



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 12:39 AM UTC

PDB ID : 2ZHP / pdb_00002zhp
Title : Crystal structure of bleomycin-binding protein from Streptoalloteichus hindustanus complexed with bleomycin derivative
Authors : Okumura, H.; Miyazaki, I.; Simizu, S.; Osada, H.
Deposited on : 2008-02-06
Resolution : 1.60 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

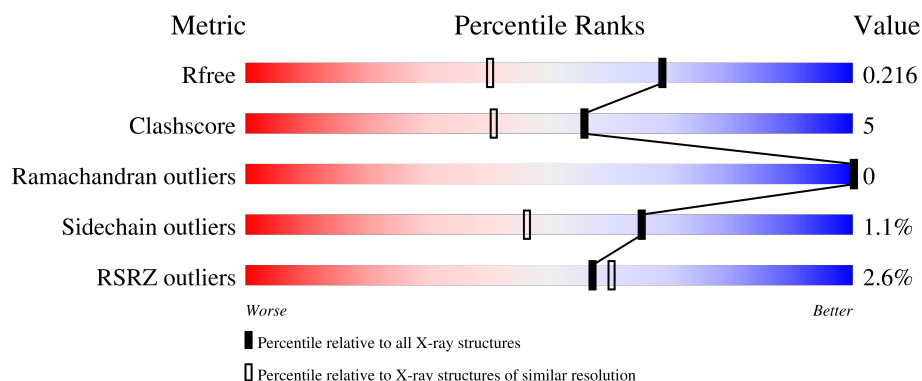
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	4673 (1.60-1.60)
Clashscore	190562	4931 (1.60-1.60)
Ramachandran outliers	187476	4831 (1.60-1.60)
Sidechain outliers	187428	4830 (1.60-1.60)
RSRZ outliers	180081	4672 (1.60-1.60)

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	A	124	 4% 83% 13% ..
1	B	124	 % 80% 10% 10%

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
4	CL	B	404	-	-	X	-

2 Entry composition [i](#)

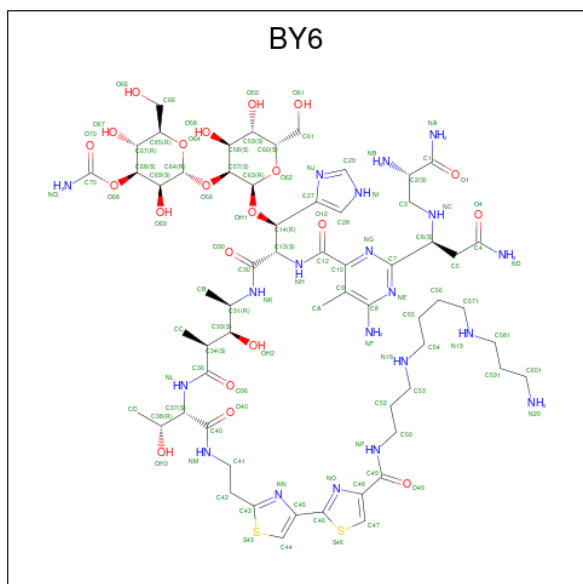
There are 5 unique types of molecules in this entry. The entry contains 2202 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bleomycin resistance protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	120	Total	C	N	O	S	0	0	0
			940	598	158	182	2			
1	B	111	Total	C	N	O	S	0	0	0
			866	551	144	169	2			

- Molecule 2 is [(2S,3R,4R,5S,6S)-2-[(3R,4R,5R,6R)-2-[(1R,2S)-2-[[6-amino-2-[(1S)-3-amino-1-[(2,3-diamino-3-oxopropyl)amino]-3-oxopropyl]-5-methylpyrimidine-4-carbonyl]amino]-3-[[[(2S,3R,4R)-5-[[[(2R,3S)-1-[2-[4-[4-[3-[4-(3-aminopropylamino)butylamino]propylcarbamoyl]-1,3-thiazol-2-yl]-1,3-thiazol-2-yl]ethylamino]-3-hydroxy-1-oxobutan-2-yl]amino]-3-hydroxy-4-methyl-5-oxopentan-2-yl]amino]-1-(3H-imidazol-4-yl)-3-oxopropoxy]-4,5-dihydroxy-6-(hydroxymethyl)oxan-3-yl]oxy-3,5-dihydroxy-6-(hydroxymethyl)oxan-4-yl] carbamate (CCD ID: BY6) (formula: $C_{60}H_{96}N_{20}O_{21}S_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	S	0	0
			91	51	17	21	2		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	C	N	O	S	0	0
			91	51	17	21	2		

- Molecule 3 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Cu	0	0
			1	1		
3	B	1	Total	Cu	0	0
			1	1		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	B	2	Total	Cl	0	0
			2	2		

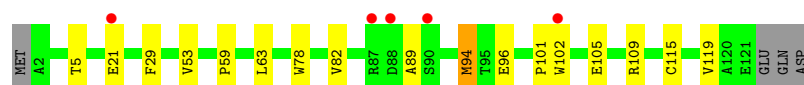
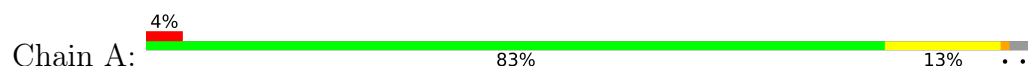
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	103	Total	O	0	0
			103	103		
5	B	106	Total	O	0	0
			106	106		

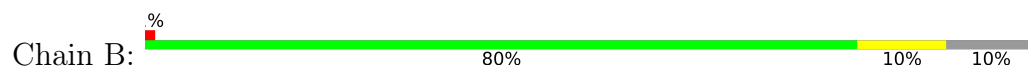
3 Residue-property plots [i](#)

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: Bleomycin resistance protein



- Molecule 1: Bleomycin resistance protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	39.99Å 53.87Å 125.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	32.11 – 1.60 32.11 – 1.60	Depositor EDS
% Data completeness (in resolution range)	99.8 (32.11-1.60) 99.9 (32.11-1.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.87 (at 1.60Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.187 , 0.216 0.187 , 0.216	Depositor DCC
R_{free} test set	1834 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	11.7	Xtriage
Anisotropy	0.293	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	2202	wwPDB-VP
Average B, all atoms (Å ²)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.48% 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: BY6, CU, CL

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	0.91	1/964 (0.1%)	1.07	2/1319 (0.2%)
1	B	0.88	0/884	1.08	1/1205 (0.1%)
All	All	0.89	1/1848 (0.1%)	1.07	3/2524 (0.1%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	MET	SD-CE	-9.94	1.54	1.79

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	69	ARG	N-CA-C	-5.28	100.64	109.24
1	A	5	THR	N-CA-C	5.17	120.24	113.88
1	A	29	PHE	N-CA-C	-5.01	103.79	110.55

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	940	0	883	9	0
1	B	866	0	820	8	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	91	0	69	0	0
2	B	91	0	69	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	1	0	0	0	0
4	B	2	0	0	5	1
5	A	103	0	0	0	0
5	B	106	0	0	0	0
All	All	2202	0	1841	18	1

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 5.

All (18) 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:25:ASP:OD2	4:B:404:CL:CL	1.98	1.18
1:B:21:GLU:OE2	4:B:404:CL:CL	2.11	1.05
1:B:21:GLU:CD	4:B:404:CL:CL	2.62	0.80
1:A:89:ALA:HA	1:A:109:ARG:NH2	2.14	0.62
1:A:101:PRO:HD2	1:A:102:TRP:CZ3	2.35	0.62
1:B:55:ASP:CG	1:B:57:VAL:HG12	2.31	0.55
1:B:21:GLU:OE1	4:B:404:CL:CL	2.62	0.55
1:A:63:LEU:HD23	1:A:115:CYS:HB3	1.89	0.54
2:B:401:BY6:H33	4:B:403:CL:CL	2.48	0.51
1:B:63:LEU:HD23	1:B:115:CYS:HB3	1.93	0.50
1:B:58:VAL:HB	1:B:59:PRO:CD	2.46	0.45
1:A:105:GLU:HB3	1:A:119:VAL:HG22	1.99	0.45
1:A:101:PRO:HD2	1:A:102:TRP:CE3	2.52	0.45
1:A:82:VAL:HB	1:A:94:MET:HE2	1.99	0.44
1:A:63:LEU:CD2	1:A:115:CYS:HB3	2.48	0.44
1:A:78:TRP:C	1:A:94:MET:HE1	2.45	0.41
1:A:53:VAL:HG21	1:A:59:PRO:HD3	2.03	0.41
1:B:74:LEU:HD11	1:B:78:TRP:CZ2	2.56	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:73:GLU:OE2	4:B:404:CL:CL[4_565]	2.08	0.12

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	118/124 (95%)	116 (98%)	2 (2%)	0	100	100
1	B	105/124 (85%)	104 (99%)	1 (1%)	0	100	100
All	All	223/248 (90%)	220 (99%)	3 (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	98/102 (96%)	96 (98%)	2 (2%)	48	24
1	B	91/102 (89%)	91 (100%)	0	100	100
All	All	189/204 (93%)	187 (99%)	2 (1%)	65	46

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	GLU
1	A	96	GLU

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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 7 ligands modelled in this entry, 5 are monoatomic - leaving 2 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	BY6	B	401	3	94,96,108	2.48	34 (36%)	119,137,149	2.72	44 (36%)
2	BY6	A	301	3	94,96,108	2.57	35 (37%)	119,137,149	2.92	50 (42%)

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	BY6	B	401	3	-	12/96/137/149	0/6/6/6
2	BY6	A	301	3	-	10/96/137/149	0/6/6/6

All (69) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	BY6	C13-C14	10.52	1.70	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	BY6	C13-C14	8.46	1.67	1.54
2	B	401	BY6	C43-NN	-6.76	1.23	1.31
2	B	401	BY6	C8-NE	6.18	1.43	1.35
2	A	301	BY6	C43-NN	-5.91	1.24	1.31
2	B	401	BY6	O62-C63	5.82	1.56	1.41
2	A	301	BY6	C8-NE	5.78	1.42	1.35
2	A	301	BY6	C49-NP	-5.69	1.25	1.33
2	A	301	BY6	C43-S43	5.18	1.80	1.72
2	A	301	BY6	O62-C63	5.18	1.55	1.41
2	A	301	BY6	C38-C37	5.17	1.68	1.53
2	B	401	BY6	C38-C37	5.11	1.68	1.53
2	A	301	BY6	C13-NH	4.94	1.56	1.45
2	B	401	BY6	C43-S43	4.89	1.80	1.72
2	B	401	BY6	OH3-C38	-4.47	1.31	1.43
2	B	401	BY6	C13-NH	4.47	1.55	1.45
2	A	301	BY6	CB-C31	4.40	1.61	1.52
2	B	401	BY6	CC-C34	4.31	1.62	1.53
2	B	401	BY6	O64-C65	4.13	1.54	1.44
2	A	301	BY6	C34-C36	-4.07	1.46	1.52
2	B	401	BY6	C12-NH	-3.98	1.25	1.34
2	A	301	BY6	C30-NK	-3.80	1.25	1.34
2	A	301	BY6	O64-C65	3.75	1.53	1.44
2	A	301	BY6	C12-NH	-3.74	1.25	1.34
2	B	401	BY6	CB-C31	3.73	1.60	1.52
2	A	301	BY6	C50-NP	3.71	1.52	1.45
2	A	301	BY6	CC-C34	3.67	1.60	1.53
2	B	401	BY6	C49-NP	-3.66	1.28	1.33
2	B	401	BY6	C50-NP	3.60	1.52	1.45
2	B	401	BY6	C7-NG	-3.50	1.28	1.34
2	B	401	BY6	C34-C36	-3.42	1.47	1.52
2	A	301	BY6	C46-NO	3.41	1.38	1.31
2	B	401	BY6	C59-C60	3.39	1.60	1.53
2	B	401	BY6	C30-NK	-3.30	1.26	1.34
2	A	301	BY6	C27-NJ	-3.24	1.33	1.39
2	A	301	BY6	O69-C69	3.17	1.50	1.43
2	B	401	BY6	O69-C69	3.05	1.50	1.43
2	A	301	BY6	OH3-C38	-3.01	1.35	1.43
2	A	301	BY6	C67-C65	2.98	1.59	1.53
2	A	301	BY6	C4-ND	2.93	1.42	1.32
2	A	301	BY6	C7-NG	-2.87	1.29	1.34
2	A	301	BY6	C28-NI	-2.87	1.29	1.36
2	B	401	BY6	C69-C68	2.65	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	BY6	O59-C59	2.65	1.49	1.43
2	A	301	BY6	O68-C68	2.65	1.48	1.44
2	B	401	BY6	C67-C65	2.63	1.58	1.53
2	B	401	BY6	C4-ND	2.62	1.41	1.32
2	A	301	BY6	C41-NM	2.62	1.52	1.46
2	A	301	BY6	C70-NQ	2.57	1.38	1.33
2	B	401	BY6	C28-NI	-2.50	1.30	1.36
2	B	401	BY6	C6-NC	-2.50	1.43	1.47
2	A	301	BY6	C59-C60	2.44	1.58	1.53
2	A	301	BY6	C59-C58	2.44	1.58	1.52
2	B	401	BY6	C27-NJ	-2.32	1.35	1.39
2	B	401	BY6	C59-C58	2.31	1.58	1.52
2	B	401	BY6	C36-NL	2.31	1.39	1.34
2	A	301	BY6	C6-NC	-2.26	1.44	1.47
2	B	401	BY6	C46-NO	2.25	1.36	1.31
2	B	401	BY6	C47-C48	2.21	1.40	1.36
2	A	301	BY6	C7-NE	-2.19	1.30	1.34
2	B	401	BY6	C46-S46	2.18	1.76	1.73
2	A	301	BY6	O59-C59	2.16	1.48	1.43
2	B	401	BY6	O68-C70	2.16	1.39	1.35
2	A	301	BY6	O40-C40	-2.14	1.19	1.23
2	B	401	BY6	O56-C57	2.13	1.49	1.43
2	A	301	BY6	C5-C6	2.12	1.57	1.53
2	A	301	BY6	CA-C9	-2.10	1.47	1.51
2	A	301	BY6	C44-C45	2.05	1.39	1.36
2	B	401	BY6	C41-NM	2.02	1.50	1.46

All (94) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	BY6	C28-C27-NJ	-12.56	101.40	109.91
2	B	401	BY6	C28-C27-NJ	-11.90	101.84	109.91
2	B	401	BY6	C63-OH1-C14	-8.46	104.44	115.23
2	A	301	BY6	C63-OH1-C14	-8.01	105.01	115.23
2	A	301	BY6	C9-C10-NG	-7.29	115.74	123.32
2	B	401	BY6	C9-C10-NG	-7.23	115.80	123.32
2	B	401	BY6	C50-NP-C49	6.73	128.76	121.90
2	A	301	BY6	O30-C30-C13	-6.73	107.08	120.75
2	A	301	BY6	NE-C7-NG	-6.06	117.65	126.15
2	B	401	BY6	O30-C30-C13	-6.03	108.50	120.75
2	B	401	BY6	NE-C7-NG	-6.02	117.71	126.15
2	A	301	BY6	C50-NP-C49	5.93	127.94	121.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	BY6	C63-C57-C58	5.65	119.60	110.37
2	A	301	BY6	C3-NC-C6	5.49	124.23	114.00
2	B	401	BY6	C3-NC-C6	5.47	124.18	114.00
2	A	301	BY6	C47-S46-C46	-5.30	85.01	89.21
2	A	301	BY6	O36-C36-C34	5.07	129.92	121.11
2	A	301	BY6	C37-C40-NM	-4.94	108.24	116.38
2	B	401	BY6	C47-S46-C46	-4.68	85.51	89.21
2	B	401	BY6	C37-C40-NM	-4.62	108.77	116.38
2	A	301	BY6	C48-C47-S46	4.53	115.47	110.24
2	B	401	BY6	C14-C13-NH	-4.45	105.41	110.20
2	A	301	BY6	C14-C13-NH	-4.31	105.56	110.20
2	A	301	BY6	C13-C30-NK	4.29	125.64	116.39
2	B	401	BY6	C63-C57-C58	4.17	117.18	110.37
2	A	301	BY6	C7-C6-NC	-4.15	105.05	112.14
2	A	301	BY6	NI-C29-NJ	-4.13	101.37	111.49
2	A	301	BY6	S46-C46-NO	4.06	119.67	114.49
2	B	401	BY6	C13-C30-NK	4.06	125.13	116.39
2	A	301	BY6	O56-C57-C63	-3.96	98.49	108.47
2	B	401	BY6	C45-C44-S43	-3.95	105.68	110.24
2	A	301	BY6	C5-C6-NC	3.86	116.53	109.25
2	B	401	BY6	O56-C57-C63	-3.85	98.76	108.47
2	B	401	BY6	C42-C41-NM	-3.77	103.97	112.00
2	B	401	BY6	NI-C29-NJ	-3.72	102.39	111.49
2	A	301	BY6	C2-C1-NA	3.71	123.09	116.75
2	A	301	BY6	C45-C44-S43	-3.70	105.96	110.24
2	B	401	BY6	O36-C36-C34	3.66	127.47	121.11
2	B	401	BY6	S46-C46-NO	3.57	119.04	114.49
2	B	401	BY6	C48-C47-S46	3.51	114.30	110.24
2	A	301	BY6	C42-C41-NM	-3.49	104.57	112.00
2	A	301	BY6	O4-C4-C5	-3.43	113.70	120.87
2	A	301	BY6	O62-C63-C57	-3.40	102.88	109.49
2	B	401	BY6	C64-O64-C65	3.35	120.25	113.72
2	A	301	BY6	C7-NE-C8	3.33	123.33	118.12
2	B	401	BY6	C2-C1-NA	3.26	122.31	116.75
2	A	301	BY6	C13-NH-C12	-3.23	115.00	121.53
2	B	401	BY6	C68-C67-C65	-3.23	102.67	109.57
2	B	401	BY6	OH3-C38-C37	-3.21	102.60	109.07
2	A	301	BY6	O58-C58-C57	3.21	118.16	109.94
2	A	301	BY6	O36-C36-NL	-3.14	117.34	122.96
2	B	401	BY6	C13-NH-C12	-3.14	115.19	121.53
2	A	301	BY6	C49-C48-NO	3.11	126.56	120.67
2	B	401	BY6	O58-C58-C57	3.09	117.86	109.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	BY6	C5-C6-NC	3.01	114.93	109.25
2	A	301	BY6	C47-C48-C49	-2.98	118.97	123.92
2	A	301	BY6	C28-NI-C29	2.93	112.59	107.36
2	A	301	BY6	C68-C67-C65	-2.87	103.46	109.57
2	A	301	BY6	C44-C45-NN	2.85	119.35	115.43
2	B	401	BY6	O36-C36-NL	-2.81	117.92	122.96
2	B	401	BY6	C7-C6-NC	-2.78	107.40	112.14
2	B	401	BY6	C28-NI-C29	2.77	112.30	107.36
2	B	401	BY6	C7-NE-C8	2.75	122.42	118.12
2	A	301	BY6	C48-NO-C46	-2.75	105.37	110.10
2	B	401	BY6	O4-C4-C5	-2.72	115.18	120.87
2	B	401	BY6	C44-C45-NN	2.64	119.07	115.43
2	B	401	BY6	O62-C63-C57	-2.54	104.55	109.49
2	A	301	BY6	OH3-C38-CD	2.47	117.08	109.68
2	A	301	BY6	C34-C36-NL	-2.46	112.71	116.35
2	A	301	BY6	O30-C30-NK	2.42	127.28	122.96
2	B	401	BY6	C45-NN-C43	2.41	111.23	110.27
2	B	401	BY6	O56-C64-C69	-2.40	102.17	108.09
2	A	301	BY6	C64-O64-C65	2.39	118.39	113.72
2	A	301	BY6	C33-C34-C36	2.36	113.08	109.53
2	A	301	BY6	OH1-C63-C57	2.36	113.67	109.09
2	B	401	BY6	OH3-C38-CD	2.34	116.70	109.68
2	B	401	BY6	C49-C48-NO	2.31	125.05	120.67
2	A	301	BY6	CD-C38-C37	-2.27	107.79	112.29
2	A	301	BY6	C2-C3-NC	2.27	117.25	112.11
2	A	301	BY6	C5-C4-ND	2.25	121.58	116.19
2	A	301	BY6	O1-C1-NA	-2.24	119.09	123.04
2	B	401	BY6	OH1-C63-C57	2.20	113.37	109.09
2	A	301	BY6	O64-C65-C67	-2.19	105.74	109.70
2	A	301	BY6	C67-C68-C69	2.15	113.85	110.86
2	A	301	BY6	C3-C2-C1	-2.13	104.24	108.70
2	B	401	BY6	C5-C4-ND	2.12	121.27	116.19
2	B	401	BY6	C33-C34-C36	2.12	112.72	109.53
2	B	401	BY6	C2-C3-NC	2.12	116.90	112.11
2	A	301	BY6	O56-C64-C69	-2.12	102.88	108.09
2	B	401	BY6	O56-C57-C58	2.11	112.59	107.23
2	B	401	BY6	O1-C1-NA	-2.08	119.36	123.04
2	A	301	BY6	C33-C31-NK	-2.06	105.58	109.56
2	B	401	BY6	C38-C37-C40	-2.06	106.83	111.32
2	A	301	BY6	C44-C45-C46	-2.01	120.51	126.44

There are no chirality outliers.

All (22) torsion outliers are listed below:

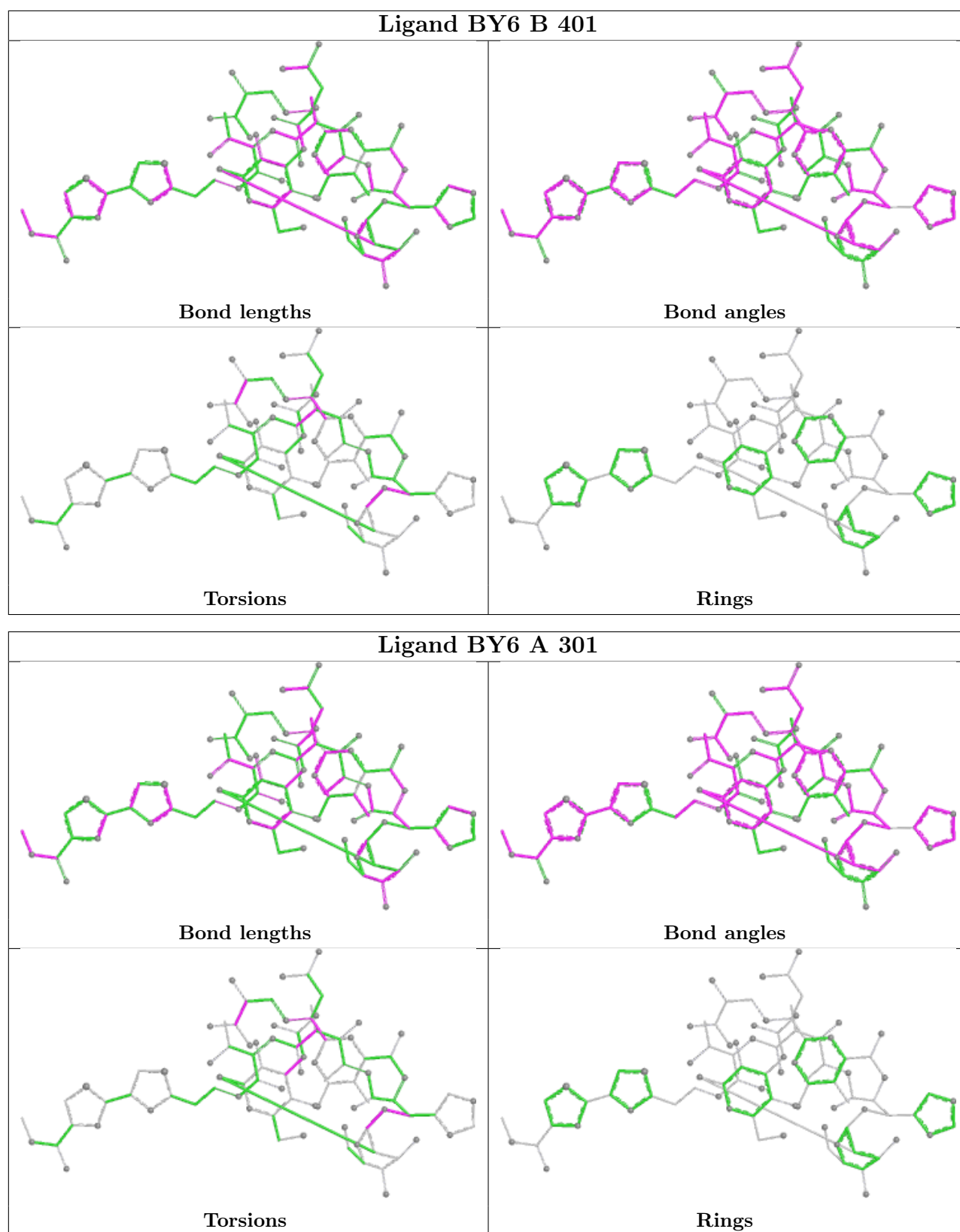
Mol	Chain	Res	Type	Atoms
2	A	301	BY6	NA-C1-C2-C3
2	A	301	BY6	O1-C1-C2-C3
2	A	301	BY6	C5-C6-NC-C3
2	A	301	BY6	C5-C6-C7-NG
2	A	301	BY6	C5-C6-C7-NE
2	A	301	BY6	CC-C34-C36-O36
2	A	301	BY6	CC-C34-C36-NL
2	B	401	BY6	C5-C6-NC-C3
2	B	401	BY6	C5-C6-C7-NG
2	B	401	BY6	C5-C6-C7-NE
2	B	401	BY6	C13-C14-OH1-C63
2	B	401	BY6	C27-C14-OH1-C63
2	B	401	BY6	CC-C34-C36-O36
2	B	401	BY6	CC-C34-C36-NL
2	B	401	BY6	O1-C1-C2-C3
2	A	301	BY6	C69-C68-O68-C70
2	A	301	BY6	C13-C14-OH1-C63
2	B	401	BY6	NA-C1-C2-C3
2	B	401	BY6	C57-C63-OH1-C14
2	A	301	BY6	C57-C63-OH1-C14
2	B	401	BY6	O1-C1-C2-NB
2	B	401	BY6	O62-C63-OH1-C14

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	BY6	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 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 ⓘ

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, 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	120/124 (96%)	-0.12	5 (4%) 40 42	6, 11, 29, 36	0
1	B	111/124 (89%)	-0.22	1 (0%) 81 84	6, 10, 24, 35	0
All	All	231/248 (93%)	-0.17	6 (2%) 57 60	6, 11, 26, 36	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	103	GLY	2.9
1	A	102	TRP	2.4
1	A	90	SER	2.4
1	A	87	ARG	2.3
1	A	21	GLU	2.2
1	A	88	ASP	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](#)

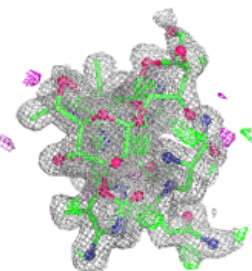
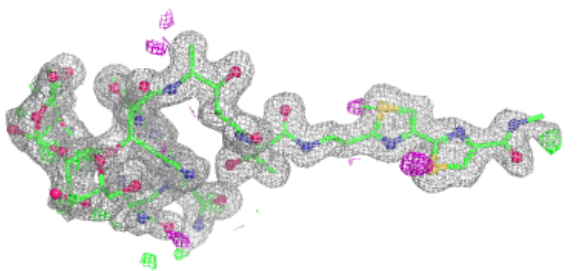
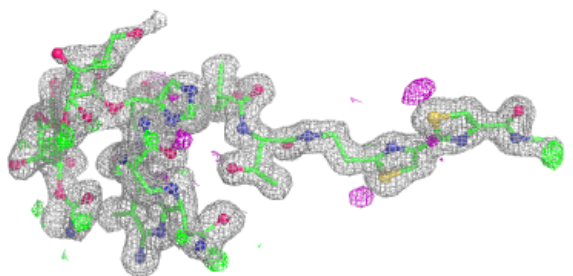
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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	BY6	B	401	91/103	0.93	0.11	8,21,42,48	0
4	CL	A	303	1/1	0.95	0.16	19,19,19,19	0
4	CL	B	404	1/1	0.95	0.05	10,10,10,10	0
2	BY6	A	301	91/103	0.96	0.06	5,11,22,29	0
4	CL	B	403	1/1	0.97	0.17	25,25,25,25	0
3	CU	B	402	1/1	0.99	0.02	17,17,17,17	0
3	CU	A	302	1/1	0.99	0.03	12,12,12,12	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.

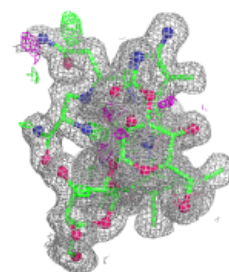
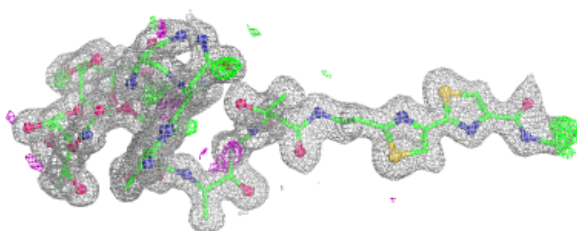
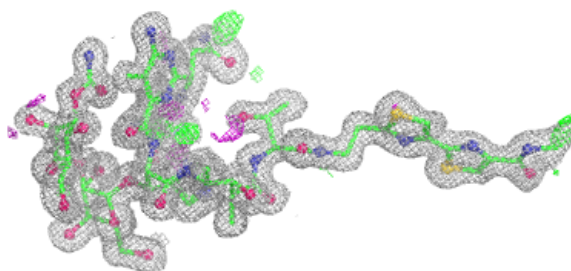
Electron density around BY6 B 401:

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 BY6 A 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.