

wwPDB X-ray Structure Validation Summary Report (i)

Jan 17, 2022 - 04:38 pm GMT

	: 6TTG	
Title	: Crystal structure of the ATP binding domain of S. aureu	us GyrB complexed
	with LMD62	
Authors	: Welin, M.; Kimbung, R.; Focht, D.	
Deposited on		
Resolution	: 1.70 Å(reported)	
Authors Deposited on	with LMD62 : Welin, M.; Kimbung, R.; Focht, D.	ıs GyrB complex

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 following versions of software and data (see references (1)) were used in the production of this report:

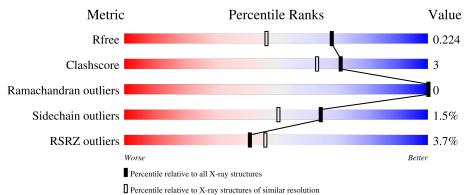
MolProbity		4 021 467
Mogul	:	1.8.4, CSD as 541 be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.24
buster-report	:	1.1.7(2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0267
CCP4	:	7.1.010 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.24

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 1.70 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	4298 (1.70-1.70)
Clashscore	141614	4695 (1.70-1.70)
Ramachandran outliers	138981	4610 (1.70-1.70)
Sidechain outliers	138945	4610 (1.70-1.70)
RSRZ outliers	127900	4222 (1.70-1.70)

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	212	2% 8 5%	5%	10%
1	В	212	4% 83% 5%	6•	12%



2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 3431 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	Λ	191	Total	С	N O S O	0	4	0		
	A	191	1558	979	269	308	2	0	4	0
1	В	187	Total	С	Ν	0	S	0	Б	0
	D	107	1542	966	266	308	2	0	5	0

• Molecule 1 is a protein called DNA gyrase subunit B.

Chain	Residue	Modelled	Actual	Comment	Reference
А	0	GLY	-	expression tag	UNP P0A0K8
А	1	SER	-	expression tag	UNP P0A0K8
А	?	-	VAL	deletion	UNP P0A0K8
А	?	-	LEU	deletion	UNP P0A0K8
А	?	-	HIS	deletion	UNP P0A0K8
А	?	-	ALA	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	LYS	deletion	UNP P0A0K8
А	?	-	PHE	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	TYR	deletion	UNP P0A0K8
А	?	-	LYS	deletion	UNP P0A0K8
А	?	-	VAL	deletion	UNP P0A0K8
А	?	-	SER	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	LEU	deletion	UNP P0A0K8
А	?	-	HIS	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
А	?	-	VAL	deletion	UNP P0A0K8
А	?	-	GLY	deletion	UNP P0A0K8
				e .	-

There are 50 discrepancies between the modelled and reference sequences:

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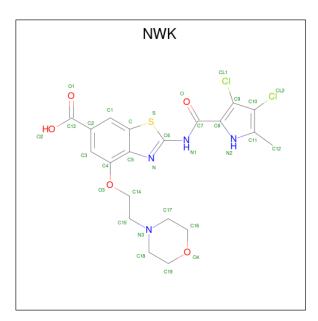


Chain	Residue	Modelled	Actual	Comment	Reference
В	0	GLY	-	expression tag	UNP P0A0K8
В	1	SER	-	expression tag	UNP P0A0K8
В	?	_	VAL	deletion	UNP P0A0K8
В	?	_	LEU	deletion	UNP P0A0K8
В	?	-	HIS	deletion	UNP P0A0K8
В	?	-	ALA	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	LYS	deletion	UNP P0A0K8
В	?	-	PHE	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	TYR	deletion	UNP P0A0K8
В	?	-	LYS	deletion	UNP P0A0K8
В	?	-	VAL	deletion	UNP P0A0K8
В	?	-	SER	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	LEU	deletion	UNP P0A0K8
В	?	-	HIS	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8
В	?	-	VAL	deletion	UNP P0A0K8
В	?	-	GLY	deletion	UNP P0A0K8

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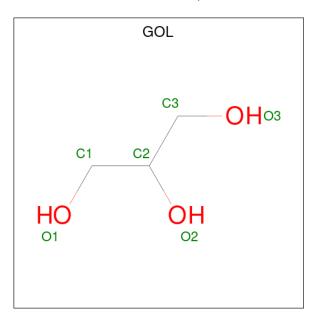
• Molecule 2 is 2-[[3,4-bis(chloranyl)-5-methyl-1 {H}-pyrrol-2-yl]carbonylamino]-4-(2-morp holin-4-ylethoxy)-1,3-benzothiazole-6-carboxylic acid (three-letter code: NWK) (formula: $C_{20}H_{20}Cl_2N_4O_5S$) (labeled as "Ligand of Interest" by depositor).





Mol	Chain	Residues	Atoms			ZeroOcc	AltConf				
2	Δ	1	Total	С	Cl	Ν	0	S	0	0	
2	Л	T	32	20	2	4	5	1	0	0	
9	В	1	Total	С	Cl	Ν	0	\mathbf{S}	0	0	
	D	1	32	20	2	4	5	1		0	

• Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).

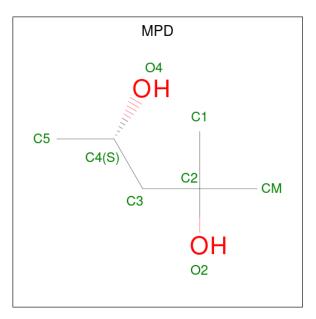


Mo	bl	Chain	Residues	Atoms		ZeroOcc	AltConf	
3		А	1	Total 6	${ m C} { m 3}$	O 3	0	0

• Molecule 4 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula:



 $C_6H_{14}O_2).$



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0
4	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0
4	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0
4	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0
4	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 8 & 6 & 2 \end{array}$	0	0

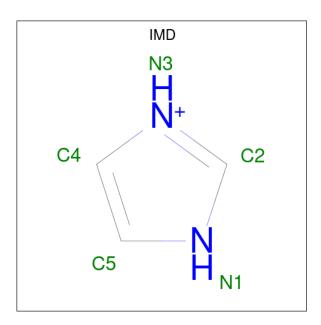
• Molecule 5 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mo	bl	Chain	Residues	Atoms		ZeroOcc	AltConf
5		А	1	Total 1	Ca 1	0	0

• Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	Total Cl 1 1	0	0
6	В	1	Total Cl 1 1	0	0





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
7	В	1	Total 5	$\begin{array}{c} \mathrm{C} \\ \mathrm{3} \end{array}$	N 2	0	0

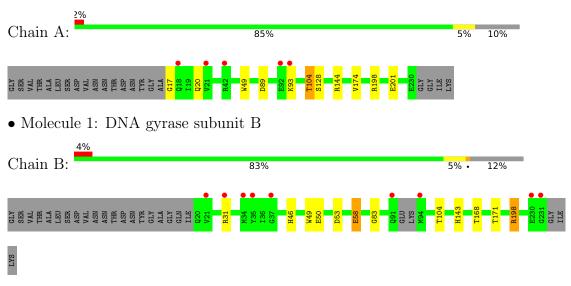
• Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	А	114	Total O 114 114	0	0
8	В	99	Total O 99 99	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: DNA gyrase subunit B



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants	142.70Å 55.36Å 50.81Å	Depositor
a, b, c, α , β , γ	90.00° 100.42° 90.00°	Depositor
Resolution (Å)	51.50 - 1.70	Depositor
	51.50 - 1.70	EDS
% Data completeness	98.0 (51.50-1.70)	Depositor
(in resolution range)	98.0(51.50-1.70)	EDS
R _{merge}	0.04	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.42 (at 1.70 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0257	Depositor
R, R_{free}	0.186 , 0.211	Depositor
It, Itfree	0.199 , 0.224	DCC
R_{free} test set	2032 reflections $(4.82%)$	wwPDB-VP
Wilson B-factor ($Å^2$)	29.9	Xtriage
Anisotropy	0.098	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	(Not available), (Not available)	EDS
L-test for twinning ²	$< L >=0.50, < L^2>=0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3431	wwPDB-VP
Average B, all atoms $(Å^2)$	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.79% 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: CL, NWK, MPD, GOL, IMD, CA

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		Bo	nd lengths	Bond angles		
	Unain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.85	1/1595~(0.1%)	0.87	2/2159~(0.1%)	
1	В	0.80	1/1572~(0.1%)	0.85	0/2127	
All	All	0.83	2/3167~(0.1%)	0.86	2/4286~(0.0%)	

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	А	104	THR	C-N	14.69	1.67	1.34
1	В	104	THR	C-N	13.83	1.65	1.34

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
1	А	144	ARG	NE-CZ-NH2	-6.64	116.98	120.30
1	А	104	THR	O-C-N	-5.19	114.40	122.70

There are no chirality outliers.

There are no planarity outliers.

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	1558	0	1541	10	0
1	В	1542	0	1506	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	А	32	0	0	1	0
2	В	32	0	0	1	0
3	А	6	0	8	0	0
4	А	24	0	42	3	0
4	В	16	0	28	0	0
5	А	1	0	0	0	0
6	А	1	0	0	0	0
6	В	1	0	0	0	0
7	В	5	0	5	0	0
8	А	114	0	0	0	0
8	В	99	0	0	3	0
All	All	3431	0	3130	22	0

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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 3.

The worst 5 of 22 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:53[A]:ASP:OD1	8:B:401:HOH:O	1.55	1.23
4:A:305:MPD:HM1	4:A:305:MPD:H52	1.70	0.72
1:A:198:ARG:HH11	1:A:201:GLU:CD	2.08	0.57
1:A:198:ARG:NH1	1:A:201:GLU:OE2	2.37	0.57
2:B:301:NWK:S	2:B:301:NWK:O	2.64	0.55

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	193/212~(91%)	190~(98%)	3~(2%)	0	100 100

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	Chain	Analysed	Favoured	Allowed	Outliers	Perce	ntiles
1	В	188/212~(89%)	180 (96%)	8 (4%)	0	100	100
All	All	381/424~(90%)	370~(97%)	11 (3%)	0	100	100

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There are no Ramachandran outliers to report.

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.

Mol	Chain	Analysed	Analysed Rotameric		Percentiles
1	А	172/183~(94%)	170~(99%)	2(1%)	71 59
1	В	169/183~(92%)	166 (98%)	3~(2%)	59 43
All	All	341/366~(93%)	336~(98%)	5(2%)	65 51

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

Mol	Chain	Res	Type
1	А	89	ASP
1	А	93	LYS
1	В	31	ARG
1	В	58	GLU
1	В	198	ARG

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

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)

Of 12 ligands modelled in this entry, 3 are monoatomic - leaving 9 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	Turne	Chain	Res	Link	Bo	ond leng	$_{\rm ths}$	B	ond ang	gles
NIOI	Type	Unam	nes		Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MPD	А	303	-	7,7,7	0.21	0	$9,\!10,\!10$	0.50	0
4	MPD	А	304	-	7,7,7	0.18	0	9,10,10	0.42	0
4	MPD	В	303	-	7,7,7	0.13	0	9,10,10	0.51	0
7	IMD	В	304	-	$3,\!5,\!5$	0.30	0	4,5,5	0.69	0
2	NWK	А	301	-	$26,\!35,\!35$	1.47	2 (7%)	$27,\!50,\!50$	2.46	14 (51%)
3	GOL	А	302	-	$5,\!5,\!5$	0.08	0	$5,\!5,\!5$	0.22	0
4	MPD	В	302	-	7,7,7	0.20	0	9,10,10	0.33	0
2	NWK	В	301	-	$26,\!35,\!35$	1.23	3 (11%)	$27,\!50,\!50$	1.54	5 (18%)
4	MPD	А	305	-	7,7,7	0.11	0	9,10,10	0.48	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
4	MPD	А	303	-	-	1/5/5/5	-
4	MPD	А	304	-	-	1/5/5/5	-
4	MPD	В	303	-	-	2/5/5/5	-
7	IMD	В	304	-	-	-	0/1/1/1
2	NWK	А	301	-	-	6/8/26/26	0/4/4/4
3	GOL	А	302	-	-	0/4/4/4	-
4	MPD	В	302	-	-	0/5/5/5	-
2	NWK	В	301	-	-	4/8/26/26	0/4/4/4
4	MPD	А	305	-	-	1/5/5/5	-



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	301	NWK	C4-C5	4.59	1.49	1.42
2	В	301	NWK	C4-C5	4.47	1.49	1.42
2	А	301	NWK	C2-C13	4.39	1.51	1.47
2	В	301	NWK	C2-C13	2.37	1.49	1.47
2	В	301	NWK	С5-С	2.03	1.49	1.42

All (5) bond length outliers are listed below:

The worst 5 of 19 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	А	301	NWK	O3-C4-C5	5.98	122.51	115.16
2	В	301	NWK	C11-C10-CL2	4.22	130.75	125.94
2	А	301	NWK	O-C7-C8	-4.18	114.33	120.59
2	А	301	NWK	C3-C2-C13	-4.05	115.03	120.36
2	А	301	NWK	C11-C10-CL2	4.05	130.55	125.94

There are no chirality outliers.

5 of 15 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	А	301	NWK	C5-C4-O3-C14
2	В	301	NWK	C14-C15-N3-C18
4	А	303	MPD	C2-C3-C4-O4
4	А	305	MPD	C2-C3-C4-O4
2	А	301	NWK	O3-C14-C15-N3

There are no ring outliers.

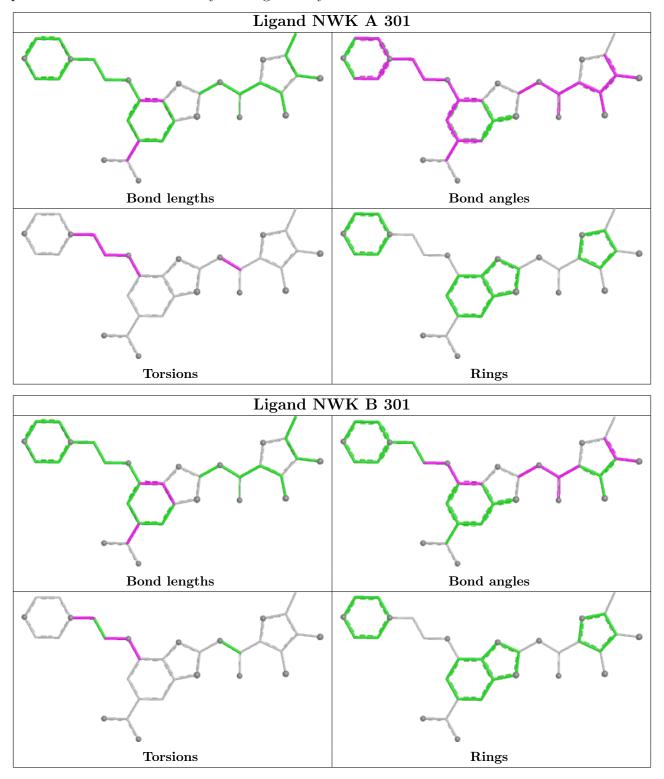
5 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	А	303	MPD	1	0
4	А	304	MPD	1	0
2	А	301	NWK	1	0
2	В	301	NWK	1	0
4	А	305	MPD	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.





5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	А	1
1	В	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	А	104:THR	С	128:SER	Ν	1.67
1	В	104:THR	С	128:SER	Ν	1.65



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	191/212~(90%)	0.08	5 (2%) 56 60	25, 34, 53, 86	0
1	В	187/212 (88%)	0.20	9 (4%) 30 34	24, 34, 63, 74	0
All	All	378/424~(89%)	0.14	14 (3%) 41 46	24, 34, 60, 86	0

The worst 5 of 14 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	34	MET	5.0
1	А	93	LYS	4.5
1	В	37	GLY	4.0
1	В	31	ARG	3.3
1	В	21	VAL	2.9

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.

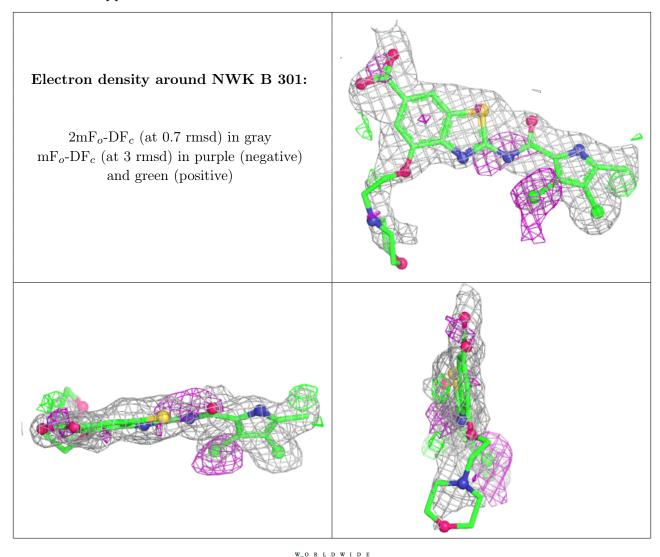
6.4 Ligands (i)

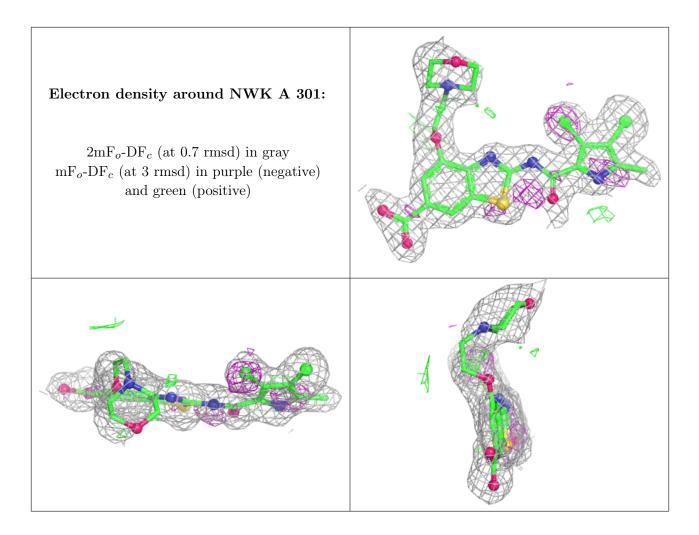
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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



Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(A^2)$	Q<0.9
4	MPD	А	305	8/8	0.75	0.23	62,69,75,80	0
4	MPD	А	304	8/8	0.83	0.18	51,57,60,60	0
4	MPD	В	302	8/8	0.83	0.18	45,49,53,54	0
2	NWK	В	301	32/32	0.85	0.21	47,56,94,96	0
7	IMD	В	304	5/5	0.85	0.13	59,59,62,62	0
4	MPD	В	303	8/8	0.89	0.18	54,59,65,65	0
3	GOL	А	302	6/6	0.89	0.12	68,71,72,74	0
4	MPD	А	303	8/8	0.91	0.20	42,53,66,67	0
6	CL	В	305	1/1	0.91	0.09	$52,\!52,\!52,\!52$	0
2	NWK	А	301	32/32	0.91	0.13	31,43,77,80	0
5	CA	А	306	1/1	0.95	0.05	100,100,100,100	1
6	CL	А	307	1/1	0.98	0.05	45,45,45,45	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.

