WORLDWIDE wwPDB PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2022 – 01:46 pm BST

PDB ID	:	7Z78
Title	:	Crystal structure of compound 4 in complex with the bromodomain of human SMARCA2 and pVHL:ElonginC:ElonginB
Authors	:	Bader, G.; Boettcher, J.; Wolkerstorfer, B.
Deposited on	:	2022-03-15
Resolution	:	1.32 Å(reported)

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

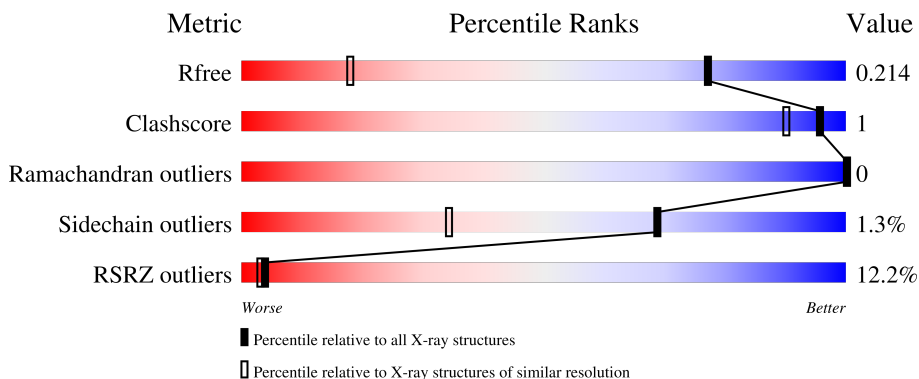|		
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.30
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.30

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 1.32 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	Percentile Ranks	Value
Rfree		0.214
Clashscore		1
Ramachandran outliers		0
Sidechain outliers		1.3%
RSRZ outliers		12.2%

Worse   *Better*

▮ Percentile relative to all X-ray structures

▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1611 (1.34-1.30)
Clashscore	141614	1667 (1.34-1.30)
Ramachandran outliers	138981	1615 (1.34-1.30)
Sidechain outliers	138945	1615 (1.34-1.30)
RSRZ outliers	127900	1580 (1.34-1.30)

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 >=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 <=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	123	7% ... 89% .. 8%
1	B	123	4% ... 89% 11%
1	C	123	22% ... 88% 5% 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	ZN	A	1503	-	-	-	X

2 Entry composition ⓘ

There are 4 unique types of molecules in this entry. The entry contains 3291 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.
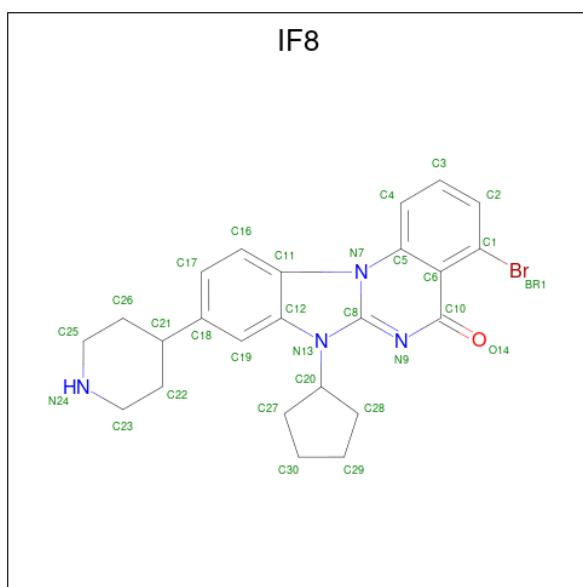
- Molecule 1 is a protein called Probable global transcription activator SNF2L2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	113	Total	C	N	O	S	0	0	0
			927	589	163	172	3			
1	B	110	Total	C	N	O	S	0	0	0
			908	578	160	167	3			
1	C	114	Total	C	N	O	S	0	0	0
			935	595	164	173	3			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1371	SER	-	expression tag	UNP P51531
A	1372	MET	-	expression tag	UNP P51531
B	1371	SER	-	expression tag	UNP P51531
B	1372	MET	-	expression tag	UNP P51531
C	1371	SER	-	expression tag	UNP P51531
C	1372	MET	-	expression tag	UNP P51531

- Molecule 2 is 4-bromanyl-7-cyclopentyl-9-piperidin-4-yl-benzimidazolo[1,2-a]quinazolin-5-one (three-letter code: IF8) (formula: $C_{24}H_{25}BrN_4O$) (labeled as "Ligand of Interest" by depositor).

IF8

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	Br	C	N	O	0	0
			30	1	24	4	1		
2	B	1	Total	Br	C	N	O	0	0
			30	1	24	4	1		
2	C	1	Total	Br	C	N	O	0	0
			30	1	24	4	1		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	B	1	Total	Zn	0	0
			1	1		
3	C	1	Total	Zn	0	0
			1	1		

- Molecule 4 is water.

| Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	161	Total	O	0	0
			161	161		
4	B	168	Total	O	0	0
			168	168		
4	C	98	Total	O	0	0
			98	98		

3 Residue-property plots ⓘ

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.
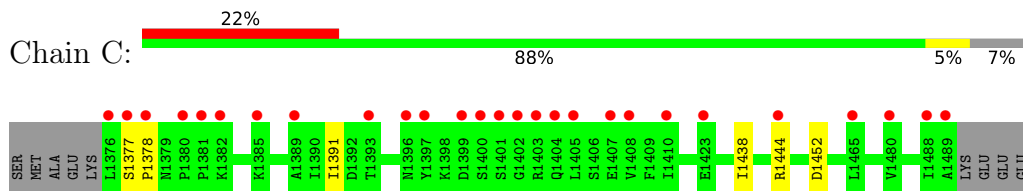
• Molecule 1: Probable global transcription activator SNF2L2
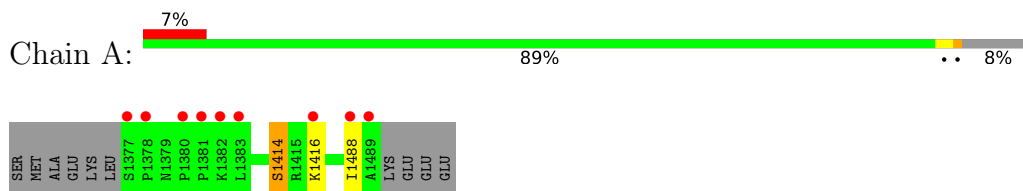
• Molecule 1: Probable global transcription activator SNF2L2
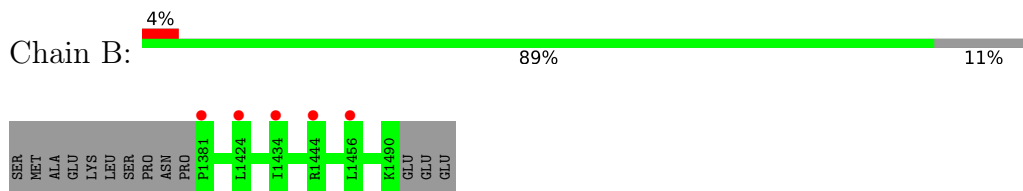
• Molecule 1: Probable global transcription activator SNF2L2

4 Data and refinement statistics ⓘ

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α, β, γ	61.06Å 61.06Å 89.28Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	26.40 – 1.32 25.93 – 1.32	Depositor EDS
% Data completeness (in resolution range)	69.7 (26.40-1.32) 70.4 (25.93-1.32)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.36 (at 1.32Å)	Xtriage
Refinement program	BUSTER 2.11.7	Depositor
R, R_{free}	0.187 , 0.210 0.189 , 0.214	Depositor DCC
R_{free} test set	3172 reflections (5.20%)	wwPDB-VP
Wilson B-factor (Å^2)	16.8	Xtriage
Anisotropy	0.024	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	(Not available) , (Not available)	EDS
L-test for twinning[2]	$< |L| > = 0.49, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.014 for -h,-k,l 0.042 for h,-h-k,-l 0.026 for -k,-h,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3291	wwPDB-VP
Average B, all atoms (Å^2)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.50% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: IF8, 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	0.42	0/943	0.52	0/1267
1	B	0.41	0/922	0.51	0/1234
1	C	0.43	0/951	0.54	0/1278
All	All	0.42	0/2816	0.52	0/3779

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	927	0	955	1	0
1	B	908	0	944	0	0
1	C	935	0	966	2	0
2	A	30	0	0	1	0
2	B	30	0	0	1	0
2	C	30	0	0	1	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
4	A	161	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	168	0	0	0	0
4	C	98	0	0	0	0
All	All	3291	0	2865	6	0

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 1.

All (6) 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:1414:SER:OG	1:A:1416:LYS:HG2	2.11	0.51
1:C:1391:ILE:HG13	1:C:1438:ILE:HD11	1.98	0.45
2:C:1501:IF8:BR1	2:C:1501:IF8:O14	2.90	0.44
2:A:1501:IF8:O14	2:A:1501:IF8:BR1	2.93	0.42
2:B:1501:IF8:O14	2:B:1501:IF8:BR1	2.93	0.41
1:C:1377:SER:H	1:C:1444:ARG:NH2	2.19	0.41

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	111/123 (90%)	111 (100%)	0	0	100	100
1	B	108/123 (88%)	108 (100%)	0	0	100	100
1	C	112/123 (91%)	110 (98%)	2 (2%)	0	100	100
All	All	331/369 (90%)	329 (99%)	2 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	106/115 (92%)	104 (98%)	2 (2%)	57	19
1	B	103/115 (90%)	103 (100%)	0	100	100
1	C	107/115 (93%)	105 (98%)	2 (2%)	57	19
All	All	316/345 (92%)	312 (99%)	4 (1%)	69	34

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1414	SER
1	A	1488	ILE
1	C	1378	PRO
1	C	1452	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\#|Z| > 2$
2	IF8	A	1501	-	35,35,35	0.37	0	41,52,52	0.68	0
2	IF8	B	1501	-	35,35,35	0.37	0	41,52,52	0.66	0
2	IF8	C	1501	-	35,35,35	0.34	0	41,52,52	0.68	0

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
2	IF8	A	1501	-	-	0/8/23/23	0/6/6/6
2	IF8	B	1501	-	-	0/8/23/23	0/6/6/6
2	IF8	C	1501	-	-	0/8/23/23	0/6/6/6

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.
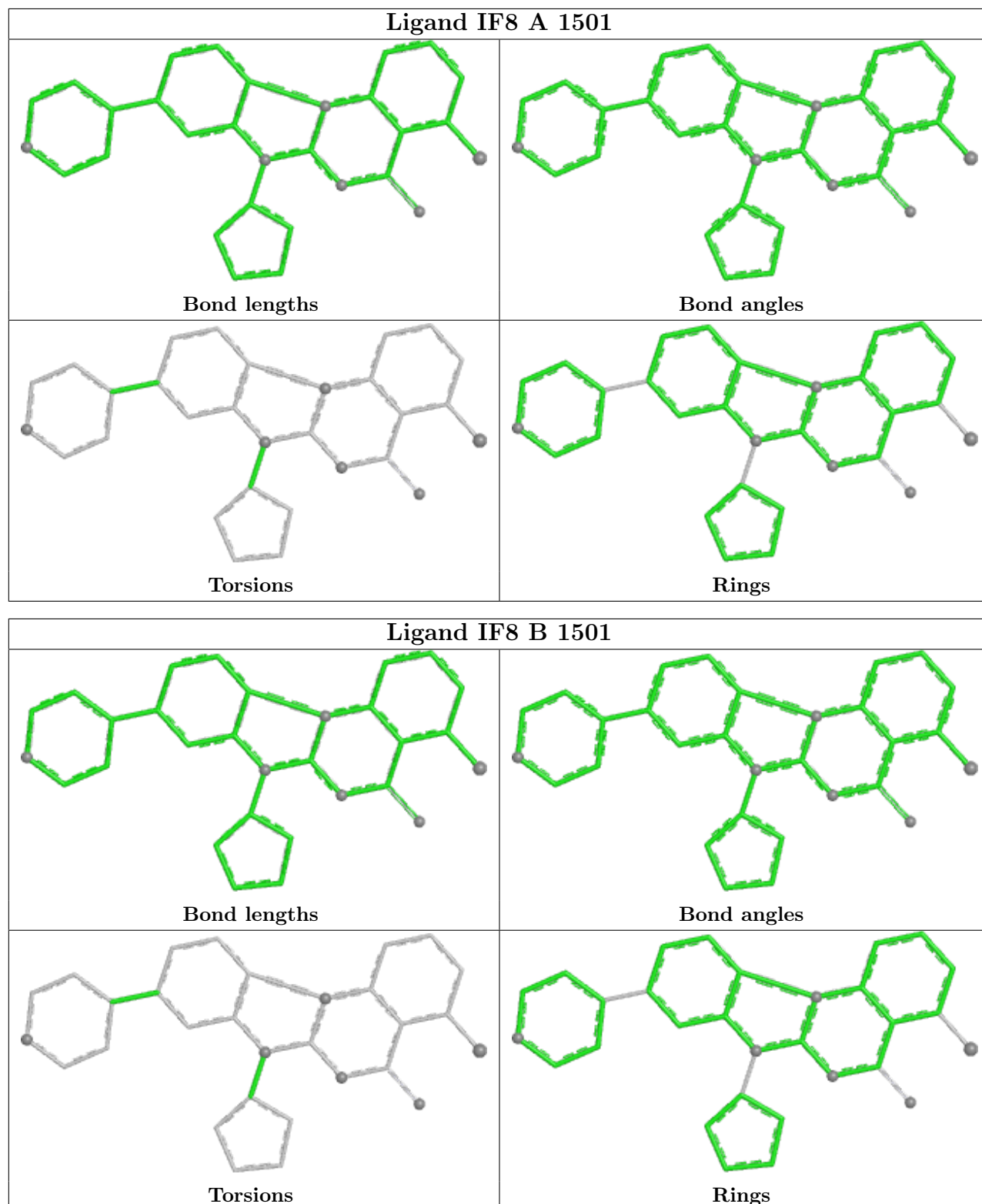
There are no torsion outliers.
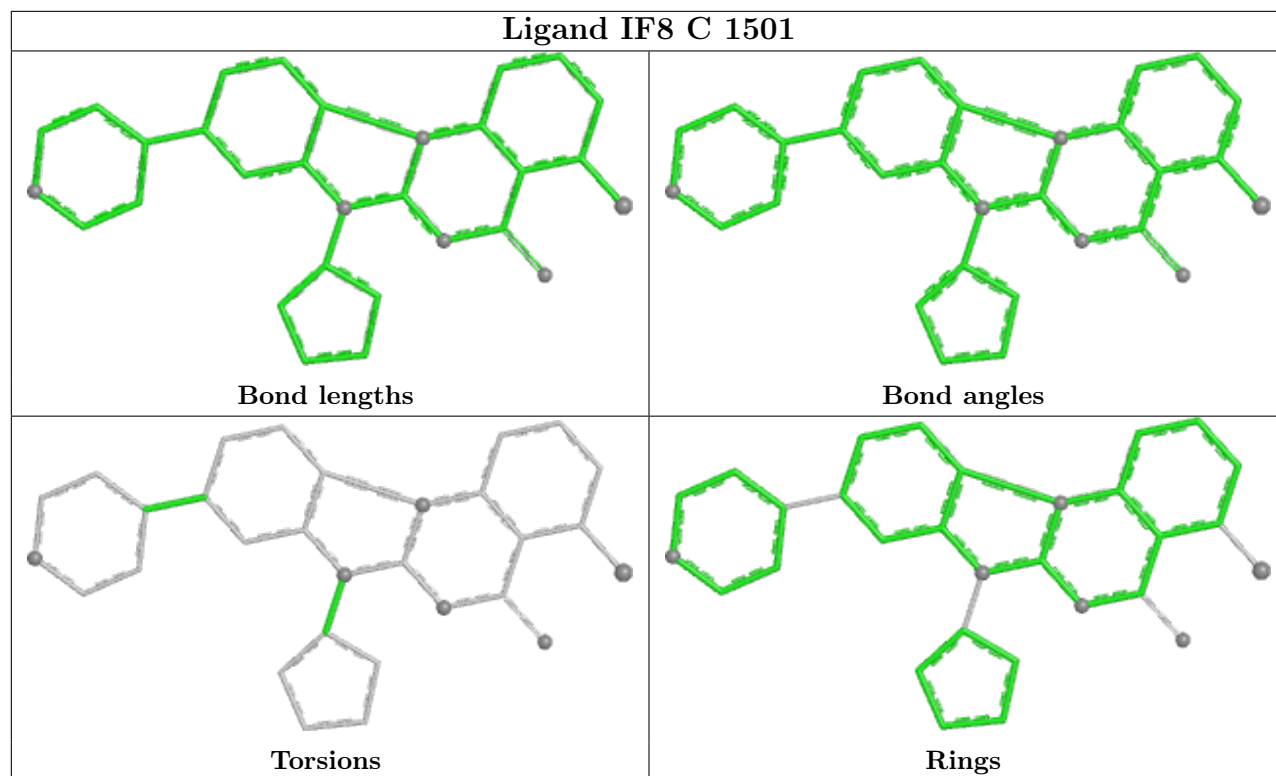
There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1501	IF8	1	0
2	B	1501	IF8	1	0
2	C	1501	IF8	1	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 then 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.

Ligand IF8 A 1501	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand IF8 B 1501	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand IF8 C 1501

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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, 95th 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^2)	Q<0.9
1	A	113/123 (91%)	0.50	9 (7%)	12	10	13, 21, 41, 64	0
1	B	110/123 (89%)	0.36	5 (4%)	33	34	13, 21, 43, 71	0
1	C	114/123 (92%)	1.16	27 (23%)	0	0	16, 33, 72, 92	0
All	All	337/369 (91%)	0.68	41 (12%)	4	3	13, 23, 55, 92	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1488	ILE	8.0
1	A	1488	ILE	6.2
1	C	1400	SER	5.9
1	C	1489	ALA	5.9
1	C	1397	TYR	5.9
1	A	1378	PRO	4.3
1	C	1396	ASN	4.2
1	B	1424	LEU	3.8
1	C	1393	THR	3.7
1	B	1381	PRO	3.5
1	C	1405	LEU	3.5
1	C	1380	PRO	3.4
1	C	1465	LEU	3.3
1	A	1489	ALA	3.3
1	C	1376	LEU	3.3
1	C	1402	GLY	3.2
1	B	1444	ARG	3.1
1	A	1381	PRO	2.8
1	C	1423	GLU	2.8
1	A	1377	SER	2.8
1	A	1380	PRO	2.8
1	C	1404	GLN	2.8
1	C	1381	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	1382	LYS	2.6
1	C	1382	LYS	2.6
1	C	1444	ARG	2.5
1	A	1383	LEU	2.5
1	C	1389	ALA	2.5
1	C	1403	ARG	2.5
1	C	1401	SER	2.5
1	C	1410	ILE	2.5
1	C	1378	PRO	2.4
1	C	1377	SER	2.2
1	C	1408	VAL	2.2
1	B	1456	LEU	2.2
1	C	1399	ASP	2.2
1	C	1385	LYS	2.2
1	C	1407	GLU	2.1
1	A	1416	LYS	2.1
1	C	1480	VAL	2.1
1	B	1434	ILE	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

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, 95th 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.
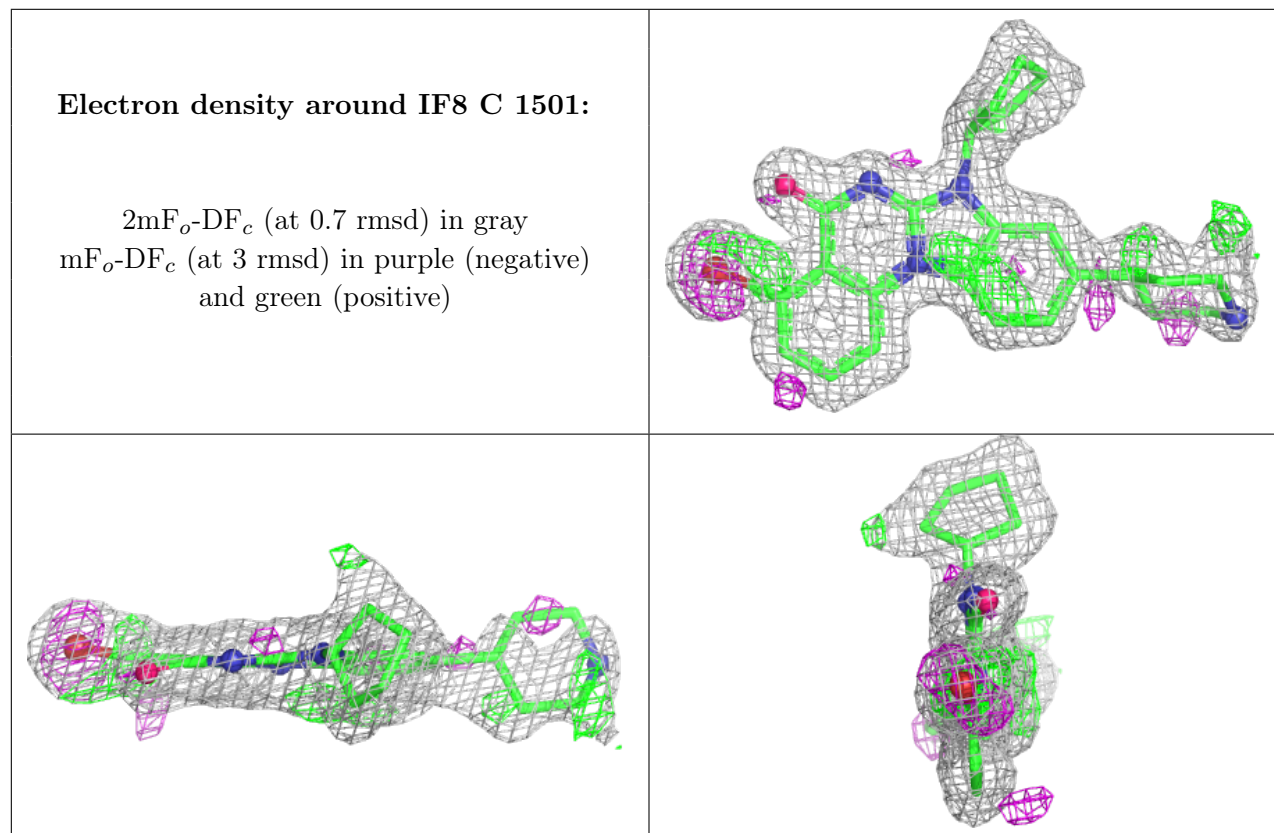
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($Å^2$)	Q<0.9
3	ZN	A	1503	1/1	0.65	0.98	255,255,255,255	0
3	ZN	A	1502	1/1	0.94	0.07	23,23,23,23	0
3	ZN	B	1502	1/1	0.96	0.05	27,27,27,27	0
2	IF8	C	1501	30/30	0.98	0.11	16,23,43,44	0
2	IF8	A	1501	30/30	0.99	0.08	11,15,34,36	0

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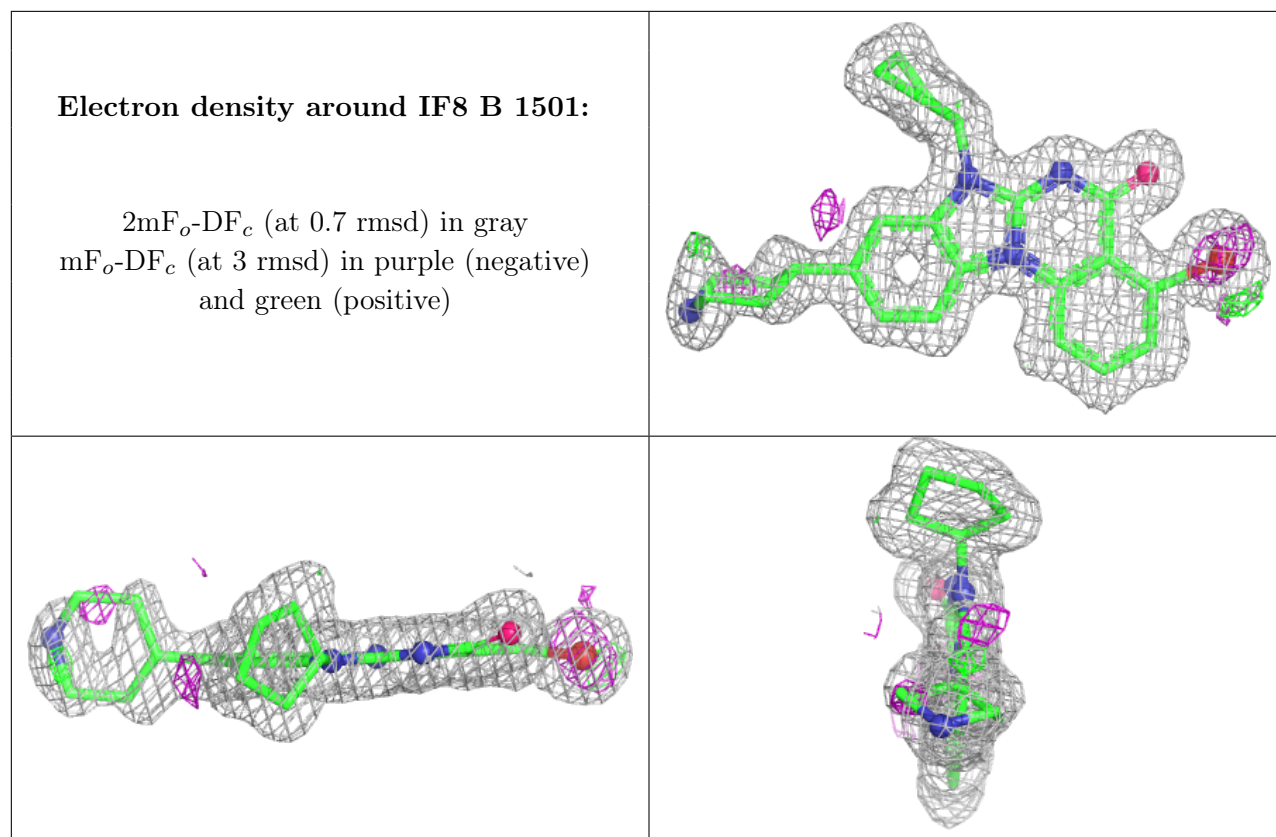| Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	IF8	B	1501	30/30	0.99	0.08	11,15,36,38	0
3	ZN	C	1502	1/1	0.99	0.04	21,21,21,21	0

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 IF8 C 1501:

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

Electron density around IF8 A 1501:

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

Electron density around IF8 B 1501:

$2mF_o$-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.