



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:05 AM UTC

PDB ID : 3IXG / pdb_00003ixg
Title : X-ray crystal structure of the extended-spectrum AmpC T70I mutant beta-lactamase with and without benzo(b)thiophene-2-boronic acid bound at 2.14 Angstrom resolution
Authors : Shoichet, B.K.; Thomas, V.L.
Deposited on : 2009-09-04
Resolution : 2.14 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
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

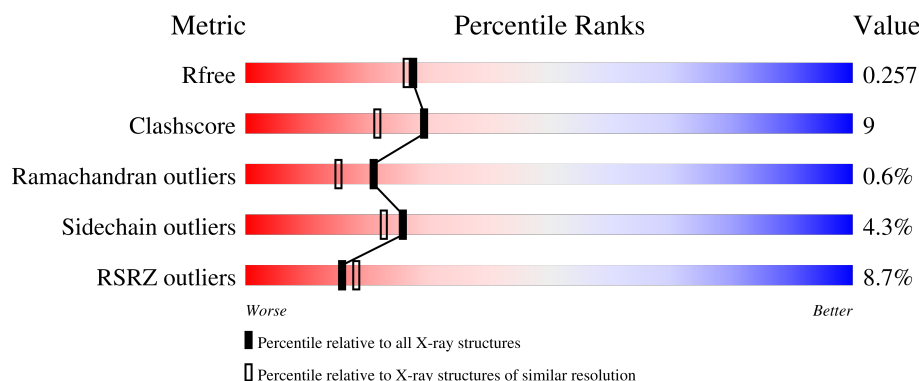
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.14 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	3689 (2.16-2.12)
Clashscore	190562	3812 (2.16-2.12)
Ramachandran outliers	187476	3773 (2.16-2.12)
Sidechain outliers	187428	3772 (2.16-2.12)
RSRZ outliers	180081	3691 (2.16-2.12)

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	358	9% 71% 16% • 11%
1	B	358	7% 80% 18% •

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5524 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.

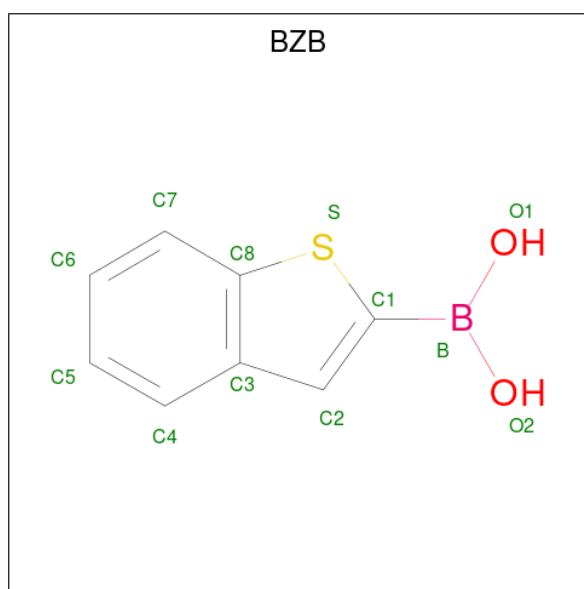
- Molecule 1 is a protein called Beta-lactamase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	317	Total	C	N	O	S	0	0	0
			2491	1616	422	447	6			
1	B	358	Total	C	N	O	S	0	0	0
			2801	1806	476	513	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	70	ILE	THR	engineered mutation	UNP P00811
B	70	ILE	THR	engineered mutation	UNP P00811

- Molecule 2 is BENZO[B]THIOPHENE-2-BORONIC ACID (CCD ID: BZB) (formula: $C_8H_7BO_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	B	1	Total	B	C	O	S	0	0
			12	1	8	2	1		

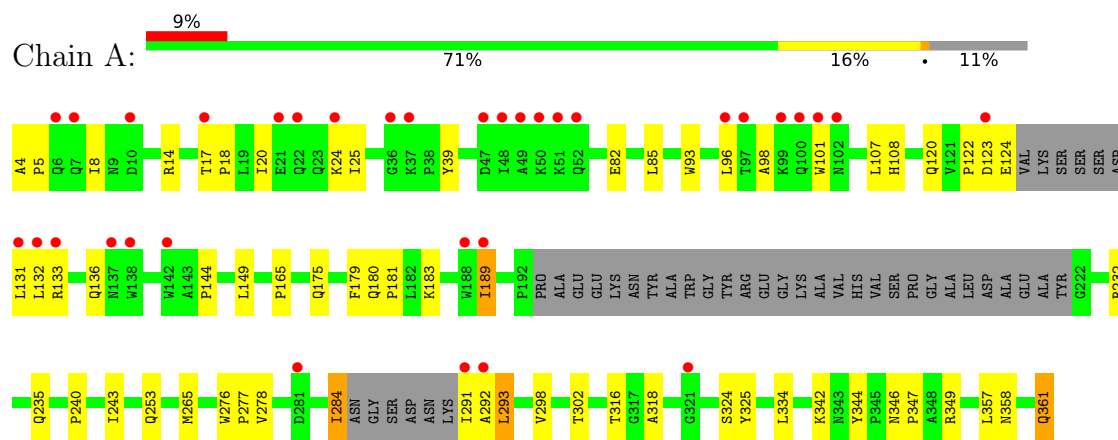
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	110	Total 110	O 110	0	0
3	B	110	Total 110	O 110	0	0

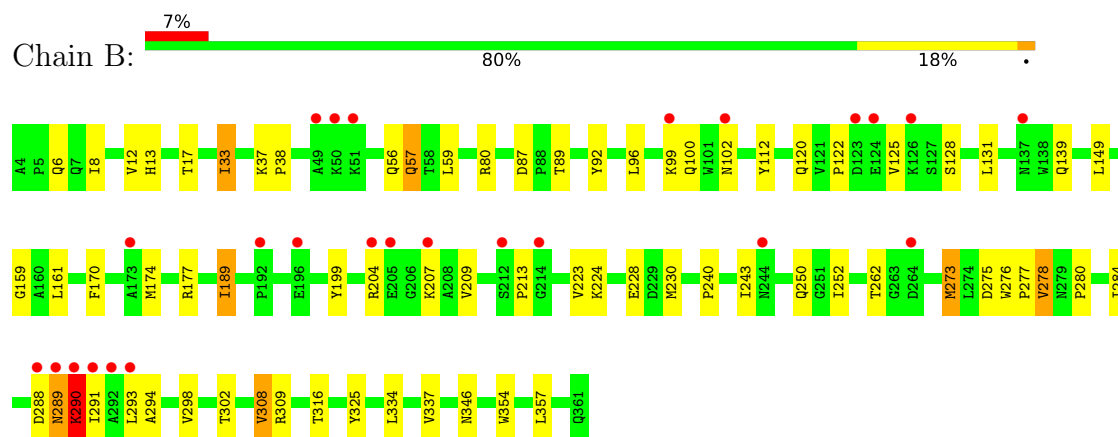
3 Residue-property plots

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: Beta-lactamase



• Molecule 1: Beta-lactamase



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	114.16Å 77.36Å 91.75Å 90.00° 122.12° 90.00°	Depositor
Resolution (Å)	30.00 – 2.14 30.00 – 2.14	Depositor EDS
% Data completeness (in resolution range)	98.4 (30.00-2.14) 98.3 (30.00-2.14)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.22 (at 2.14Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.197 , 0.245 (Not available) , 0.257	Depositor DCC
R_{free} test set	1844 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	26.1	Xtriage
Anisotropy	0.494	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 58.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5524	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.60% 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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.64	0/2560	0.85	1/3494 (0.0%)
1	B	0.74	0/2881	0.89	0/3934
All	All	0.70	0/5441	0.87	1/7428 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	318	ALA	N-CA-C	5.79	118.34	111.33

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	A	2491	0	2497	40	0
1	B	2801	0	2782	50	0
2	B	12	0	7	1	0
3	A	110	0	0	7	0
3	B	110	0	0	4	0
All	All	5524	0	5286	89	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 9.

All (89) 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:308:VAL:HG22	1:B:309:ARG:HG2	1.61	0.81
1:A:149:LEU:HD11	1:A:293:LEU:HG	1.66	0.77
1:B:262:THR:CG2	1:B:298:VAL:HG12	2.15	0.77
1:B:120:GLN:HG2	2:B:362:BZB:HC51	1.69	0.74
1:B:262:THR:HG22	1:B:298:VAL:HG12	1.68	0.73
1:B:290:LYS:HD3	1:B:294:ALA:HB2	1.69	0.72
3:A:471:HOH:O	1:B:250:GLN:HG2	1.90	0.72
1:B:100:GLN:NE2	3:B:405:HOH:O	2.25	0.69
1:A:132:LEU:HG	3:A:389:HOH:O	1.92	0.69
1:A:334:LEU:HG	1:A:357:LEU:HD22	1.75	0.69
1:B:288:ASP:O	1:B:289:ASN:HB2	1.95	0.66
1:A:5:PRO:HG2	1:A:8:ILE:HD12	1.80	0.63
1:B:122:PRO:HB2	1:B:125:VAL:HG23	1.79	0.63
1:B:290:LYS:HD3	1:B:294:ALA:CB	2.28	0.63
1:A:14:ARG:NH2	3:A:398:HOH:O	2.31	0.62
1:B:56:GLN:NE2	1:B:228:GLU:OE2	2.32	0.62
1:B:57:GLN:HE21	1:B:57:GLN:HA	1.66	0.61
1:A:316:THR:HG22	1:A:325:TYR:CE1	2.38	0.59
1:B:189:ILE:HD13	1:B:223:VAL:HG22	1.84	0.58
1:B:276:TRP:CD2	1:B:277:PRO:HA	2.37	0.58
1:B:284:ILE:HG23	1:B:288:ASP:OD2	2.04	0.57
1:B:289:ASN:C	1:B:291:ILE:H	2.12	0.57
1:B:334:LEU:HG	1:B:357:LEU:HD22	1.86	0.56
1:A:240:PRO:HG2	1:A:253:GLN:HE21	1.72	0.54
1:A:17:THR:HB	1:A:18:PRO:HD3	1.89	0.54
1:A:316:THR:HG21	1:A:346:ASN:HD21	1.73	0.54
1:B:59:LEU:HB2	1:B:199:TYR:HA	1.89	0.54
1:A:232:ARG:HA	1:A:235:GLN:HG2	1.90	0.53
1:B:243:ILE:HD13	1:B:252:ILE:HD12	1.91	0.52
1:B:280:PRO:HD3	1:B:354:TRP:CE2	2.44	0.52
1:A:20:ILE:HA	1:A:25:ILE:HD12	1.93	0.51
1:B:128:SER:HA	1:B:131:LEU:HD12	1.92	0.50
1:A:232:ARG:O	1:A:235:GLN:HG2	2.12	0.50
1:B:102:ASN:HA	3:B:484:HOH:O	2.12	0.50
1:A:136:GLN:HG3	3:A:389:HOH:O	2.12	0.50
1:A:131:LEU:N	3:A:475:HOH:O	2.45	0.49
1:B:33:ILE:HD12	1:B:33:ILE:N	2.27	0.49
1:B:89:THR:HG22	1:B:96:LEU:HD23	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:89:THR:CG2	1:B:96:LEU:HD23	2.44	0.48
1:A:180:GLN:HB2	1:A:181:PRO:HD3	1.95	0.47
1:A:82:GLU:HB3	1:A:165:PRO:HB2	1.97	0.47
1:B:89:THR:HG23	1:B:161:LEU:CD1	2.45	0.47
1:A:107:LEU:HD22	1:B:302:THR:HG21	1.97	0.47
1:A:344:TYR:CZ	1:A:349:ARG:HG2	2.50	0.46
1:A:324:SER:HB3	3:A:468:HOH:O	2.14	0.46
1:B:289:ASN:O	1:B:291:ILE:N	2.49	0.46
1:A:302:THR:HG22	3:B:424:HOH:O	2.16	0.46
1:B:8:ILE:O	1:B:12:VAL:HG23	2.16	0.46
1:B:316:THR:HG22	1:B:325:TYR:CD1	2.51	0.46
1:A:276:TRP:CD2	1:A:277:PRO:HA	2.51	0.45
1:A:292:ALA:O	1:A:293:LEU:C	2.59	0.45
1:A:175:GLN:HG3	1:A:189:ILE:HD11	1.98	0.45
1:B:284:ILE:O	1:B:288:ASP:HB2	2.17	0.45
1:B:288:ASP:O	1:B:289:ASN:CB	2.64	0.45
1:B:288:ASP:OD1	1:B:346:ASN:HB3	2.16	0.45
1:A:179:PHE:HE2	1:A:189:ILE:HG13	1.81	0.45
1:B:243:ILE:CD1	1:B:252:ILE:HD12	2.47	0.45
1:A:316:THR:HG22	1:A:325:TYR:HE1	1.79	0.45
1:B:275:ASP:O	1:B:278:VAL:HG22	2.17	0.44
1:B:87:ASP:OD2	1:B:92:TYR:OH	2.33	0.44
1:A:316:THR:HG22	1:A:325:TYR:CD1	2.53	0.44
1:B:80:ARG:HH21	1:B:177:ARG:HG2	1.83	0.44
1:A:5:PRO:HD2	1:A:39:TYR:CE2	2.53	0.44
1:A:284:ILE:HG21	1:A:347:PRO:HA	2.00	0.44
1:B:159:GLY:HA3	1:B:170:PHE:CZ	2.53	0.43
1:A:24:LYS:HB2	1:A:342:LYS:NZ	2.32	0.43
1:B:99:LYS:O	1:B:102:ASN:HB2	2.18	0.43
1:A:122:PRO:HB2	1:A:124:GLU:HG2	2.01	0.43
1:B:13:HIS:O	1:B:17:THR:HB	2.19	0.43
1:B:240:PRO:HA	1:B:243:ILE:HD12	2.01	0.43
1:B:289:ASN:C	1:B:291:ILE:N	2.76	0.42
1:A:175:GLN:NE2	3:A:402:HOH:O	2.53	0.42
1:B:37:LYS:HG3	1:B:38:PRO:HD2	2.01	0.42
1:A:183:LYS:HD2	1:A:232:ARG:NH2	2.35	0.42
1:B:243:ILE:HD13	1:B:252:ILE:CD1	2.49	0.42
1:B:273:MET:HE2	1:B:273:MET:HB2	1.82	0.42
1:A:358:ASN:HA	1:A:361:GLN:HE21	1.85	0.41
1:B:262:THR:HG22	1:B:298:VAL:CG1	2.44	0.41
1:A:240:PRO:HA	1:A:243:ILE:HD13	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:TYR:HB3	1:B:149:LEU:O	2.20	0.41
1:A:98:ALA:HB3	1:A:101:TRP:HD1	1.84	0.41
1:A:85:LEU:HD23	1:A:85:LEU:HA	1.88	0.41
1:A:93:TRP:CE3	1:A:96:LEU:HD22	2.56	0.41
1:B:230:MET:HB3	1:B:337:VAL:HG11	2.02	0.41
1:B:288:ASP:OD1	1:B:346:ASN:CB	2.69	0.41
1:A:108:HIS:CE1	1:A:144:PRO:HB2	2.56	0.41
1:B:224:LYS:HE2	3:B:396:HOH:O	2.19	0.41
1:A:284:ILE:H	1:A:284:ILE:HG12	1.70	0.40
1:A:4:ALA:HA	1:A:5:PRO:HD2	1.94	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	A	309/358 (86%)	300 (97%)	8 (3%)	1 (0%)	36	33
1	B	356/358 (99%)	344 (97%)	9 (2%)	3 (1%)	16	9
All	All	665/716 (93%)	644 (97%)	17 (3%)	4 (1%)	21	15

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	289	ASN
1	A	293	LEU
1	B	213	PRO
1	B	290	LYS

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	261/292 (89%)	251 (96%)	10 (4%)	29	28
1	B	292/292 (100%)	278 (95%)	14 (5%)	23	18
All	All	553/584 (95%)	529 (96%)	24 (4%)	26	23

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	120	GLN
1	A	123	ASP
1	A	133	ARG
1	A	189	ILE
1	A	265	MET
1	A	278	VAL
1	A	284	ILE
1	A	291	ILE
1	A	298	VAL
1	A	361	GLN
1	B	6	GLN
1	B	33	ILE
1	B	57	GLN
1	B	139	GLN
1	B	174	MET
1	B	189	ILE
1	B	204	ARG
1	B	207	LYS
1	B	209	VAL
1	B	273	MET
1	B	278	VAL
1	B	290	LYS
1	B	293	LEU
1	B	308	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	35	GLN
1	A	137	ASN
1	A	180	GLN
1	A	185	ASN
1	A	190	ASN
1	A	235	GLN
1	A	250	GLN
1	A	253	GLN
1	A	341	ASN
1	A	358	ASN
1	A	361	GLN
1	B	7	GLN
1	B	9	ASN
1	B	35	GLN
1	B	56	GLN
1	B	57	GLN
1	B	102	ASN
1	B	120	GLN
1	B	180	GLN
1	B	285	ASN
1	B	346	ASN

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

1 ligand is 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	BZB	B	362	-	12,13,13	5.15	4 (33%)	12,18,18	1.31	1 (8%)

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	BZB	B	362	-	-	1/2/4/4	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	362	BZB	B-O2	12.27	1.54	1.36
2	B	362	BZB	B-O1	11.67	1.53	1.36
2	B	362	BZB	C3-C8	3.89	1.48	1.41
2	B	362	BZB	C8-S	-3.29	1.67	1.74

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	362	BZB	C4-C3-C8	-2.46	116.00	119.05

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	362	BZB	O2-B-C1-C2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	362	BZB	1	0

5.7 Other polymers

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

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	A	317/358 (88%)	0.57	34 (10%)	11 13	13, 25, 50, 61	12 (3%)
1	B	358/358 (100%)	0.40	25 (6%)	22 26	12, 25, 43, 49	14 (3%)
All	All	675/716 (94%)	0.48	59 (8%)	16 18	12, 25, 46, 61	26 (3%)

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	50	LYS	11.5
1	B	126	LYS	7.5
1	B	288	ASP	7.4
1	A	51	LYS	7.0
1	B	50	LYS	6.5
1	A	22	GLN	6.3
1	B	205	GLU	6.1
1	B	123	ASP	6.1
1	B	207	LYS	5.9
1	B	289	ASN	5.8
1	A	48	ILE	5.6
1	B	124	GLU	5.5
1	B	51	LYS	5.2
1	B	290	LYS	5.0
1	B	99	LYS	5.0
1	A	49	ALA	4.8
1	A	47	ASP	4.8
1	A	188	TRP	4.6
1	B	102	ASN	4.6
1	A	131	LEU	4.5
1	A	291	ILE	4.4
1	B	196	GLU	4.4
1	A	21	GLU	4.2
1	B	49	ALA	4.2

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Mol	Chain	Res	Type	RSRZ
1	B	214	GLY	4.0
1	A	97	THR	3.9
1	A	281	ASP	3.5
1	A	132	LEU	3.4
1	A	96	LEU	3.3
1	A	99	LYS	3.2
1	A	123	ASP	3.1
1	A	100	GLN	3.1
1	B	264	ASP	3.1
1	B	244	ASN	3.0
1	A	7	GLN	3.0
1	A	321	GLY	3.0
1	B	137	ASN	2.9
1	B	192	PRO	2.8
1	B	293	LEU	2.8
1	A	37	LYS	2.8
1	B	291	ILE	2.6
1	A	6	GLN	2.5
1	A	17	THR	2.5
1	A	142	TRP	2.5
1	A	24	LYS	2.5
1	A	292	ALA	2.4
1	A	102	ASN	2.4
1	A	52	GLN	2.4
1	B	204	ARG	2.4
1	A	101	TRP	2.4
1	A	137	ASN	2.3
1	B	292	ALA	2.3
1	A	133	ARG	2.2
1	A	36	GLY	2.2
1	B	212	SER	2.1
1	A	10	ASP	2.1
1	A	189	ILE	2.1
1	A	138	TRP	2.1
1	B	173	ALA	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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	BZB	B	362	12/12	0.82	0.13	24,34,35,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.