



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 05:16 AM UTC

PDB ID : 6J09 / pdb_00006j09
Title : Crystal structure of Haemophilus Influenzae BamA POTRA1-4
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Deposited on : 2018-12-21
Resolution : 3.00 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
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

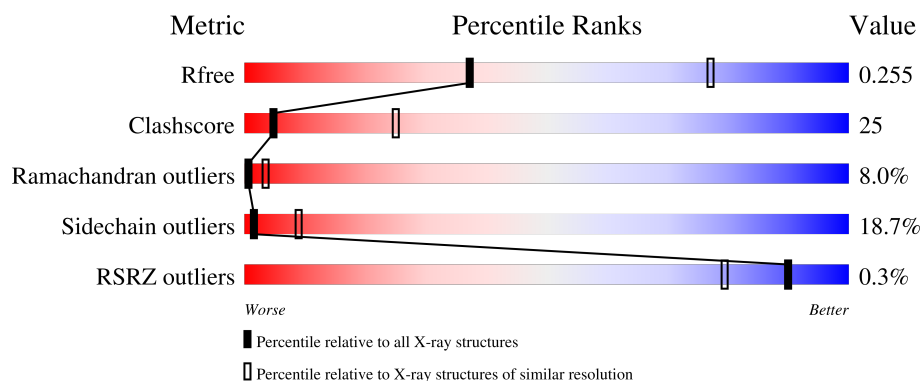
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	333	41% 38% 16% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2581 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	329	Total	C	N	O	S	0	0	0
			2552	1592	452	506	2			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	expression tag	UNP A0A2X1PZ23
A	2	SER	-	expression tag	UNP A0A2X1PZ23
A	3	HIS	-	expression tag	UNP A0A2X1PZ23
A	4	MET	-	expression tag	UNP A0A2X1PZ23

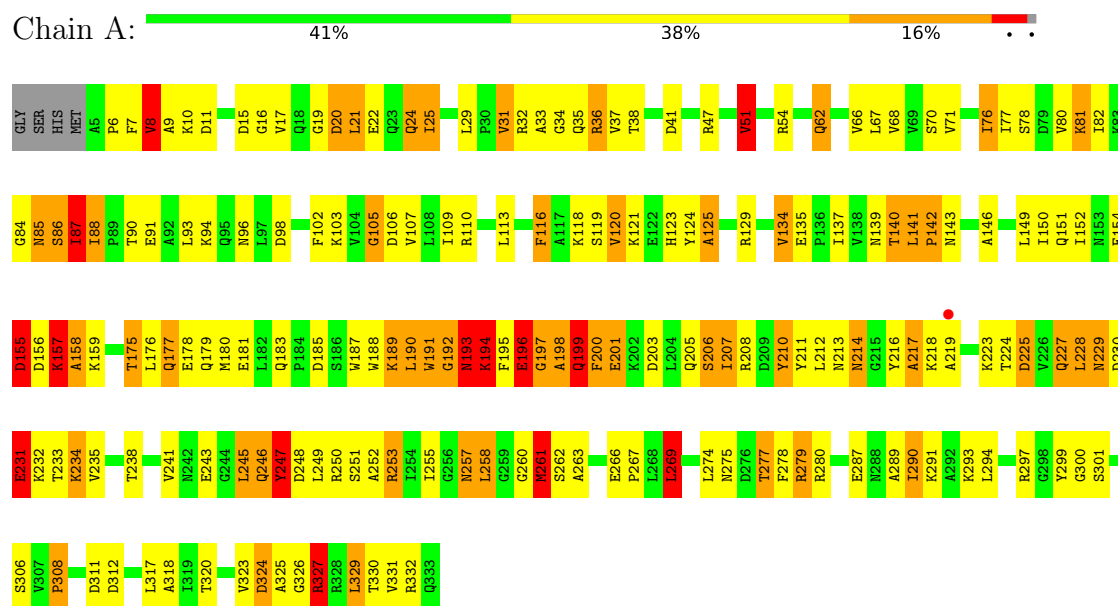
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	29	Total	O	0	0
			29	29		

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: Outer membrane protein assembly factor BamA



4 Data and refinement statistics

Property	Value	Source
Space group	P 62 2 2	Depositor
Cell constants a, b, c, α , β , γ	127.34Å 127.34Å 136.84Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	58.10 – 3.00 58.10 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (58.10-3.00) 99.8 (58.10-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.20	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.46 (at 3.01Å)	Xtriage
Refinement program	REFMAC 5.8.0218	Depositor
R, R_{free}	0.191 , 0.252 0.200 , 0.255	Depositor DCC
R_{free} test set	682 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	98.7	Xtriage
Anisotropy	0.366	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 107.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	2581	wwPDB-VP
Average B, all atoms (Å ²)	122.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.85% 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	1.42	16/2585 (0.6%)	1.62	44/3500 (1.3%)

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	A	0	10

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	157	LYS	CA-C	-9.17	1.40	1.52
1	A	88	ILE	N-CA	-8.42	1.38	1.46
1	A	158	ALA	CA-C	-6.32	1.45	1.52
1	A	185	ASP	CA-C	-6.31	1.44	1.52
1	A	142	PRO	N-CA	-6.06	1.39	1.47
1	A	199	GLN	CA-C	6.03	1.60	1.52
1	A	332	ARG	N-CA	5.97	1.53	1.46
1	A	306	SER	CA-C	5.77	1.59	1.53
1	A	70	SER	C-O	-5.75	1.16	1.23
1	A	125	ALA	N-CA	-5.63	1.38	1.46
1	A	80	VAL	CA-CB	-5.58	1.47	1.53
1	A	78	SER	C-O	-5.30	1.19	1.23
1	A	191	TRP	CA-C	-5.24	1.46	1.53
1	A	194	LYS	CA-CB	5.23	1.62	1.53
1	A	41	ASP	CA-C	-5.12	1.46	1.52
1	A	6	PRO	N-CA	5.04	1.53	1.47

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	88	ILE	N-CA-C	-12.09	93.94	107.73
1	A	82	ILE	N-CA-C	9.27	122.02	107.28
1	A	109	ILE	CB-CA-C	-8.16	99.88	111.31
1	A	82	ILE	CB-CA-C	-8.03	100.55	110.99
1	A	200	PHE	N-CA-C	-7.72	103.66	113.23
1	A	125	ALA	N-CA-C	-7.48	104.77	114.04
1	A	51	VAL	CB-CA-C	-7.39	101.26	112.05
1	A	231	GLU	N-CA-C	-7.31	100.69	110.55
1	A	134	VAL	CB-CA-C	-7.28	103.88	111.23
1	A	141	LEU	CA-C-N	-7.11	113.38	120.21
1	A	141	LEU	C-N-CA	-7.11	113.38	120.21
1	A	80	VAL	N-CA-C	6.83	117.48	107.37
1	A	110	ARG	N-CA-C	6.71	119.43	111.71
1	A	16	GLY	N-CA-C	6.49	122.29	113.99
1	A	152	ILE	N-CA-C	6.33	117.14	107.77
1	A	331	VAL	N-CA-C	-6.29	99.18	109.12
1	A	68	VAL	N-CA-C	6.21	117.65	109.21
1	A	290	ILE	CB-CA-C	-6.17	103.96	112.04
1	A	155	ASP	CA-C-N	-6.16	112.97	122.09
1	A	155	ASP	C-N-CA	-6.16	112.97	122.09
1	A	80	VAL	CB-CA-C	-6.14	102.66	111.19
1	A	21	LEU	N-CA-C	-6.09	104.83	112.93
1	A	210	TYR	N-CA-C	6.07	119.79	112.38
1	A	158	ALA	O-C-N	6.04	129.32	123.29
1	A	306	SER	N-CA-C	5.82	118.65	109.39
1	A	332	ARG	N-CA-C	5.71	117.55	107.61
1	A	135	GLU	CA-C-N	5.70	126.14	119.99
1	A	135	GLU	C-N-CA	5.70	126.14	119.99
1	A	98	ASP	N-CA-C	-5.69	105.06	112.23
1	A	102	PHE	N-CA-C	-5.57	103.87	111.56
1	A	41	ASP	N-CA-C	-5.55	105.31	111.36
1	A	261	MET	N-CA-C	5.55	122.62	110.80
1	A	120	VAL	N-CA-C	-5.39	105.28	110.72
1	A	253	ARG	N-CA-C	5.38	118.39	109.46
1	A	25	ILE	O-C-N	5.36	127.48	121.90
1	A	199	GLN	N-CA-C	5.34	122.18	110.80
1	A	247	TYR	N-CA-C	5.29	117.16	110.33
1	A	150	ILE	CB-CA-C	-5.27	104.56	111.25
1	A	68	VAL	CB-CA-C	-5.20	103.95	111.34
1	A	8	VAL	CB-CA-C	-5.18	103.88	110.98
1	A	157	LYS	O-C-N	5.17	129.47	122.59
1	A	155	ASP	N-CA-C	-5.08	102.63	109.95
1	A	116	PHE	N-CA-C	-5.04	105.87	111.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	312	ASP	N-CA-C	-5.01	105.50	110.97

There are no chirality outliers.

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	GLY	Peptide
1	A	155	ASP	Peptide
1	A	157	LYS	Peptide
1	A	190	LEU	Peptide
1	A	191	TRP	Peptide
1	A	193	ASN	Peptide
1	A	194	LYS	Peptide
1	A	217	ALA	Peptide
1	A	269	LEU	Peptide
1	A	87	ILE	Peptide

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	2552	0	2549	130	2
2	A	29	0	0	7	0
All	All	2581	0	2549	130	2

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 25.

All (130) 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:261:MET:HE1	1:A:293:LYS:HG3	1.42	1.01
1:A:229:ASN:HB3	1:A:231:GLU:O	1.62	0.99
1:A:280:ARG:NH2	1:A:287:GLU:OE2	2.02	0.91
1:A:188:TRP:CD1	1:A:188:TRP:O	2.27	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:261:MET:CE	1:A:293:LYS:HG3	2.09	0.81
1:A:294:LEU:O	1:A:299:TYR:HB2	1.80	0.81
1:A:193:ASN:OD1	1:A:194:LYS:N	2.16	0.78
1:A:47:ARG:O	1:A:51:VAL:HG23	1.84	0.77
1:A:280:ARG:HD2	2:A:409:HOH:O	1.84	0.76
1:A:195:PHE:O	1:A:196:GLU:O	2.04	0.75
1:A:245:LEU:HB2	1:A:247:TYR:CE2	2.25	0.72
1:A:279:ARG:HB2	1:A:279:ARG:HH11	1.55	0.71
1:A:85:ASN:OD1	1:A:90:THR:OG1	2.05	0.70
1:A:85:ASN:ND2	1:A:88:ILE:O	2.24	0.70
1:A:201:GLU:HA	1:A:201:GLU:OE1	1.89	0.70
1:A:252:ALA:O	1:A:269:LEU:CD2	2.40	0.70
1:A:326:GLY:O	1:A:327:ARG:HG3	1.91	0.70
1:A:198:ALA:O	1:A:201:GLU:HB3	1.93	0.69
1:A:76:ILE:HD12	1:A:105:GLY:O	1.94	0.67
1:A:20:ASP:OD2	1:A:143:ASN:ND2	2.26	0.67
1:A:253:ARG:HH12	1:A:320:THR:HG23	1.60	0.66
1:A:124:TYR:CE1	1:A:154:GLU:HG2	2.30	0.66
1:A:88:ILE:HD12	1:A:124:TYR:CZ	2.31	0.65
1:A:177:GLN:NE2	1:A:183:GLN:OE1	2.28	0.65
1:A:250:ARG:O	1:A:274:LEU:HD13	1.97	0.65
1:A:158:ALA:CB	1:A:233:THR:HA	2.27	0.65
1:A:197:GLY:HA3	2:A:412:HOH:O	1.96	0.64
1:A:297:ARG:HD3	2:A:416:HOH:O	1.97	0.64
1:A:326:GLY:C	1:A:327:ARG:HG3	2.23	0.64
1:A:19:GLY:O	1:A:21:LEU:N	2.31	0.64
1:A:245:LEU:HB2	1:A:247:TYR:HE2	1.63	0.63
1:A:76:ILE:HG13	1:A:77:ILE:N	2.13	0.63
1:A:123:HIS:HB3	2:A:423:HOH:O	2.00	0.62
1:A:260:GLY:O	1:A:262:SER:N	2.32	0.62
1:A:188:TRP:O	1:A:188:TRP:HD1	1.81	0.59
1:A:250:ARG:NH2	1:A:318:ALA:HB2	2.17	0.59
1:A:32:ARG:O	1:A:33:ALA:C	2.45	0.59
1:A:213:ASN:O	1:A:214:ASN:CG	2.45	0.59
1:A:207:ILE:O	1:A:208:ARG:C	2.47	0.57
1:A:250:ARG:HG3	1:A:251:SER:N	2.20	0.57
1:A:158:ALA:HB3	1:A:233:THR:HA	1.88	0.56
1:A:197:GLY:O	1:A:200:PHE:N	2.37	0.56
1:A:176:LEU:O	1:A:179:GLN:HB2	2.06	0.55
1:A:195:PHE:O	1:A:196:GLU:C	2.49	0.55
1:A:233:THR:O	1:A:234:LYS:O	2.25	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:196:GLU:O	1:A:198:ALA:N	2.41	0.54
1:A:208:ARG:CZ	1:A:208:ARG:HB2	2.38	0.53
1:A:308:PRO:HB2	1:A:317:LEU:HD21	1.90	0.53
1:A:261:MET:O	1:A:262:SER:C	2.51	0.53
1:A:188:TRP:O	1:A:189:LYS:C	2.52	0.53
1:A:19:GLY:O	1:A:20:ASP:C	2.51	0.52
1:A:201:GLU:OE1	1:A:201:GLU:CA	2.55	0.52
1:A:299:TYR:CD1	1:A:325:ALA:HB1	2.44	0.52
1:A:248:ASP:OD2	1:A:277:THR:HA	2.09	0.52
1:A:87:ILE:HG12	1:A:154:GLU:OE2	2.10	0.51
1:A:88:ILE:HG21	1:A:93:LEU:HD21	1.93	0.51
1:A:96:ASN:HD22	1:A:116:PHE:HE1	1.58	0.51
1:A:180:MET:HG2	1:A:203:ASP:HB3	1.93	0.50
1:A:188:TRP:O	1:A:188:TRP:CG	2.61	0.50
1:A:17:VAL:HG11	1:A:21:LEU:HD23	1.93	0.50
1:A:190:LEU:HB2	2:A:403:HOH:O	2.12	0.50
1:A:297:ARG:CD	2:A:416:HOH:O	2.58	0.49
1:A:326:GLY:C	1:A:327:ARG:CG	2.85	0.49
1:A:325:ALA:O	1:A:327:ARG:N	2.45	0.49
1:A:7:PHE:CZ	1:A:62:GLN:HB2	2.47	0.49
1:A:156:ASP:O	1:A:157:LYS:CB	2.60	0.49
1:A:289:ALA:O	1:A:290:ILE:C	2.56	0.48
1:A:90:THR:O	1:A:91:GLU:C	2.57	0.48
1:A:326:GLY:O	1:A:327:ARG:O	2.32	0.48
1:A:140:THR:O	1:A:141:LEU:HD23	2.13	0.48
1:A:139:ASN:O	1:A:146:ALA:HB1	2.14	0.48
1:A:140:THR:HA	1:A:146:ALA:HB2	1.96	0.47
1:A:155:ASP:OD2	1:A:157:LYS:O	2.32	0.47
1:A:308:PRO:CB	1:A:317:LEU:HD21	2.44	0.47
1:A:278:PHE:CD2	1:A:279:ARG:N	2.83	0.47
1:A:84:GLY:O	1:A:85:ASN:C	2.58	0.47
1:A:258:LEU:N	1:A:258:LEU:HD13	2.29	0.47
1:A:207:ILE:O	1:A:210:TYR:N	2.48	0.47
1:A:257:ASN:HD22	1:A:257:ASN:N	2.13	0.47
1:A:124:TYR:CE1	1:A:154:GLU:CG	2.97	0.46
1:A:249:LEU:O	1:A:275:ASN:N	2.47	0.46
1:A:31:VAL:HG22	1:A:35:GLN:CD	2.40	0.46
1:A:253:ARG:O	1:A:320:THR:HA	2.14	0.46
1:A:123:HIS:C	1:A:125:ALA:H	2.23	0.46
1:A:247:TYR:N	1:A:247:TYR:HD2	2.13	0.46
1:A:250:ARG:HG3	1:A:251:SER:H	1.79	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:LEU:HA	1:A:24:GLN:HG3	1.97	0.45
1:A:54:ARG:O	1:A:54:ARG:HD2	2.16	0.45
1:A:81:LYS:O	1:A:149:LEU:HA	2.15	0.45
1:A:190:LEU:O	1:A:192:GLY:N	2.49	0.45
1:A:299:TYR:O	1:A:301:SER:N	2.47	0.45
1:A:211:TYR:HE1	1:A:243:GLU:HG2	1.81	0.45
1:A:247:TYR:N	1:A:247:TYR:CD2	2.84	0.45
1:A:257:ASN:N	1:A:257:ASN:ND2	2.65	0.45
1:A:88:ILE:HD12	1:A:124:TYR:CE1	2.52	0.44
1:A:193:ASN:HB2	1:A:194:LYS:HB2	1.99	0.44
1:A:103:LYS:HB2	1:A:106:ASP:OD2	2.17	0.44
1:A:205:GLN:O	1:A:208:ARG:N	2.51	0.44
1:A:205:GLN:O	1:A:206:SER:C	2.61	0.44
1:A:262:SER:O	1:A:266:GLU:HG2	2.18	0.44
1:A:287:GLU:O	1:A:291:LYS:HB2	2.18	0.44
1:A:9:ALA:O	1:A:34:GLY:N	2.49	0.44
1:A:129:ARG:HB3	1:A:154:GLU:O	2.18	0.44
1:A:21:LEU:O	1:A:22:GLU:C	2.61	0.43
1:A:212:LEU:CD2	1:A:217:ALA:HA	2.47	0.43
1:A:25:ILE:O	1:A:29:LEU:HD13	2.18	0.43
1:A:227:GLN:HE21	1:A:227:GLN:HA	1.83	0.43
1:A:157:LYS:HA	1:A:190:LEU:CD2	2.49	0.42
1:A:181:GLU:OE2	1:A:199:GLN:OE1	2.36	0.42
1:A:266:GLU:N	1:A:267:PRO:HD2	2.35	0.42
1:A:199:GLN:O	1:A:199:GLN:HG3	2.18	0.42
1:A:192:GLY:O	1:A:193:ASN:C	2.62	0.42
1:A:278:PHE:CD2	1:A:278:PHE:C	2.97	0.42
1:A:8:VAL:HG13	1:A:36:ARG:HB2	2.02	0.42
1:A:86:SER:C	1:A:87:ILE:HG23	2.44	0.41
1:A:158:ALA:HB3	1:A:233:THR:CA	2.48	0.41
1:A:113:LEU:HA	1:A:113:LEU:HD12	1.82	0.41
1:A:141:LEU:HB3	1:A:142:PRO:HD2	2.01	0.41
1:A:216:TYR:CZ	1:A:246:GLN:HG3	2.54	0.41
1:A:67:LEU:C	1:A:67:LEU:HD12	2.46	0.41
1:A:17:VAL:HG21	1:A:25:ILE:HD12	2.03	0.41
1:A:36:ARG:HG2	1:A:37:VAL:N	2.35	0.41
1:A:120:VAL:HG12	1:A:134:VAL:HG21	2.02	0.41
1:A:158:ALA:HA	1:A:187:TRP:CE3	2.55	0.41
1:A:86:SER:OG	1:A:87:ILE:HG23	2.22	0.40
1:A:245:LEU:N	2:A:402:HOH:O	2.49	0.40
1:A:151:GLN:OE1	1:A:232:LYS:HE2	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:THR:HA	1:A:178:GLU:OE2	2.21	0.40
1:A:257:ASN:O	1:A:324:ASP:HA	2.21	0.40
1:A:279:ARG:HH11	1:A:279:ARG:CB	2.29	0.40

All (2) 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:A:224:THR:O	1:A:329:LEU:O[11_555]	2.17	0.03
1:A:225:ASP:OD1	1:A:330:THR:OG1[11_555]	2.18	0.02

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	327/333 (98%)	257 (79%)	44 (14%)	26 (8%)	1 3

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	20	ASP
1	A	193	ASN
1	A	196	GLU
1	A	197	GLY
1	A	198	ALA
1	A	214	ASN
1	A	219	ALA
1	A	234	LYS
1	A	261	MET
1	A	263	ALA
1	A	327	ARG
1	A	85	ASN

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Mol	Chain	Res	Type
1	A	157	LYS
1	A	159	LYS
1	A	192	GLY
1	A	194	LYS
1	A	207	ILE
1	A	230	ASP
1	A	199	GLN
1	A	206	SER
1	A	228	LEU
1	A	308	PRO
1	A	87	ILE
1	A	229	ASN
1	A	269	LEU
1	A	300	GLY

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	278/283 (98%)	226 (81%)	52 (19%)	1 9

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	VAL
1	A	10	LYS
1	A	11	ASP
1	A	15	ASP
1	A	24	GLN
1	A	31	VAL
1	A	36	ARG
1	A	38	THR
1	A	51	VAL
1	A	62	GLN
1	A	66	VAL
1	A	71	VAL

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Mol	Chain	Res	Type
1	A	76	ILE
1	A	81	LYS
1	A	86	SER
1	A	94	LYS
1	A	107	VAL
1	A	118	LYS
1	A	119	SER
1	A	121	LYS
1	A	137	ILE
1	A	140	THR
1	A	155	ASP
1	A	175	THR
1	A	177	GLN
1	A	189	LYS
1	A	194	LYS
1	A	196	GLU
1	A	201	GLU
1	A	218	LYS
1	A	223	LYS
1	A	225	ASP
1	A	227	GLN
1	A	228	LEU
1	A	231	GLU
1	A	235	VAL
1	A	238	THR
1	A	241	VAL
1	A	245	LEU
1	A	246	GLN
1	A	247	TYR
1	A	255	ILE
1	A	257	ASN
1	A	258	LEU
1	A	261	MET
1	A	277	THR
1	A	279	ARG
1	A	311	ASP
1	A	323	VAL
1	A	324	ASP
1	A	327	ARG
1	A	329	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	44	ASN
1	A	62	GLN
1	A	114	ASN
1	A	131	ASN
1	A	153	ASN
1	A	199	GLN
1	A	214	ASN
1	A	227	GLN
1	A	257	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 [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	329/333 (98%)	-0.44	1 (0%) 90 79	71, 123, 172, 215	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	219	ALA	2.1

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.