



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 18, 2026 – 02:37 PM UTC

PDB ID : 3L2Z / pdb_00003l2z
Title : Crystal structure of hydrated Biotin Protein Ligase from M. tuberculosis
Authors : Gupta, V.; Gupta, R.K.; Khare, G.; Salunke, D.M.; Tyagi, A.K.
Deposited on : 2009-12-16
Resolution : 2.80 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

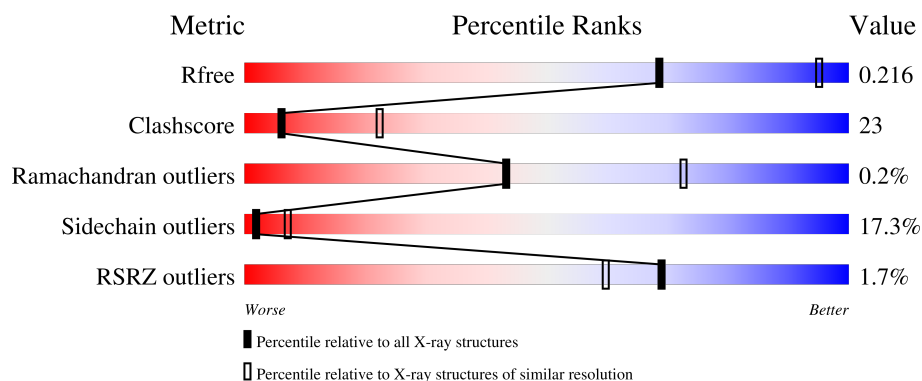
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

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	266	2% 47% 32% 11% 10%
1	B	266	2% 45% 35% 11% 9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3561 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 BirA bifunctional protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	240	Total	C	N	O	S	0	0	0
			1742	1095	318	328	1			
1	B	243	Total	C	N	O	S	0	1	0
			1793	1123	337	332	1			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	11	Total	O	0	0
			11	11		
2	B	15	Total	O	0	0
			15	15		

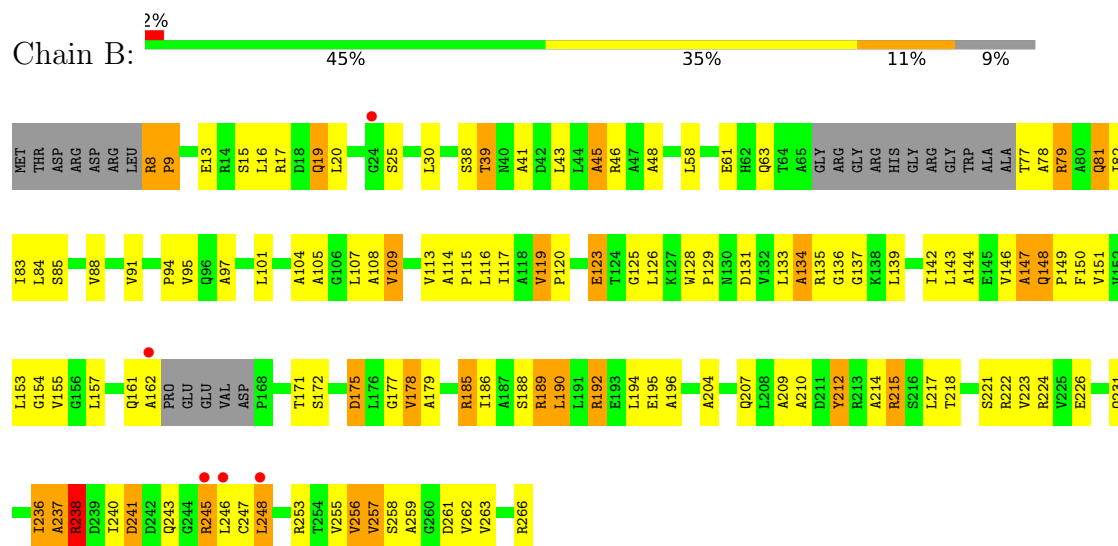
3 Residue-property plots [i](#)

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: BirA bifunctional protein



• Molecule 1: BirA bifunctional protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	79.69Å 62.77Å 105.75Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	14.13 – 2.80 14.13 – 2.80	Depositor EDS
% Data completeness (in resolution range)	99.3 (14.13-2.80) 98.3 (14.13-2.80)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.18 (at 2.81Å)	Xtriage
Refinement program	PHENIX (phenix.refine)	Depositor
R, R_{free}	0.174 , 0.231 0.194 , 0.216	Depositor DCC
R_{free} test set	1338 reflections (10.00%)	wwPDB-VP
Wilson B-factor (Å ²)	50.3	Xtriage
Anisotropy	0.158	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3561	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.06% 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 ⓘ

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	2.24	41/1765 (2.3%)	1.19	8/2412 (0.3%)
1	B	2.19	40/1819 (2.2%)	1.22	8/2480 (0.3%)
All	All	2.22	81/3584 (2.3%)	1.20	16/4892 (0.3%)

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	B	0	1

All (81) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	237	ALA	CA-CB	-7.73	1.43	1.53
1	B	240	ILE	C-O	-6.88	1.16	1.24
1	A	155	VAL	C-O	-6.55	1.17	1.24
1	A	233	VAL	C-O	-6.52	1.17	1.24
1	A	83	ILE	C-O	-6.49	1.17	1.24
1	B	204	ALA	CA-CB	-6.48	1.43	1.53
1	A	237	ALA	CA-CB	-6.25	1.44	1.53
1	A	209	ALA	CA-CB	-6.22	1.43	1.53
1	B	257	VAL	C-O	-6.19	1.17	1.24
1	A	240	ILE	C-O	-6.17	1.17	1.24
1	B	147	ALA	CA-CB	-6.14	1.46	1.53
1	B	209	ALA	CA-CB	-6.14	1.43	1.53
1	A	223	VAL	C-O	-6.02	1.17	1.23
1	A	210	ALA	CA-CB	-6.02	1.44	1.53
1	A	88	VAL	C-O	-5.91	1.18	1.24
1	A	263	VAL	C-O	-5.85	1.18	1.24
1	A	180	ALA	C-O	-5.78	1.20	1.25
1	A	59	ILE	C-O	-5.75	1.18	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	209	ALA	C-O	-5.75	1.17	1.24
1	B	108	ALA	CA-CB	-5.73	1.44	1.53
1	B	255	VAL	C-O	-5.72	1.18	1.24
1	A	247	CYS	C-O	-5.71	1.17	1.24
1	A	187	ALA	CA-CB	-5.70	1.44	1.53
1	B	263	VAL	C-O	-5.70	1.18	1.24
1	B	154	GLY	CA-C	-5.66	1.48	1.52
1	A	200	GLN	C-O	-5.61	1.17	1.24
1	B	104	ALA	CA-CB	-5.59	1.44	1.53
1	B	194	LEU	C-O	-5.58	1.17	1.24
1	B	153	LEU	C-N	-5.56	1.29	1.33
1	A	110	LEU	C-O	-5.53	1.17	1.24
1	B	109	VAL	C-O	-5.52	1.18	1.24
1	A	235	GLY	C-O	-5.50	1.18	1.23
1	B	45	ALA	CA-CB	-5.50	1.44	1.53
1	B	256	VAL	C-O	-5.50	1.18	1.24
1	A	249	ASP	C-O	-5.46	1.17	1.24
1	A	105	ALA	C-O	-5.45	1.17	1.24
1	A	257	VAL	C-O	-5.43	1.18	1.24
1	B	97	ALA	CA-CB	-5.41	1.44	1.53
1	A	187	ALA	C-O	-5.41	1.17	1.24
1	A	126	LEU	C-O	-5.40	1.17	1.24
1	B	155	VAL	C-O	-5.39	1.18	1.24
1	A	152	VAL	C-O	-5.38	1.18	1.24
1	A	194	LEU	C-O	-5.36	1.17	1.24
1	B	231	GLN	C-O	-5.36	1.17	1.24
1	B	261	ASP	C-O	-5.33	1.17	1.24
1	A	108	ALA	CA-CB	-5.29	1.45	1.53
1	A	113	VAL	C-O	-5.27	1.19	1.24
1	A	31	ASP	C-O	-5.27	1.17	1.23
1	B	214	ALA	CA-CB	-5.25	1.45	1.53
1	A	212	TYR	C-O	-5.24	1.18	1.24
1	B	212	TYR	C-O	-5.21	1.18	1.24
1	B	238	ARG	C-N	-5.21	1.26	1.33
1	B	236	ILE	C-O	-5.20	1.18	1.24
1	B	144	ALA	CA-CB	-5.19	1.45	1.53
1	A	186	ILE	CA-CB	-5.19	1.47	1.54
1	B	196	ALA	CA-CB	-5.18	1.45	1.53
1	A	105	ALA	CA-CB	-5.18	1.45	1.53
1	B	94	PRO	C-O	-5.17	1.17	1.23
1	B	195	GLU	C-O	-5.17	1.18	1.24
1	B	210	ALA	CA-CB	-5.17	1.45	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	107	LEU	C-O	-5.17	1.18	1.24
1	A	192	ARG	C-O	-5.16	1.18	1.24
1	B	142	ILE	C-O	-5.16	1.18	1.24
1	B	155	VAL	CA-CB	-5.15	1.48	1.54
1	B	134	ALA	CA-CB	-5.13	1.45	1.53
1	A	211	ASP	C-O	-5.12	1.18	1.24
1	B	153	LEU	C-O	-5.11	1.18	1.24
1	A	210	ALA	C-O	-5.09	1.18	1.24
1	B	192	ARG	C-O	-5.09	1.18	1.24
1	B	9	PRO	C-O	-5.09	1.20	1.24
1	A	125	GLY	C-O	-5.09	1.17	1.23
1	A	248	LEU	C-O	-5.08	1.17	1.24
1	A	255	VAL	C-O	-5.07	1.18	1.24
1	B	88	VAL	C-O	-5.06	1.19	1.24
1	B	236	ILE	CA-C	-5.05	1.46	1.52
1	A	14	ARG	C-O	-5.03	1.18	1.24
1	A	143	LEU	C-O	-5.03	1.18	1.24
1	A	190	LEU	C-O	-5.03	1.18	1.24
1	A	57	VAL	C-O	-5.03	1.18	1.24
1	B	143	LEU	C-O	-5.03	1.18	1.24
1	A	47	ALA	CA-CB	-5.02	1.45	1.53

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	19	GLN	N-CA-C	9.04	121.21	111.36
1	B	179	ALA	N-CA-C	9.04	121.21	111.36
1	B	139	LEU	N-CA-C	7.78	119.83	111.36
1	A	218	THR	N-CA-C	6.73	118.70	111.36
1	B	218	THR	N-CA-C	6.17	118.09	111.36
1	B	186	ILE	N-CA-C	6.04	116.78	110.62
1	A	139	LEU	N-CA-C	5.96	117.85	111.36
1	B	101	LEU	N-CA-C	5.88	117.77	111.36
1	A	101	LEU	N-CA-C	5.78	117.66	111.36
1	A	179	ALA	N-CA-C	5.58	117.04	111.07
1	A	131	ASP	N-CA-C	5.46	118.30	109.40
1	B	131	ASP	N-CA-C	5.41	118.22	109.40
1	B	20	LEU	N-CA-C	5.40	120.55	112.94
1	A	37	GLY	N-CA-C	5.36	118.98	112.49
1	A	130	ASN	N-CA-C	5.33	120.65	113.72
1	B	178	VAL	N-CA-C	-5.15	101.98	108.93

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	238	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	A	1742	0	1768	71	0
1	B	1793	0	1847	94	1
2	A	11	0	0	2	1
2	B	15	0	0	2	0
All	All	3561	0	3615	161	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 23.

All (161) 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:119:VAL:HG23	1:B:120:PRO:CD	1.71	1.20
1:A:189:ARG:HH11	1:A:189:ARG:CG	1.58	1.11
1:B:119:VAL:HG23	1:B:120:PRO:HD2	1.09	1.08
1:A:189:ARG:HG2	1:A:189:ARG:NH1	1.46	1.06
1:B:246:LEU:O	1:B:248:LEU:HD23	1.59	1.01
1:A:11:LEU:HG	1:A:183:ARG:HH12	1.30	0.97
1:A:260:GLY:HA3	1:B:253:ARG:NH1	1.78	0.97
1:A:194:LEU:O	1:A:198:ILE:HG13	1.69	0.93
1:B:17:ARG:NH2	1:B:30:LEU:O	2.01	0.93
1:B:258:SER:O	1:B:259:ALA:C	2.13	0.88
1:A:260:GLY:HA3	1:B:253:ARG:CZ	2.03	0.86
1:B:17:ARG:CZ	1:B:30:LEU:O	2.23	0.86
1:B:117:ILE:CG2	1:B:119:VAL:CG1	2.55	0.85
1:B:245:ARG:HD2	1:B:256:VAL:HG12	1.59	0.84
1:B:117:ILE:CG2	1:B:119:VAL:HG13	2.08	0.84
1:A:224:ARG:HD2	1:A:226:GLU:OE2	1.78	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:105:ALA:O	1:B:109:VAL:HG23	1.79	0.82
1:B:117:ILE:HG22	1:B:119:VAL:CG1	2.10	0.81
1:B:241:ASP:OD2	1:B:243:GLN:N	2.14	0.81
1:B:246:LEU:HG	1:B:248:LEU:HD21	1.63	0.80
1:A:189:ARG:HH11	1:A:189:ARG:HG2	0.68	0.79
1:B:266:ARG:CG	1:B:266:ARG:O	2.29	0.79
1:A:182:ASP:O	1:A:186:ILE:HG13	1.83	0.79
1:B:8:ARG:HG2	1:B:79:ARG:HB3	1.64	0.78
1:B:236:ILE:HG21	1:B:238:ARG:HH11	1.48	0.78
1:B:266:ARG:O	1:B:266:ARG:HG2	1.84	0.77
1:B:119:VAL:CG2	1:B:120:PRO:CD	2.61	0.76
1:B:241:ASP:OD1	1:B:245:ARG:NE	2.17	0.76
1:A:22:GLY:O	1:A:23:ALA:O	2.04	0.75
1:A:46:ARG:O	1:A:51:ALA:HB3	1.86	0.74
1:A:203:ASN:O	1:A:204:ALA:C	2.29	0.74
1:B:246:LEU:O	1:B:248:LEU:CD2	2.35	0.74
1:B:245:ARG:HD2	1:B:256:VAL:CG1	2.19	0.73
1:B:161:GLN:O	1:B:162:ALA:HB3	1.88	0.72
1:A:11:LEU:HG	1:A:183:ARG:NH1	2.03	0.72
1:A:81:GLN:NE2	1:A:158:ASN:OD1	2.22	0.72
1:B:117:ILE:HG21	1:B:119:VAL:CG1	2.19	0.71
1:B:241:ASP:OD2	1:B:241:ASP:C	2.30	0.70
1:B:237:ALA:HA	1:B:248:LEU:HD22	1.73	0.69
1:A:11:LEU:HG	1:A:61:GLU:OE2	1.93	0.68
1:B:117:ILE:HG22	1:B:119:VAL:HG13	1.74	0.67
1:A:29:GLN:NE2	1:A:31:ASP:OD2	2.26	0.67
1:B:188:SER:O	1:B:192:ARG:HG3	1.94	0.67
1:B:247:CYS:C	1:B:248:LEU:HD23	2.20	0.67
1:A:260:GLY:CA	1:B:253:ARG:NH1	2.58	0.65
1:B:117:ILE:HG22	1:B:119:VAL:HG12	1.77	0.65
1:B:13:GLU:OE2	1:B:17:ARG:NH2	2.25	0.65
1:B:248:LEU:HD23	1:B:248:LEU:N	2.12	0.64
1:B:247:CYS:C	1:B:248:LEU:CD2	2.70	0.64
1:A:241:ASP:C	1:A:241:ASP:OD2	2.38	0.63
1:B:175:ASP:OD2	1:B:175:ASP:N	2.30	0.63
1:A:264:HIS:O	1:A:265:LEU:HD23	1.99	0.62
1:B:8:ARG:N	1:B:9:PRO:HD3	2.15	0.62
1:B:161:GLN:O	1:B:162:ALA:CB	2.49	0.61
1:B:78:ALA:O	1:B:79:ARG:HB2	1.98	0.61
1:B:223:VAL:HG21	1:B:248:LEU:HD11	1.83	0.60
1:A:81:GLN:HG2	1:A:158:ASN:HA	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:ASP:OD2	1:A:175:ASP:N	2.30	0.60
1:A:182:ASP:HB3	1:A:185:ARG:HB3	1.83	0.60
1:B:15:SER:O	1:B:19:GLN:CG	2.50	0.60
1:B:117:ILE:HG21	1:B:119:VAL:HG13	1.82	0.60
1:B:248:LEU:CD2	1:B:248:LEU:N	2.63	0.60
1:A:133:LEU:HD12	1:A:138:LYS:HA	1.84	0.60
1:A:39:THR:HB	1:A:85:SER:OG	2.03	0.59
1:B:15:SER:O	1:B:19:GLN:HG2	2.02	0.58
1:B:120:PRO:HD2	1:B:123:GLU:HB2	1.84	0.58
1:A:22:GLY:O	1:A:23:ALA:C	2.47	0.58
1:A:174:LEU:HD12	1:A:178:VAL:O	2.05	0.56
1:A:194:LEU:O	1:A:198:ILE:CG1	2.49	0.56
1:A:46:ARG:O	1:A:51:ALA:CB	2.53	0.56
1:B:126:LEU:HD13	1:B:212:TYR:CE1	2.40	0.56
1:B:223:VAL:CG2	1:B:248:LEU:HD11	2.35	0.56
1:A:126:LEU:HD13	1:A:212:TYR:CE1	2.41	0.56
1:A:17:ARG:NH1	1:A:30:LEU:O	2.38	0.56
1:A:215:ARG:HG2	1:A:215:ARG:HH11	1.71	0.56
1:B:8:ARG:N	1:B:9:PRO:CD	2.68	0.56
1:A:29:GLN:HG3	1:A:30:LEU:N	2.21	0.55
1:B:128:TRP:CH2	1:B:259:ALA:HA	2.42	0.55
1:B:223:VAL:CG2	1:B:248:LEU:CD1	2.84	0.55
1:B:134:ALA:O	1:B:135:ARG:C	2.47	0.54
1:B:224:ARG:NH2	1:B:226:GLU:OE2	2.39	0.54
1:A:11:LEU:CG	1:A:61:GLU:OE2	2.55	0.54
1:B:171:THR:OG1	1:B:172:SER:N	2.40	0.54
1:A:189:ARG:CG	1:A:189:ARG:NH1	2.31	0.53
1:B:61:GLU:OE1	1:B:79:ARG:HD2	2.08	0.53
1:B:222:ARG:H	1:B:266:ARG:HB3	1.73	0.53
1:B:246:LEU:CG	1:B:248:LEU:HD21	2.36	0.53
1:A:148:GLN:HA	1:A:150:PHE:N	2.25	0.52
1:A:13:GLU:O	1:A:17:ARG:HG3	2.10	0.52
1:A:264:HIS:C	1:A:265:LEU:HD23	2.35	0.51
1:B:114:ALA:N	1:B:115:PRO:CD	2.72	0.51
1:B:17:ARG:NH1	1:B:30:LEU:O	2.44	0.51
1:B:241:ASP:CG	1:B:245:ARG:HG3	2.36	0.51
1:B:81:GLN:HB3	1:B:157:LEU:O	2.11	0.50
1:B:241:ASP:OD1	1:B:245:ARG:HG3	2.11	0.50
1:A:128:TRP:CH2	1:A:259:ALA:HA	2.48	0.49
1:B:125:GLY:HA3	1:B:217:LEU:HG	1.93	0.49
1:B:215:ARG:HD2	2:B:270:HOH:O	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:128:TRP:CE3	1:B:129:PRO:HD3	2.48	0.49
1:A:82:ILE:HD12	1:A:159:VAL:HG21	1.95	0.49
1:A:215:ARG:HH11	1:A:215:ARG:CG	2.26	0.48
1:B:15:SER:O	1:B:19:GLN:HG3	2.13	0.48
1:A:120:PRO:O	1:A:121:PRO:C	2.57	0.48
1:A:80:ALA:HB1	1:A:160:THR:O	2.13	0.47
1:B:247:CYS:C	1:B:248:LEU:HD22	2.38	0.47
1:B:123:GLU:HG3	1:B:135:ARG:HG3	1.95	0.47
1:A:123:GLU:O	2:A:268:HOH:O	2.20	0.47
1:A:131:ASP:OD1	1:A:138:LYS:NZ	2.36	0.47
1:B:84:LEU:O	1:B:84:LEU:HD12	2.14	0.47
1:B:247:CYS:O	1:B:248:LEU:HD22	2.14	0.47
1:A:258:SER:O	1:A:259:ALA:C	2.55	0.46
1:B:16:LEU:HA	1:B:19:GLN:HG3	1.97	0.46
1:B:149:PRO:HG2	1:B:150:PHE:CE1	2.50	0.46
1:A:114:ALA:HB3	1:A:115:PRO:HD3	1.97	0.46
1:B:84:LEU:HD12	1:B:84:LEU:C	2.39	0.46
1:B:119:VAL:HG23	1:B:120:PRO:N	2.08	0.46
1:B:190:LEU:C	1:B:190:LEU:CD2	2.89	0.46
1:A:114:ALA:N	1:A:115:PRO:CD	2.78	0.46
1:A:245:ARG:HB3	1:A:257:VAL:O	2.16	0.46
1:A:81:GLN:HA	1:A:159:VAL:HG23	1.98	0.46
1:A:82:ILE:HD12	1:A:159:VAL:CG2	2.46	0.45
1:B:78:ALA:O	1:B:79:ARG:CB	2.65	0.45
1:A:215:ARG:NE	2:A:276:HOH:O	2.41	0.45
1:B:39:THR:HB	1:B:85:SER:OG	2.16	0.45
1:B:45:ALA:O	1:B:48:ALA:HB3	2.17	0.45
1:B:147:ALA:O	1:B:148:GLN:C	2.58	0.45
1:B:246:LEU:C	1:B:248:LEU:HD23	2.37	0.45
1:B:177:GLY:O	1:B:178:VAL:C	2.56	0.45
1:B:236:ILE:O	1:B:248:LEU:HA	2.17	0.44
1:B:148:GLN:HA	1:B:149:PRO:C	2.44	0.43
1:B:148:GLN:HA	1:B:150:PHE:N	2.34	0.43
1:B:246:LEU:HG	1:B:248:LEU:CD2	2.43	0.43
1:A:100:TRP:CH2	1:A:209:ALA:HA	2.54	0.43
1:B:39:THR:HG22	1:B:58:LEU:HG	2.00	0.43
1:A:148:GLN:HA	1:A:150:PHE:H	1.82	0.43
1:B:136:GLY:N	2:B:268:HOH:O	2.39	0.42
1:A:146:VAL:HB	1:A:151:VAL:HG22	2.01	0.42
1:A:225:VAL:HG22	1:A:262:VAL:HG12	2.00	0.42
1:B:172:SER:OG	1:B:175:ASP:OD2	2.37	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:11:LEU:CG	1:A:183:ARG:NH1	2.76	0.42
1:A:148:GLN:H	1:A:148:GLN:HG2	1.28	0.42
1:A:114:ALA:N	1:A:115:PRO:HD2	2.35	0.42
1:A:224:ARG:NH2	1:A:226:GLU:OE2	2.28	0.41
1:A:117:ILE:HD12	1:A:117:ILE:HA	1.48	0.41
1:B:189:ARG:HA	1:B:192:ARG:HD3	2.01	0.41
1:B:190:LEU:CD2	1:B:190:LEU:O	2.69	0.41
1:A:198:ILE:HD12	1:A:198:ILE:HG21	1.79	0.41
1:A:224:ARG:HG3	1:A:234:VAL:HG22	2.02	0.41
1:A:240:ILE:HG21	1:A:240:ILE:HD12	1.81	0.41
1:A:215:ARG:CG	1:A:215:ARG:NH1	2.84	0.41
1:A:255:VAL:HG22	1:B:257:VAL:HG12	2.03	0.41
1:B:38:SER:HB3	1:B:41:ALA:HB3	2.02	0.41
1:A:23:ALA:HA	1:A:24:GLY:HA2	1.68	0.41
1:A:174:LEU:HD12	1:A:174:LEU:HA	1.68	0.41
1:A:224:ARG:CD	1:A:226:GLU:OE2	2.61	0.41
1:B:241:ASP:C	1:B:243:GLN:H	2.28	0.41
1:A:11:LEU:HD23	1:A:11:LEU:HA	1.63	0.41
1:A:168:PRO:HB2	1:A:169:ASP:H	1.58	0.40
1:B:135:ARG:C	1:B:137:GLY:H	2.28	0.40
1:A:199:ILE:HG22	1:A:200:GLN:N	2.33	0.40
1:B:117:ILE:HG22	1:B:119:VAL:H	1.86	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:185:ARG:CD	2:A:277:HOH:O[3_545]	2.10	0.10

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	234/266 (88%)	228 (97%)	5 (2%)	1 (0%)	30	60
1	B	238/266 (90%)	230 (97%)	8 (3%)	0	100	100
All	All	472/532 (89%)	458 (97%)	13 (3%)	1 (0%)	43	72

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	23	ALA

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	176/202 (87%)	147 (84%)	29 (16%)	2	8
1	B	183/202 (91%)	150 (82%)	33 (18%)	2	6
All	All	359/404 (89%)	297 (83%)	62 (17%)	2	7

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	8	ARG
1	A	17	ARG
1	A	39	THR
1	A	63	GLN
1	A	82	ILE
1	A	88	VAL
1	A	102	SER
1	A	113	VAL
1	A	116	LEU
1	A	117	ILE
1	A	127	LYS
1	A	133	LEU
1	A	146	VAL
1	A	148	GLN
1	A	159	VAL

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Mol	Chain	Res	Type
1	A	171	THR
1	A	175	ASP
1	A	189	ARG
1	A	192	ARG
1	A	194	LEU
1	A	195	GLU
1	A	198	ILE
1	A	199	ILE
1	A	202	ARG
1	A	215	ARG
1	A	221	SER
1	A	226	GLU
1	A	238	ARG
1	A	263	VAL
1	B	8	ARG
1	B	19	GLN
1	B	25	SER
1	B	39	THR
1	B	43	LEU
1	B	46	ARG
1	B	63	GLN
1	B	77	THR
1	B	79	ARG
1	B	81	GLN
1	B	82	ILE
1	B	83	ILE
1	B	91	VAL
1	B	95	VAL
1	B	113	VAL
1	B	116	LEU
1	B	119	VAL
1	B	123	GLU
1	B	133	LEU
1	B	146	VAL
1	B	148	GLN
1	B	151	VAL
1	B	175	ASP
1	B	185	ARG
1	B	189	ARG
1	B	190	LEU
1	B	207	GLN
1	B	215	ARG

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Mol	Chain	Res	Type
1	B	221	SER
1	B	241	ASP
1	B	245	ARG
1	B	248	LEU
1	B	262	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	A	130	ASN
1	A	148	GLN
1	B	19	GLN
1	B	35	GLN
1	B	81	GLN
1	B	148	GLN

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 ⓘ

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	240/266 (90%)	-0.51	3 (1%) 75 66	20, 43, 72, 95	0
1	B	243/266 (91%)	-0.41	5 (2%) 63 54	24, 42, 76, 109	1 (0%)
All	All	483/532 (90%)	-0.46	8 (1%) 69 60	20, 43, 72, 109	1 (0%)

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	161	GLN	4.1
1	B	24	GLY	4.0
1	B	248	LEU	3.6
1	A	61	GLU	3.0
1	B	245	ARG	2.8
1	B	246	LEU	2.6
1	A	160	THR	2.6
1	B	162	ALA	2.3

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.