



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 17, 2024 – 07:47 AM EST

PDB ID : 4I0F  
Title : Design and Synthesis of Thiophene Dihydroisoquinolins as Novel BACE-1 Inhibitors  
Authors : Yao, N.; Brecht, E.  
Deposited on : 2012-11-16  
Resolution : 1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

---

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

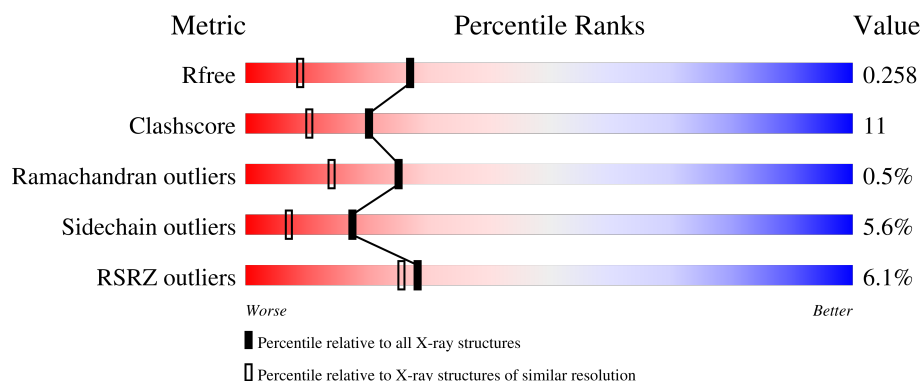
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	164625	7108 (1.80-1.80)
Clashscore	180529	8162 (1.80-1.80)
Ramachandran outliers	177936	8077 (1.80-1.80)
Sidechain outliers	177891	8076 (1.80-1.80)
RSRZ outliers	164620	7108 (1.80-1.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	406	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3478 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	0	0
			3099	1977	523	585	14			

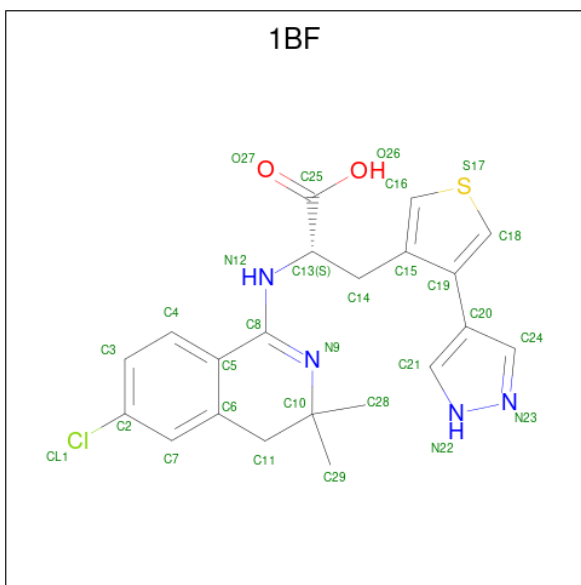
There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	56	MET	-	expression tag	UNP P56817
A	454	ARG	-	expression tag	UNP P56817
A	455	SER	-	expression tag	UNP P56817
A	456	HIS	-	expression tag	UNP P56817
A	457	HIS	-	expression tag	UNP P56817
A	458	HIS	-	expression tag	UNP P56817
A	459	HIS	-	expression tag	UNP P56817
A	460	HIS	-	expression tag	UNP P56817
A	461	HIS	-	expression tag	UNP P56817

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Zn	0	0
			3	3		

- Molecule 3 is N-(6-chloro-3,3-dimethyl-3,4-dihydroisoquinolin-1-yl)-3-[4-(1H-pyrazol-4-yl)thiophen-3-yl]-L-alanine (three-letter code: 1BF) (formula: C<sub>21</sub>H<sub>21</sub>ClN<sub>4</sub>O<sub>2</sub>S).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	Cl	N	O	S	0	0
			29	21	1	4	2	1		

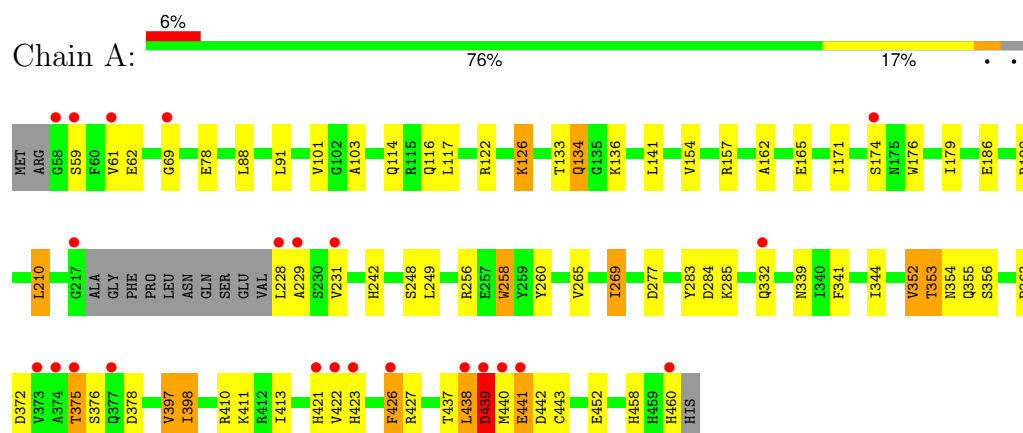
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	347	Total	O	0	0
			347	347		

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: Beta-secretase 1



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	75.17Å 103.53Å 100.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.00 – 1.80 60.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	81.2 (60.00-1.80) 85.6 (60.00-1.80)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.60 (at 1.76Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, $R_{free}$	0.208 , 0.266 0.204 , 0.258	Depositor DCC
$R_{free}$ test set	1853 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	30.7	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 42.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.54$ , $\langle L^2 \rangle = 0.38$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3478	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.44% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 1BF, ZN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	1.11	3/3180 (0.1%)	0.97	6/4322 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	352	VAL	CB-CG1	5.73	1.64	1.52
1	A	78	GLU	CB-CG	-5.26	1.42	1.52
1	A	356	SER	C-O	5.10	1.33	1.23

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	277	ASP	CB-CG-OD1	7.40	124.96	118.30
1	A	210	LEU	CA-CB-CG	6.87	131.11	115.30
1	A	157	ARG	NE-CZ-NH1	6.23	123.42	120.30
1	A	427	ARG	NE-CZ-NH2	-5.76	117.42	120.30
1	A	122	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	A	367	LEU	CA-CB-CG	5.67	128.33	115.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	258	TRP	Peptide

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3099	0	2990	66	1
2	A	3	0	0	0	0
3	A	29	0	20	4	0
4	A	347	0	0	15	2
All	All	3478	0	3010	68	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (68) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:437:THR:HG22	1:A:438:LEU:N	1.53	1.18
1:A:437:THR:CG2	1:A:438:LEU:H	1.54	1.17
1:A:397:VAL:HG12	1:A:398:ILE:HD13	1.31	1.10
1:A:439:ASP:O	1:A:441:GLU:HG2	1.51	1.09
1:A:441:GLU:H	1:A:442:ASP:HA	1.32	0.95
1:A:114:GLN:HG2	1:A:116:GLN:HE22	1.33	0.93
1:A:440:MET:HA	1:A:441:GLU:HG3	1.49	0.93
1:A:59:SER:O	4:A:734:HOH:O	1.90	0.88
1:A:423:HIS:HB3	4:A:856:HOH:O	1.77	0.83
1:A:437:THR:HG22	1:A:438:LEU:H	0.69	0.82
1:A:423:HIS:CB	4:A:856:HOH:O	2.30	0.79
1:A:440:MET:HA	1:A:441:GLU:CG	2.13	0.78
1:A:441:GLU:H	1:A:442:ASP:CA	1.96	0.78
1:A:62:GLU:HB2	4:A:734:HOH:O	1.88	0.74
1:A:441:GLU:N	1:A:442:ASP:HA	2.03	0.72
1:A:439:ASP:C	1:A:441:GLU:HG2	2.12	0.69
1:A:440:MET:N	1:A:441:GLU:HA	2.07	0.69
1:A:439:ASP:O	1:A:440:MET:SD	2.52	0.68
1:A:134:GLN:HG3	3:A:504:1BF:C3	2.26	0.65

*Continued on next page...*



*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:397:VAL:CG1	1:A:398:ILE:HD13	2.18	0.65
1:A:376:SER:HB3	4:A:707:HOH:O	1.97	0.65
1:A:437:THR:CG2	1:A:438:LEU:N	2.29	0.64
1:A:265:VAL:HG11	1:A:440:MET:HB3	1.79	0.64
1:A:133:THR:HG22	1:A:134:GLN:NE2	2.14	0.63
1:A:441:GLU:HB2	1:A:443:CYS:SG	2.42	0.59
1:A:339:ASN:HB3	1:A:426:PHE:CZ	2.38	0.59
1:A:126:LYS:HG2	1:A:141:LEU:HD12	1.85	0.58
1:A:242:HIS:HE1	4:A:641:HOH:O	1.87	0.57
1:A:88:LEU:HD13	1:A:101:VAL:HG11	1.89	0.55
1:A:260:TYR:HB3	1:A:413:ILE:HD11	1.90	0.54
1:A:397:VAL:HG12	1:A:398:ILE:CD1	2.20	0.54
3:A:504:1BF:H16	3:A:504:1BF:H21	1.91	0.52
1:A:458:HIS:HE1	4:A:693:HOH:O	1.92	0.52
1:A:176:TRP:HZ2	3:A:504:1BF:H18	1.74	0.52
1:A:375:THR:HB	1:A:378:ASP:HB2	1.92	0.51
1:A:438:LEU:O	1:A:439:ASP:HB2	2.12	0.50
1:A:332:GLN:HG2	4:A:741:HOH:O	2.12	0.49
1:A:269:ILE:HD13	1:A:344:ILE:HG12	1.94	0.49
1:A:186:GLU:OE2	1:A:256:ARG:NH1	2.36	0.49
1:A:440:MET:HG2	4:A:930:HOH:O	2.13	0.49
1:A:154:VAL:HG21	1:A:205:THR:HG21	1.96	0.48
1:A:355:GLN:C	1:A:440:MET:HE2	2.33	0.48
1:A:353:THR:HG23	1:A:354:ASN:ND2	2.29	0.48
1:A:365:GLN:O	1:A:397:VAL:HG22	2.14	0.48
1:A:458:HIS:HB3	1:A:460:HIS:CE1	2.49	0.47
1:A:114:GLN:HG2	1:A:116:GLN:NE2	2.14	0.47
1:A:341:PHE:HB3	1:A:363:PRO:HB3	1.96	0.47
1:A:411:LYS:HD2	4:A:892:HOH:O	2.14	0.47
1:A:59:SER:OG	1:A:61:VAL:HG12	2.15	0.47
1:A:126:LYS:CG	1:A:141:LEU:HD12	2.44	0.47
1:A:103:ALA:CB	1:A:162:ALA:HB1	2.45	0.46
1:A:203:LYS:NZ	1:A:203:LYS:HB3	2.30	0.46
1:A:242:HIS:CE1	4:A:641:HOH:O	2.64	0.46
1:A:285:LYS:NZ	4:A:910:HOH:O	2.50	0.45
1:A:171:ILE:HG13	1:A:174:SER:HB3	1.98	0.45
1:A:440:MET:CA	1:A:441:GLU:CG	2.91	0.43
1:A:423:HIS:HB2	4:A:856:HOH:O	2.08	0.43
1:A:114:GLN:HB3	1:A:117:LEU:HG	2.01	0.42
1:A:228:LEU:HB3	1:A:229:ALA:H	1.53	0.42
1:A:91:LEU:HD23	1:A:179:ILE:HG13	2.01	0.42

*Continued on next page...*

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:372:ASP:HB2	4:A:699:HOH:O	2.20	0.42
3:A:504:1BF:H16	3:A:504:1BF:C24	2.49	0.42
1:A:69:GLY:C	1:A:231:VAL:HG22	2.39	0.42
1:A:363:PRO:HA	1:A:366:TYR:CE2	2.54	0.42
1:A:352:VAL:HG12	1:A:353:THR:O	2.21	0.41
1:A:441:GLU:H	1:A:443:CYS:N	2.19	0.41
1:A:410:ARG:NE	4:A:802:HOH:O	2.33	0.40
1:A:283:TYR:HA	1:A:284:ASP:HA	1.69	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:828:HOH:O	4:A:844:HOH:O[4_555]	1.72	0.48
1:A:165:GLU:OE1	4:A:828:HOH:O[4_555]	2.08	0.12

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	389/406 (96%)	377 (97%)	10 (3%)	2 (0%)	25 14

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	439	ASP
1	A	441	GLU

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	337/348 (97%)	318 (94%)	19 (6%)	17 7

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	126	LYS
1	A	134	GLN
1	A	136	LYS
1	A	189	ARG
1	A	210	LEU
1	A	248	SER
1	A	249	LEU
1	A	258	TRP
1	A	269	ILE
1	A	353	THR
1	A	375	THR
1	A	397	VAL
1	A	398	ILE
1	A	421	HIS
1	A	422	VAL
1	A	426	PHE
1	A	438	LEU
1	A	439	ASP
1	A	452	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	116	GLN
1	A	134	GLN
1	A	153	ASN
1	A	242	HIS
1	A	355	GLN
1	A	387	GLN

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	456	HIS
1	A	458	HIS
1	A	460	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

### 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

### 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

### 5.6 Ligand geometry ⓘ

Of 4 ligands modelled in this entry, 3 are monoatomic - leaving 1 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
3	1BF	A	504	-	29,32,32	2.48	8 (27%)	31,47,47	8.37	17 (54%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	1BF	A	504	-	-	1/13/30/30	0/4/4/4

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	504	1BF	C18-C19	6.96	1.41	1.37
3	A	504	1BF	C16-C15	6.16	1.40	1.37
3	A	504	1BF	C19-C15	4.90	1.52	1.40
3	A	504	1BF	C5-C6	3.79	1.45	1.40
3	A	504	1BF	C5-C8	3.56	1.50	1.46
3	A	504	1BF	C16-S17	3.00	1.74	1.70
3	A	504	1BF	N23-N22	2.74	1.43	1.37
3	A	504	1BF	C18-S17	2.05	1.73	1.70

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	504	1BF	C19-C18-S17	-33.54	103.17	112.53
3	A	504	1BF	C18-C19-C15	26.35	118.95	111.69
3	A	504	1BF	C15-C16-S17	-14.48	108.49	112.53
3	A	504	1BF	C18-C19-C20	-4.55	118.62	125.52
3	A	504	1BF	N12-C8-N9	4.18	125.62	119.71
3	A	504	1BF	C18-S17-C16	3.80	100.11	92.37
3	A	504	1BF	C3-C4-C5	-3.45	115.17	120.86
3	A	504	1BF	C2-C7-C6	-3.20	115.75	119.96
3	A	504	1BF	C4-C5-C6	3.20	122.78	119.41
3	A	504	1BF	C4-C3-C2	2.92	122.18	119.24
3	A	504	1BF	O26-C25-O27	-2.64	118.09	124.08
3	A	504	1BF	C14-C15-C16	-2.38	122.95	127.46
3	A	504	1BF	C21-C20-C24	2.24	111.87	105.13
3	A	504	1BF	C3-C2-C7	2.08	124.25	121.53
3	A	504	1BF	C29-C10-N9	2.01	111.89	107.16
3	A	504	1BF	C25-C13-N12	-2.01	105.91	110.57
3	A	504	1BF	C6-C5-C8	-2.00	116.31	118.92

There are no chirality outliers.

All (1) torsion outliers are listed below:

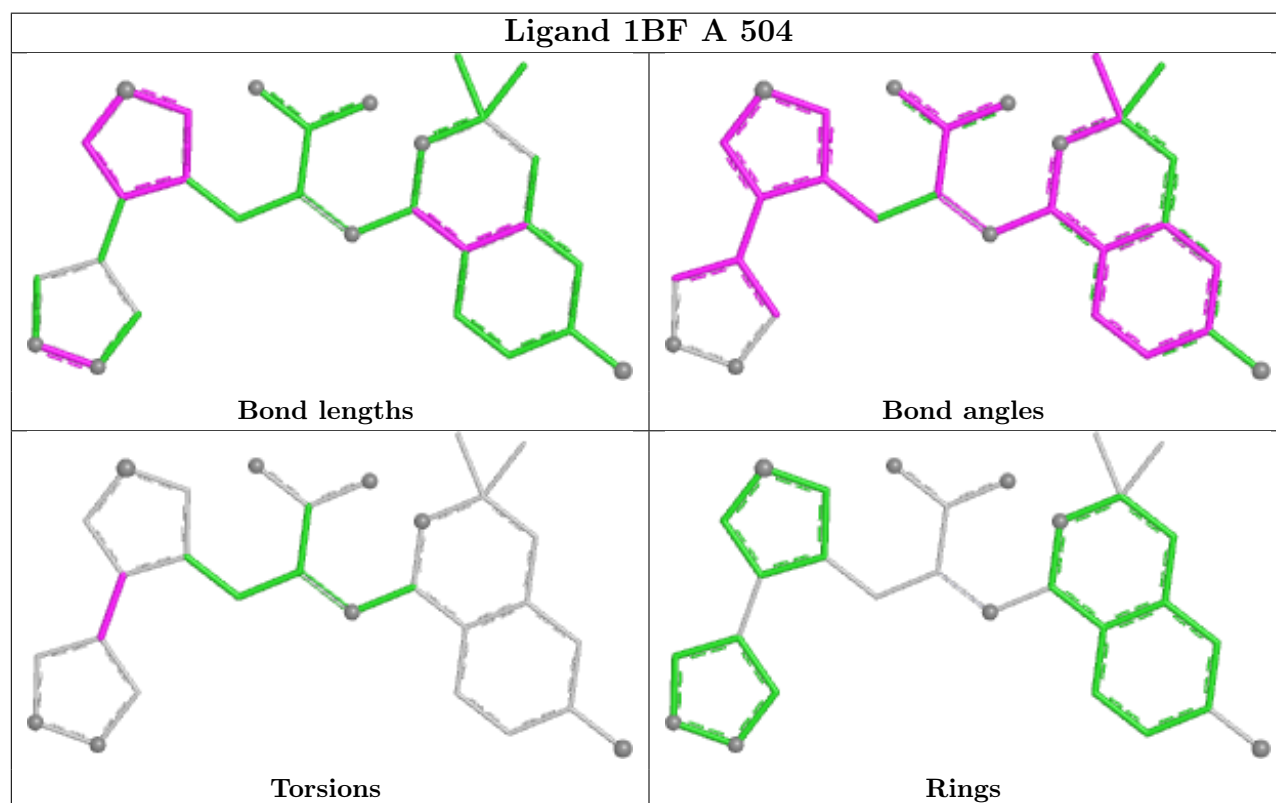
Mol	Chain	Res	Type	Atoms
3	A	504	1BF	C18-C19-C20-C21

There are no ring outliers.

1 monomer is involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	504	1BF	4	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å <sup>2</sup> )	Q<0.9
1	A	393/406 (96%)	0.23	24 (6%)	28 26	15, 27, 52, 69	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	438	LEU	5.2
1	A	439	ASP	4.2
1	A	373	VAL	4.2
1	A	426	PHE	4.2
1	A	375	THR	3.8
1	A	231	VAL	3.5
1	A	440	MET	3.4
1	A	228	LEU	3.3
1	A	460	HIS	3.3
1	A	58	GLY	3.2
1	A	422	VAL	3.1
1	A	374	ALA	3.0
1	A	423	HIS	3.0
1	A	229	ALA	2.9
1	A	441	GLU	2.8
1	A	206	HIS	2.8
1	A	69	GLY	2.7
1	A	377	GLN	2.6
1	A	174	SER	2.5
1	A	61	VAL	2.4
1	A	332	GLN	2.4
1	A	59	SER	2.3
1	A	421	HIS	2.2
1	A	217	GLY	2.1

## 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

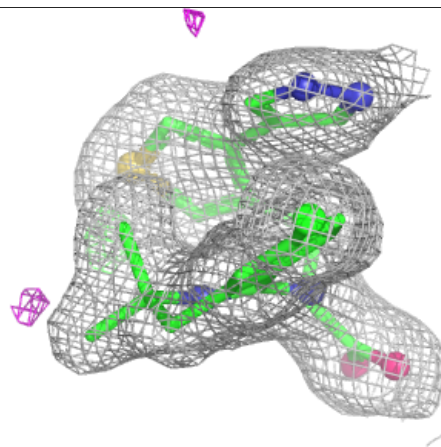
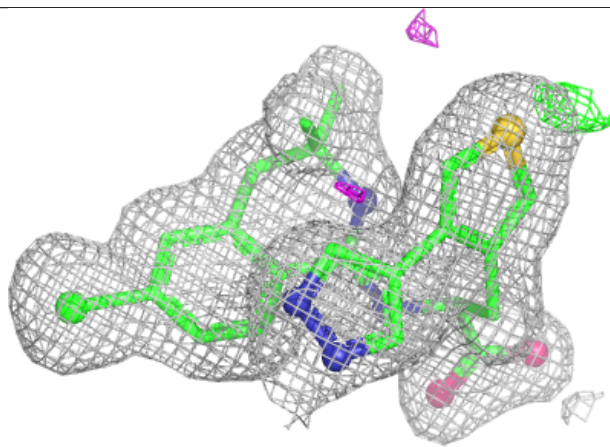
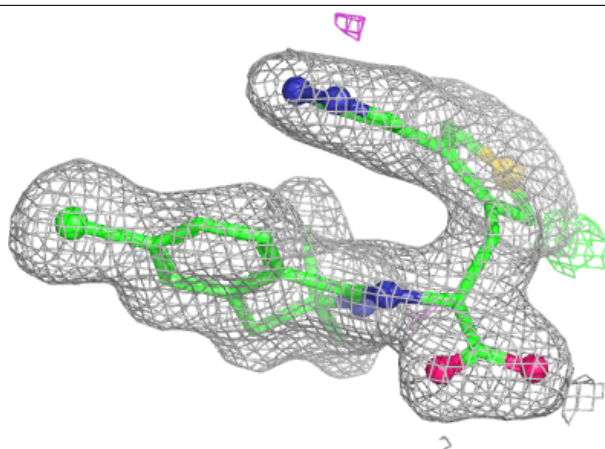
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	ZN	A	502	1/1	0.95	0.08	47,47,47,47	0
3	1BF	A	504	29/29	0.96	0.07	17,22,34,37	0
2	ZN	A	503	1/1	0.98	0.06	37,37,37,37	0
2	ZN	A	501	1/1	1.00	0.03	30,30,30,30	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



**Electron density around 1BF A 504:**

$2mF_o - DF_c$  (at 0.7 rmsd) in gray  
 $mF_o - DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers ⓘ

There are no such residues in this entry.