

wwPDB X-ray Structure Validation Summary Report (i)

Apr 23, 2024 – 12:20 PM JST

PDB ID	:	8HKN
Title	:	Mutated human ADP-ribosyltransferase 2 (PARP2) catalytic domain bound
		to approved drug Fluzoparib
Authors	:	Wang, X.Y.; Zhou, J.; Xu, B.L.
Deposited on	:	2022-11-27
Resolution	:	2.50 Å(reported)

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

We welcome your comments at 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 (i) 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 (1)) were used in the production of this report:

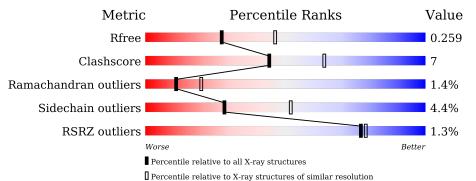
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 2.50 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R _{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559(2.50-2.50)

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	А	353	2% 8 6%	13%	:
1	В	353	% 	18%	•



8HKN

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 5913 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Р	353	Total	С	Ν	0	S	0	0	0
	D	202	2802	1777	486	520	19	0		
1	Δ	353	Total	С	Ν	0	S	0	0	0
	A	ამა	2802	1777	486	520	19	0	U	U

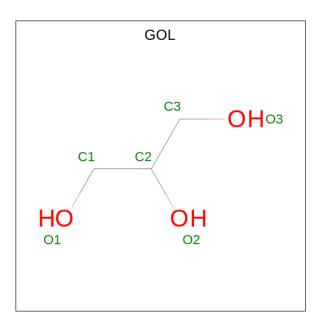
• Molecule 1 is a protein called Poly [ADP-ribose] polymerase 2.

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	229	GLY	-	expression tag	UNP Q9UGN5
В	349	SER	THR	engineered mutation	UNP Q9UGN5
В	351	ARG	LEU	engineered mutation	UNP Q9UGN5
В	353	GLY	SER	engineered mutation	UNP Q9UGN5
В	354	LEU	PRO	engineered mutation	UNP Q9UGN5
А	229	GLY	-	expression tag	UNP Q9UGN5
А	349	SER	THR	engineered mutation	UNP Q9UGN5
А	351	ARG	LEU	engineered mutation	UNP Q9UGN5
А	353	GLY	SER	engineered mutation	UNP Q9UGN5
А	354	LEU	PRO	engineered mutation	UNP Q9UGN5

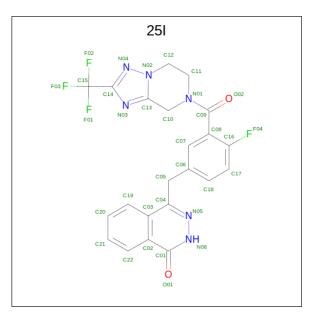
• Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0
2	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 6 & 3 & 3 \end{array}$	0	0

• Molecule 3 is Fluzoparib (three-letter code: 25I) (formula: $C_{22}H_{16}F_4N_6O_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	В	1	Total 34	C 22	F 4	N 6	O 2	0	0

Continued on next page...



Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
2	Λ	1	Total	С	F	Ν	Ο	0	0
3	A	1	34	22	4	6	2	0	0

• Molecule 4 is water.

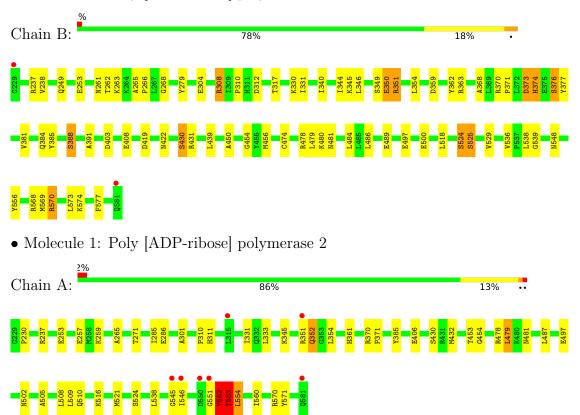
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	В	110	Total O 110 110	0	0
4	А	119	Total O 119 119	0	0



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: Poly [ADP-ribose] polymerase 2





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	41.01Å 85.58Å 95.41Å	Deperitor
a, b, c, α , β , γ	90.00° 90.03° 90.00°	Depositor
Resolution (Å)	36.98 - 2.50	Depositor
Resolution (A)	95.41 - 2.50	EDS
% Data completeness	99.3 (36.98-2.50)	Depositor
(in resolution range)	99.3 (95.41 - 2.50)	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$6.43 (at 2.51 \text{\AA})$	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
D D	0.181 , 0.261	Depositor
R, R_{free}	0.179 , 0.259	DCC
R_{free} test set	1102 reflections (4.84%)	wwPDB-VP
Wilson B-factor $(Å^2)$	40.2	Xtriage
Anisotropy	0.026	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35 , 49.9	EDS
L-test for twinning ²	$< L > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.024 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5913	wwPDB-VP
Average B, all atoms $(Å^2)$	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 5.97% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



¹Intensities estimated from amplitudes.

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: GOL, $25\mathrm{I}$

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		
	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.41	0/2861	0.75	0/3860	
1	В	0.42	0/2861	0.75	0/3860	
All	All	0.41	0/5722	0.75	0/7720	

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	В	0	3

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	В	308	ARG	Sidechain
1	В	431	ARG	Sidechain
1	В	478	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	А	2802	0	2808	29	0
1	В	2802	0	2808	45	0
2	А	6	0	8	1	0
2	В	6	0	8	0	0
3	А	34	0	0	1	0
3	В	34	0	0	0	0
4	А	119	0	0	2	0
4	В	110	0	0	3	0
All	All	5913	0	5632	75	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 7.

The worst 5 of 75 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:B:384:GLN:O	1:B:388:SER:HB3	1.62	0.99
1:B:439:LEU:HD23	1:B:569:MET:HE1	1.66	0.78
1:A:502:ASN:ND2	1:A:508:LEU:HD11	2.00	0.76
1:A:370:ARG:NH1	1:A:406:GLU:OE2	2.20	0.74
1:B:374:HIS:O	4:B:701:HOH:O	2.06	0.73

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	А	351/353~(99%)	332~(95%)	14 (4%)	5(1%)	11 20
1	В	351/353~(99%)	328~(93%)	18 (5%)	5(1%)	11 20
All	All	702/706~(99%)	660 (94%)	32~(5%)	10 (1%)	11 20

5 of 10 Ramachandran outliers are listed below:



Mol	Chain	Res	Type
1	В	350	GLU
1	В	376	SER
1	В	525	SER
1	А	352	GLN
1	А	552	TYR

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent side chain 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	Percentiles		
1	А	305/305~(100%)	295~(97%)	10 (3%)	38 64	
1	В	305/305~(100%)	288 (94%)	17 (6%)	21 40	
All	All	610/610~(100%)	583~(96%)	27~(4%)	28 52	

5 of 27 residues with a non-rotameric side chain are listed below:

Mol	Chain	\mathbf{Res}	Type
1	В	480	LYS
1	А	271	THR
1	А	552	TYR
1	В	570	ARG
1	А	345	LYS

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such side chains are listed below:

Mol	Chain	Res	Type
1	В	422	ASN
1	А	361	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.



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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

4 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			
	Type	Chain	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	25I	В	602	-	$35,\!38,\!38$	1.18	5 (14%)	42,57,57	1.75	9 (21%)
2	GOL	А	601	-	$5,\!5,\!5$	1.21	0	$5,\!5,\!5$	0.53	0
3	25I	А	602	-	35,38,38	1.14	1 (2%)	42,57,57	1.96	6 (14%)
2	GOL	В	601	-	$5,\!5,\!5$	1.33	1 (20%)	5,5,5	0.70	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
3	25I	В	602	-	-	0/18/27/27	0/4/5/5
2	GOL	А	601	-	-	3/4/4/4	-
3	25I	А	602	-	-	0/18/27/27	0/4/5/5
2	GOL	В	601	-	-	0/4/4/4	-

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
3	А	602	25I	C04-N05	4.59	1.32	1.29
3	В	602	25I	C13-N03	-2.45	1.30	1.35

Continued on next page...



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	$\mathrm{Ideal}(\mathrm{\AA})$
3	В	602	25I	C21-C20	2.33	1.44	1.38
3	В	602	25I	C05-C06	-2.32	1.47	1.51
2	В	601	GOL	O2-C2	2.11	1.49	1.43

Continued from previous page...

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
3	А	602	25I	C14-N03-C13	-9.76	100.28	104.05
3	В	602	25I	C14-N03-C13	-5.56	101.90	104.05
3	В	602	25I	C17-C16-C08	-4.15	118.59	123.11
3	В	602	25I	C11-C12-N02	-3.73	101.47	109.06
3	В	602	25I	C18-C17-C16	3.08	122.77	119.05

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	А	601	GOL	C1-C2-C3-O3
2	А	601	GOL	O2-C2-C3-O3
2	А	601	GOL	O1-C1-C2-O2

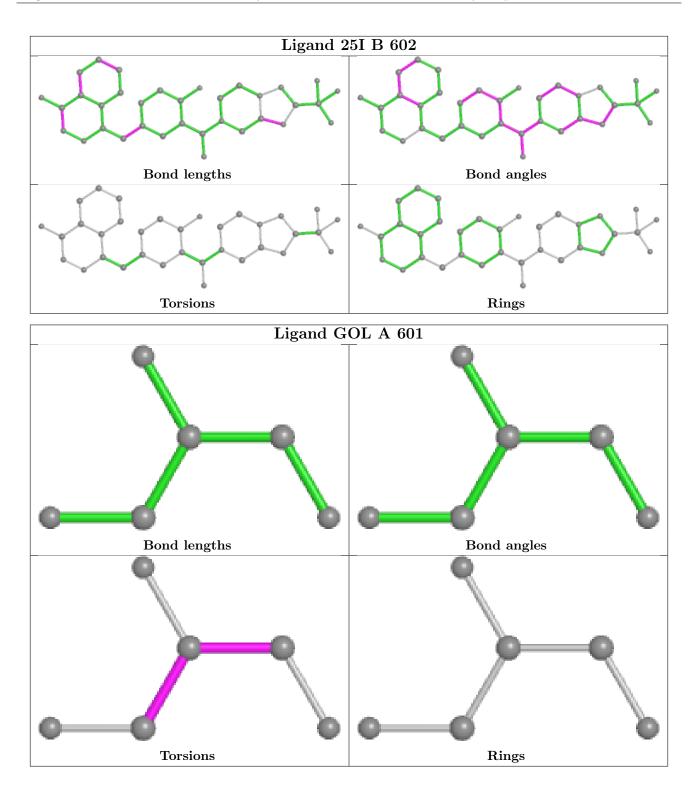
There are no ring outliers.

2 monomers are involved in 1 short contact:

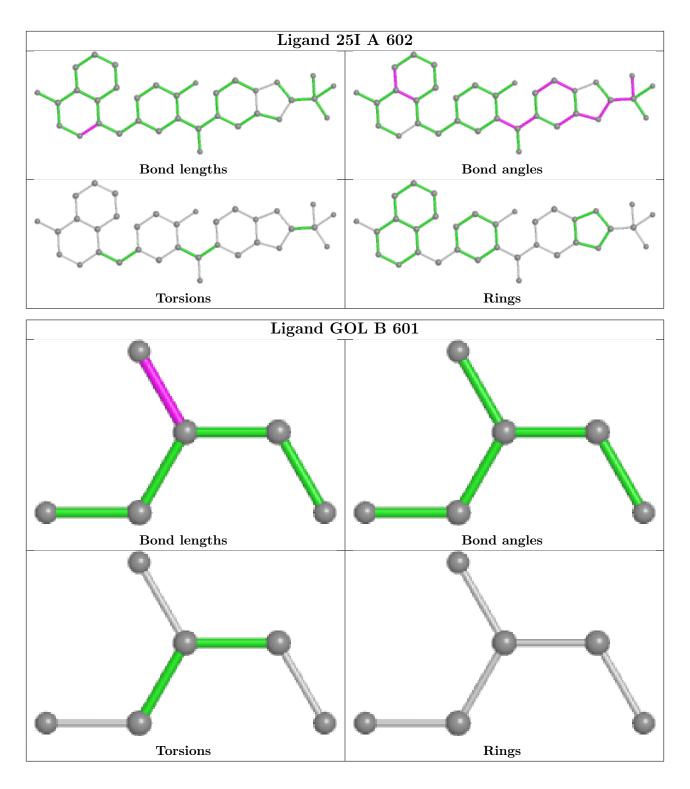
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	А	601	GOL	1	0
3	А	602	25I	1	0

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 sufficient 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 (i)

6.1 Protein, DNA and RNA chains (i)

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^{th} 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	$\langle RSRZ \rangle$	#RSRZ>2		$OWAB(Å^2)$	Q<0.9	
1	А	353/353~(100%)	-0.29	7 (1%)	65	68	24, 41, 72, 126	0
1	В	353/353~(100%)	-0.27	2 (0%)	89	90	21, 42, 75, 106	0
All	All	706/706~(100%)	-0.28	9 (1%)	77	79	21, 42, 75, 126	0

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	581	GLN	4.2
1	А	351	ARG	3.4
1	В	229	GLY	3.3
1	А	551	GLY	3.2
1	А	545	GLY	2.3

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.

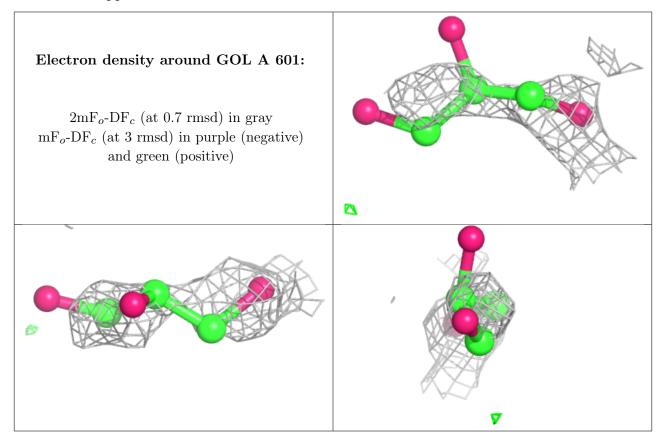
6.4 Ligands (i)

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^{th} 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.

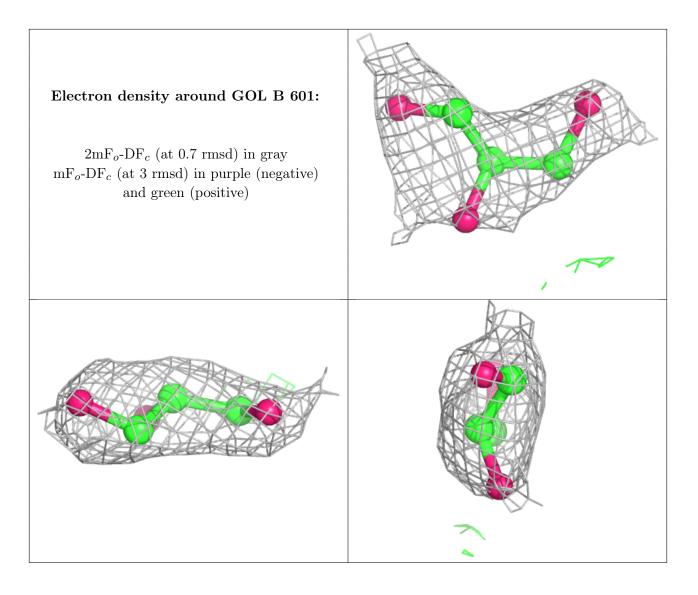


Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q < 0.9
2	GOL	А	601	6/6	0.89	0.52	57,80,84,93	0
2	GOL	В	601	6/6	0.90	0.18	53,58,60,62	0
3	25I	А	602	34/34	0.97	0.11	24,28,39,44	0
3	25I	В	602	34/34	0.98	0.11	21,25,43,45	0

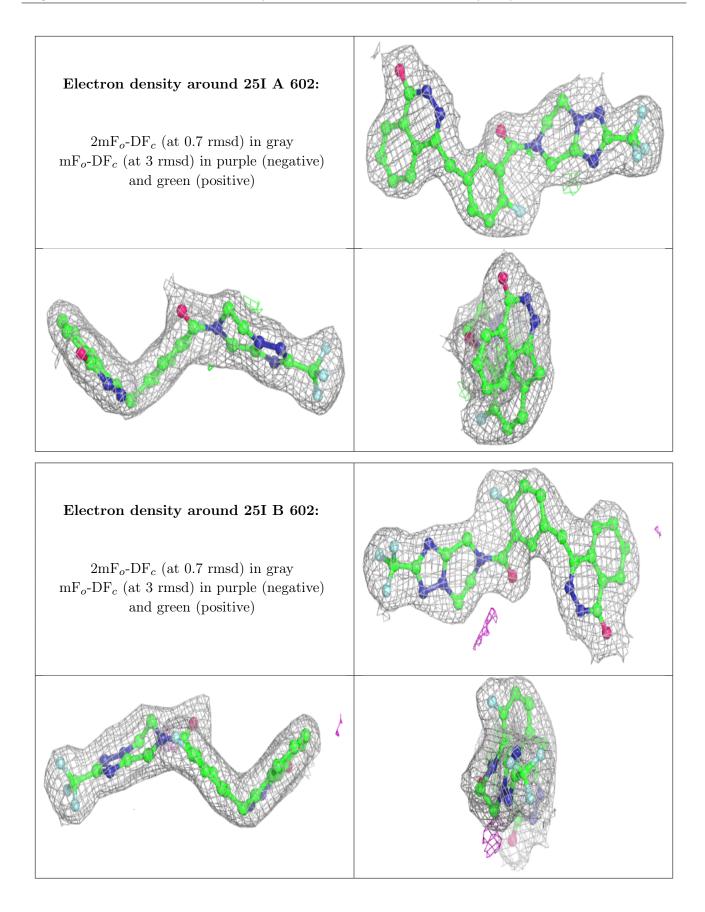
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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













6.5 Other polymers (i)

There are no such residues in this entry.

