILES: CC(C)Cn1cc(c2cccc2c1=O)NC(=O)N3CCN(CC3)c4cccc4F

RUN: RUN81

DDG (kcal/mol): -1.52

dDDG (kcal/mol): 0.20

RUB-POS-1325a9ea-2_1



CID: RUB-POS-1325a9ea-2_1

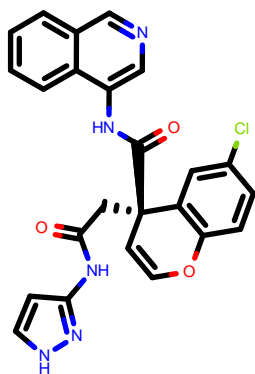
SMILES: c1cc(cc(c1)Cl)CC(=O)Nc2cncc3c2c(ccc3)F

RUN: RUN3606

DDG (kcal/mol): -1.51

dDDG (kcal/mol): 0.20

LON-WEI-4d77710c-50_2



CID: LON-WEI-4d77710c-50_2

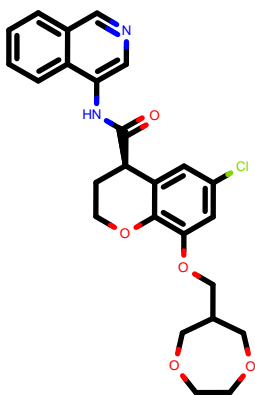
SMILES: CCCC[N@H+]([C]N(C(=O)Nc1cn(c(=O)c2c1cccc2)C)Cc3ccco3

RUN: RUN246

DDG (kcal/mol): -1.51

dDDG (kcal/mol): 0.26

JOH-UNI-ea72002d-4_1



CID: JOH-UNI-ea72002d-4_1

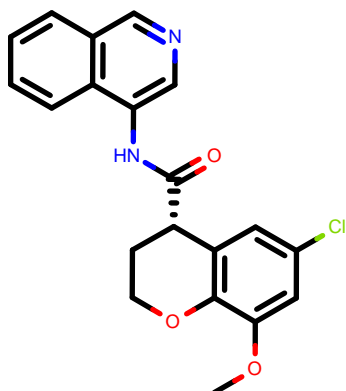
SMILES: c1ccc2c(c1)ncnc2[C@@]([C@H]3CCOC4C3CC(C4)C)(N5C(=O)C=CC5=O)F

RUN: RUN2488

DDG (kcal/mol): -1.50

dDDG (kcal/mol): 0.43

LON-WEI-4d77710c-21_1



CID: LON-WEI-4d77710c-21_1

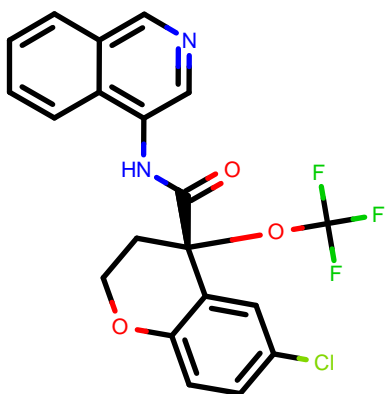
SMILES: Cc1cccnc1NC(=O)Nc2cn(c(=O)c3c2cccc3)CC(C)C

RUN: RUN207

DDG (kcal/mol): -1.50

dDDG (kcal/mol): 0.13

MAT-POS-5369c344-3_2



CID: MAT-POS-5369c344-3_2

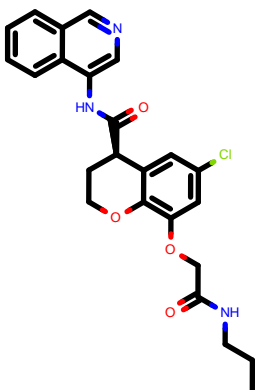
SMILES: c1ccc(cc1)COc2cc(cc3c2OCC[C@H]3C(=O)Nc4cnc5c4cccc5)Cl

RUN: RUN3619

DDG (kcal/mol): -1.50

dDDG (kcal/mol): 0.08

EDJ-MED-9e38fd34-4_1



CID: EDJ-MED-9e38fd34-4_1

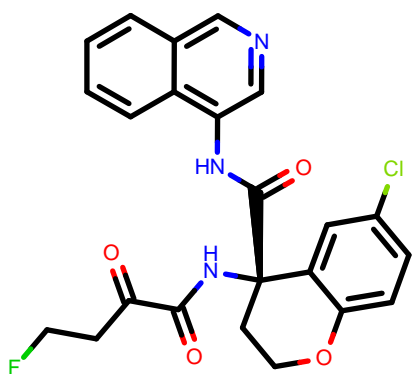
SMILES: C[C@@]1(c2cc(c(cc2NC1=O)Cl)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN2349

DDG (kcal/mol): -1.50

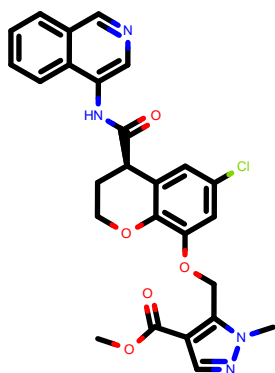
dDDG (kcal/mol): 0.34

DAR-DIA-0587064e-13_2



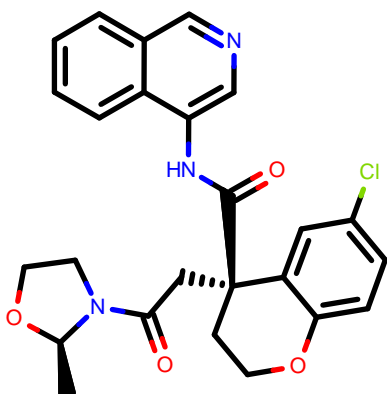
CID:	DAR-DIA-0587064e-13_2
SMILES:	<chem>CCCOc1cc(cc2c1NCC[C@H]2C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN3361
DDG (kcal/mol):	-1.49
dDDG (kcal/mol):	0.12

EDJ-MED-d203f206-34_1



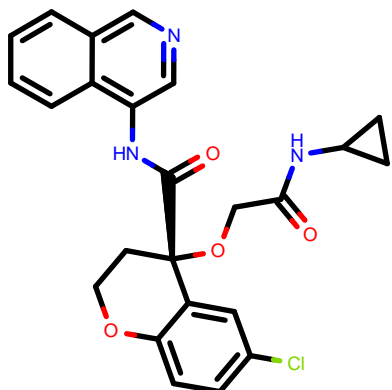
CID:	EDJ-MED-d203f206-34_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(c4)Cl)CC(=O)N[C@@H]5C[C@H]5CO</chem>
RUN:	RUN2597
DDG (kcal/mol):	-1.49
dDDG (kcal/mol):	0.34

ALP-POS-d3acb8cc-1_1



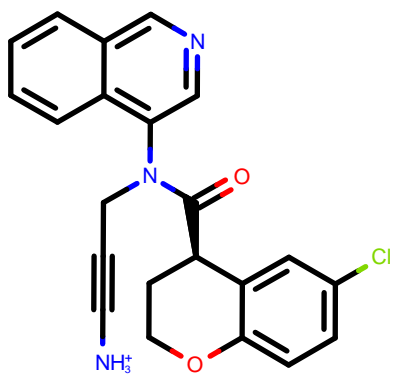
CID:	ALP-POS-d3acb8cc-1_1
SMILES:	<chem>C[C@H]1C[C@H](c2cc(ccc2O1)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN1092
DDG (kcal/mol):	-1.49
dDDG (kcal/mol):	0.26

MIC-UNK-d854bf4c-1_1



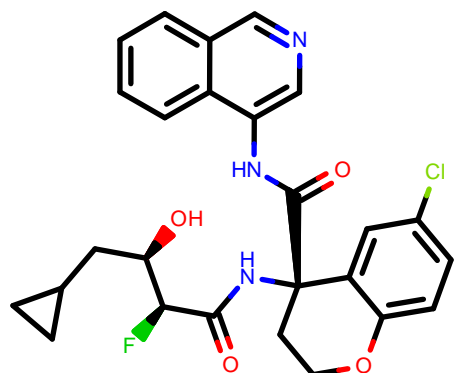
CID:	MIC-UNK-d854bf4c-1_1
SMILES:	<chem>CC(=O)N1CCC2(CC1)CN(C(=O)[C@@H]2c3cccc(c3)Cl)c4cncc5c4cccc5</chem>
RUN:	RUN3326
DDG (kcal/mol):	-1.49
dDDG (kcal/mol):	0.12

DAR-DIA-ecdbc7dd-8_2



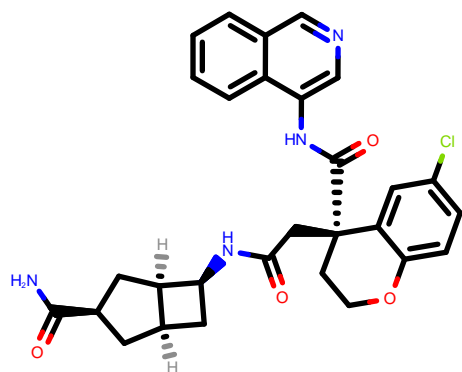
CID:	DAR-DIA-ecdbc7dd-8_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]3(CCNc4c3cc(cc4)Cl)C[NH+]5CCCC5</chem>
RUN:	RUN2888
DDG (kcal/mol):	-1.48
dDDG (kcal/mol):	0.13

DAR-DIA-0d514e7d-35_1



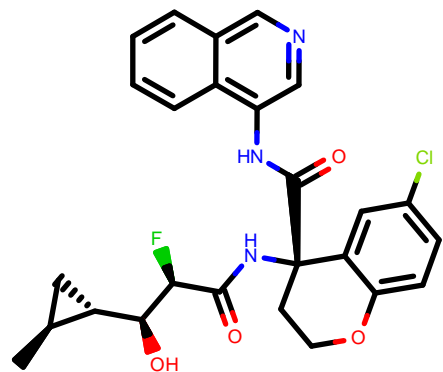
CID:	DAR-DIA-0d514e7d-35_1
SMILES:	<chem>C[C@H]1[C@@H](c2cc(ccc2O1)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN887
DDG (kcal/mol):	-1.48
dDDG (kcal/mol):	0.29

DAR-DIA-9e4459de-15_12



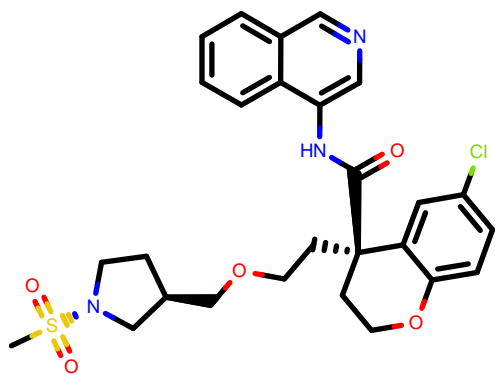
CID:	DAR-DIA-9e4459de-15_12
SMILES:	<chem>c1cc2c(c1)NC(=O)CCCC(=O)Nc3cccc(c3)cncc4NC(=O)[C@]5(CCNc6c4cc(cc6)Cl)C[nH+]7CCCC1=O)NC7=O</chem>
RUN:	RUN1455
DDG (kcal/mol):	-1.48
dDDG (kcal/mol):	0.27

EDJ-MED-15e90dfc-3_1



CID:	EDJ-MED-15e90dfc-3_1
SMILES:	<chem>COCC[NH2+][C][C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3441
DDG (kcal/mol):	-1.47
dDDG (kcal/mol):	0.14

MAK-UNK-83e0a0b4-1_2



CID: MAK-UNK-83e0a0b4-1_2

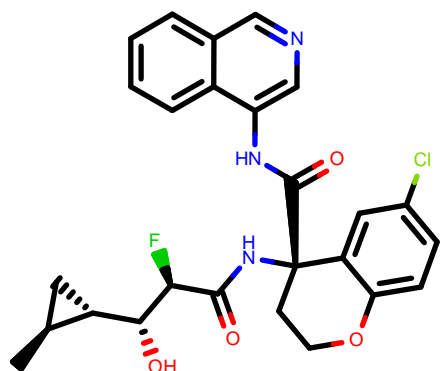
SMILES: CCCC1ccc2c(c1)[C@H](C)CCO2Cl-O[N]C(C)C3C(NH2+)[C]3c4ccc5c4ccc(c5)C[N+]2+

RUN: RUN733

DDG (kcal/mol): -1.47

dDDG (kcal/mol): 0.37

MIC-UNK-5a93dd5f-3_6



CID: MIC-UNK-5a93dd5f-3_6

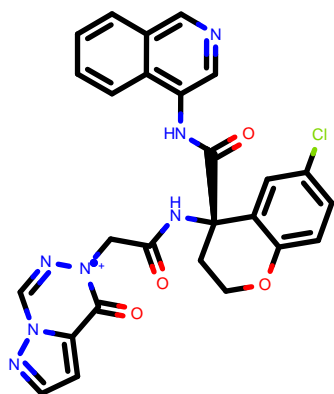
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@H](c3ccccc3)Cl[N+]@@H+4CC[C@H]5CCCC[C@H]5C4

RUN: RUN745

DDG (kcal/mol): -1.46

dDDG (kcal/mol): 0.28

RAL-THA-8416115c-1_4



CID: RAL-THA-8416115c-1_4

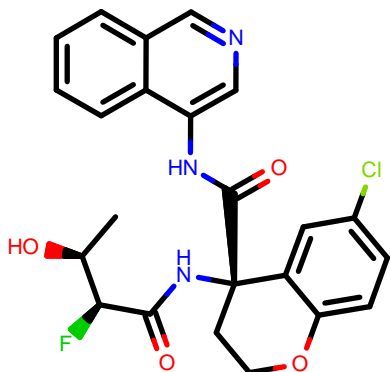
SMILES: c1ccc(cc1)CN2CC[C@H](c3c2ccc(c3)Cl)C(=O)Nc4cnc5c4cccc5

RUN: RUN1250

DDG (kcal/mol): -1.46

dDDG (kcal/mol): 0.34

MIC-UNK-cdc2493e-8_7



CID: MIC-UNK-cdc2493e-8_7

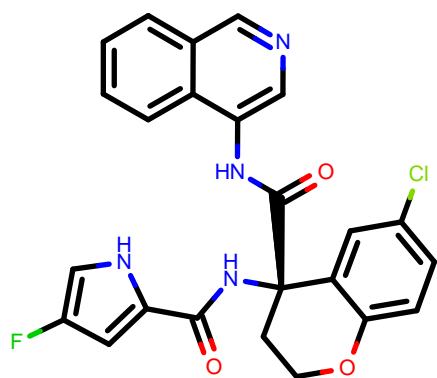
SMILES: c1ccc2c(c1)ncnc2NC(=O)N(c3ccccc3)Cl[C@H]4CC[C@H]5CCCC[C@H]5C4

RUN: RUN542

DDG (kcal/mol): -1.46

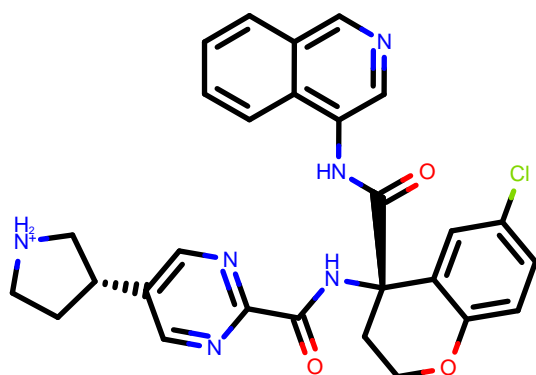
dDDG (kcal/mol): 0.24

JAG-UCB-706446eb-5_2



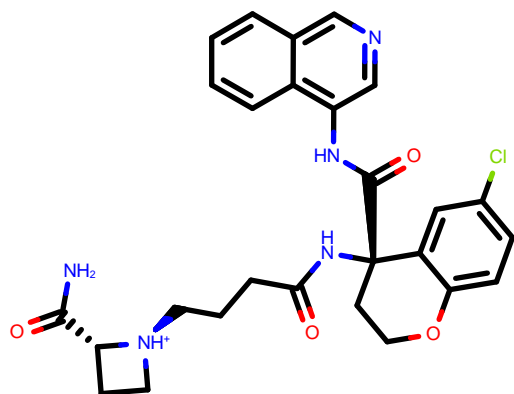
CID:	JAG-UCB-706446eb-5_2
SMILES:	<chem>COCN([C@H]1CCCOC1)C(=O)C[C@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4ccc5c4cccc5</chem>
RUN:	RUN625
DDG (kcal/mol):	-1.46
dDDG (kcal/mol):	0.25

MIC-UNK-8758c41d-1_2



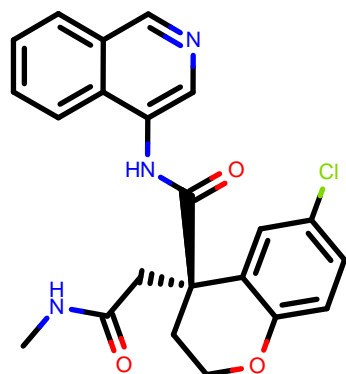
CID:	MIC-UNK-8758c41d-1_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCS(=O)(=O)c4c3cc(c(c4)Cl)Cl</chem>
RUN:	RUN3314
DDG (kcal/mol):	-1.46
dDDG (kcal/mol):	0.16

MAT-POS-0c8fa4a7-1_1



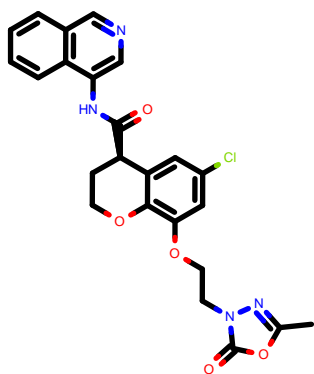
CID:	MAT-POS-0c8fa4a7-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)C3=CCCc4c3cc(cc4)Cl</chem>
RUN:	RUN517
DDG (kcal/mol):	-1.46
dDDG (kcal/mol):	0.37

MAK-UNK-c749d764-3_5



CID:	MAK-UNK-c749d764-3_5
SMILES:	<chem>CC1(C[NH2+])C1OCN(c2ccc3c2cccc3)C(=O)C[C@@H]4CC[C@@H]([C@H]4O)C(F)F</chem>
RUN:	RUN909
DDG (kcal/mol):	-1.45
dDDG (kcal/mol):	0.30

ALP-UNI-3496895b-1_2



CID: ALP-UNI-3496895b-1_2

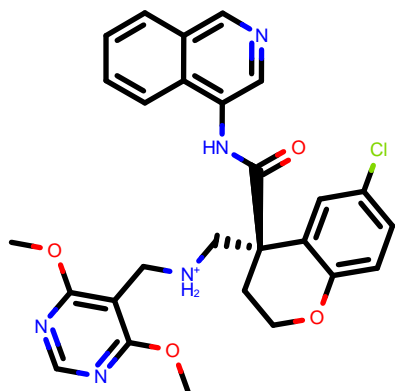
SMILES: c1ccc2c(c1)ncnc2NC(=O)C@@@3[C@@H](COCc4cc(Cl)cc(C1=O)N(CCS1=O)=O)N3

RUN: RUN2500

DDG (kcal/mol): -1.45

dDDG (kcal/mol): 0.36

ALP-POS-347519b5-1_13



CID: ALP-POS-347519b5-1_13

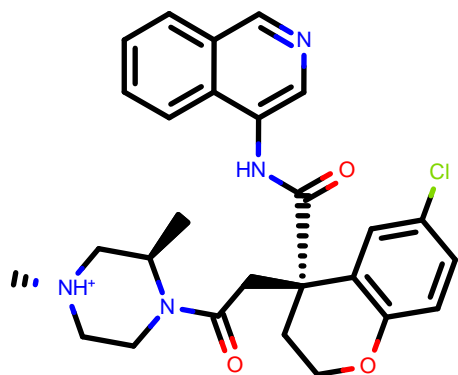
SMILES: CS1=O=O)N@@1C[C@@H]2[C@@H]3CC[C@@H]3(C)C@@H2[C@@H]1C1=O)Nc4ncoc5c4ccccc5

RUN: RUN4236

DDG (kcal/mol): -1.45

dDDG (kcal/mol): 0.30

BEN-BAS-c2bc0d80-3_1



CID: BEN-BAS-c2bc0d80-3_1

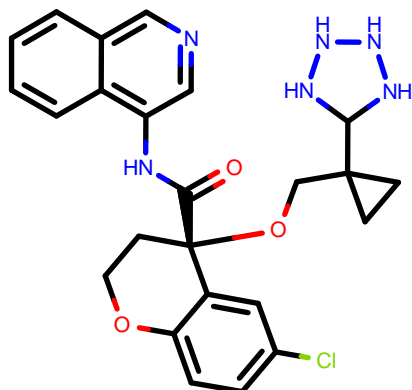
SMILES: c1ccc2c(c1)cncc2N3C(=O)[C@@@4(COCc5c4cc(Cl)cc5)CN3

RUN: RUN1134

DDG (kcal/mol): -1.45

dDDG (kcal/mol): 0.52

ALP-POS-fe871b40-10_1



CID: ALP-POS-fe871b40-10_1

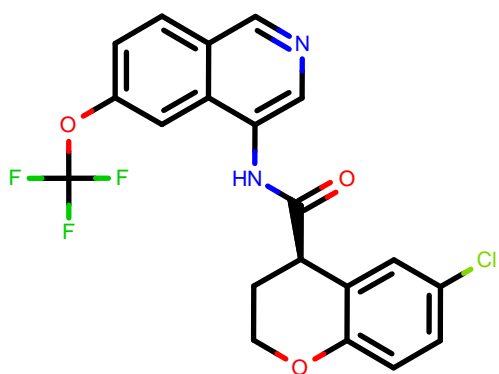
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@@H]3CCNc4c3cc(c(c4)F)Cl

RUN: RUN3123

DDG (kcal/mol): -1.45

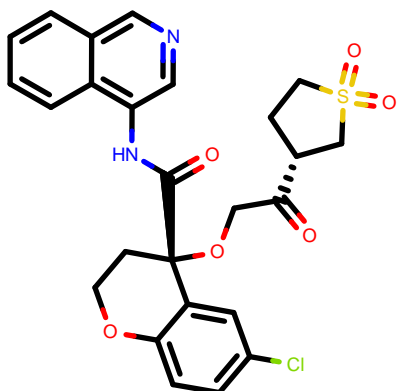
dDDG (kcal/mol): 0.22

EDJ-MED-e4b030d8-12_1



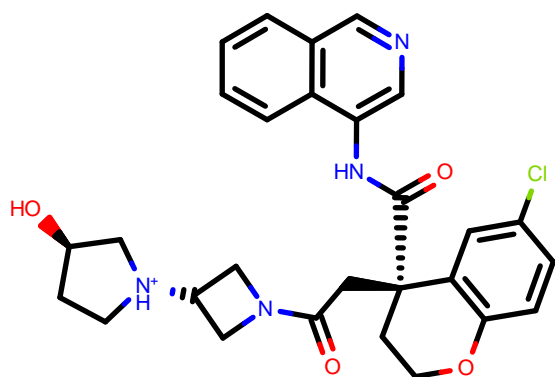
CID:	EDJ-MED-e4b030d8-12_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3Cc4cc(ccc4OC3)Cl</chem>
RUN:	RUN292
DDG (kcal/mol):	-1.45
dDDG (kcal/mol):	0.22

DAR-DIA-f6ee7aeb-5_2



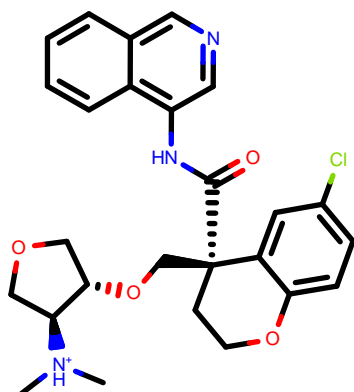
CID:	DAR-DIA-f6ee7aeb-5_2
SMILES:	<chem>CCCCc1cc(cc(c1)Cl)[C@H]2CC(=O)N(C)C@@H]2c3d[nH]c(=O)[nH]c3=O)c4cncc5c4cccc5</chem>
RUN:	RUN3415
DDG (kcal/mol):	-1.44
dDDG (kcal/mol):	0.20

LON-WEI-5e7d1b3e-34_2



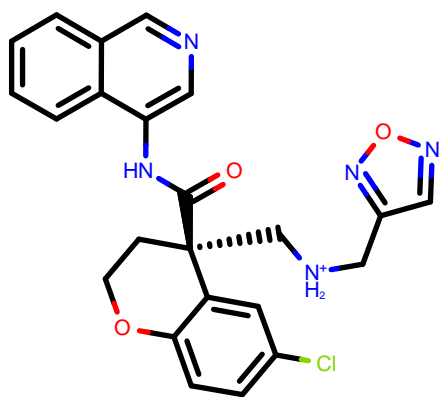
CID:	LON-WEI-5e7d1b3e-34_2
SMILES:	<chem>CC(C)Cn1cc(cc2ccccc2c1=O)NC(=O)N3CC[C@H](C3)c4cccc4</chem>
RUN:	RUN1336
DDG (kcal/mol):	-1.43
dDDG (kcal/mol):	0.29

EDJ-MED-15e90dfc-7_2



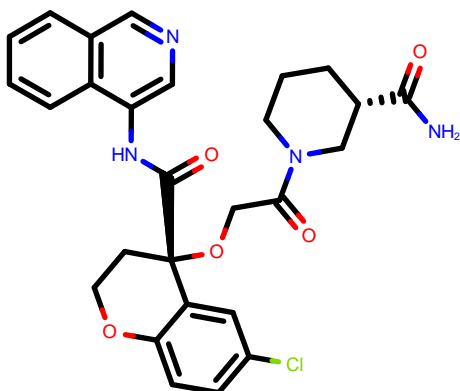
CID:	EDJ-MED-15e90dfc-7_2
SMILES:	<chem>CS(=O)(=O)CCNC[C@]1(C)C(CO)c2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3458
DDG (kcal/mol):	-1.43
dDDG (kcal/mol):	0.35

EDG-MED-90036822-94_2



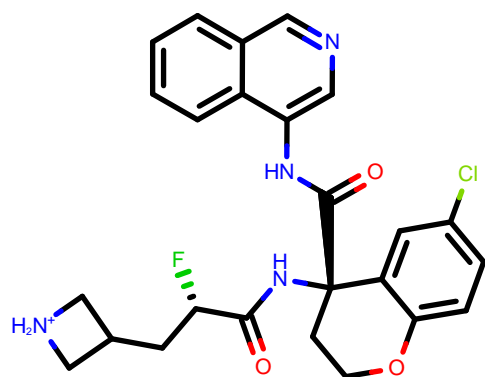
CID:	EDG-MED-90036822-94_2
SMILES:	<chem>C1C=NC(=O)N[C@@H]1C(=O)N[C@@H]2C=NC(=O)N=C2</chem>
RUN:	RUN1805
DDG (kcal/mol):	-1.43
dDDG (kcal/mol):	0.41

PET-UNK-29afea89-1_1



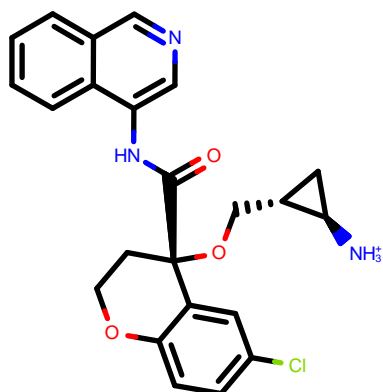
CID:	PET-UNK-29afea89-1_1
SMILES:	<chem>C#C[C@]1(COC(=O)N)C(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN657
DDG (kcal/mol):	-1.43
dDDG (kcal/mol):	0.34

ADA-UCB-dc2b944c-12_1



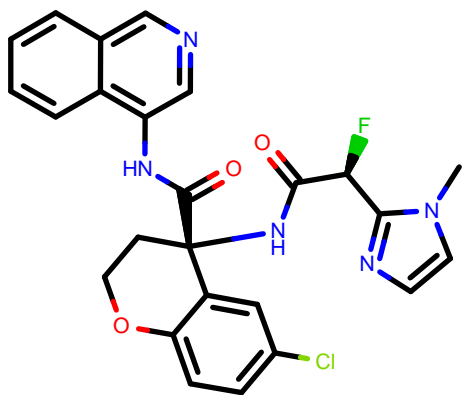
CID:	ADA-UCB-dc2b944c-12_1
SMILES:	<chem>Cc1c2cccc2c(cn1)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN609
DDG (kcal/mol):	-1.42
dDDG (kcal/mol):	0.32

DAR-DIA-23e5a6a0-1_2



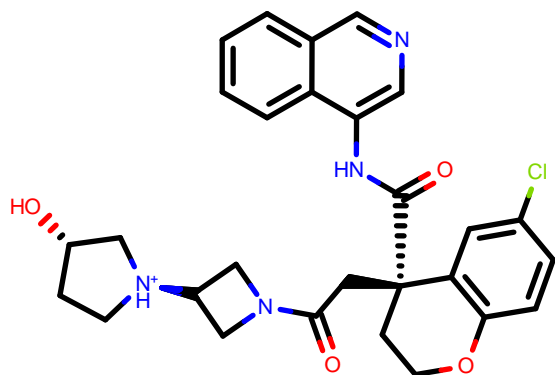
CID:	DAR-DIA-23e5a6a0-1_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN402
DDG (kcal/mol):	-1.42
dDDG (kcal/mol):	0.25

DAR-DIA-0587064e-17_1



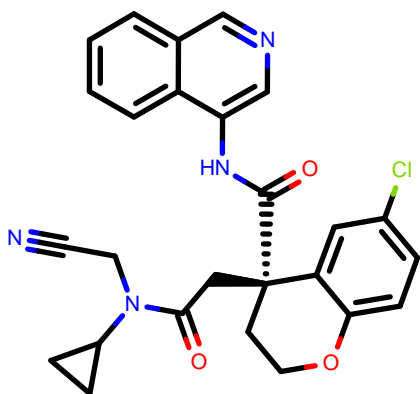
CID:	DAR-DIA-0587064e-17_1
SMILES:	<chem>COCCOC1cc(cc2c1OCC[C@@H]2C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN3369
DDG (kcal/mol):	-1.42
dDDG (kcal/mol):	0.16

ERI-UCB-d6de1f3c-2_1



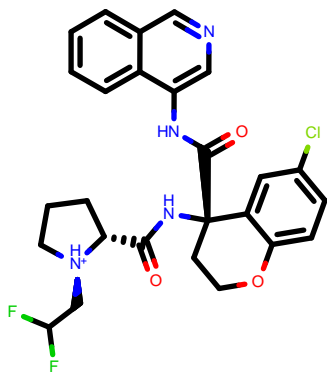
CID:	ERI-UCB-d6de1f3c-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2C(=O)N3CCN(C(=O)C3)c4cccc(c4)Cl</chem>
RUN:	RUN1091
DDG (kcal/mol):	-1.42
dDDG (kcal/mol):	0.33

DAR-DIA-5ff57136-12_1



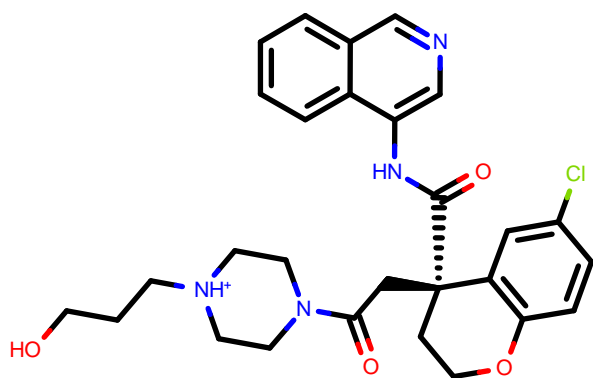
CID:	DAR-DIA-5ff57136-12_1
SMILES:	<chem>c1ccc2c(c1)cncc2N(C#CC[C@H]3CC[C@H]3)C(=O)[C@@H]4CCOC5c4ccc(c5)Cl</chem>
RUN:	RUN1382
DDG (kcal/mol):	-1.41
dDDG (kcal/mol):	0.17

MAK-UNK-ffc90da7-5_1



CID:	MAK-UNK-ffc90da7-5_1
SMILES:	<chem>C[C@@](CO)(c1ccc2c(c1)cncc2NC(=O)Cc3cccc(c3)Cl)[NH2+][C4CC4</chem>
RUN:	RUN702
DDG (kcal/mol):	-1.41
dDDG (kcal/mol):	0.35

ED_-GRI-5b13fbe2-57_1



CID: ED_-GRI-5b13fbe2-57_1

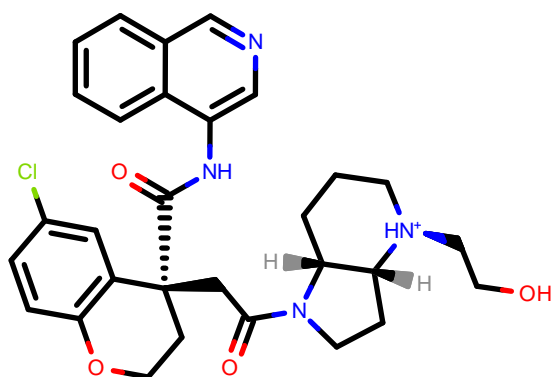
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OCc5c[nH]5

RUN: RUN1604

DDG (kcal/mol): -1.40

dDDG (kcal/mol): 0.31

ED_-GRI-5b13fbe2-42_1



CID: ED_-GRI-5b13fbe2-42_1

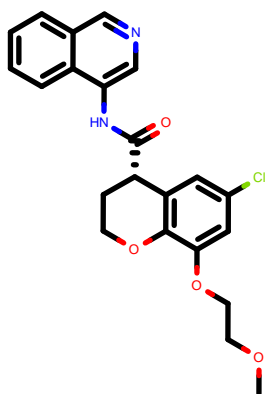
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OC[C@H](CF)[NH3+]

RUN: RUN1576

DDG (kcal/mol): -1.40

dDDG (kcal/mol): 0.60

EDJ-MED-d08626de-3_1



CID: EDJ-MED-d08626de-3_1

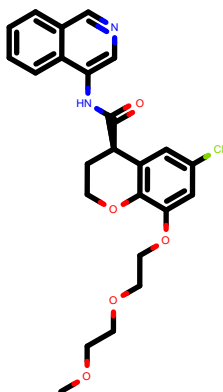
SMILES: CO[C@@]1(CCOc2c1cc(c(c2)Cl)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN2341

DDG (kcal/mol): -1.40

dDDG (kcal/mol): 0.24

KAD-UNI-8a629cb0-12_1



CID: KAD-UNI-8a629cb0-12_1

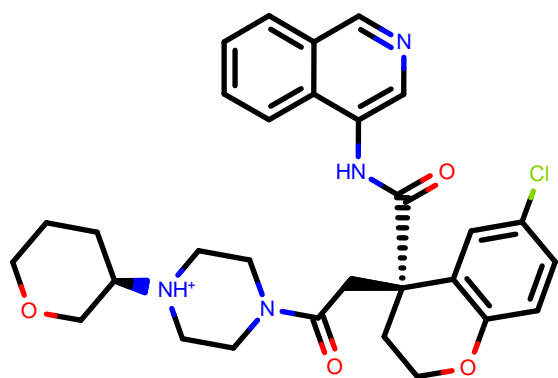
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CNC(=O)[C@H]5C6[nH+]cnc6C5

RUN: RUN2094

DDG (kcal/mol): -1.40

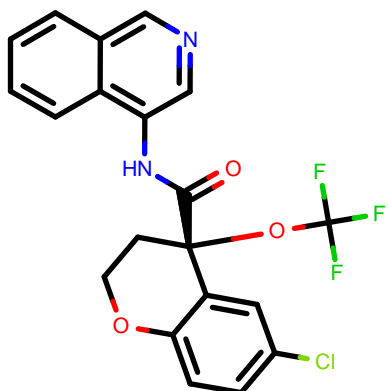
dDDG (kcal/mol): 0.36

BEN-DND-f2e727cd-6_2



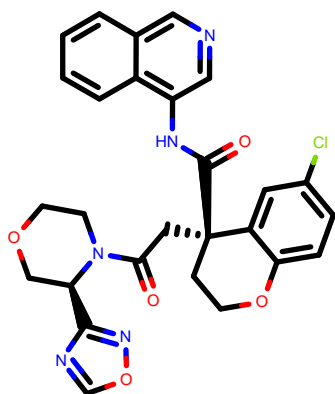
CID:	BEN-DND-f2e727cd-6_2
SMILES:	<chem>C[N@H+]1Cc2ccc(cc2[C@@H](C1)C(=O)Nc3ccc4c3cccc4)Cl</chem>
RUN:	RUN1199
DDG (kcal/mol):	-1.40
dDDG (kcal/mol):	0.26

MAT-POS-5369c344-5_2



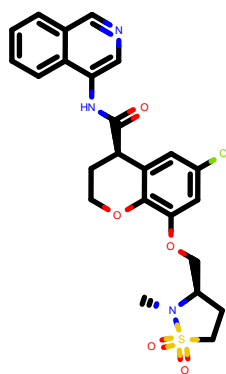
CID:	MAT-POS-5369c344-5_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H](C1)C(=O)Nc3ccc4c3cccc4)Cl</chem>
RUN:	RUN3623
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.07

KAD-UNI-b13decd3-7_1



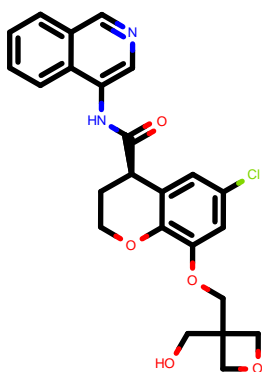
CID:	KAD-UNI-b13decd3-7_1
SMILES:	<chem>CC(=O)OCCn1cc(en1)C[NH2+]C[C@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4ccc5c4cccc5</chem>
RUN:	RUN3792
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.15

ALP-UNI-76695c4f-8_2



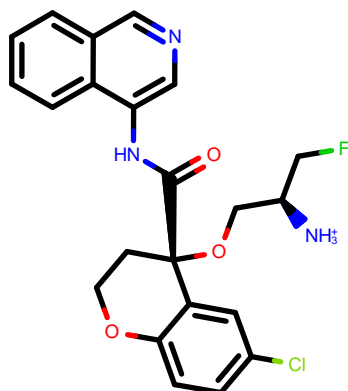
CID:	ALP-UNI-76695c4f-8_2
SMILES:	<chem>C1N@E)C1CN(C1)Cl(=O)C[C@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4ccc5c4cccc5)Si(=O)(=O)C</chem>
RUN:	RUN2179
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.43

KAD-UNI-8a629cb0-22_1



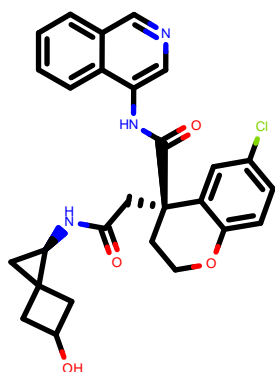
CID:	KAD-UNI-8a629cb0-22_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CCNC(=O)c5cc6c([nH]c(=O)[nH]6)nc5</chem>
RUN:	RUN2104
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.41

MAT-POS-8a69d52e-5_1



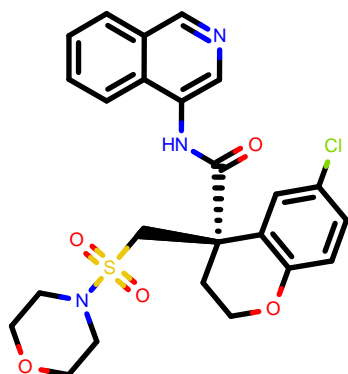
CID:	MAT-POS-8a69d52e-5_1
SMILES:	<chem>C[C@H]1COc2ccc(cc2[C@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN372
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.52

MAK-UNK-c749d764-24_1



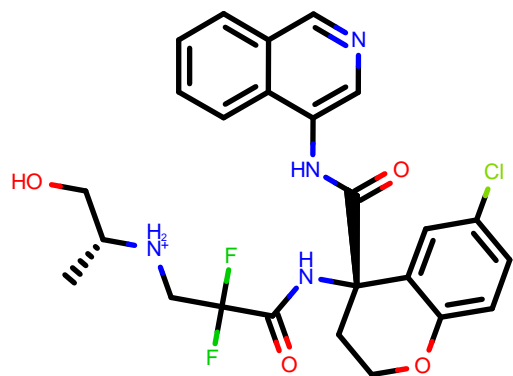
CID:	MAK-UNK-c749d764-24_1
SMILES:	<chem>CCN(c1cncc2c1cccc2)C(=O)C[C@@H]3CCC[C@@H]([C@@H]3O)C(F)F</chem>
RUN:	RUN1030
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.32

LAU-MED-88a3970a-6_1



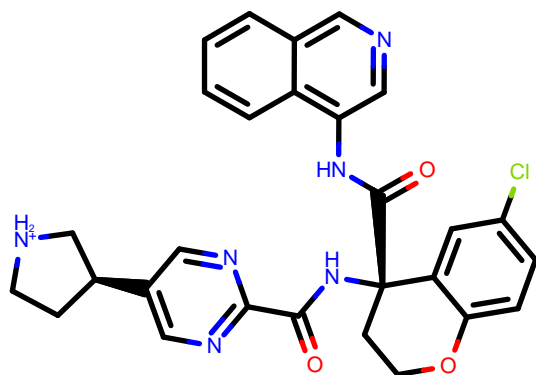
CID:	LAU-MED-88a3970a-6_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCOCc4c3cc(cc4CCCC)Cl</chem>
RUN:	RUN1500
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.18

MIC-UNK-cdc2493e-24_1



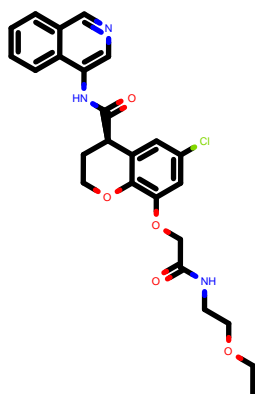
CID:	MIC-UNK-cdc2493e-24_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)N(CCn3cncn3)c4cccc(c4)Cl</chem>
RUN:	RUN575
DDG (kcal/mol):	-1.39
dDDG (kcal/mol):	0.43

PET-UNK-bb7ffe78-3_1



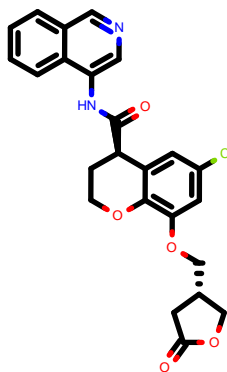
CID:	PET-UNK-bb7ffe78-3_1
SMILES:	<chem>CCCC1CC(Cc2c(Cl)CC(=O)Nc2cncc3c2ccccc3</chem>
RUN:	RUN3330
DDG (kcal/mol):	-1.38
dDDG (kcal/mol):	0.17

MAT-POS-fce787c2-6_1



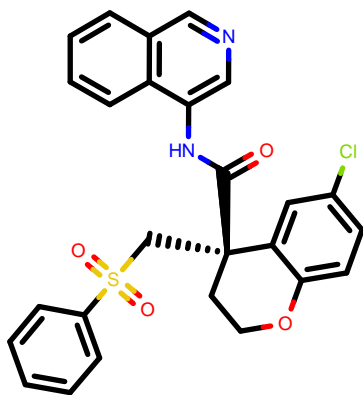
CID:	MAT-POS-fce787c2-6_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)C(c3ccc(c(c3)Cl)Cl)(F)F</chem>
RUN:	RUN2151
DDG (kcal/mol):	-1.38
dDDG (kcal/mol):	0.43

ALP-POS-2da19ca7-7_7



CID:	ALP-POS-2da19ca7-7_7
SMILES:	<chem>C[C@H]1CN(C)CC[C@H]1NC(=O)C(C)C2(COC2C3C(C)C)C1=O)Nc4nc5c4cccc5(S(=O)(=O)C</chem>
RUN:	RUN2388
DDG (kcal/mol):	-1.38
dDDG (kcal/mol):	0.48

ED_-GRI-5b13fbe2-45_1



CID: ED_-GRI-5b13fbe2-45_1

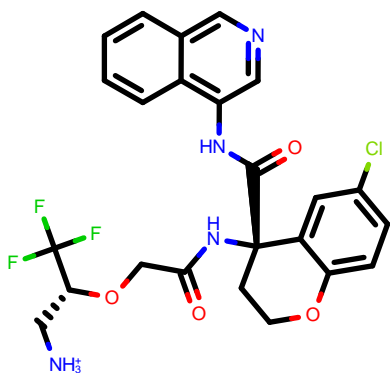
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@H]3[C@@H](COCc4c3cc(c4)C)OC[C@@H]3C5c6c(nc6)COC5C7=CC=CC=C7

RUN: RUN1584

DDG (kcal/mol): -1.38

dDDG (kcal/mol): 0.32

EDG-MED-70ae9412-2_2



CID: EDG-MED-70ae9412-2_2

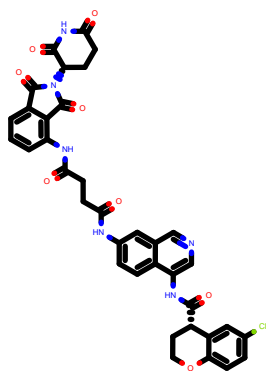
SMILES: C[NH+](C)CC(=O)NC[C@]1(C)COCc2c1cc(cc2)C(=O)Nc3cnc4c3cccc4

RUN: RUN3168

DDG (kcal/mol): -1.37

dDDG (kcal/mol): 0.39

EDJ-MED-50011917-1_2



CID: EDJ-MED-50011917-1_2

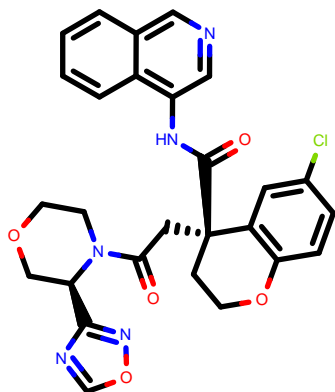
SMILES: c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)C[C@@H]4CC(=O)N4

RUN: RUN371

DDG (kcal/mol): -1.37

dDDG (kcal/mol): 0.31

MAK-UNK-c749d764-24_5



CID: MAK-UNK-c749d764-24_5

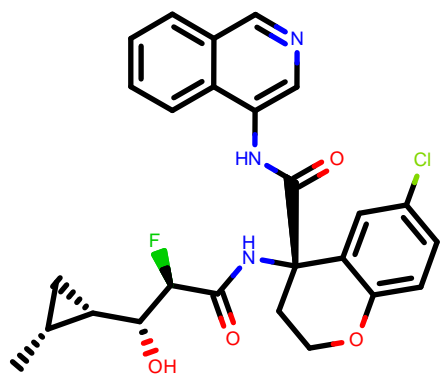
SMILES: CCN(c1cnc2c1cccc2)C(=O)C[C@@H]3CCC[C@@H]3C[C@@H]4C(F)F

RUN: RUN1034

DDG (kcal/mol): -1.37

dDDG (kcal/mol): 0.36

ADA-UCB-dc2b944c-10_1



CID: ADA-UCB-dc2b944c-10_1

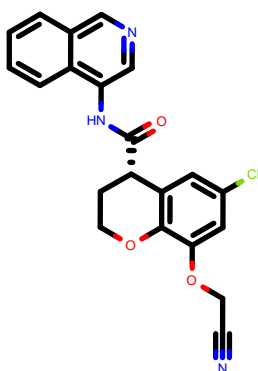
SMILES: Cc1c(c2ccccc2cn1)NC(=O)[C@@H]3CCOCc4ccc(cc4)Cl

RUN: RUN608

DDG (kcal/mol): -1.36

dDDG (kcal/mol): 0.25

ALF-EVA-07677224-4_11



CID: ALF-EVA-07677224-4_11

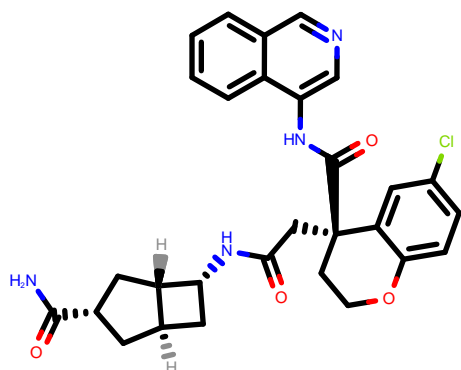
SMILES: C[C@@H]1[C@H](CO)S(=O)(=O)N1[C@@H]2C=CC(=O)N2[C@@H]3C=CC(=O)N3

RUN: RUN4918

DDG (kcal/mol): -1.36

dDDG (kcal/mol): 0.29

MAK-UNK-c749d764-18_5



CID: MAK-UNK-c749d764-18_5

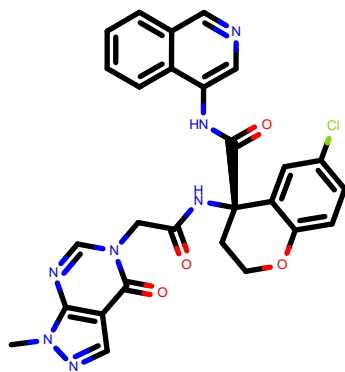
SMILES: CC(C)SCN(c1ccc2c1ccc2)C(=O)C[C@@H]3CCC[C@@H]1[C@@H]3O1C(F)F

RUN: RUN994

DDG (kcal/mol): -1.36

dDDG (kcal/mol): 0.36

LON-WEI-5e7d1b3e-43_1



CID: LON-WEI-5e7d1b3e-43_1

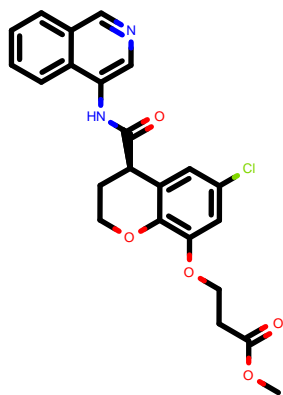
SMILES: Cc1cc(c(cc1Cl)OC)NC(=O)Nc2cn(c(=O)c3c2cccc3)CC(C)C

RUN: RUN1350

DDG (kcal/mol): -1.36

dDDG (kcal/mol): 0.28

EDJ-MED-40433386-8_1



CID: EDJ-MED-40433386-8_1

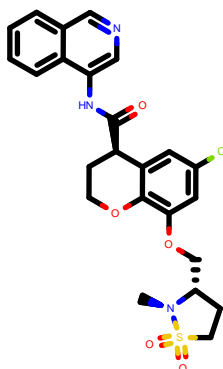
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CNC(=O)CC@5(CCOCC5)C#N

RUN: RUN2559

DDG (kcal/mol): -1.35

dDDG (kcal/mol): 0.35

CHO-MSK-a31cca77-1_1



CID: CHO-MSK-a31cca77-1_1

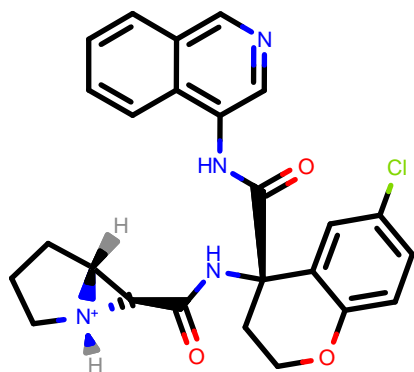
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CC(=O)Nc5nncoc5

RUN: RUN2191

DDG (kcal/mol): -1.35

dDDG (kcal/mol): 0.39

LON-WEI-9739a092-10_1



CID: LON-WEI-9739a092-10_1

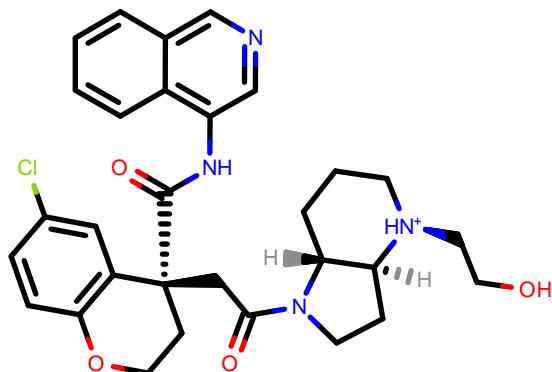
SMILES: COc1c(cc(cn1)Br)Nc2cc(cc(c2)Cl)CC(=O)Nc3ncc4c3cccc4

RUN: RUN3276

DDG (kcal/mol): -1.34

dDDG (kcal/mol): 0.16

DAR-DIA-5ff57136-9_1



CID: DAR-DIA-5ff57136-9_1

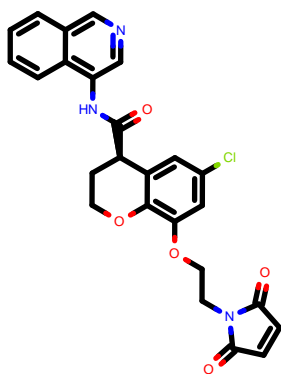
SMILES: COC(=O)/C=C/C(=O)N(c1cncc2c1cccc2)C(=O)[C@@]3(CCOc4c3cc(cc4)Cl)C

RUN: RUN1378

DDG (kcal/mol): -1.33

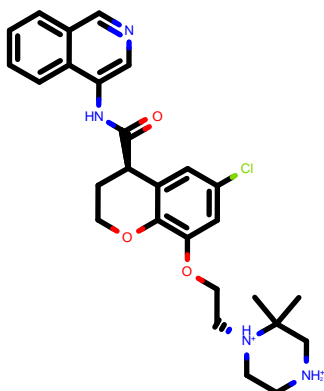
dDDG (kcal/mol): 0.53

ALF-EVA-07677224-1_2



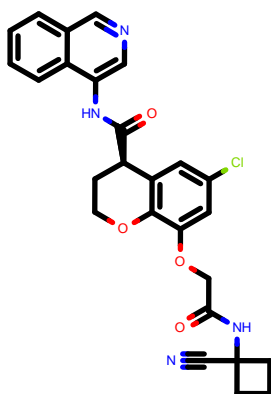
CID:	ALF-EVA-07677224-1_2
SMILES:	<chem>Cc1nc(ns1)S(=O)(=O)[N@]2Cc3ccc(cc3[C@@H](C2)C(=O)Nc4cncc5c4cccc5)Cl</chem>
RUN:	RUN4895
DDG (kcal/mol):	-1.33
dDDG (kcal/mol):	0.19

MAT-POS-de59a476-2_2



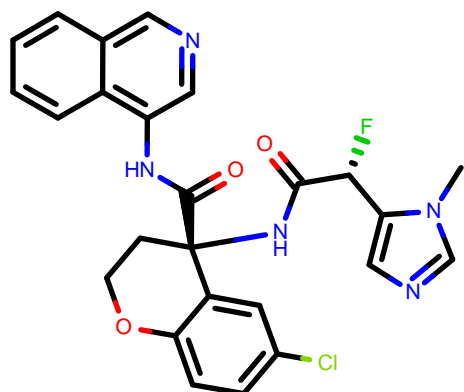
CID:	MAT-POS-de59a476-2_2
SMILES:	<chem>CO[C@H](c1ccc(c(c1)Cl)Cl)C(=O)Nc2cncc3c2cccc3</chem>
RUN:	RUN2218
DDG (kcal/mol):	-1.33
dDDG (kcal/mol):	0.55

EDG-MED-ba1ac7b9-3_1



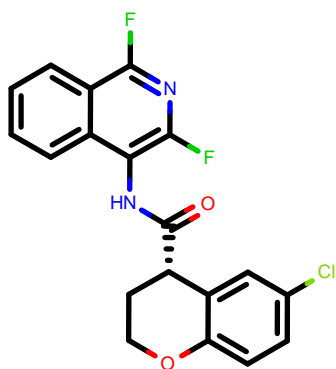
CID:	EDG-MED-ba1ac7b9-3_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(c4)Cl)CC(=O)N(CCO)C5CC5</chem>
RUN:	RUN2624
DDG (kcal/mol):	-1.33
dDDG (kcal/mol):	0.51

EDJ-MED-00143744-1_1



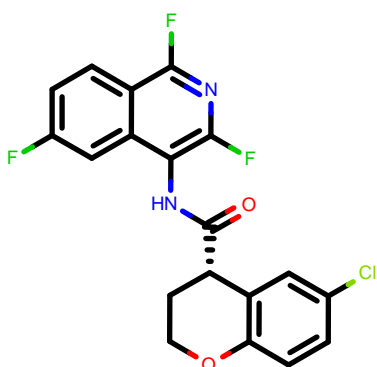
CID:	EDJ-MED-00143744-1_1
SMILES:	<chem>Cn1c(nnn1)COC[C@@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4cncc5c4cccc5</chem>
RUN:	RUN3166
DDG (kcal/mol):	-1.33
dDDG (kcal/mol):	0.16

LON-WEI-4d77710c-1_1



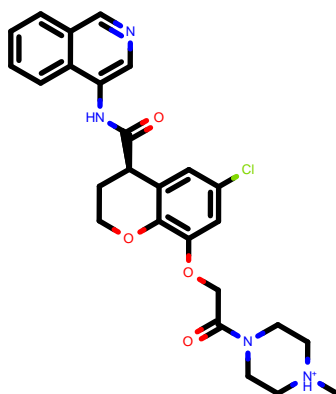
CID:	LON-WEI-4d77710c-1_1
SMILES:	<chem>Cn1cc(c2ccccc2c1=O)NC(=O)NCc3cc(ccc3OC)OC</chem>
RUN:	RUN193
DDG (kcal/mol):	-1.33
dDDG (kcal/mol):	0.11

VLA-UCB-1dbca3b4-15_2



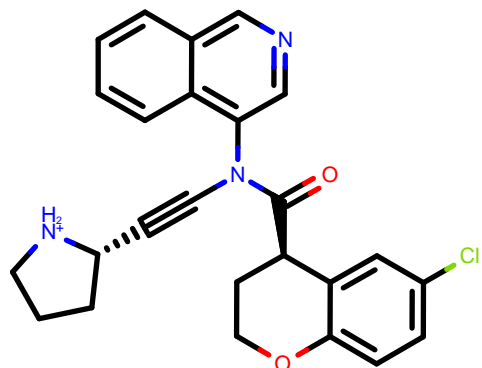
CID:	VLA-UCB-1dbca3b4-15_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN172
DDG (kcal/mol):	-1.32
dDDG (kcal/mol):	0.11

EDJ-MED-d203f206-24_1



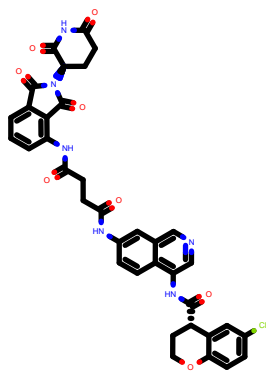
CID:	EDJ-MED-d203f206-24_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCO[C@]6H[C@]5(C)Cn6ccnn6</chem>
RUN:	RUN2588
DDG (kcal/mol):	-1.32
dDDG (kcal/mol):	0.40

ALP-UNI-8e43a71e-2_5



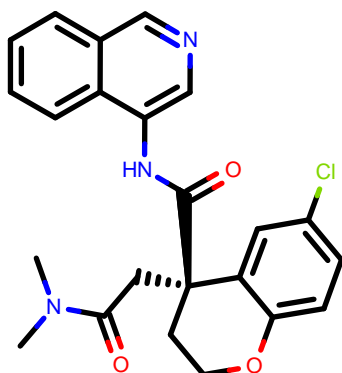
CID:	ALP-UNI-8e43a71e-2_5
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCO[C@]6H[C@]5(C)Cn6ccnn6</chem>
RUN:	RUN2928
DDG (kcal/mol):	-1.32
dDDG (kcal/mol):	0.18

ALP-UNI-8e43a71e-5_3



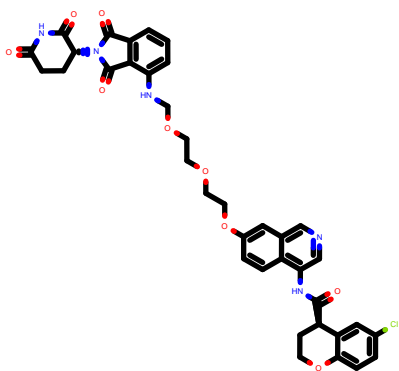
CID:	ALP-UNI-8e43a71e-5_3
SMILES:	<chem>C1C@H](CN@)CC1C@H1NC(=O)C1C2(CCO3C2CC(C3)C1=O)N4NCC5C4CC5S(=O)(=O)C</chem>
RUN:	RUN2946
DDG (kcal/mol):	-1.32
dDDG (kcal/mol):	0.42

MAK-UNK-ffc90da7-6_2



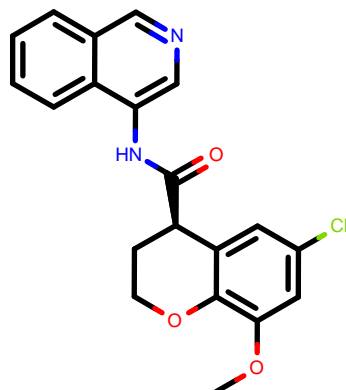
CID:	MAK-UNK-ffc90da7-6_2
SMILES:	<chem>C1C@H](C[NH2+]CCCO)c1ccc2nccc(c2c1)NC(=O)Cc3ccccc3Cl</chem>
RUN:	RUN706
DDG (kcal/mol):	-1.31
dDDG (kcal/mol):	0.25

JAG-UCB-f37eaa14-1_1



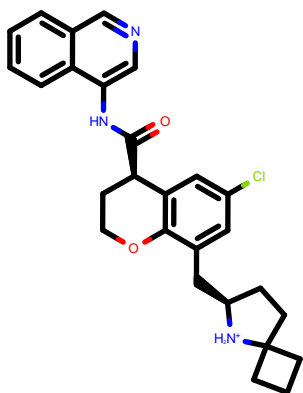
CID:	JAG-UCB-f37eaa14-1_1
SMILES:	<chem>c1ccc2c(c1)cnc2N3CC[C@@]4(C3=O)COc5c4cc(c(c5)F)Cl</chem>
RUN:	RUN3055
DDG (kcal/mol):	-1.31
dDDG (kcal/mol):	0.30

ERI-UCB-ce40166b-10_2



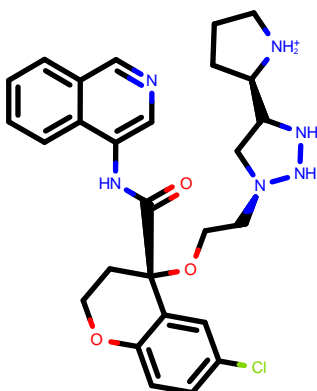
CID:	ERI-UCB-ce40166b-10_2
SMILES:	<chem>c1ccc2c(c1)cnc2CC(=O)Nc3cc(cc(c3)O)[C@H]4CCC(=O)N4)C#N</chem>
RUN:	RUN49
DDG (kcal/mol):	-1.31
dDDG (kcal/mol):	0.24

MIC-UNK-66895286-5_1



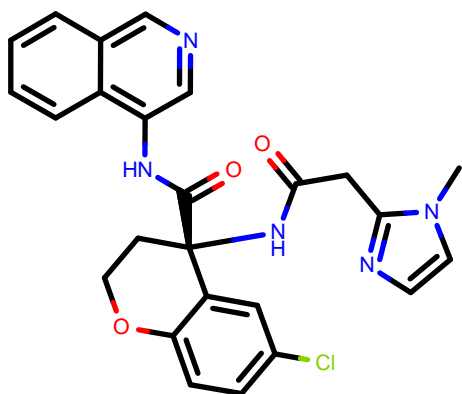
CID:	MIC-UNK-66895286-5_1
SMILES:	<chem>c1ccc2c(c1)c(c[nH]c2=O)CC(=O)Nc3cccc(c3)Cl</chem>
RUN:	RUN31
DDG (kcal/mol):	-1.31
dDDG (kcal/mol):	0.43

MAK-UNK-8be7dca9-1_2



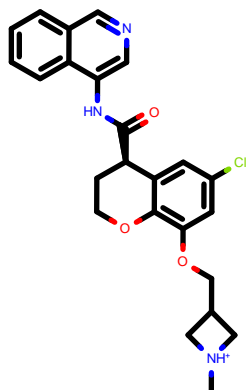
CID:	MAK-UNK-8be7dca9-1_2
SMILES:	<chem>c1cc2c(cc1CN3CC[NH2+][CC3]cnc2NC(=O)[C@@H]4CCOc5c4cc(cc5)Cl</chem>
RUN:	RUN493
DDG (kcal/mol):	-1.30
dDDG (kcal/mol):	0.43

EDJ-MED-f893e2a1-6_1



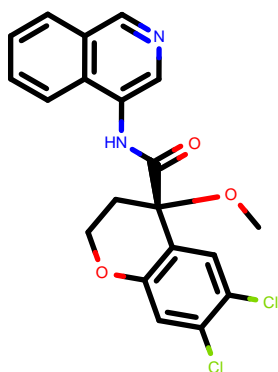
CID:	EDJ-MED-f893e2a1-6_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)C[NH2+][C]C@@H]5CC(=O)NC5</chem>
RUN:	RUN3203
DDG (kcal/mol):	-1.29
dDDG (kcal/mol):	0.14

ALF-EVA-5b152d2f-4_1



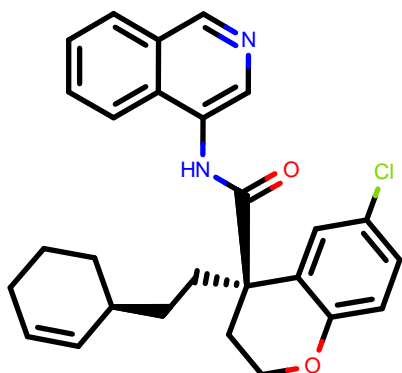
CID:	ALF-EVA-5b152d2f-4_1
SMILES:	<chem>c1ccc2c(c1)cnc2NNC(=O)[C@@]3(CCOc4c3cc(cc4)Cl</chem>
RUN:	RUN2354
DDG (kcal/mol):	-1.29
dDDG (kcal/mol):	0.37

DAR-DIA-0d514e7d-31_14



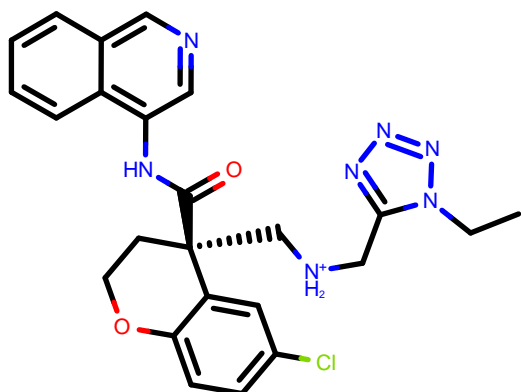
CID:	DAR-DIA-0d514e7d-31_14
SMILES:	<chem>C[C@H]1CCO[C@@H]2C=CC(=C[C@H]2[C@H]1C(=O)Nc3ccc(Cl)c(Cl)c3</chem>
RUN:	RUN847
DDG (kcal/mol):	-1.29
dDDG (kcal/mol):	0.25

DAR-DIA-0cde14eb-57_2



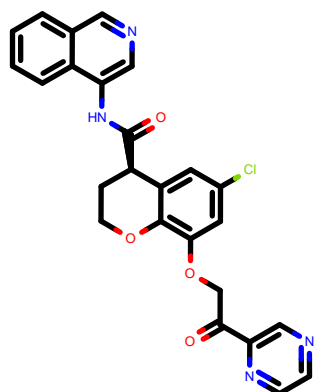
CID:	DAR-DIA-0cde14eb-57_2
SMILES:	<chem>C[C@H](c1cccc(c1)C2(CC2)F)C(=O)Nc3ccc(Cl)cc3</chem>
RUN:	RUN26
DDG (kcal/mol):	-1.28
dDDG (kcal/mol):	0.27

MIC-UNK-ea4eb352-6_1



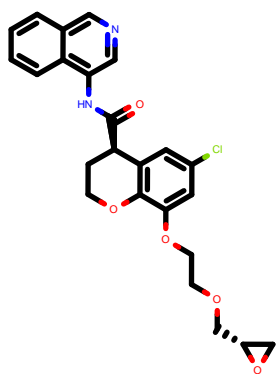
CID:	MIC-UNK-ea4eb352-6_1
SMILES:	<chem>CCNCCN[C@@H]1CCO[C@@H]2C=CC(=C[C@H]2[C@H]1C(=O)Nc3ccc(Cl)cc3</chem>
RUN:	RUN4646
DDG (kcal/mol):	-1.27
dDDG (kcal/mol):	0.29

ALP-UNI-3496895b-15_3



CID:	ALP-UNI-3496895b-15_3
SMILES:	<chem>c1ccc2c(c1)ncnc2N[C@@H]1CCO[C@@H]2C=CC(=C[C@H]2[C@H]1C(=O)Nc3ccc(Cl)cc3</chem>
RUN:	RUN2533
DDG (kcal/mol):	-1.27
dDDG (kcal/mol):	0.31

MAT-POS-f9802937-5_1



CID: MAT-POS-f9802937-5_1

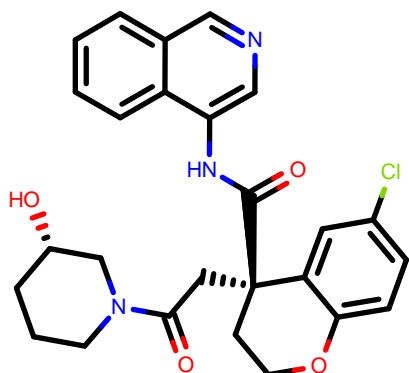
SMILES: CNC(=O)C[C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cnc4c3ccc4

RUN: RUN2396

DDG (kcal/mol): -1.27

dDDG (kcal/mol): 0.37

BEN-BAS-5c03e89e-1_1



CID: BEN-BAS-5c03e89e-1_1

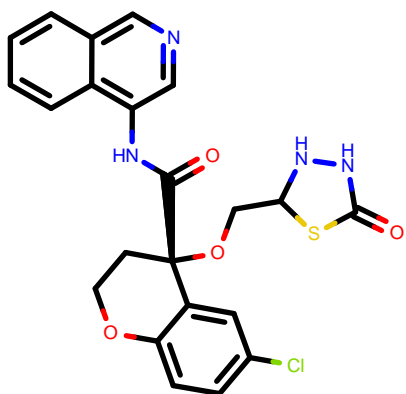
SMILES: CC(C)([C@@H]1COc2ccc(cc2[C@@H]1C(=O)Nc3cnc4c3ccc4)Cl)O

RUN: RUN1143

DDG (kcal/mol): -1.27

dDDG (kcal/mol): 0.27

NIR-THE-af15c15d-1_2



CID: NIR-THE-af15c15d-1_2

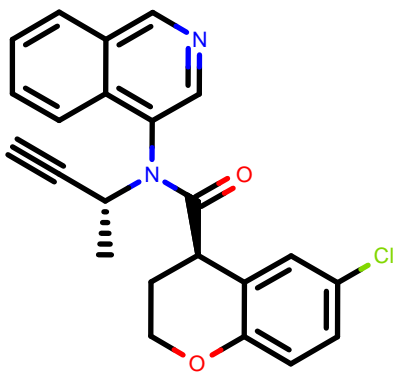
SMILES: CC#CC(=O)N(c1cnc2c1ccc2)C(=O)[C@H]3COc4c3cc(cc4)Cl

RUN: RUN3320

DDG (kcal/mol): -1.27

dDDG (kcal/mol): 0.13

ALP-UNI-8e43a71e-15_22



CID: ALP-UNI-8e43a71e-15_22

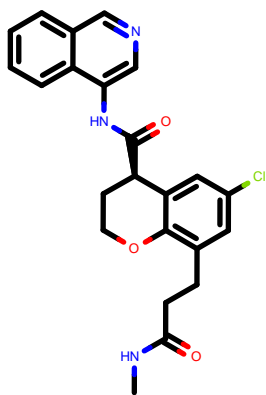
SMILES: c1ccc2c(c1)cnc2NCl(-O)[C@]3(CCOc4ccc(cc4)Cl)CC(-O)NC@@H5C[C@H]6[C@]3([C@H]5)C6(-)N

RUN: RUN3003

DDG (kcal/mol): -1.27

dDDG (kcal/mol): 0.12

DAR-DIA-6a508060-5_2



CID: DAR-DIA-6a508060-5_2

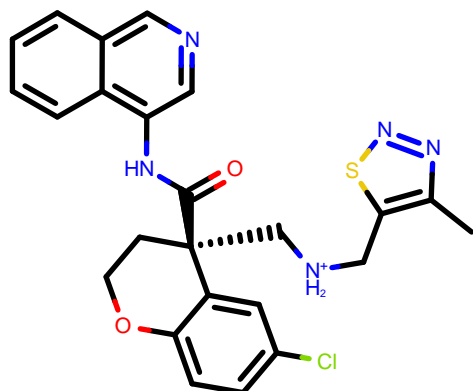
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOC4c3cc(cc4C)[C@@H]5CC(=O)N5Cl

RUN: RUN341

DDG (kcal/mol): -1.26

dDDG (kcal/mol): 0.30

MAT-POS-5f1400cf-1_6



CID: MAT-POS-5f1400cf-1_6

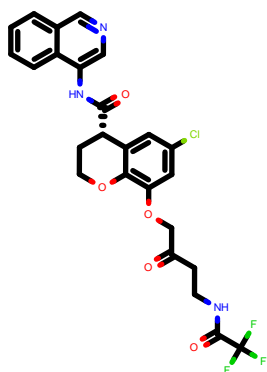
SMILES: Cc1cc(en1)N2CC[C@@H](C2)C[NH2+][C@@]3(CCO4c3cc(cc4)Cl)C(=O)Nc5ncc6c5ccc6

RUN: RUN4511

DDG (kcal/mol): -1.26

dDDG (kcal/mol): 0.25

MAT-POS-e9e99895-8_4



CID: MAT-POS-e9e99895-8_4

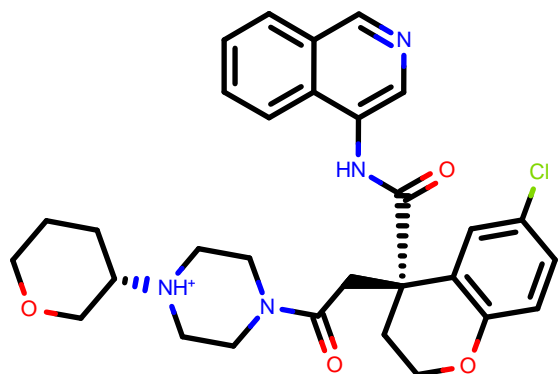
SMILES: Cc1nc2n(n1)C[C@@H](CC2)C(=O)N[C@@]3(C)C3cc(c(c3)Cl)C(=O)Nc4ncc5c4cccc5

RUN: RUN2257

DDG (kcal/mol): -1.26

dDDG (kcal/mol): 0.26

BEN-DND-c852c98b-4_1



CID: BEN-DND-c852c98b-4_1

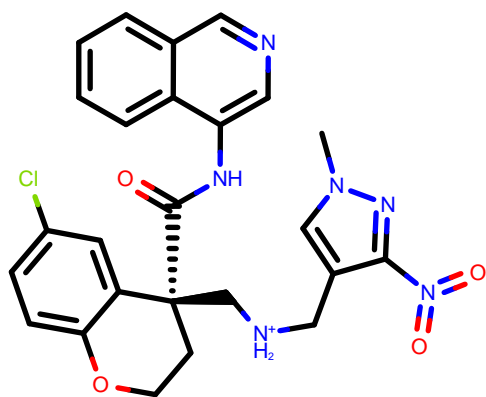
SMILES: c1cc2cncc(c2cc1O)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl

RUN: RUN1207

DDG (kcal/mol): -1.25

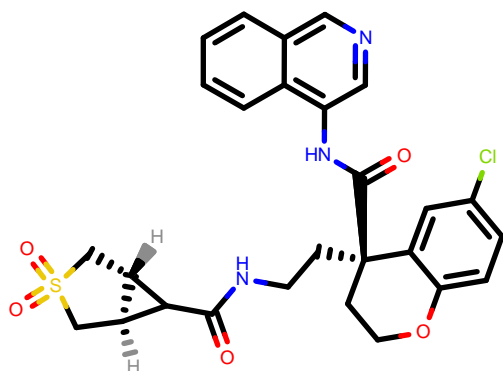
dDDG (kcal/mol): 0.31

EDG-MED-90036822-36_2



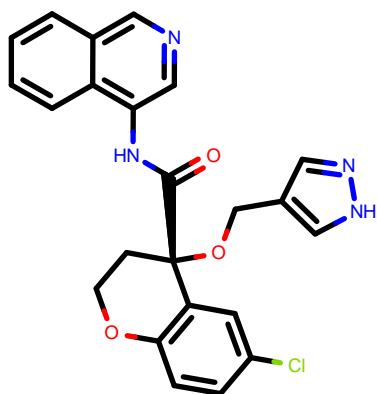
CID:	EDG-MED-90036822-36_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)C)NC(=O)[C@H](n5cnc5)F</chem>
RUN:	RUN1711
DDG (kcal/mol):	-1.25
dDDG (kcal/mol):	0.46

DAR-DIA-f6ee7aeb-5_1



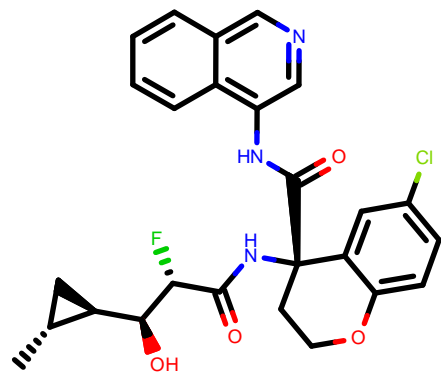
CID:	DAR-DIA-f6ee7aeb-5_1
SMILES:	<chem>CCCOc1cc(cc(c1)Cl)[C@@]2C(=O)N[C]C@@[H]2c3c[nH]c(=O)[nH]c3=O)c4cnc5c4cccc5</chem>
RUN:	RUN3414
DDG (kcal/mol):	-1.25
dDDG (kcal/mol):	0.17

MAT-POS-3b92565d-7_1



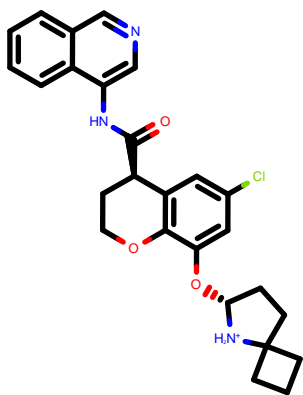
CID:	MAT-POS-3b92565d-7_1
SMILES:	<chem>Cn1cc(c2ccccc2c1=O)NC(=O)Cc3cc(ccc3OC)Cl</chem>
RUN:	RUN113
DDG (kcal/mol):	-1.25
dDDG (kcal/mol):	0.30

LON-WEI-9739a092-4_2



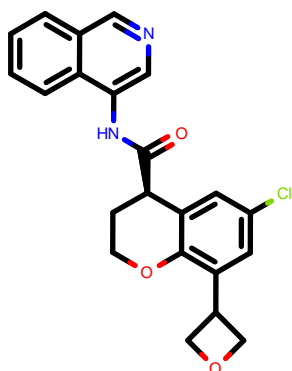
CID:	LON-WEI-9739a092-4_2
SMILES:	<chem>CCN(Cc1ccc(cc1)F)c2cc(cc(c2)Cl)CC(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN3272
DDG (kcal/mol):	-1.24
dDDG (kcal/mol):	0.15

MAT-POS-b5746674-102_1



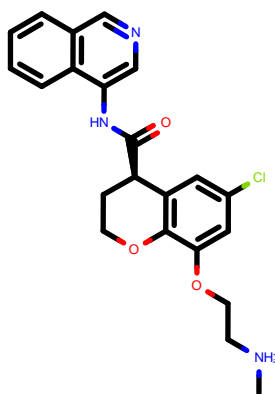
CID:	MAT-POS-b5746674-102_1
SMILES:	<chem>CCCCN(CCCNC(=O)Nc1cn(c(=O)c2c1cccc2)CC(C)C)c3cccc3</chem>
RUN:	RUN83
DDG (kcal/mol):	-1.24
dDDG (kcal/mol):	0.41

DAR-DIA-0cde14eb-55_1



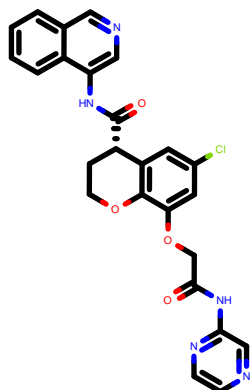
CID:	DAR-DIA-0cde14eb-55_1
SMILES:	<chem>C[C@@H](c1cccc(c1)C2(CC2)C)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN21
DDG (kcal/mol):	-1.24
dDDG (kcal/mol):	0.24

KAD-UNI-8a629cb0-11_1



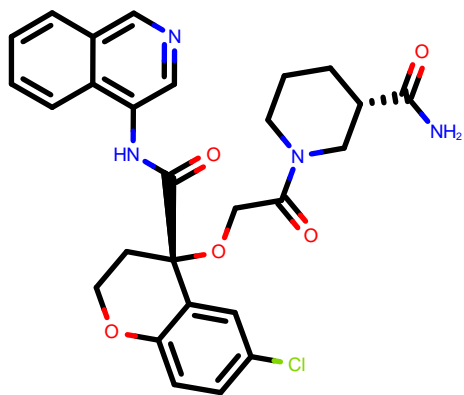
CID:	KAD-UNI-8a629cb0-11_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(C)CCOc4c3cc(cc4)C)CCNC(=O)Cc5c(=O)[nH]enc5[O-]</chem>
RUN:	RUN2096
DDG (kcal/mol):	-1.24
dDDG (kcal/mol):	0.55

MAT-POS-2905de8c-3_2



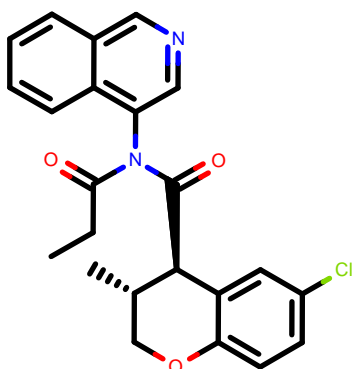
CID:	MAT-POS-2905de8c-3_2
SMILES:	<chem>C[NH+](C)[C@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN2232
DDG (kcal/mol):	-1.23
dDDG (kcal/mol):	0.23

MIC-UNK-d854bf4c-4_1



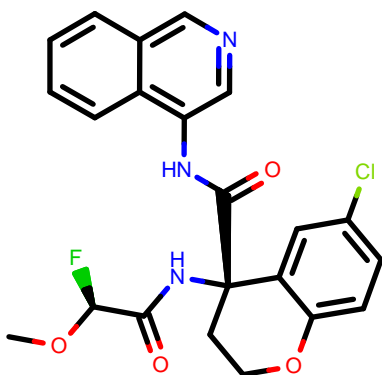
CID:	MIC-UNK-d854bf4c-4_1
SMILES:	<chem>CC(=O)N1CCC2(CC1)CCN(C(=O)[C@@H]2c3ccc(c(c3)Cl)Cl)c4ccc5c4ccc5</chem>
RUN:	RUN3337
DDG (kcal/mol):	-1.23
dDDG (kcal/mol):	0.21

ALP-UNI-8e43a71e-1_4



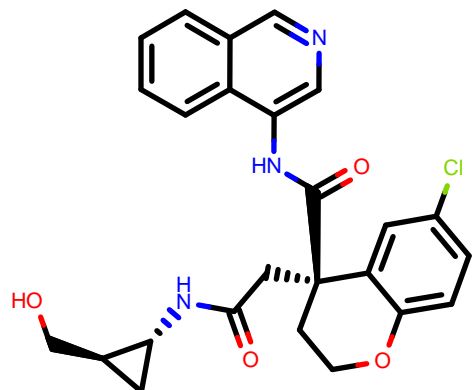
CID:	ALP-UNI-8e43a71e-1_4
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@H]3[C@@H](COCc4ccc(cc4)Cl)C(=O)N5CCCC5</chem>
RUN:	RUN2922
DDG (kcal/mol):	-1.23
dDDG (kcal/mol):	0.13

MAT-POS-fb82b63d-1_1



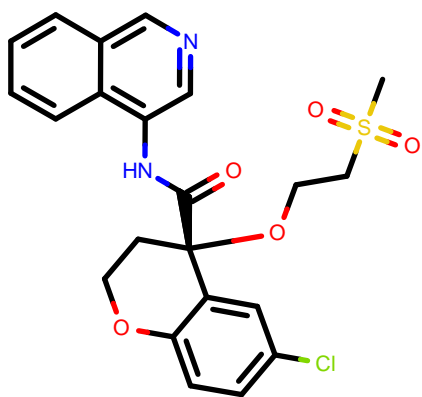
CID:	MAT-POS-fb82b63d-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3c4cc(ccc4CC[NH2+])3]Cl</chem>
RUN:	RUN3163
DDG (kcal/mol):	-1.23
dDDG (kcal/mol):	0.10

MAK-UNK-c749d764-33_7



CID:	MAK-UNK-c749d764-33_7
SMILES:	<chem>CS(=O)(=O)N(c1ccc2c1cccc2)C(=O)C[C@@H]3CCC[C@H]1[C@H]3O)C(F)F</chem>
RUN:	RUN1085
DDG (kcal/mol):	-1.23
dDDG (kcal/mol):	0.31

MAT-POS-4223bc15-41_2



CID: MAT-POS-4223bc15-41_2

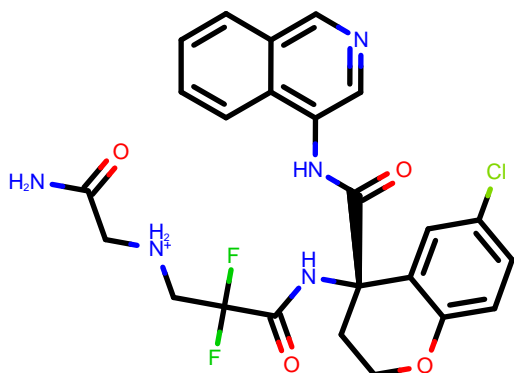
SMILES: Cn1cnc1C(=O)N2Cc3ccc(cc3[C@H](C2)C(=O)Nc4cnc5c4cccc5)Cl

RUN: RUN4164

DDG (kcal/mol): -1.23

dDDG (kcal/mol): 0.16

MIC-UNK-bcd487e9-6_2



CID: MIC-UNK-bcd487e9-6_2

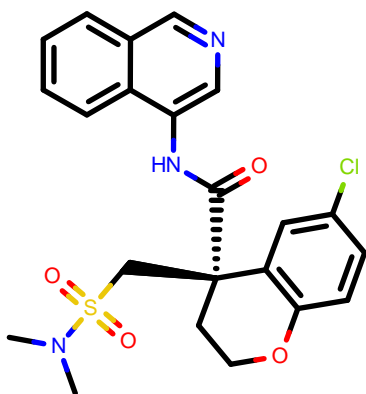
SMILES: c1ccc2c(c1)cnc2NC(=O)N[C@H](C)C3CCCS3(=O)=O)c4cccc(c4)Cl

RUN: RUN593

DDG (kcal/mol): -1.22

dDDG (kcal/mol): 0.42

LAU-MED-88a3970a-2_1



CID: LAU-MED-88a3970a-2_1

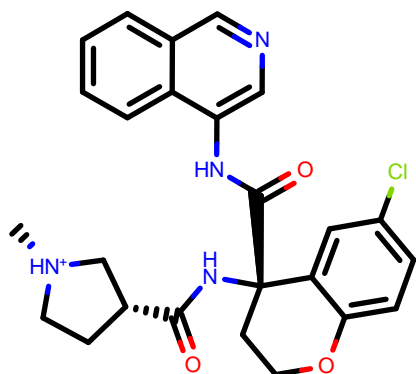
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H](C)C3CCO4c3cc(cc4OCCO)Cl

RUN: RUN1498

DDG (kcal/mol): -1.22

dDDG (kcal/mol): 0.16

RAL-THA-c9f97604-1_1



CID: RAL-THA-c9f97604-1_1

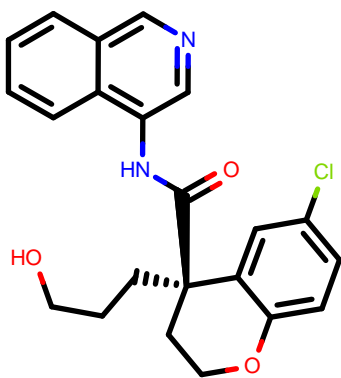
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H](C)C3c4ccc4[C@H]5CC[C@H]3O5)Cl

RUN: RUN584

DDG (kcal/mol): -1.22

dDDG (kcal/mol): 0.36

ALP-UNI-8e43a71e-5_6



CID: ALP-UNI-8e43a71e-5_6

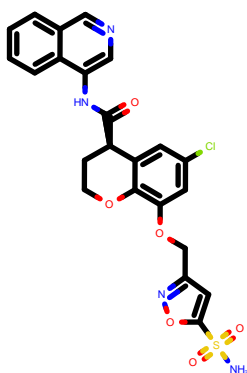
SMILES: C1C@H]1CN@@]CC[C@H]1NC(-O)C[C@@]2(CCOc3c2cc(c3)C)C1-O)Nc4nc5c4ccc5)S(-O)(-O)C

RUN: RUN2950

DDG (kcal/mol): -1.22

dDDG (kcal/mol): 0.10

EDJ-MED-fcba3f31-6_1



CID: EDJ-MED-fcba3f31-6_1

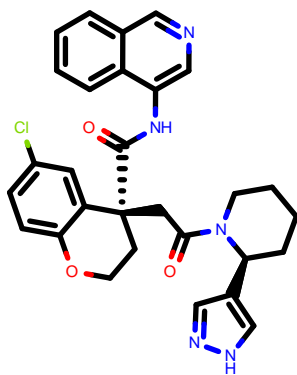
SMILES: c1ccc2c(c1)ncoc2NC(=O)[C@@]3(CCOc4c3cc(c4)C)COC[C@H]5CNCN5=O

RUN: RUN2542

DDG (kcal/mol): -1.22

dDDG (kcal/mol): 0.38

BEN-DND-c852c98b-3_2



CID: BEN-DND-c852c98b-3_2

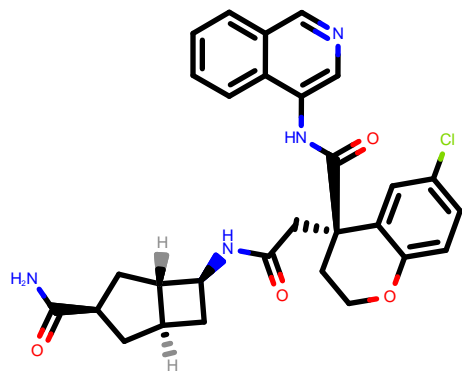
SMILES: CC(C)(c1ccc2cnccc(c2c1)NC(=O)[C@H]3CCOc4c3cc(cc4)Cl)O

RUN: RUN1209

DDG (kcal/mol): -1.21

dDDG (kcal/mol): 0.41

LON-WEI-5e7d1b3e-35_1



CID: LON-WEI-5e7d1b3e-35_1

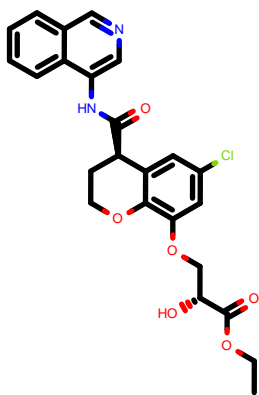
SMILES: Cc1ccc(cc1)C[N@@H]2CC[C@@H]1(C2)CNC(=O)Nc3cn(c(=O)c4c3ccc4)CC(C)C

RUN: RUN1342

DDG (kcal/mol): -1.21

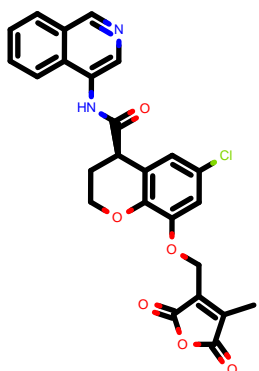
dDDG (kcal/mol): 0.37

CHO-MSK-a31cca77-4_2



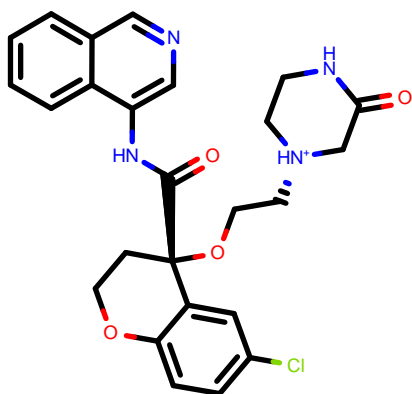
CID:	CHO-MSK-a31cca77-4_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CC(=O)Nc5ncon5</chem>
RUN:	RUN2198
DDG (kcal/mol):	-1.21
dDDG (kcal/mol):	0.44

ALP-UNI-3496895b-5_6



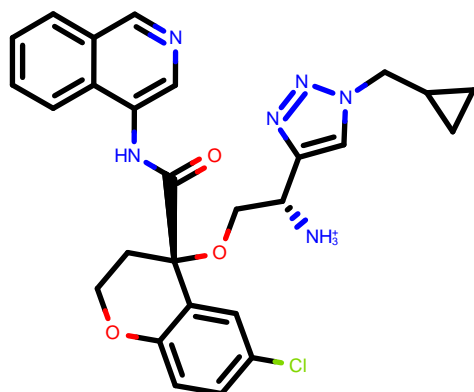
CID:	ALP-UNI-3496895b-5_6
SMILES:	<chem>C[C@H]1C[N@@](C)C[C@@H]1NC(=O)[C@]2(COC3C2cc(c3)Cl)C(=O)Nc4nc5c4cccc5(Si(=O)(=O)C</chem>
RUN:	RUN2517
DDG (kcal/mol):	-1.21
dDDG (kcal/mol):	0.36

DAR-DIA-6a508060-11_2



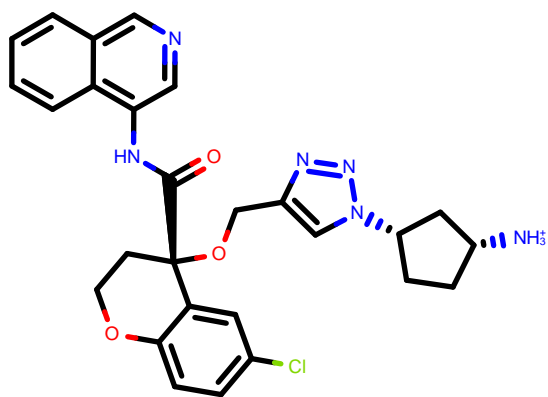
CID:	DAR-DIA-6a508060-11_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)C3CC(C3)[C@H]4C[C@@H]4Cl</chem>
RUN:	RUN347
DDG (kcal/mol):	-1.21
dDDG (kcal/mol):	0.43

MAT-POS-8a69d52e-4_1



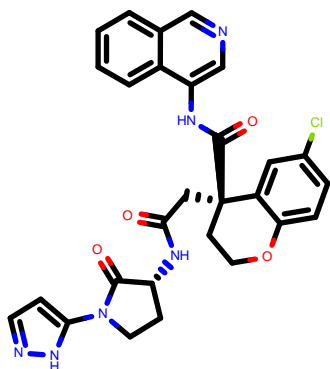
CID:	MAT-POS-8a69d52e-4_1
SMILES:	<chem>C[C@@H]1COc2ccc(cc2[C@@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN366
DDG (kcal/mol):	-1.20
dDDG (kcal/mol):	0.46

ALP-POS-fe871b40-2_2



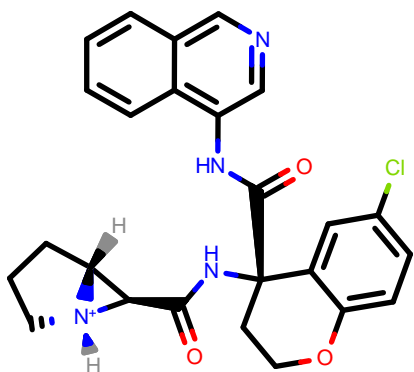
CID:	ALP-POS-fe871b40-2_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@H]3CCNc4c3cc(cc4C#N)Cl</chem>
RUN:	RUN3108
DDG (kcal/mol):	-1.20
dDDG (kcal/mol):	0.22

LON-WEI-5e7d1b3e-35_4



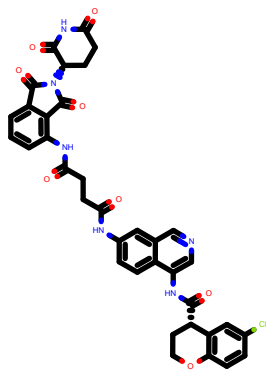
CID:	LON-WEI-5e7d1b3e-35_4
SMILES:	<chem>Cc1ccc(cc1)C[N@H]2CC[C@H](C2)CNC(=O)Nc3cn(c(=O)c4c3cccc4)CC(C)C</chem>
RUN:	RUN1344
DDG (kcal/mol):	-1.20
dDDG (kcal/mol):	0.37

MIC-UNK-d854bf4c-2_1



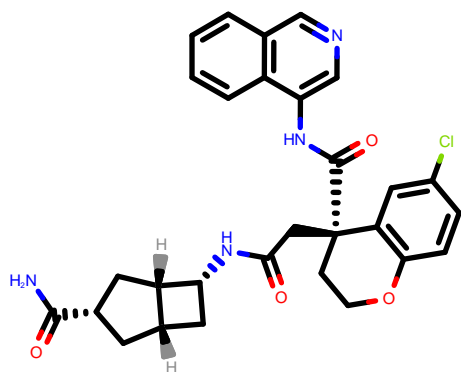
CID:	MIC-UNK-d854bf4c-2_1
SMILES:	<chem>CC(=O)N1CCC2(CC1)CN(C(=O)[C@@H]2c3cccc(c3)Cl)Cl)c4cccc5c4cccc5</chem>
RUN:	RUN3332
DDG (kcal/mol):	-1.20
dDDG (kcal/mol):	0.17

MIC-UNK-91acba05-3_1



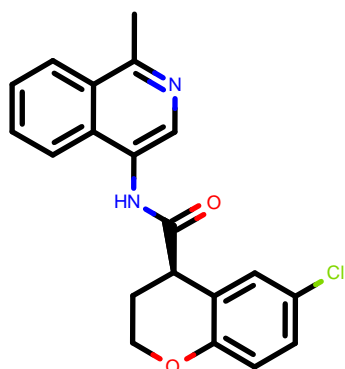
CID:	MIC-UNK-91acba05-3_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCS(=O)(=O)c4c3cc(cc4)Cl</chem>
RUN:	RUN472
DDG (kcal/mol):	-1.19
dDDG (kcal/mol):	0.32

LON-WEI-5e7d1b3e-49_1



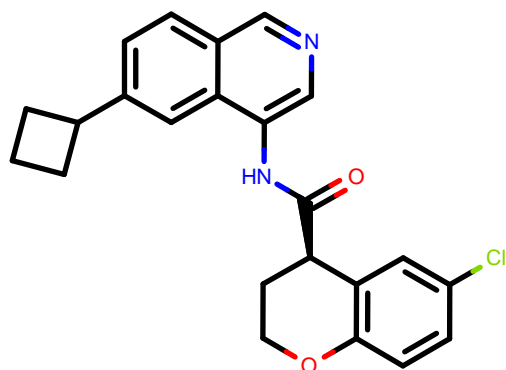
CID:	LON-WEI-5e7d1b3e-49_1
SMILES:	<chem>Cn1cc(c2ccccc2c1=O)NC(=O)NCc3ccccc3</chem>
RUN:	RUN1357
DDG (kcal/mol):	-1.18
dDDG (kcal/mol):	0.28

MAT-POS-b5746674-98_1



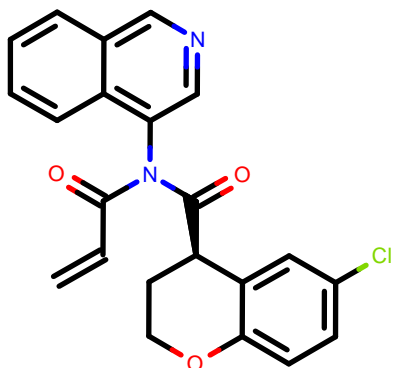
CID:	MAT-POS-b5746674-98_1
SMILES:	<chem>Cn1cc(c2ccccc2c1=O)NC(=O)NCCC[N@@H]3CCc4ccccc4C3</chem>
RUN:	RUN73
DDG (kcal/mol):	-1.18
dDDG (kcal/mol):	0.20

MAK-UNK-c749d764-20_6



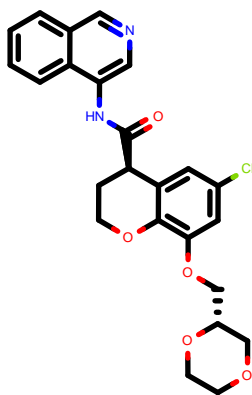
CID:	MAK-UNK-c749d764-20_6
SMILES:	<chem>CCCN(c1nccc2c1cccc2)C(=O)C[C@H]3CCC[C@@H]([C@H]3O)C(F)F</chem>
RUN:	RUN1011
DDG (kcal/mol):	-1.18
dDDG (kcal/mol):	0.21

EDG-MED-ba1ac7b9-27_3



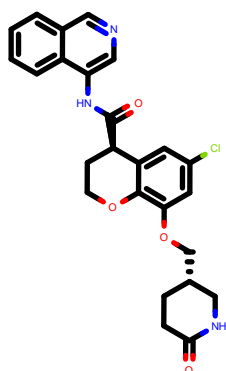
CID:	EDG-MED-ba1ac7b9-27_3
SMILES:	<chem>C/C=C/C(=O)Nc1ccc2c(c1)ncnc2C3=CC=C(Cl)C=C3</chem>
RUN:	RUN2725
DDG (kcal/mol):	-1.18
dDDG (kcal/mol):	0.16

CHO-MSK-a31cca77-2_2



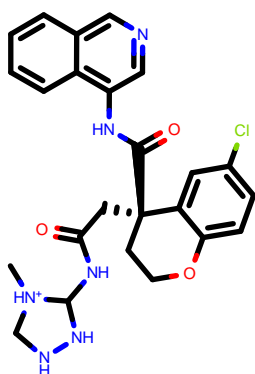
CID:	CHO-MSK-a31cca77-2_2
SMILES:	<chem>Cn1cnc1NC(=O)C[C@]2(CCOC3c2cc(cc3)Cl)C(=O)Nc4cnc5c4cccc5</chem>
RUN:	RUN2195
DDG (kcal/mol):	-1.17
dDDG (kcal/mol):	0.34

DAR-DIA-e7614d05-1_2



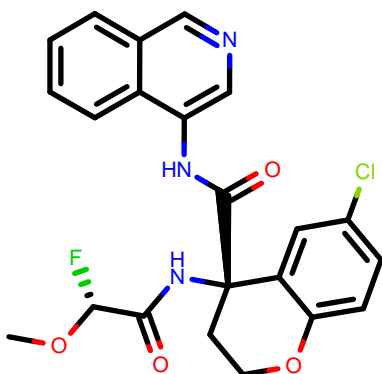
CID:	DAR-DIA-e7614d05-1_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@]3(CCOC4c3cc(cc4)Cl)OC(F)(F)F</chem>
RUN:	RUN2479
DDG (kcal/mol):	-1.17
dDDG (kcal/mol):	0.37

RAL-THA-e002e396-11_2



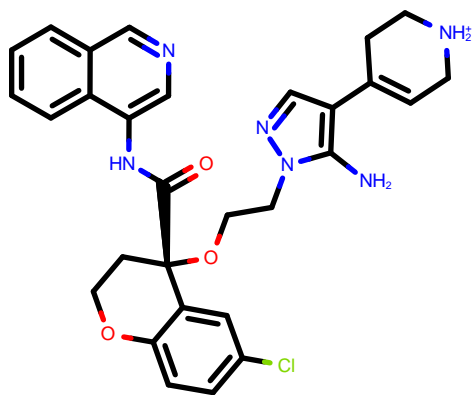
CID:	RAL-THA-e002e396-11_2
SMILES:	<chem>COC[C@]1(CCOC2c1cc(cc2)Cl)C(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN3470
DDG (kcal/mol):	-1.16
dDDG (kcal/mol):	0.12

EDG-MED-70ae9412-1_2



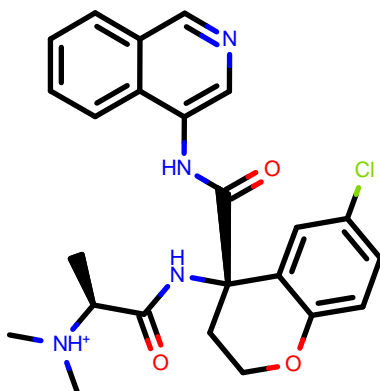
CID:	EDG-MED-70ae9412-1_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@]3(CCOC4c3cc(cc4)Cl)CNC(=O)Cn5cnc5</chem>
RUN:	RUN3164
DDG (kcal/mol):	-1.16
dDDG (kcal/mol):	0.11

MAT-POS-8a69d52e-6_1



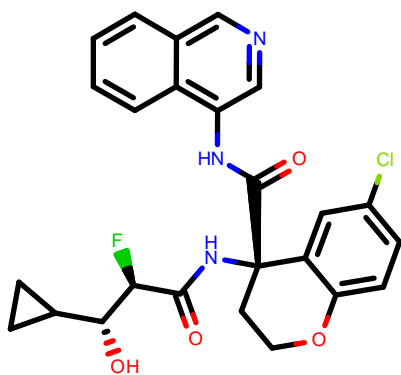
CID:	MAT-POS-8a69d52e-6_1
SMILES:	<chem>C[C@@H]1COc2ccc(cc2[C@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN374
DDG (kcal/mol):	-1.15
dDDG (kcal/mol):	0.38

MAK-UNK-8be7dca9-4_2



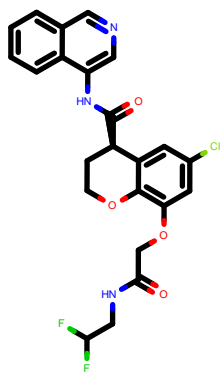
CID:	MAK-UNK-8be7dca9-4_2
SMILES:	<chem>c1cc2cncc(c2cc1N3CC[NH2+][CC3])NC(=O)[C@H]4CCOc5c4cc(cc5)Cl</chem>
RUN:	RUN505
DDG (kcal/mol):	-1.15
dDDG (kcal/mol):	0.33

DAR-DIA-0f2f46c9-9_2



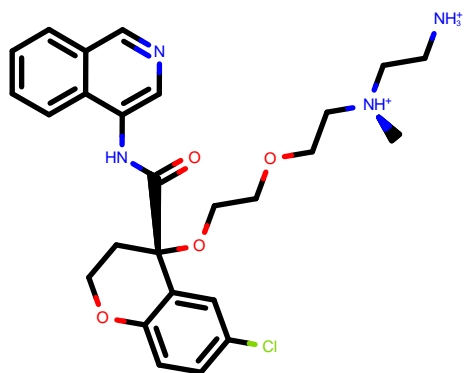
CID:	DAR-DIA-0f2f46c9-9_2
SMILES:	<chem>CNS(=O)(=O)[N@]1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3243
DDG (kcal/mol):	-1.15
dDDG (kcal/mol):	0.15

ALP-UNI-3496895b-5_3



CID:	ALP-UNI-3496895b-5_3
SMILES:	<chem>C[C@@H]1C[N@]1CC[C@@H]1NC(=O)C[C@]2(COC3Ck2cc(c3)Cl)C(=O)Nc4nc5c4cccc5)S(=O)(=O)C</chem>
RUN:	RUN2514
DDG (kcal/mol):	-1.15
dDDG (kcal/mol):	0.39

DAR-DIA-23e5a6a0-7_1



CID: DAR-DIA-23e5a6a0-7_1

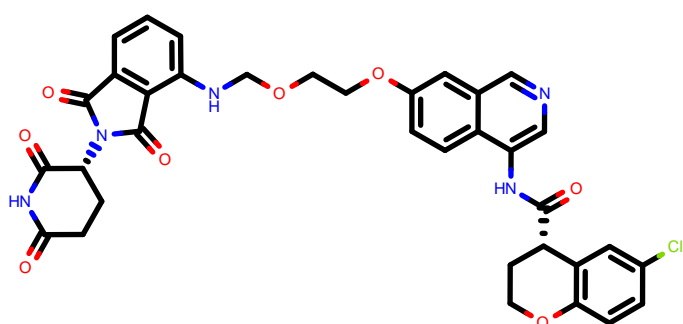
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H](C)COCc3ccccc3c4c(c1)C@H](C)COCc5[nH]2+5)CCc6)Cl

RUN: RUN413

DDG (kcal/mol): -1.15

dDDG (kcal/mol): 0.37

DAR-DIA-5d6f1b43-10_1



CID: DAR-DIA-5d6f1b43-10_1

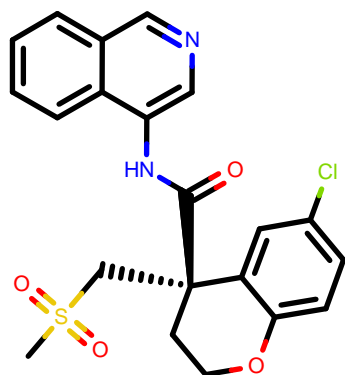
SMILES: c1ccc2c(c1)cnc2N3CCCC4(CCCCC4)N(C3=O)c5ccccc5)Cl

RUN: RUN487

DDG (kcal/mol): -1.15

dDDG (kcal/mol): 0.32

PET-UNK-12d8d43f-2_1



CID: PET-UNK-12d8d43f-2_1

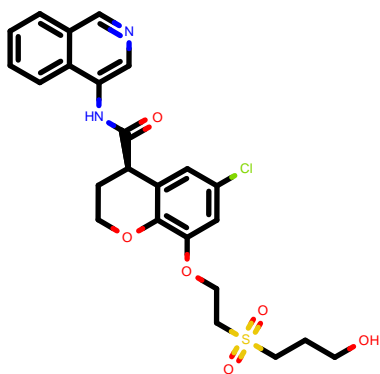
SMILES: COC(=O)N(c1cnc2c1cccc2)C(=O)Cc3ccccc3)Cl

RUN: RUN1496

DDG (kcal/mol): -1.15

dDDG (kcal/mol): 0.25

JOH-UNI-ea72002d-3_4



CID: JOH-UNI-ea72002d-3_4

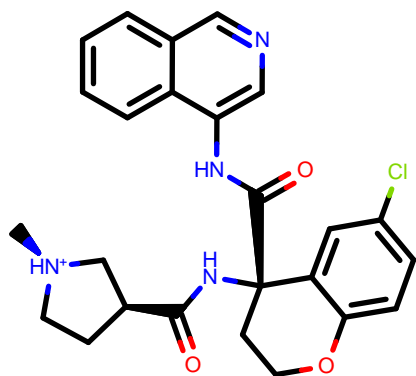
SMILES: c1ccc2c(c1)cnc2[C@@H](C(=O)[C@@H](C)COCc3ccccc3)N5C(=O)C=CC5=O

RUN: RUN2487

DDG (kcal/mol): -1.14

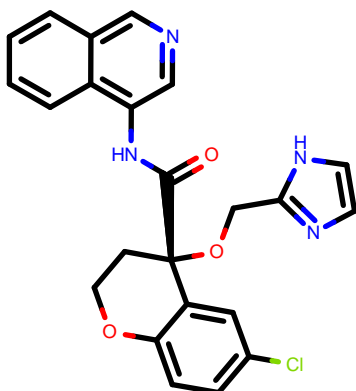
dDDG (kcal/mol): 0.57

MIC-UNK-cdc2493e-19_1



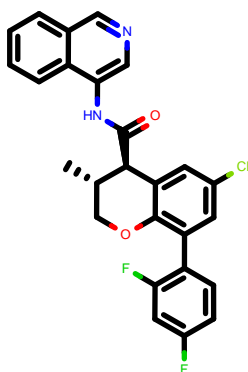
CID:	MIC-UNK-cdc2493e-19_1
SMILES:	<chem>CC(C)CCCN(c1cccc(c1)Cl)C(=O)Nc2cncc3c2cccc3</chem>
RUN:	RUN567
DDG (kcal/mol):	-1.14
dDDG (kcal/mol):	0.40

PET-UNK-689df078-4_2



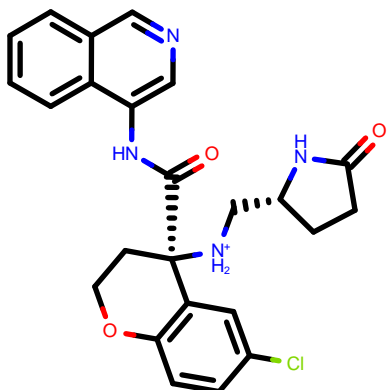
CID:	PET-UNK-689df078-4_2
SMILES:	<chem>C[N@H]1CC(=O)N(C(=O)[C@@H]1c2cccc(c2)Cl)c3cncc4c3cccc4</chem>
RUN:	RUN3106
DDG (kcal/mol):	-1.14
dDDG (kcal/mol):	0.12

MIC-UNK-c66144cb-3_2



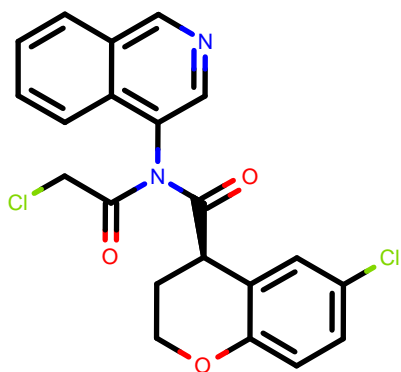
CID:	MIC-UNK-c66144cb-3_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](CCCCCCCC3)c4cccc(c4)Cl</chem>
RUN:	RUN134
DDG (kcal/mol):	-1.14
dDDG (kcal/mol):	0.28

EDG-MED-90036822-29_1



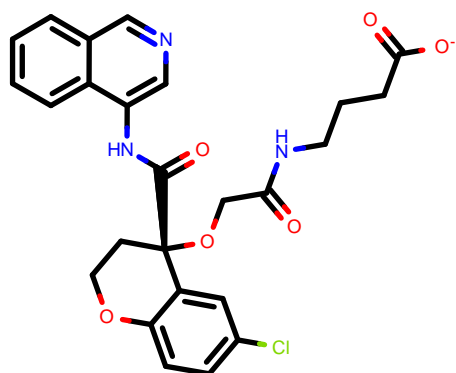
CID:	EDG-MED-90036822-29_1
SMILES:	<chem>C[C@H](C(=O)N[C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4)O</chem>
RUN:	RUN1698
DDG (kcal/mol):	-1.12
dDDG (kcal/mol):	0.28

EDG-MED-ba1ac7b9-23_3



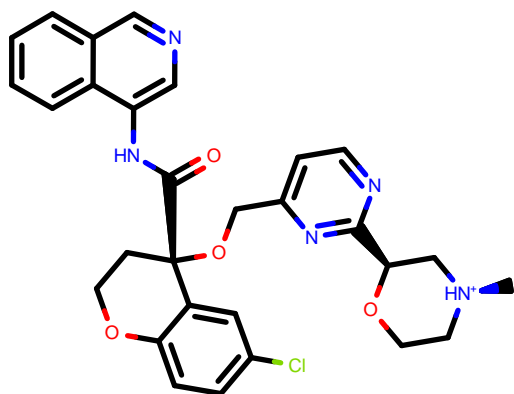
CID:	EDG-MED-ba1ac7b9-23_3
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)C(Cl)C(=O)Nc3ccc4c3ccoc4</chem>
RUN:	RUN2709
DDG (kcal/mol):	-1.11
dDDG (kcal/mol):	0.15

MIC-UNK-cdc2493e-4_2



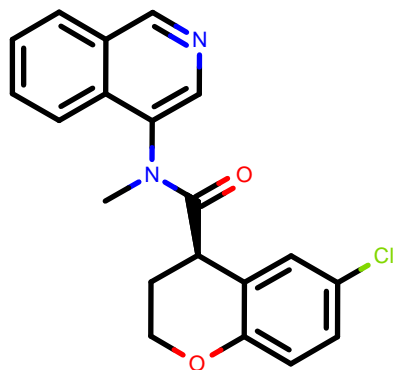
CID:	MIC-UNK-cdc2493e-4_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)N(c3cccc(c3)Cl)C4C[C@H]5CCCC[C@H]5C4</chem>
RUN:	RUN528
DDG (kcal/mol):	-1.10
dDDG (kcal/mol):	0.34

PET-UNK-dd44aeb6-1_1



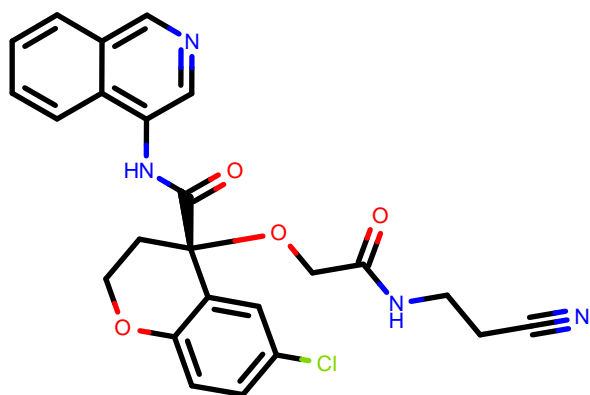
CID:	PET-UNK-dd44aeb6-1_1
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)Cc3cc(ccc3F)Cl</chem>
RUN:	RUN426
DDG (kcal/mol):	-1.10
dDDG (kcal/mol):	0.36

ALF-EVA-ced740bd-1_2



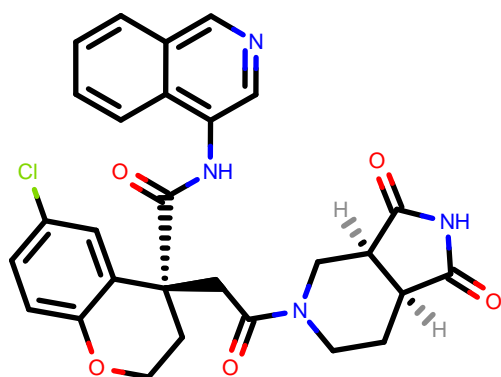
CID:	ALF-EVA-ced740bd-1_2
SMILES:	<chem>COc1c(ccc2c1c(nc2)NC(=O)[C@H]3CCOC4c3cc(cc4)Cl)Cl</chem>
RUN:	RUN2789
DDG (kcal/mol):	-1.10
dDDG (kcal/mol):	0.09

PET-UNK-bb7ffe78-2_1



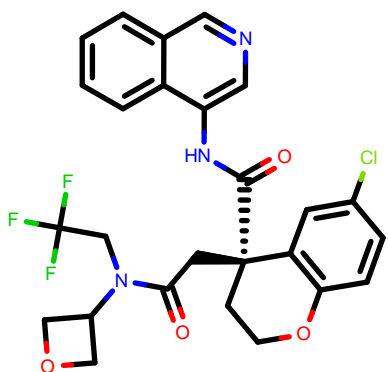
CID:	PET-UNK-bb7ffe78-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)OC(F)(F)F</chem>
RUN:	RUN3331
DDG (kcal/mol):	-1.10
dDDG (kcal/mol):	0.16

VLA-UNK-5b0345c3-1_1



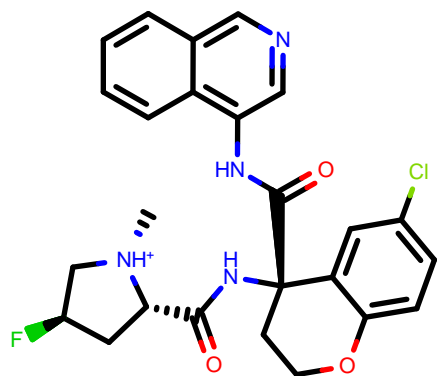
CID:	VLA-UNK-5b0345c3-1_1
SMILES:	<chem>COc1cccc2c1cc([nH]2)C(=O)N3Cc4ccc(cc4[C@@H](C3)C(=O)Nc5ccc6c5ccc6)Cl</chem>
RUN:	RUN3883
DDG (kcal/mol):	-1.09
dDDG (kcal/mol):	0.16

RAL-THA-8416115c-5_3



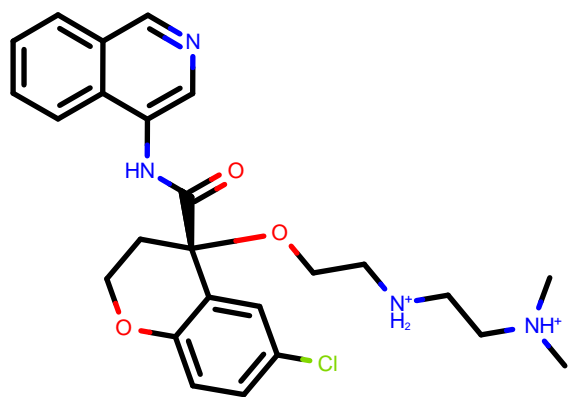
CID:	RAL-THA-8416115c-5_3
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](C3CCN(c4c3cc(cc4)Cl)Cc5[nH]cn5</chem>
RUN:	RUN1265
DDG (kcal/mol):	-1.08
dDDG (kcal/mol):	0.21

DAR-DIA-0f2f46c9-7_1



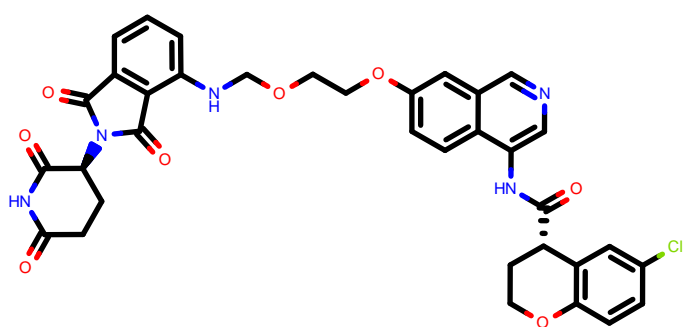
CID:	DAR-DIA-0f2f46c9-7_1
SMILES:	<chem>CNS(=O)(=O)[N@@]1CC[C@@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3ccc4</chem>
RUN:	RUN3234
DDG (kcal/mol):	-1.08
dDDG (kcal/mol):	0.23

MAT-POS-8a69d52e-4_2



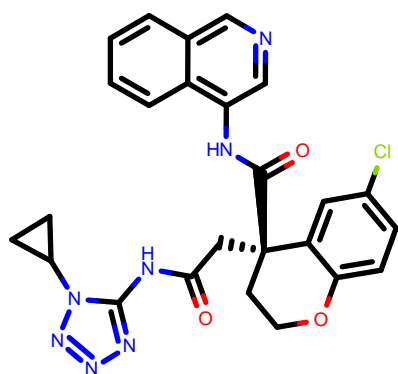
CID:	MAT-POS-8a69d52e-4_2
SMILES:	<chem>C[C@H]1COc2ccc(cc2[C@@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN367
DDG (kcal/mol):	-1.08
dDDG (kcal/mol):	0.46

EDJ-MED-e4b030d8-4_1



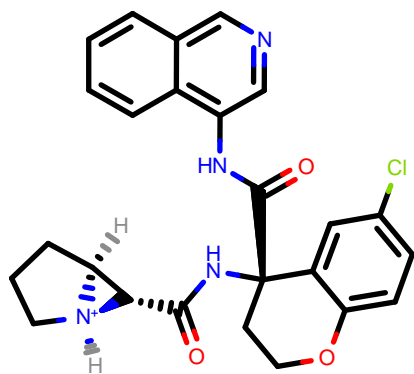
CID:	EDJ-MED-e4b030d8-4_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOc4c3cc(cc4OC5COC5)Cl</chem>
RUN:	RUN287
DDG (kcal/mol):	-1.08
dDDG (kcal/mol):	0.39

KAD-UNI-cb0f2bbc-21_1



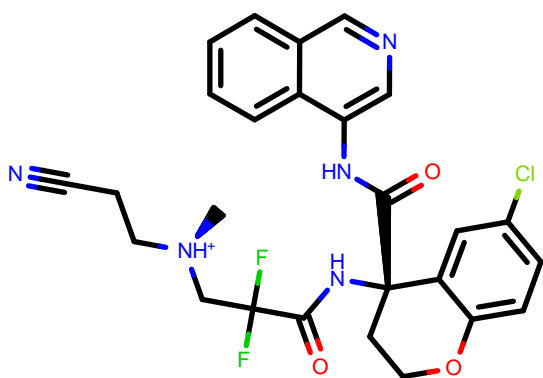
CID:	KAD-UNI-cb0f2bbc-21_1
SMILES:	<chem>Cn1cc(n2e1c(n2)C#N)C(NH2)C[C@@]3(CCOc4c3cc(cc4)O)C(=O)Nc5cncc6c5cccc6</chem>
RUN:	RUN3707
DDG (kcal/mol):	-1.08
dDDG (kcal/mol):	0.17

MAT-POS-78e1d523-1_2



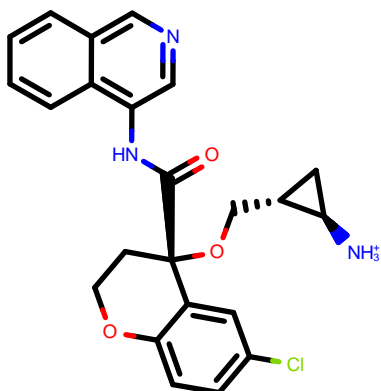
CID:	MAT-POS-78e1d523-1_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCSc4c3cc(cc4)Cl</chem>
RUN:	RUN3279
DDG (kcal/mol):	-1.07
dDDG (kcal/mol):	0.16

MIC-UNK-5a93dd5f-2_5



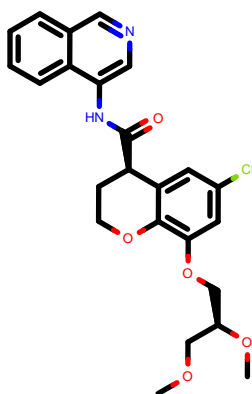
CID:	MIC-UNK-5a93dd5f-2_5
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@@H](c3cccc(c3)Cl)N[+]C(C)(F)F5CCCCC(C)H5C4</chem>
RUN:	RUN738
DDG (kcal/mol):	-1.07
dDDG (kcal/mol):	0.33

VLA-UNK-f49ebb87-1_2



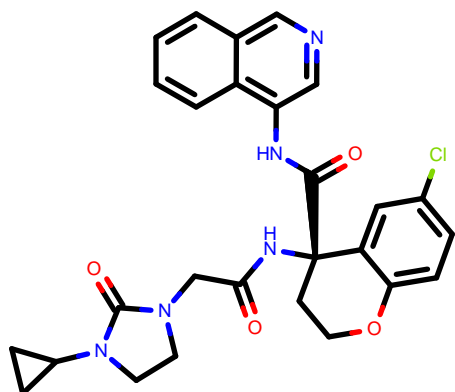
CID:	VLA-UNK-f49ebb87-1_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H](c3ccccc3)NCC(=O)O3Cl</chem>
RUN:	RUN3102
DDG (kcal/mol):	-1.06
dDDG (kcal/mol):	0.16

ALF-EVA-5b152d2f-8_1



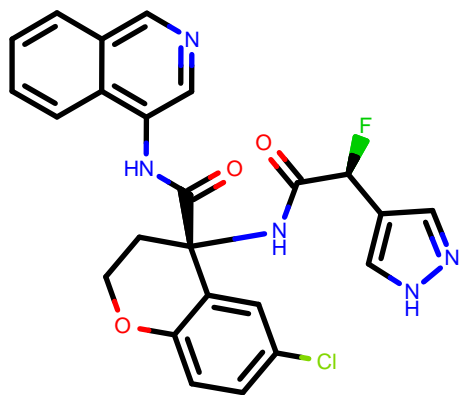
CID:	ALF-EVA-5b152d2f-8_1
SMILES:	<chem>c1cc2c(cc1CC3CC3)ncnc2NC(=O)[C@@H](c4ccccc4)C(=O)O3Cl</chem>
RUN:	RUN2362
DDG (kcal/mol):	-1.06
dDDG (kcal/mol):	0.48

RAL-THA-8416115c-6_2



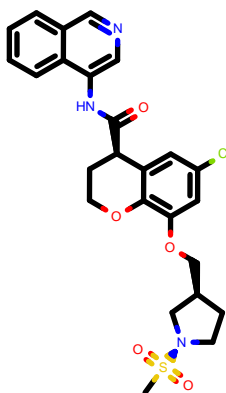
CID:	RAL-THA-8416115c-6_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@H](c3ccccc3)NCC(=O)O3Cl</chem>
RUN:	RUN1268
DDG (kcal/mol):	-1.06
dDDG (kcal/mol):	0.34

ALP-POS-fe871b40-11_1



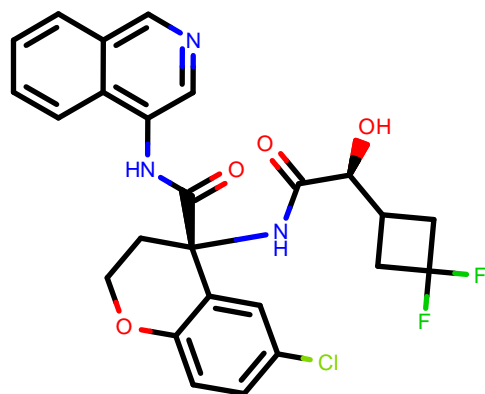
CID:	ALP-POS-fe871b40-11_1
SMILES:	<chem>CO[C@@]1(CCOc2c1cc(c(c2)F)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3129
DDG (kcal/mol):	-1.06
dDDG (kcal/mol):	0.15

KAD-UNI-80f122c8-3_1



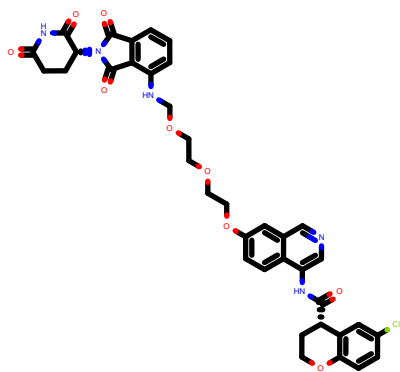
CID:	KAD-UNI-80f122c8-3_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(c(c4)Cl)CC(=O)Nc5cnc6c5cccc6</chem>
RUN:	RUN2288
DDG (kcal/mol):	-1.06
dDDG (kcal/mol):	0.46

VLA-UNK-70dd90ef-2_1



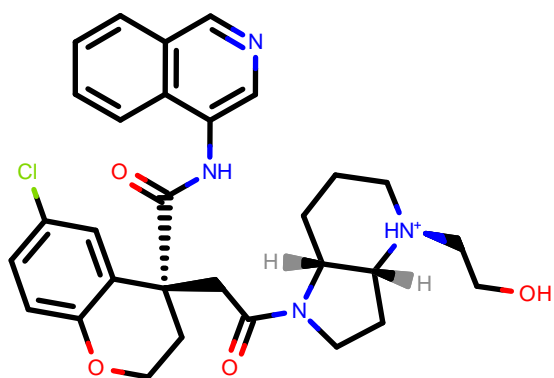
CID:	VLA-UNK-70dd90ef-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3C(=O)[C@@]4(C5CC(CCC5NC4=O)Cl)NC3=O</chem>
RUN:	RUN3303
DDG (kcal/mol):	-1.06
dDDG (kcal/mol):	0.16

FRA-DIA-0fa076fe-1_1



CID:	FRA-DIA-0fa076fe-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3COc4c3cccc4Cl</chem>
RUN:	RUN417
DDG (kcal/mol):	-1.06
dDDG (kcal/mol):	0.48

MAT-POS-4223bc15-24_2



CID: MAT-POS-4223bc15-24_2

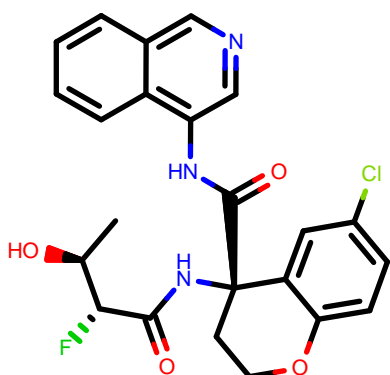
SMILES: CN(C)C(=O)C[N@H+]1Cc2ccc(cc2[C@@H](C1)C(=O)Nc3cncc4c3cccc4)Cl

RUN: RUN4110

DDG (kcal/mol): -1.05

dDDG (kcal/mol): 0.27

NIR-THE-3fc2bec4-1_1



CID: NIR-THE-3fc2bec4-1_1

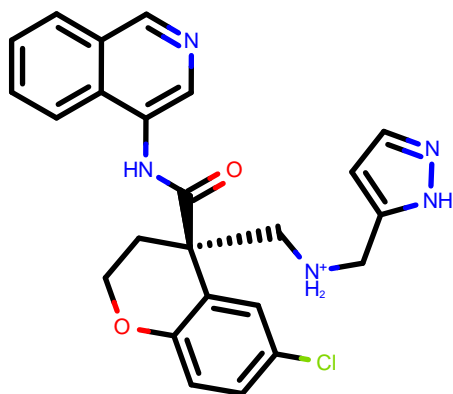
SMILES: CO[C@]1(CCNc2c1cc(c(c2)F)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN3275

DDG (kcal/mol): -1.05

dDDG (kcal/mol): 0.11

EDG-MED-90036822-91_1



CID: EDG-MED-90036822-91_1

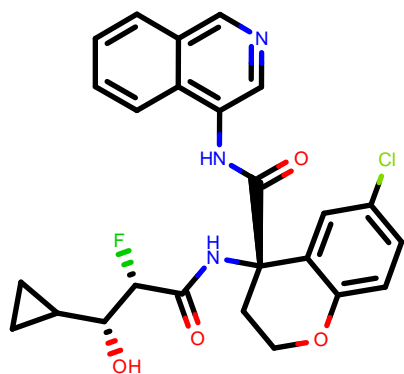
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)NC(=O)[C@@]@H(c5cccc(c5)F)[NH3+]

RUN: RUN1799

DDG (kcal/mol): -1.05

dDDG (kcal/mol): 0.41

DAR-DIA-0f2f46c9-10_3



CID: DAR-DIA-0f2f46c9-10_3

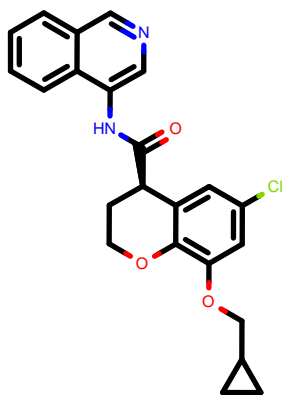
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]@H3CC[N@](c4c3cc(cc4)Cl)S(=O)(=O)[O-]

RUN: RUN3244

DDG (kcal/mol): -1.04

dDDG (kcal/mol): 0.17

DAR-DIA-9e4459de-15_10



CID: DAR-DIA-9e4459de-15_10

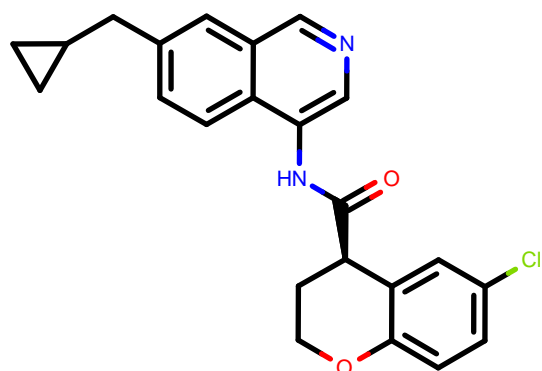
SMILES: c1cc2c(c1)NC(=O)CCCl(=O)Nc3ccc4c(c3)ncoc4NC(=O)C@H5CCOC6c5cc6c7C|c1c(c2O)[C@H]7CCCl(=O)NC7=O|O

RUN: RUN1450

DDG (kcal/mol): -1.03

dDDG (kcal/mol): 0.32

DAR-DIA-0d514e7d-32_9



CID: DAR-DIA-0d514e7d-32_9

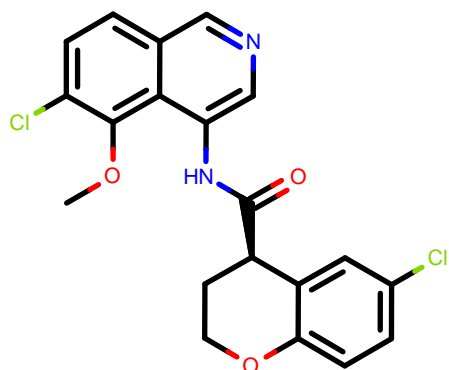
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3[C@@H]4C[C@@H]4CO[C@H]5[C@@H]3C=C(C=C5)Cl

RUN: RUN858

DDG (kcal/mol): -1.03

dDDG (kcal/mol): 0.23

LON-WEI-5e7d1b3e-34_1



CID: LON-WEI-5e7d1b3e-34_1

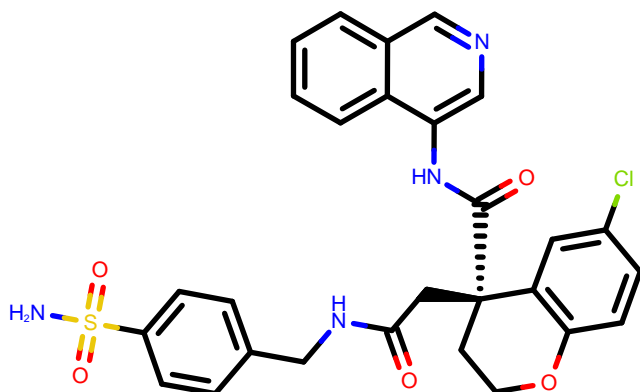
SMILES: CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)N3CC[C@@H](C3)c4ccccc4

RUN: RUN1335

DDG (kcal/mol): -1.03

dDDG (kcal/mol): 0.20

NAU-LAT-4ce8bf23-1_1



CID: NAU-LAT-4ce8bf23-1_1

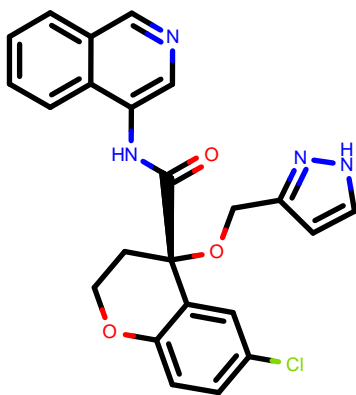
SMILES: CC(=O)NC[C@@H](c1cncc2c1cccc2)C(=O)Nc3cccc(c3)Cl

RUN: RUN1389

DDG (kcal/mol): -1.03

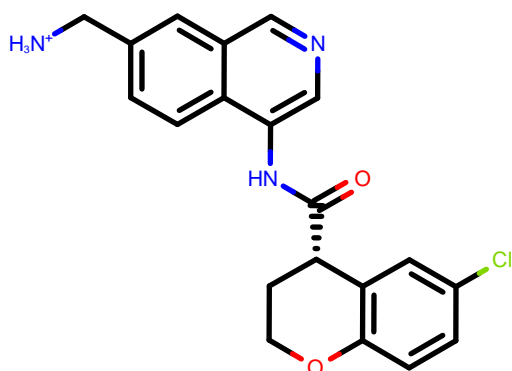
dDDG (kcal/mol): 0.25

MAT-POS-3b97339c-2_1



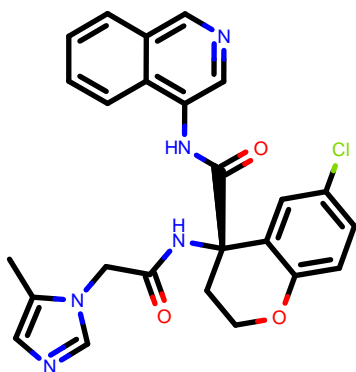
CID:	MAT-POS-3b97339c-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)[NH3+]</chem>
RUN:	RUN3296
DDG (kcal/mol):	-1.03
dDDG (kcal/mol):	0.11

ERI-UCB-ce40166b-12_1



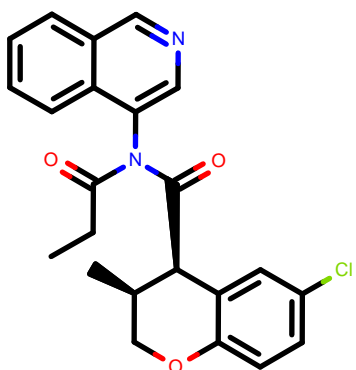
CID:	ERI-UCB-ce40166b-12_1
SMILES:	<chem>c1ccc2c(c1)cncc2CC(=O)Nc3cc(cc(c3)O)c4cccnc4C#N</chem>
RUN:	RUN51
DDG (kcal/mol):	-1.03
dDDG (kcal/mol):	0.16

VLA-UNK-3a43cd95-4_1



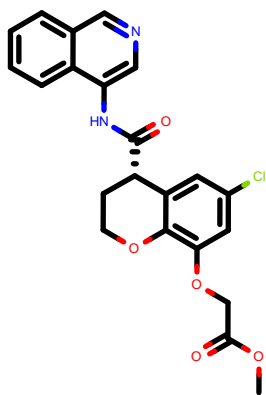
CID:	VLA-UNK-3a43cd95-4_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3C(=O)C[C@@]4(C3=O)COc5c4cc(c(c5)F)Cl</chem>
RUN:	RUN3191
DDG (kcal/mol):	-1.02
dDDG (kcal/mol):	0.14

LON-WEI-4d77710c-27_1



CID:	LON-WEI-4d77710c-27_1
SMILES:	<chem>CC(C)Cn1cc(c2cccc2c1=O)NC(=O)NC3CCCCC3</chem>
RUN:	RUN211
DDG (kcal/mol):	-1.02
dDDG (kcal/mol):	0.22

MAT-POS-e9e99895-13_7



CID: MAT-POS-e9e99895-13_7

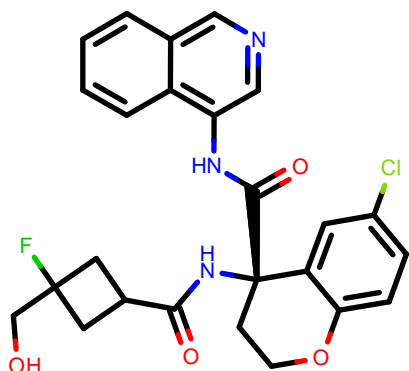
SMILES: C[C@@H](c1ccc(c(c1)Cl)Cl)(C=O)Nc2nccc3c2ccc3NC(=O)[C@H]4CC[N@H+](C4)C

RUN: RUN2273

DDG (kcal/mol): -1.02

dDDG (kcal/mol): 0.23

LON-WEI-9739a092-3_1



CID: LON-WEI-9739a092-3_1

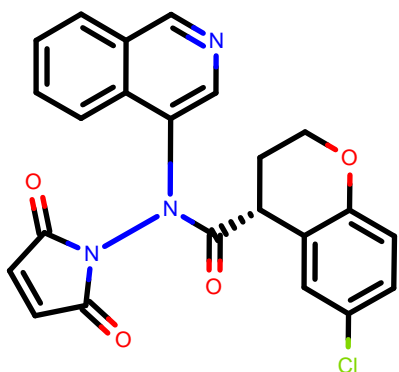
SMILES: COc1cc(ccc1Nc2cc(cc(c2)Cl)Cl)CC(=O)Nc3cncc4c3cccc4)Br

RUN: RUN3267

DDG (kcal/mol): -1.01

dDDG (kcal/mol): 0.14

MAK-UNK-c749d764-10_3



CID: MAK-UNK-c749d764-10_3

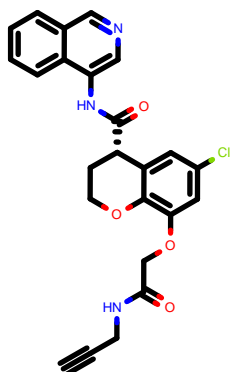
SMILES: CSN(c1cncc2c1cccc2)C(=O)[C@H](C@@H)3CCC[C@H]([C@@H]3O)C(F)F

RUN: RUN943

DDG (kcal/mol): -1.01

dDDG (kcal/mol): 0.30

JOH-UNI-ea72002d-4_2



CID: JOH-UNI-ea72002d-4_2

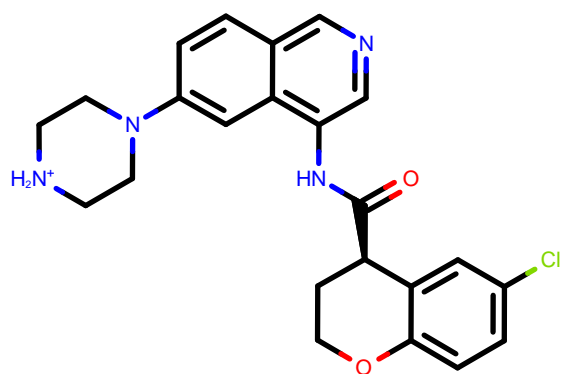
SMILES: c1ccc2c(c1)cncc2[C@H](C(=O)[C@H]3CCOCc4c3cc(cc4)Cl)(N5C(=O)C=CC5=O)F

RUN: RUN2490

DDG (kcal/mol): -1.00

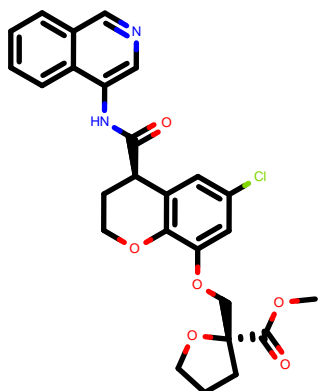
dDDG (kcal/mol): 0.29

EDG-MED-ba1ac7b9-30_2



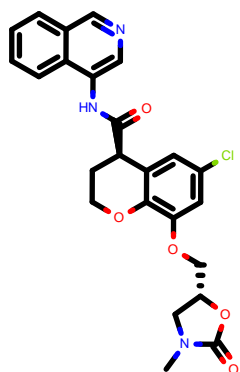
CID:	EDG-MED-ba1ac7b9-30_2
SMILES:	<chem>C[NH+]1CCC(CC1)N(CCO)C(=O)C[C@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4ccc5c4cccc5</chem>
RUN:	RUN2741
DDG (kcal/mol):	-1.00
dDDG (kcal/mol):	0.10

MAT-POS-2905de8c-2_2



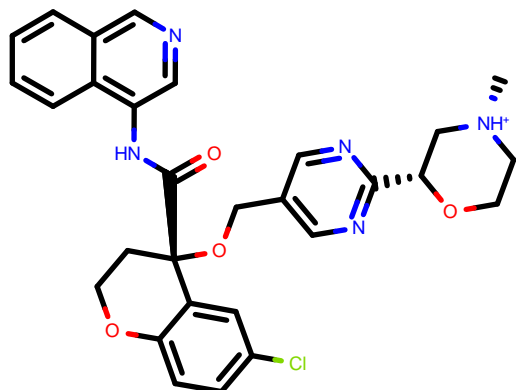
CID:	MAT-POS-2905de8c-2_2
SMILES:	<chem>C[NH2+][C@]1(CCOC2c1cc(cc2)Cl)C(=O)Nc3ccc4c3cccc4</chem>
RUN:	RUN2230
DDG (kcal/mol):	-1.00
dDDG (kcal/mol):	0.48

ALP-UNI-0676e700-5_1



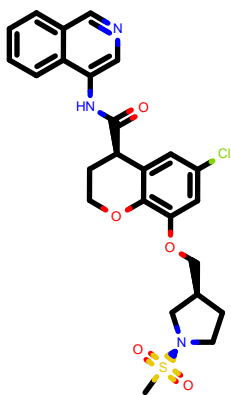
CID:	ALP-UNI-0676e700-5_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CNC(=O)c5cccc(c5)OCC(=O)N</chem>
RUN:	RUN2449
DDG (kcal/mol):	-1.00
dDDG (kcal/mol):	0.37

JAG-UCB-f37eaa14-9_2



CID:	JAG-UCB-f37eaa14-9_2
SMILES:	<chem>c1ccc2c(c1)cncc2N3CC[C@@]4(C3=O)C[N@](c5c4cc(cc5)Cl)C6CC(=O)O6</chem>
RUN:	RUN3071
DDG (kcal/mol):	-1.00
dDDG (kcal/mol):	0.21

LEE-CAM-7ab9b158-1_3



CID: LEE-CAM-7ab9b158-1_3

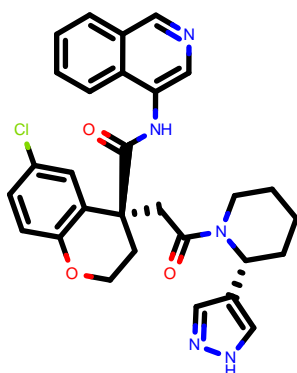
SMILES: C[NH+]([C]([C]([H]1COC[C@H]1OC[C@@]2(CCOc3c2cc(cc3)C)C(=O)Nc4ccc5c4cccc5

RUN: RUN2202

DDG (kcal/mol): -1.00

dDDG (kcal/mol): 0.44

RAL-THA-8416115c-2_3



CID: RAL-THA-8416115c-2_3

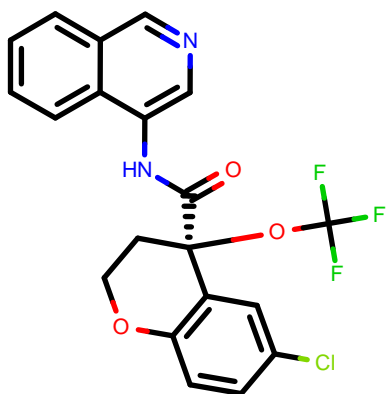
SMILES: CN1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN1253

DDG (kcal/mol): -1.00

dDDG (kcal/mol): 0.25

MAK-UNK-c749d764-15_1



CID: MAK-UNK-c749d764-15_1

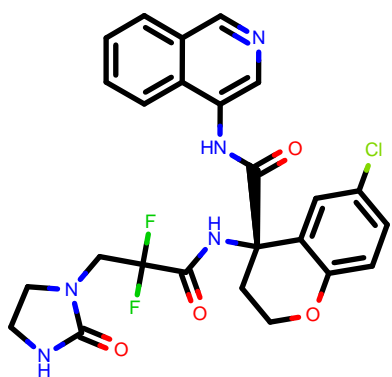
SMILES: C[C@@H](Nc1cnc2c1ccc2Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN959

DDG (kcal/mol): -0.99

dDDG (kcal/mol): 0.08

MIC-UNK-cdc2493e-14_2



CID: MIC-UNK-cdc2493e-14_2

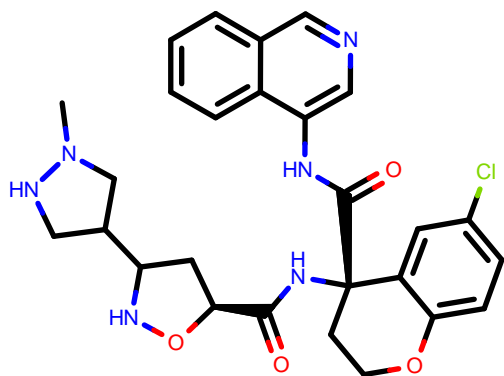
SMILES: c1ccc2c(c1)cnc2N(C(=O)N(c3cccc(c3)Cl)C@H)C[C@@H](C@H)(C4)NH5CCCC5

RUN: RUN556

DDG (kcal/mol): -0.99

dDDG (kcal/mol): 0.33

EDJ-MED-009f762b-4_2



CID: EDJ-MED-009f762b-4_2

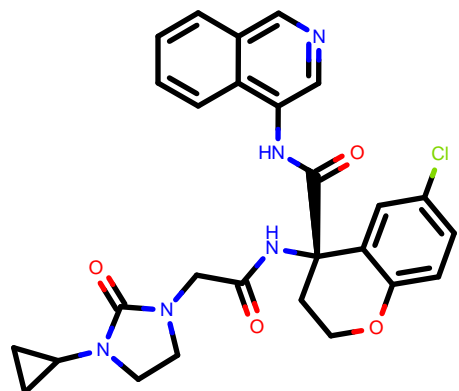
SMILES: Cc1[nH]c(c[nH+]1)C[N@H]2Cc3ccc(cc3[C@@H](C2)C(=O)Nc4ccc5c4cc(cc5)F)Cl

RUN: RUN3913

DDG (kcal/mol): -0.99

dDDG (kcal/mol): 0.14

ALP-POS-a0a4abd7-1_2



CID: ALP-POS-a0a4abd7-1_2

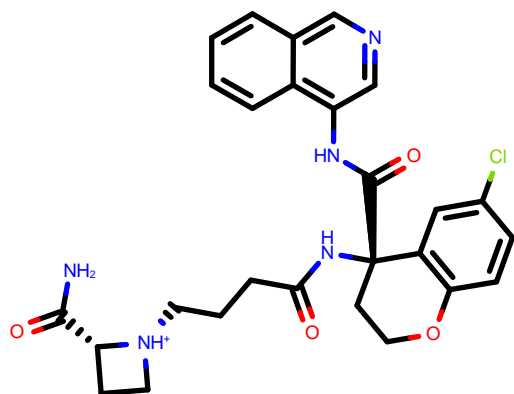
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@]3(CCOc4c3ccc(c4)Cl)Cn5c[nH]c(=O)[nH]5=O

RUN: RUN3562

DDG (kcal/mol): -0.99

dDDG (kcal/mol): 0.18

ADA-UCB-dc2b944c-16_1



CID: ADA-UCB-dc2b944c-16_1

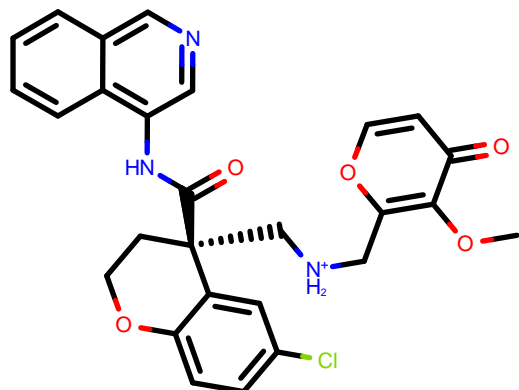
SMILES: CC1(C[C@H](c2cc(ccc2O1)Cl)C(=O)Nc3cncc4c3cccc4)C

RUN: RUN614

DDG (kcal/mol): -0.98

dDDG (kcal/mol): 0.33

ALP-POS-a577c8a2-3_1



CID: ALP-POS-a577c8a2-3_1

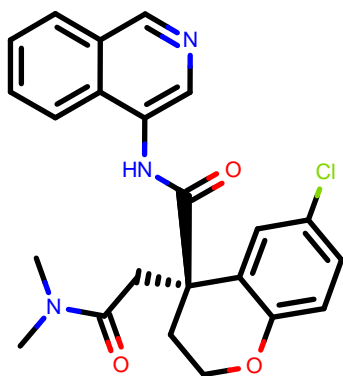
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3C[N@@]4(S(=O)(=O)c4c3cc(cc4)Cl)C[C5[nH]ccn5

RUN: RUN4628

DDG (kcal/mol): -0.98

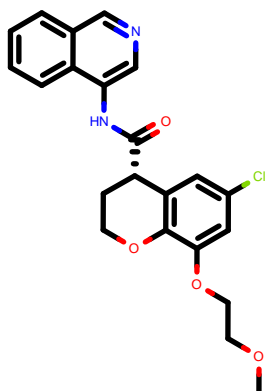
dDDG (kcal/mol): 0.27

MAK-UNK-c749d764-18_6



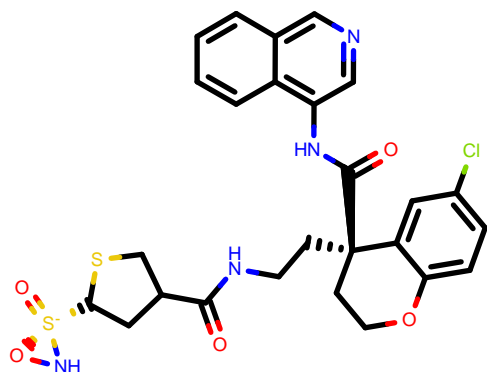
CID:	MAK-UNK-c749d764-18_6
SMILES:	<chem>CC(C)SCN(c1cncc2c1cccc2)C(=O)C[C@H]3CCC[C@@H]([C@H]3O)C(F)F</chem>
RUN:	RUN995
DDG (kcal/mol):	-0.97
dDDG (kcal/mol):	0.31

EDG-MED-90036822-24_1



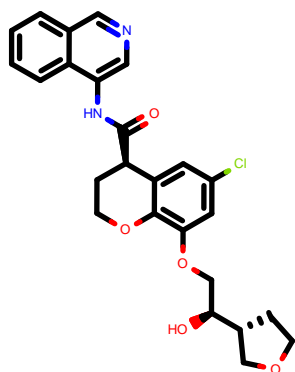
CID:	EDG-MED-90036822-24_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)NC(=O)C5(COC5)F</chem>
RUN:	RUN1688
DDG (kcal/mol):	-0.97
dDDG (kcal/mol):	0.22

DAR-DIA-0587064e-27_2



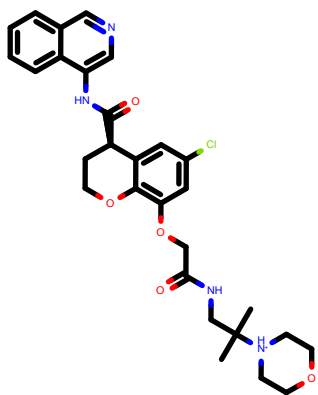
CID:	DAR-DIA-0587064e-27_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)OCc5ccc(cc5Cl)F)F)Cl</chem>
RUN:	RUN3388
DDG (kcal/mol):	-0.96
dDDG (kcal/mol):	0.21

EDJ-MED-d203f206-35_1



CID:	EDJ-MED-d203f206-35_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)O)CC(O)N6CC[C@H]([C@@H]3C5(C)O)N6=O</chem>
RUN:	RUN2596
DDG (kcal/mol):	-0.95
dDDG (kcal/mol):	0.40

ALP-POS-5bb456a5-9_2



CID: ALP-POS-5bb456a5-9_2

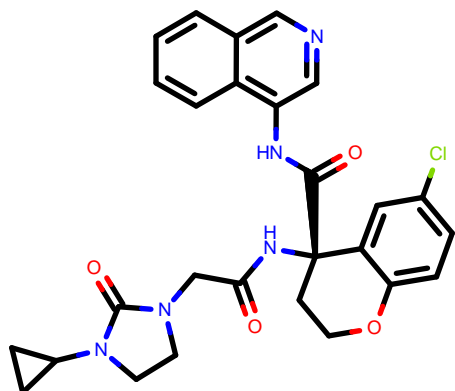
SMILES: CCNC(=O)C[NH+]1CCN(CC1)C(=O)C[C@]2(CCOc3c2cc(cc3)C)C(=O)Nc4ncc5c4cccc5

RUN: RUN2445

DDG (kcal/mol): -0.95

dDDG (kcal/mol): 0.40

MAT-POS-4223bc15-9_10



CID: MAT-POS-4223bc15-9_10

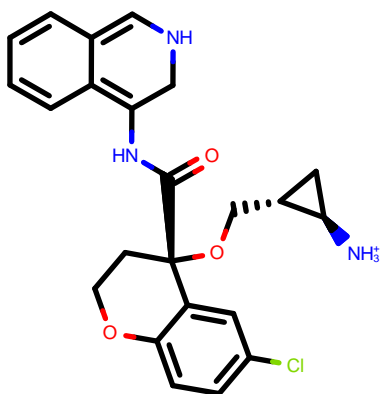
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3C[N@@]4)Cc4c3cc(cc4)C(S(=O)(=O)C)C#N

RUN: RUN4018

DDG (kcal/mol): -0.95

dDDG (kcal/mol): 0.17

VLA-UNK-9a7dc93f-4_1



CID: VLA-UNK-9a7dc93f-4_1

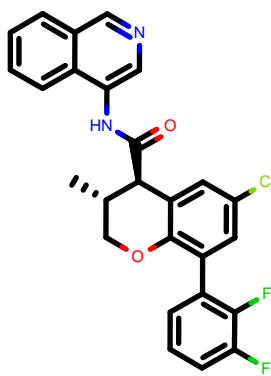
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCOCc4c3cc(c(c4F)F)C#N

RUN: RUN3086

DDG (kcal/mol): -0.95

dDDG (kcal/mol): 0.16

DAR-DIA-53551c05-11_1



CID: DAR-DIA-53551c05-11_1

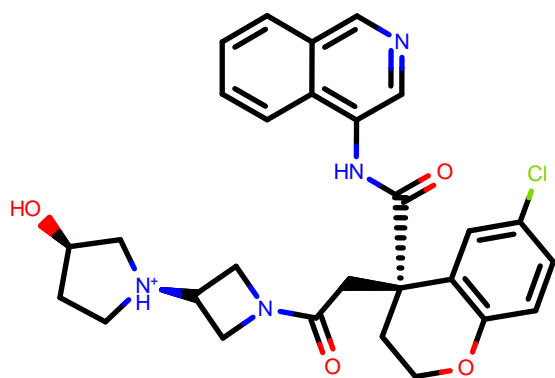
SMILES: c1ccc2c(c1)cnc2NC3c4cc(ccc4C(=O)C3=O)c5cc(cc(c5)C)O[C@@H]6CC(=O)N6

RUN: RUN144

DDG (kcal/mol): -0.95

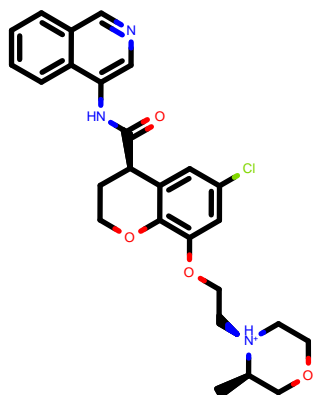
dDDG (kcal/mol): 0.29

ERI-UCB-d6de1f3c-7_2



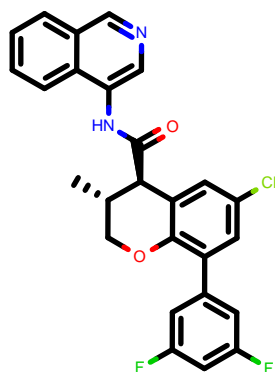
CID:	ERI-UCB-d6de1f3c-7_2
SMILES:	<chem>C[C@H]1CN(CC(=O)N1c2cccc(c2)Cl)C(=O)c3cnc4c3cccc4</chem>
RUN:	RUN1100
DDG (kcal/mol):	-0.94
dDDG (kcal/mol):	0.25

EDJ-MED-fcba3f31-8_1



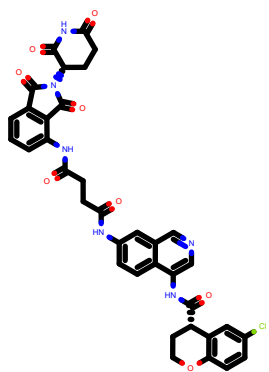
CID:	EDJ-MED-fcba3f31-8_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(C)CCOc4c3cc(cc4)Cl)CO[C@H]5CCOC5=O</chem>
RUN:	RUN2545
DDG (kcal/mol):	-0.94
dDDG (kcal/mol):	0.50

MIC-UNK-c66144cb-1_2



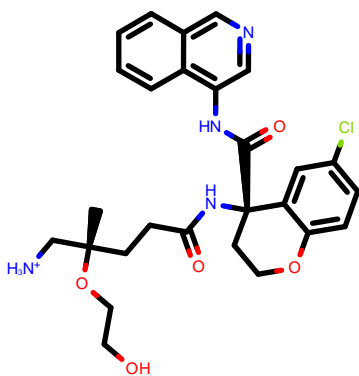
CID:	MIC-UNK-c66144cb-1_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@H](CCc3cccc(c3)F)c4cccc(c4)Cl</chem>
RUN:	RUN129
DDG (kcal/mol):	-0.94
dDDG (kcal/mol):	0.29

VLA-UCB-05e51b3f-14_1



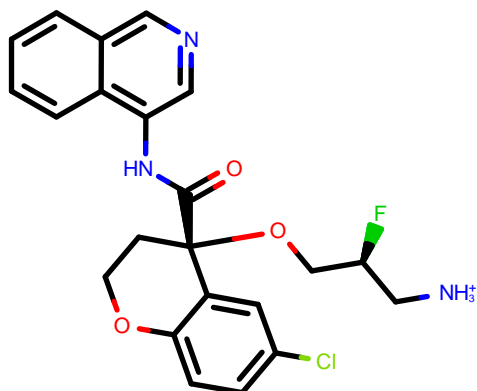
CID:	VLA-UCB-05e51b3f-14_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)N(CCC3CCCCC3)c4cccc(c4)Cl</chem>
RUN:	RUN324
DDG (kcal/mol):	-0.94
dDDG (kcal/mol):	0.33

MAT-POS-fb82b63d-4_1



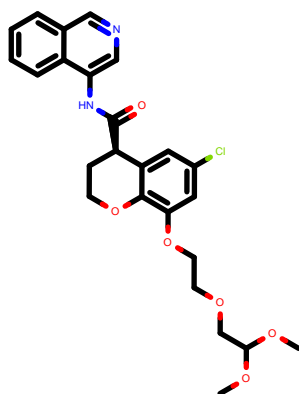
CID:	MAT-POS-fb82b63d-4_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](N)CCc1ccc4CC[N@@H+](3CC5CC5)Cl</chem>
RUN:	RUN3177
DDG (kcal/mol):	-0.93
dDDG (kcal/mol):	0.28

ERI-UCB-b3e6b0c2-13_2



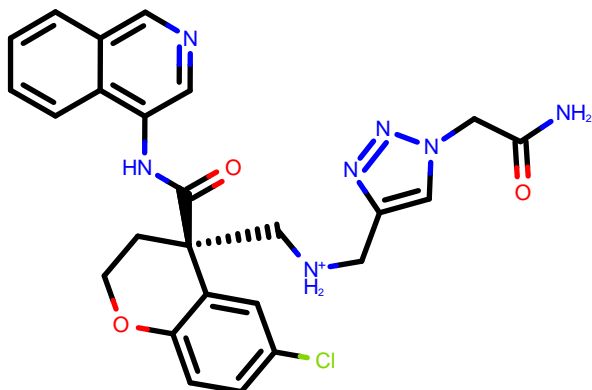
CID:	ERI-UCB-b3e6b0c2-13_2
SMILES:	<chem>C[N@]1C[C@H](c2c1ccc(c2)Cl)C(=O)Nc3ccc4c3ccc(c4)CN5CC[NH2+]CC5</chem>
RUN:	RUN3049
DDG (kcal/mol):	-0.93
dDDG (kcal/mol):	0.25

ALP-UNI-3496895b-2_7



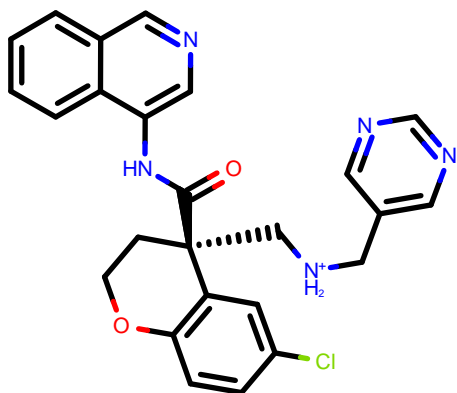
CID:	ALP-UNI-3496895b-2_7
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](N)CCOC3CCOC4CC(=O)N5CC[C@@H](C)C5[NH2+]COC4</chem>
RUN:	RUN2508
DDG (kcal/mol):	-0.93
dDDG (kcal/mol):	0.38

MAT-POS-61f37a1a-15_1



CID:	MAT-POS-61f37a1a-15_1
SMILES:	<chem>Cn1c[nC]1c(-O)[nH]c1(-O)n2C[C]N12+)[C]C(=O)Nc3cc4c3ccc(cc4)Cl)C1=CN=CN1CC(=O)N</chem>
RUN:	RUN4624
DDG (kcal/mol):	-0.93
dDDG (kcal/mol):	0.32

ED_-GRI-5b13fbe2-51_1



CID: ED_-GRI-5b13fbe2-51_1

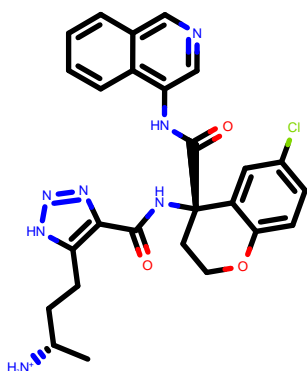
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OCC5(COC5)COC6C[NH2+]C6

RUN: RUN1598

DDG (kcal/mol): -0.92

dDDG (kcal/mol): 0.37

MAT-POS-fb82b63d-4_4



CID: MAT-POS-fb82b63d-4_4

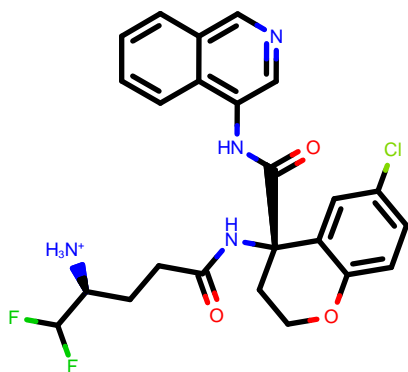
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3c4cc(ccc4CC[N@H+]3CC5CC5)Cl

RUN: RUN3183

DDG (kcal/mol): -0.92

dDDG (kcal/mol): 0.20

PET-UNK-824b5c6a-3_1



CID: PET-UNK-824b5c6a-3_1

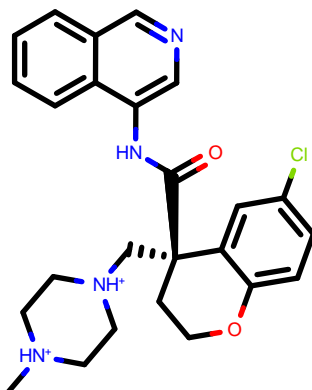
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OC5CC5

RUN: RUN3291

DDG (kcal/mol): -0.91

dDDG (kcal/mol): 0.23

RAL-THA-4aa06b95-7_4



CID: RAL-THA-4aa06b95-7_4

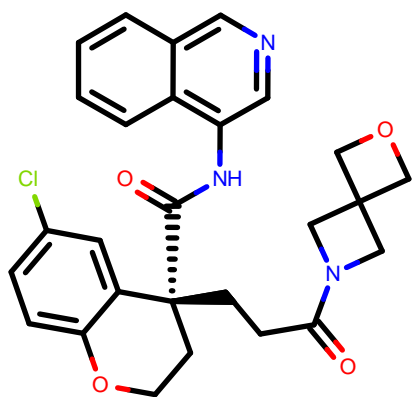
SMILES: COCCN1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN1246

DDG (kcal/mol): -0.91

dDDG (kcal/mol): 0.29

RAL-THA-e002e396-2_1



CID: RAL-THA-e002e396-2_1

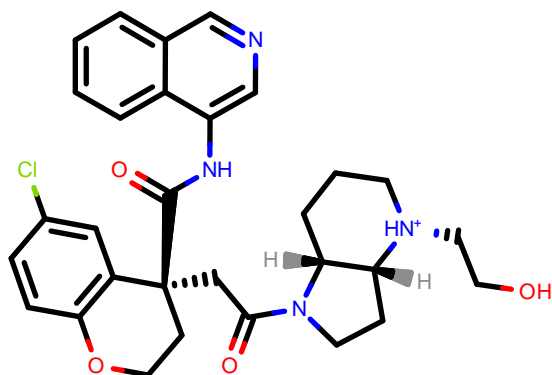
SMILES: CCS(=O)(=O)C[C@@]1(C)C(CO)C2C1CC(C2)C(=O)Nc3cnc4c3cccc4

RUN: RUN3453

DDG (kcal/mol): -0.91

dDDG (kcal/mol): 0.15

RAL-THA-8416115c-11_4



CID: RAL-THA-8416115c-11_4

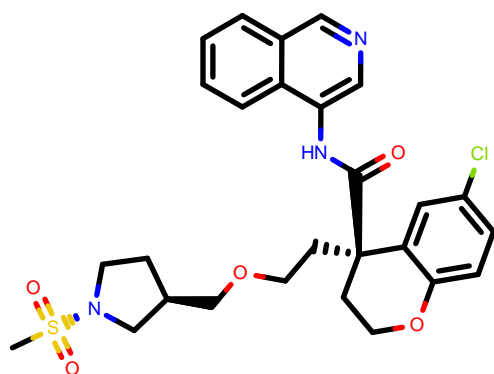
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H]3CCN(c4c3cc(cc4)Cl)C5cnc[nH]5

RUN: RUN1289

DDG (kcal/mol): -0.91

dDDG (kcal/mol): 0.53

MIC-UNK-5a93dd5f-2_2



CID: MIC-UNK-5a93dd5f-2_2

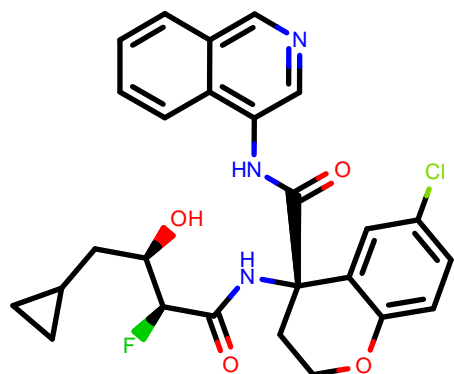
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H](C3CCCC3)N4[C@@H](C5CCCC[C@H]5C4)S(=O)(=O)N

RUN: RUN735

DDG (kcal/mol): -0.91

dDDG (kcal/mol): 0.41

DAR-DIA-f6ee7aeb-2_2



CID: DAR-DIA-f6ee7aeb-2_2

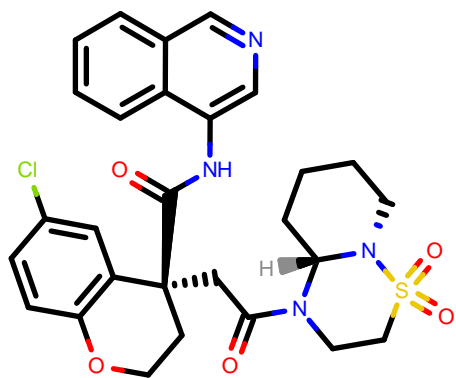
SMILES: c1ccc2c(c1)cnc2N3C[C@H]([C@@H](C3=O)c4cc(cc4)Cl)OCC(F)(F)F5CCCC5C#N

RUN: RUN3405

DDG (kcal/mol): -0.90

dDDG (kcal/mol): 0.18

BEN-DND-a7517465-1_1



CID: BEN-DND-a7517465-1_1

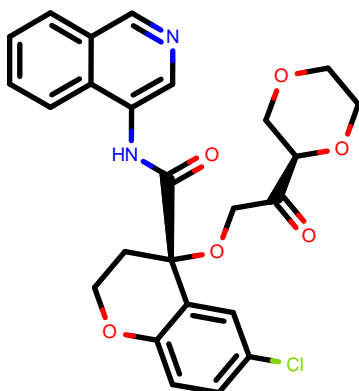
SMILES: Cc1cccc2c1c(cnc2)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl

RUN: RUN1477

DDG (kcal/mol): -0.90

dDDG (kcal/mol): 0.26

MAK-UNK-ffc90da7-9_4



CID: MAK-UNK-ffc90da7-9_4

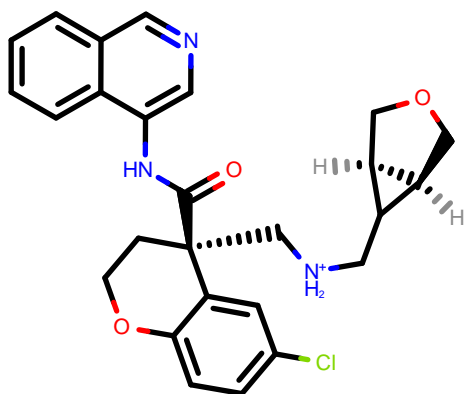
SMILES: C[C@H]1[C@H](CCO1)SC[C@@H](C)c2ccc3c(c2)ncc3NC(=O)C4CCCC4Cl

RUN: RUN715

DDG (kcal/mol): -0.90

dDDG (kcal/mol): 0.28

EDG-MED-90036822-108_1



CID: EDG-MED-90036822-108_1

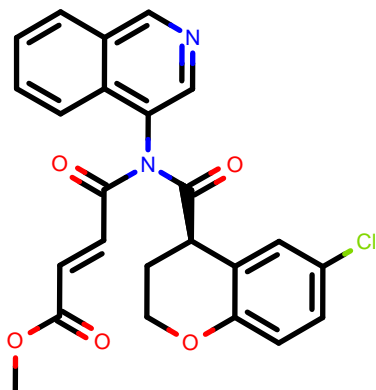
SMILES: c1ccc2c(c1)ncc2NC(=O)[C@@H]3[C@@H](C)CCOC4c3cc(cc4)C(=O)C[C@@H]5[C@@H](C)CC5O

RUN: RUN1846

DDG (kcal/mol): -0.90

dDDG (kcal/mol): 0.34

MIC-UNK-9582b2c5-2_4



CID: MIC-UNK-9582b2c5-2_4

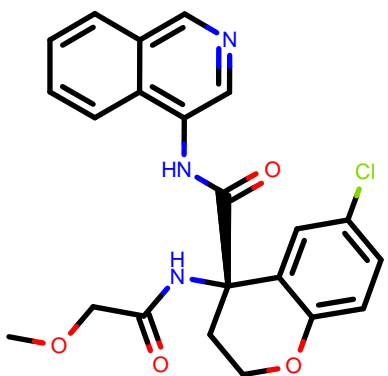
SMILES: CC(=O)N1CC[C@H]2[C@H](C1)C[C@@H](C(=O)N2c3cccc(c3)Cl)c4cccc5c4ccoc5

RUN: RUN265

DDG (kcal/mol): -0.90

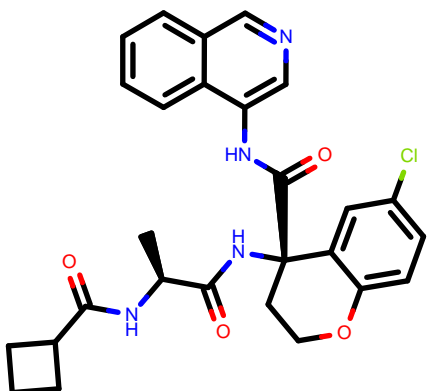
dDDG (kcal/mol): 0.36

EDJ-MED-4f704dc9-1_2



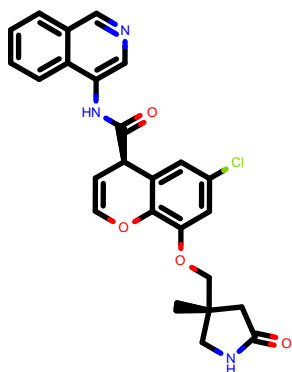
CID:	EDJ-MED-4f704dc9-1_2
SMILES:	<chem>CO[C@]1(CCNC2c1cc(c(c2)Cl)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3160
DDG (kcal/mol):	-0.90
dDDG (kcal/mol):	0.10

RAL-THA-4aa06b95-3_2



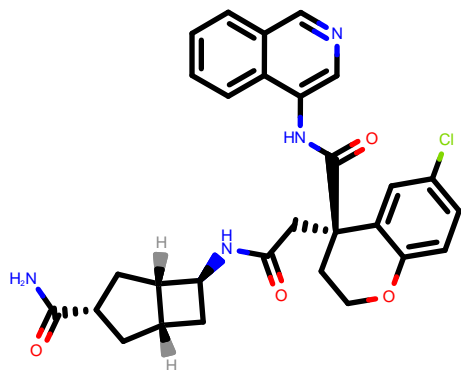
CID:	RAL-THA-4aa06b95-3_2
SMILES:	<chem>CNC(=O)N1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN1234
DDG (kcal/mol):	-0.89
dDDG (kcal/mol):	0.28

MAT-POS-2905de8c-2_1



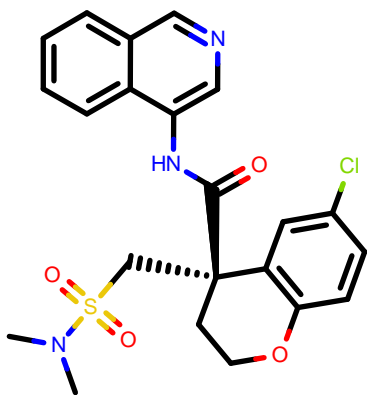
CID:	MAT-POS-2905de8c-2_1
SMILES:	<chem>C[NH2+][C@@]1(CCOC2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN2229
DDG (kcal/mol):	-0.89
dDDG (kcal/mol):	0.38

MAK-UNK-c749d764-33_3



CID:	MAK-UNK-c749d764-33_3
SMILES:	<chem>CS(=O)(=O)N(c1cncc2c1cccc2)C(=O)C[C@@H]3CCC[C@H](C[C@@H]3O)C(F)F</chem>
RUN:	RUN1081
DDG (kcal/mol):	-0.89
dDDG (kcal/mol):	0.43

ED_-GRI-5b13fbe2-4_1



CID: ED_-GRI-5b13fbe2-4_1

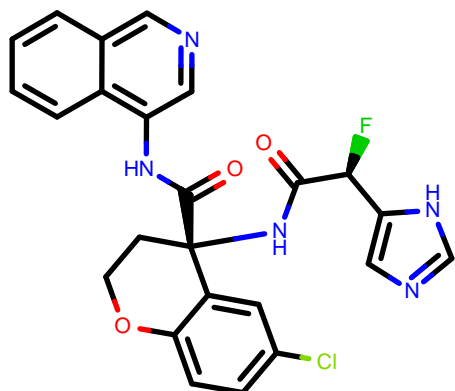
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3(CCOc4ccc(Cl)OC)C@H](c5[nH]n5)[NH3+]

RUN: RUN1525

DDG (kcal/mol): -0.88

dDDG (kcal/mol): 0.27

VLA-UNK-ba665ac8-2_1



CID: VLA-UNK-ba665ac8-2_1

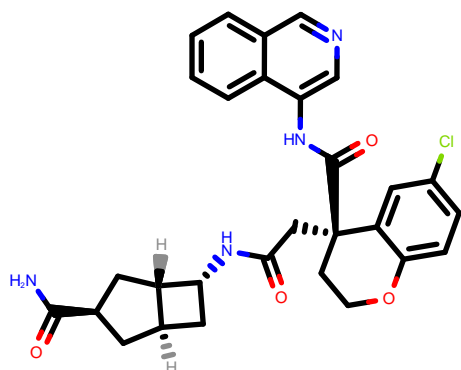
SMILES: C[N@@H]1CC(=O)N(C(=O)[C@@]12CCOC3c2cc(Cl)C)C4=CC=CC=C4

RUN: RUN3301

DDG (kcal/mol): -0.88

dDDG (kcal/mol): 0.17

DAR-DIA-9e4459de-13_16



CID: DAR-DIA-9e4459de-13_16

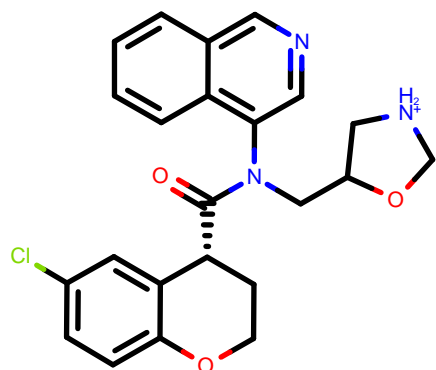
SMILES: c1cc2c(c1)NC(=O)C3c4c(c3)ncnc4N(C(=O)[C@@]5(CCOc6ccc(Cl)OC)C@H)(c2O)[C@@H]7C(=O)NC7=O

RUN: RUN1440

DDG (kcal/mol): -0.88

dDDG (kcal/mol): 0.36

ALF-EVA-ced740bd-3_2



CID: ALF-EVA-ced740bd-3_2

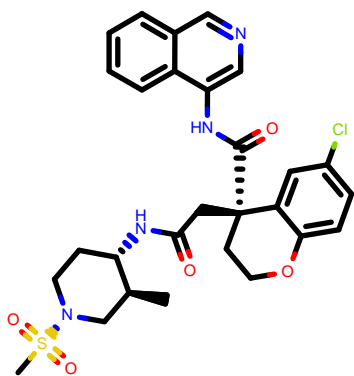
SMILES: c1cc2cnc(c2cc1F)NC(=O)[C@H]3COc4c3cc(cc4)F

RUN: RUN2793

DDG (kcal/mol): -0.88

dDDG (kcal/mol): 0.14

MAK-UNK-c749d764-21_3



CID: MAK-UNK-c749d764-21_3

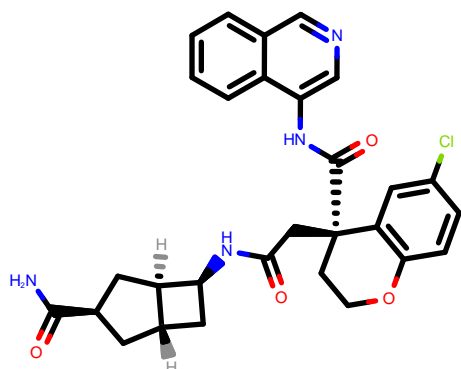
SMILES: CCCCN(c1cncc2c1cccc2)C(=O)C[C@@H]3CCC[C@H]([C@@H]3O)C(F)F

RUN: RUN1016

DDG (kcal/mol): -0.86

dDDG (kcal/mol): 0.23

LON-WEI-5e7d1b3e-40_1



CID: LON-WEI-5e7d1b3e-40_1

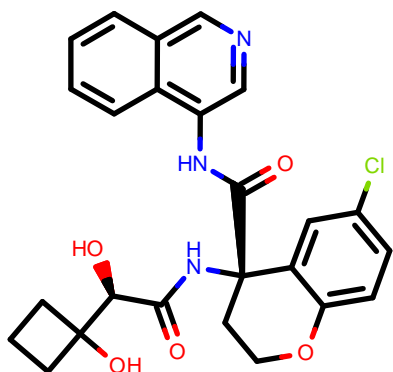
SMILES: CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)NCCC3=CCCC3

RUN: RUN1352

DDG (kcal/mol): -0.85

dDDG (kcal/mol): 0.22

KAD-UNI-877d7bed-14_1



CID: KAD-UNI-877d7bed-14_1

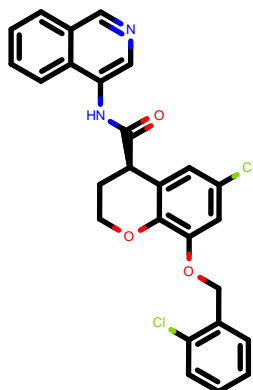
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4c3cc(cc4OCC(=O)CNC(=O)C(F)F)O

RUN: RUN3760

DDG (kcal/mol): -0.85

dDDG (kcal/mol): 0.16

DAR-DIA-9e4459de-15_15



CID: DAR-DIA-9e4459de-15_15

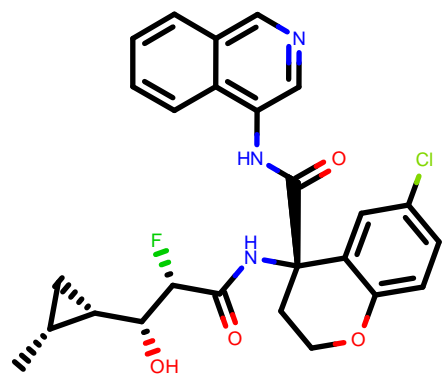
SMILES: c1cc2c(c1)NC(=O)C(Cc1ccc(Cl)cc1)O[C@@H]3CCOCc4c3cc(Cl)c1c(c2O)[C@@H]7CCCl=O)NC7=O

RUN: RUN1453

DDG (kcal/mol): -0.85

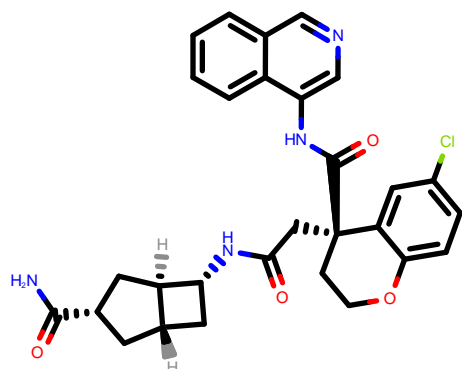
dDDG (kcal/mol): 0.46

PET-UNK-7a31b064-2_1



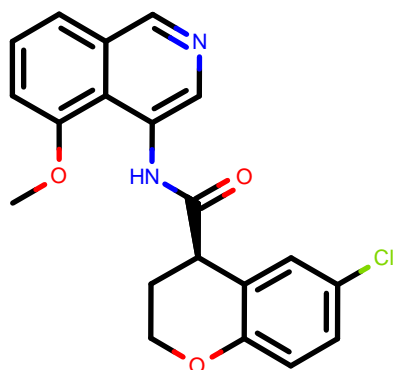
CID:	PET-UNK-7a31b064-2_1
SMILES:	<chem>CCN1C[C@H](C(=O)N(C1=O)c2cncc3c2cccc3)c4cccc(c4)Cl</chem>
RUN:	RUN3290
DDG (kcal/mol):	-0.85
dDDG (kcal/mol):	0.17

NAU-LAT-4ce8bf23-1_2



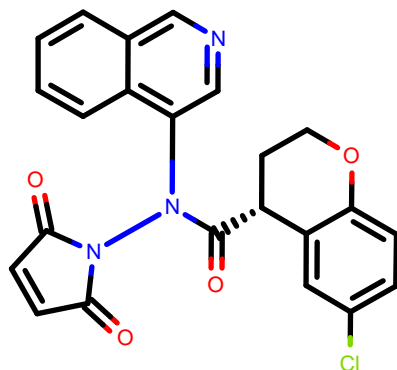
CID:	NAU-LAT-4ce8bf23-1_2
SMILES:	<chem>CC(=O)NC[C@H](c1cncc2c1cccc2)C(=O)Nc3cccc(c3)Cl</chem>
RUN:	RUN1394
DDG (kcal/mol):	-0.85
dDDG (kcal/mol):	0.41

MAT-POS-8a69d52e-1_3



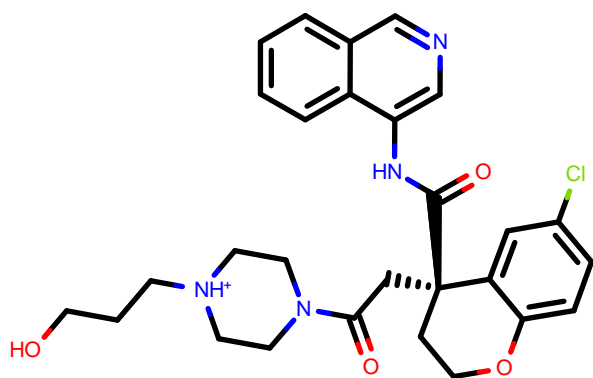
CID:	MAT-POS-8a69d52e-1_3
SMILES:	<chem>C[C@@H]1C[C@H](c2cc(ccc2O1)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN362
DDG (kcal/mol):	-0.85
dDDG (kcal/mol):	0.24

BEN-DND-34fc7f90-8_1



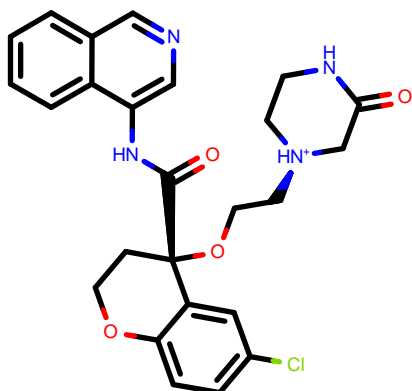
CID:	BEN-DND-34fc7f90-8_1
SMILES:	<chem>C[N@@H+][C]2cc(c(cc2[C@@H](C1)C(=O)Nc3cncc4c3cccc4)Cl)Cl</chem>
RUN:	RUN3674
DDG (kcal/mol):	-0.84
dDDG (kcal/mol):	0.12

BEN-DND-c852c98b-6_2



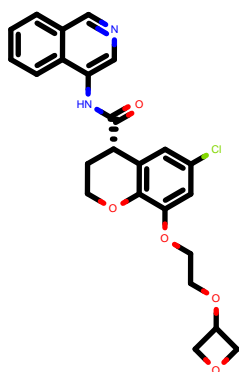
CID:	BEN-DND-c852c98b-6_2
SMILES:	<chem>c1cc2nccc(c2cc1OC(F)F)NC(=O)[C@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN1216
DDG (kcal/mol):	-0.84
dDDG (kcal/mol):	0.34

DAR-DIA-6a508060-10_1



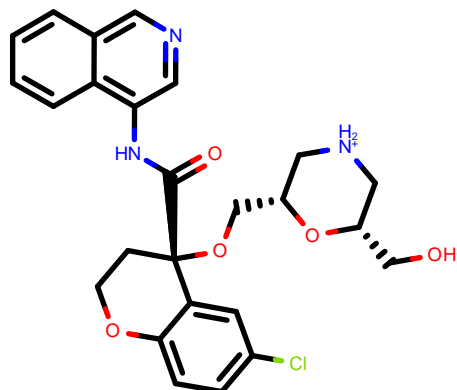
CID:	DAR-DIA-6a508060-10_1
SMILES:	<chem>C[C@@]1(CCOC2c1cc(cc2)Cl)C(=O)Nc3nccc4c3cccc4</chem>
RUN:	RUN349
DDG (kcal/mol):	-0.84
dDDG (kcal/mol):	0.36

MAT-POS-f9802937-4_1



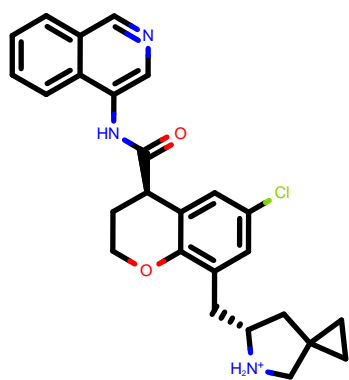
CID:	MAT-POS-f9802937-4_1
SMILES:	<chem>CN(C)C(=O)C[C@@]1(CCOC2c1cc(cc2)Cl)C(=O)Nc3nccc4c3cccc4</chem>
RUN:	RUN2397
DDG (kcal/mol):	-0.84
dDDG (kcal/mol):	0.29

DAR-DIA-0f2f46c9-1_2



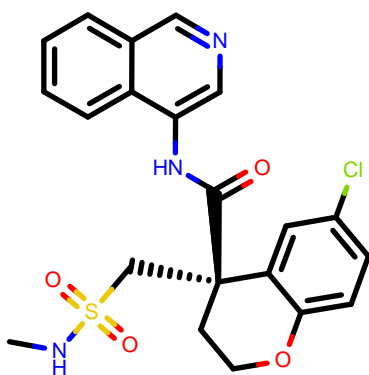
CID:	DAR-DIA-0f2f46c9-1_2
SMILES:	<chem>c1ccc2c(c1)nccc2NC(=O)[C@H]3CC[N@@]4(c4c3cc(cc4)Cl)S(=O)(=O)N</chem>
RUN:	RUN3219
DDG (kcal/mol):	-0.84
dDDG (kcal/mol):	0.23

ALP-POS-3b848b35-4_1



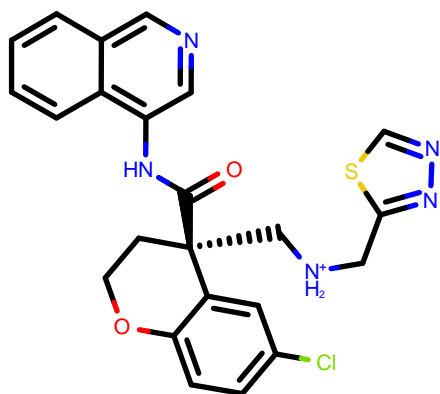
CID:	ALP-POS-3b848b35-4_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](CCC3CCCCC3)c4cccc(c4)Cl</chem>
RUN:	RUN61
DDG (kcal/mol):	-0.83
dDDG (kcal/mol):	0.34

ED_-GRI-5b13fbe2-12_2



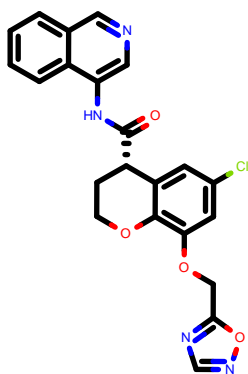
CID:	ED_-GRI-5b13fbe2-12_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](C(=O)N)OCC[N@H](CO)CCOCCO</chem>
RUN:	RUN1539
DDG (kcal/mol):	-0.83
dDDG (kcal/mol):	0.24

EDJ-MED-1981ceba-3_4



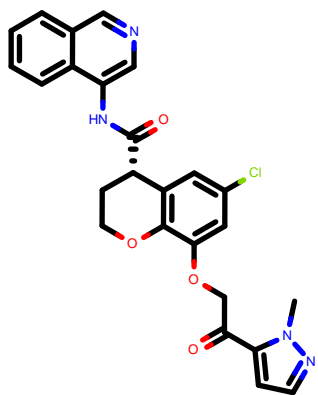
CID:	EDJ-MED-1981ceba-3_4
SMILES:	<chem>COC1CN(C1)S(=O)(=O)[N@]2Cc3ccc(cc3[C@@H](C2)C(=O)Nc4cncc5c4cccc5)Cl</chem>
RUN:	RUN4692
DDG (kcal/mol):	-0.82
dDDG (kcal/mol):	0.24

EDG-MED-5d232de5-5_2



CID:	EDG-MED-5d232de5-5_2
SMILES:	<chem>CN1CC[C@@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN2368
DDG (kcal/mol):	-0.82
dDDG (kcal/mol):	0.24

ALP-UNI-3496895b-2_5



CID: ALP-UNI-3496895b-2_5

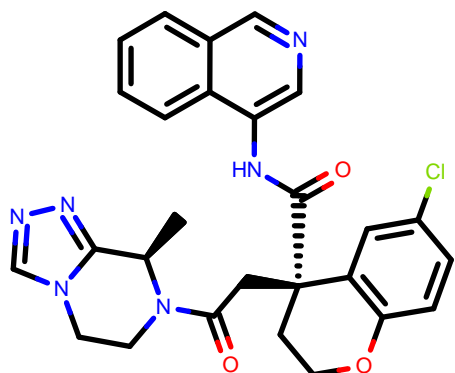
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@H](Cl)C3C(=O)N(C)C3C(=O)N(C)C3C(=O)N(C)C3

RUN: RUN2506

DDG (kcal/mol): -0.82

dDDG (kcal/mol): 0.25

BEN-DND-f2e727cd-5_1



CID: BEN-DND-f2e727cd-5_1

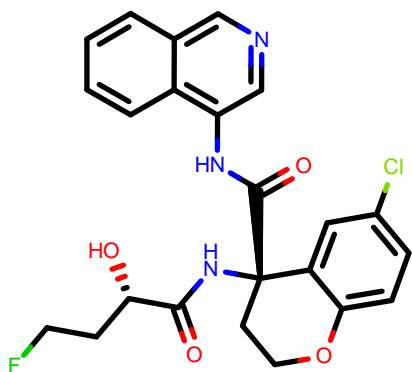
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@H](Cl)C3C(=O)N(C)C3C(=O)N(C)C3

RUN: RUN1196

DDG (kcal/mol): -0.82

dDDG (kcal/mol): 0.22

MIC-UNK-b9827f26-4_1



CID: MIC-UNK-b9827f26-4_1

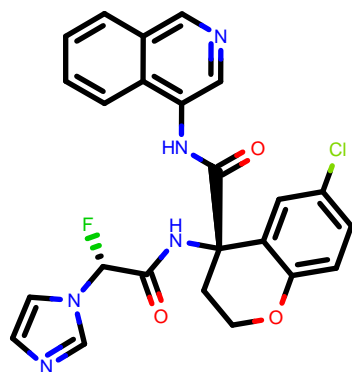
SMILES: COc1ccc(cc1N2CCN(CC=O)C(=O)c3ncnc4c3ccc4)Cl

RUN: RUN3255

DDG (kcal/mol): -0.81

dDDG (kcal/mol): 0.15

MAT-POS-fb82b63d-3_1



CID: MAT-POS-fb82b63d-3_1

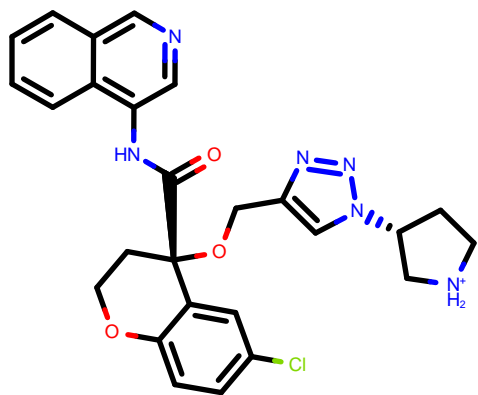
SMILES: C[N@@H+]1CCc2ccc(cc2[C@@H]1C(=O)Nc3ncnc4c3ccc4)Cl

RUN: RUN3173

DDG (kcal/mol): -0.81

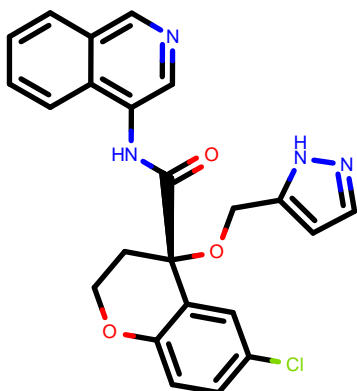
dDDG (kcal/mol): 0.14

JAG-UCB-f37eaa14-8_1



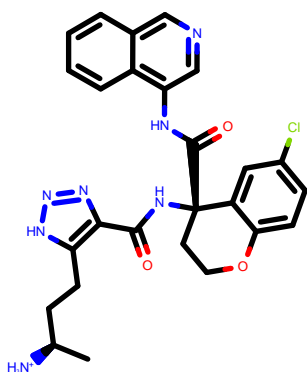
CID:	JAG-UCB-f37eaa14-8_1
SMILES:	<chem>c1ccc2c(c1)ncnc2N3CC[C@@H](C3=O)C[N@@](c5c4cc(cc5)C)CCn6ccc6</chem>
RUN:	RUN3068
DDG (kcal/mol):	-0.81
dDDG (kcal/mol):	0.20

ALP-UNI-ba800595-1_2



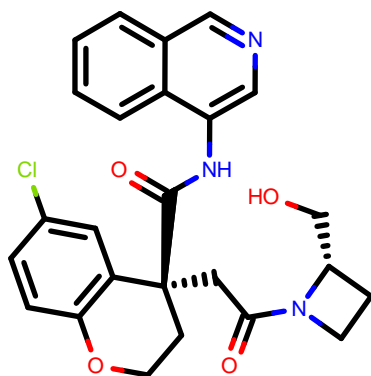
CID:	ALP-UNI-ba800595-1_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@H]3CCNc4c3cc(cc4Cl)Cl</chem>
RUN:	RUN3073
DDG (kcal/mol):	-0.81
dDDG (kcal/mol):	0.11

EDJ-MED-4f704dc9-1_1



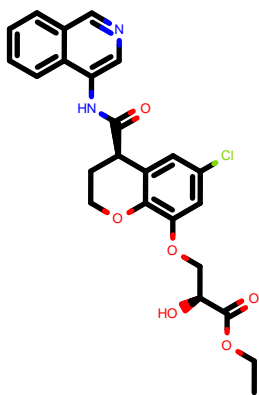
CID:	EDJ-MED-4f704dc9-1_1
SMILES:	<chem>CO[C@@]1(CCNc2c1cc(c(c2)Cl)Cl)C(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN3159
DDG (kcal/mol):	-0.81
dDDG (kcal/mol):	0.23

RAL-THA-8416115c-10_3



CID:	RAL-THA-8416115c-10_3
SMILES:	<chem>CNC(=O)CN1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN1285
DDG (kcal/mol):	-0.81
dDDG (kcal/mol):	0.21

KAD-UNI-8a629cb0-38_2



CID: KAD-UNI-8a629cb0-38_2

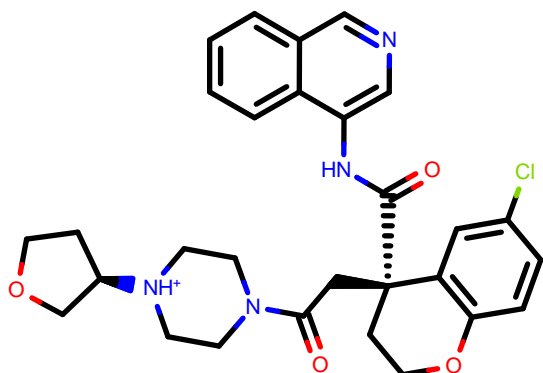
SMILES: C[N@@]1[C@@H](CCS1(=O)=O)COC[C@]2(CCOc3c2cc(cc3)C)C(=O)Nc4ncc5c4cccc5

RUN: RUN2124

DDG (kcal/mol): -0.81

dDDG (kcal/mol): 0.41

MAK-UNK-c749d764-15_15



CID: MAK-UNK-c749d764-15_15

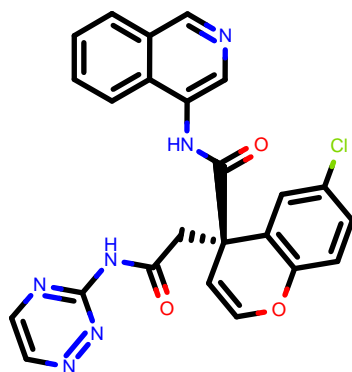
SMILES: C[C@@H](N(c1cncc2c1ccc2)C(=O)C)C[C@H]3CCC@H3O(C)F)OCC4CCCC4

RUN: RUN972

DDG (kcal/mol): -0.80

dDDG (kcal/mol): 0.26

MIC-UNK-5a93dd5f-11_2



CID: MIC-UNK-5a93dd5f-11_2

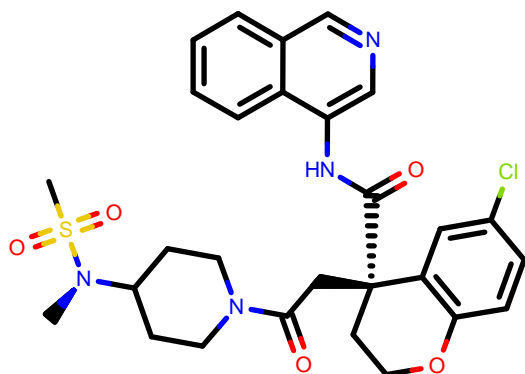
SMILES: c1ccc2c(c1)cncc2N(C(=O)[C@H](c3ccccc(c3)C)N)N4CCC(CC4)N5C5CCCC5

RUN: RUN789

DDG (kcal/mol): -0.80

dDDG (kcal/mol): 0.31

JOH-UNI-a38a7bdd-5_4



CID: JOH-UNI-a38a7bdd-5_4

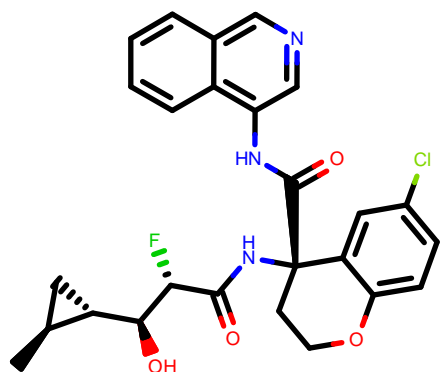
SMILES: c1ccc2c(c1)cncc2N(C(=O)Cc3ccccc(c3)C)C(=O)[C@H]4C[C@H]4F

RUN: RUN1486

DDG (kcal/mol): -0.80

dDDG (kcal/mol): 0.46

MIC-UNK-0a05c952-4_2



CID: MIC-UNK-0a05c952-4_2

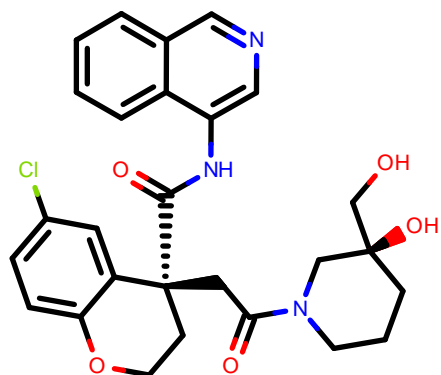
SMILES: c1ccc2c(c1)cncc2N3[C@H](CC[C@H](C3=O)c4ccc(c(c4)Cl)Cl)[C@H]5CO5

RUN: RUN3522

DDG (kcal/mol): -0.80

dDDG (kcal/mol): 0.17

MAK-UNK-c749d764-26_8



CID: MAK-UNK-c749d764-26_8

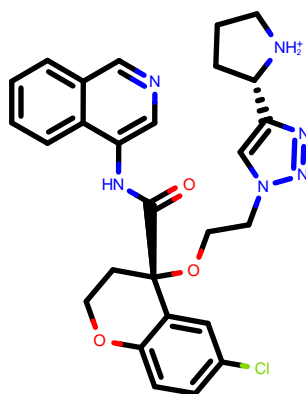
SMILES: c1ccc2c(c1)cncc2NC(=O)C[C@H]3CCCC[C@H]1([C@H]3O)Cl

RUN: RUN1053

DDG (kcal/mol): -0.80

dDDG (kcal/mol): 0.45

MIC-UNK-bcd487e9-6_1



CID: MIC-UNK-bcd487e9-6_1

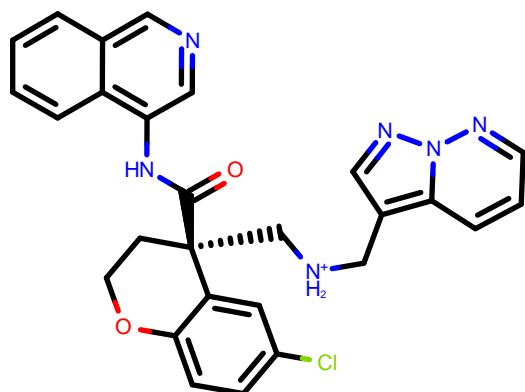
SMILES: c1ccc2c(c1)cncc2NC(=O)N(C[C@H]3CCCCS3(=O)=O)c4cccc(c4)Cl

RUN: RUN592

DDG (kcal/mol): -0.80

dDDG (kcal/mol): 0.40

EDG-MED-4c68219f-14_1



CID: EDG-MED-4c68219f-14_1

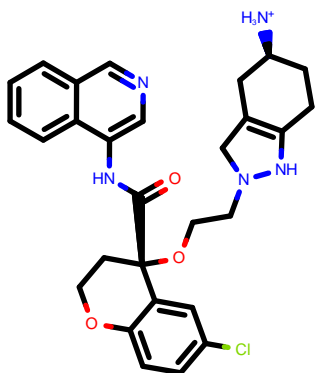
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)O[C@H]5C[C@H]([NH2+][C5])CO

RUN: RUN1649

DDG (kcal/mol): -0.79

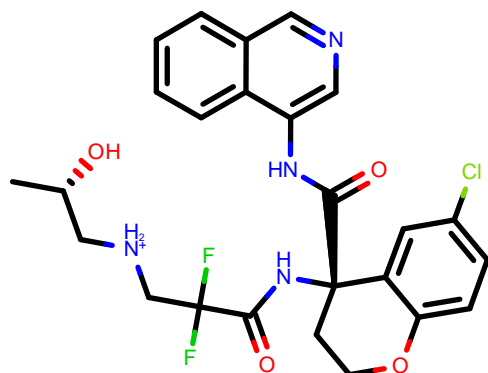
dDDG (kcal/mol): 0.40

FRA-DIA-b66f7109-2_2



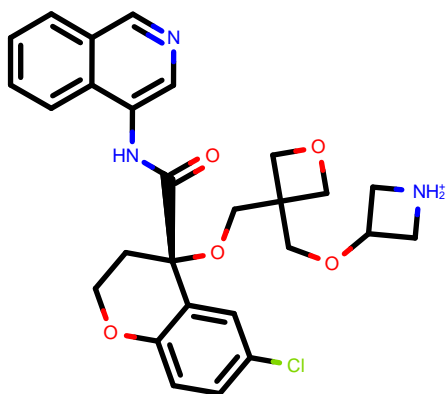
CID:	FRA-DIA-b66f7109-2_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3COc4c3cc(cc4Cl)O[C@H]5CC(=O)N5</chem>
RUN:	RUN394
DDG (kcal/mol):	-0.79
dDDG (kcal/mol):	0.37

MAK-UNK-8be7dca9-10_2



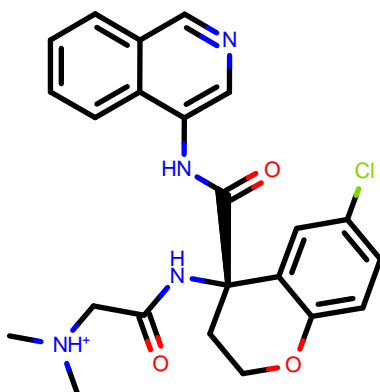
CID:	MAK-UNK-8be7dca9-10_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3C[C@@H](Oc4c3cc(cc4Cl)C(=O)O)[O-]</chem>
RUN:	RUN512
DDG (kcal/mol):	-0.79
dDDG (kcal/mol):	0.38

FRA-DIA-b66f7109-1_1



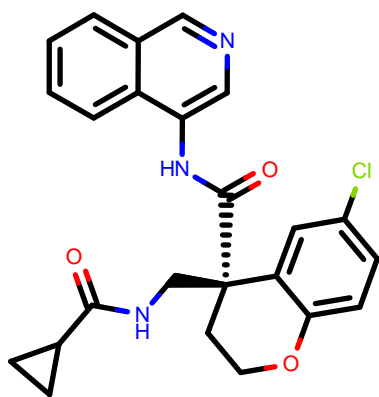
CID:	FRA-DIA-b66f7109-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CNc4c3cc(cc4Cl)O[C@@H]5CC(=O)N5</chem>
RUN:	RUN391
DDG (kcal/mol):	-0.79
dDDG (kcal/mol):	0.33

ALP-UNI-44c99a80-2_1



CID:	ALP-UNI-44c99a80-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2N(Cc3cnc[nH]3)C(=O)[C@H]4CCOc5c4cc(cc5)Cl</chem>
RUN:	RUN463
DDG (kcal/mol):	-0.79
dDDG (kcal/mol):	0.34

RAL-THA-8416115c-5_2



CID: RAL-THA-8416115c-5_2

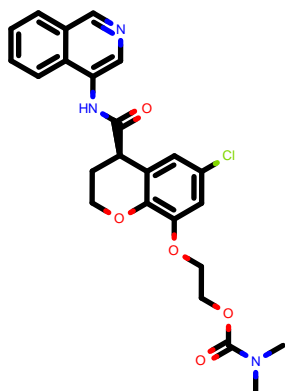
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H]3CCN(c4c3cc(cc4)Cl)Cc5[nH]cn5

RUN: RUN1264

DDG (kcal/mol): -0.79

dDDG (kcal/mol): 0.23

ALP-POS-5bb456a5-1_5



CID: ALP-POS-5bb456a5-1_5

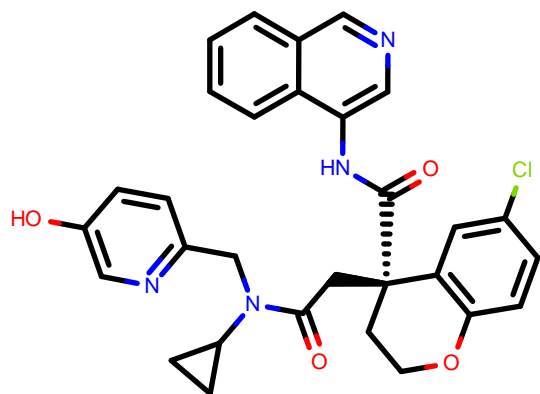
SMILES: CN(C)COCOC1=CC=C(Cl)C2=CC=C(NC(=O)C3=CN=C4C=CC=CC34)O2

RUN: RUN2410

DDG (kcal/mol): -0.79

dDDG (kcal/mol): 0.42

BEN-DND-c852c98b-6_1



CID: BEN-DND-c852c98b-6_1

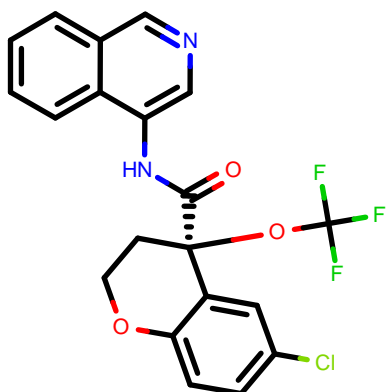
SMILES: Oc1ccc(nc1)CN(C)C(=O)C[C@H]23CCOC4C3CC(=O)N2C

RUN: RUN1213

DDG (kcal/mol): -0.79

dDDG (kcal/mol): 0.28

MAK-UNK-c749d764-10_4



CID: MAK-UNK-c749d764-10_4

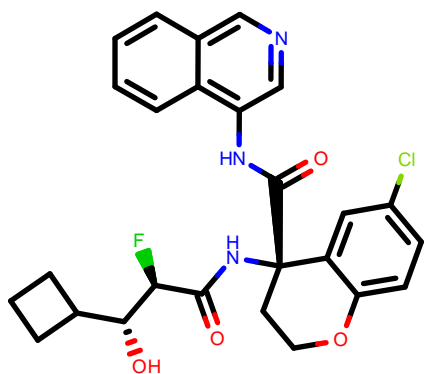
SMILES: CSN(c1cnc2c1cccc2)C(=O)C[C@H]3CCC[C@H]1[C@@H]3O(C(F)(F)F)

RUN: RUN944

DDG (kcal/mol): -0.79

dDDG (kcal/mol): 0.08

DAR-DIA-0587064e-1_1



CID: DAR-DIA-0587064e-1_1

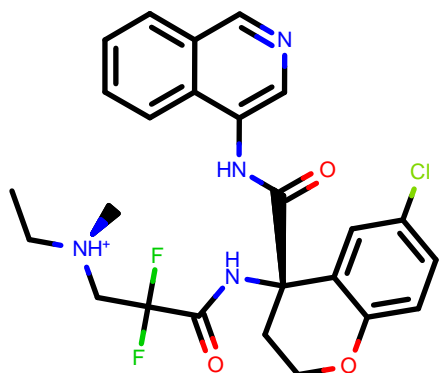
SMILES: c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)OCCC(F)(F)F

RUN: RUN3348

DDG (kcal/mol): -0.78

dDDG (kcal/mol): 0.15

ROB-UNI-611831f5-2_2



CID: ROB-UNI-611831f5-2_2

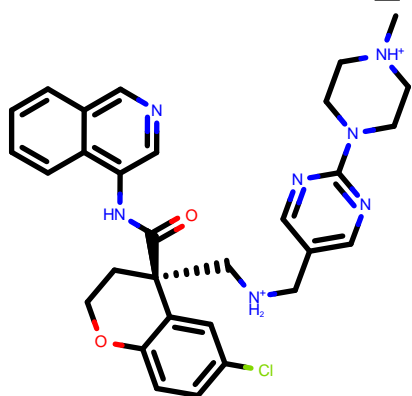
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3CCNc4c3cc(cn4)Cl

RUN: RUN3324

DDG (kcal/mol): -0.78

dDDG (kcal/mol): 0.18

ALP-POS-347519b5-3_58



CID: ALP-POS-347519b5-3_58

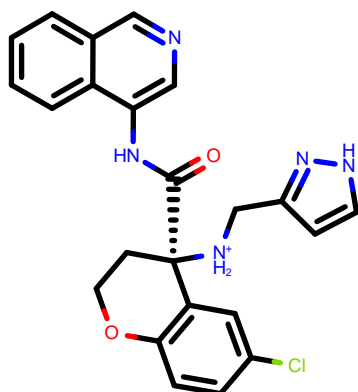
SMILES: CS(=O)(=O)N(B)1CC@@H2C@@H3CC[C@H]1[C@H]2C@H1(C1C(=O)N4cccc54cccc5)O3

RUN: RUN4325

DDG (kcal/mol): -0.78

dDDG (kcal/mol): 0.32

EDG-MED-90036822-61_1



CID: EDG-MED-90036822-61_1

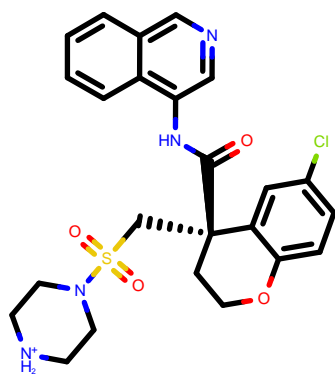
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)NC(=O)C[NH+]5CCCS5

RUN: RUN1745

DDG (kcal/mol): -0.77

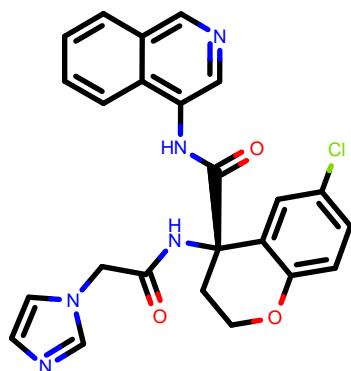
dDDG (kcal/mol): 0.26

ALP-POS-347519b5-1_35



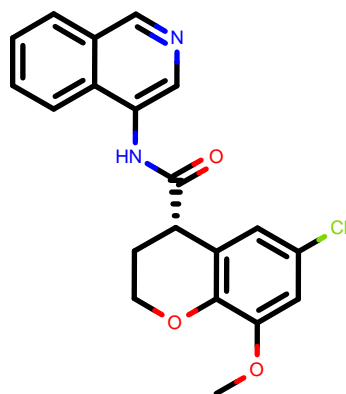
CID:	ALP-POS-347519b5-1_35
SMILES:	<chem>CS(=O)(=O)N[C@@H]1C[C@@H]2[C@@H]3CC[C@@H]3C[C@@H]2[C@@H]1C(=O)Nc4ccc5c4ccc5</chem>
RUN:	RUN4250
DDG (kcal/mol):	-0.77
dDDG (kcal/mol):	0.24

EDJ-MED-37aac4bd-4_1



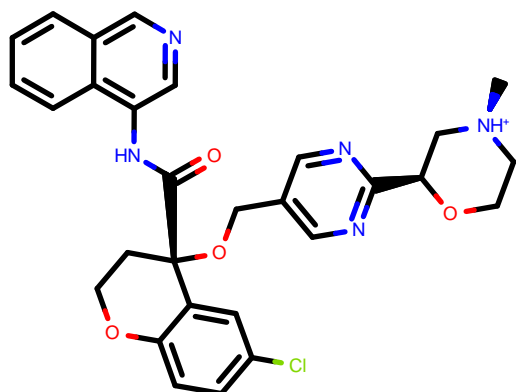
CID:	EDJ-MED-37aac4bd-4_1
SMILES:	<chem>CO[C@@H]1(CCOc2c1cc(cc2F)F)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3145
DDG (kcal/mol):	-0.77
dDDG (kcal/mol):	0.13

MIC-UNK-5a93dd5f-2_3



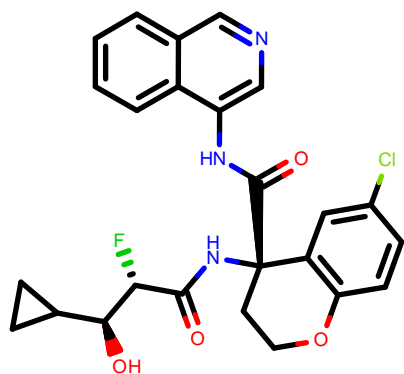
CID:	MIC-UNK-5a93dd5f-2_3
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](c3cccc(c3)Cl)N[C@@H]4C[C@@H]5CCCC[C@@H]5C4</chem>
RUN:	RUN736
DDG (kcal/mol):	-0.77
dDDG (kcal/mol):	0.14

EDJ-MED-f893e2a1-2_1



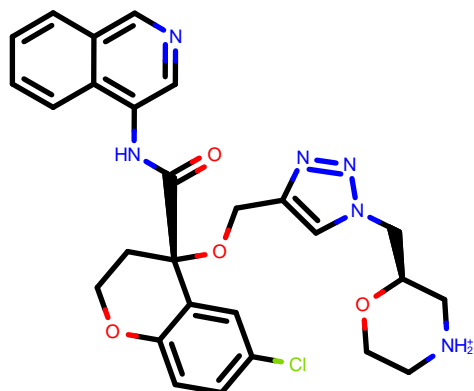
CID:	EDJ-MED-f893e2a1-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](c3(CCOc4c3cc(cc4)Cl)CNC[C@@H]5(CCOc5)C#N</chem>
RUN:	RUN3195
DDG (kcal/mol):	-0.77
dDDG (kcal/mol):	0.27

PET-UNK-a692de38-1_1



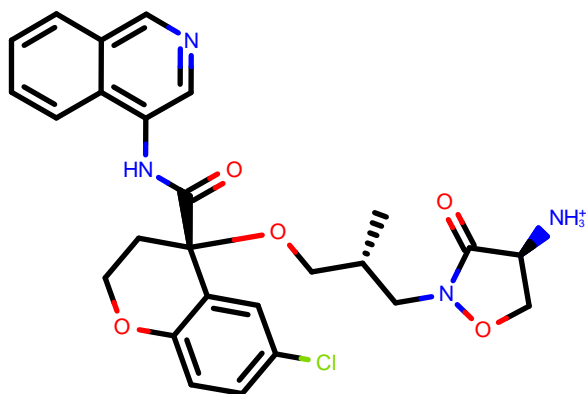
CID:	PET-UNK-a692de38-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2N(C(=O)Cc3ccccc(c3)Cl)NC#N</chem>
RUN:	RUN583
DDG (kcal/mol):	-0.76
dDDG (kcal/mol):	0.26

WIL-UCB-7ba4ac3a-2_1



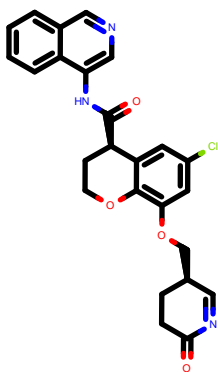
CID:	WIL-UCB-7ba4ac3a-2_1
SMILES:	<chem>CN(c1cncc2c1cccc2)C(=O)[C@@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN3024
DDG (kcal/mol):	-0.76
dDDG (kcal/mol):	0.23

ALP-POS-fe871b40-3_1



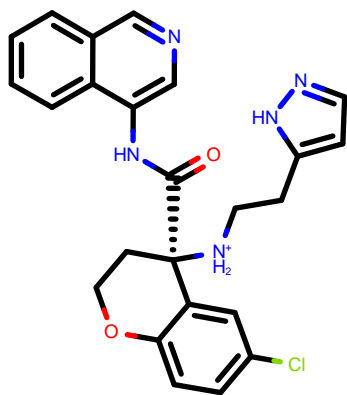
CID:	ALP-POS-fe871b40-3_1
SMILES:	<chem>CO[C@@@]1(CCOc2c1cc(cc2C#N)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3113
DDG (kcal/mol):	-0.76
dDDG (kcal/mol):	0.22

JOH-UNI-ea72002d-3_3



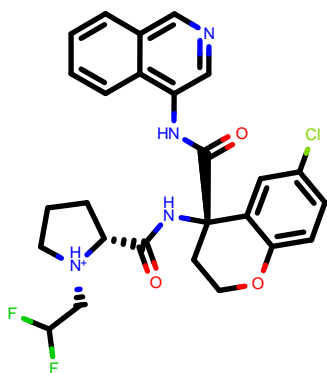
CID:	JOH-UNI-ea72002d-3_3
SMILES:	<chem>c1ccc2c(c1)cncc2[C@@H](C(=O)[C@@H]3CCOC4c3cc(cc4)Cl)N5C(=O)C=C5=O</chem>
RUN:	RUN2486
DDG (kcal/mol):	-0.76
dDDG (kcal/mol):	0.52

ALP-POS-ce760d3f-7_2



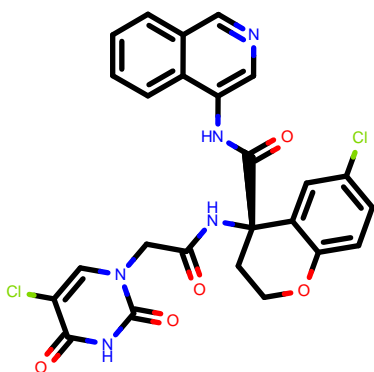
CID:	ALP-POS-ce760d3f-7_2
SMILES:	<chem>c1cc2cncc(c2cc1F)NC(=O)[C@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN1464
DDG (kcal/mol):	-0.76
dDDG (kcal/mol):	0.29

MAT-POS-dd3ad2b5-5_2



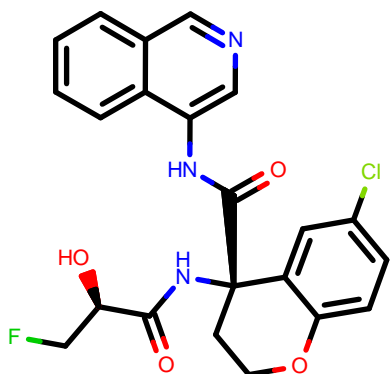
CID:	MAT-POS-dd3ad2b5-5_2
SMILES:	<chem>CNC(=O)N1Cc2ccc(cc2[C@H](C1)C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN3544
DDG (kcal/mol):	-0.75
dDDG (kcal/mol):	0.16

MIC-UNK-5a93dd5f-3_8



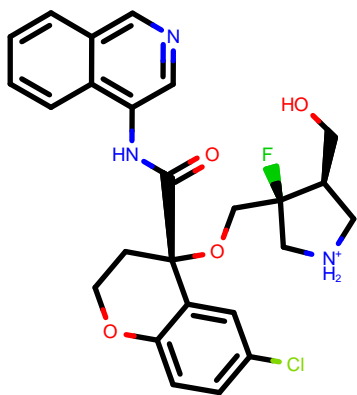
CID:	MIC-UNK-5a93dd5f-3_8
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H](c3cccc(c3)Cl)[N@H]4CC[C@H]5CCCC[C@]54</chem>
RUN:	RUN746
DDG (kcal/mol):	-0.75
dDDG (kcal/mol):	0.33

VLA-UCB-34f3ed0c-21_1



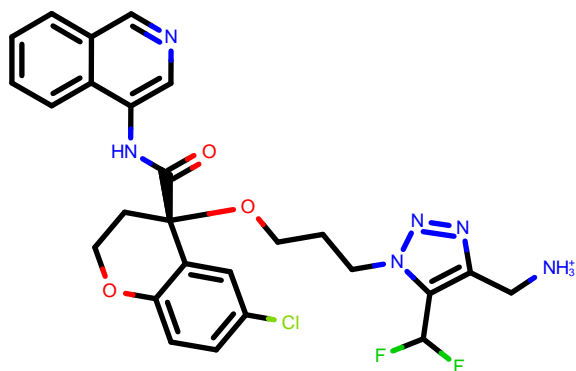
CID:	VLA-UCB-34f3ed0c-21_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(Cc4cc4c3cc(cc4)Cl)C(=O)CC[C@@]3[C@@H]5CC(=O)Nc5</chem>
RUN:	RUN648
DDG (kcal/mol):	-0.75
dDDG (kcal/mol):	0.29

VLA-UCB-50c39ae8-7_1



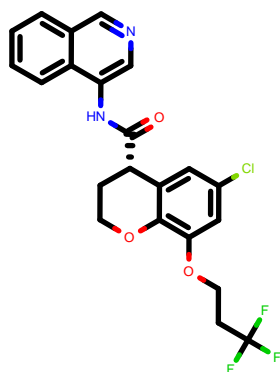
CID:	VLA-UCB-50c39ae8-7_1
SMILES:	<chem>C=CC(=O)N(c1cncc2c1cccc2)C(=O)[C@@H]3CCOc4c3cc(cc4)Cl</chem>
RUN:	RUN386
DDG (kcal/mol):	-0.75
dDDG (kcal/mol):	0.30

DAR-DIA-0587064e-21_2



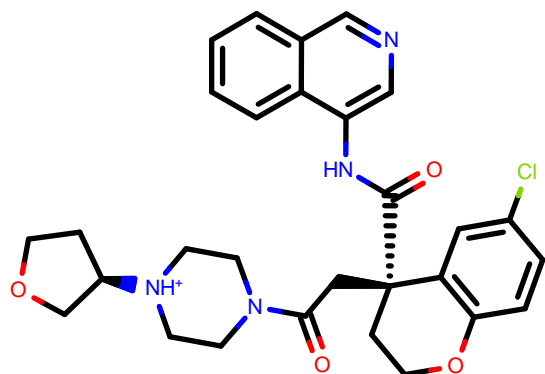
CID:	DAR-DIA-0587064e-21_2
SMILES:	<chem>c1ccc(cc1)CCOc2cc(cc3c2NCC[C@H]3C(=O)Nc4cncc5c4cccc5)Cl</chem>
RUN:	RUN3376
DDG (kcal/mol):	-0.75
dDDG (kcal/mol):	0.22

JOH-UNI-a38a7bdd-6_4



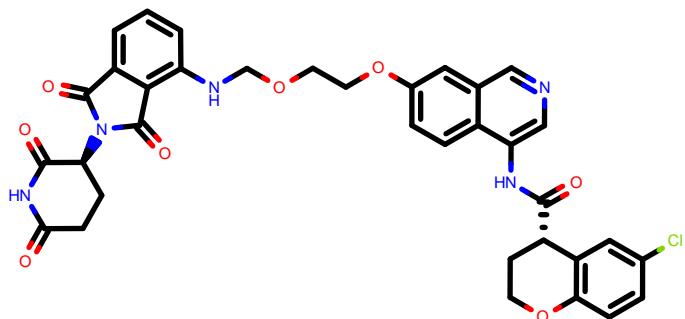
CID:	JOH-UNI-a38a7bdd-6_4
SMILES:	<chem>c1ccc2c(c1)cncc2N(C(=O)Cc3cccc(c3)Cl)C(=O)[C@H]4C[C@H]4C(F)(F)F</chem>
RUN:	RUN1494
DDG (kcal/mol):	-0.75
dDDG (kcal/mol):	0.31

MAK-UNK-c749d764-3_7



CID:	MAK-UNK-c749d764-3_7
SMILES:	<chem>CC1(C[NH2+]C1)OCN(c2ncc3c2cccc3)C(=O)[C@@H]4CCC[C@H]4(C)[F]F</chem>
RUN:	RUN911
DDG (kcal/mol):	-0.74
dDDG (kcal/mol):	0.27

EDJ-MED-e4b030d8-5_1



CID: EDJ-MED-e4b030d8-5_1

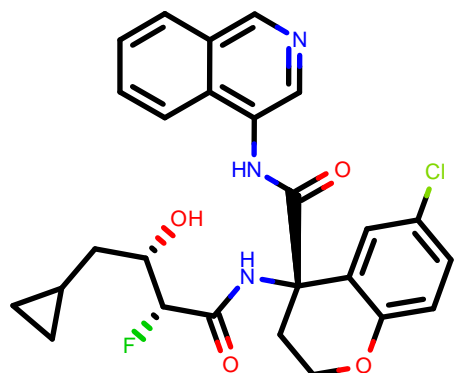
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCOC4c3cc(cc4C5COC5)Cl

RUN: RUN286

DDG (kcal/mol): -0.74

dDDG (kcal/mol): 0.37

DAR-DIA-b4e9dd8d-5_2



CID: DAR-DIA-b4e9dd8d-5_2

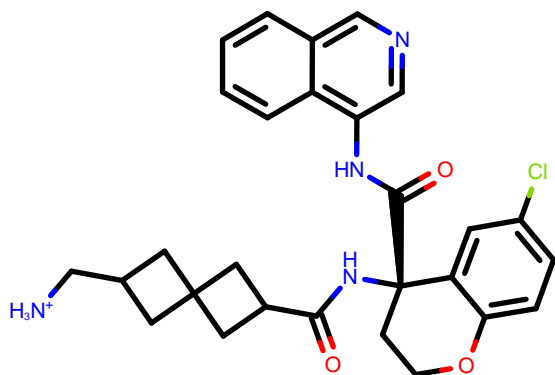
SMILES: COC(=O)/C=C/C(=O)N(c1cnc2c1ccc2)C(=O)[C@@H]3CNc4c3cc(cc4)Cl

RUN: RUN3400

DDG (kcal/mol): -0.73

dDDG (kcal/mol): 0.15

MIC-UNK-8758c41d-1_1



CID: MIC-UNK-8758c41d-1_1

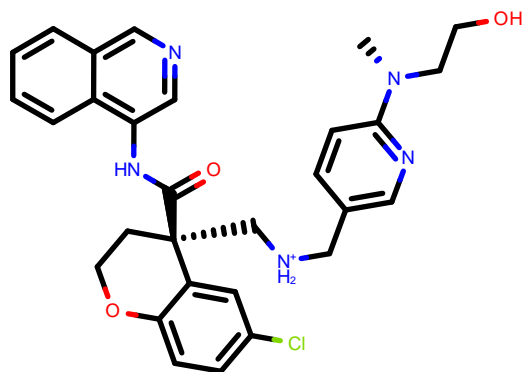
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCS(=O)(=O)c4c3cc(c(c4)Cl)Cl

RUN: RUN3312

DDG (kcal/mol): -0.73

dDDG (kcal/mol): 0.18

EDJ-MED-b7309adf-2_1



CID: EDJ-MED-b7309adf-2_1

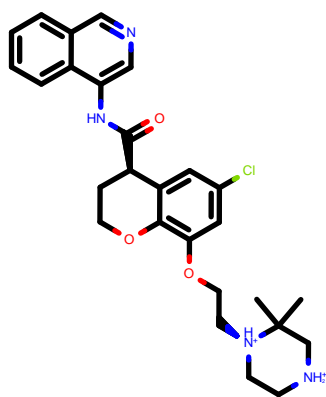
SMILES: c1cc2c(cc1Cl)cnc2NC(=O)[C@@H]3CCS(=O)(=O)c4c3cc(cc4)Cl

RUN: RUN4526

DDG (kcal/mol): -0.73

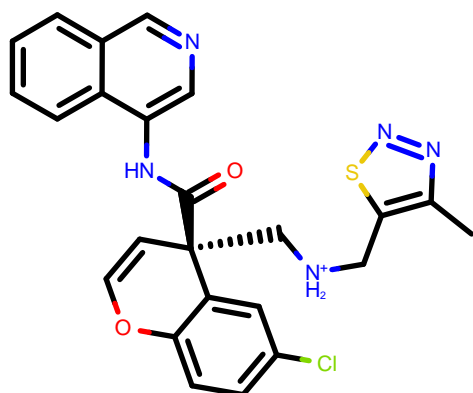
dDDG (kcal/mol): 0.45

MAT-POS-de59a476-4_1



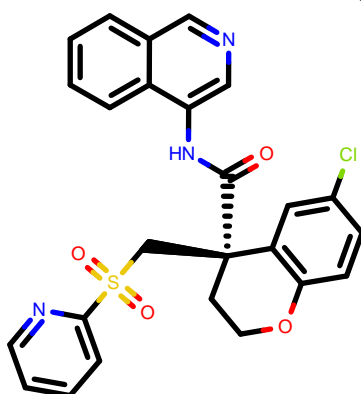
CID:	MAT-POS-de59a476-4_1
SMILES:	<chem>COCCO[C@@H](c1ccc(c(c1)Cl)Cl)C(=O)Nc2cnc3c2cccc3</chem>
RUN:	RUN2222
DDG (kcal/mol):	-0.73
dDDG (kcal/mol):	0.48

EDG-MED-90036822-17_1



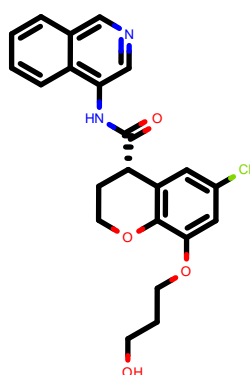
CID:	EDG-MED-90036822-17_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3[C@@H](COc4c3cc(cc4)Cl)NC(=O)[C@@H](CF)O</chem>
RUN:	RUN1680
DDG (kcal/mol):	-0.73
dDDG (kcal/mol):	0.42

MIC-UNK-08fa0751-1_2



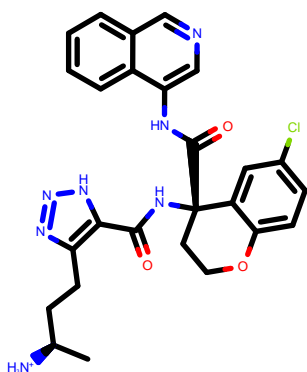
CID:	MIC-UNK-08fa0751-1_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3[C@@H](CCOc4c3cc(cc4C)O)(F)F)Cl</chem>
RUN:	RUN1520
DDG (kcal/mol):	-0.72
dDDG (kcal/mol):	0.19

VLA-UNK-f702bf1c-2_1



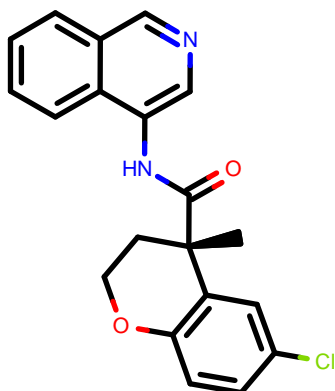
CID:	VLA-UNK-f702bf1c-2_1
SMILES:	<chem>c1ccc2c(c1)cnc2N3C(=O)[C@@H]4(CCOc5c4cc(cc5)Cl)N(C3=O)C6ccn[nH]6</chem>
RUN:	RUN2306
DDG (kcal/mol):	-0.72
dDDG (kcal/mol):	0.31

PET-UNK-55f647aa-2_1



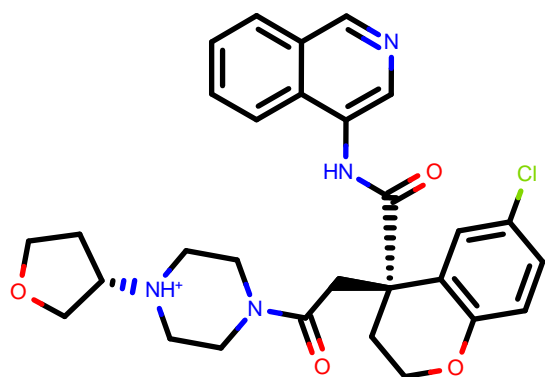
CID:	PET-UNK-55f647aa-2_1
SMILES:	<chem>CN(C)c1ccc(en1)N(Cc2ccsc2)C(=O)Cc3cnc4c3cccc4</chem>
RUN:	RUN570
DDG (kcal/mol):	-0.72
dDDG (kcal/mol):	0.33

NAU-LAT-356bd3c2-7_1



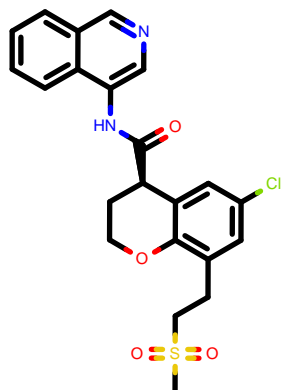
CID:	NAU-LAT-356bd3c2-7_1
SMILES:	<chem>c1ccc2c(c1)cnc2OC(=O)c3cccc4c3cc[nH]4</chem>
RUN:	RUN58
DDG (kcal/mol):	-0.72
dDDG (kcal/mol):	0.13

BEN-DND-c852c98b-2_1



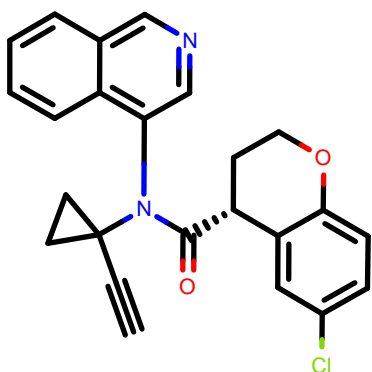
CID:	BEN-DND-c852c98b-2_1
SMILES:	<chem>c1cc2nccc(e2cc1OC(F)(F)F)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN1204
DDG (kcal/mol):	-0.72
dDDG (kcal/mol):	0.28

MAT-POS-afd4d4fd-1_1



CID:	MAT-POS-afd4d4fd-1_1
SMILES:	<chem>Cc1c(cccc1Cl)CC(=O)Nc2cnc3c2cccc3</chem>
RUN:	RUN303
DDG (kcal/mol):	-0.72
dDDG (kcal/mol):	0.33

ALP-UNI-8e43a71e-2_12



CID: ALP-UNI-8e43a71e-2_12

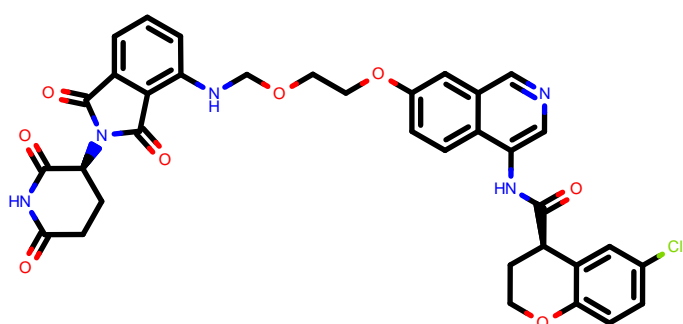
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@H]3[C@@H](COCc4ccc(Cl)cc4)CC(=O)N5CC[C@H]6[C@@H](C#C)N#H5

RUN: RUN2933

DDG (kcal/mol): -0.71

dDDG (kcal/mol): 0.12

ROB-UNI-322e8f70-2_2



CID: ROB-UNI-322e8f70-2_2

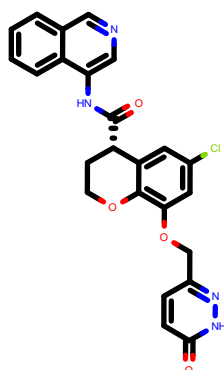
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@H]3CS(=O)(=O)Nc4c3cc(cc4)Cl

RUN: RUN3156

DDG (kcal/mol): -0.71

dDDG (kcal/mol): 0.28

EDJ-MED-40433386-4_1



CID: EDJ-MED-40433386-4_1

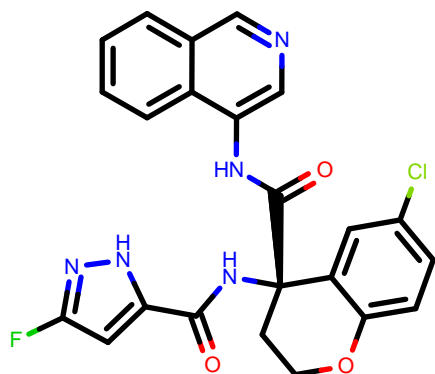
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@H]3[C@@H](COCc4ccc(Cl)cc4)CNC(=O)C5(CCC5)CO

RUN: RUN2556

DDG (kcal/mol): -0.70

dDDG (kcal/mol): 0.24

EDJ-MED-f893e2a1-10_1



CID: EDJ-MED-f893e2a1-10_1

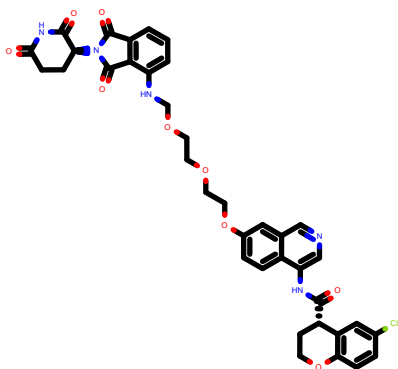
SMILES: CC(C)(CNC[C@@H]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3ccc4c3ccc4)C#N

RUN: RUN3214

DDG (kcal/mol): -0.70

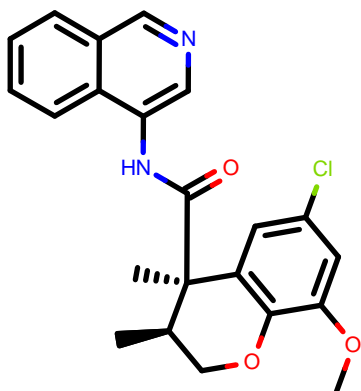
dDDG (kcal/mol): 0.11

MAT-POS-8a69d52e-3_1



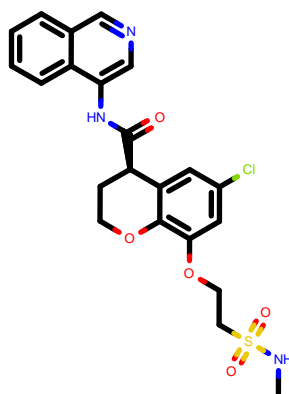
CID:	MAT-POS-8a69d52e-3_1
SMILES:	<chem>C[C@H]1C[C@@H](c2cc(ccc2O1)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN365
DDG (kcal/mol):	-0.70
dDDG (kcal/mol):	0.40

LON-WEI-4d77710c-7_2



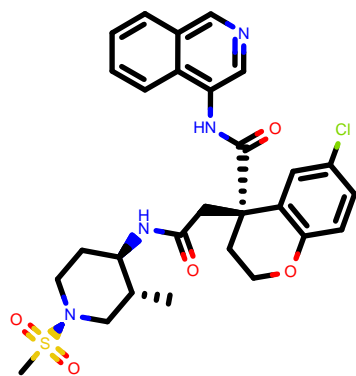
CID:	LON-WEI-4d77710c-7_2
SMILES:	<chem>Cn1cc(c2cccc2c1=O)NC(=O)NCCC[N@H+]3CCc4cccc4C3</chem>
RUN:	RUN198
DDG (kcal/mol):	-0.70
dDDG (kcal/mol):	0.22

KAD-UNI-8a629cb0-13_1



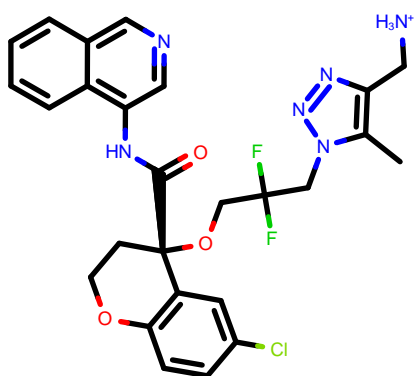
CID:	KAD-UNI-8a629cb0-13_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(Cc4cc3cc(cc4)Cl)NC(=O)CN5CCN(C5=O)O6CC6</chem>
RUN:	RUN2095
DDG (kcal/mol):	-0.69
dDDG (kcal/mol):	0.34

MAT-POS-4223bc15-14_2



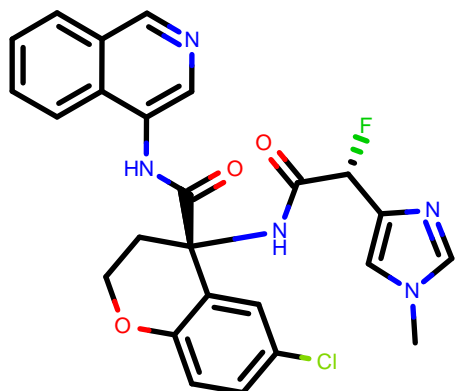
CID:	MAT-POS-4223bc15-14_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3[C@H](N@)4(Cc4c3cc(cc4)Cl)S(=O)(=O)CC5C(NH2+)C5</chem>
RUN:	RUN4067
DDG (kcal/mol):	-0.69
dDDG (kcal/mol):	0.34

ERI-UCB-b3e6b0c2-12_1



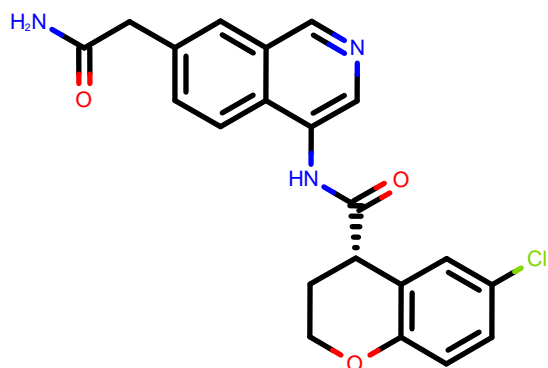
CID:	ERI-UCB-b3e6b0c2-12_1
SMILES:	<chem>C[N+]([O-])C1C[C@H](c2c1ccc(c2)Cl)C(=O)Nc3ncnc4c3ccc(c4)C[NH3+]</chem>
RUN:	RUN3046
DDG (kcal/mol):	-0.69
dDDG (kcal/mol):	0.21

EDJ-MED-4f704dc9-2_2



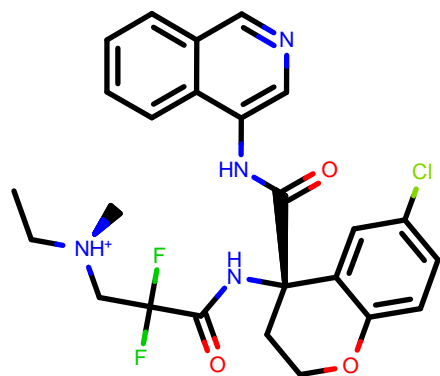
CID:	EDJ-MED-4f704dc9-2_2
SMILES:	<chem>COC[C@]1(CC(=O)Nc2c1cc(c(c2)Cl)C)C(=O)Nc3ncnc4c3ccc4</chem>
RUN:	RUN3162
DDG (kcal/mol):	-0.69
dDDG (kcal/mol):	0.14

MAT-POS-b5746674-37_2



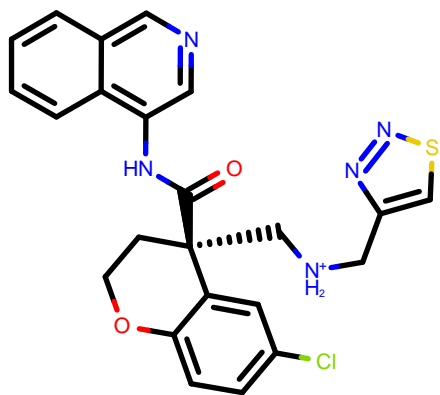
CID:	MAT-POS-b5746674-37_2
SMILES:	<chem>Cn1cc(c2ccccc2c1=O)NC(=O)N(Cc3ccco3)C[C@H]4CCCCO4</chem>
RUN:	RUN67
DDG (kcal/mol):	-0.69
dDDG (kcal/mol):	0.13

DAR-DIA-0d514e7d-33_3



CID:	DAR-DIA-0d514e7d-33_3
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)C@H]3c4cc(ccc4O)C@@H]5[C@H]3C5)Cl</chem>
RUN:	RUN884
DDG (kcal/mol):	-0.69
dDDG (kcal/mol):	0.34

ALP-POS-347519b5-1_3



CID: ALP-POS-347519b5-1_3

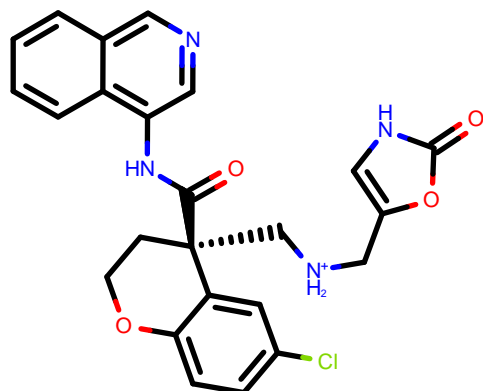
SMILES: CS(-O)(N)C1C2C@H3CC(C@H)C3(C@H)2C@H(C1)C(-O)N4ncc5c4ccc5

RUN: RUN4233

DDG (kcal/mol): -0.69

dDDG (kcal/mol): 0.20

ALP-POS-347519b5-3_24



CID: ALP-POS-347519b5-3_24

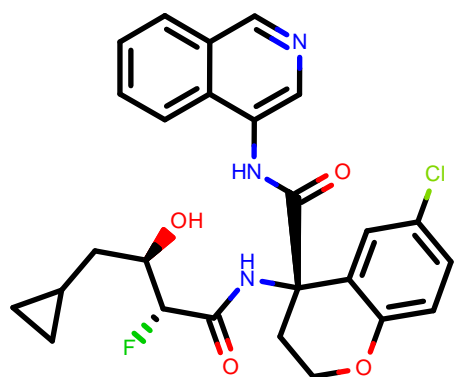
SMILES: CS(-O)(N)C1C2C@H3CC(C@H)C3(C@H)2C@H(C1)C(-O)N4ncc5c4ccc5O

RUN: RUN4307

DDG (kcal/mol): -0.68

dDDG (kcal/mol): 0.32

MIC-UNK-0a05c952-1_8



CID: MIC-UNK-0a05c952-1_8

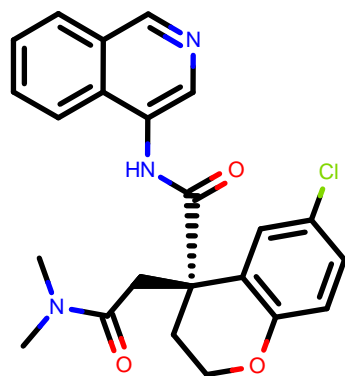
SMILES: c1ccc2c(c1)ncc2N3[C@H](C[C@H](C3=O)c4cccc(c4)Cl)[C@H]5CO5

RUN: RUN3504

DDG (kcal/mol): -0.68

dDDG (kcal/mol): 0.18

MAK-UNK-ffc90da7-6_1



CID: MAK-UNK-ffc90da7-6_1

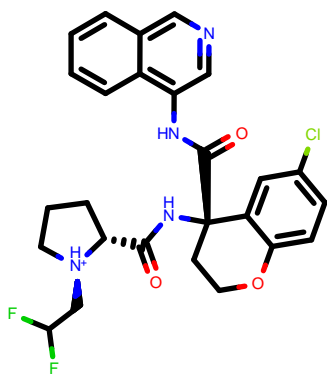
SMILES: C[C@H](C[NH2+])CCCOc1ccc2nccc(c2c1)NC(=O)Cc3ccccc(c3)Cl

RUN: RUN703

DDG (kcal/mol): -0.68

dDDG (kcal/mol): 0.18

MIC-UNK-5a93dd5f-9_4



CID: MIC-UNK-5a93dd5f-9_4

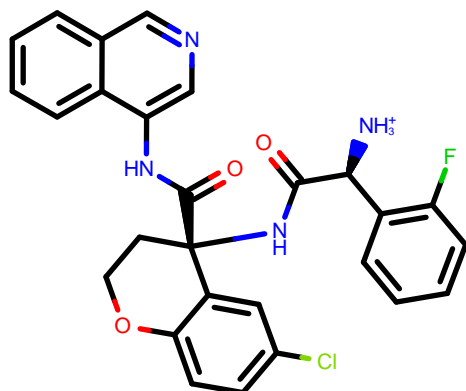
SMILES: CN(C)[C@H]1CC[N@H+]1C1[C@@H](c2cccc(c2)Cl)C(=O)Nc3ncc4c3cccc4

RUN: RUN781

DDG (kcal/mol): -0.68

dDDG (kcal/mol): 0.28

ADA-UCB-dc2b944c-1_2



CID: ADA-UCB-dc2b944c-1_2

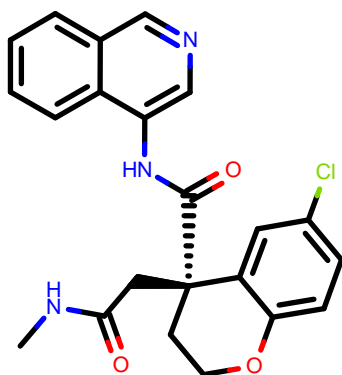
SMILES: c1ccc2c(c1)cncc2N3C(=O)C[N@H+]1[C@@H](C3=O)CCOc5c4cc(cc5)C1)CC6CCCC6

RUN: RUN598

DDG (kcal/mol): -0.67

dDDG (kcal/mol): 0.31

DAR-DIA-0d514e7d-32_1



CID: DAR-DIA-0d514e7d-32_1

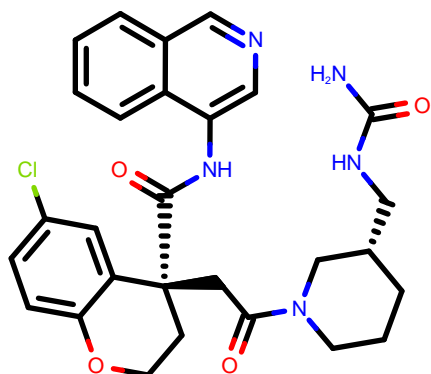
SMILES: c1ccc2c(c1)cncc2N3C(=O)[C@@H](C3=O)C@H4C(C3=O)CCOc5c4cc(cc5)C1)CC6CCCC6

RUN: RUN850

DDG (kcal/mol): -0.67

dDDG (kcal/mol): 0.16

NAU-LAT-2fed8305-6_1



CID: NAU-LAT-2fed8305-6_1

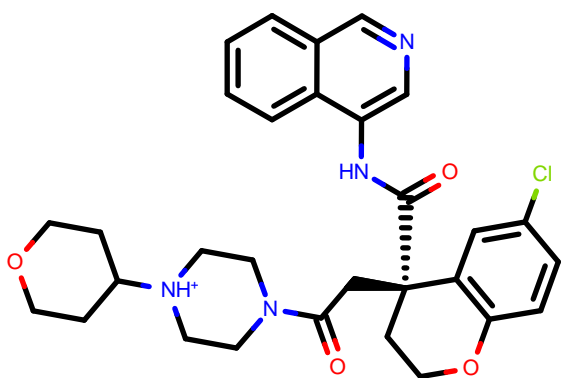
SMILES: c1ccc2c(c1)cncc2NC(=O)Cc3cccc(s3)Cl

RUN: RUN1113

DDG (kcal/mol): -0.67

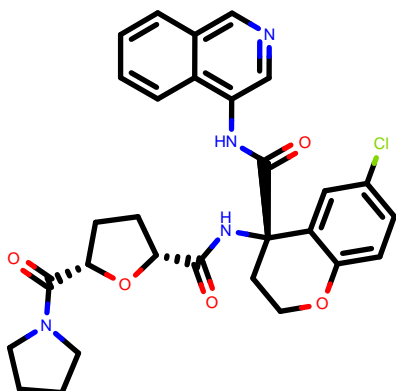
dDDG (kcal/mol): 0.41

ALP-POS-ce760d3f-1_2



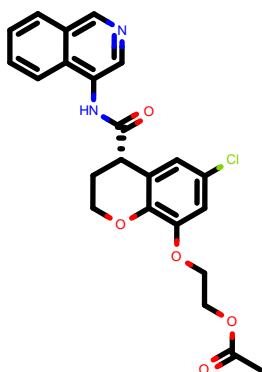
CID:	ALP-POS-ce760d3f-1_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCOCc4c3cc(cc4O)Cl</chem>
RUN:	RUN1462
DDG (kcal/mol):	-0.66
dDDG (kcal/mol):	0.33

ALP-POS-966f8da6-1_1



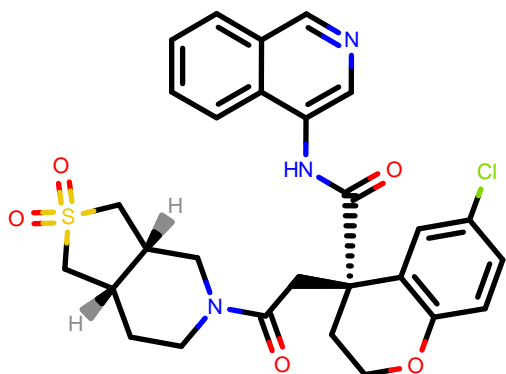
CID:	ALP-POS-966f8da6-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCN(c4c3cc(cc4)Cl)Cc5cn[nH]5</chem>
RUN:	RUN1218
DDG (kcal/mol):	-0.66
dDDG (kcal/mol):	0.33

EDG-MED-ba1ac7b9-1_4



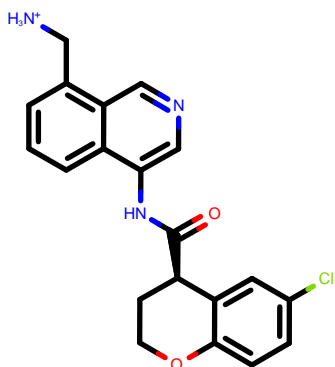
CID:	EDG-MED-ba1ac7b9-1_4
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CC[C@H]5C(=O)N</chem>
RUN:	RUN2618
DDG (kcal/mol):	-0.66
dDDG (kcal/mol):	0.24

DAR-DIA-0d514e7d-12_1



CID:	DAR-DIA-0d514e7d-12_1
SMILES:	<chem>C[C@H]1COc2c(cc(cc2[C@@H]1C(=O)Nc3ncc4c3ccc4)Cl)c5ccc(cc5)S(=O)(=O)F</chem>
RUN:	RUN816
DDG (kcal/mol):	-0.66
dDDG (kcal/mol):	0.28

DAR-DIA-0cde14eb-57_1



CID: DAR-DIA-0cde14eb-57_1

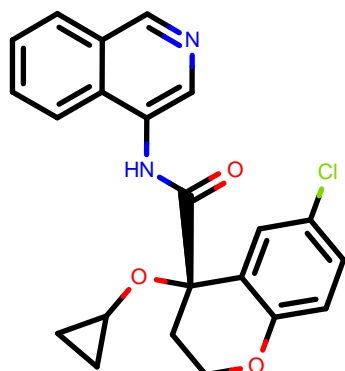
SMILES: C[C@@H](c1cccc(c1)C2(CC2)F)C(=O)Nc3cncc4c3cccc4

RUN: RUN25

DDG (kcal/mol): -0.66

dDDG (kcal/mol): 0.23

MAT-POS-4223bc15-31_3



CID: MAT-POS-4223bc15-31_3

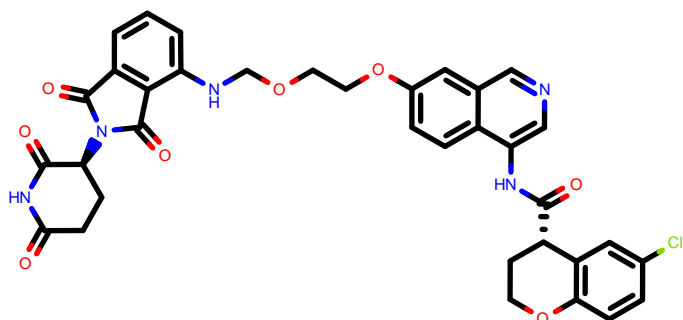
SMILES: Cn1ncc(n1)S(=O)(=O)N@@2Cc3ccc(cc3[C@@H](C2)C(=O)Nc4cncc5c4cccc5)Cl

RUN: RUN4140

DDG (kcal/mol): -0.65

dDDG (kcal/mol): 0.08

EDJ-MED-e4b030d8-9_1



CID: EDJ-MED-e4b030d8-9_1

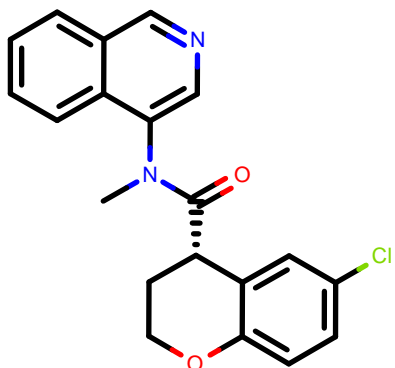
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4c3nc(cc4)Cl

RUN: RUN290

DDG (kcal/mol): -0.65

dDDG (kcal/mol): 0.35

LON-WEI-4d77710c-35_1



CID: LON-WEI-4d77710c-35_1

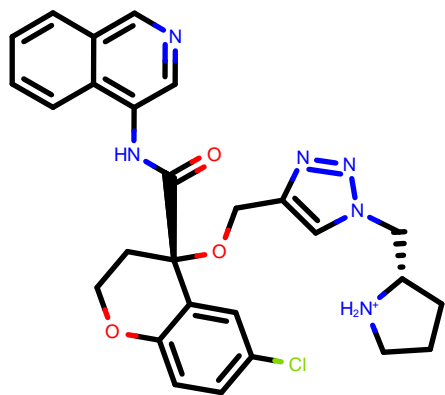
SMILES: Cc1ccc(cc1)C[N@@H+]2CC[C@@H](C2)CNC(=O)Nc3cn(cc3=O)c4c3cccc4)CC(C)C

RUN: RUN222

DDG (kcal/mol): -0.65

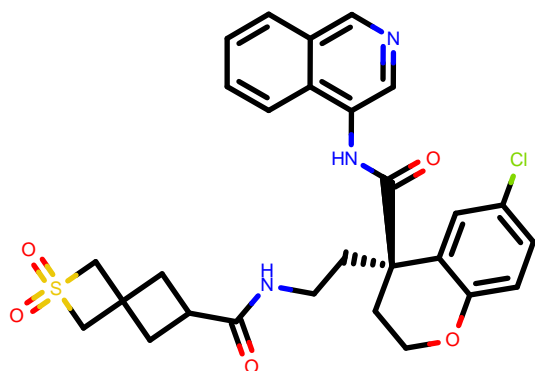
dDDG (kcal/mol): 0.22

ALP-POS-fe871b40-12_2



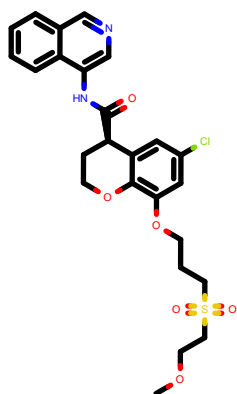
CID:	ALP-POS-fe871b40-12_2
SMILES:	<chem>CO[C@J]1(CCNC2c1cc(c(c2)F)Cl)C(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN3128
DDG (kcal/mol):	-0.65
dDDG (kcal/mol):	0.20

DAR-DIA-f6ee7aeb-3_3



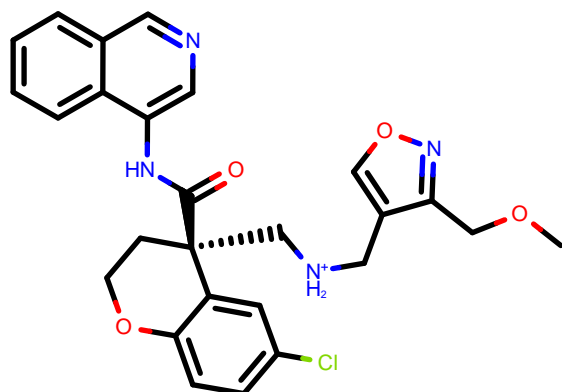
CID:	DAR-DIA-f6ee7aeb-3_3
SMILES:	<chem>CCCOc1cc(cc(c1)Cl)C@@H]2CC(=O)N(C)C@H]2c3cccc3C#N)c4cnc5c4cccc5</chem>
RUN:	RUN3409
DDG (kcal/mol):	-0.64
dDDG (kcal/mol):	0.20

EDG-MED-ba1ac7b9-8_2



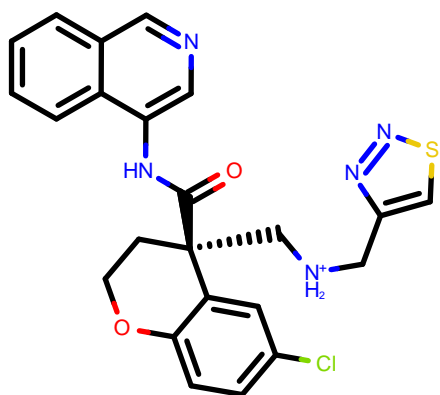
CID:	EDG-MED-ba1ac7b9-8_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)C@]3(CCOc4c3cc(cc4)Cl)CC(=O)N(CC#N)C5O5</chem>
RUN:	RUN2645
DDG (kcal/mol):	-0.64
dDDG (kcal/mol):	0.51

MIC-UNK-460e637d-2_6



CID:	MIC-UNK-460e637d-2_6
SMILES:	<chem>C[N@H+]1CCN(C)C@@H]1CO[C@H]2CC(=O)N2)C(=O)Cc3cnc4c3cccc4</chem>
RUN:	RUN1887
DDG (kcal/mol):	-0.64
dDDG (kcal/mol):	0.52

ED_-GRI-5b13fbe2-24_1



CID: ED_-GRI-5b13fbe2-24_1

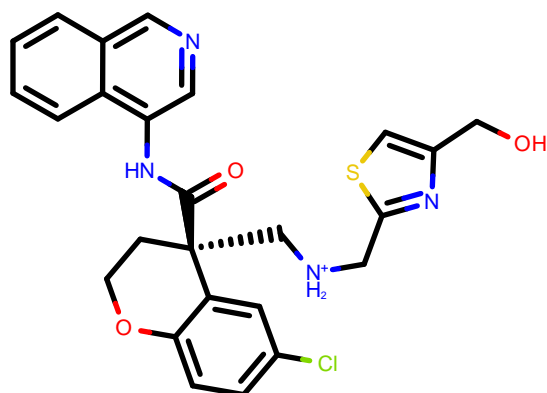
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3(CCOc4c3cc(c4)Cl)OCc5cn(n5)CC6C[NH2+]C6

RUN: RUN1555

DDG (kcal/mol): -0.64

dDDG (kcal/mol): 0.36

EDJ-MED-93390d0c-4_1



CID: EDJ-MED-93390d0c-4_1

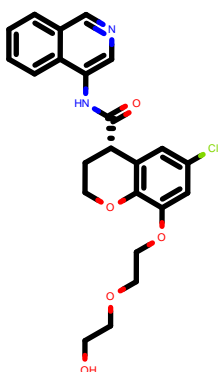
SMILES: CO[C@@]1(CCOc2c1cc(c2)F)Cl)C(=O)Nc3ncc4c3cc(c4)S(=O)(=O)C

RUN: RUN4525

DDG (kcal/mol): -0.64

dDDG (kcal/mol): 0.25

ALP-POS-2da19ca7-8_1



CID: ALP-POS-2da19ca7-8_1

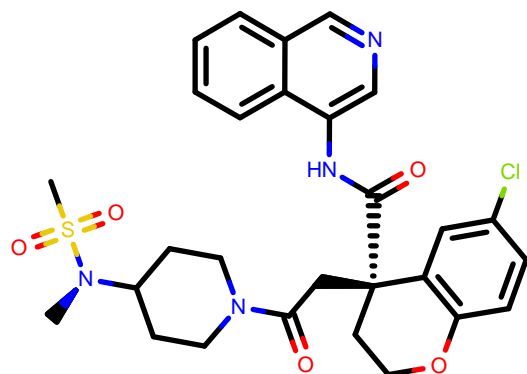
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3(CCOc4c3cc(c4)Cl)CC1=O)N[C@@H](CCO)N(C)COC(C)C

RUN: RUN2394

DDG (kcal/mol): -0.64

dDDG (kcal/mol): 0.28

LON-WEI-5e7d1b3e-21_1



CID: LON-WEI-5e7d1b3e-21_1

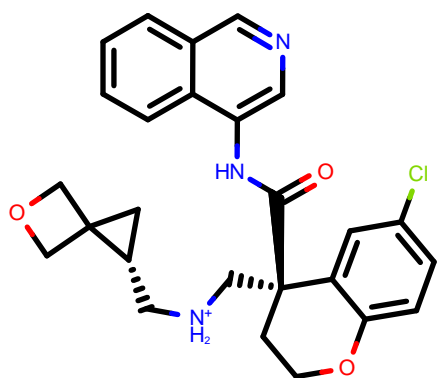
SMILES: Cc1cccnc1NC(=O)Nc2cn(c(=O)c3c2cccc3)CC(C)C

RUN: RUN1326

DDG (kcal/mol): -0.63

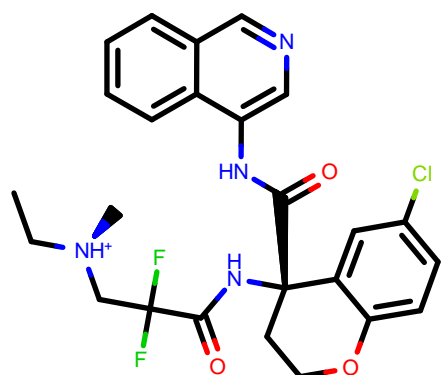
dDDG (kcal/mol): 0.67

RAL-THA-05e671eb-20_1



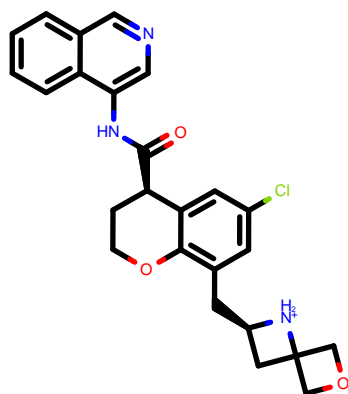
CID:	RAL-THA-05e671eb-20_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](C)CCOCc4c3cc(cc4C#N)Cl</chem>
RUN:	RUN2044
DDG (kcal/mol):	-0.63
dDDG (kcal/mol):	0.33

MAK-UNK-ffc90da7-4_5



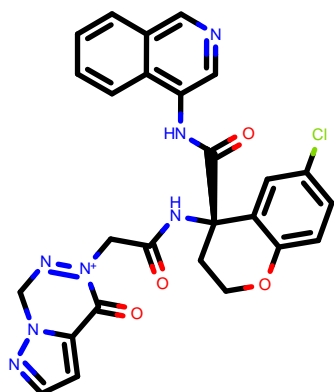
CID:	MAK-UNK-ffc90da7-4_5
SMILES:	<chem>C[C@@H](C)[NH2+][C]C@@H1CCCCO1[C@H](c2ccc(c2)Cl)C(=O)Nc3nccc4c3ccc4</chem>
RUN:	RUN700
DDG (kcal/mol):	-0.63
dDDG (kcal/mol):	0.29

DAR-DIA-0cde14eb-60_2



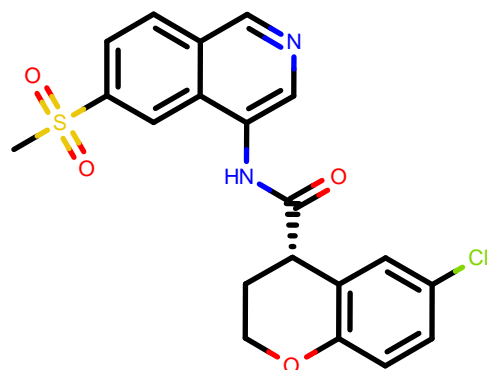
CID:	DAR-DIA-0cde14eb-60_2
SMILES:	<chem>C[C@H](c1cccc(c1)C2(CC2)C#N)C(=O)Nc3nccc4c3cccc4</chem>
RUN:	RUN32
DDG (kcal/mol):	-0.63
dDDG (kcal/mol):	0.38

RAL-THA-6e4c80cf-3_1



CID:	RAL-THA-6e4c80cf-3_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](C)CN(Cc4c3cc(cc4)Cl)C(=O)OCc5ncs5</chem>
RUN:	RUN3898
DDG (kcal/mol):	-0.62
dDDG (kcal/mol):	0.18

MAT-POS-b3e365b9-2_1



CID: MAT-POS-b3e365b9-2_1

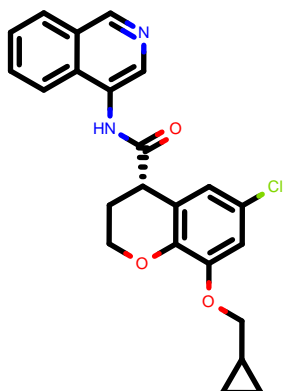
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3CCOCc4c3cc(cc4)Cl

RUN: RUN277

DDG (kcal/mol): -0.62

dDDG (kcal/mol): 0.12

DAR-DIA-9e4459de-15_4



CID: DAR-DIA-9e4459de-15_4

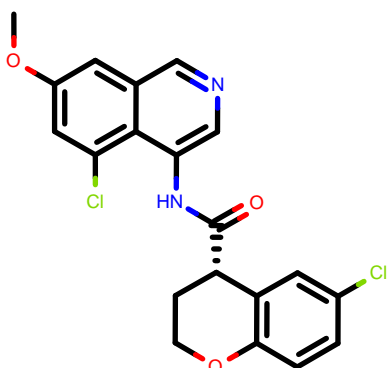
SMILES: c1cc2c(c1)NC(=O)COC1=O)Nc3ccc4c(c3)cncc4NC(=O)[C@H]5COCc6cc(cc6)Cl)c7c(c2)O[C@H]7COC1=O)NC7=O)O

RUN: RUN1446

DDG (kcal/mol): -0.62

dDDG (kcal/mol): 0.24

RAL-THA-4aa06b95-7_3



CID: RAL-THA-4aa06b95-7_3

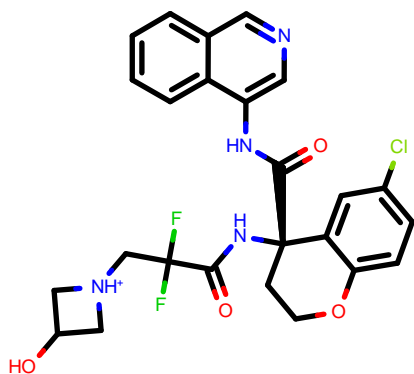
SMILES: COCCN1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN1245

DDG (kcal/mol): -0.61

dDDG (kcal/mol): 0.13

DAR-DIA-2964957d-9_1



CID: DAR-DIA-2964957d-9_1

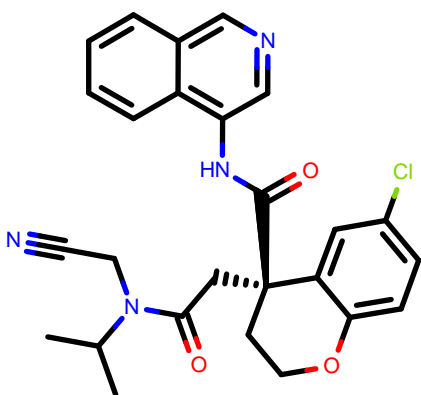
SMILES: CN(c1cccc(c1)Cl)c2c(c(=O)c2=O)Nc3cncc4c3cccc4

RUN: RUN518

DDG (kcal/mol): -0.61

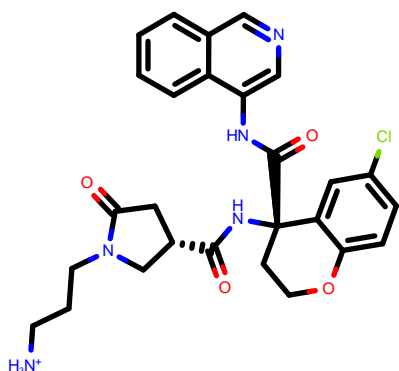
dDDG (kcal/mol): 0.29

ERI-UCB-d6de1f3c-7_1



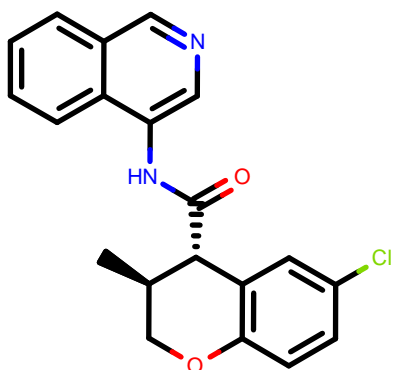
CID:	ERI-UCB-d6de1f3c-7_1
SMILES:	<chem>C[C@@H]1CN(CC(=O)N1c2cccc(c2)Cl)C(=O)c3cncc4c3cccc4</chem>
RUN:	RUN1099
DDG (kcal/mol):	-0.61
dDDG (kcal/mol):	0.36

ALP-POS-64a710fa-2_1



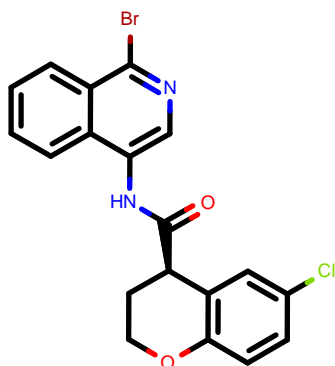
CID:	ALP-POS-64a710fa-2_1
SMILES:	<chem>C[NH+](C)CCCN(Cc1cccc1)C(=O)C2cnc3c2cccc3</chem>
RUN:	RUN465
DDG (kcal/mol):	-0.61
dDDG (kcal/mol):	0.35

MAT-POS-b5746674-34_1



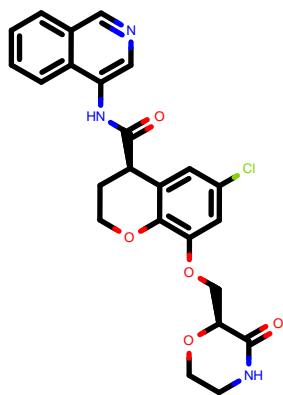
CID:	MAT-POS-b5746674-34_1
SMILES:	<chem>CCCC[N@@H+](CCNC(=O)Nc1cn(c(=O)c2c1cccc2)C)Cc3cccc3</chem>
RUN:	RUN66
DDG (kcal/mol):	-0.61
dDDG (kcal/mol):	0.12

MAT-POS-b5746674-109_1



CID:	MAT-POS-b5746674-109_1
SMILES:	<chem>CC(C)Cn1cc(c2cccc2c1=O)NC(=O)NCC[NH+]3CCOCC3</chem>
RUN:	RUN92
DDG (kcal/mol):	-0.60
dDDG (kcal/mol):	0.23

ALP-POS-5bb456a5-2_8



CID: ALP-POS-5bb456a5-2_8

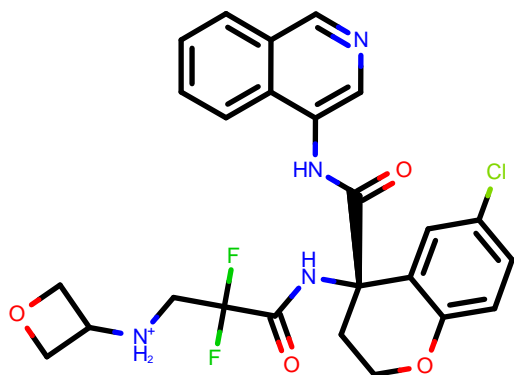
SMILES: C1C@H1C[N@]([C@@H]1NC(=O)C[C@@]2(COC3C2cc3C)C)C(=O)N4ncc5c4cccc5[S(=O)2](=O)C

RUN: RUN2429

DDG (kcal/mol): -0.60

dDDG (kcal/mol): 0.42

PET-UNK-55f647aa-1_1



CID: PET-UNK-55f647aa-1_1

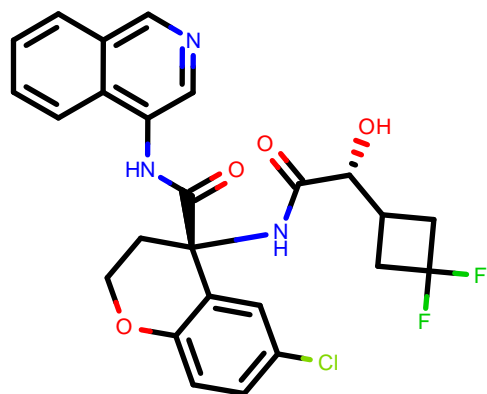
SMILES: CN(C)c1ccc(nc1)N(Cc2ccsc2)C(=O)Cc3cnc4c3cccc4

RUN: RUN569

DDG (kcal/mol): -0.60

dDDG (kcal/mol): 0.42

EDJ-MED-6d9ff7d0-8_1



CID: EDJ-MED-6d9ff7d0-8_1

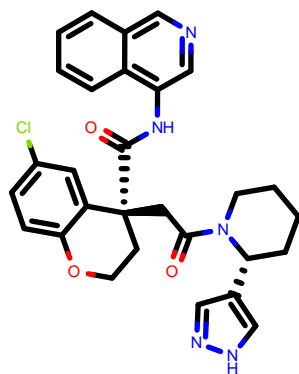
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4C)O)N[C@@H]3C[C@@]4H5CCOCC(=O)N5

RUN: RUN3433

DDG (kcal/mol): -0.60

dDDG (kcal/mol): 0.15

EDJ-MED-670ad2ee-6_2



CID: EDJ-MED-670ad2ee-6_2

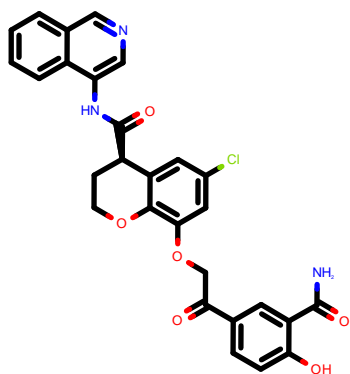
SMILES: c1cc2ncc(c2cc1F)NC(=O)[C@@]3(C)C[N@]([C@@H]3C)C(=O)S(=O)(=O)C5CC5

RUN: RUN3870

DDG (kcal/mol): -0.60

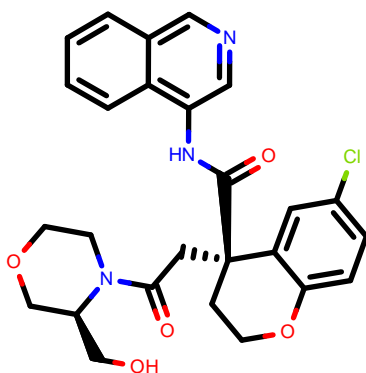
dDDG (kcal/mol): 0.22

EDG-MED-ba1ac7b9-12_2



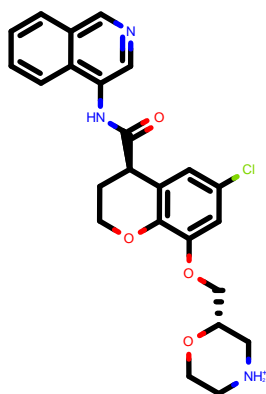
CID:	EDG-MED-ba1ac7b9-12_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3[C@@O]c4c3cc(cc4)C)CC(=O)N5CCC[C@H]5c6r(n-)nn6</chem>
RUN:	RUN2659
DDG (kcal/mol):	-0.60
dDDG (kcal/mol):	0.43

KAD-UNI-877d7bed-11_2



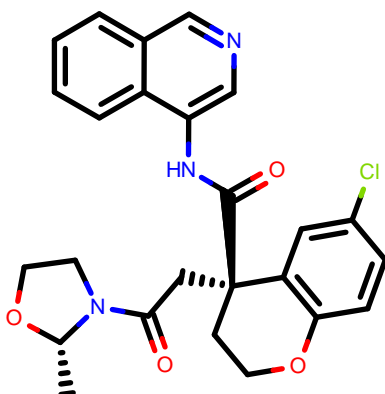
CID:	KAD-UNI-877d7bed-11_2
SMILES:	<chem>CC(C)[C@H]1CCN(C1)C(=O)COc2cc(cc3c2OCC[C@H]3C(=O)Nc4cncc5c4ccc5)C)O</chem>
RUN:	RUN3748
DDG (kcal/mol):	-0.59
dDDG (kcal/mol):	0.21

MAT-POS-24589f88-12_2



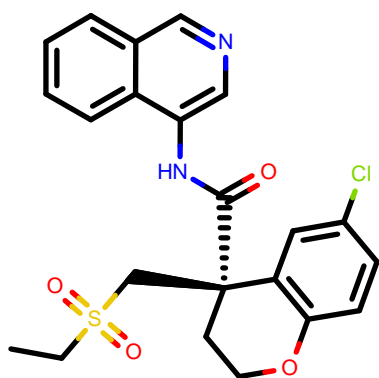
CID:	MAT-POS-24589f88-12_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3[C@@O]c4c3cc(cc4)C)C[NH2+]Cc5ccc(cc5)S(=O)(=O)N</chem>
RUN:	RUN4799
DDG (kcal/mol):	-0.59
dDDG (kcal/mol):	0.21

MAT-POS-8d5af1ef-1_2



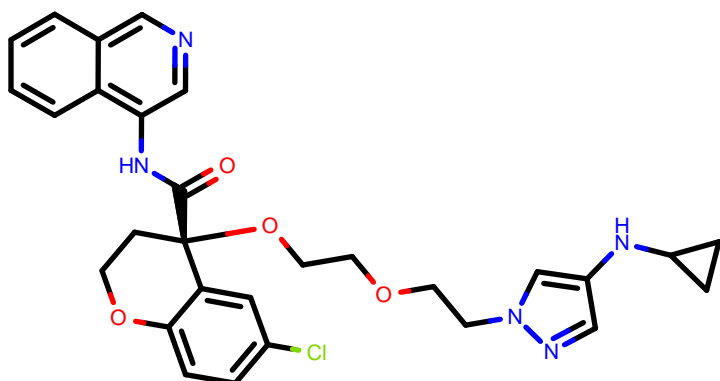
CID:	MAT-POS-8d5af1ef-1_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CC(=O)Nc4c3cc(cc4)Br</chem>
RUN:	RUN1129
DDG (kcal/mol):	-0.58
dDDG (kcal/mol):	0.26

LAU-MED-88a3970a-7_1



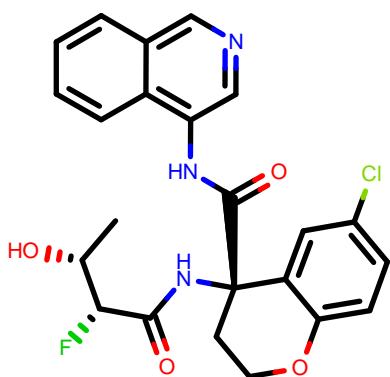
CID:	LAU-MED-88a3970a-7_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOc4c3cc(cc4CCCCO)Cl</chem>
RUN:	RUN1502
DDG (kcal/mol):	-0.58
dDDG (kcal/mol):	0.16

ALP-POS-fe871b40-7_1



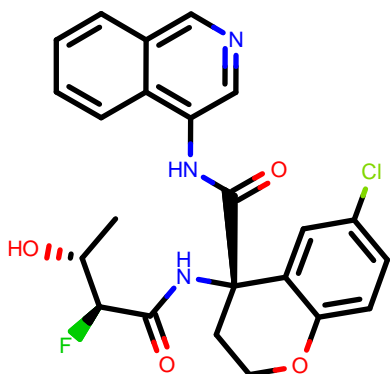
CID:	ALP-POS-fe871b40-7_1
SMILES:	<chem>CO[C@@]1(CCOC2c1cc(cn2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3119
DDG (kcal/mol):	-0.58
dDDG (kcal/mol):	0.27

MAT-POS-fb82b63d-3_3



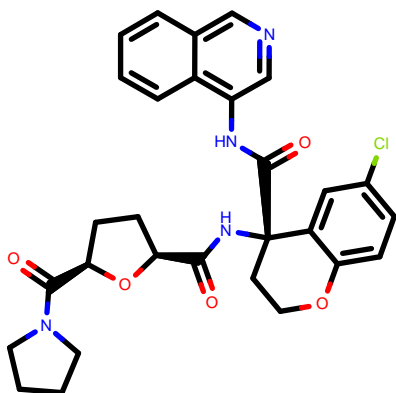
CID:	MAT-POS-fb82b63d-3_3
SMILES:	<chem>C[N@@H+]1CCc2ccc(cc2[C@H]1C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN3175
DDG (kcal/mol):	-0.58
dDDG (kcal/mol):	0.12

MAT-POS-78e1d523-5_1



CID:	MAT-POS-78e1d523-5_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCNc4c3cc(c(c4)Cl)Cl</chem>
RUN:	RUN3282
DDG (kcal/mol):	-0.58
dDDG (kcal/mol):	0.13

MAT-POS-4223bc15-9_4



CID: MAT-POS-4223bc15-9_4

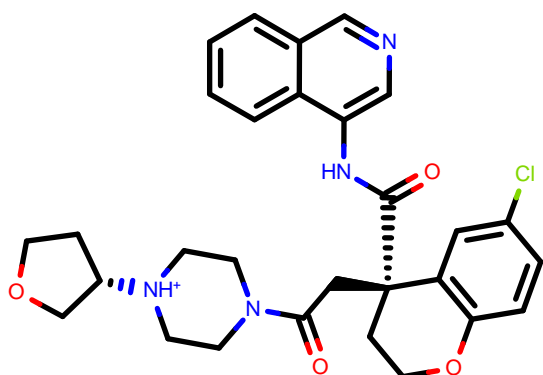
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@H]3C[N@@]3C4c3ccc(cc4)Cl5(-O)-O)C@@H5COC3C@@H5O

RUN: RUN4012

DDG (kcal/mol): -0.57

dDDG (kcal/mol): 0.16

MAK-UNK-c749d764-20_1



CID: MAK-UNK-c749d764-20_1

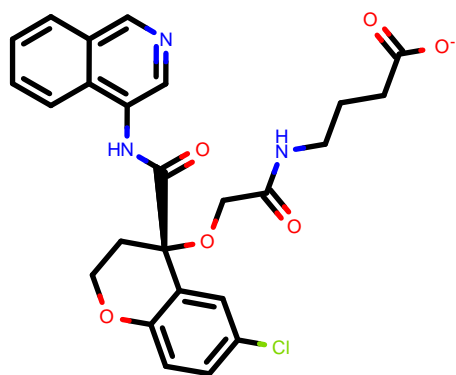
SMILES: CCCN(c1cnc2c1cccc2)C(=O)C[C@@H]3CCCC[C@@H]([C@@H]3O)C(F)F

RUN: RUN1007

DDG (kcal/mol): -0.57

dDDG (kcal/mol): 0.29

ALP-POS-fe871b40-11_2



CID: ALP-POS-fe871b40-11_2

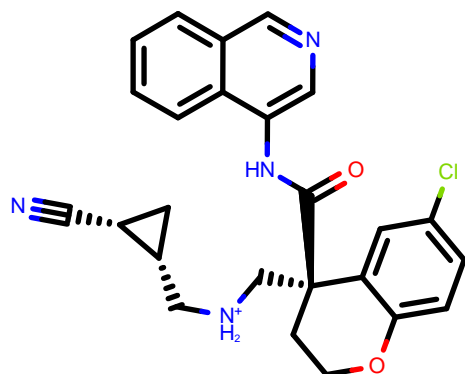
SMILES: CO[C@]1(CCOc2c1cc(c(c2)F)Cl)C(=O)Nc3nccc4c3cccc4

RUN: RUN3130

DDG (kcal/mol): -0.57

dDDG (kcal/mol): 0.19

EDG-MED-90036822-45_1



CID: EDG-MED-90036822-45_1

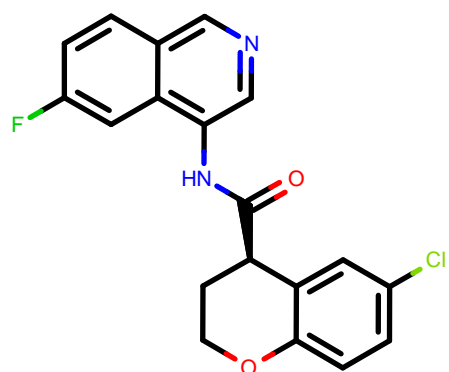
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@]3(CCOc4c3ccc(cc4)Cl)NC(=O)c5[nH]cn5(F)F

RUN: RUN1723

DDG (kcal/mol): -0.57

dDDG (kcal/mol): 0.34

DAR-DIA-0f7b7cd9-8_1



CID: DAR-DIA-0f7b7cd9-8_1

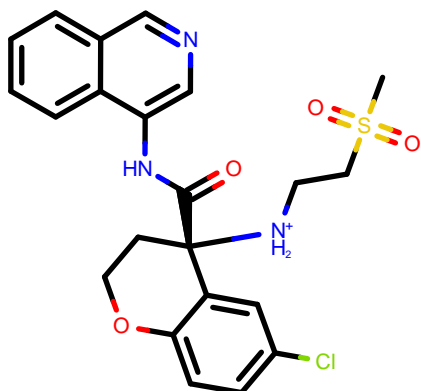
SMILES: c1ccc2c(c1)cncc2N3C(=O)[C@@]4(CCOc5c4cc(cc5)Cl)C(=O)S3

RUN: RUN3016

DDG (kcal/mol): -0.57

dDDG (kcal/mol): 0.07

EDG-MED-971238d3-4_1



CID: EDG-MED-971238d3-4_1

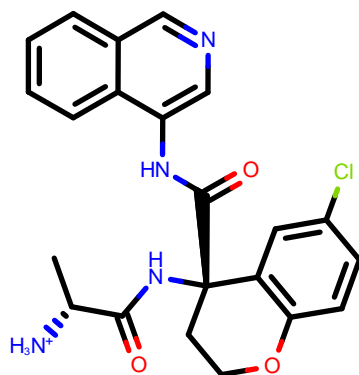
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)[NH3+]

RUN: RUN1469

DDG (kcal/mol): -0.57

dDDG (kcal/mol): 0.35

RAL-THA-1d44ff04-11_1



CID: RAL-THA-1d44ff04-11_1

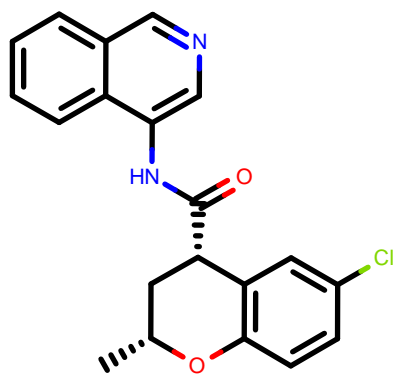
SMILES: CNS(=O)(=O)c1cc(cc(c1)Cl)CC(=O)Nc2cncc3c2cccc3

RUN: RUN446

DDG (kcal/mol): -0.57

dDDG (kcal/mol): 0.26

LON-WEI-4d77710c-42_2



CID: LON-WEI-4d77710c-42_2

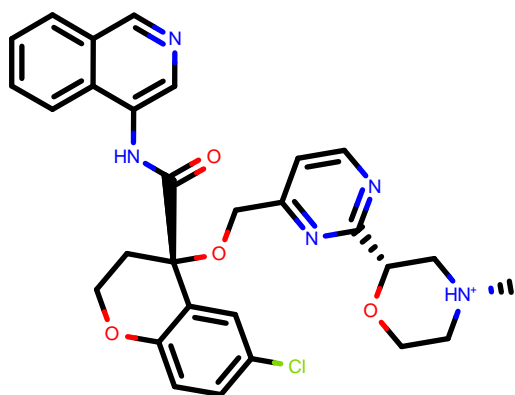
SMILES: CC(C)Cn1cc(c2cccc2c1=O)NC(=O)N[C@H]3CCCC4c3cccc4

RUN: RUN230

DDG (kcal/mol): -0.56

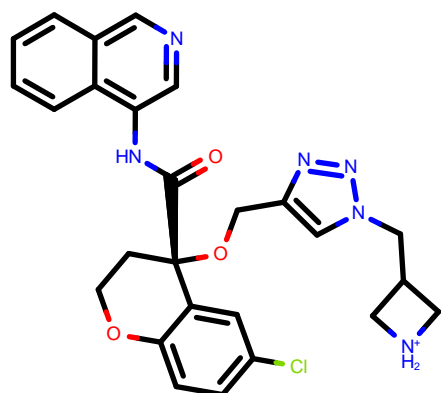
dDDG (kcal/mol): 0.12

MIC-UNK-644c43c7-1_1



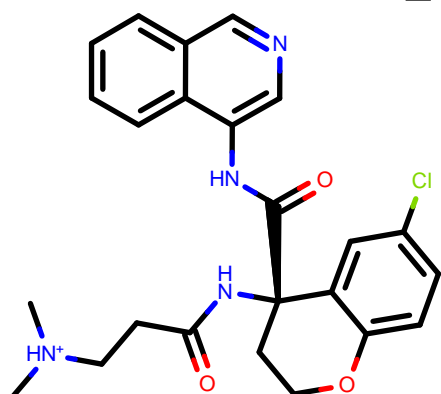
CID:	MIC-UNK-644c43c7-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)CCNCC(F)(F)F</chem>
RUN:	RUN433
DDG (kcal/mol):	-0.56
dDDG (kcal/mol):	0.33

JAG-UCB-f37eaa14-7_1



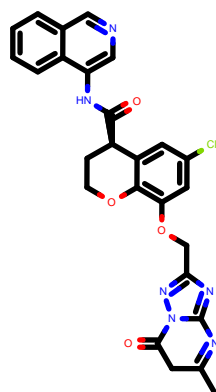
CID:	JAG-UCB-f37eaa14-7_1
SMILES:	<chem>Cc1c[nH]1c2cc[nH]2c3cc(cc3)Cl)c4cc5c6c7c8c9c5c6c7c8c9c4</chem>
RUN:	RUN3066
DDG (kcal/mol):	-0.55
dDDG (kcal/mol):	0.18

MAT-POS-53907a1c-3_2



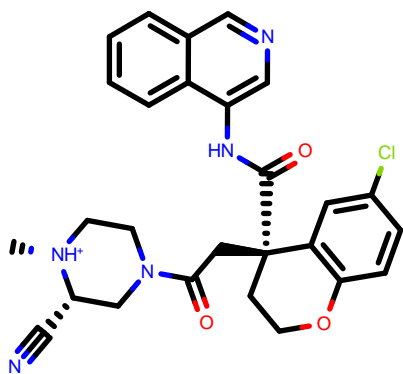
CID:	MAT-POS-53907a1c-3_2
SMILES:	<chem>c1ccc2c(c1)cncc2N(C(=O)N(CCC3CCCC3)c4cc(cc(c4)Cl)O)C@H]5CC(=O)N5</chem>
RUN:	RUN526
DDG (kcal/mol):	-0.55
dDDG (kcal/mol):	0.43

ALP-POS-e0fe77e5-1_1



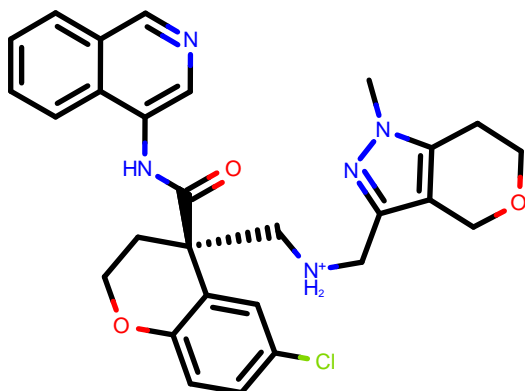
CID:	ALP-POS-e0fe77e5-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3CCCC[C@@H](C3=O)c4ccc(cc4)Cl</chem>
RUN:	RUN2318
DDG (kcal/mol):	-0.55
dDDG (kcal/mol):	0.48

JOH-SUS-a69c159d-5_1



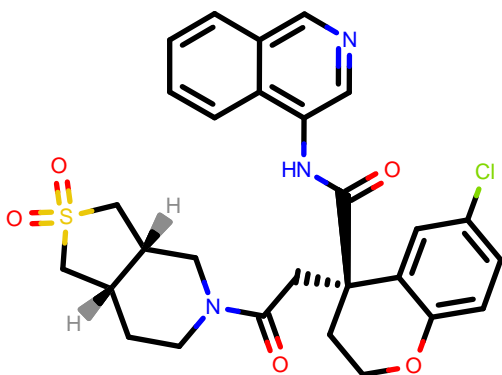
CID:	JOH-SUS-a69c159d-5_1
SMILES:	<chem>c1cc2c(cc1F)c(cnc2F)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl</chem>
RUN:	RUN1125
DDG (kcal/mol):	-0.55
dDDG (kcal/mol):	0.25

RAL-THA-05e671eb-28_2



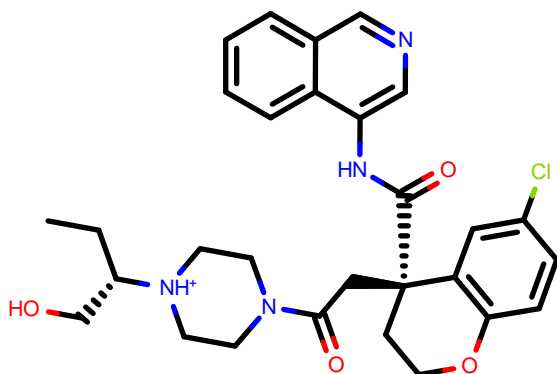
CID:	RAL-THA-05e671eb-28_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCOC4c3c(c(cc4)Cl)F</chem>
RUN:	RUN2061
DDG (kcal/mol):	-0.55
dDDG (kcal/mol):	0.62

MIC-UNK-0a05c952-1_2



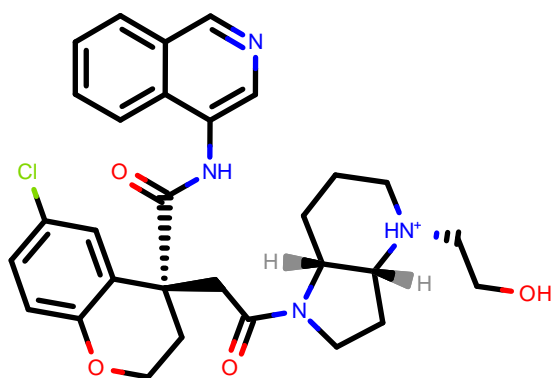
CID:	MIC-UNK-0a05c952-1_2
SMILES:	<chem>c1ccc2c(c1)cnc2N3[C@H](C[C@@H](C3=O)c4cccc(c4)Cl)[C@@H]5COC5</chem>
RUN:	RUN3497
DDG (kcal/mol):	-0.55
dDDG (kcal/mol):	0.18

RAL-THA-8416115c-2_2



CID:	RAL-THA-8416115c-2_2
SMILES:	<chem>CN1CC[C@@H](c2c1ccc(c2)Cl)C(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN1252
DDG (kcal/mol):	-0.55
dDDG (kcal/mol):	0.38

PET-UNK-8c422e11-1_1



CID: PET-UNK-8c422e11-1_1

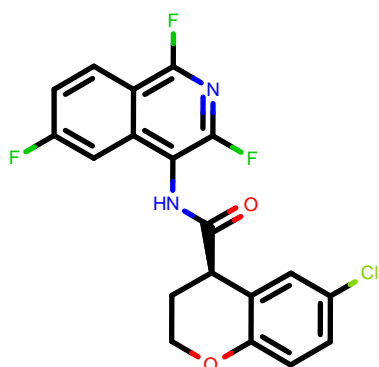
SMILES: CC(c1cc(cc(c1)Cl)CC(=O)Nc2cncc3c2cccc3)(F)F

RUN: RUN3712

DDG (kcal/mol): -0.54

dDDG (kcal/mol): 0.45

EDJ-MED-8c98ee63-1_1



CID: EDJ-MED-8c98ee63-1_1

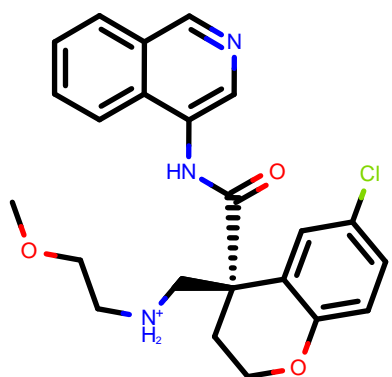
SMILES: COC[C][NH2+][C][C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN2849

DDG (kcal/mol): -0.54

dDDG (kcal/mol): 0.07

EDG-MED-90036822-55_1



CID: EDG-MED-90036822-55_1

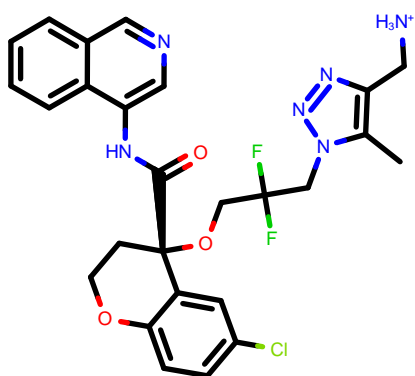
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)NC(=O)c5ccc(nc5)N6CC(C6)[NH3+]

RUN: RUN1735

DDG (kcal/mol): -0.54

dDDG (kcal/mol): 0.45

FRA-DIA-b66f7109-2_1



CID: FRA-DIA-b66f7109-2_1

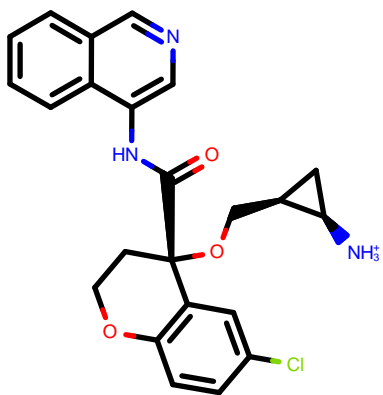
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3COc4c3cc(cc4Cl)O[C@@H]5CC(=O)N5

RUN: RUN393

DDG (kcal/mol): -0.54

dDDG (kcal/mol): 0.36

DAR-DIA-23e5a6a0-4_1



CID: DAR-DIA-23e5a6a0-4_1

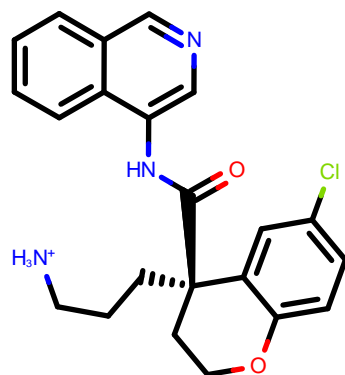
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4c3cc(cc4O)[C@@H]5CC6(CC6)C[NH2+]5Cl

RUN: RUN407

DDG (kcal/mol): -0.54

dDDG (kcal/mol): 0.29

ALP-UNI-8e43a71e-5_4



CID: ALP-UNI-8e43a71e-5_4

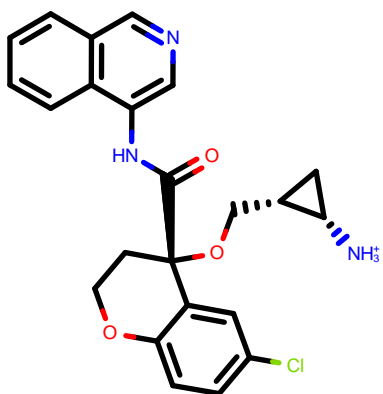
SMILES: ClC1=CC=C(C=C1)C2=CC=CC=C2NC(=O)[C@@H]3CCOCc4c3cc(cc4O)[C@@H]5CC6(CC6)C[NH2+]5Cl

RUN: RUN2948

DDG (kcal/mol): -0.53

dDDG (kcal/mol): 0.13

PET-UNK-c5865d42-1_2



CID: PET-UNK-c5865d42-1_2

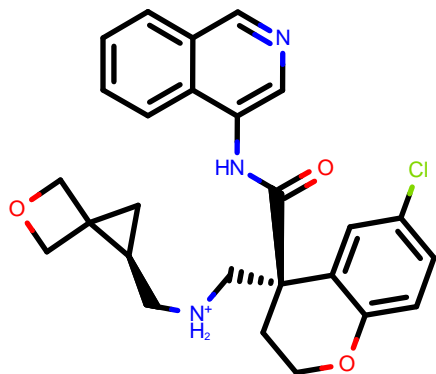
SMILES: c1ccc2c(c1)cncc2NC(=O)Cc3cc(cc(c3)Cl)C[C@H]4CC[NH2+]4

RUN: RUN424

DDG (kcal/mol): -0.53

dDDG (kcal/mol): 0.31

RAL-THA-05e671eb-20_2



CID: RAL-THA-05e671eb-20_2

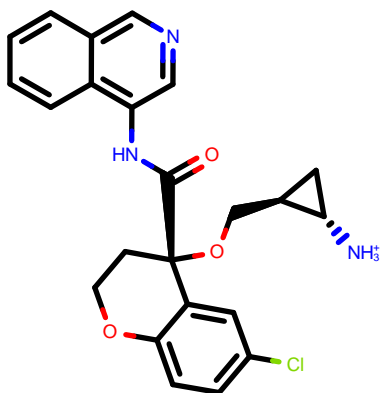
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4c3cc(cc4C#N)Cl

RUN: RUN2045

DDG (kcal/mol): -0.53

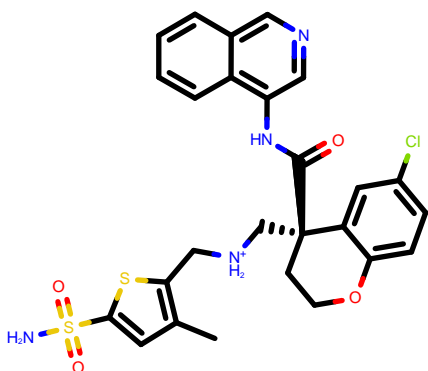
dDDG (kcal/mol): 0.44

ALP-POS-fe871b40-15_1



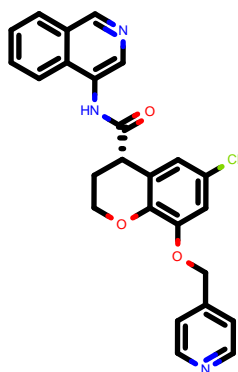
CID:	ALP-POS-fe871b40-15_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CC(=O)Nc4c3cc(c(c4)F)Cl</chem>
RUN:	RUN3135
DDG (kcal/mol):	-0.53
dDDG (kcal/mol):	0.15

EDJ-MED-1981ceba-4_4



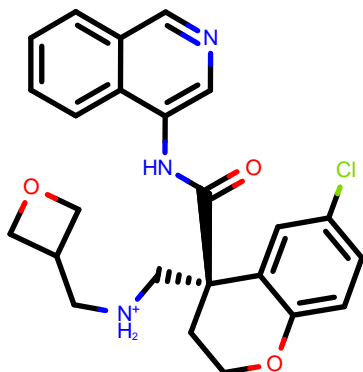
CID:	EDJ-MED-1981ceba-4_4
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3C[N@](C)C4C3cc(cc4)Cl)S(=O)(=O)N5CC(C5)C#N</chem>
RUN:	RUN4695
DDG (kcal/mol):	-0.53
dDDG (kcal/mol):	0.29

EDG-MED-90036822-60_1



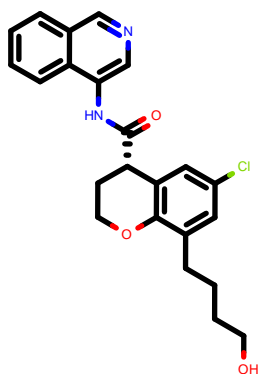
CID:	EDG-MED-90036822-60_1
SMILES:	<chem>C[C@@H](C(=O)N[C@@]1(C)COC2C1C(C)C(=O)Nc3ccc4c3ccc4)n5ccn5</chem>
RUN:	RUN1738
DDG (kcal/mol):	-0.52
dDDG (kcal/mol):	0.24

EDG-MED-90036822-90_1



CID:	EDG-MED-90036822-90_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(C)COC4C3cc(cc4)Cl)N(C(=O)[C@@H](c5ccccc5F))N[HS+]</chem>
RUN:	RUN1795
DDG (kcal/mol):	-0.52
dDDG (kcal/mol):	0.31

ALP-POS-477dc5b7-2_1



CID: ALP-POS-477dc5b7-2_1

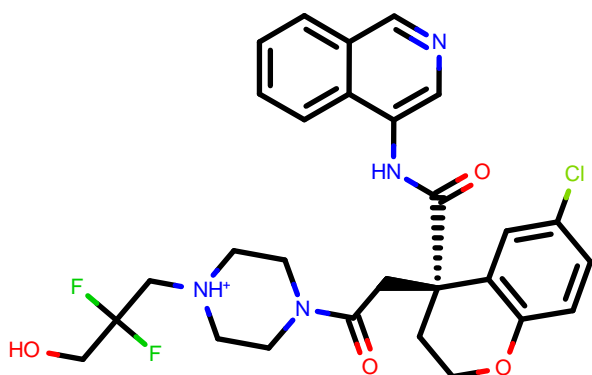
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCNc4c3cc(cc4)Cl

RUN: RUN299

DDG (kcal/mol): -0.52

dDDG (kcal/mol): 0.23

LON-WEI-5e7d1b3e-33_1



CID: LON-WEI-5e7d1b3e-33_1

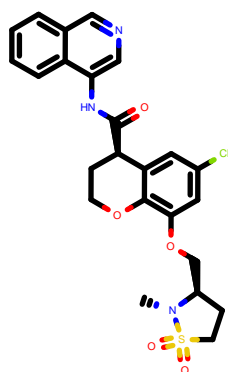
SMILES: C[C@@H]1c2cc(c(cc2CCN1C(=O)Nc3cn(c(=O)c4c3cccc4)C)OC)OC

RUN: RUN1334

DDG (kcal/mol): -0.52

dDDG (kcal/mol): 0.39

ALP-UNI-76695c4f-10_1



CID: ALP-UNI-76695c4f-10_1

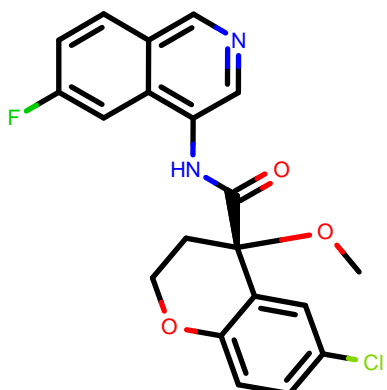
SMILES: C[C@@]12CN[C@@]1(C1=O)NC2=O(C)C(=O)C[C@@]3(CCOc4c3cc(cc4)Cl)C(=O)Nc5ncnc6s5ccc6

RUN: RUN2174

DDG (kcal/mol): -0.51

dDDG (kcal/mol): 0.36

PET-UNK-83d689b6-2_1



CID: PET-UNK-83d689b6-2_1

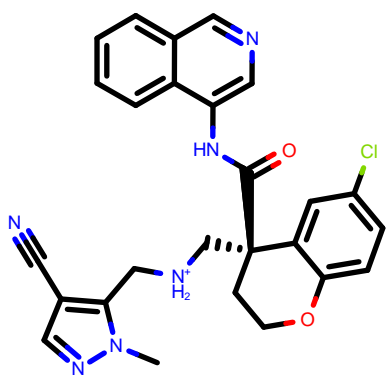
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OCCc5cccn5

RUN: RUN4194

DDG (kcal/mol): -0.51

dDDG (kcal/mol): 0.07

LEE-CAM-7ab9b158-4_4



CID: LEE-CAM-7ab9b158-4_4

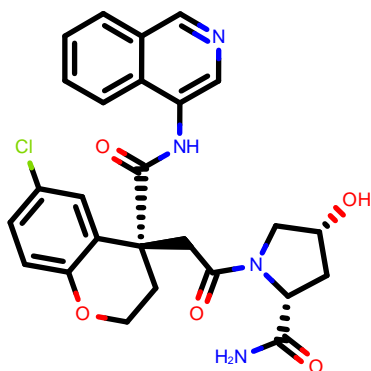
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@]3(CCOc4c3cc(cc4)C)COC[C@@H]5CN6C[C@@H](O5)CCC6=O

RUN: RUN2213

DDG (kcal/mol): -0.51

dDDG (kcal/mol): 0.44

ERI-UCB-d6de1f3c-5_2



CID: ERI-UCB-d6de1f3c-5_2

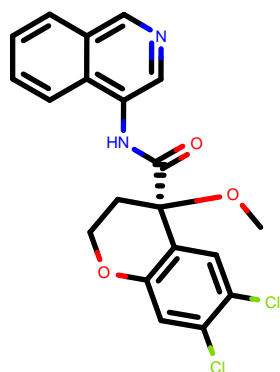
SMILES: c1ccc2c(c1)ncnc2C(=O)N3C[C@H](N(C(=O)C3)c4cccc(c4)C)CC5CCCC5

RUN: RUN1096

DDG (kcal/mol): -0.51

dDDG (kcal/mol): 0.33

DAR-DIA-9e4459de-11_11



CID: DAR-DIA-9e4459de-11_11

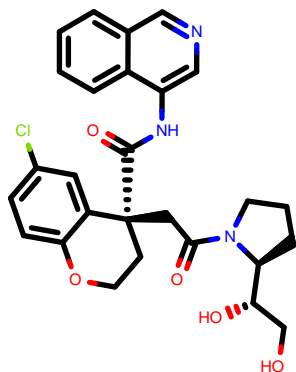
SMILES: c1cc2c(c1)ncnc2C(=O)C3[C@H](N(C(=O)C3)C4=C(C=C(C=C4)Cl)Cl)C5=C(C=C(C=C5)OC)C

RUN: RUN1419

DDG (kcal/mol): -0.51

dDDG (kcal/mol): 0.12

BEN-BAS-5c03e89e-2_4



CID: BEN-BAS-5c03e89e-2_4

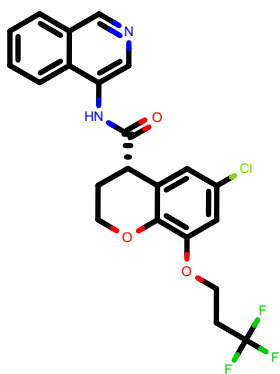
SMILES: CC(C)([C@H]1C=Nc2ccc(cc2[C@H]1C(=O)Nc3cnc4c3ccc4)Cl)O

RUN: RUN1146

DDG (kcal/mol): -0.50

dDDG (kcal/mol): 0.40

EDG-MED-90036822-5_1



CID: EDG-MED-90036822-5_1

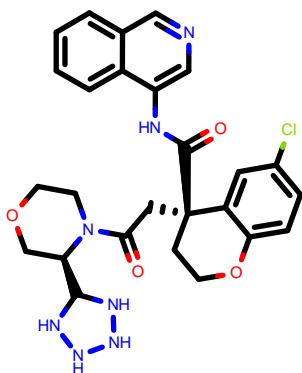
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOC4c3cc(cc4)Cl)NC(=O)[C@@H](CF)(NH3+)

RUN: RUN1660

DDG (kcal/mol): -0.49

dDDG (kcal/mol): 0.30

JOH-SUS-a69c159d-7_1



CID: JOH-SUS-a69c159d-7_1

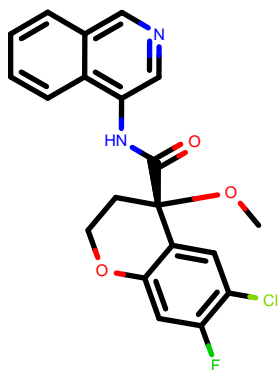
SMILES: c1cc2c(cc1F)c(c(nc2F)F)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl

RUN: RUN1130

DDG (kcal/mol): -0.48

dDDG (kcal/mol): 0.38

MAT-POS-4223bc15-23_3



CID: MAT-POS-4223bc15-23_3

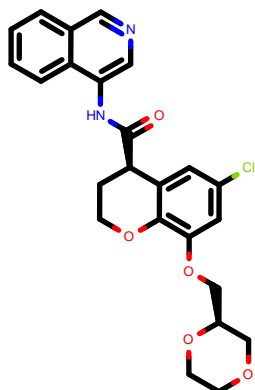
SMILES: CNC(=O)C[N@@H+]1Cc2ccc(cc2[C@@H](C1)C(=O)Nc3ncc4c3ccc4)Cl

RUN: RUN4108

DDG (kcal/mol): -0.48

dDDG (kcal/mol): 0.06

PET-UNK-2c6614b6-6_1



CID: PET-UNK-2c6614b6-6_1

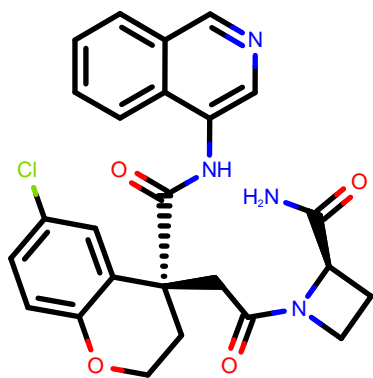
SMILES: c1cc2ncc(c2cc1F)NC(=O)[C@@H]3CCS(=O)(=O)c4c3cc(cc4)Cl

RUN: RUN4811

DDG (kcal/mol): -0.48

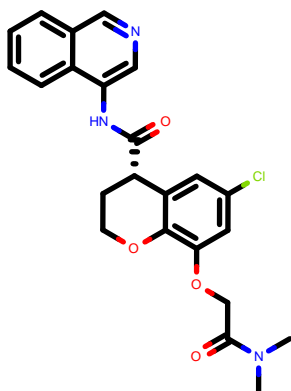
dDDG (kcal/mol): 0.17

BEN-BAS-c2bc0d80-7_1



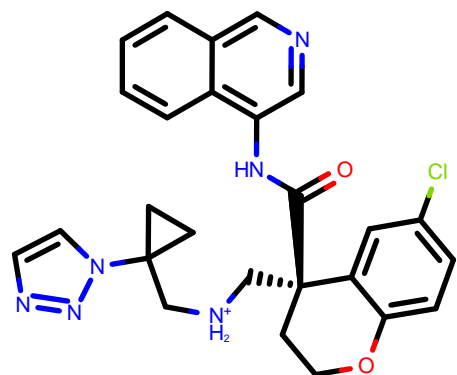
CID:	BEN-BAS-c2bc0d80-7_1
SMILES:	<chem>CN1C(=O)N(C(=O))[C@@]12CCOCc3c2cc(cc3)Cl)c4ncc5c4cccc5</chem>
RUN:	RUN1138
DDG (kcal/mol):	-0.48
dDDG (kcal/mol):	0.27

VLA-UNK-f702bf1c-4_1



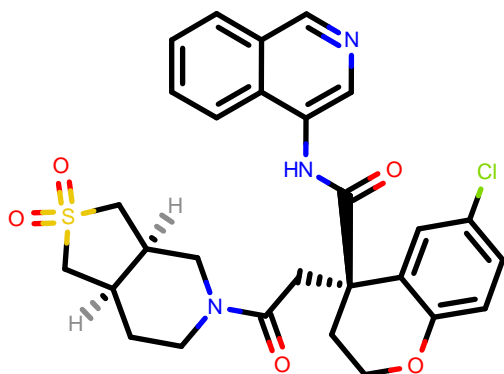
CID:	VLA-UNK-f702bf1c-4_1
SMILES:	<chem>c1ccc2c(c1)ncc2N3C(=O)[C@@]4(CCOc5c4cc(cc5)Cl)N(C3=O)C[C@@]4Hj6CC(=O)N(C)C</chem>
RUN:	RUN2313
DDG (kcal/mol):	-0.48
dDDG (kcal/mol):	0.26

EDJ-MED-1981ceba-3_3



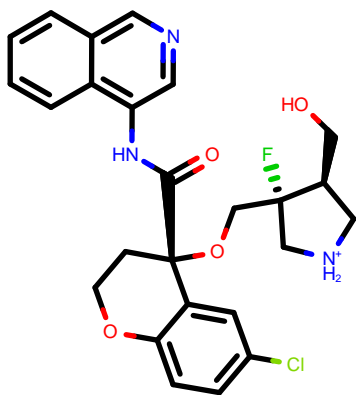
CID:	EDJ-MED-1981ceba-3_3
SMILES:	<chem>OOC1CN(C1)S(=O)(=O)N@@]2Cc3ccc(cc3[C@H](C2)C(=O)Nc4ncc5c4cccc5)Cl</chem>
RUN:	RUN4691
DDG (kcal/mol):	-0.48
dDDG (kcal/mol):	0.28

MIC-UNK-5a93dd5f-12_2



CID:	MIC-UNK-5a93dd5f-12_2
SMILES:	<chem>c1ccc2c(c1)ncc2NC(=O)[C@H](c3cccc(c3)Cl)N@@]4CC[C@@]4H(C4)NH+5CCCCC5</chem>
RUN:	RUN791
DDG (kcal/mol):	-0.48
dDDG (kcal/mol):	0.28

JAG-UCB-f37eaa14-6_1



CID: JAG-UCB-f37eaa14-6_1

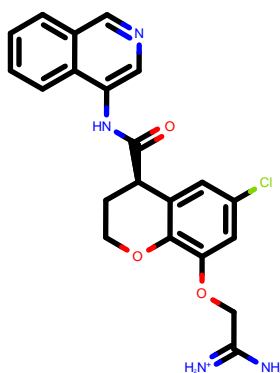
SMILES: c1ccc2c(c1)cnc2N3CC[C@@H](C3=O)C[N@@]([c5c4cc(cc5)Cl])[C@@H]6CC(=O)N6

RUN: RUN3062

DDG (kcal/mol): -0.48

dDDG (kcal/mol): 0.21

PET-UNK-1320d94d-3_1



CID: PET-UNK-1320d94d-3_1

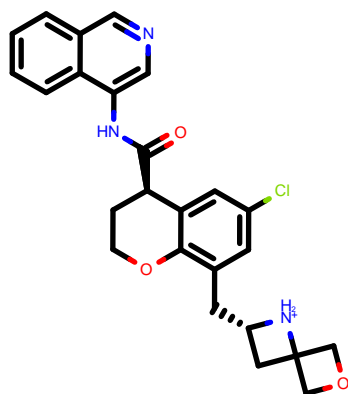
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]([c3ccc(c3)Cl]N[C@@]([C@@H]4CC[C@@H]4)O)N5CCCC5

RUN: RUN4974

DDG (kcal/mol): -0.48

dDDG (kcal/mol): 0.18

DAR-DIA-0cde14eb-58_1



CID: DAR-DIA-0cde14eb-58_1

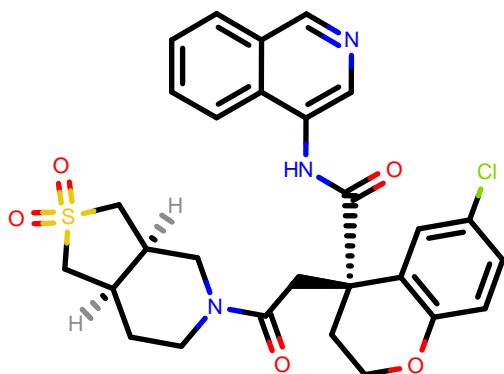
SMILES: C[C@@H]([c1cccc(c1)C2(CC2)I])C(=O)Nc3cnc4c3cccc4

RUN: RUN28

DDG (kcal/mol): -0.47

dDDG (kcal/mol): 0.45

MAK-UNK-c749d764-21_7



CID: MAK-UNK-c749d764-21_7

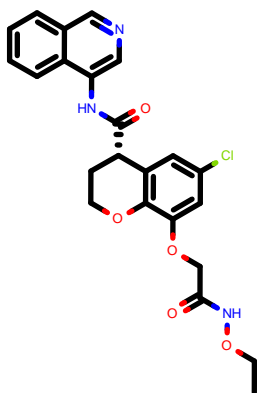
SMILES: CCCCN(c1cnc2c1cccc2)C(=O)C[C@@H]3CCC[C@H]([C@H]3O)C(F)F

RUN: RUN1021

DDG (kcal/mol): -0.47

dDDG (kcal/mol): 0.24

EDJ-MED-d203f206-10_1



CID: EDJ-MED-d203f206-10_1

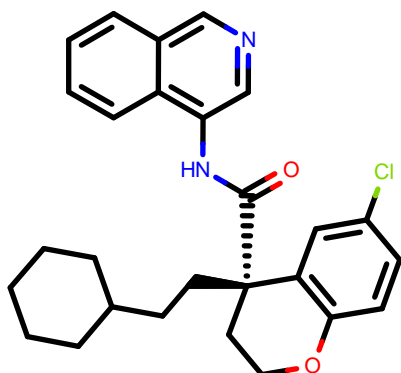
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@H](C)COC3COCc4cc3cc4)C(=O)N5CCCC6(C5)C@H](C)C@H)6O

RUN: RUN2572

DDG (kcal/mol): -0.47

dDDG (kcal/mol): 0.28

ALP-UNI-dbbfd3db-15_2



CID: ALP-UNI-dbbfd3db-15_2

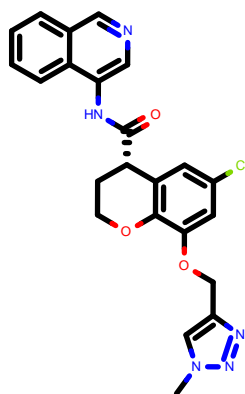
SMILES: C[C@H](C(=O)N[C@@H]1(CCOc2c1cc(cc2)C)C(=O)Nc3ccc4c3ccc4)NC(=O)C5CCCC5

RUN: RUN2786

DDG (kcal/mol): -0.47

dDDG (kcal/mol): 0.29

ALP-POS-5bb456a5-1_2



CID: ALP-POS-5bb456a5-1_2

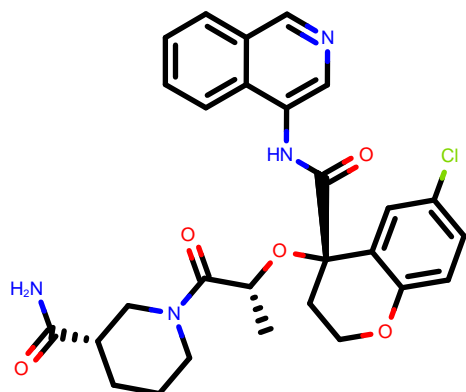
SMILES: C[C@H](C)N[C@@H](C)C(=O)N[C@@H]1(CCOc2c1cc(cc2)C)C(=O)Nc3ccc4c3ccc4)S(=O)(=O)C

RUN: RUN2407

DDG (kcal/mol): -0.47

dDDG (kcal/mol): 0.25

MIC-UNK-5a93dd5f-2_6



CID: MIC-UNK-5a93dd5f-2_6

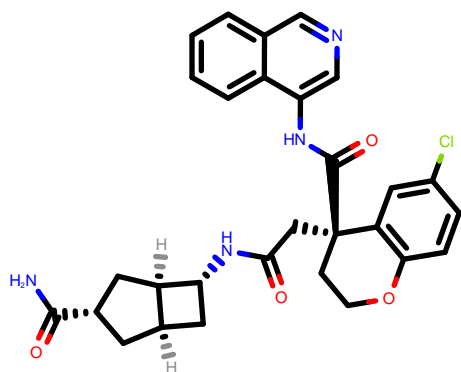
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@H](C)COC3COCc4cc3cc4)C(=O)N5CCCC6(C5)C@H](C)C@H)6O

RUN: RUN739

DDG (kcal/mol): -0.47

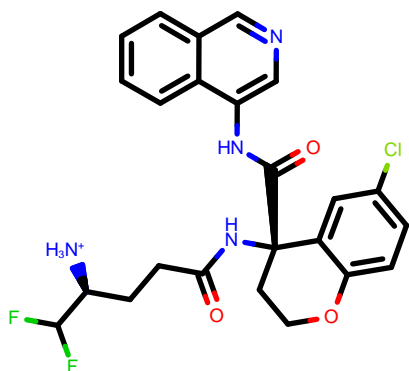
dDDG (kcal/mol): 0.29

LAU-MED-88a3970a-17_1



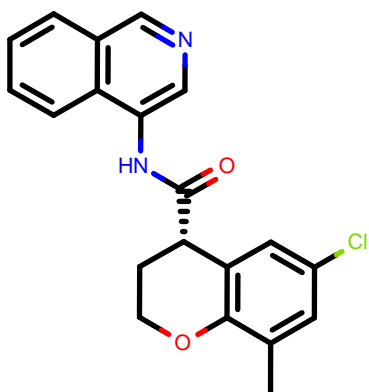
CID:	LAU-MED-88a3970a-17_1
SMILES:	<chem>CNC(=O)CCc1cc(cc2c1OCC[C@H]2C(=O)Nc3ncc4c3cccc4)Cl</chem>
RUN:	RUN1513
DDG (kcal/mol):	-0.47
dDDG (kcal/mol):	0.56

MIC-UNK-cdc2493e-4_3



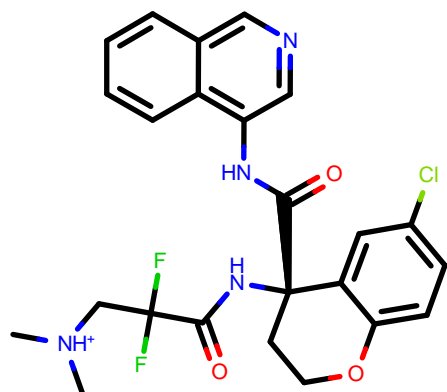
CID:	MIC-UNK-cdc2493e-4_3
SMILES:	<chem>c1ccc2c(c1)nccc2NC(=O)N(c3cccc(c3)Cl)C4C[C@H]5CCCC[C@H]5C4</chem>
RUN:	RUN529
DDG (kcal/mol):	-0.46
dDDG (kcal/mol):	0.41

MAK-UNK-ffc90da7-3_2



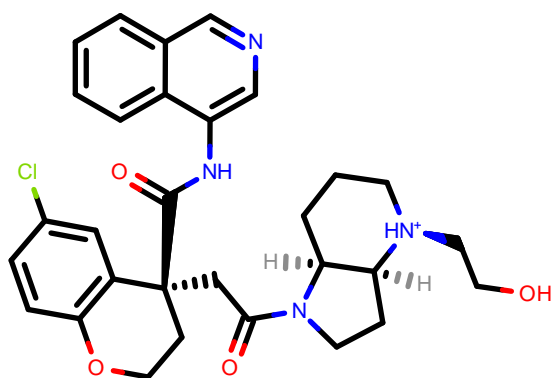
CID:	MAK-UNK-ffc90da7-3_2
SMILES:	<chem>CC(C)OC[C@H](c1ccc2c(c1)nccc2NC(=O)Cc3cccc(c3)Cl)[NH2+]C</chem>
RUN:	RUN696
DDG (kcal/mol):	-0.46
dDDG (kcal/mol):	0.12

MAT-POS-3b97339c-3_2



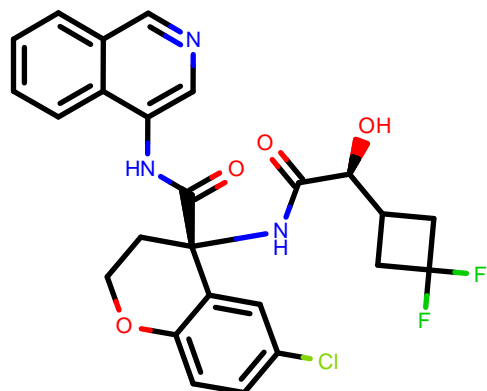
CID:	MAT-POS-3b97339c-3_2
SMILES:	<chem>c1ccc2c(c1)nccc2NC(=O)[C@]3(CCOc4c3cc(cc4)Cl)CNC(=O)c5cn[nH]c5S(=O)(=O)N</chem>
RUN:	RUN3308
DDG (kcal/mol):	-0.46
dDDG (kcal/mol):	0.16

LON-WEI-5e7d1b3e-50_2



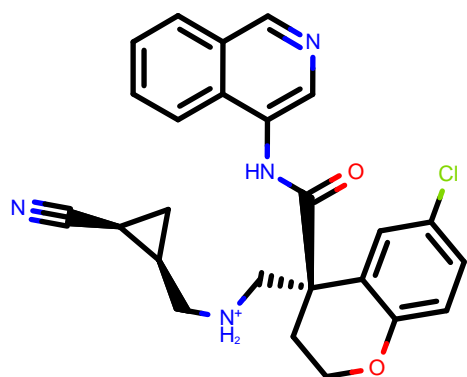
CID:	LON-WEI-5e7d1b3e-50_2
SMILES:	<chem>CCCC[N@H+](CCNC(=O)Nc1cn(c(=O)c2c1cccc2)C)Cc3ccco3</chem>
RUN:	RUN1365
DDG (kcal/mol):	-0.45
dDDG (kcal/mol):	0.48

JAG-UCB-706446eb-4_1



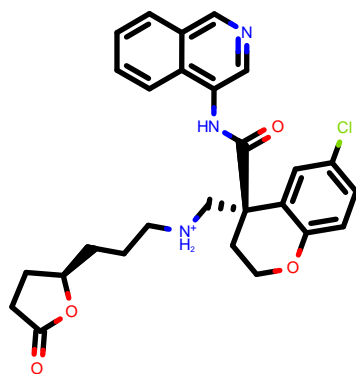
CID:	JAG-UCB-706446eb-4_1
SMILES:	<chem>CCOC(=O)CC[C@]1(C)C(CO)c2c1cc(cc2)C(C(=O)Nc3cncc4c3ccoc4)</chem>
RUN:	RUN619
DDG (kcal/mol):	-0.45
dDDG (kcal/mol):	0.26

EDG-MED-90036822-38_2



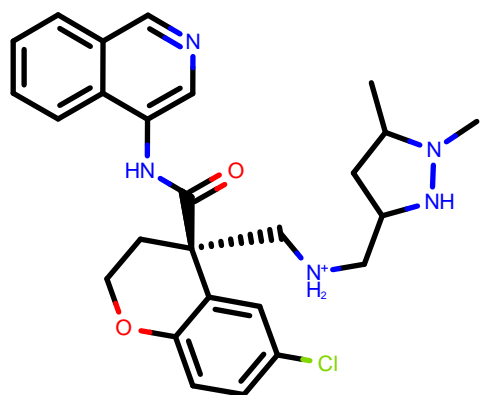
CID:	EDG-MED-90036822-38_2
SMILES:	<chem>C[C@H](CNCC(C(=O)N[C@]1(C)C(CO)c2c1cc(cc2)C(C(=O)Nc3cncc4c3ccoc4)(F)F)O</chem>
RUN:	RUN1720
DDG (kcal/mol):	-0.45
dDDG (kcal/mol):	0.42

MAT-POS-61f37a1a-8_2



CID:	MAT-POS-61f37a1a-8_2
SMILES:	<chem>Cc1c(c(n1)C)N2CCOCC2)C(NH2+)C[C@]3(C)C(CO)c4c3cc(cc4)C(C(=O)Nc5cncc6c5ccoc6)</chem>
RUN:	RUN4602
DDG (kcal/mol):	-0.45
dDDG (kcal/mol):	0.24

EDG-MED-90036822-52_1



CID: EDG-MED-90036822-52_1

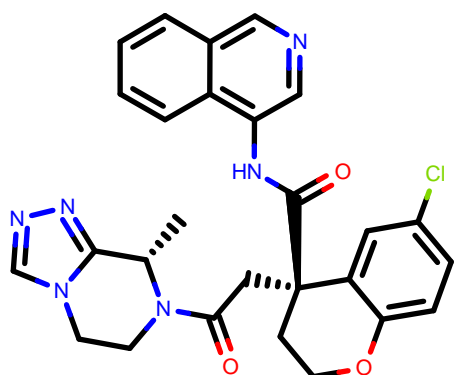
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)C)NC(=O)C(CNCCCO)(F)F

RUN: RUN1737

DDG (kcal/mol): -0.45

dDDG (kcal/mol): 0.37

EDJ-MED-841e0cf0-2_1



CID: EDJ-MED-841e0cf0-2_1

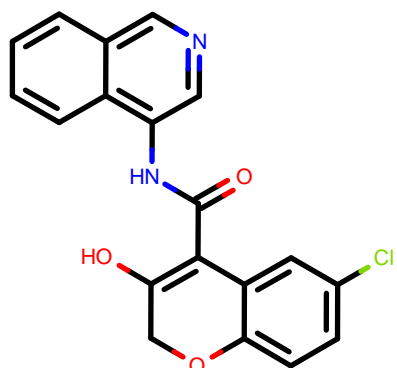
SMILES: CO[C@@]1(C)N@@(C)C2c1cc(c(c2)F)Cl)S(=O)(=O)C(=O)C(=O)Nc3ncc4c3ccoc4

RUN: RUN3830

DDG (kcal/mol): -0.45

dDDG (kcal/mol): 0.15

EDG-MED-ba1ac7b9-18_1



CID: EDG-MED-ba1ac7b9-18_1

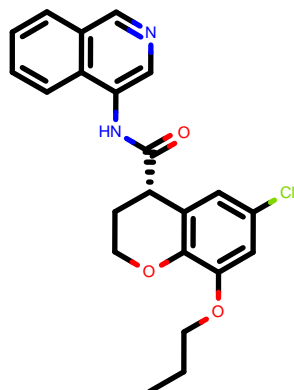
SMILES: C[C@@]1(C)N2c(ccn2)CN1C(=O)C[C@@]3(CCOc4c3cc(cc4)C)C(=O)Nc5ncc6c5ccoc6

RUN: RUN2682

DDG (kcal/mol): -0.44

dDDG (kcal/mol): 0.23

EDG-MED-971238d3-6_1



CID: EDG-MED-971238d3-6_1

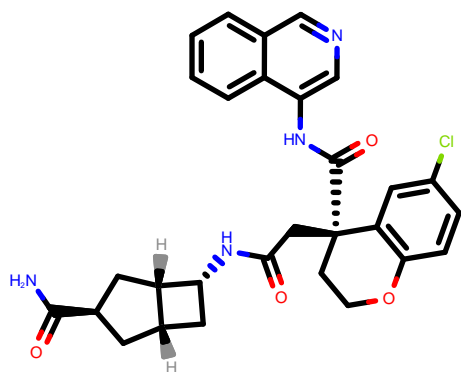
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)C)CC[NH3+]

RUN: RUN1471

DDG (kcal/mol): -0.44

dDDG (kcal/mol): 0.26

LON-WEI-5e7d1b3e-35_2



CID: LON-WEI-5e7d1b3e-35_2

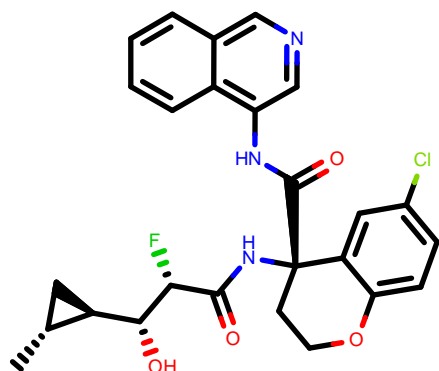
SMILES: Cc1ccc(cc1)C[N@H]2CC[C@@H](C2)CNC(=O)Nc3cn(c(O)c4c3cccc4)CC(C)C

RUN: RUN1341

DDG (kcal/mol): -0.44

dDDG (kcal/mol): 0.27

MAT-POS-3b97339c-3_1



CID: MAT-POS-3b97339c-3_1

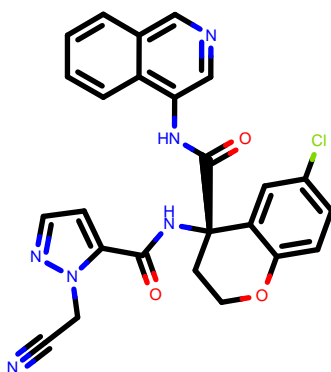
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3(CCOc4c3cc(c4)Cl)CNC(=O)c5cn[nH]c5S(=O)(=O)N

RUN: RUN3307

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.18

KAD-UNI-877d7bed-18_2



CID: KAD-UNI-877d7bed-18_2

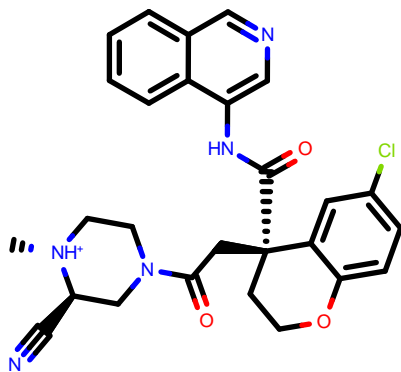
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3(CCOc4c3cc(c4)OCCS(=O)(=O)CCCO)Cl

RUN: RUN3771

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.14

DAR-DIA-076fb6ea-8_1



CID: DAR-DIA-076fb6ea-8_1

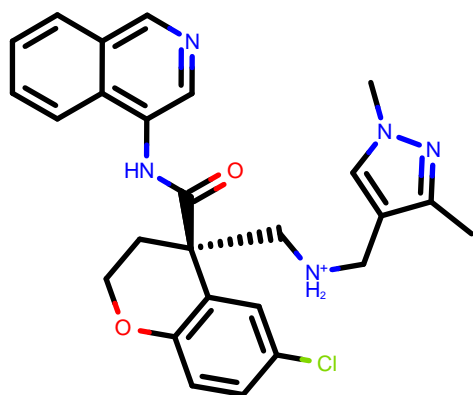
SMILES: C=C=CC(=O)N(c1cnc2c1cccc2)C(=O)Cc3cccc(c3)Cl

RUN: RUN1400

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.23

ED_-GRI-5b13fbe2-74_4



CID: ED_-GRI-5b13fbe2-74_4

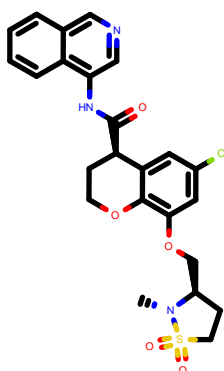
SMILES: C[N@H+]1CCO[C@H](C1)c2nccc(n2)CO[C@]3(CCOc4c3cc(cc4)Cl)C(=O)Nc5ncc6c5cccc6

RUN: RUN1634

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.51

ALP-POS-67d5babe-1_3



CID: ALP-POS-67d5babe-1_3

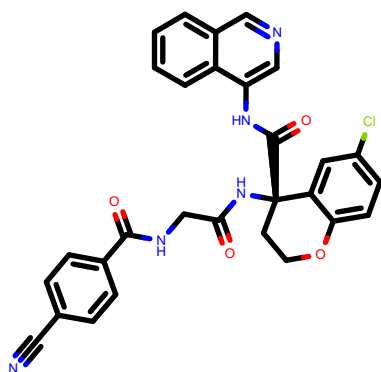
SMILES: c1ccc2c(c1)nccc2NC(=O)[C@@]3(COCC4=CC(=O)C(C1)CQ[N@H]5CC[C@@H]5C(S(=O)(=O)C)C@H)3C1=CC=C(Cl)C=C1

RUN: RUN4868

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.19

ALP-POS-966f8da6-1_4



CID: ALP-POS-966f8da6-1_4

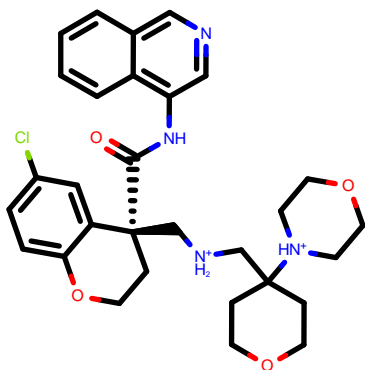
SMILES: c1ccc2c(c1)nccc2NC(=O)[C@H]3CCN(c4c3cc(cc4)Cl)Cc5cnc[nH]5

RUN: RUN1223

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.31

MIC-UNK-f792ef5d-1_1



CID: MIC-UNK-f792ef5d-1_1

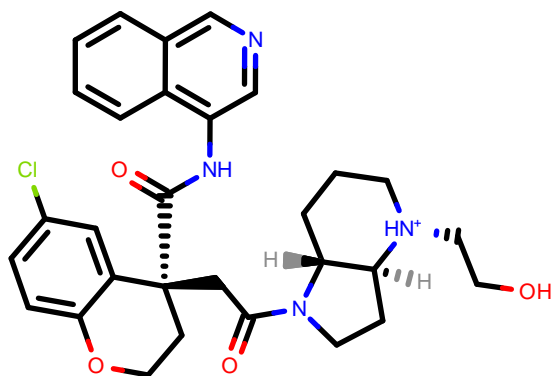
SMILES: C[N@@H+]1[C@H]2COc3ccc(cc3[C@]2(C1)C(=O)Nc4cnc5c4cccc5)Cl

RUN: RUN4665

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.32

MAT-POS-4223bc15-7_1



CID: MAT-POS-4223bc15-7_1

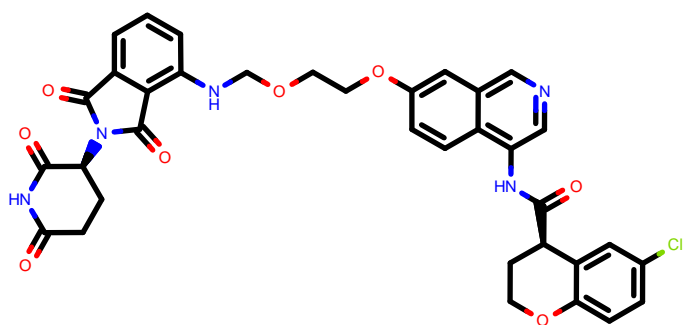
SMILES: c1ccc2c(c1)ncnc2NC(=O)[C@@H]3C[N@@]4(Cc4c3ccc(cc4)Cl)S(=O)(=O)C5COCS

RUN: RUN3997

DDG (kcal/mol): -0.43

dDDG (kcal/mol): 0.35

MIC-UNK-25b9c114-1_1



CID: MIC-UNK-25b9c114-1_1

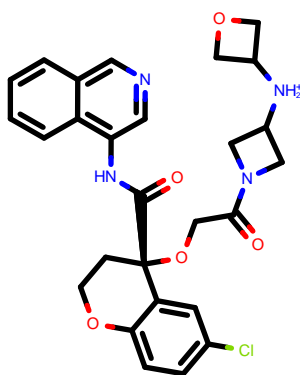
SMILES: CS(=O)(=O)NC[C@@H]1CN(C(=O)CN1C(=O)c2ccc3c2ccc3c4ccc(cc4)Cl

RUN: RUN3259

DDG (kcal/mol): -0.42

dDDG (kcal/mol): 0.28

EDJ-MED-15e90dfc-7_1



CID: EDJ-MED-15e90dfc-7_1

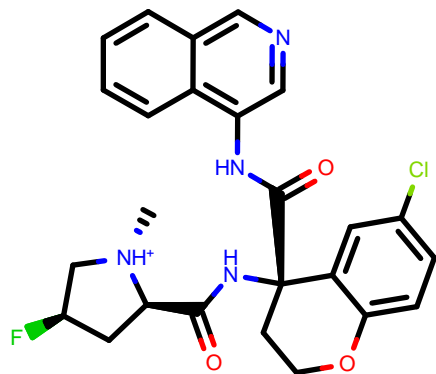
SMILES: CS(=O)(=O)CCNC[C@@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3ccc4c3ccc4

RUN: RUN3456

DDG (kcal/mol): -0.42

dDDG (kcal/mol): 0.24

MIC-UNK-cdc2493e-12_1



CID: MIC-UNK-cdc2493e-12_1

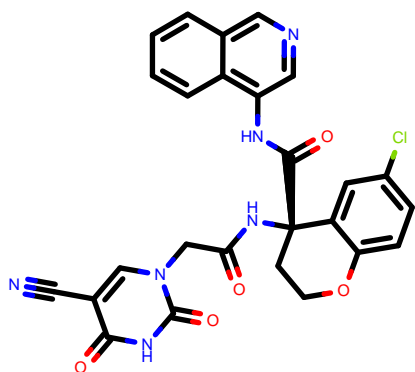
SMILES: C[NH+]1C[C@@H]1CC[C@@H]1C1N(c2ccc(c2)Cl)C(=O)Nc3ccc4c3ccc4

RUN: RUN551

DDG (kcal/mol): -0.42

dDDG (kcal/mol): 0.31

MAT-POS-4223bc15-14_3



CID: MAT-POS-4223bc15-14_3

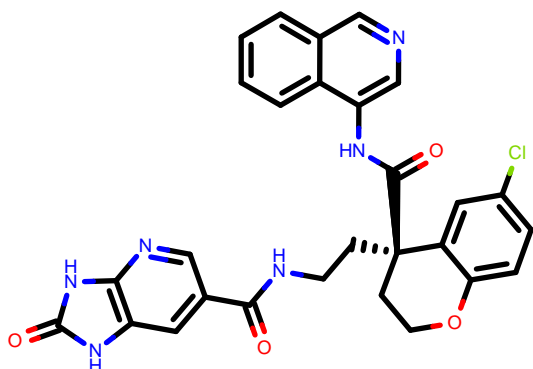
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3C[N@@]([C@4c3cc(cc4)Cl]S(=O)(=O)CC5C[NH2+]C5

RUN: RUN4068

DDG (kcal/mol): -0.42

dDDG (kcal/mol): 0.16

DAR-DIA-0d514e7d-31_16



CID: DAR-DIA-0d514e7d-31_16

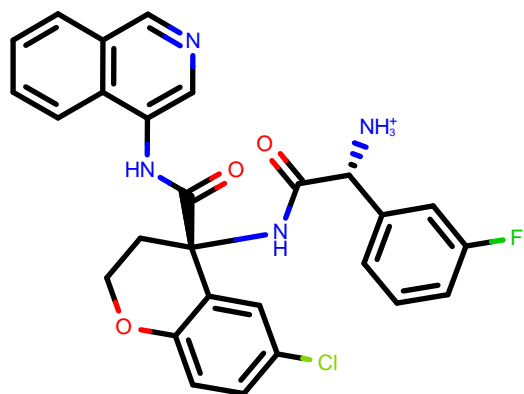
SMILES: C[C@H]1CCO[C@H]2C=CC(=C[C@H]2[C@H]1C(=O)Nc3cnc4c3cccc4)Cl

RUN: RUN849

DDG (kcal/mol): -0.42

dDDG (kcal/mol): 0.29

MIC-UNK-bcd487e9-1_1



CID: MIC-UNK-bcd487e9-1_1

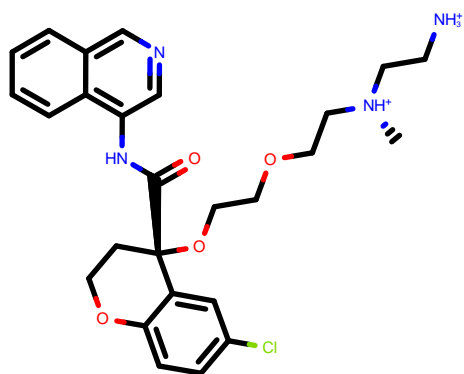
SMILES: c1ccc(cc1)CN(c2cccc(c2)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN586

DDG (kcal/mol): -0.42

dDDG (kcal/mol): 0.31

WIL-UCB-7ba4ac3a-3_1



CID: WIL-UCB-7ba4ac3a-3_1

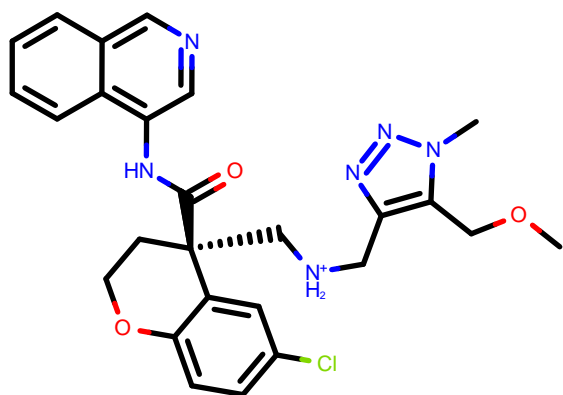
SMILES: c1ccc2c(c1)cnc2NS(=O)(=O)[C@@H]3CCOC4c3cc(cc4)Cl

RUN: RUN3020

DDG (kcal/mol): -0.42

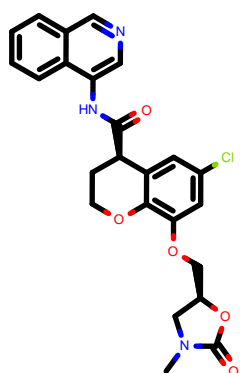
dDDG (kcal/mol): 0.30

EDG-MED-10fcb19e-1_2



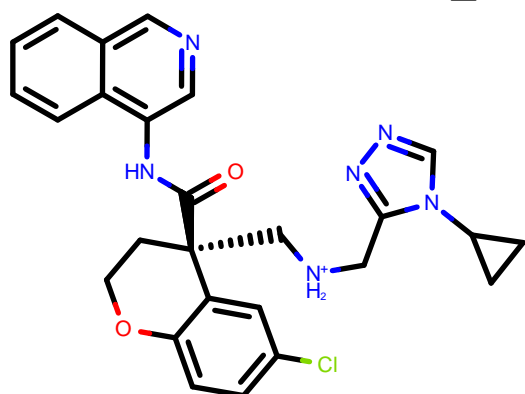
CID:	EDG-MED-10fcb19e-1_2
SMILES:	<chem>CO[C@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cc(cc4)NS(=O)(=O)C</chem>
RUN:	RUN4497
DDG (kcal/mol):	-0.42
dDDG (kcal/mol):	0.27

EDG-MED-ba1ac7b9-4_4



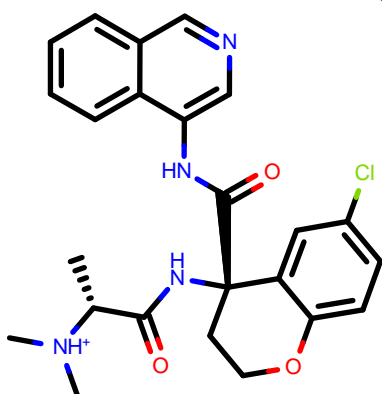
CID:	EDG-MED-ba1ac7b9-4_4
SMILES:	<chem>c1ccc2c(c1)ncnc2N(C=O)[C@]3(CCOc4c3cc(cc4)Cl)C(=O)N5CC(C5)N#H#CC(C)C@H(C)O</chem>
RUN:	RUN2629
DDG (kcal/mol):	-0.41
dDDG (kcal/mol):	0.34

EDJ-MED-c82a5324-1_7



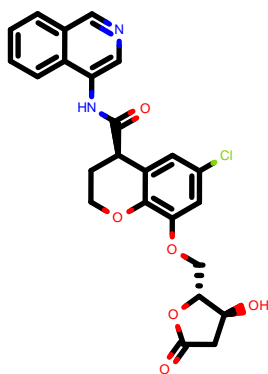
CID:	EDJ-MED-c82a5324-1_7
SMILES:	<chem>CNC(=O)[C@@H](CO)[N@H]1C2cocc(cc2[C@H](C1)C(=O)Nc3cncc4c3ccc4)Cl</chem>
RUN:	RUN4726
DDG (kcal/mol):	-0.41
dDDG (kcal/mol):	0.22

MAK-UNK-83e0a0b4-1_1



CID:	MAK-UNK-83e0a0b4-1_1
SMILES:	<chem>CCCC1ccc2c(c1)[C@@H](CO)[N@H]1C2cocc(cc2[C@H](C1)C(=O)Nc3cncc4c3ccc4)COC</chem>
RUN:	RUN727
DDG (kcal/mol):	-0.41
dDDG (kcal/mol):	0.36

ALP-POS-5bb456a5-6_1



CID: ALP-POS-5bb456a5-6_1

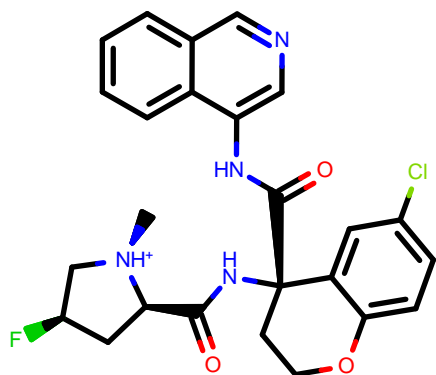
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCC(CC5)c6n[nH]n6

RUN: RUN2436

DDG (kcal/mol): -0.40

dDDG (kcal/mol): 0.45

MIC-UNK-cdc2493e-12_3



CID: MIC-UNK-cdc2493e-12_3

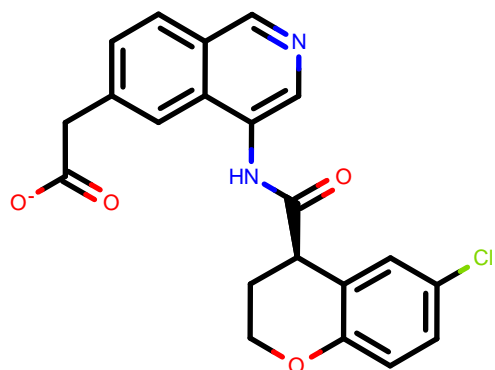
SMILES: C[NH+](C)[C@@H]1CC[C@H](C1)N(c2cccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN552

DDG (kcal/mol): -0.39

dDDG (kcal/mol): 0.28

MAT-POS-b5746674-35_2



CID: MAT-POS-b5746674-35_2

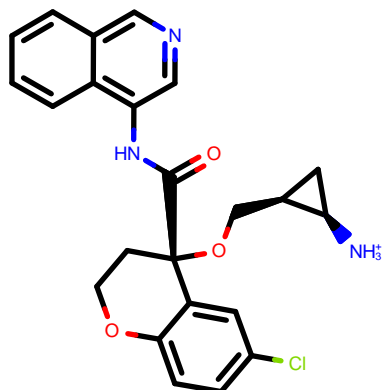
SMILES: Cn1cc(c2cccc2c1=O)NC(=O)NCC[N@H+](Cc3cccc3)C4CCCC4

RUN: RUN70

DDG (kcal/mol): -0.39

dDDG (kcal/mol): 0.23

VLA-UNK-9a7dc93f-4_2



CID: VLA-UNK-9a7dc93f-4_2

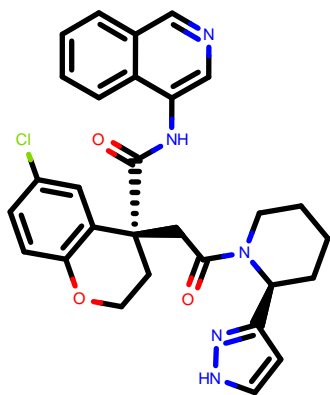
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3CCOC4c3cc(c(c4F)F)C#N

RUN: RUN3087

DDG (kcal/mol): -0.39

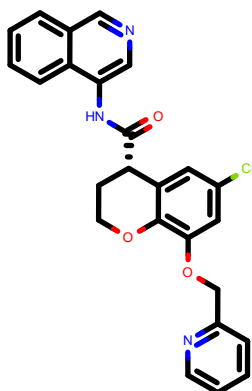
dDDG (kcal/mol): 0.20

BEN-BAS-c2bc0d80-4_1



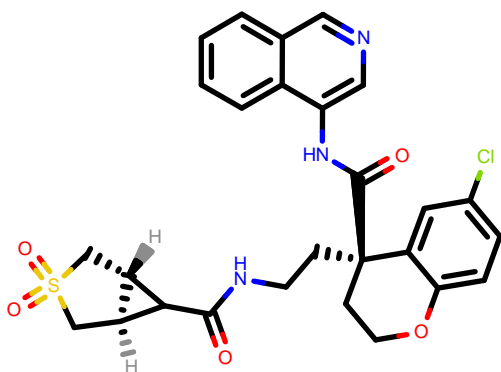
CID:	BEN-BAS-c2bc0d80-4_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CC=Nc4c3[nH]c(cc4=O)Cl</chem>
RUN:	RUN1139
DDG (kcal/mol):	-0.38
dDDG (kcal/mol):	0.32

ED_-GRI-5b13fbe2-17_1



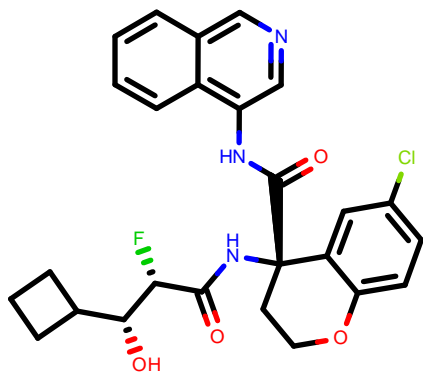
CID:	ED_-GRI-5b13fbe2-17_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)OC[C@H](c5cn(m5)CC6C6)[NH3+]</chem>
RUN:	RUN1549
DDG (kcal/mol):	-0.38
dDDG (kcal/mol):	0.26

MIC-UNK-5a93dd5f-8_2



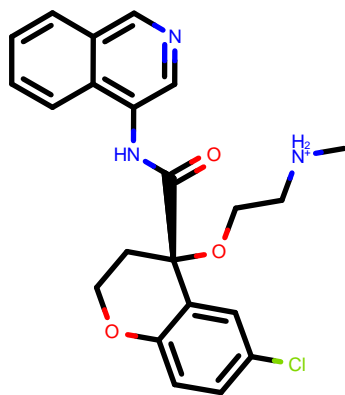
CID:	MIC-UNK-5a93dd5f-8_2
SMILES:	<chem>C[NH+](C)C1CC[NH+](CC1)[C@H](c2ccccc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN775
DDG (kcal/mol):	-0.38
dDDG (kcal/mol):	0.29

RAL-THA-e002e396-5_2



CID:	RAL-THA-e002e396-5_2
SMILES:	<chem>CNS(=O)(=O)C[C@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3462
DDG (kcal/mol):	-0.38
dDDG (kcal/mol):	0.19

DAR-DIA-0f2f46c9-4_3



CID: DAR-DIA-0f2f46c9-4_3

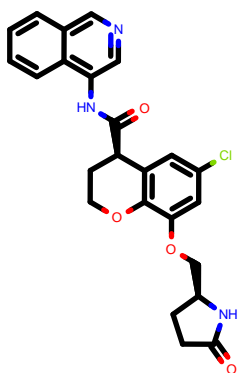
SMILES: CS(=O)(=O)N@@1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3ccc4c3ccc4

RUN: RUN3228

DDG (kcal/mol): -0.38

dDDG (kcal/mol): 0.30

MAT-POS-a3f7f96a-7_4



CID: MAT-POS-a3f7f96a-7_4

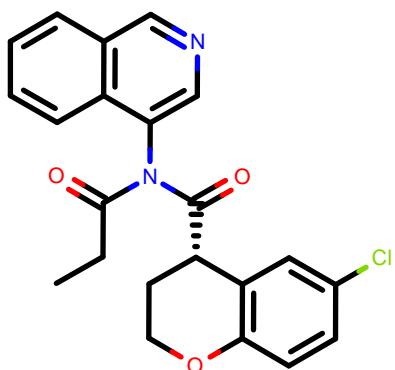
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H]3CCOC4=CC(=O)C(C)C(NH2)C5=CC=C5)C@@H6CCS1=O1=O)C6

RUN: RUN5045

DDG (kcal/mol): -0.38

dDDG (kcal/mol): 0.20

ERI-UCB-ce40166b-1_1



CID: ERI-UCB-ce40166b-1_1

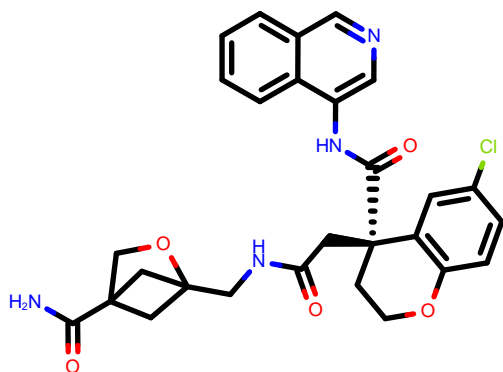
SMILES: c1ccc2c(c1)cnc2CC(=O)Nc3ccc(c3)O[C@@H]4CC(=O)N4

RUN: RUN33

DDG (kcal/mol): -0.38

dDDG (kcal/mol): 0.20

MAK-UNK-c749d764-1_4



CID: MAK-UNK-c749d764-1_4

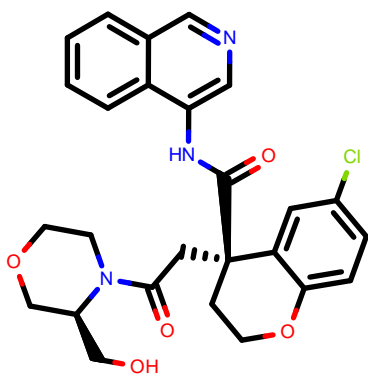
SMILES: c1ccc2c(c1)cnc2NC(=O)C[C@H]3CCC[C@H]([C@@H]3O)C4CC4

RUN: RUN892

DDG (kcal/mol): -0.38

dDDG (kcal/mol): 0.26

MAK-UNK-c749d764-31_3



CID: MAK-UNK-c749d764-31_3

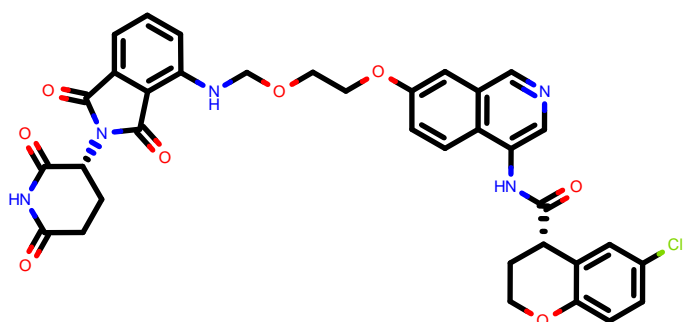
SMILES: CC/C=C(c1cncc2c1cccc2)/C(=O)C[C@@H]3CC[C@H](C[C@@H]3O)C(F)F

RUN: RUN1072

DDG (kcal/mol): -0.37

dDDG (kcal/mol): 0.34

MIC-UNK-9582b2c5-2_8



CID: MIC-UNK-9582b2c5-2_8

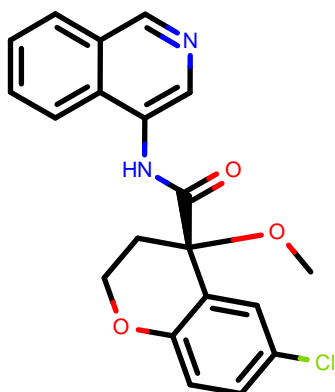
SMILES: CC(=O)N1C[C@H]2[C@H](C1)C[C@H](C(=O)N2c3cccc(c3)Cl)c4cncc5c4cccc5

RUN: RUN268

DDG (kcal/mol): -0.37

dDDG (kcal/mol): 0.36

ALP-UNI-dbbfd3db-13_2



CID: ALP-UNI-dbbfd3db-13_2

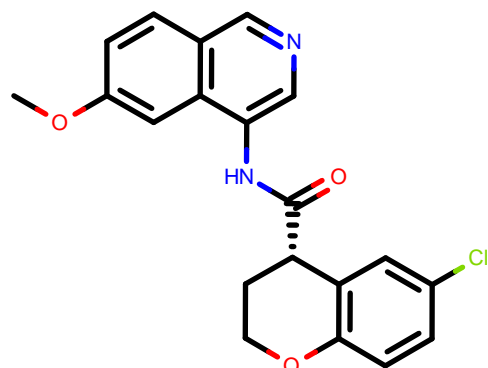
SMILES: Cn1cc(cn1)C2=NO[C@H](C2)C(=O)N[C@@H]3(CCOc4c3cc(cc4)Cl)C(=O)Nc5cncc6c5cccc6

RUN: RUN2782

DDG (kcal/mol): -0.36

dDDG (kcal/mol): 0.07

DAR-DIA-0d514e7d-32_20



CID: DAR-DIA-0d514e7d-32_20

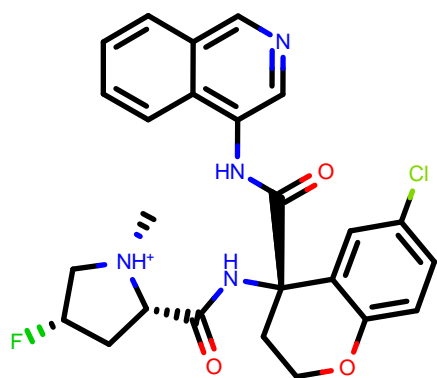
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3[C@H]4C[C@@H]4CO[C@@H]5[C@H]3C=C(C=C5)Cl

RUN: RUN868

DDG (kcal/mol): -0.36

dDDG (kcal/mol): 0.10

DAR-DIA-0f2f46c9-6_1



CID: DAR-DIA-0f2f46c9-6_1

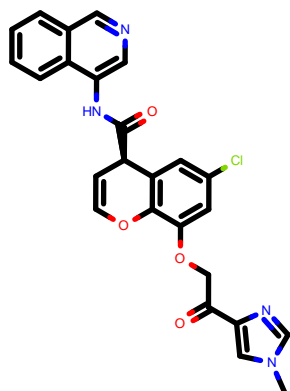
SMILES: CS(=O)(=O)N@@1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cnc4c3cccc4

RUN: RUN3232

DDG (kcal/mol): -0.36

dDDG (kcal/mol): 0.16

MAT-POS-e9e99895-12_2



CID: MAT-POS-e9e99895-12_2

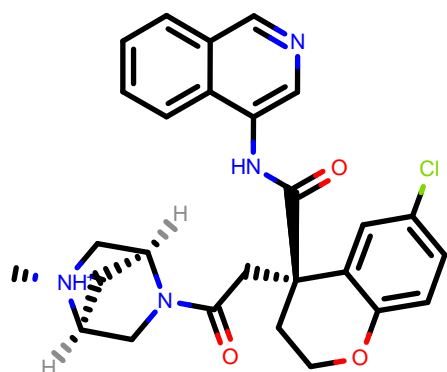
SMILES: C[C@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2cnc3c2cccc3)NC(=O)CC[NH+](C)C

RUN: RUN2266

DDG (kcal/mol): -0.36

dDDG (kcal/mol): 0.40

JOH-SUS-a69c159d-3_1



CID: JOH-SUS-a69c159d-3_1

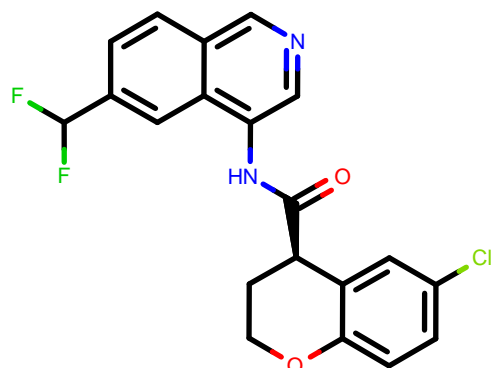
SMILES: c1ccc2c(c1)c(cnc2C(F)F)NC(=O)[C@@H]3CCOC4c3cc(cc4)Cl

RUN: RUN1120

DDG (kcal/mol): -0.36

dDDG (kcal/mol): 0.36

DAR-DIA-ecdbc7dd-10_2



CID: DAR-DIA-ecdbc7dd-10_2

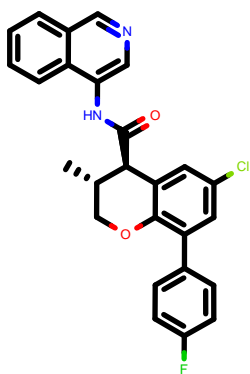
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCNC4c3cc(cc4)Cl)C[NH+]5CCOCC5

RUN: RUN2895

DDG (kcal/mol): -0.36

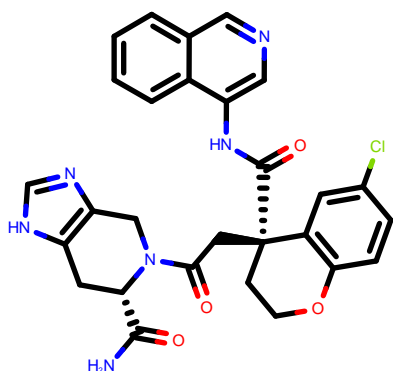
dDDG (kcal/mol): 0.08

LON-WEI-4d77710c-35_2



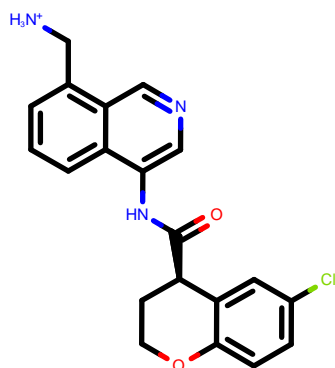
CID:	LON-WEI-4d77710c-35_2
SMILES:	<chem>Cc1ccc(cc1)C[N+](=O)[C@@H](C2)CNC(=O)Nc3cn(c(O)c4c3cccc4)CC(C)C</chem>
RUN:	RUN223
DDG (kcal/mol):	-0.35
dDDG (kcal/mol):	0.31

RAL-THA-4aa06b95-6_1



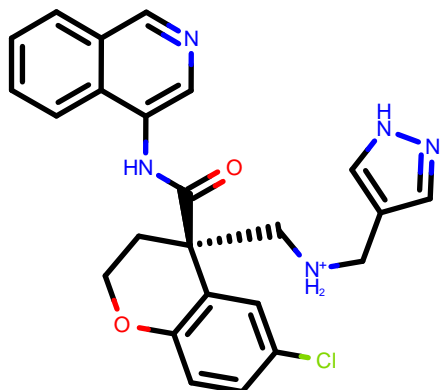
CID:	RAL-THA-4aa06b95-6_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](C3)CCN(c4c3cc(cc4)Cl)CCO</chem>
RUN:	RUN1239
DDG (kcal/mol):	-0.35
dDDG (kcal/mol):	0.25

ERI-UCB-ce40166b-11_1



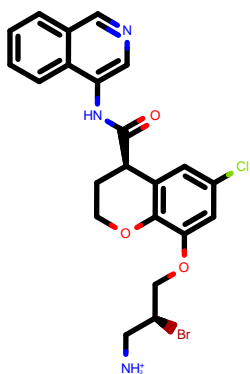
CID:	ERI-UCB-ce40166b-11_1
SMILES:	<chem>c1ccc2c(c1)cncc2CC(=O)Nc3cc(cc(c3)Oc4cccc(O)[nH]4)C#N</chem>
RUN:	RUN50
DDG (kcal/mol):	-0.35
dDDG (kcal/mol):	0.25

KAD-UNI-8a629cb0-34_1



CID:	KAD-UNI-8a629cb0-34_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H](C3)CCO4c3cc(cc4)CNC(=O)[C@@H](C5)C#N</chem>
RUN:	RUN2116
DDG (kcal/mol):	-0.35
dDDG (kcal/mol):	0.37

MAT-POS-24589f88-6_2



CID: MAT-POS-24589f88-6_2

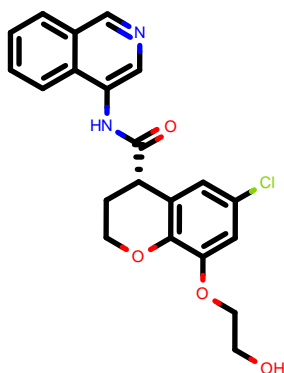
SMILES: c1ccc2c(c1)ncnc2N(C(=O)C[C@H]3CCOCc4cc(Cl)C[C@H]3C)C(=O)NCCBr

RUN: RUN4777

DDG (kcal/mol): -0.35

dDDG (kcal/mol): 0.31

VLA-UCB-05e51b3f-9_1



CID: VLA-UCB-05e51b3f-9_1

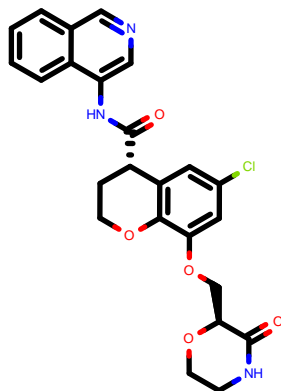
SMILES: c1ccc2c(c1)ncnc2N(C(=O)C[C@H]3CCOCc4cc(Cl)C[C@H]3C)C(=O)CCO

RUN: RUN317

DDG (kcal/mol): -0.34

dDDG (kcal/mol): 0.21

ALP-POS-5bb456a5-2_7



CID: ALP-POS-5bb456a5-2_7

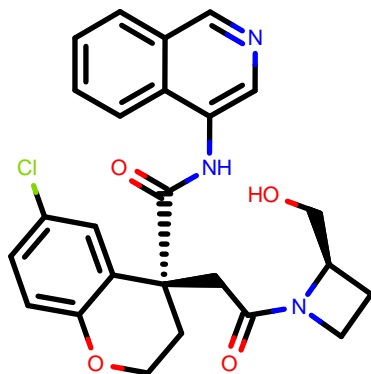
SMILES: C1=CN(C(=O)C[C@H]2CCOCc3cc(Cl)C[C@H]2C)C(=O)NCC(=O)N1

RUN: RUN2428

DDG (kcal/mol): -0.34

dDDG (kcal/mol): 0.28

MIC-UNK-67d4a29a-1_1



CID: MIC-UNK-67d4a29a-1_1

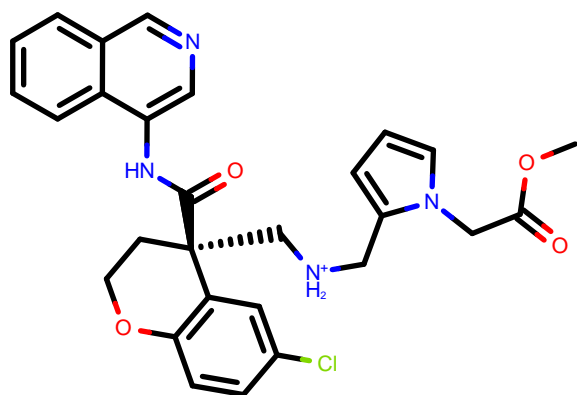
SMILES: CN1C(=O)N(C(=O)C[C@H]2CCOCc3cc(Cl)C[C@H]2C)C1O

RUN: RUN1086

DDG (kcal/mol): -0.34

dDDG (kcal/mol): 0.31

ED_-GRI-5b13fbe2-60_2



CID: ED_-GRI-5b13fbe2-60_2

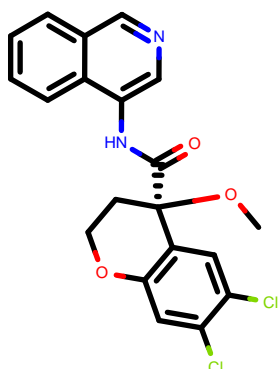
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)C)OCCn5cc6c(n5)CC[C@@H](C6)[NH3+]

RUN: RUN1607

DDG (kcal/mol): -0.34

dDDG (kcal/mol): 0.50

DAR-DIA-5ff57136-17_1



CID: DAR-DIA-5ff57136-17_1

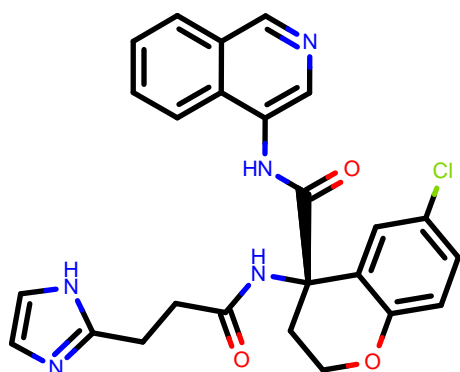
SMILES: c1ccc2c(c1)cncc2N(CC#CBr)C(=O)[C@@]3(CCOc4c3cc(cc4)Cl

RUN: RUN1390

DDG (kcal/mol): -0.34

dDDG (kcal/mol): 0.13

NAU-LAT-a5c7d7cb-8_1



CID: NAU-LAT-a5c7d7cb-8_1

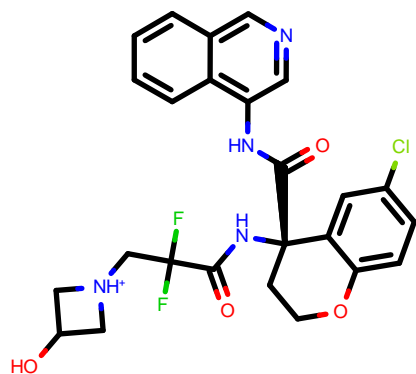
SMILES: c1ccc2c(c1)cncc2CNC(=O)COc3cccc(c3)Cl

RUN: RUN578

DDG (kcal/mol): -0.34

dDDG (kcal/mol): 0.32

EDJ-MED-946e547c-1_1



CID: EDJ-MED-946e547c-1_1

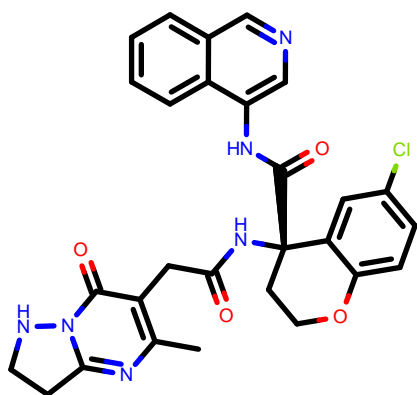
SMILES: CO[C@@]1(CCNc2c1cc(c2)F)C(=O)Nc3ncc4c3cc(cc4)S(=O)(=O)C

RUN: RUN3316

DDG (kcal/mol): -0.33

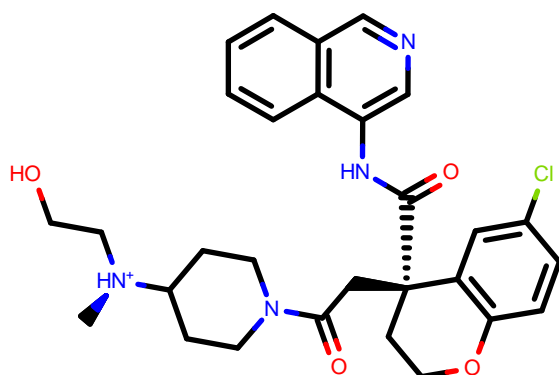
dDDG (kcal/mol): 0.17

RAL-THA-c11c1343-1_1



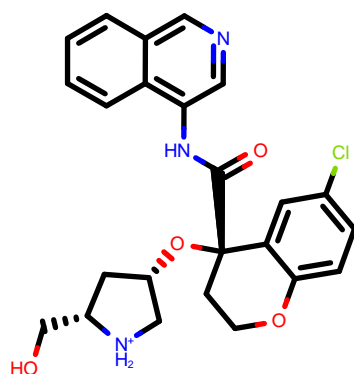
CID:	RAL-THA-c11c1343-1_1
SMILES:	<chem>CC(=O)N1CC[C@@H](c2c1ccc(c2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN1227
DDG (kcal/mol):	-0.33
dDDG (kcal/mol):	0.29

BEN-DND-f2e727cd-4_2



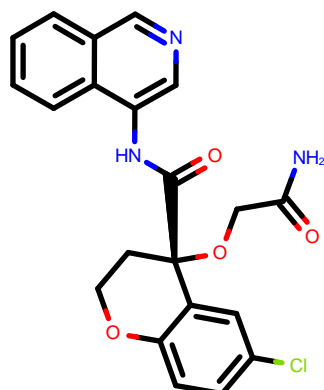
CID:	BEN-DND-f2e727cd-4_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3COCc4c3cc(cc4)Cl</chem>
RUN:	RUN1194
DDG (kcal/mol):	-0.33
dDDG (kcal/mol):	0.48

DAR-DIA-0f2f46c9-2_2



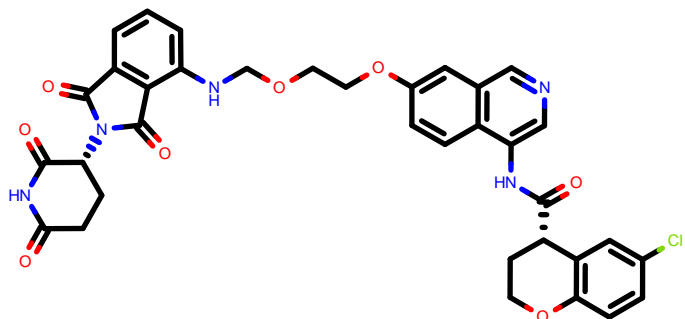
CID:	DAR-DIA-0f2f46c9-2_2
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CC[N@](c4c3cc(cc4)Cl)S(=O)(=O)N</chem>
RUN:	RUN3223
DDG (kcal/mol):	-0.33
dDDG (kcal/mol):	0.23

ALP-POS-7c6e02c7-3_1



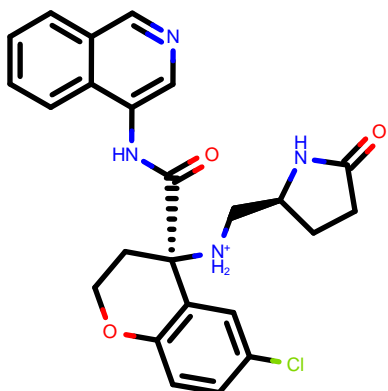
CID:	ALP-POS-7c6e02c7-3_1
SMILES:	<chem>CN(C)c1ccc(nc1)N(Cc2ccc(c(c2)Cl)F)C(=O)Cc3cncc4c3cccc4</chem>
RUN:	RUN3100
DDG (kcal/mol):	-0.33
dDDG (kcal/mol):	0.11

DAR-DIA-23e5a6a0-2_1



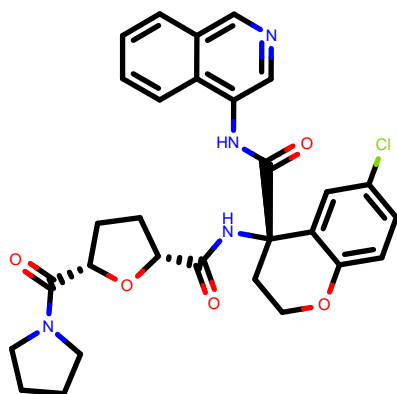
CID:	DAR-DIA-23e5a6a0-2_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCOCc4cc3cc(c4O)[C@@H]5CC6([NH2+][5]COC6)Cl</chem>
RUN:	RUN403
DDG (kcal/mol):	-0.33
dDDG (kcal/mol):	0.42

ED_-GRI-5b13fbe2-7_2



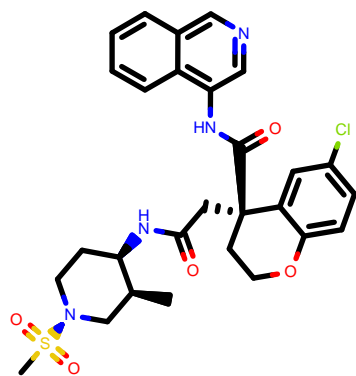
CID:	ED_-GRI-5b13fbe2-7_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(c4)O)C5c6n(n5)C[C@@H]6C[NH2+][COC6]</chem>
RUN:	RUN1532
DDG (kcal/mol):	-0.33
dDDG (kcal/mol):	0.30

MAT-POS-4223bc15-9_9



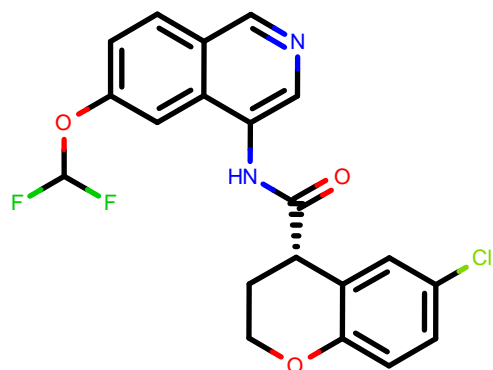
CID:	MAT-POS-4223bc15-9_9
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3C[N@@]4(Cc4c3cc(c4)Cl)S(=O)(=O)[C@@H]5CC[C@@H]5O</chem>
RUN:	RUN4017
DDG (kcal/mol):	-0.32
dDDG (kcal/mol):	0.18

MAT-POS-4223bc15-9_2



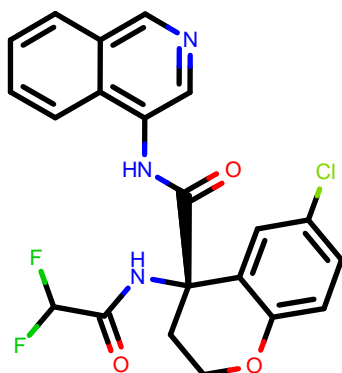
CID:	MAT-POS-4223bc15-9_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3C[N@@]4(Cc4c3cc(c4)Cl)S(=O)(=O)[C@@H]5CC[C@@H]5O</chem>
RUN:	RUN4010
DDG (kcal/mol):	-0.32
dDDG (kcal/mol):	0.25

LON-WEI-4d77710c-20_1



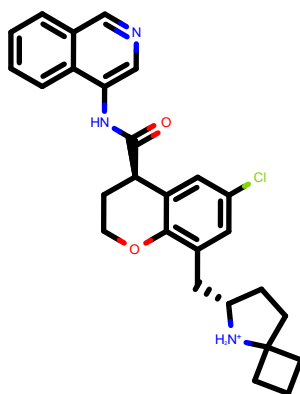
CID:	LON-WEI-4d77710c-20_1
SMILES:	<chem>CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)NCCc3c[nH]c4c3ccccc4</chem>
RUN:	RUN208
DDG (kcal/mol):	-0.31
dDDG (kcal/mol):	0.11

EDJ-MED-f893e2a1-6_3



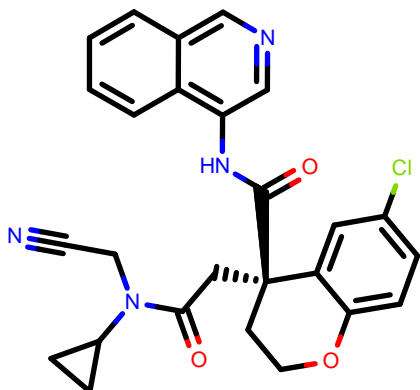
CID:	EDJ-MED-f893e2a1-6_3
SMILES:	<chem>c1ccc2c(c1)ncoc2NC(=O)[C@@]3(C)C(CO)C4c3cc(cc4)C]C[NH2+]C[C@H]5CC(=O)NC5</chem>
RUN:	RUN3205
DDG (kcal/mol):	-0.31
dDDG (kcal/mol):	0.08

ERI-UCB-ce40166b-3_1



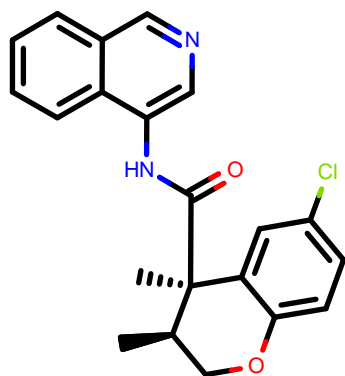
CID:	ERI-UCB-ce40166b-3_1
SMILES:	<chem>c1ccc2c(c1)ncoc2CC(=O)Nc3ccccc(c3)Oc4cccc(=O)[nH]4</chem>
RUN:	RUN38
DDG (kcal/mol):	-0.31
dDDG (kcal/mol):	0.38

KAD-UNI-b13decd3-1_4



CID:	KAD-UNI-b13decd3-1_4
SMILES:	<chem>Cn1cc(en1)N2CC[C@H](C2)C[NH2+]C[C@]3(C)C(CO)C4c3cc(cc4)C]C(=O)Nc5ccc6c5ccccc6</chem>
RUN:	RUN3782
DDG (kcal/mol):	-0.31
dDDG (kcal/mol):	0.12

EDJ-MED-ee07cf00-6_4



CID: EDJ-MED-ee07cf00-6_4

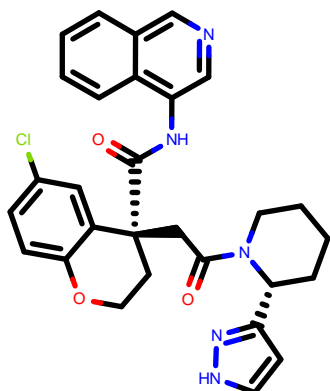
SMILES: Cn1cc(en1)C2=NO[C@H](C2)C(=O)N[C@H](c3ccccc3)C(C=O)Nc4nccc5c4ccccc5

RUN: RUN2819

DDG (kcal/mol): -0.31

dDDG (kcal/mol): 0.07

KAD-UNI-b13decd3-10_1



CID: KAD-UNI-b13decd3-10_1

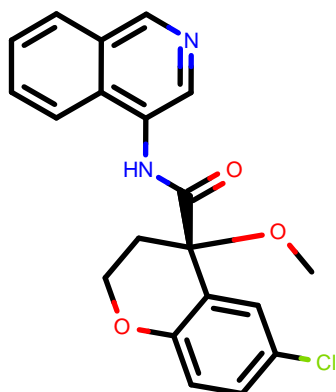
SMILES: COc1nc2ncc(en2n1)C[NH2+]C[C@]3(CCOc4c3cc(cc4)C)C(=O)Nc5nccc6c5ccccc6

RUN: RUN3791

DDG (kcal/mol): -0.31

dDDG (kcal/mol): 0.15

MAT-POS-932d1078-3_1



CID: MAT-POS-932d1078-3_1

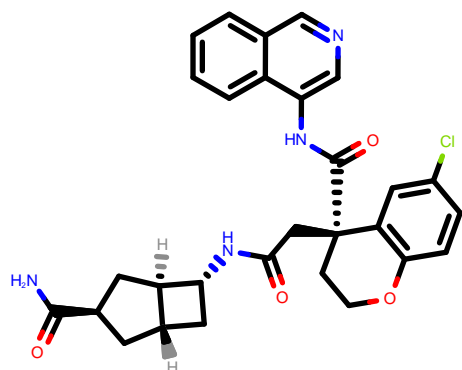
SMILES: CO[C@]1(CCOc2c1cc(cc2F)F)C(=O)Nc3nccc4c3ccccc4

RUN: RUN3595

DDG (kcal/mol): -0.30

dDDG (kcal/mol): 0.06

ED_-GRI-5b13fbe2-58_2



CID: ED_-GRI-5b13fbe2-58_2

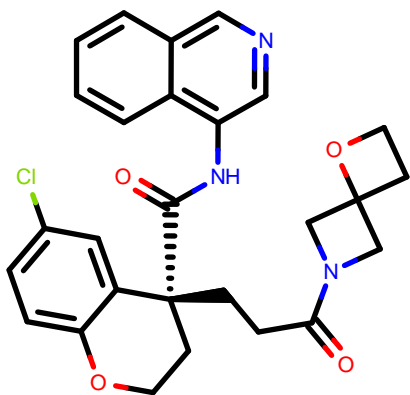
SMILES: c1ccc2c(c1)encc2NC(=O)[C@@]3(CCOc4c3cc(cc4)C)OCCn5cc(m5)[C@H]6CCCC[NH2+]6

RUN: RUN1608

DDG (kcal/mol): -0.30

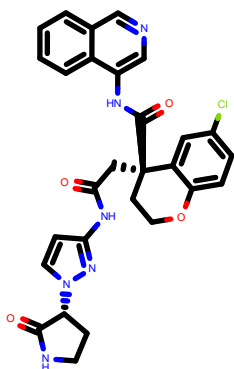
dDDG (kcal/mol): 0.27

DAR-DIA-f6ee7aeb-4_2



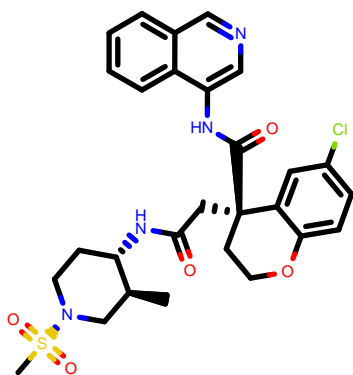
CID:	DAR-DIA-f6ee7aeb-4_2
SMILES:	<chem>c1ccc2c(c1)ncnc2N3C[C@@H](C[C@@H](CC3=O)4ccc5c(c4)OCC(C(F)F)C5)H[C@@H](O)[H]5=O</chem>
RUN:	RUN3413
DDG (kcal/mol):	-0.30
dDDG (kcal/mol):	0.16

MAK-UNK-c749d764-16_15



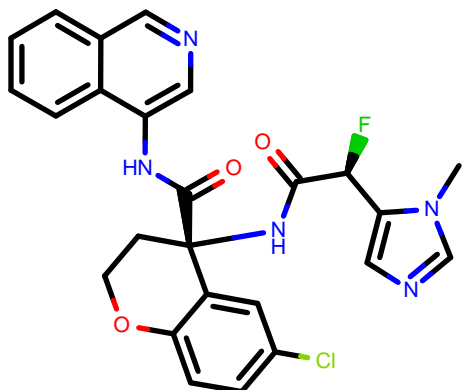
CID:	MAK-UNK-c749d764-16_15
SMILES:	<chem>C[C@@H](N)(c1ncnc2c1ccc2)C(=O)C[C@H]3CCC[C@H]([C@H]3O)C(F)F</chem>
RUN:	RUN988
DDG (kcal/mol):	-0.30
dDDG (kcal/mol):	0.47

MAK-UNK-c749d764-1_6



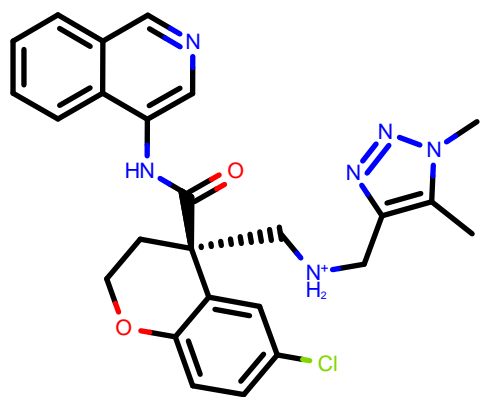
CID:	MAK-UNK-c749d764-1_6
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)C[C@H]3CCC[C@@H]([C@H]3O)C4CC4</chem>
RUN:	RUN894
DDG (kcal/mol):	-0.29
dDDG (kcal/mol):	0.34

MAK-UNK-8be7dca9-10_4



CID:	MAK-UNK-8be7dca9-10_4
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@H]3C[C@H](Oc4c3cc(cc4)C)C(=O)[O-]</chem>
RUN:	RUN514
DDG (kcal/mol):	-0.28
dDDG (kcal/mol):	0.29

EDG-MED-90036822-96_2



CID: EDG-MED-90036822-96_2

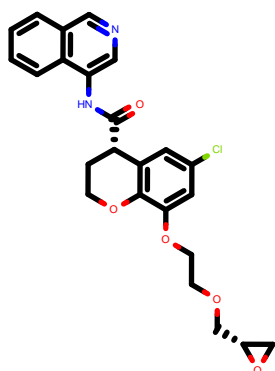
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4C)NC(=O)CC[C@H](CCF)N)H3+

RUN: RUN1824

DDG (kcal/mol): -0.28

dDDG (kcal/mol): 0.36

MAT-POS-f9802937-6_1



CID: MAT-POS-f9802937-6_1

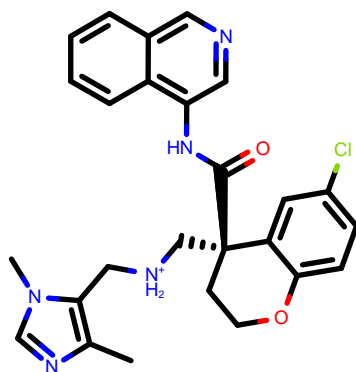
SMILES: CNC(=O)C[C@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cnc4c3ccccc4

RUN: RUN2399

DDG (kcal/mol): -0.27

dDDG (kcal/mol): 0.29

EDG-MED-90036822-106_1



CID: EDG-MED-90036822-106_1

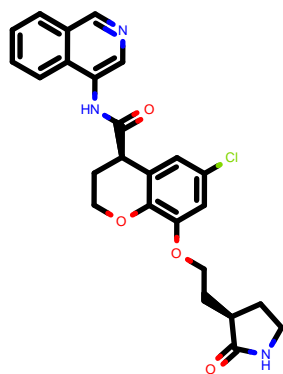
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4C)C)NC(=O)CC(C[NH3+])F

RUN: RUN1843

DDG (kcal/mol): -0.27

dDDG (kcal/mol): 0.50

KAD-UNI-8a629cb0-47_1



CID: KAD-UNI-8a629cb0-47_1

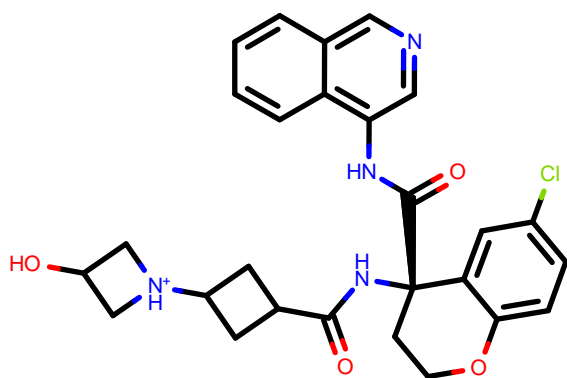
SMILES: Cn1cc(nn1)CC(=O)NCC[C@]2(CCOc3c2cc(cc3)Cl)C(=O)Nc4cnc5c4ccccc5

RUN: RUN2139

DDG (kcal/mol): -0.27

dDDG (kcal/mol): 0.45

VLA-UNK-c65c1026-5_1



CID: VLA-UNK-c65c1026-5_1

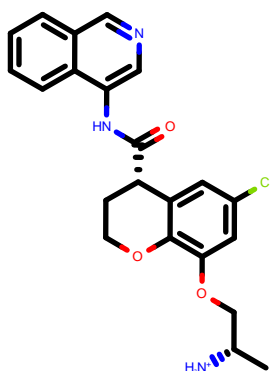
SMILES: c1ccc2c(c1)cncc2N3C(=O)C[NH2+][C@@]4(C3=O)COc5c4cc(cc5)Cl

RUN: RUN3186

DDG (kcal/mol): -0.26

dDDG (kcal/mol): 0.18

MAT-POS-e9e99895-7_2



CID: MAT-POS-e9e99895-7_2

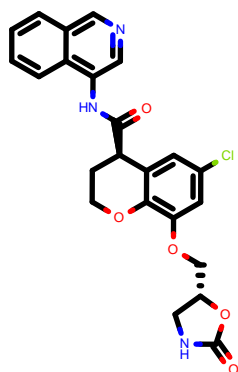
SMILES: C[C@](c1ccc(cc1)Cl)(C(=O)Nc2ccc3c2cccc3)NC(=O)Cc4ccc(cc4)n5cnm5

RUN: RUN2254

DDG (kcal/mol): -0.26

dDDG (kcal/mol): 0.29

MIC-UNK-06e5f114-2_1



CID: MIC-UNK-06e5f114-2_1

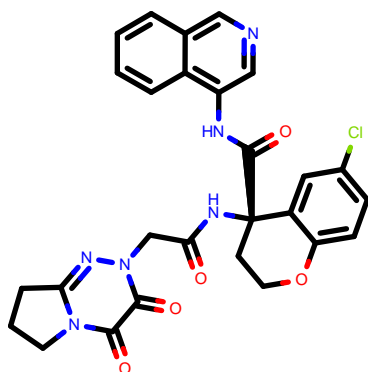
SMILES: CS(=O)(=O)c1cccc(c1)NCc2ccc(cc2CC(=O)Nc3nccc4c3cccc4)Cl

RUN: RUN5061

DDG (kcal/mol): -0.26

dDDG (kcal/mol): 0.19

DAR-DIA-076fb6ea-1_1



CID: DAR-DIA-076fb6ea-1_1

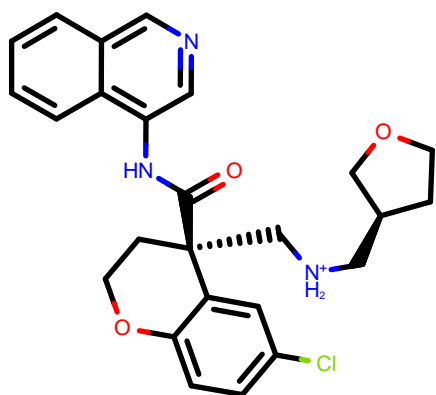
SMILES: C=CC(=O)N(c1cncc2c1cccc2)C(=O)Cc3cccc(c3)Cl

RUN: RUN1397

DDG (kcal/mol): -0.26

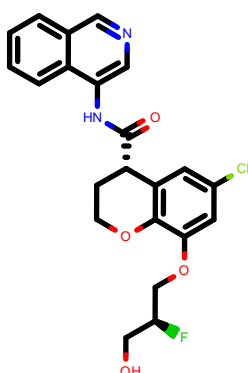
dDDG (kcal/mol): 0.40

PET-UNK-9b23ef84-7_1



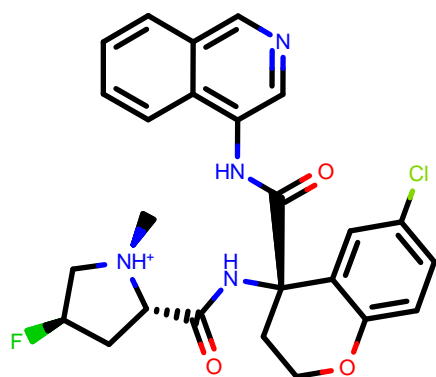
CID:	PET-UNK-9b23ef84-7_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CN(Cc4c3cc(c(c4)F)Cl)c5nnc5</chem>
RUN:	RUN4433
DDG (kcal/mol):	-0.26
dDDG (kcal/mol):	0.21

MAT-POS-1f3f1a6f-2_2



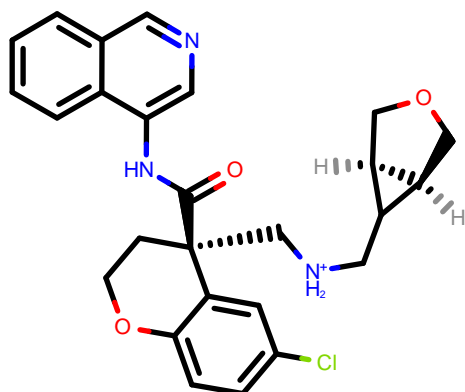
CID:	MAT-POS-1f3f1a6f-2_2
SMILES:	<chem>C[NH+](C)[C@H](c1ccc(c(c1)Cl)Cl)C(=O)Nc2cnc3c2cccc3</chem>
RUN:	RUN2278
DDG (kcal/mol):	-0.25
dDDG (kcal/mol):	0.27

DAR-DIA-0f2f46c9-7_3



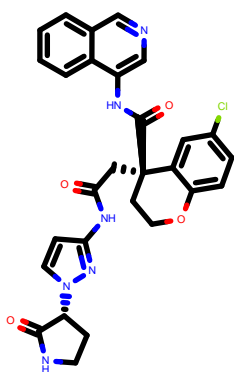
CID:	DAR-DIA-0f2f46c9-7_3
SMILES:	<chem>CNS(=O)(=O)[N@@]1CC[C@H](c2c1ccc(c2)Cl)C(=O)Nc3cnc4c3cccc4</chem>
RUN:	RUN3236
DDG (kcal/mol):	-0.25
dDDG (kcal/mol):	0.16

EDJ-MED-92e193ae-2_1



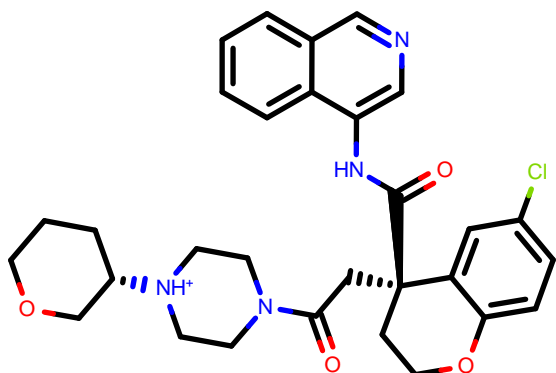
CID:	EDJ-MED-92e193ae-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3CCNc4c3cc(cc4)Cl</chem>
RUN:	RUN1853
DDG (kcal/mol):	-0.24
dDDG (kcal/mol):	0.39

ED_-GRI-5b13fbe2-40_1



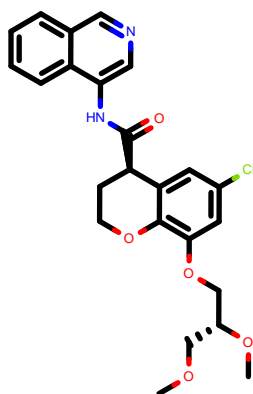
CID:	ED_-GRI-5b13fbe2-40_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OC[C@H](C[NH3+])F</chem>
RUN:	RUN1574
DDG (kcal/mol):	-0.24
dDDG (kcal/mol):	0.42

BEN-DND-c852c98b-2_2



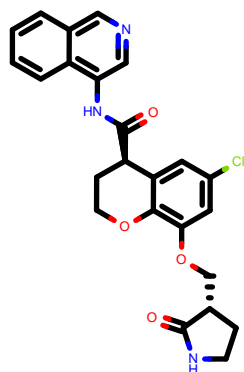
CID:	BEN-DND-c852c98b-2_2
SMILES:	<chem>c1cc2cnc(c2cc1OC(F)(F)F)NC(=O)[C@H]3CCOCc4c3cc(cc4)Cl</chem>
RUN:	RUN1206
DDG (kcal/mol):	-0.24
dDDG (kcal/mol):	0.37

EDJ-MED-9e38fd34-5_1



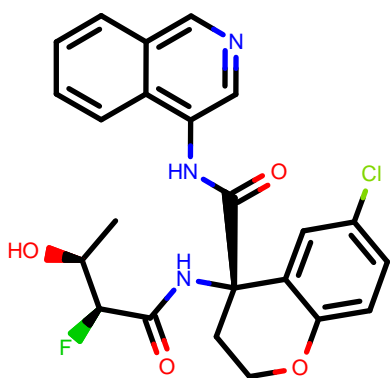
CID:	EDJ-MED-9e38fd34-5_1
SMILES:	<chem>C[C@]1(c2cc(c(cc2NC1=O)Cl)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN2351
DDG (kcal/mol):	-0.24
dDDG (kcal/mol):	0.40

MAT-POS-8293a91a-4_2



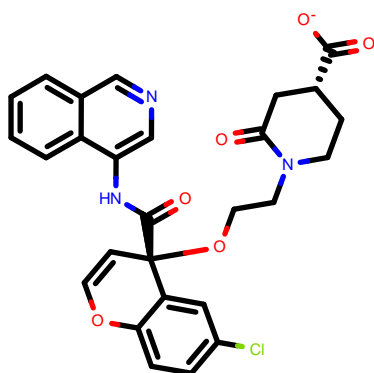
CID:	MAT-POS-8293a91a-4_2
SMILES:	<chem>CNC(=O)C[N@H+]1Cc2ccc(cc2[C@@H](C1)C(=O)Nc3cncc4c3cccc4)Cl</chem>
RUN:	RUN5083
DDG (kcal/mol):	-0.24
dDDG (kcal/mol):	0.19

MAT-POS-fb82b63d-4_2



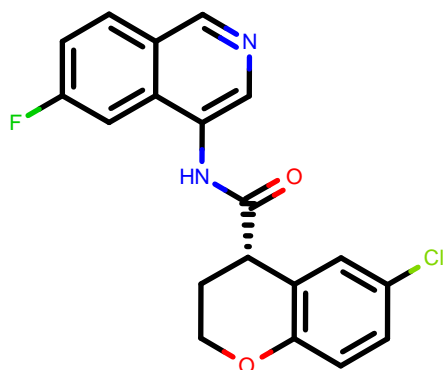
CID:	MAT-POS-fb82b63d-4_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@H]3c4cc(ccc4CC[N+](H+)(3)CC5CC5)Cl</chem>
RUN:	RUN3181
DDG (kcal/mol):	-0.23
dDDG (kcal/mol):	0.14

MIC-UNK-cdc2493e-12_4



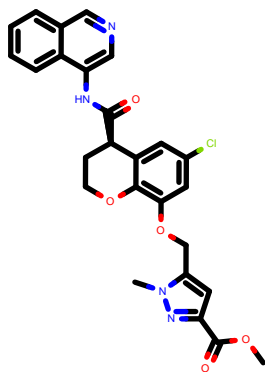
CID:	MIC-UNK-cdc2493e-12_4
SMILES:	<chem>C[NH+](C)[C@H]1CC[C@H](C1)N(c2cccc(c2)Cl)C(=O)Nc3ncc4c3cccc4</chem>
RUN:	RUN553
DDG (kcal/mol):	-0.22
dDDG (kcal/mol):	0.34

MAK-UNK-c749d764-15_6



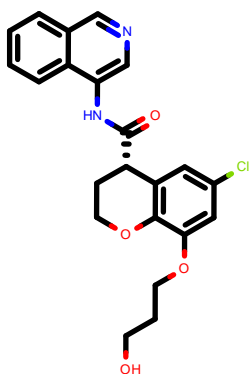
CID:	MAK-UNK-c749d764-15_6
SMILES:	<chem>C[C@H](N)(c1cnc2c1cccc2)C(=O)C[C@@H]3CC[C@H](C3)C(F)F)OCC4CCccc4</chem>
RUN:	RUN963
DDG (kcal/mol):	-0.22
dDDG (kcal/mol):	0.10

EDJ-MED-d203f206-36_1



CID:	EDJ-MED-d203f206-36_1
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CC[C@@](C5)(C#N)C(=O)N</chem>
RUN:	RUN2598
DDG (kcal/mol):	-0.22
dDDG (kcal/mol):	0.39

DAR-DIA-23e5a6a0-9_2



CID: DAR-DIA-23e5a6a0-9_2

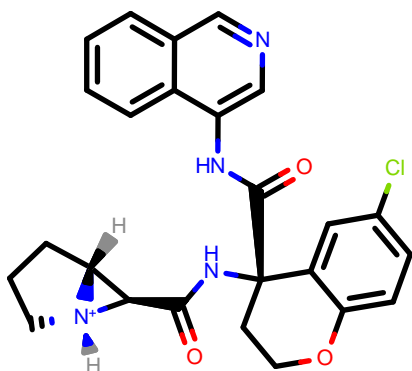
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4cc(Cc5c6(Cc6)C[NH2+]5)Cl

RUN: RUN419

DDG (kcal/mol): -0.22

dDDG (kcal/mol): 0.23

MIC-UNK-cdc2493e-16_1



CID: MIC-UNK-cdc2493e-16_1

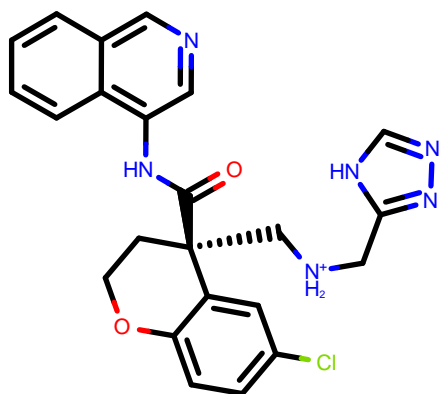
SMILES: c1ccc2c(c1)cncc2NC(=O)N(CCC3CCCC3)c4cccc(c4)Cl

RUN: RUN561

DDG (kcal/mol): -0.22

dDDG (kcal/mol): 0.29

EDJ-MED-139368ae-5_2



CID: EDJ-MED-139368ae-5_2

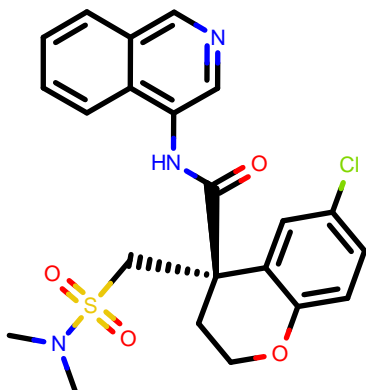
SMILES: C[NH+]1CCN(CC1)S(=O)(=O)[N@]2Cc3cccc3[C@@H](C2)C(=O)Nc4ncc5c4cccc5

RUN: RUN4549

DDG (kcal/mol): -0.21

dDDG (kcal/mol): 0.22

JOH-UNI-21fd6073-1_2



CID: JOH-UNI-21fd6073-1_2

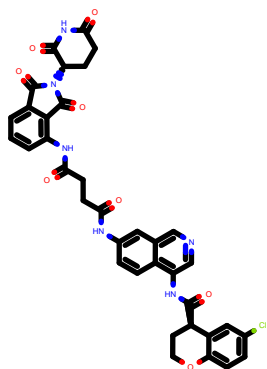
SMILES: c1ccc(cc1)S(=O)(=O)N(CC#N)C(=O)CN(c2ncc3c2cccc3)C(=O)[C@H]4COc5c4cc(oc5)Cl

RUN: RUN4183

DDG (kcal/mol): -0.21

dDDG (kcal/mol): 0.12

JAG-UCB-f37eaa14-4_2



CID: JAG-UCB-f37eaa14-4_2

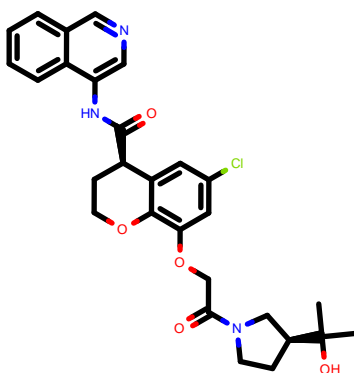
SMILES: c1ccc2c(c1)cncc2N3CC[C@]4(C3=O)C[N@]5c5c4cc(cc5Cl)CC[NH2+][C6CC6

RUN: RUN3061

DDG (kcal/mol): -0.21

dDDG (kcal/mol): 0.22

ALF-EVA-07677224-1_3



CID: ALF-EVA-07677224-1_3

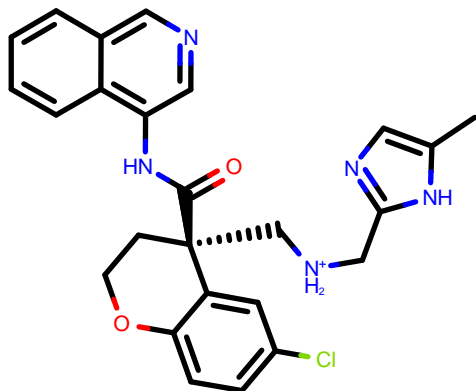
SMILES: Cc1nc(ns1)S(=O)(=O)[N@]2Cc3ccc(cc3[C@H](C2)C(=O)Nc4cccc5c4cccc5)Cl

RUN: RUN4898

DDG (kcal/mol): -0.21

dDDG (kcal/mol): 0.20

MAT-POS-dc2604c4-3_1



CID: MAT-POS-dc2604c4-3_1

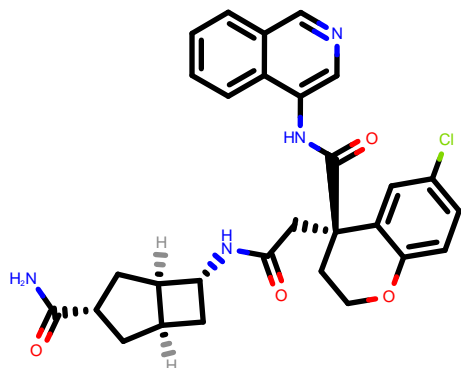
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3[C@H]([C@]4(C3)C(=O)S(=O)(=O)CC5(CNC5)C#N

RUN: RUN4711

DDG (kcal/mol): -0.21

dDDG (kcal/mol): 0.18

RAL-THA-8416115c-13_2



CID: RAL-THA-8416115c-13_2

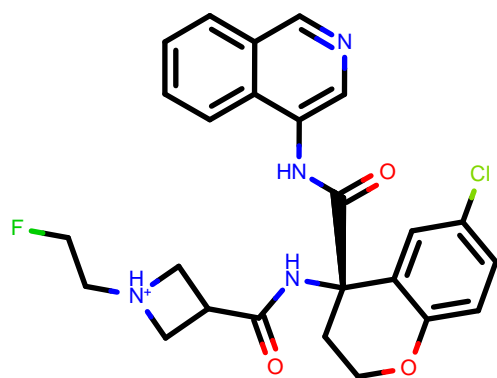
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3CCN(c4c3cc(cc4Cl)C)c5[nH]ncn5

RUN: RUN1296

DDG (kcal/mol): -0.21

dDDG (kcal/mol): 0.46

DAR-DIA-0d514e7d-17_1



CID: DAR-DIA-0d514e7d-17_1

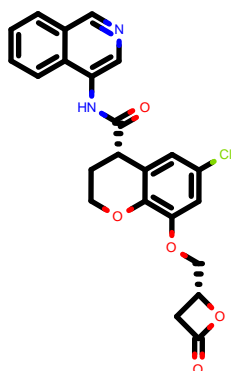
SMILES: C[C@H]1COc2c(cc(cc2[C@@H]1C(=O)Nc3encc4c3cccc4Cl)c5cc(cc5)F)F

RUN: RUN820

DDG (kcal/mol): -0.20

dDDG (kcal/mol): 0.40

DAR-DIA-6be260fc-4_1



CID: DAR-DIA-6be260fc-4_1

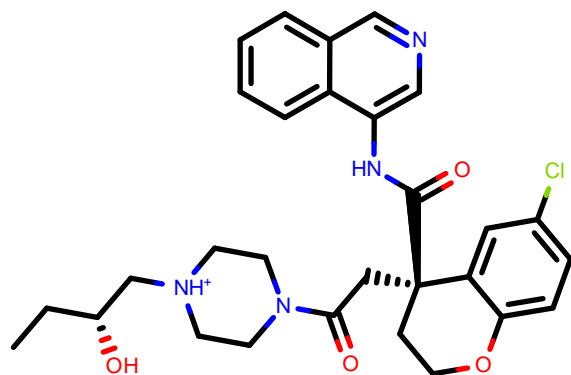
SMILES: CC(C)[C@H]1CN(C(=O)[C@@]12CNc3c2cc(cc3Cl)c4cncc5c4cccc5

RUN: RUN2135

DDG (kcal/mol): -0.20

dDDG (kcal/mol): 0.27

MIC-UNK-5a93dd5f-3_12



CID: MIC-UNK-5a93dd5f-3_12

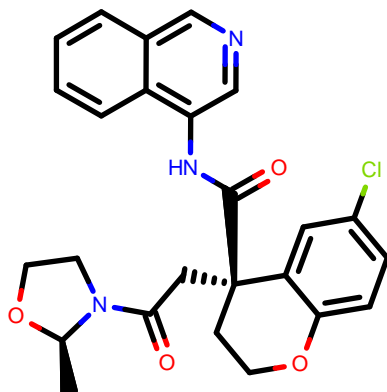
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H](c3cccc(c3)Cl)[N@H]4CC[C@H]5CCCC[C@H]5C4

RUN: RUN751

DDG (kcal/mol): -0.19

dDDG (kcal/mol): 0.49

KAD-UNI-b13decd3-1_2



CID: KAD-UNI-b13decd3-1_2

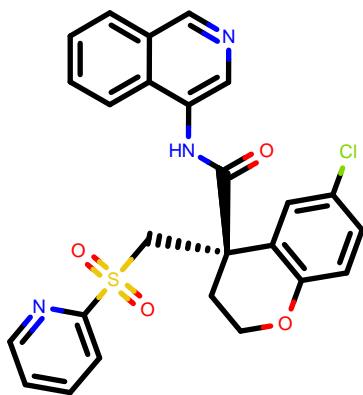
SMILES: Cn1cc(en1)N2CC[C@@H](C2)C[NH2+]C[C@@]3(CCOc4c3cc(cc4Cl)C(=O)Nc5encc6c5cccc6

RUN: RUN3779

DDG (kcal/mol): -0.19

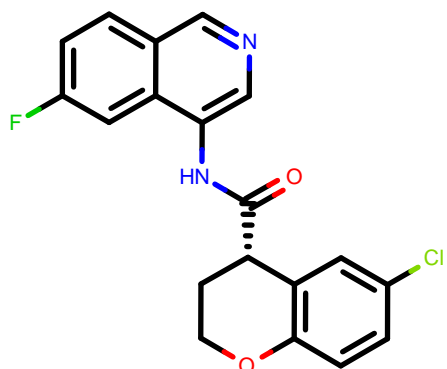
dDDG (kcal/mol): 0.10

PET-UNK-12d8d43f-1_1



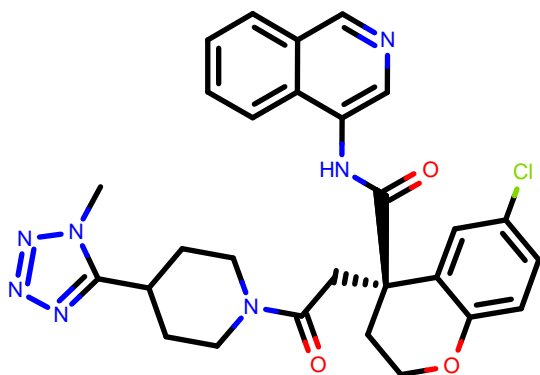
CID:	PET-UNK-12d8d43f-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2N(C=O)C(=O)Cc3ccccc(c3)Cl</chem>
RUN:	RUN1495
DDG (kcal/mol):	-0.19
dDDG (kcal/mol):	0.32

MIC-UNK-9582b2c5-3_4



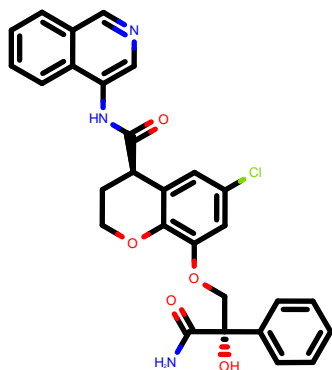
CID:	MIC-UNK-9582b2c5-3_4
SMILES:	<chem>CC(=O)N1CC[C@H]2[C@H](C1)CN(C(=O)N2c3ccccc(c3)Cl)c4cncc5c4cccc5</chem>
RUN:	RUN272
DDG (kcal/mol):	-0.19
dDDG (kcal/mol):	0.11

MIC-UNK-5a93dd5f-5_8



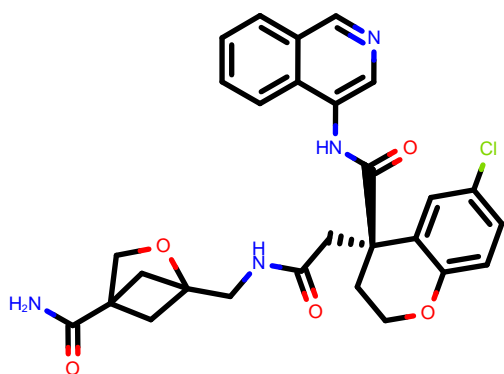
CID:	MIC-UNK-5a93dd5f-5_8
SMILES:	<chem>CC(=O)N[C@H]1CC[N@H](C1)[C@H](C2c3ccccc(c2)Cl)C(=O)Nc3cnnc4c3cccc4</chem>
RUN:	RUN764
DDG (kcal/mol):	-0.18
dDDG (kcal/mol):	0.32

MAT-POS-e9e99895-3_2



CID:	MAT-POS-e9e99895-3_2
SMILES:	<chem>C[C@](c1ccc(c(c1)Cl)Cl)(C(=O)Nc2ncc3c2cccc3)NC(=O)[C@@H]4CCC(=O)NC4</chem>
RUN:	RUN2244
DDG (kcal/mol):	-0.18
dDDG (kcal/mol):	0.46

MIC-UNK-5a93dd5f-5_7



CID: MIC-UNK-5a93dd5f-5_7

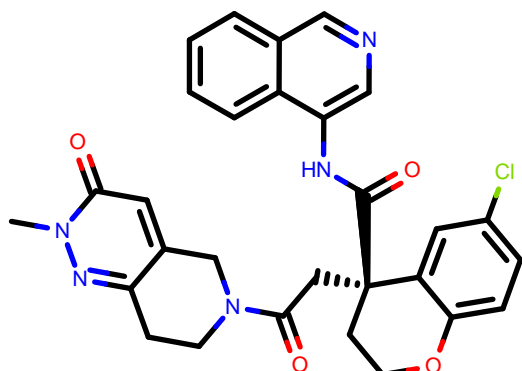
SMILES: CC(=O)N[C@@H]1CC[N@H+](C1)[C@H](c2ccoc(c2)Cl)C(=O)Nc3ccc4c3ccoc4

RUN: RUN765

DDG (kcal/mol): -0.18

dDDG (kcal/mol): 0.31

MAK-UNK-c749d764-18_2



CID: MAK-UNK-c749d764-18_2

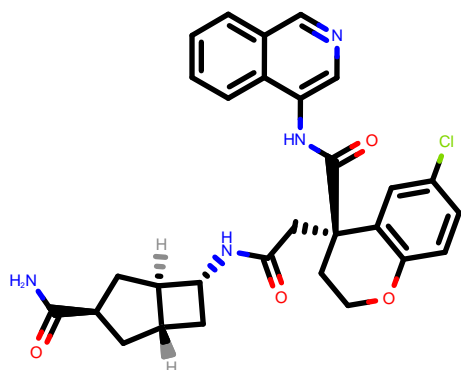
SMILES: CC(C)SCN(c1cnc2c1cccc2)C(=O)C[C@H]3CC[C@@H](C3)[C@@H](C3O)C(F)F

RUN: RUN991

DDG (kcal/mol): -0.18

dDDG (kcal/mol): 0.31

NAU-LAT-4ce8bf23-3_1



CID: NAU-LAT-4ce8bf23-3_1

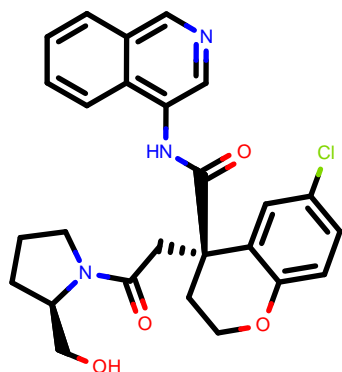
SMILES: CC(=O)N[C@@H](c1cnc2c1cccc2)C(=O)Nc3ccccc(c3)Cl

RUN: RUN1393

DDG (kcal/mol): -0.17

dDDG (kcal/mol): 0.38

DAR-DIA-4987d2cd-4_4



CID: DAR-DIA-4987d2cd-4_4

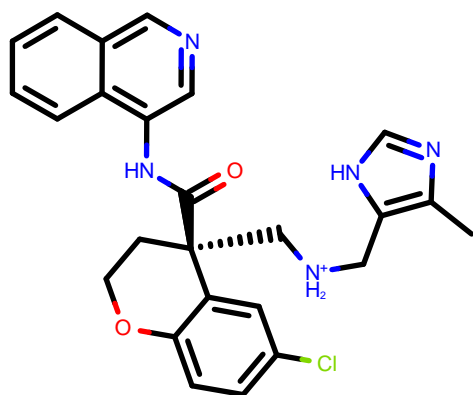
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@H]3c4cc(ccc4CN3c5c(c(=O)c5=O)[O-])Cl

RUN: RUN3817

DDG (kcal/mol): -0.17

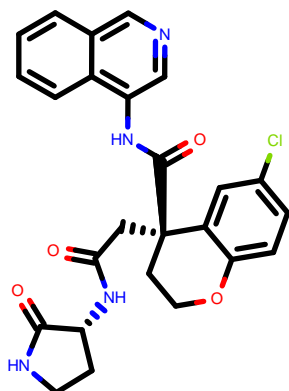
dDDG (kcal/mol): 0.16

RAL-THA-b9d6aec1-2_1



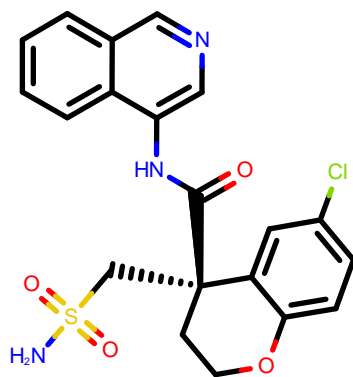
CID:	RAL-THA-b9d6aec1-2_1
SMILES:	<chem>CNS(=O)(=O)c1ccc2ncnc(c2c1)NC(=O)[C@@H]3CCOCc4cc(cc4)Cl</chem>
RUN:	RUN4490
DDG (kcal/mol):	-0.17
dDDG (kcal/mol):	0.33

KAD-UNI-877d7bed-8_1



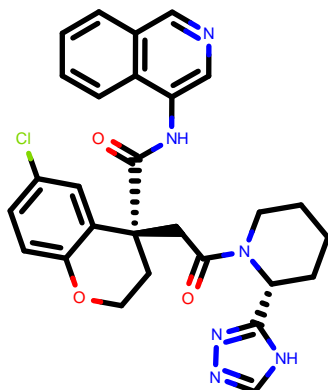
CID:	KAD-UNI-877d7bed-8_1
SMILES:	<chem>CC(=O)N1CCN(CC1)C(=O)COC2cc(cc3c2OCC[C@@H]3C(=O)Nc4cnc5c4ccoc5)Cl</chem>
RUN:	RUN3735
DDG (kcal/mol):	-0.17
dDDG (kcal/mol):	0.15

JOH-UNI-a38a7bdd-6_3



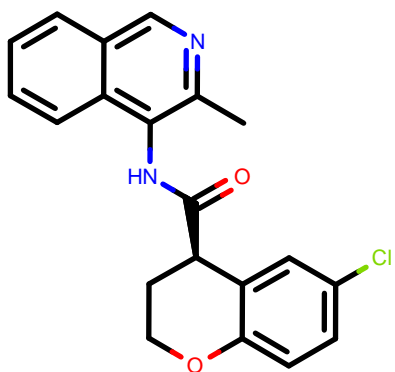
CID:	JOH-UNI-a38a7bdd-6_3
SMILES:	<chem>c1ccc2c(c1)cnc2N(C(=O)Cc3ccccc3)C(=O)[C@@H]4C[C@@H]4C(F)(F)F</chem>
RUN:	RUN1492
DDG (kcal/mol):	-0.16
dDDG (kcal/mol):	0.22

RAL-THA-4aa06b95-1_2



CID:	RAL-THA-4aa06b95-1_2
SMILES:	<chem>c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCN(c4c3cc(cc4)Cl)C(=O)N</chem>
RUN:	RUN1230
DDG (kcal/mol):	-0.16
dDDG (kcal/mol):	0.37

ALP-UNI-dbbfd3db-10_1



CID: ALP-UNI-dbbfd3db-10_1

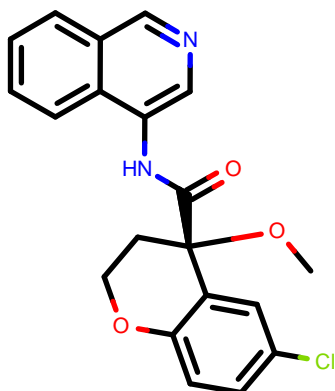
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4ccc(cc4)Cl)NC(=O)Cn5cc(c(=O)[nH]5=O)Cl

RUN: RUN2778

DDG (kcal/mol): -0.16

dDDG (kcal/mol): 0.11

MIC-UNK-0a05c952-3_8



CID: MIC-UNK-0a05c952-3_8

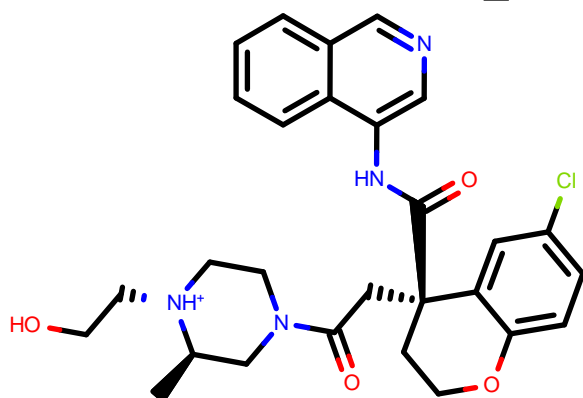
SMILES: c1ccc2c(c1)cncc2N3[C@H](C[C@H](C3=O)c4ccc(c(c4)Cl)Cl)[C@H]5CO5

RUN: RUN3518

DDG (kcal/mol): -0.15

dDDG (kcal/mol): 0.06

ALP-POS-c3a96089-2_1



CID: ALP-POS-c3a96089-2_1

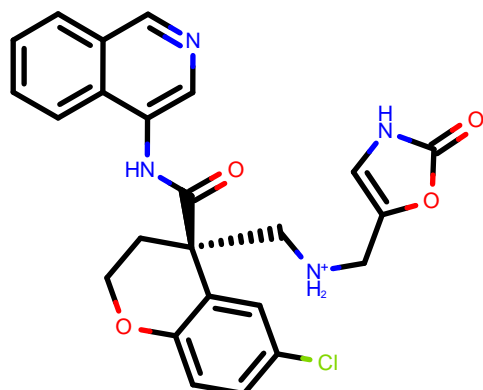
SMILES: Cc1ccc(nc1)N(Cc2cscn2)C(=O)Cc3cncc4c3cccc4

RUN: RUN1182

DDG (kcal/mol): -0.15

dDDG (kcal/mol): 0.58

DAR-DIA-9e4459de-13_14



CID: DAR-DIA-9e4459de-13_14

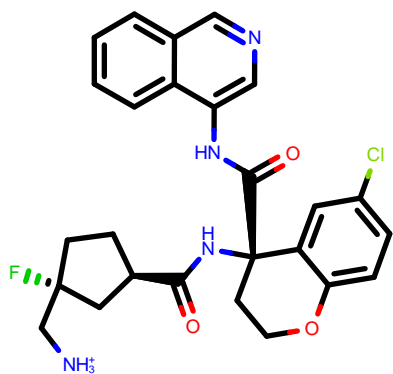
SMILES: c1cc2c(c1)NC(=O)C3CC4C(C3)NCC4NC(=O)C[C@H]5C(=O)NC(=O)N5

RUN: RUN1437

DDG (kcal/mol): -0.15

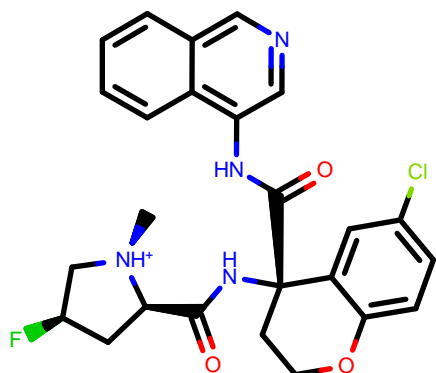
dDDG (kcal/mol): 0.37

MAK-UNK-ffc90da7-8_1



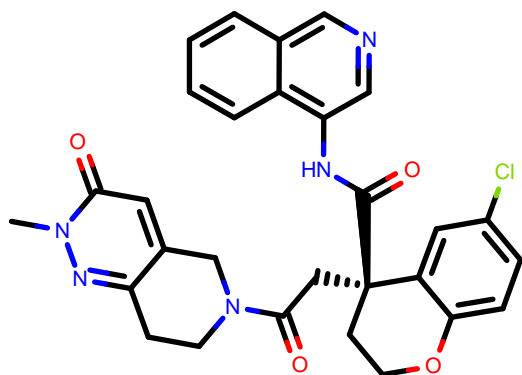
CID:	MAK-UNK-ffc90da7-8_1
SMILES:	<chem>CC(C)OCCCC1ccc2c(c1)cncc2NC(=O)Cc3cccc(c3)Cl</chem>
RUN:	RUN711
DDG (kcal/mol):	-0.15
dDDG (kcal/mol):	0.42

PET-UNK-824b5c6a-2_1



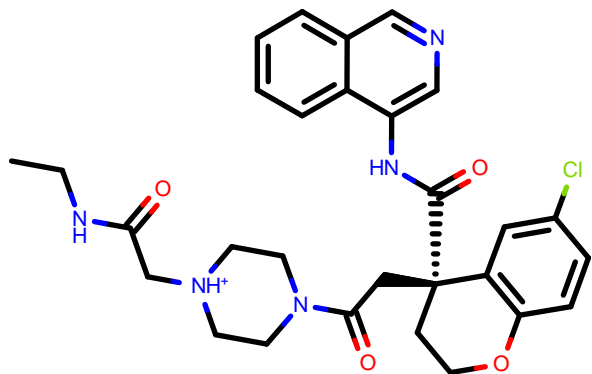
CID:	PET-UNK-824b5c6a-2_1
SMILES:	<chem>CCCO[C@]1(CCOC2c1cc(cc2)Cl)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3288
DDG (kcal/mol):	-0.14
dDDG (kcal/mol):	0.15

JOH-UNI-21fd6073-4_2



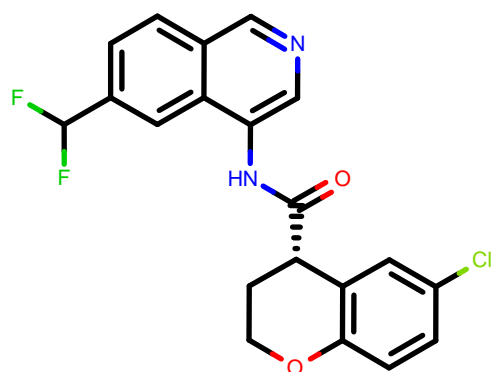
CID:	JOH-UNI-21fd6073-4_2
SMILES:	<chem>c1ccc2c(c1)cncc2N(CC(=O)O)n3c(ccc3O)O)C(=O)C[C@H]4COc5c4cc(cc5)Cl</chem>
RUN:	RUN4184
DDG (kcal/mol):	-0.14
dDDG (kcal/mol):	0.16

MAK-UNK-c749d764-19_6



CID:	MAK-UNK-c749d764-19_6
SMILES:	<chem>c1ccc2c(c1)cncc2N(C[NH3+])C(=O)C[C@H]3CCC[C@H]4[C@H]3O)C(F)F</chem>
RUN:	RUN1003
DDG (kcal/mol):	-0.13
dDDG (kcal/mol):	0.31

LON-WEI-4d77710c-16_1



CID: LON-WEI-4d77710c-16_1

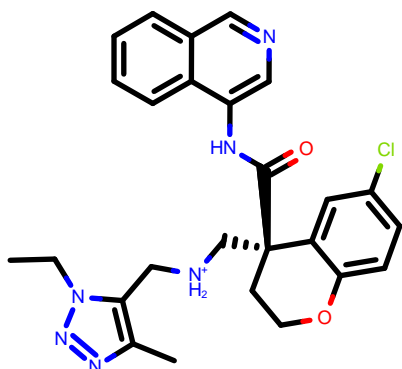
SMILES: CC(C)Cn1cc(c2ccccc2c1=O)NC(=O)NCC[NH+]3CCOCC3

RUN: RUN204

DDG (kcal/mol): -0.13

dDDG (kcal/mol): 0.11

RAL-THA-05e671eb-24_2



CID: RAL-THA-05e671eb-24_2

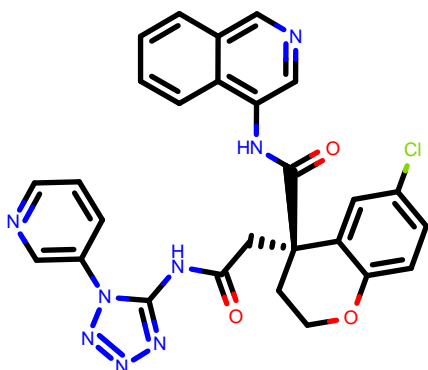
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3CCOC4c3ccccc4C#N

RUN: RUN2053

DDG (kcal/mol): -0.13

dDDG (kcal/mol): 0.36

KAD-UNI-877d7bed-1_2



CID: KAD-UNI-877d7bed-1_2

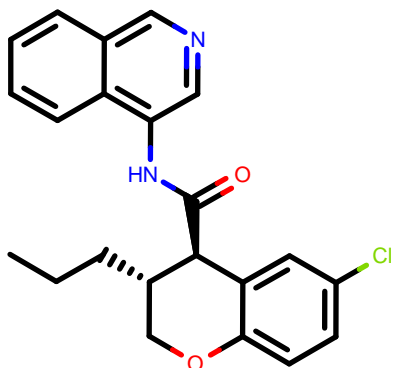
SMILES: c1ccc(cc1)[C@]([C@@]2OC2c(c3c2OCC[C@@]3[H]3C(=O)Nc4ncc5c4cccc5)Cl)C(=O)N

RUN: RUN3720

DDG (kcal/mol): -0.13

dDDG (kcal/mol): 0.17

EDG-MED-ba1ac7b9-20_3



CID: EDG-MED-ba1ac7b9-20_3

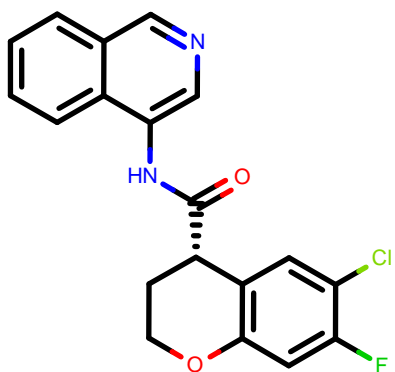
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)CC(=O)N5CCOC[C@H]5C(F)F

RUN: RUN2696

DDG (kcal/mol): -0.12

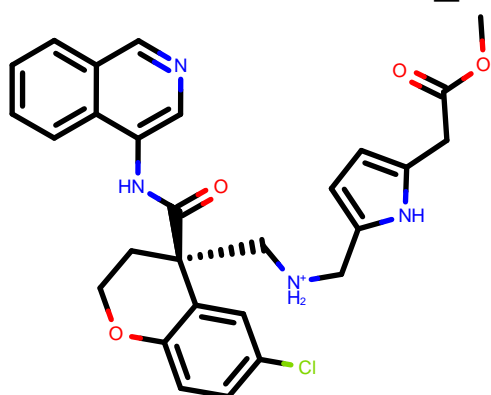
dDDG (kcal/mol): 0.09

LAU-MED-88a3970a-15_1



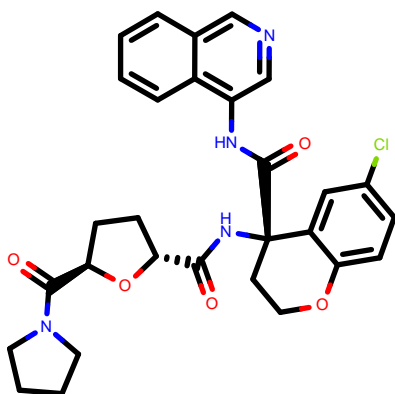
CID:	LAU-MED-88a3970a-15_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3CCOCc4cc(cc4CCS(=O)(=O)N)Cl</chem>
RUN:	RUN1511
DDG (kcal/mol):	-0.12
dDDG (kcal/mol):	0.10

JOH-UNI-ee5ed7c8-10_1



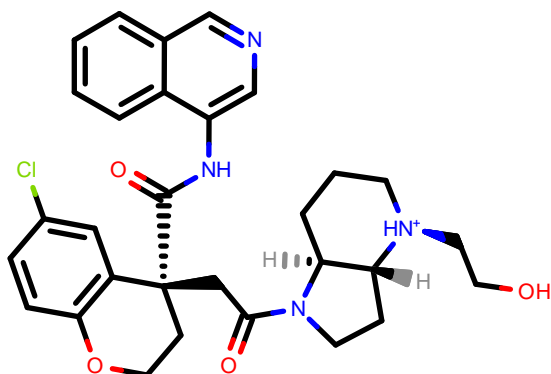
CID:	JOH-UNI-ee5ed7c8-10_1
SMILES:	<chem>CN(c1c2cccc2cnc1CC(F)(F)F)C(=O)Cc3cccc(c3)Cl</chem>
RUN:	RUN1909
DDG (kcal/mol):	-0.12
dDDG (kcal/mol):	0.37

MAT-POS-4223bc15-5_4



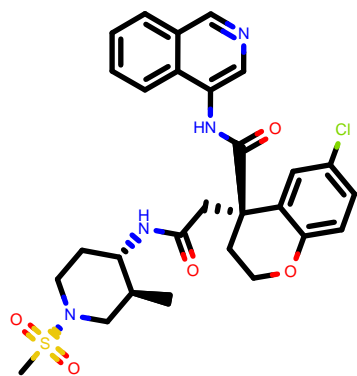
CID:	MAT-POS-4223bc15-5_4
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@H]3C[N@](C4c3cc(cc4)Cl)S(=O)(=O)CCO</chem>
RUN:	RUN3992
DDG (kcal/mol):	-0.12
dDDG (kcal/mol):	0.19

PET-UNK-9bf1291a-5_2



CID:	PET-UNK-9bf1291a-5_2
SMILES:	<chem>CO[C@]1(C[N@](Cc2c1cc(cc2)Cl)CC#N)C(=O)Nc3cncc4c3cccc4</chem>
RUN:	RUN3970
DDG (kcal/mol):	-0.12
dDDG (kcal/mol):	0.31

MAK-UNK-c749d764-16_14



CID: MAK-UNK-c749d764-16_14

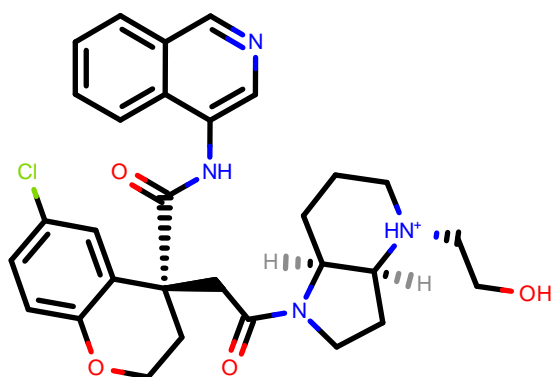
SMILES: C[C@H](N(c1cnc2c1cccc2)C(=O)C)[C@@H]3CCCC[C@H]([C@H]3O)C(F)O

RUN: RUN987

DDG (kcal/mol): -0.11

dDDG (kcal/mol): 0.40

MAK-UNK-c749d764-29_8



CID: MAK-UNK-c749d764-29_8

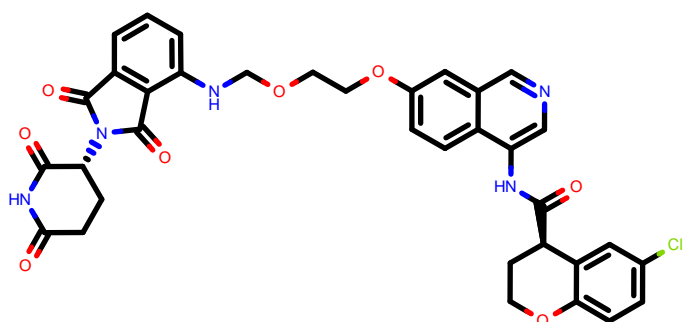
SMILES: c1ccc2c(c1)cncc2N(C(=O)C)[C@H]3CCCC[C@H]([C@H]3O)C(F)C(=O)ON

RUN: RUN1070

DDG (kcal/mol): -0.11

dDDG (kcal/mol): 0.45

ALP-UNI-8e43a71e-2_16



CID: ALP-UNI-8e43a71e-2_16

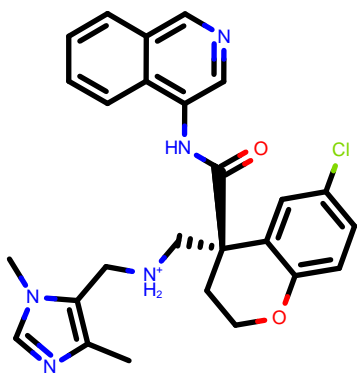
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3[C@@H]4CCOC(=O)C1=CC=C(C=C1)O[C@H]3C[C@@H]4N

RUN: RUN2938

DDG (kcal/mol): -0.11

dDDG (kcal/mol): 0.29

EDJ-MED-93390d0c-1_1



CID: EDJ-MED-93390d0c-1_1

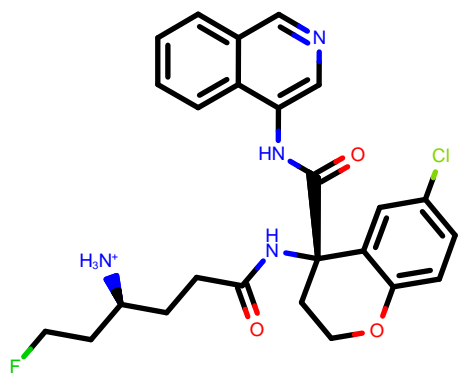
SMILES: CS(=O)(=O)c1ccc2cncc(c2c1)NC(=O)[C@H]3CCCC(=O)O[C@H]4C3CC(=O)C1=CC=C(C=C1)Cl

RUN: RUN4516

DDG (kcal/mol): -0.11

dDDG (kcal/mol): 0.26

MAK-UNK-c749d764-3_3



CID: MAK-UNK-c749d764-3_3

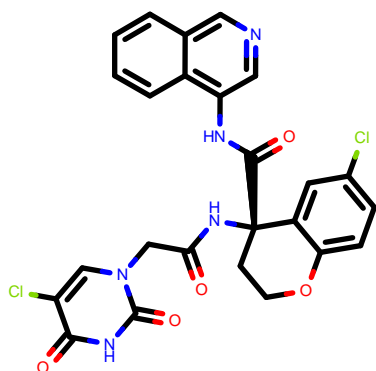
SMILES: CC1(C[NH2+](C1)OCN(c2cncc3c2cccc3)C(=O)C[C@@H]4CC[C@H]([C@@H]4O)C(F)F

RUN: RUN907

DDG (kcal/mol): -0.11

dDDG (kcal/mol): 0.33

DAR-DIA-f6ee7aeb-6_4



CID: DAR-DIA-f6ee7aeb-6_4

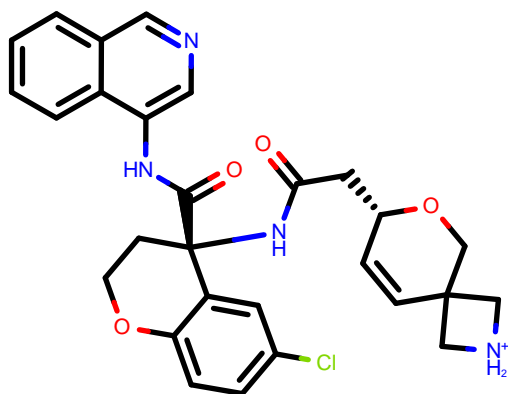
SMILES: c1ccc2c(c1)cncc2N3C[C@H]([C@H](CC3=O)c4cccc(c4)Cl)c5c[nH]c(=O)[nH]5=O

RUN: RUN3422

DDG (kcal/mol): -0.10

dDDG (kcal/mol): 0.15

MIC-UNK-bcd487e9-2_1



CID: MIC-UNK-bcd487e9-2_1

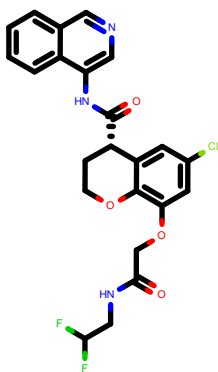
SMILES: c1ccc2c(c1)cncc2NC(=O)N(CC3CCCCC3)c4cccc(c4)Cl

RUN: RUN587

DDG (kcal/mol): -0.10

dDDG (kcal/mol): 0.31

MAT-POS-f39f51fd-1_1



CID: MAT-POS-f39f51fd-1_1

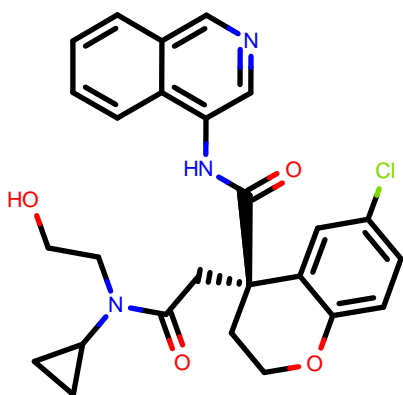
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@H]3CC(=O)Nc4c3cc(cc4Cl)Cl

RUN: RUN2404

DDG (kcal/mol): -0.10

dDDG (kcal/mol): 0.29

EDJ-MED-841e0cf0-7_3



CID: EDJ-MED-841e0cf0-7_3

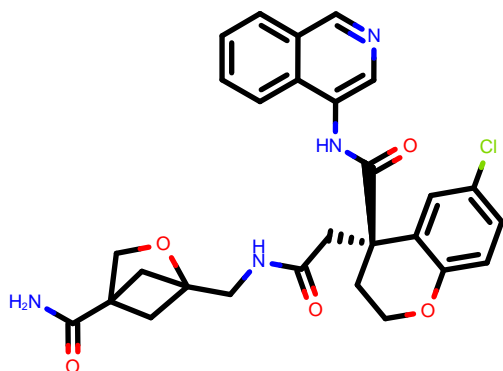
SMILES: CNS(=O)(=O)[N@@]1Cc2ccc(cc2[C@H](C1)C(=O)Nc3cncc4c3cc(cc4)F)Cl

RUN: RUN3846

DDG (kcal/mol): -0.10

dDDG (kcal/mol): 0.15

MIC-UNK-5a93dd5f-6_1



CID: MIC-UNK-5a93dd5f-6_1

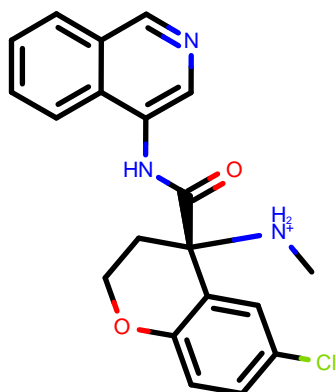
SMILES: CC(=O)N(C)C1CC[NH+]([C@]1[C@@H](c2cccc(c2)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN766

DDG (kcal/mol): -0.10

dDDG (kcal/mol): 0.49

MAK-UNK-c749d764-1_2



CID: MAK-UNK-c749d764-1_2

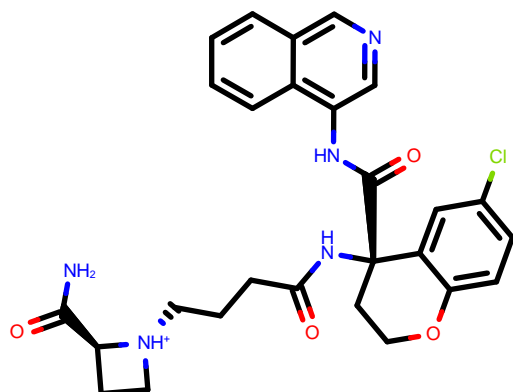
SMILES: c1ccc2c(c1)cncc2NC(=O)C[C@H]3CCCC[C@@]3([C@@H]3O)C4CC4

RUN: RUN889

DDG (kcal/mol): -0.09

dDDG (kcal/mol): 0.31

EDJ-MED-37aac4bd-1_2



CID: EDJ-MED-37aac4bd-1_2

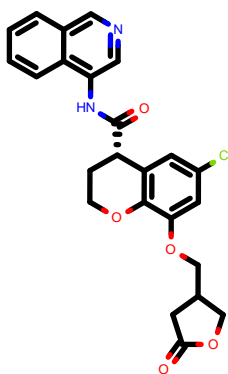
SMILES: CO[C@@]1(CCOc2c1cc(c(c2)Cl)Cl)C(=O)Nc3cncc4c3cccc4

RUN: RUN3140

DDG (kcal/mol): -0.09

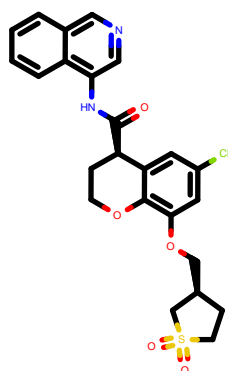
dDDG (kcal/mol): 0.21

ALP-POS-2da19ca7-2_2



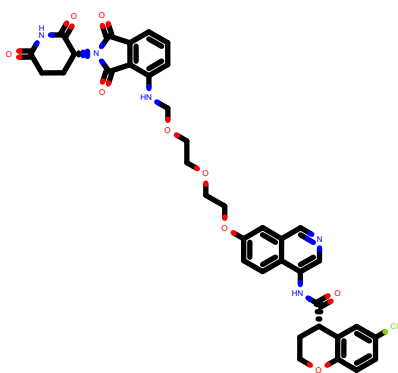
CID:	ALP-POS-2da19ca7-2_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@H]3[C@@H](C)CCOC4=CC(=O)NCC(NH)CC5=C[C@H]6COC(=O)C5</chem>
RUN:	RUN2376
DDG (kcal/mol):	-0.09
dDDG (kcal/mol):	0.23

MIC-UNK-37660950-3_2



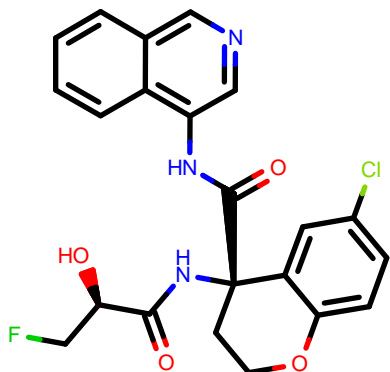
CID:	MIC-UNK-37660950-3_2
SMILES:	<chem>CS(=O)(=O)[N@H]1CC[C@H](C1)CNc2ccc(cc2CC(=O)Nc3cncoc4c3ccccc4)Cl</chem>
RUN:	RUN5113
DDG (kcal/mol):	-0.08
dDDG (kcal/mol):	0.20

MAT-POS-8a69d52e-4_3



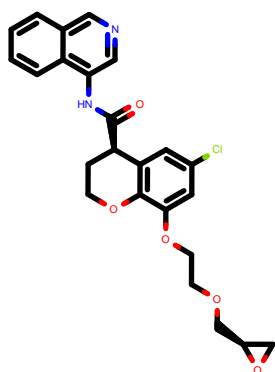
CID:	MAT-POS-8a69d52e-4_3
SMILES:	<chem>C[C@H]1COc2ccc(cc2[C@H]1C(=O)Nc3cncoc4c3ccccc4)Cl</chem>
RUN:	RUN368
DDG (kcal/mol):	-0.08
dDDG (kcal/mol):	0.38

ALP-POS-fe871b40-15_2



CID:	ALP-POS-fe871b40-15_2
SMILES:	<chem>c1ccc2c(c1)ncnc2NC(=O)[C@H]3CC(=O)Nc4c3cc(c(c4)F)Cl</chem>
RUN:	RUN3137
DDG (kcal/mol):	-0.08
dDDG (kcal/mol):	0.16

BEN-DND-11faade0-1_2



CID: BEN-DND-11faade0-1_2

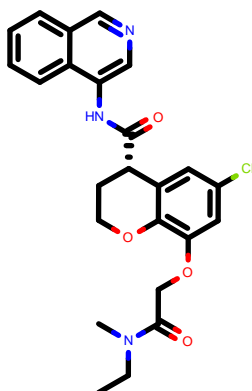
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@H]3CNS(=O)(=O)c4c3cccc4

RUN: RUN5075

DDG (kcal/mol): -0.07

dDDG (kcal/mol): 0.20

EDJ-MED-9e38fd34-4_2



CID: EDJ-MED-9e38fd34-4_2

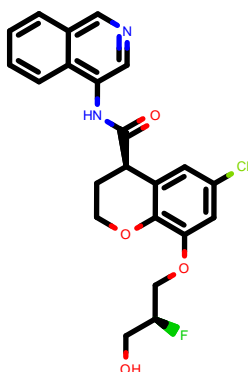
SMILES: C[C@]1(c2cc(c(cc2NC1=O)Cl)C)C(=O)Nc3cncc4c3cccc4

RUN: RUN2350

DDG (kcal/mol): -0.06

dDDG (kcal/mol): 0.28

ALF-EVA-a24cc7ce-2_2



CID: ALF-EVA-a24cc7ce-2_2

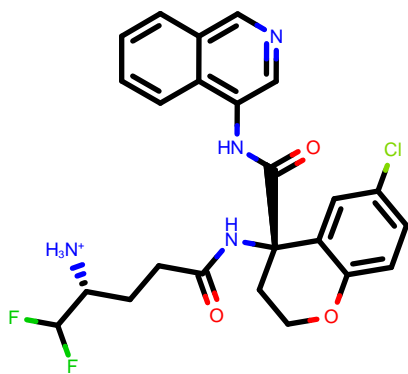
SMILES: CN(c1cncc2c1cccc2)C(=O)[C@H]3COc4c3cc(cc4)Cl

RUN: RUN4956

DDG (kcal/mol): -0.06

dDDG (kcal/mol): 0.22

DAR-DIA-0587064e-20_2



CID: DAR-DIA-0587064e-20_2

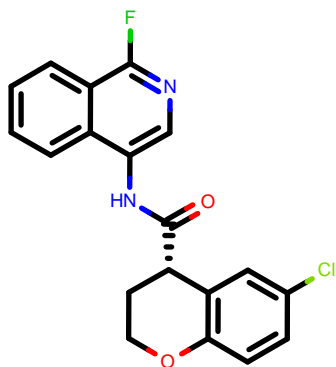
SMILES: c1ccc(cc1)CCOc2cc(cc3c2OCC[C@H]3C(=O)Nc4cncc5c4cccc5)Cl

RUN: RUN3378

DDG (kcal/mol): -0.06

dDDG (kcal/mol): 0.24

ALP-POS-6747fa38-1_1



CID: ALP-POS-6747fa38-1_1

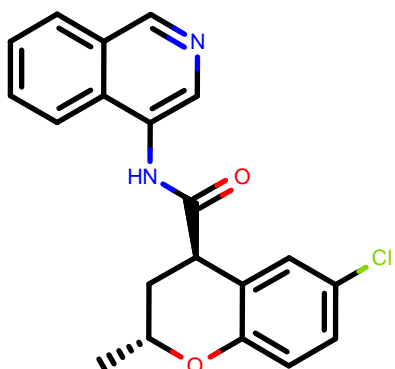
SMILES: CC(=O)N1CC[NH+](CC1)CC(=O)Nc2cncc3c2cccc3

RUN: RUN188

DDG (kcal/mol): -0.06

dDDG (kcal/mol): 0.10

EDG-MED-ba1ac7b9-17_2



CID: EDG-MED-ba1ac7b9-17_2

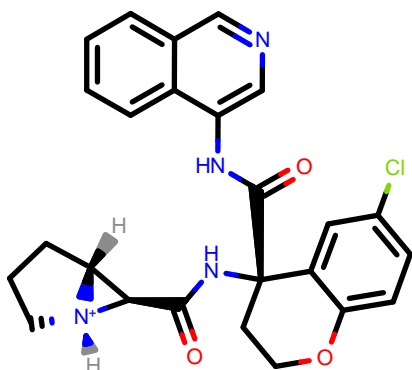
SMILES: Cn1ccnc1CN(C2CC2)C(=O)C[C@@]3(CCOc4c3cc(cc4)Cl)C(=O)Nc5cncc6c5cccc6

RUN: RUN2689

DDG (kcal/mol): -0.05

dDDG (kcal/mol): 0.08

NAU-LAT-0543f7f2-10_1



CID: NAU-LAT-0543f7f2-10_1

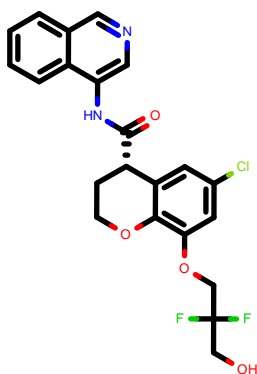
SMILES: CC(=O)NCCOC1cc(cc2c1OCC[C@@]3(CCOc4c3cc(cc4)Cl)C(=O)Nc5cncc4c3cccc4)Cl

RUN: RUN659

DDG (kcal/mol): -0.05

dDDG (kcal/mol): 0.27

VLA-UNK-f702bf1c-8_1



CID: VLA-UNK-f702bf1c-8_1

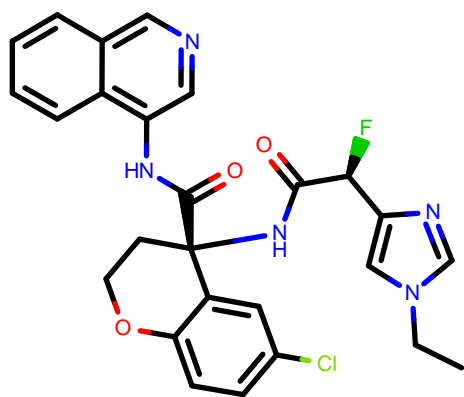
SMILES: c1ccc2c(c1)cncc2N3C(=O)[C@@]4(CCOc5c4cc(cc5)Cl)N(C3=O)CCc6[nH]nnc6

RUN: RUN2322

DDG (kcal/mol): -0.05

dDDG (kcal/mol): 0.28

EDJ-MED-f893e2a1-7_1



CID: EDJ-MED-f893e2a1-7_1

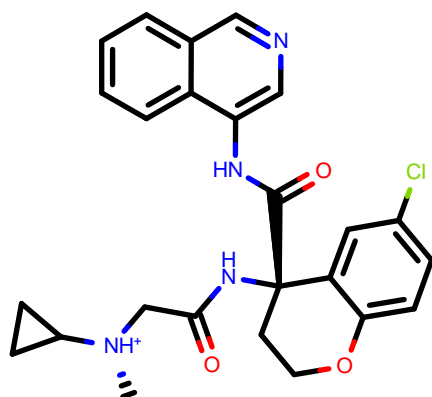
SMILES: c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)C([NH2+][C5c[nH]c(=O)o5

RUN: RUN3206

DDG (kcal/mol): -0.03

dDDG (kcal/mol): 0.18

MIC-UNK-bcd487e9-5_1



CID: MIC-UNK-bcd487e9-5_1

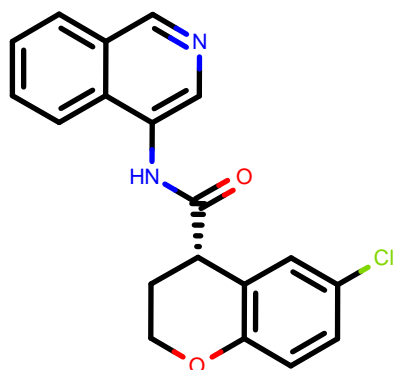
SMILES: c1ccc2c(c1)cncc2NC(=O)N(Cc3ccccc3)c4cccc(c4)Cl

RUN: RUN591

DDG (kcal/mol): -0.02

dDDG (kcal/mol): 0.39

DAR-DIA-0cde14eb-54_1



CID: DAR-DIA-0cde14eb-54_1

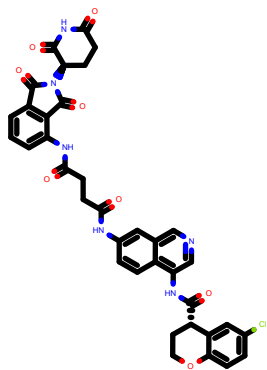
SMILES: c1ccc2c(c1)cncc2NC(=O)Nc3ccccc(c3)C4(CC4)C#N

RUN: RUN19

DDG (kcal/mol): -0.02

dDDG (kcal/mol): 0.09

ALP-POS-7c6e02c7-1_1



CID: ALP-POS-7c6e02c7-1_1

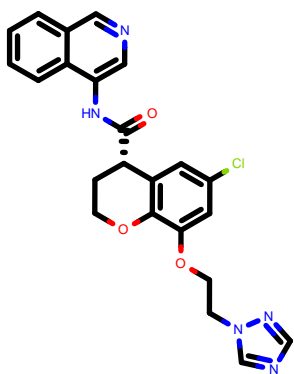
SMILES: CN(C)c1cnc(nc1)N(Cc2ccc(c(c2)Cl)Cl)C(=O)Cc3cnc4c3ccccc4

RUN: RUN3099

DDG (kcal/mol): -0.02

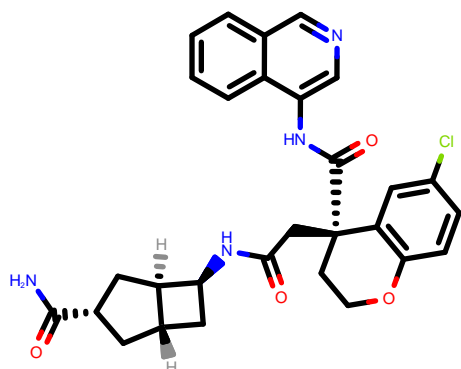
dDDG (kcal/mol): 0.42

ALP-POS-5bb456a5-1_13



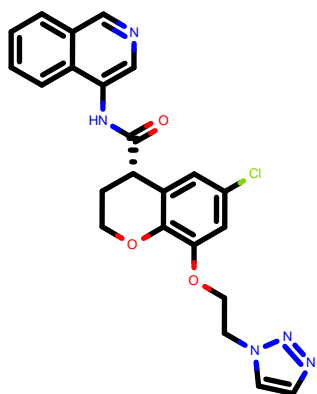
CID:	ALP-POS-5bb456a5-1_13
SMILES:	<chem>C1C=CN=C2C=CC=C1C(=O)N(C2)C(=O)C3=CC(OC3)C(Cl)=O</chem>
RUN:	RUN2417
DDG (kcal/mol):	-0.02
dDDG (kcal/mol):	0.25

DAR-DIA-5ff57136-7_1



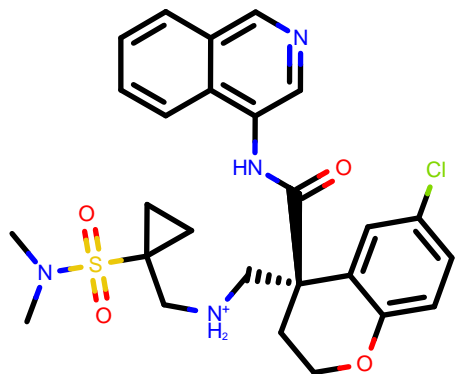
CID:	DAR-DIA-5ff57136-7_1
SMILES:	<chem>C=CC(=O)N(c1cncc2c1cccc2)C(=O)C3=CCOc4c3cc(cc4)Cl</chem>
RUN:	RUN1376
DDG (kcal/mol):	-0.02
dDDG (kcal/mol):	0.24

ALF-EVA-5b152d2f-7_1



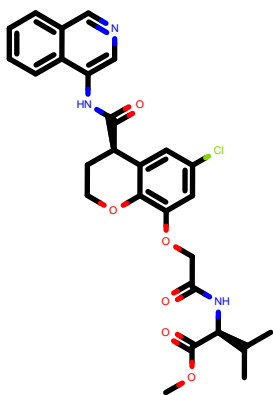
CID:	ALF-EVA-5b152d2f-7_1
SMILES:	<chem>c1cc2ncc(c2cc1C3CCC3)NC(=O)[C@@H]4CCOc5c4cc(cc5)Cl</chem>
RUN:	RUN2361
DDG (kcal/mol):	-0.02
dDDG (kcal/mol):	0.27

MAT-POS-ec6d90b7-2_1



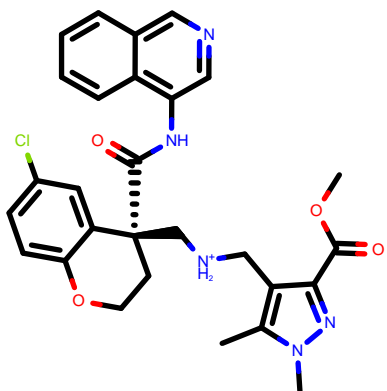
CID:	MAT-POS-ec6d90b7-2_1
SMILES:	<chem>c1ccc2c(c1)cncc2N3C(=O)[C@]4(CCOc5c4cc(cc5)Cl)OC3=O</chem>
RUN:	RUN4452
DDG (kcal/mol):	-0.02
dDDG (kcal/mol):	0.45

MAT-POS-2905de8c-1_1



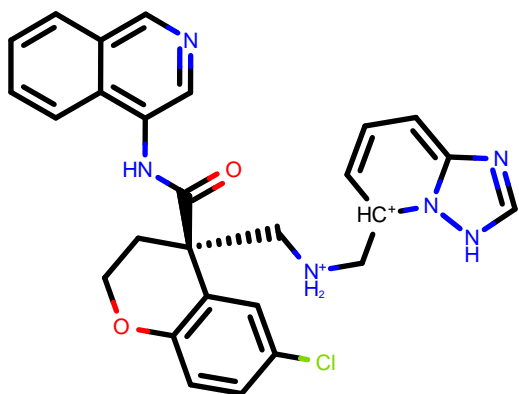
CID:	MAT-POS-2905de8c-1_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)[NH3+]</chem>
RUN:	RUN2225
DDG (kcal/mol):	-0.02
dDDG (kcal/mol):	0.36

EDG-MED-90036822-22_1



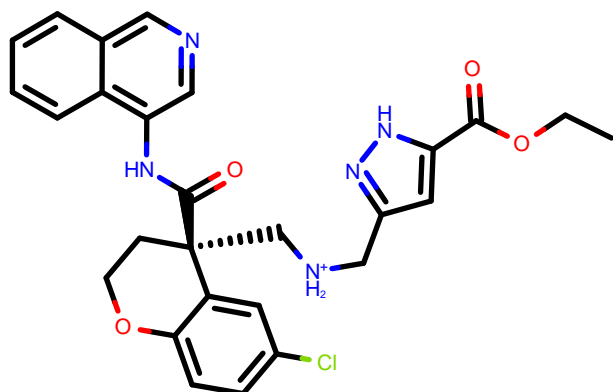
CID:	EDG-MED-90036822-22_1
SMILES:	<chem>C[C@@]([CC](=O)N[C@]1(CCOc2c1cc(cc2)Cl)C(=O)Nc3cncc4c3ccccc4)[C](NH3+)]OCCO</chem>
RUN:	RUN1691
DDG (kcal/mol):	-0.02
dDDG (kcal/mol):	0.46

ED_-GRI-5b13fbe2-54_1



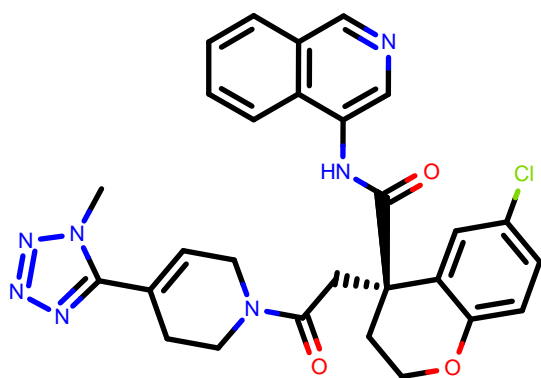
CID:	ED_-GRI-5b13fbe2-54_1
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@]3(CCOc4c3cc(cc4)Cl)OCCCN5c(c1nn5)[C](NH3+)]C(F)F</chem>
RUN:	RUN1599
DDG (kcal/mol):	-0.02
dDDG (kcal/mol):	0.41

MAT-POS-dc2604c4-3_3



CID:	MAT-POS-dc2604c4-3_3
SMILES:	<chem>c1ccc2c(c1)cncc2NC(=O)[C@@H]3C[N@](C)C4c3cc(cc4)Cl)S(=O)(=O)CC5(CNC5)C#N</chem>
RUN:	RUN4714
DDG (kcal/mol):	-0.01
dDDG (kcal/mol):	0.25

MAK-UNK-c749d764-12_5



CID: MAK-UNK-c749d764-12_5

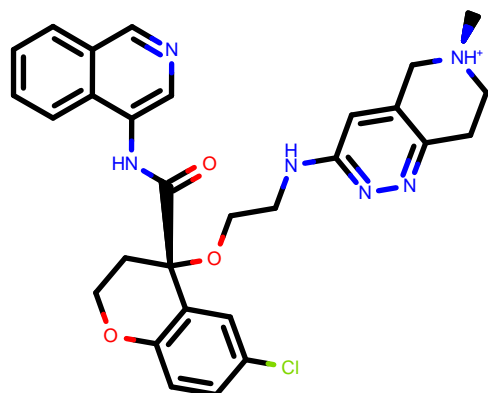
SMILES: c1ccc2c(c1)cnc2NC(=O)C[C@@H]3CCCC[C@@H]([C@@H]3O)Cl

RUN: RUN954

DDG (kcal/mol): -0.01

dDDG (kcal/mol): 0.29

DAR-DIA-23e5a6a0-6_2



CID: DAR-DIA-23e5a6a0-6_2

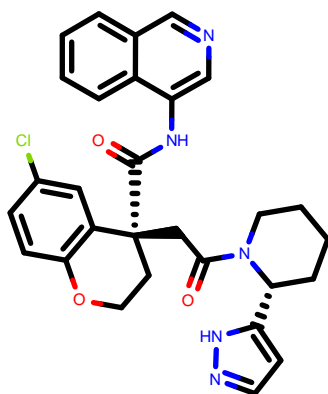
SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3CCOCc4c3cc(cc4C)[C@H]5CCCC6([NH2+][5]O)C6)Cl

RUN: RUN412

DDG (kcal/mol): -0.01

dDDG (kcal/mol): 0.41

JOH-SUS-a69c159d-2_2



CID: JOH-SUS-a69c159d-2_2

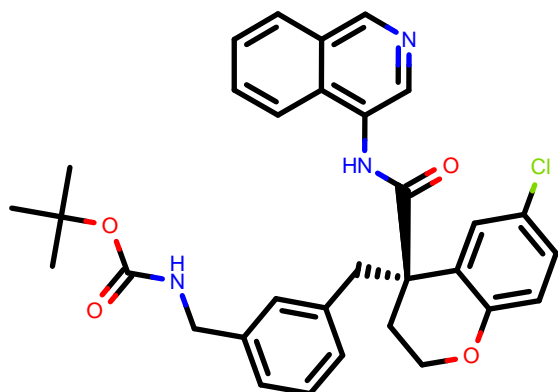
SMILES: c1ccc2c(c1)c(cnc2F)NC(=O)[C@@H]3CCOCc4c3cc(cc4)Cl

RUN: RUN1119

DDG (kcal/mol): -0.01

dDDG (kcal/mol): 0.28

DAR-DIA-ecdbc7dd-4_2



CID: DAR-DIA-ecdbc7dd-4_2

SMILES: c1ccc2c(c1)cnc2NC(=O)[C@@H]3(CCOc4c3cc(cc4)Cl)C[NH+]5CCCC5

RUN: RUN2882

DDG (kcal/mol): -0.01

dDDG (kcal/mol): 0.16