



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:38 PM UTC

PDB ID : 9GVP / pdb\_00009gvp  
Title : Human PPAR-gamma ligand binding domain in complex with WG115  
Authors : Oerlemans, G.J.M.; Koops, A.A.; Brunsveld, L.  
Deposited on : 2024-09-25  
Resolution : 2.05 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>.

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

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

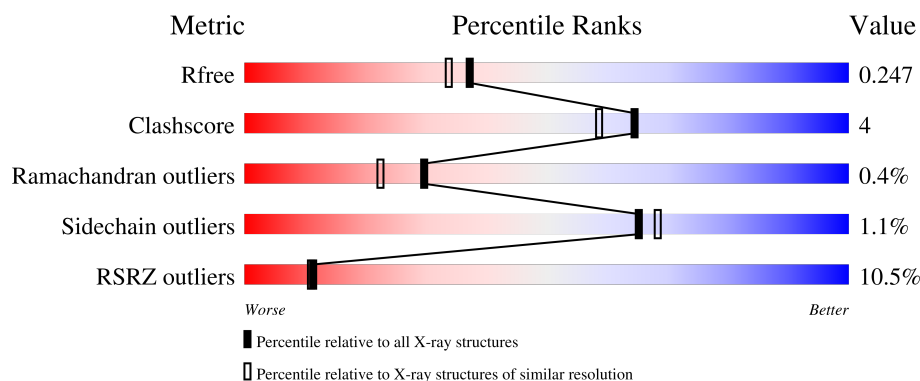
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*



The reported resolution of this entry is 2.05 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))
$R_{free}$	180053	2260 (2.04-2.04)
Clashscore	190562	2333 (2.04-2.04)
Ramachandran outliers	187476	2318 (2.04-2.04)
Sidechain outliers	187428	2318 (2.04-2.04)
RSRZ outliers	180081	2260 (2.04-2.04)

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	273	 9% 88% 8% . .
1	B	273	 11% 86% 9% . .

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4222 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.

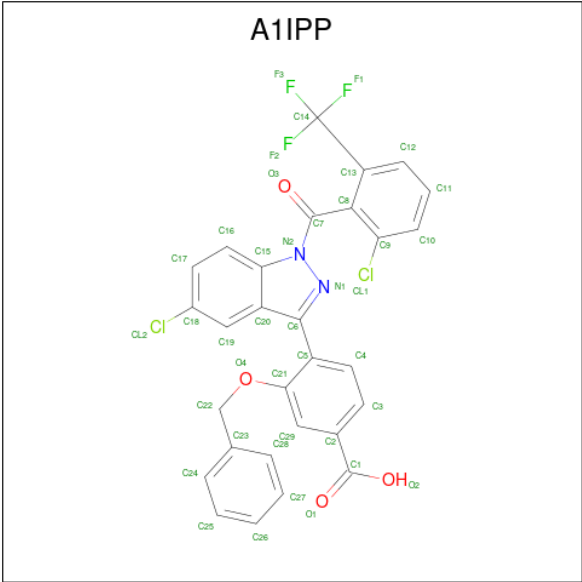
- Molecule 1 is a protein called Peroxisome proliferator-activated receptor gamma.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	262	Total	C	N	O	S	0	0	0
			2024	1308	329	376	11			
1	A	263	Total	C	N	O	S	0	0	0
			2063	1328	334	390	11			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	231	GLY	-	expression tag	UNP P37231
B	232	SER	-	expression tag	UNP P37231
B	233	HIS	-	expression tag	UNP P37231
B	234	MET	-	expression tag	UNP P37231
A	231	GLY	-	expression tag	UNP P37231
A	232	SER	-	expression tag	UNP P37231
A	233	HIS	-	expression tag	UNP P37231
A	234	MET	-	expression tag	UNP P37231

- Molecule 2 is 4-[5-chloranyl-1-[2-chloranyl-6-(trifluoromethyl)phenyl]carbonyl-indazol-3-yl]-3-phenylmethoxy-benzoic acid (CCD ID: A1IPP) (formula: C<sub>29</sub>H<sub>17</sub>Cl<sub>2</sub>F<sub>3</sub>N<sub>2</sub>O<sub>4</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	
2	B	1	Total	C	Cl	F	N	O	0	0
			40	29	2	3	2	4		
2	A	1	Total	C	Cl	F	N	O	0	0
			40	29	2	3	2	4		

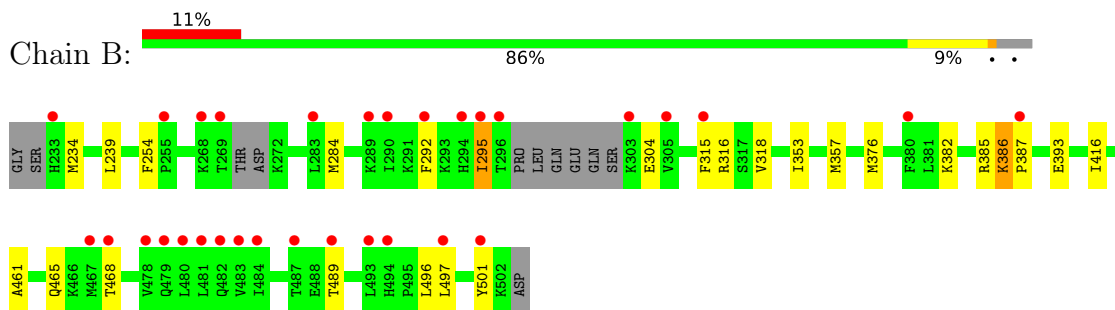
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	27	Total	O	0	0
			27	27		
3	A	28	Total	O	0	0
			28	28		

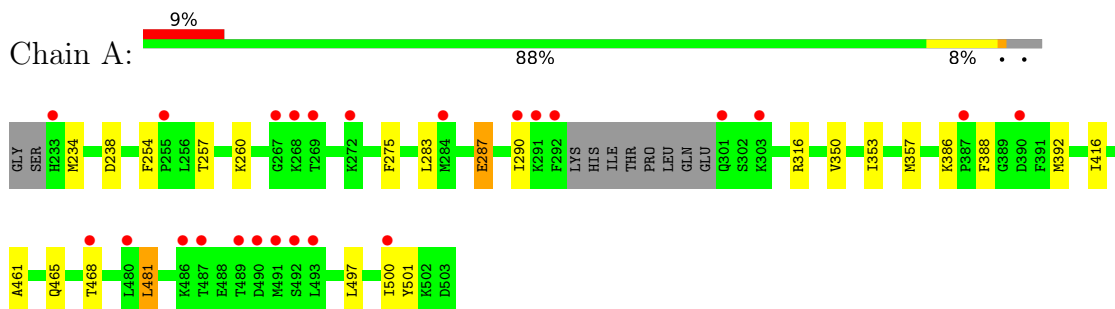
### 3 Residue-property plots [i](#)

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: Peroxisome proliferator-activated receptor gamma



- Molecule 1: Peroxisome proliferator-activated receptor gamma



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.11Å 61.73Å 119.41Å 90.00° 101.37° 90.00°	Depositor
Resolution (Å)	58.53 – 2.05 58.53 – 2.05	Depositor EDS
% Data completeness (in resolution range)	98.1 (58.53-2.05) 98.2 (58.53-2.05)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.33 (at 2.05Å)	Xtriage
Refinement program	PHENIX 1.21.2-5419	Depositor
R, $R_{free}$	0.223 , 0.249 (Not available) , 0.247	Depositor DCC
$R_{free}$ test set	2065 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	41.7	Xtriage
Anisotropy	0.305	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 47.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	4222	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.91% 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 [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: A1IPP

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	0.36	0/2098	0.53	0/2835
1	B	0.36	0/2058	0.54	0/2785
All	All	0.36	0/4156	0.54	0/5620

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
1	B	0	1
All	All	0	2

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	316	ARG	Sidechain
1	B	316	ARG	Sidechain

### 5.2 Too-close contacts [i](#)

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	2063	0	2059	15	0
1	B	2024	0	1998	18	0
2	A	40	0	0	0	0
2	B	40	0	0	0	0
3	A	28	0	0	0	0
3	B	27	0	0	0	0
All	All	4222	0	4057	32	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 4.

All (32) 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:353:ILE:HG23	1:A:416:ILE:HD12	1.73	0.71
1:B:353:ILE:HG23	1:B:416:ILE:HD12	1.82	0.60
1:B:465:GLN:O	1:B:468:THR:HB	2.03	0.59
1:B:304:GLU:OE1	1:B:385:ARG:NH2	2.35	0.58
1:A:234:MET:HE2	1:A:238:ASP:HB3	1.86	0.57
1:B:468:THR:HA	1:A:468:THR:HG23	1.86	0.56
1:A:461:ALA:O	1:A:465:GLN:HG3	2.06	0.55
1:A:257:THR:HG23	1:A:260:LYS:H	1.74	0.52
1:A:465:GLN:O	1:A:468:THR:HB	2.09	0.52
1:B:382:LYS:HE2	1:B:393:GLU:CD	2.35	0.51
1:B:304:GLU:OE2	1:B:385:ARG:NE	2.35	0.51
1:B:461:ALA:O	1:B:465:GLN:HG3	2.11	0.51
1:A:257:THR:CG2	1:A:260:LYS:H	2.26	0.48
1:A:254:PHE:CG	1:A:357:MET:HE1	2.49	0.48
1:B:234:MET:HE3	1:B:239:LEU:HD23	1.95	0.48
1:B:292:PHE:HB2	1:B:295:ILE:HG12	1.95	0.47
1:B:385:ARG:HH11	1:B:386:LYS:CB	2.28	0.47
1:B:318:VAL:HG21	1:B:501:TYR:HD2	1.79	0.47
1:A:350:VAL:HG21	1:A:500:ILE:HD13	1.97	0.46
1:A:283:LEU:O	1:A:287:GLU:HG2	2.16	0.46
1:A:388:PHE:O	1:A:392:MET:HG2	2.15	0.46
1:B:254:PHE:CD1	1:B:357:MET:HE1	2.52	0.45
1:B:318:VAL:HG21	1:B:501:TYR:CD2	2.51	0.45
1:B:497:LEU:HD23	1:B:497:LEU:HA	1.77	0.44
1:A:497:LEU:HD23	1:A:497:LEU:HA	1.80	0.44
1:B:386:LYS:CB	1:B:387:PRO:CD	2.97	0.43
1:A:481:LEU:HD23	1:A:481:LEU:HA	1.83	0.42
1:B:284:MET:HE2	1:B:284:MET:HB3	1.88	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:481:LEU:HD12	1:A:501:TYR:CZ	2.55	0.42
1:B:376:MET:HE3	1:B:376:MET:HB2	1.94	0.42
1:A:275:PHE:HB3	1:A:290:ILE:HD11	2.02	0.41
1:B:315:PHE:CD2	1:B:315:PHE:C	3.00	0.40

There are no symmetry-related clashes.

## 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	259/273 (95%)	256 (99%)	2 (1%)	1 (0%)	30	22
1	B	256/273 (94%)	253 (99%)	2 (1%)	1 (0%)	30	22
All	All	515/546 (94%)	509 (99%)	4 (1%)	2 (0%)	30	22

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	386	LYS
1	A	386	LYS

### 5.3.2 Protein sidechains [i](#)

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	225/245 (92%)	223 (99%)	2 (1%)	70	75
1	B	214/245 (87%)	211 (99%)	3 (1%)	59	60
All	All	439/490 (90%)	434 (99%)	5 (1%)	65	68

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	295	ILE
1	B	489	THR
1	B	496	LEU
1	A	287	GLU
1	A	481	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	233	HIS
1	B	440	ASN
1	B	458	GLN
1	B	498	GLN
1	A	373	GLN

### 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

2 ligands are modelled in this entry.

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	A1IPP	B	601	-	44,44,44	0.52	0	61,65,65	0.76	2 (3%)
2	A1IPP	A	601	-	44,44,44	0.56	0	61,65,65	0.76	1 (1%)

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	A1IPP	B	601	-	-	0/27/27/27	0/5/5/5
2	A1IPP	A	601	-	-	2/27/27/27	0/5/5/5

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	601	A1IPP	C20-C15-N2	-2.83	105.22	106.40
2	B	601	A1IPP	C20-C15-N2	-2.77	105.24	106.40
2	B	601	A1IPP	C9-C8-C7	-2.50	118.19	121.19

There are no chirality outliers.

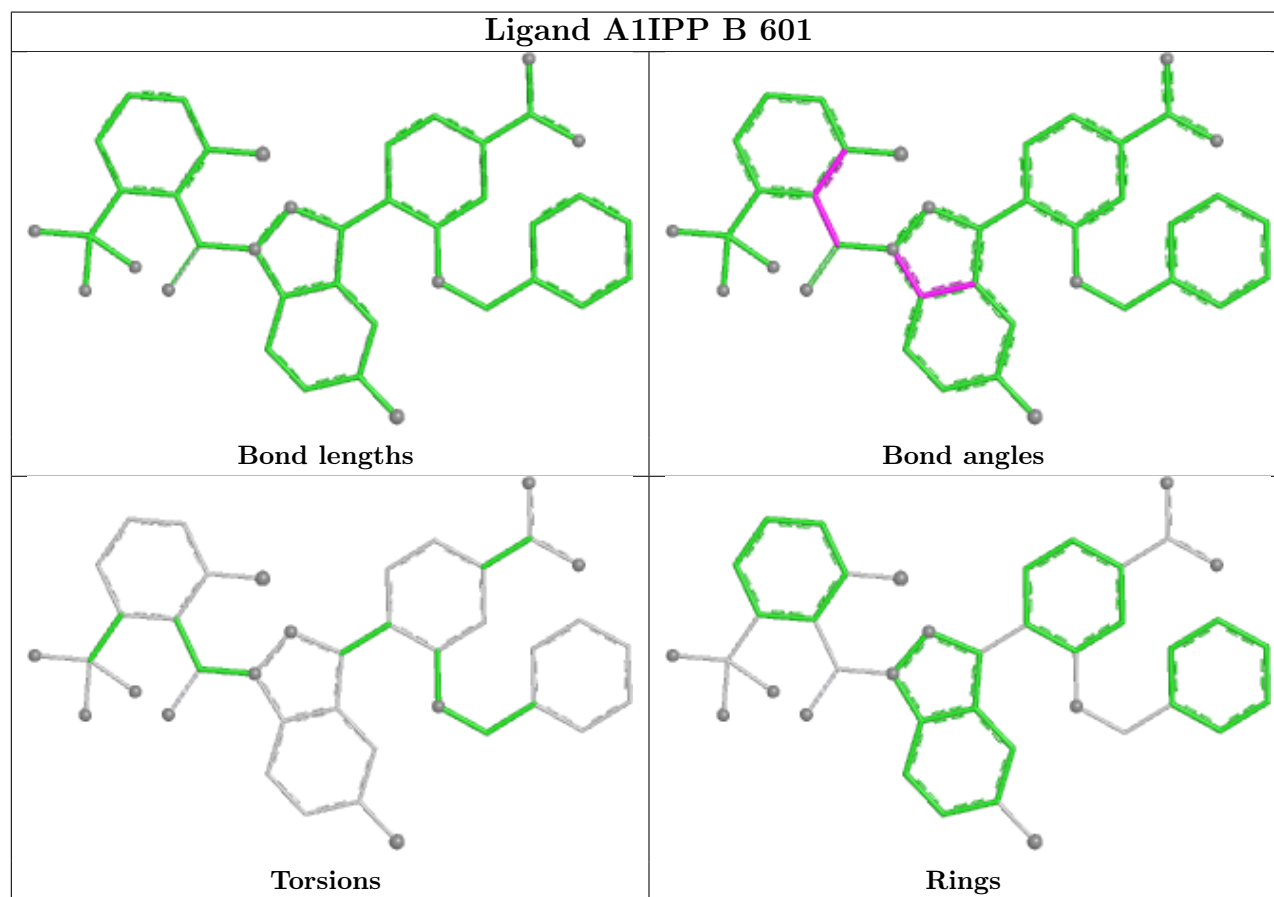
All (2) torsion outliers are listed below:

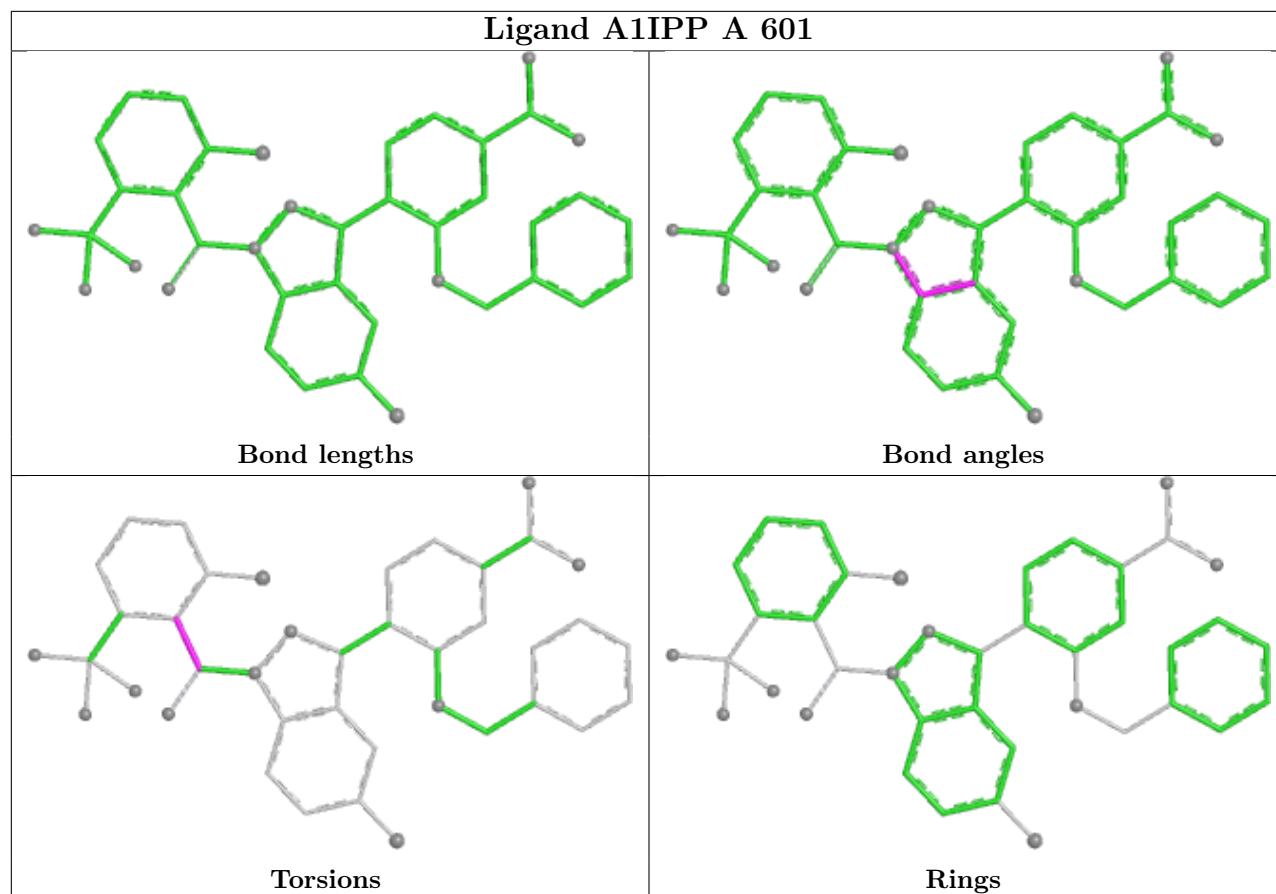
Mol	Chain	Res	Type	Atoms
2	A	601	A1IPP	N2-C7-C8-C9
2	A	601	A1IPP	O3-C7-C8-C9

There are no ring outliers.

No monomer is involved in short contacts.

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	263/273 (96%)	0.61	24 (9%) 15 15	35, 51, 91, 133	0
1	B	262/273 (95%)	0.67	31 (11%) 9 8	33, 53, 100, 141	0
All	All	525/546 (96%)	0.64	55 (10%) 11 11	33, 52, 98, 141	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	292	PHE	6.4
1	B	295	ILE	4.4
1	B	489	THR	4.4
1	B	468	THR	4.3
1	A	284	MET	4.2
1	B	483	VAL	4.2
1	B	290	ILE	4.1
1	B	292	PHE	4.0
1	B	315	PHE	4.0
1	B	294	HIS	3.9
1	B	269	THR	3.9
1	A	291	LYS	3.8
1	A	290	ILE	3.7
1	A	490	ASP	3.7
1	B	480	LEU	3.6
1	B	467	MET	3.6
1	A	303	LYS	3.4
1	B	493	LEU	3.4
1	A	269	THR	3.3
1	B	289	LYS	3.3
1	A	468	THR	3.3
1	A	486	LYS	3.3
1	B	482	GLN	3.3
1	B	305	VAL	3.2

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Mol	Chain	Res	Type	RSRZ
1	B	296	THR	3.2
1	B	268	LYS	3.1
1	A	255	PRO	3.0
1	B	479	GLN	3.0
1	A	487	THR	3.0
1	B	233	HIS	2.9
1	B	484	ILE	2.9
1	B	501	TYR	2.9
1	B	387	PRO	2.8
1	A	489	THR	2.7
1	B	255	PRO	2.6
1	A	390	ASP	2.6
1	A	267	GLY	2.6
1	A	491	MET	2.6
1	A	268	LYS	2.5
1	B	303	LYS	2.5
1	B	481	LEU	2.4
1	A	500	ILE	2.3
1	A	493	LEU	2.3
1	B	478	VAL	2.2
1	B	494	HIS	2.2
1	A	301	GLN	2.1
1	B	487	THR	2.1
1	A	233	HIS	2.1
1	B	497	LEU	2.1
1	A	387	PRO	2.1
1	B	283	LEU	2.1
1	B	380	PHE	2.1
1	A	480	LEU	2.0
1	A	272	LYS	2.0
1	A	492	SER	2.0

## 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](#)

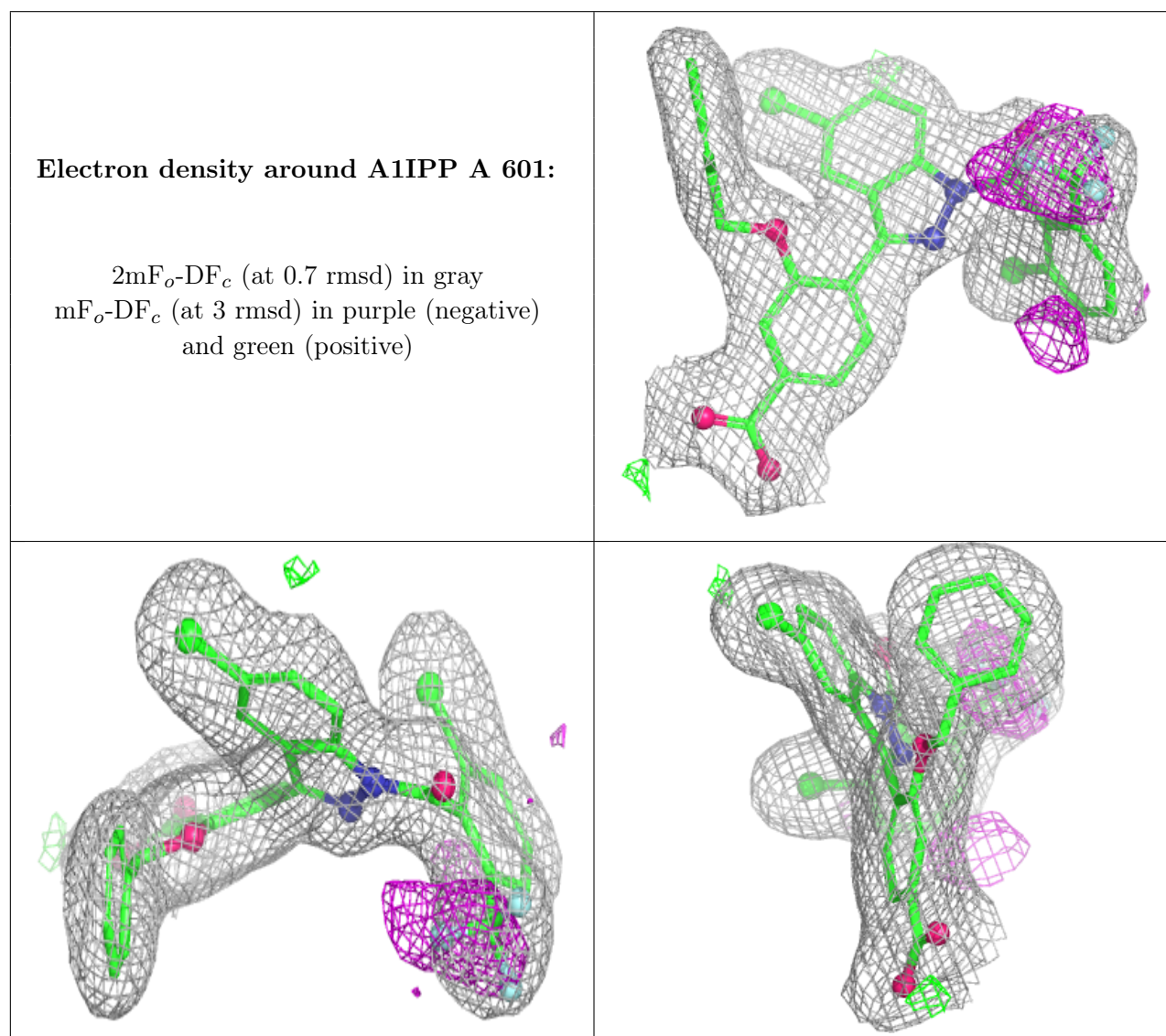
There are no oligosaccharides in this entry.

## 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	A1IPP	A	601	40/40	0.93	0.10	46,51,64,65	0
2	A1IPP	B	601	40/40	0.94	0.09	50,61,69,74	0

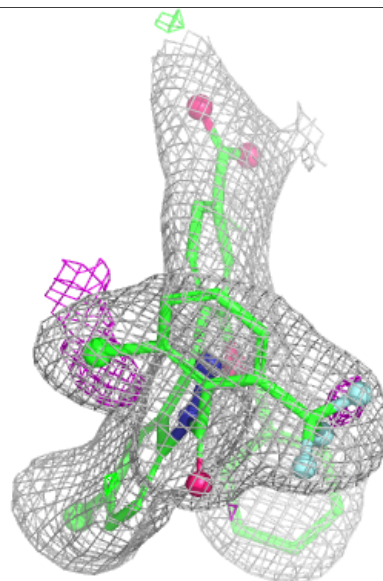
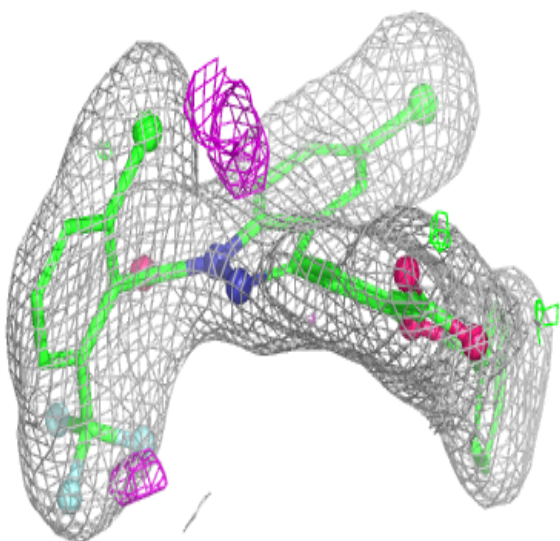
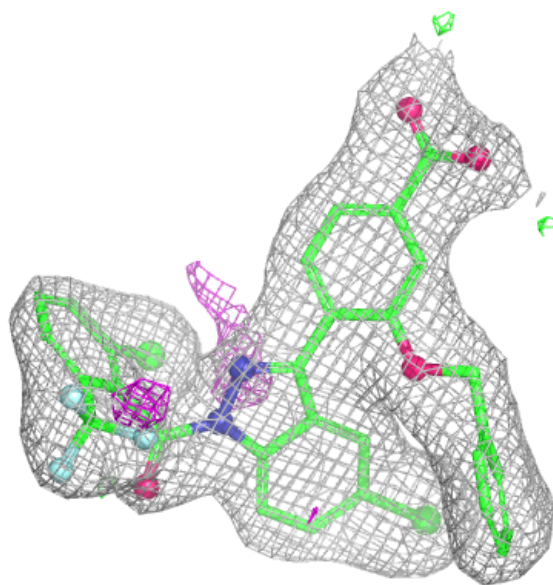
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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





**Electron density around A1IPP B 601:**

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



## 6.5 Other polymers [i](#)

There are no such residues in this entry.