



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 11:51 AM UTC

PDB ID : 8RIY / pdb_00008riy
Title : Human NUDT5 with ibrutinib derivative
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Deposited on : 2023-12-19
Resolution : 2.29 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

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 : 4-5-2 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

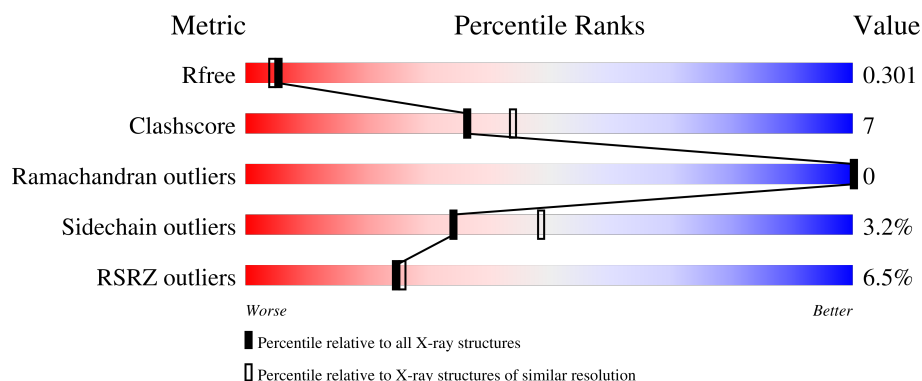
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.29 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	9078 (2.30-2.26)
Clashscore	190562	9802 (2.30-2.26)
Ramachandran outliers	187476	9690 (2.30-2.26)
Sidechain outliers	187428	9691 (2.30-2.26)
RSRZ outliers	180081	9085 (2.30-2.26)

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 $\leq 5\%$. 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	AAA	209	5% 80%13%7%
1	BBB	209	7% 76%14%8%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 3015 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.

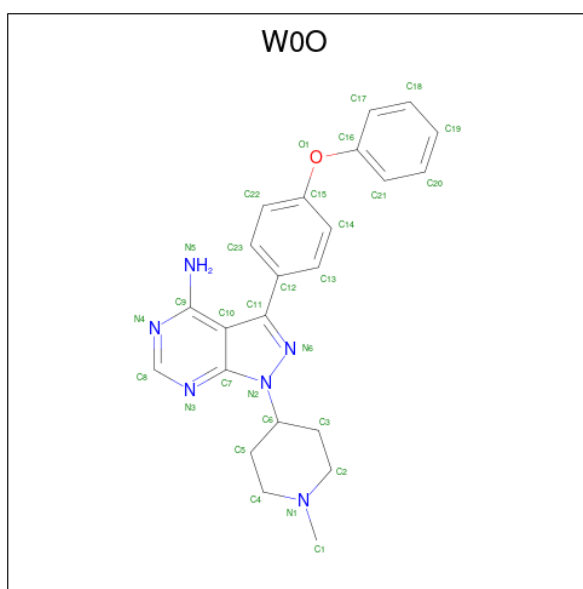
- Molecule 1 is a protein called ADP-sugar pyrophosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AAA	195	Total	C	N	O	S	1	0	0
			1466	926	242	290	8			
1	BBB	192	Total	C	N	O	S	0	0	0
			1456	918	239	291	8			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AAA	0	SER	-	expression tag	UNP Q9UKK9
BBB	0	SER	-	expression tag	UNP Q9UKK9

- Molecule 2 is 1-(1-methylpiperidin-4-yl)-3-(4-phenoxyphenyl)pyrazolo[3,4-d]pyrimidin-4-amine (CCD ID: W0O) (formula: C₂₃H₂₄N₆O) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	AAA	1	Total	C	N	O	0	0
			30	23	6	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	BBB	1	Total	C	N	O	0	0
			30	23	6	1		

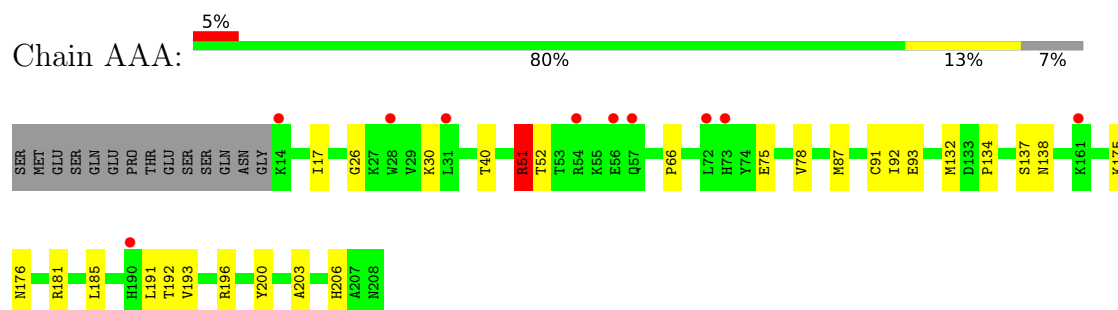
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	AAA	11	Total	O	0	0
			11	11		
3	BBB	22	Total	O	0	0
			22	22		

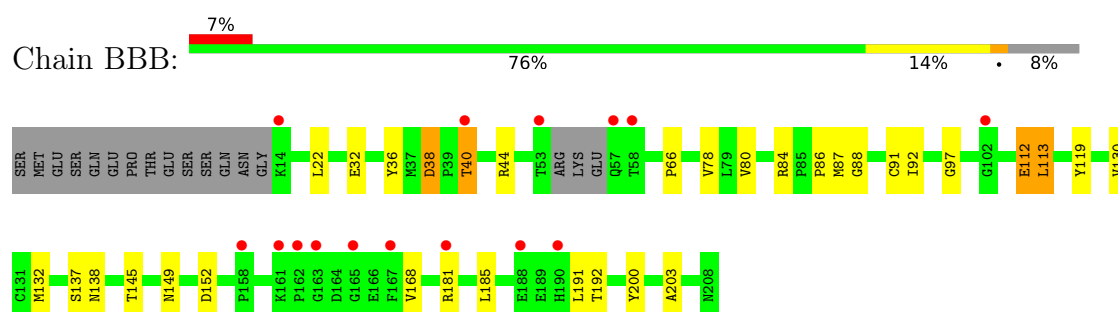
3 Residue-property plots [i](#)

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: ADP-sugar pyrophosphatase



- Molecule 1: ADP-sugar pyrophosphatase



4 Data and refinement statistics

Property	Value	Source
Space group	P 61 2 2	Depositor
Cell constants a, b, c, α , β , γ	60.60Å 60.60Å 502.61Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	83.77 – 2.29 83.77 – 2.29	Depositor EDS
% Data completeness (in resolution range)	35.5 (83.77-2.29) 35.5 (83.77-2.29)	Depositor EDS
R_{merge}	0.34	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.52 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.226 , 0.295 0.226 , 0.301	Depositor DCC
R_{free} test set	450 reflections (1.70%)	wwPDB-VP
Wilson B-factor (Å ²)	43.9	Xtriage
Anisotropy	0.179	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3015	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.70% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: W00

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	AAA	1.90	3/1494 (0.2%)	1.58	3/2037 (0.1%)
1	BBB	0.99	0/1484	1.36	0/2024
All	All	1.52	3/2978 (0.1%)	1.47	3/4061 (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	AAA	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	AAA	51	ARG	CZ-NH1	45.44	1.96	1.32
1	AAA	51	ARG	NE-CZ	-31.12	0.98	1.33
1	AAA	51	ARG	CZ-NH2	29.26	1.71	1.33

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	AAA	51	ARG	NE-CZ-NH1	-24.86	96.64	121.50
1	AAA	51	ARG	CD-NE-CZ	18.94	150.92	124.40
1	AAA	51	ARG	NH1-CZ-NH2	16.43	140.66	119.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	AAA	51	ARG	Sidechain

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	AAA	1466	0	1418	26	0
1	BBB	1456	0	1405	31	0
2	AAA	30	0	0	0	0
2	BBB	30	0	0	0	0
3	AAA	11	0	0	0	0
3	BBB	22	0	0	0	0
All	All	3015	0	2823	42	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.

All (42) 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:AAA:51:ARG:NE	1:AAA:51:ARG:NH1	2.29	0.79
1:AAA:203:ALA:HB3	1:BBB:203:ALA:HB3	1.65	0.79
1:AAA:51:ARG:NE	1:AAA:51:ARG:NH2	2.34	0.76
1:AAA:193:VAL:HG23	1:BBB:130:VAL:HG12	1.78	0.66
1:AAA:138:ASN:HB3	1:BBB:86:PRO:O	1.99	0.63
1:AAA:75:GLU:OE2	1:AAA:175:LYS:HE2	1.99	0.62
1:AAA:17:ILE:HD11	1:BBB:88:GLY:HA2	1.87	0.55
1:AAA:134:PRO:HB2	1:BBB:84:ARG:HG3	1.87	0.55
1:BBB:66:PRO:HB3	1:BBB:113:LEU:CD1	2.38	0.54
1:AAA:138:ASN:HA	1:BBB:87:MET:SD	2.50	0.52
1:AAA:92:ILE:HD11	1:AAA:191:LEU:HD13	1.92	0.52
1:AAA:203:ALA:CB	1:BBB:203:ALA:HB3	2.39	0.52
1:BBB:92:ILE:HD11	1:BBB:191:LEU:HD13	1.92	0.52
1:BBB:97:GLY:HA3	1:BBB:112:GLU:OE2	2.11	0.51
1:AAA:203:ALA:HB3	1:BBB:203:ALA:CB	2.38	0.51
1:BBB:66:PRO:HB3	1:BBB:113:LEU:HD11	1.94	0.50
1:AAA:66:PRO:HG2	1:AAA:78:VAL:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BBB:113:LEU:HD23	1:BBB:119:TYR:O	2.11	0.50
1:BBB:66:PRO:HG2	1:BBB:78:VAL:O	2.12	0.49
1:AAA:26:GLY:O	1:AAA:52:THR:HG21	2.11	0.49
1:BBB:38:ASP:OD1	1:BBB:40:THR:HG23	2.13	0.49
1:AAA:193:VAL:CG2	1:BBB:130:VAL:HG12	2.41	0.48
1:BBB:181:ARG:O	1:BBB:185:LEU:HD13	2.13	0.48
1:AAA:203:ALA:HB2	1:BBB:200:TYR:CD1	2.49	0.47
1:AAA:196:ARG:NH2	1:BBB:132:MET:O	2.47	0.46
1:AAA:200:TYR:CD1	1:BBB:203:ALA:HB2	2.52	0.45
1:BBB:149:ASN:OD1	1:BBB:152:ASP:HB2	2.17	0.45
1:AAA:51:ARG:NH2	1:AAA:51:ARG:HE	2.12	0.44
1:BBB:112:GLU:OE1	1:BBB:112:GLU:HA	2.17	0.44
1:BBB:91:CYS:HA	1:BBB:192:THR:O	2.18	0.44
1:AAA:87:MET:SD	1:BBB:138:ASN:HA	2.58	0.44
1:BBB:40:THR:O	1:BBB:40:THR:OG1	2.36	0.43
1:BBB:38:ASP:OD1	1:BBB:40:THR:CG2	2.66	0.43
1:BBB:80:VAL:HB	1:BBB:168:VAL:HG13	2.00	0.43
1:AAA:181:ARG:O	1:AAA:185:LEU:HD13	2.19	0.43
1:BBB:36:TYR:CZ	1:BBB:44:ARG:HB2	2.54	0.43
1:AAA:17:ILE:CD1	1:BBB:88:GLY:HA2	2.48	0.43
1:BBB:22:LEU:HA	1:BBB:32:GLU:HG2	2.00	0.43
1:AAA:91:CYS:HA	1:AAA:192:THR:O	2.20	0.42
1:AAA:51:ARG:NH1	1:AAA:51:ARG:CD	2.83	0.41
1:AAA:132:MET:HE3	1:AAA:196:ARG:CZ	2.51	0.41
1:AAA:206:HIS:CE1	1:BBB:145:THR:HG21	2.56	0.40

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	AAA	193/209 (92%)	182 (94%)	11 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BBB	188/209 (90%)	179 (95%)	9 (5%)	0	100	100
All	All	381/418 (91%)	361 (95%)	20 (5%)	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	AAA	153/179 (86%)	148 (97%)	5 (3%)	33	47
1	BBB	155/179 (87%)	150 (97%)	5 (3%)	34	49
All	All	308/358 (86%)	298 (97%)	10 (3%)	34	49

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	AAA	30	LYS
1	AAA	40	THR
1	AAA	93	GLU
1	AAA	137	SER
1	AAA	176	ASN
1	BBB	38	ASP
1	BBB	40	THR
1	BBB	112	GLU
1	BBB	113	LEU
1	BBB	137	SER

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 [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	W0O	BBB	301	-	34,34,34	0.52	0	48,48,48	1.08	3 (6%)
2	W0O	AAA	301	-	34,34,34	0.43	0	48,48,48	1.02	3 (6%)

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	W0O	BBB	301	-	-	0/12/22/22	0/5/5/5
2	W0O	AAA	301	-	-	4/12/22/22	0/5/5/5

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	BBB	301	W0O	C11-N6-N2	3.53	108.83	106.97
2	AAA	301	W0O	C11-N6-N2	2.80	108.44	106.97
2	BBB	301	W0O	C10-C11-C12	2.31	134.02	131.28
2	AAA	301	W0O	C16-O1-C15	2.15	123.69	118.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	AAA	301	WOO	C7-N2-N6	-2.05	109.58	110.90
2	BBB	301	WOO	C16-O1-C15	2.02	123.40	118.78

There are no chirality outliers.

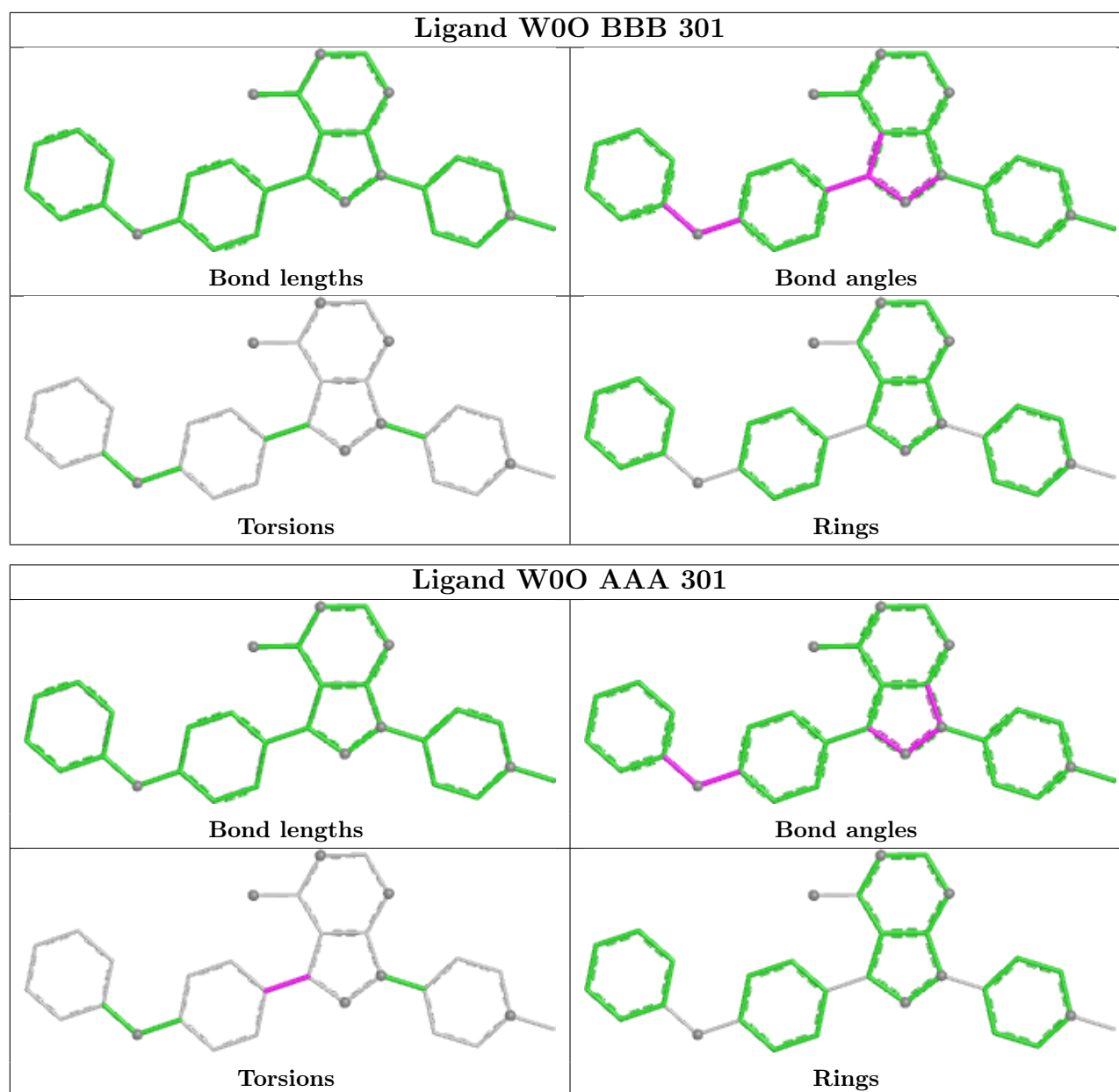
All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	AAA	301	WOO	N6-C11-C12-C13
2	AAA	301	WOO	N6-C11-C12-C23
2	AAA	301	WOO	C10-C11-C12-C13
2	AAA	301	WOO	C10-C11-C12-C23

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, 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	AAA	195/209 (93%)	0.46	10 (5%) 33 34	31, 45, 79, 93	1 (0%)
1	BBB	192/209 (91%)	0.45	15 (7%) 19 20	28, 44, 84, 113	0
All	All	387/418 (92%)	0.45	25 (6%) 25 26	28, 44, 83, 113	1 (0%)

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	AAA	57	GLN	4.2
1	BBB	162	PRO	3.8
1	BBB	57	GLN	3.4
1	BBB	53	THR	3.4
1	AAA	73	HIS	3.1
1	BBB	58	THR	3.0
1	AAA	56	GLU	2.9
1	BBB	181	ARG	2.9
1	AAA	54	ARG	2.8
1	BBB	190	HIS	2.6
1	AAA	72	LEU	2.6
1	BBB	14	LYS	2.5
1	AAA	28	TRP	2.5
1	AAA	161	LYS	2.5
1	BBB	167	PHE	2.4
1	AAA	190	HIS	2.3
1	BBB	102	GLY	2.2
1	BBB	161	LYS	2.2
1	BBB	40	THR	2.1
1	BBB	163	GLY	2.1
1	BBB	165	GLY	2.1
1	AAA	14	LYS	2.1
1	BBB	188	GLU	2.0
1	AAA	31	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	BBB	158	PRO	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](#)

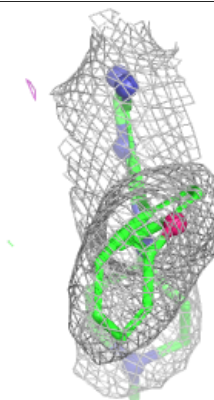
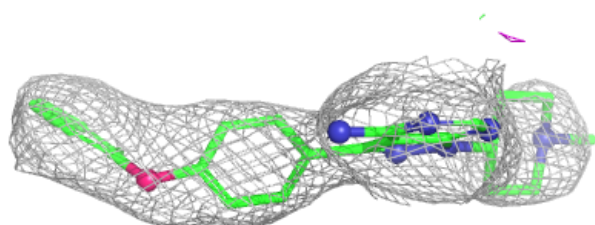
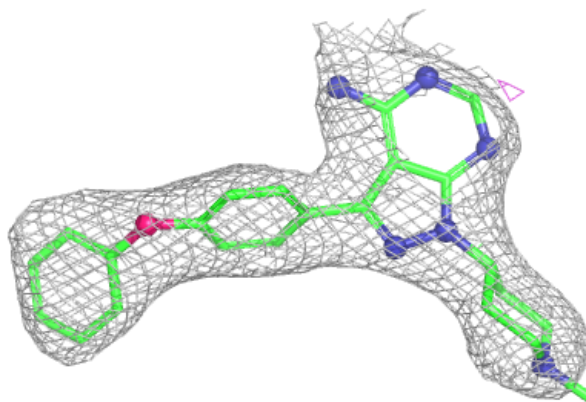
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	W0O	AAA	301	30/30	0.93	0.09	40,47,58,59	0
2	W0O	BBB	301	30/30	0.94	0.09	34,40,53,60	0

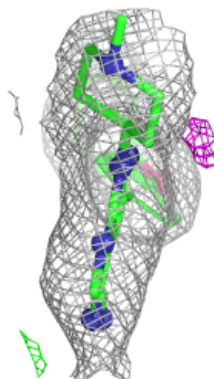
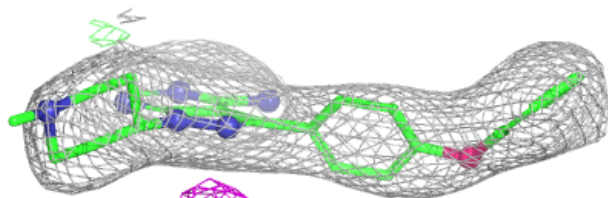
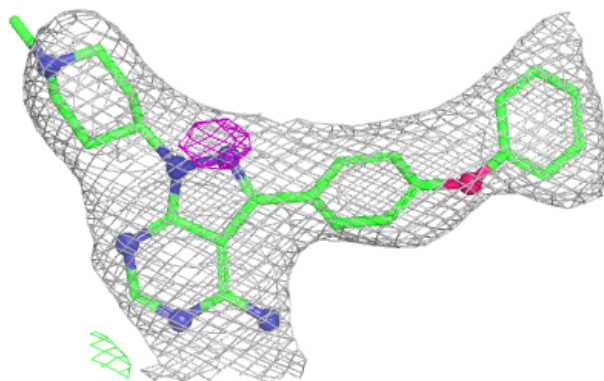
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around W0O AAA 301:

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

**Electron density around W0O BBB 301:**

$2mF_o-DF_c$ (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.