



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:03 PM UTC

PDB ID : 9CX4 / pdb_00009cx4
Title : Acinetobacter baumannii BamA POTRAs 1-4, space group P1
Authors : Overly Cottom, C.; Noinaj, N.
Deposited on : 2024-07-30
Resolution : 2.78 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

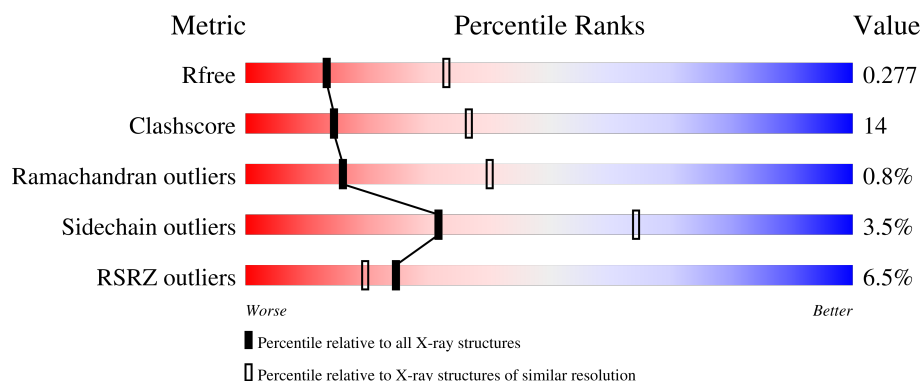
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.78 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	5248 (2.80-2.76)
Clashscore	190562	5693 (2.80-2.76)
Ramachandran outliers	187476	5590 (2.80-2.76)
Sidechain outliers	187428	5592 (2.80-2.76)
RSRZ outliers	180081	5251 (2.80-2.76)

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	328	6% 66% 33% .
1	B	328	7% 74% 24% .

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 4978 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 Outer membrane protein assembly factor BamA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	Se	0	1	0
			2486	1581	433	467	5			
1	B	328	Total	C	N	O	Se	0	0	0
			2459	1567	425	462	5			

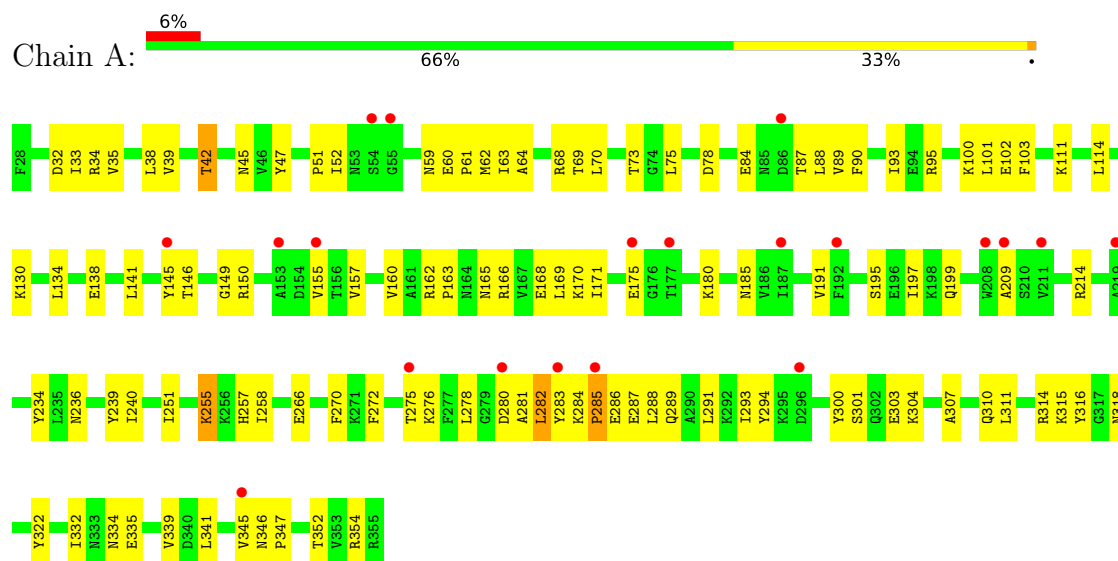
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	19	Total	O	0	0
			19	19		
2	B	14	Total	O	0	0
			14	14		

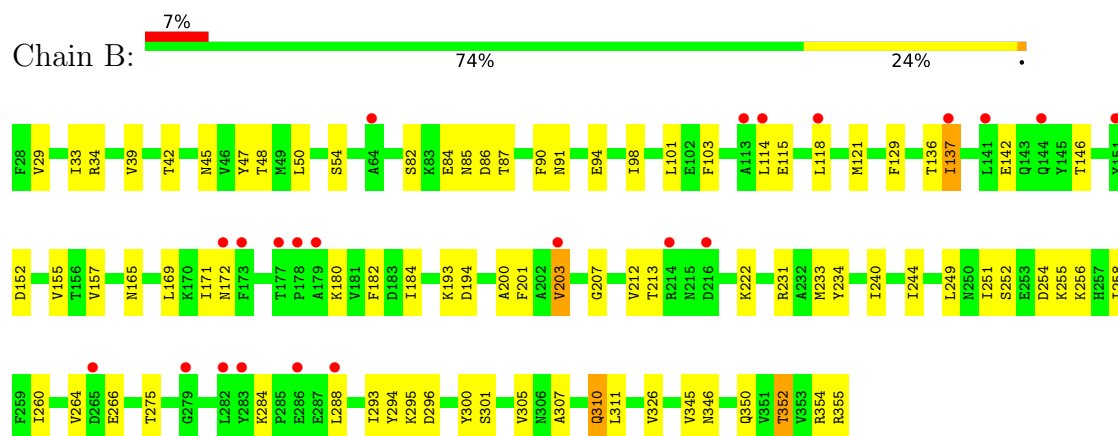
3 Residue-property plots [i](#)

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: Outer membrane protein assembly factor BamA



- Molecule 1: Outer membrane protein assembly factor BamA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	56.52Å 61.51Å 85.90Å 97.59° 107.34° 112.45°	Depositor
Resolution (Å)	39.37 – 2.78 39.37 – 2.78	Depositor EDS
% Data completeness (in resolution range)	94.0 (39.37-2.78) 94.0 (39.37-2.78)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 2.77Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.235 , 0.277 0.236 , 0.277	Depositor DCC
R_{free} test set	1988 reflections (8.09%)	wwPDB-VP
Wilson B-factor (Å ²)	63.0	Xtriage
Anisotropy	0.268	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 72.2	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	0.93	EDS
Total number of atoms	4978	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.63% 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](#)

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	0.18	0/2521	0.42	1/3421 (0.0%)
1	B	0.15	0/2495	0.36	0/3388
All	All	0.17	0/5016	0.39	1/6809 (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	146	THR	N-CA-C	-7.61	105.95	114.62

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	2486	0	2395	85	0
1	B	2459	0	2352	56	0
2	A	19	0	0	6	0
2	B	14	0	0	3	0
All	All	4978	0	4747	136	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 14.

All (136) 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:A:35:VAL:HG22	1:A:90:PHE:HB2	1.52	0.90
1:A:73:THR:HB	1:A:75:LEU:HD12	1.53	0.90
1:B:346:ASN:ND2	2:B:401:HOH:O	2.12	0.83
1:A:70:LEU:O	1:A:73:THR:OG1	1.96	0.82
1:A:103:PHE:HE1	1:A:171:ILE:HD12	1.49	0.78
1:A:68:ARG:O	2:A:401:HOH:O	2.00	0.78
1:A:78:ASP:HB3	1:A:93:ILE:HD12	1.68	0.76
1:B:157:VAL:HG22	1:B:171:ILE:HG22	1.68	0.75
1:A:42:THR:HG22	1:A:45:ASN:H	1.51	0.74
1:A:59:ASN:HD22	1:A:61:PRO:HD2	1.55	0.71
1:A:352:THR:OG1	2:A:403:HOH:O	2.07	0.71
1:A:63:ILE:HD12	1:A:64:ALA:N	2.05	0.71
1:B:180:LYS:HD3	1:B:256:LYS:HA	1.75	0.69
1:A:180:LYS:HE3	1:A:209:ALA:HA	1.77	0.66
1:B:266:GLU:O	2:B:402:HOH:O	2.13	0.66
1:B:251:ILE:HD11	1:B:255:LYS:HA	1.77	0.65
1:B:115:GLU:HA	1:B:118:LEU:HD12	1.78	0.63
1:A:102:GLU:HB2	1:A:170:LYS:HA	1.81	0.62
1:B:355:ARG:NH2	2:B:404:HOH:O	2.32	0.61
1:A:294:TYR:HB3	1:A:304:LYS:HD2	1.82	0.61
1:A:160:VAL:HG22	1:A:168:GLU:HB3	1.81	0.61
1:A:163:PRO:HB2	1:B:91:ASN:HB3	1.82	0.61
1:B:180:LYS:HA	1:B:180:LYS:HE2	1.83	0.61
1:A:294:TYR:OH	2:A:404:HOH:O	2.10	0.60
1:A:157:VAL:HG22	1:A:171:ILE:HG23	1.84	0.60
1:B:182:PHE:HE1	1:B:258:ILE:HD11	1.67	0.60
1:A:287:GLU:HA	1:A:291:LEU:HD23	1.85	0.58
1:A:35:VAL:HG12	1:A:38:LEU:HD21	1.86	0.58
1:A:150:ARG:HH12	1:A:175:GLU:HB3	1.69	0.58
1:A:52:ILE:HG12	1:A:62:MSE:SE	2.53	0.58
1:A:301:SER:HB3	1:A:304:LYS:HG3	1.87	0.57
1:A:280:ASP:HB2	1:A:345:VAL:O	2.05	0.56
1:A:316:TYR:CE2	1:A:347:PRO:HB3	2.40	0.56
1:B:182:PHE:CE1	1:B:258:ILE:HD11	2.40	0.56
1:B:103:PHE:CE2	1:B:171:ILE:HD11	2.40	0.56
1:A:322:TYR:CE1	1:B:231:ARG:HD2	2.41	0.56
1:B:121:MSE:HE1	1:B:137:ILE:HA	1.87	0.55
1:B:155:VAL:HA	1:B:172:ASN:O	2.06	0.55
1:B:33:ILE:O	1:B:47:TYR:OH	2.12	0.55
1:B:252:SER:OG	1:B:254:ASP:OD2	2.24	0.55
1:A:60:GLU:HA	1:A:63:ILE:HG13	1.90	0.54
1:A:314[B]:ARG:O	1:A:318:ASN:ND2	2.41	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:157:VAL:HG13	1:A:171:ILE:HG12	1.89	0.54
1:A:280:ASP:C	1:A:282:LEU:HD12	2.34	0.53
1:A:39:VAL:HG22	1:B:39:VAL:HG22	1.91	0.53
1:A:240:ILE:HD13	1:A:300:TYR:CE2	2.43	0.53
1:A:314[A]:ARG:O	1:A:318:ASN:ND2	2.42	0.53
1:A:103:PHE:CE1	1:A:171:ILE:HD12	2.37	0.53
1:B:249:LEU:HD12	1:B:260:ILE:HD11	1.91	0.52
1:A:134:LEU:HD11	1:A:157:VAL:HB	1.91	0.52
1:A:332:ILE:HG12	1:A:339:VAL:HG23	1.91	0.52
1:B:244:ILE:HD13	1:B:264:VAL:HG12	1.91	0.52
1:B:207:GLY:HA3	1:B:222:LYS:HE3	1.92	0.52
1:A:51:PRO:HD3	1:A:69:THR:HG21	1.92	0.51
1:A:281:ALA:O	1:A:283:TYR:N	2.44	0.51
1:B:42:THR:HB	1:B:45:ASN:ND2	2.26	0.51
1:A:276:LYS:HE3	1:A:278:LEU:HG	1.92	0.51
1:A:354:ARG:NH1	2:A:402:HOH:O	2.05	0.51
1:B:193:LYS:HG3	1:B:194:ASP:N	2.26	0.50
1:A:234:TYR:CE1	1:A:266:GLU:HG2	2.46	0.50
1:A:100:LYS:HD3	1:A:102:GLU:OE2	2.12	0.49
1:A:335:GLU:OE1	1:A:335:GLU:N	2.43	0.49
1:B:293:ILE:HG22	1:B:294:TYR:HD1	1.77	0.49
1:A:197:ILE:H	1:A:197:ILE:HD12	1.77	0.49
1:A:33:ILE:HD12	1:A:88:LEU:O	2.12	0.49
1:A:64:ALA:O	1:A:68:ARG:HG3	2.13	0.49
1:B:182:PHE:CE2	1:B:203:VAL:HG23	2.47	0.48
1:A:195:SER:O	1:A:199:GLN:HG3	2.14	0.48
1:A:315:LYS:O	1:A:315:LYS:HD2	2.14	0.47
1:A:35:VAL:HG12	1:A:38:LEU:CD2	2.44	0.47
1:B:42:THR:HG22	1:B:45:ASN:H	1.80	0.47
1:B:182:PHE:HD1	1:B:258:ILE:HG13	1.79	0.47
1:B:295:LYS:HG2	1:B:296:ASP:N	2.29	0.47
1:A:293:ILE:HG22	1:A:294:TYR:HD2	1.80	0.46
1:B:50:LEU:HD21	1:B:90:PHE:CE2	2.50	0.46
1:A:255:LYS:H	1:A:255:LYS:HG3	1.52	0.46
1:A:257:HIS:C	1:A:258:ILE:HD13	2.40	0.46
1:B:101:LEU:HD23	1:B:169:LEU:HB2	1.97	0.46
1:A:150:ARG:NH1	1:A:175:GLU:HB3	2.31	0.46
1:A:102:GLU:HB2	1:A:170:LYS:HG2	1.98	0.45
1:A:346:ASN:ND2	2:A:411:HOH:O	2.49	0.45
1:A:102:GLU:HG3	1:A:170:LYS:HG2	1.99	0.45
1:B:251:ILE:HA	1:B:258:ILE:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:39:VAL:N	1:B:94:GLU:OE1	2.34	0.45
1:B:34:ARG:HH11	1:B:34:ARG:HG3	1.82	0.45
1:B:251:ILE:HD12	1:B:252:SER:H	1.81	0.45
1:B:142:GLU:HG3	1:B:155:VAL:HG22	1.99	0.45
1:B:193:LYS:HG3	1:B:194:ASP:H	1.81	0.45
1:B:326:VAL:HG22	1:B:345:VAL:HG12	1.98	0.45
1:A:63:ILE:HD12	1:A:64:ALA:H	1.82	0.44
1:A:101:LEU:HD23	1:A:169:LEU:HB2	1.98	0.44
1:A:191:VAL:HG21	1:A:239:TYR:CE2	2.53	0.44
1:A:236:ASN:ND2	1:A:303:GLU:HG3	2.33	0.44
1:A:138:GLU:HG2	1:A:157:VAL:HG23	1.98	0.44
1:B:275:THR:HG21	1:B:294:TYR:CE2	2.52	0.44
1:A:33:ILE:HG13	1:A:47:TYR:HE2	1.83	0.43
1:A:335:GLU:H	1:A:335:GLU:CD	2.26	0.43
1:A:149:GLY:C	1:A:150:ARG:HD2	2.43	0.43
1:A:289:GLN:O	1:A:293:ILE:HG12	2.19	0.43
1:A:141:LEU:HB2	1:A:155:VAL:HG21	2.00	0.43
1:B:39:VAL:HG11	1:B:165:ASN:HB2	2.01	0.43
1:B:307:ALA:HA	1:B:310:GLN:HG2	2.01	0.43
1:B:98:ILE:HG13	1:B:129:PHE:HB2	2.01	0.43
1:B:200:ALA:HB2	1:B:233:MSE:SE	2.69	0.43
1:A:162:ARG:HH21	1:A:166:ARG:HH22	1.67	0.42
1:B:212:VAL:HG13	1:B:213:THR:HG23	2.01	0.42
1:A:32:ASP:OD1	1:A:87:THR:HG22	2.19	0.42
1:A:150:ARG:HH11	1:A:150:ARG:HG3	1.85	0.42
1:A:272:PHE:HE1	1:A:341:LEU:HD11	1.85	0.42
1:A:95:ARG:NH1	1:A:130:LYS:HE2	2.35	0.42
1:A:214:ARG:HH11	1:A:214:ARG:HG2	1.85	0.42
1:B:201:PHE:CE2	1:B:260:ILE:HG21	2.54	0.42
1:A:251:ILE:HG23	1:A:258:ILE:HD12	2.01	0.42
1:B:288:LEU:HD23	1:B:288:LEU:HA	1.92	0.42
1:B:350:GLN:OE1	1:B:352:THR:HB	2.20	0.42
1:A:240:ILE:HG21	1:A:332:ILE:HD13	2.02	0.41
1:B:184:ILE:HD12	1:B:260:ILE:HB	2.01	0.41
1:B:240:ILE:HD13	1:B:300:TYR:CE2	2.55	0.41
1:A:307:ALA:O	1:A:311:LEU:HG	2.20	0.41
1:A:39:VAL:HG11	1:A:165:ASN:HB2	2.02	0.41
1:A:111:LYS:HA	1:A:114:LEU:HD11	2.03	0.41
1:B:84:GLU:O	1:B:86:ASP:N	2.53	0.41
1:A:334:ASN:ND2	2:A:413:HOH:O	2.52	0.41
1:B:310:GLN:HG3	1:B:311:LEU:N	2.35	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:240:ILE:O	1:B:354:ARG:NH2	2.54	0.41
1:A:307:ALA:HA	1:A:310:GLN:HG2	2.02	0.41
1:B:121:MSE:CE	1:B:136:THR:HG22	2.51	0.41
1:A:285:PRO:O	1:A:286:GLU:HB2	2.21	0.41
1:A:34:ARG:HG3	1:A:89:VAL:HG13	2.03	0.41
1:A:284:LYS:HA	1:A:284:LYS:HD2	1.78	0.41
1:B:234:TYR:CE1	1:B:266:GLU:HG2	2.56	0.41
1:A:240:ILE:HG13	1:B:354:ARG:NH2	2.36	0.41
1:A:240:ILE:HG22	1:A:270:PHE:CD1	2.56	0.40
1:B:114:LEU:O	1:B:114:LEU:HD13	2.21	0.40
1:A:84:GLU:O	1:A:87:THR:OG1	2.31	0.40
1:B:284:LYS:HD2	1:B:284:LYS:HA	1.74	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	327/328 (100%)	306 (94%)	18 (6%)	3 (1%)	14	37
1	B	326/328 (99%)	307 (94%)	17 (5%)	2 (1%)	21	47
All	All	653/656 (100%)	613 (94%)	35 (5%)	5 (1%)	16	40

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	282	LEU
1	B	152	ASP
1	B	85	ASN
1	A	288	LEU
1	A	285	PRO

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	247/278 (89%)	242 (98%)	5 (2%)	48	77
1	B	242/278 (87%)	230 (95%)	12 (5%)	22	51
All	All	489/556 (88%)	472 (96%)	17 (4%)	32	64

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	THR
1	A	145	TYR
1	A	185	ASN
1	A	255	LYS
1	A	275	THR
1	B	29	VAL
1	B	48	THR
1	B	54	SER
1	B	82	SER
1	B	87	THR
1	B	137	ILE
1	B	146	THR
1	B	203	VAL
1	B	301	SER
1	B	305	VAL
1	B	310	GLN
1	B	352	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	ASN
1	A	236	ASN
1	A	250	ASN
1	A	274	GLN
1	A	346	ASN

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Mol	Chain	Res	Type
1	B	245	ASN
1	B	327	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](#)

There are no ligands in this entry.

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	A	323/328 (98%)	0.66	20 (6%) 26 21	35, 81, 128, 158	1 (0%)
1	B	323/328 (98%)	0.58	22 (6%) 23 18	34, 77, 132, 161	0
All	All	646/656 (98%)	0.62	42 (6%) 25 20	34, 79, 129, 161	1 (0%)

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	283	TYR	6.5
1	B	141	LEU	5.0
1	B	283	TYR	4.3
1	A	285	PRO	3.7
1	B	282	LEU	3.6
1	B	286	GLU	3.4
1	B	177	THR	3.2
1	B	178	PRO	3.1
1	A	175	GLU	3.1
1	B	118	LEU	3.0
1	B	114	LEU	2.9
1	A	280	ASP	2.9
1	B	137	ILE	2.9
1	A	209	ALA	2.8
1	B	113	ALA	2.7
1	A	208	TRP	2.7
1	A	211	VAL	2.7
1	B	203	VAL	2.7
1	A	86	ASP	2.7
1	A	192	PHE	2.6
1	A	54	SER	2.5
1	B	172	ASN	2.4
1	B	179	ALA	2.4
1	B	151	TYR	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	219	ALA	2.3
1	B	64	ALA	2.3
1	A	145	TYR	2.3
1	B	288	LEU	2.3
1	A	177	THR	2.3
1	B	265	ASP	2.3
1	A	55	GLY	2.2
1	A	275	THR	2.2
1	A	296	ASP	2.2
1	B	173	PHE	2.2
1	A	187	ILE	2.1
1	A	345	VAL	2.1
1	B	216	ASP	2.1
1	A	155	VAL	2.1
1	B	279	GLY	2.0
1	B	214	ARG	2.0
1	A	153	ALA	2.0
1	B	144	GLN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.