



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:01 AM UTC

PDB ID : 7AG9 / pdb_00007ag9
Title : Structure of the Kar9 protein
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Deposited on : 2020-09-22
Resolution : 3.30 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

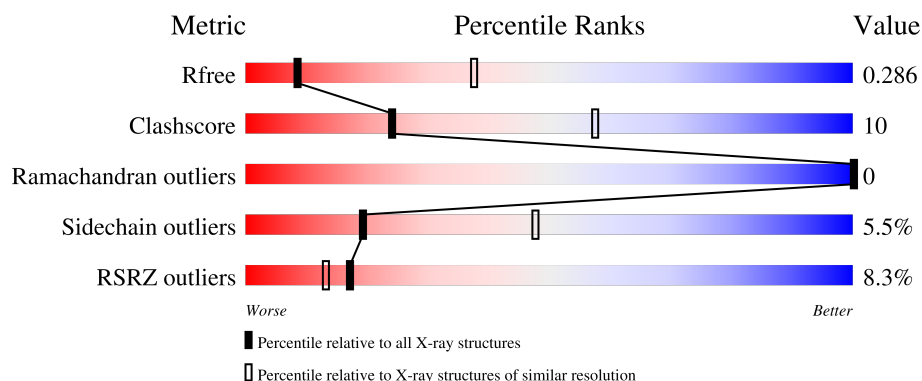
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	428	9% 69% 18% • 12%
1	B	428	6% 69% 19% • 10%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 5948 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 Kar9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	375	Total	C	N	O	S	0	0	0
			2894	1852	490	541	11			
1	B	384	Total	C	N	O	S	0	0	0
			3054	1955	513	576	10			

There are 38 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-17	MET	-	initiating methionine	UNP G0VE12
A	-16	ALA	-	expression tag	UNP G0VE12
A	-15	HIS	-	expression tag	UNP G0VE12
A	-14	HIS	-	expression tag	UNP G0VE12
A	-13	HIS	-	expression tag	UNP G0VE12
A	-12	HIS	-	expression tag	UNP G0VE12
A	-11	HIS	-	expression tag	UNP G0VE12
A	-10	HIS	-	expression tag	UNP G0VE12
A	-9	SER	-	expression tag	UNP G0VE12
A	-8	ALA	-	expression tag	UNP G0VE12
A	-7	ALA	-	expression tag	UNP G0VE12
A	-6	LEU	-	expression tag	UNP G0VE12
A	-5	GLU	-	expression tag	UNP G0VE12
A	-4	VAL	-	expression tag	UNP G0VE12
A	-3	LEU	-	expression tag	UNP G0VE12
A	-2	PHE	-	expression tag	UNP G0VE12
A	-1	GLN	-	expression tag	UNP G0VE12
A	0	GLY	-	expression tag	UNP G0VE12
A	1	PRO	-	expression tag	UNP G0VE12
B	-17	MET	-	initiating methionine	UNP G0VE12
B	-16	ALA	-	expression tag	UNP G0VE12
B	-15	HIS	-	expression tag	UNP G0VE12
B	-14	HIS	-	expression tag	UNP G0VE12
B	-13	HIS	-	expression tag	UNP G0VE12
B	-12	HIS	-	expression tag	UNP G0VE12

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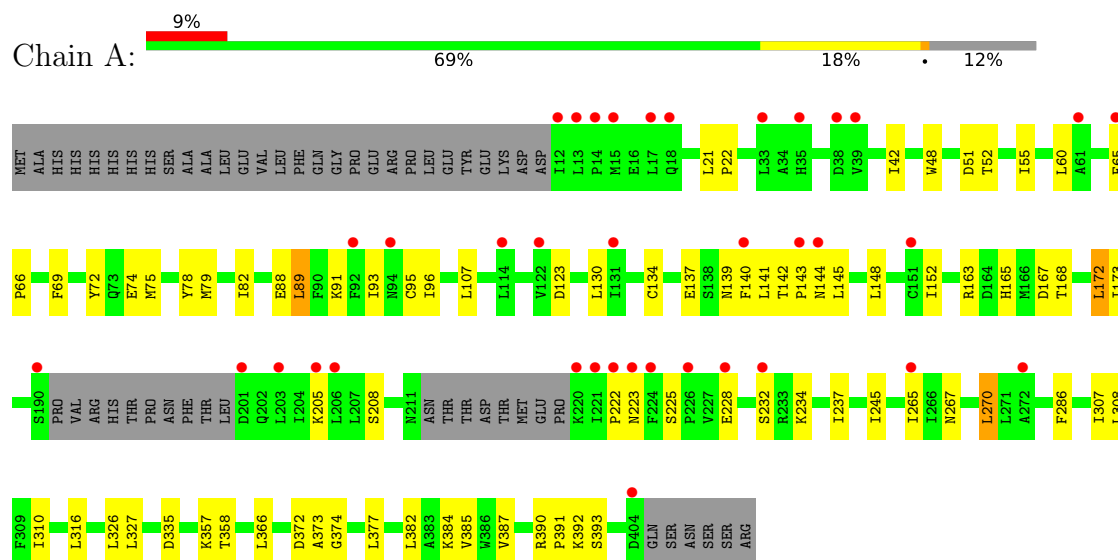
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Chain	Residue	Modelled	Actual	Comment	Reference
B	-11	HIS	-	expression tag	UNP G0VE12
B	-10	HIS	-	expression tag	UNP G0VE12
B	-9	SER	-	expression tag	UNP G0VE12
B	-8	ALA	-	expression tag	UNP G0VE12
B	-7	ALA	-	expression tag	UNP G0VE12
B	-6	LEU	-	expression tag	UNP G0VE12
B	-5	GLU	-	expression tag	UNP G0VE12
B	-4	VAL	-	expression tag	UNP G0VE12
B	-3	LEU	-	expression tag	UNP G0VE12
B	-2	PHE	-	expression tag	UNP G0VE12
B	-1	GLN	-	expression tag	UNP G0VE12
B	0	GLY	-	expression tag	UNP G0VE12
B	1	PRO	-	expression tag	UNP G0VE12

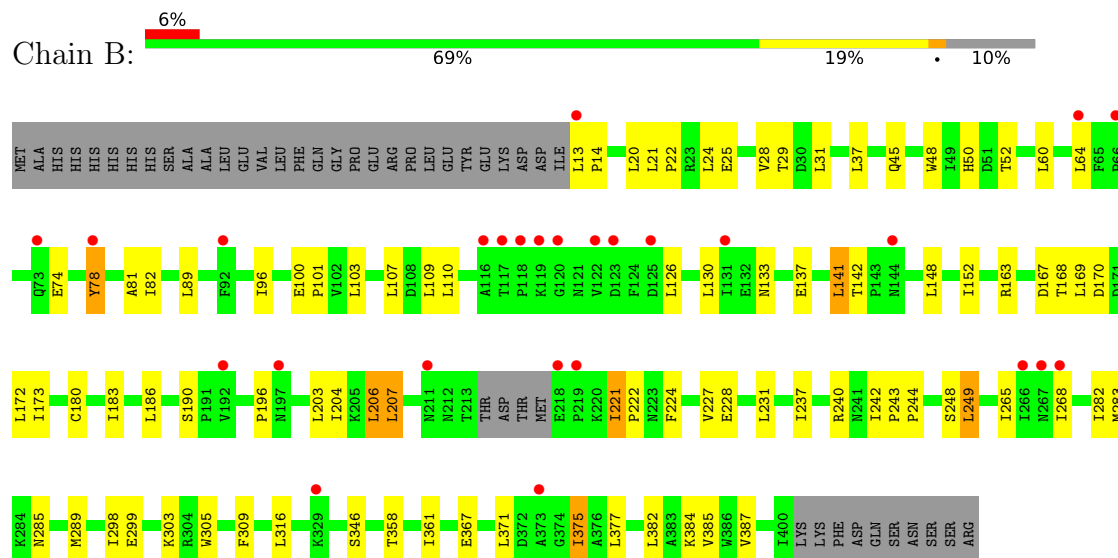
3 Residue-property plots

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: Kar9



• Molecule 1: Kar9



4 Data and refinement statistics

Property	Value	Source
Space group	P 41	Depositor
Cell constants a, b, c, α , β , γ	125.62Å 125.62Å 165.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.86 – 3.30 46.86 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.86-3.30) 99.9 (46.86-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 3.32Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.264 , 0.289 0.262 , 0.286	Depositor DCC
R_{free} test set	2574 reflections (6.68%)	wwPDB-VP
Wilson B-factor (Å ²)	141.5	Xtriage
Anisotropy	0.071	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 125.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.068 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5948	wwPDB-VP
Average B, all atoms (Å ²)	160.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.27% 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	1.01	0/2937	1.66	4/3984 (0.1%)
1	B	1.00	0/3104	1.62	0/4212
All	All	1.00	0/6041	1.64	4/8196 (0.0%)

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	91	LYS	CA-C-N	5.34	127.70	120.65
1	A	91	LYS	C-N-CA	5.34	127.70	120.65
1	A	223	ASN	CA-C-N	5.12	127.10	120.44
1	A	223	ASN	C-N-CA	5.12	127.10	120.44

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	2894	0	2780	64	0
1	B	3054	0	3016	55	0
All	All	5948	0	5796	117	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 10.

All (117) 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:65:PHE:O	1:A:78:TYR:OH	1.67	1.13
1:A:93:ILE:HG12	1:A:145:LEU:HG	1.42	1.01
1:A:225:SER:OG	1:A:228:GLU:HB2	1.84	0.77
1:A:93:ILE:HG23	1:A:145:LEU:HD23	1.67	0.74
1:A:78:TYR:CE1	1:A:82:ILE:HD11	2.24	0.72
1:A:60:LEU:HD12	1:A:144:ASN:ND2	2.04	0.72
1:A:96:ILE:HD11	1:A:145:LEU:HD21	1.74	0.70
1:A:51:ASP:O	1:A:55:ILE:HG13	1.92	0.69
1:B:169:LEU:HD23	1:B:249:LEU:CD1	2.23	0.69
1:A:139:ASN:O	1:A:143:PRO:HD2	1.93	0.67
1:B:107:LEU:HG	1:B:130:LEU:HD23	1.75	0.67
1:A:172:LEU:HG	1:A:245:ILE:HD11	1.77	0.66
1:B:309:PHE:CD2	1:B:371:LEU:HD11	2.31	0.65
1:B:309:PHE:CG	1:B:371:LEU:HD11	2.31	0.65
1:A:65:PHE:O	1:A:78:TYR:CZ	2.51	0.64
1:B:384:LYS:O	1:B:387:VAL:HG22	1.99	0.63
1:B:316:LEU:HD13	1:B:358:THR:HG21	1.81	0.63
1:A:139:ASN:O	1:A:143:PRO:CD	2.48	0.62
1:A:60:LEU:HD12	1:A:144:ASN:HD22	1.62	0.62
1:A:75:MET:SD	1:A:267:ASN:HB3	2.40	0.61
1:A:384:LYS:O	1:A:387:VAL:HG22	2.01	0.61
1:A:60:LEU:C	1:A:60:LEU:HD23	2.25	0.61
1:B:190:SER:OG	1:B:227:VAL:HG11	2.02	0.59
1:A:93:ILE:HA	1:A:145:LEU:HD21	1.86	0.57
1:B:221:ILE:HD12	1:B:221:ILE:H	1.70	0.57
1:A:78:TYR:CZ	1:A:82:ILE:HD11	2.40	0.56
1:A:107:LEU:HG	1:A:130:LEU:HD13	1.86	0.56
1:B:228:GLU:N	1:B:228:GLU:OE1	2.38	0.56
1:A:96:ILE:CD1	1:A:145:LEU:HD21	2.36	0.56
1:B:221:ILE:HD12	1:B:221:ILE:N	2.21	0.56
1:A:372:ASP:OD1	1:A:373:ALA:N	2.39	0.55
1:A:69:PHE:HA	1:A:72:TYR:CD1	2.42	0.55
1:B:96:ILE:HG22	1:B:141:LEU:HD23	1.89	0.55
1:B:24:LEU:O	1:B:28:VAL:HG23	2.08	0.54
1:A:48:TRP:O	1:A:52:THR:HG23	2.08	0.54
1:B:13:LEU:HB3	1:B:64:LEU:HD13	1.89	0.54
1:B:21:LEU:N	1:B:22:PRO:HD2	2.23	0.54
1:A:134:CYS:SG	1:B:240:ARG:NH2	2.81	0.53
1:A:142:THR:HB	1:A:143:PRO:HD3	1.88	0.53
1:B:169:LEU:HD23	1:B:249:LEU:HD13	1.91	0.53
1:B:163:ARG:O	1:B:167:ASP:HB2	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:74:GLU:OE1	1:A:74:GLU:HA	2.10	0.52
1:B:173:ILE:HG21	1:B:285:ASN:OD1	2.10	0.52
1:B:25:GLU:O	1:B:29:THR:HG23	2.10	0.51
1:B:382:LEU:O	1:B:385:VAL:HG22	2.10	0.51
1:A:88:GLU:HA	1:A:88:GLU:OE1	2.10	0.51
1:A:234:LYS:O	1:A:237:ILE:HG13	2.11	0.51
1:A:316:LEU:HD13	1:A:358:THR:HB	1.94	0.50
1:A:107:LEU:HG	1:A:130:LEU:CD1	2.41	0.50
1:B:173:ILE:CG2	1:B:285:ASN:OD1	2.60	0.49
1:B:20:LEU:C	1:B:22:PRO:HD2	2.38	0.49
1:B:227:VAL:HG12	1:B:231:LEU:HD13	1.95	0.49
1:A:60:LEU:HD23	1:A:60:LEU:O	2.13	0.48
1:B:48:TRP:CZ2	1:B:52:THR:HG21	2.49	0.48
1:B:100:GLU:HB3	1:B:101:PRO:HD3	1.96	0.48
1:A:93:ILE:HG12	1:A:145:LEU:CG	2.29	0.48
1:B:371:LEU:HD22	1:B:375:ILE:HD11	1.95	0.48
1:A:316:LEU:HD13	1:A:358:THR:CG2	2.43	0.48
1:B:82:ILE:HG21	1:B:152:ILE:HG12	1.94	0.47
1:B:240:ARG:O	1:B:243:PRO:HD2	2.14	0.47
1:B:299:GLU:O	1:B:303:LYS:HB3	2.14	0.47
1:B:13:LEU:HB2	1:B:14:PRO:HD3	1.95	0.47
1:A:372:ASP:OD1	1:A:374:GLY:N	2.36	0.47
1:B:31:LEU:HD21	1:B:45:GLN:HE21	1.78	0.47
1:A:382:LEU:O	1:A:385:VAL:HG22	2.15	0.47
1:B:249:LEU:HD23	1:B:283:MET:SD	2.55	0.47
1:A:75:MET:HE1	1:A:270:LEU:HB3	1.97	0.47
1:A:307:ILE:HG13	1:A:308:LEU:N	2.30	0.47
1:B:173:ILE:HG23	1:B:289:MET:HE3	1.96	0.46
1:A:316:LEU:HD13	1:A:358:THR:HG21	1.98	0.46
1:B:50:HIS:HB2	1:B:133:ASN:OD1	2.15	0.46
1:A:335:ASP:OD1	1:A:335:ASP:C	2.58	0.46
1:B:203:LEU:O	1:B:207:LEU:HB2	2.16	0.46
1:B:170:ASP:N	1:B:282:ILE:HD11	2.31	0.46
1:B:242:ILE:CG1	1:B:243:PRO:HD3	2.46	0.46
1:B:110:LEU:HD21	1:B:126:LEU:HD12	1.97	0.46
1:A:372:ASP:OD1	1:A:372:ASP:C	2.59	0.45
1:B:31:LEU:CD2	1:B:45:GLN:HE21	2.29	0.45
1:A:65:PHE:HB2	1:A:66:PRO:HD3	1.98	0.45
1:A:96:ILE:HD11	1:A:145:LEU:HD11	1.99	0.45
1:A:173:ILE:HD13	1:A:286:PHE:HA	1.99	0.45
1:B:316:LEU:HB3	1:B:382:LEU:HD23	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:222:PRO:HG3	1:B:305:TRP:CE2	2.52	0.45
1:A:148:LEU:O	1:A:152:ILE:HG13	2.18	0.44
1:A:327:LEU:HD21	1:A:393:SER:HA	1.99	0.44
1:A:75:MET:HB3	1:A:79:MET:HE2	2.00	0.43
1:B:78:TYR:O	1:B:82:ILE:HG13	2.18	0.43
1:A:89:LEU:N	1:A:89:LEU:HD23	2.32	0.43
1:B:305:TRP:CE2	1:B:367:GLU:HG2	2.53	0.43
1:A:137:GLU:O	1:A:141:LEU:CB	2.65	0.43
1:B:243:PRO:HB2	1:B:244:PRO:HD3	2.00	0.43
1:B:204:ILE:HG23	1:B:361:ILE:CD1	2.49	0.43
1:B:21:LEU:C	1:B:21:LEU:HD23	2.43	0.43
1:B:78:TYR:O	1:B:81:ALA:HB3	2.19	0.42
1:A:163:ARG:O	1:A:167:ASP:HB2	2.19	0.42
1:B:196:PRO:HG3	1:B:224:PHE:HD2	1.83	0.42
1:B:206:LEU:C	1:B:206:LEU:HD12	2.44	0.42
1:A:96:ILE:HG21	1:A:140:PHE:HE2	1.83	0.42
1:A:373:ALA:O	1:A:377:LEU:HD12	2.20	0.42
1:B:137:GLU:O	1:B:141:LEU:HB2	2.19	0.42
1:A:205:LYS:O	1:A:208:SER:N	2.53	0.42
1:A:222:PRO:CG	1:A:232:SER:HB2	2.49	0.42
1:B:107:LEU:CG	1:B:130:LEU:HD23	2.46	0.42
1:A:366:LEU:HD23	1:A:366:LEU:HA	1.87	0.41
1:B:169:LEU:HD21	1:B:248:SER:HB3	2.02	0.41
1:A:137:GLU:O	1:A:141:LEU:N	2.35	0.41
1:A:96:ILE:HD12	1:A:140:PHE:CE2	2.55	0.41
1:A:390:ARG:N	1:A:391:PRO:CD	2.84	0.41
1:B:37:LEU:HD11	1:B:109:LEU:C	2.46	0.41
1:A:21:LEU:N	1:A:22:PRO:CD	2.84	0.41
1:B:204:ILE:HD11	1:B:358:THR:OG1	2.21	0.41
1:A:82:ILE:CG2	1:A:152:ILE:HG12	2.51	0.40
1:A:48:TRP:CH2	1:A:52:THR:HG21	2.57	0.40
1:A:95:CYS:SG	1:A:96:ILE:N	2.94	0.40
1:A:139:ASN:O	1:A:143:PRO:HG2	2.21	0.40
1:A:107:LEU:HD13	1:B:240:ARG:HD3	2.04	0.40
1:A:205:LYS:HD2	1:A:357:LYS:HE2	2.03	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	369/428 (86%)	350 (95%)	19 (5%)	0	100	100
1	B	380/428 (89%)	362 (95%)	18 (5%)	0	100	100
All	All	749/856 (88%)	712 (95%)	37 (5%)	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	298/406 (73%)	287 (96%)	11 (4%)	30	58
1	B	335/406 (82%)	311 (93%)	24 (7%)	13	40
All	All	633/812 (78%)	598 (94%)	35 (6%)	19	48

All (35) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	42	ILE
1	A	89	LEU
1	A	123	ASP
1	A	165	HIS
1	A	168	THR
1	A	172	LEU
1	A	265	ILE
1	A	270	LEU

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Mol	Chain	Res	Type
1	A	310	ILE
1	A	326	LEU
1	A	392	LYS
1	B	60	LEU
1	B	74	GLU
1	B	78	TYR
1	B	89	LEU
1	B	103	LEU
1	B	141	LEU
1	B	142	THR
1	B	148	LEU
1	B	168	THR
1	B	172	LEU
1	B	180	CYS
1	B	183	ILE
1	B	186	LEU
1	B	206	LEU
1	B	207	LEU
1	B	221	ILE
1	B	237	ILE
1	B	249	LEU
1	B	265	ILE
1	B	268	ILE
1	B	298	ILE
1	B	346	SER
1	B	375	ILE
1	B	377	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	159	ASN
1	A	202	GLN
1	A	331	ASN
1	A	338	GLN
1	B	45	GLN
1	B	136	HIS
1	B	176	ASN
1	B	241	ASN
1	B	333	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	375/428 (87%)	0.64	37 (9%) 13 10	91, 160, 228, 283	0
1	B	384/428 (89%)	0.37	26 (6%) 23 16	100, 159, 213, 294	0
All	All	759/856 (88%)	0.50	63 (8%) 17 13	91, 159, 225, 294	0

All (63) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	117	THR	6.8
1	A	38	ASP	5.5
1	A	15	MET	5.3
1	B	13	LEU	5.0
1	A	13	LEU	4.7
1	A	114	LEU	4.7
1	A	17	LEU	4.3
1	A	35	HIS	4.2
1	A	265	ILE	4.1
1	A	92	PHE	3.9
1	B	268	ILE	3.8
1	B	73	GLN	3.8
1	A	221	ILE	3.8
1	A	190	SER	3.7
1	B	266	ILE	3.7
1	A	220	LYS	3.6
1	A	201	ASP	3.6
1	B	131	ILE	3.5
1	B	197	ASN	3.5
1	A	224	PHE	3.5
1	A	144	ASN	3.5
1	A	39	VAL	3.5
1	B	267	ASN	3.3
1	B	119	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	125	ASP	3.2
1	A	140	PHE	3.1
1	B	64	LEU	3.1
1	A	222	PRO	2.9
1	A	61	ALA	2.8
1	A	12	ILE	2.8
1	B	218	GLU	2.8
1	A	203	LEU	2.7
1	A	131	ILE	2.7
1	A	223	ASN	2.7
1	B	144	ASN	2.7
1	A	122	VAL	2.7
1	A	18	GLN	2.6
1	B	373	ALA	2.6
1	B	211	ASN	2.6
1	B	116	ALA	2.6
1	A	14	PRO	2.6
1	A	33	LEU	2.5
1	A	226	PRO	2.5
1	B	118	PRO	2.4
1	B	123	ASP	2.4
1	B	122	VAL	2.4
1	B	120	GLY	2.3
1	A	232	SER	2.3
1	A	228	GLU	2.3
1	A	272	ALA	2.3
1	B	92	PHE	2.3
1	A	94	ASN	2.3
1	B	192	VAL	2.3
1	A	143	PRO	2.2
1	B	66	PRO	2.2
1	A	151	CYS	2.2
1	B	219	PRO	2.1
1	A	404	ASP	2.1
1	A	206	LEU	2.1
1	A	205	LYS	2.1
1	B	78	TYR	2.0
1	A	65	PHE	2.0
1	B	329	LYS	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.