



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 25, 2026 – 01:32 AM EDT

PDB ID : 1R8G / pdb_00001r8g
Title : Structure and function of YbdK
Authors : Lehmann, C.; Doseeva, V.; Pullalarevu, S.; Krajewski, W.; Howard, A.; Herzberg, O.; Structure 2 Function Project (S2F)
Deposited on : 2003-10-24
Resolution : 2.15 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

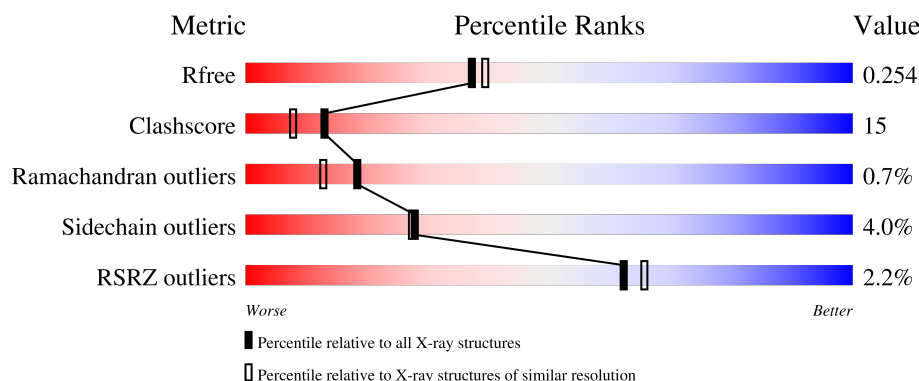
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

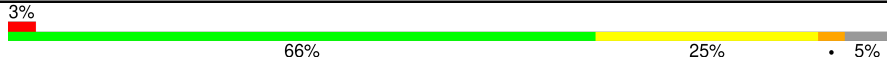
The reported resolution of this entry is 2.15 Å.

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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	372	
1	B	372	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6234 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 Hypothetical protein ybdK.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	352	Total	C	N	O	S	Se	0	0	0
			2763	1758	480	509	6	10			
1	B	358	Total	C	N	O	S	Se	0	0	0
			2815	1790	491	518	6	10			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	modified residue	UNP P77213
A	20	MSE	MET	modified residue	UNP P77213
A	36	MSE	MET	modified residue	UNP P77213
A	58	MSE	MET	modified residue	UNP P77213
A	79	MSE	MET	modified residue	UNP P77213
A	171	MSE	MET	modified residue	UNP P77213
A	195	MSE	MET	modified residue	UNP P77213
A	216	MSE	MET	modified residue	UNP P77213
A	241	MSE	MET	modified residue	UNP P77213
A	253	MSE	MET	modified residue	UNP P77213
A	348	MSE	MET	modified residue	UNP P77213
B	1	MSE	MET	modified residue	UNP P77213
B	20	MSE	MET	modified residue	UNP P77213
B	36	MSE	MET	modified residue	UNP P77213
B	58	MSE	MET	modified residue	UNP P77213
B	79	MSE	MET	modified residue	UNP P77213
B	171	MSE	MET	modified residue	UNP P77213
B	195	MSE	MET	modified residue	UNP P77213
B	216	MSE	MET	modified residue	UNP P77213
B	241	MSE	MET	modified residue	UNP P77213
B	253	MSE	MET	modified residue	UNP P77213
B	348	MSE	MET	modified residue	UNP P77213

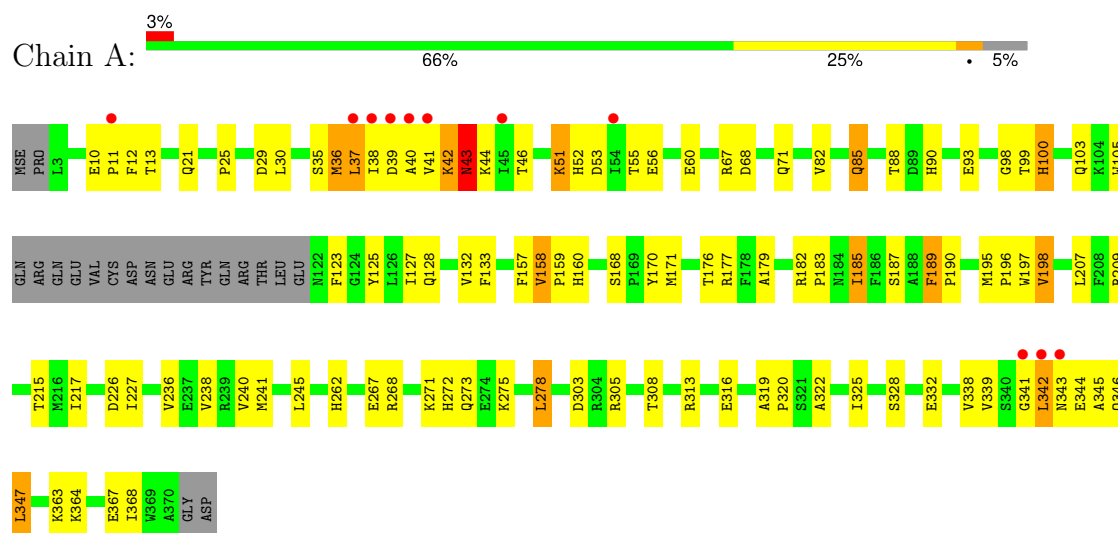
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	324	Total 324	O 324	0	0
2	B	332	Total 332	O 332	0	0

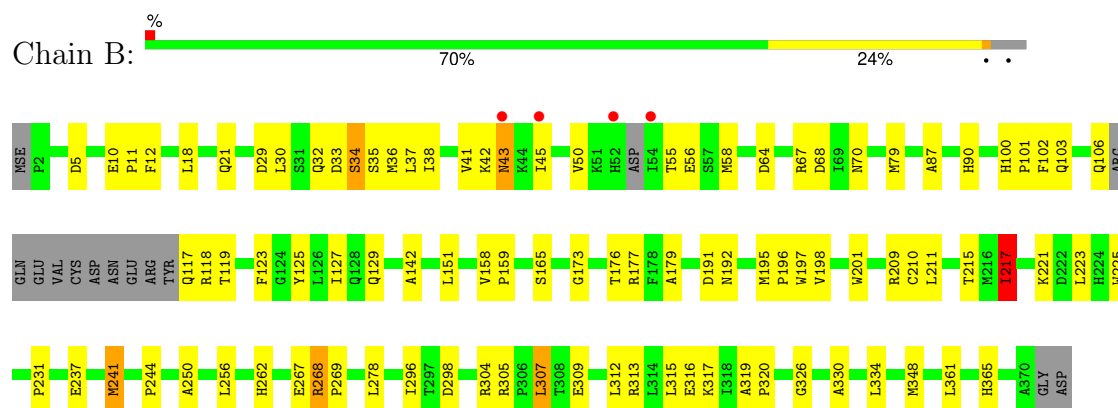
3 Residue-property plots [i](#)

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: Hypothetical protein ybdK



• Molecule 1: Hypothetical protein ybdK



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	49.47Å 100.32Å 71.37Å 90.00° 99.94° 90.00°	Depositor
Resolution (Å)	20.00 – 2.15 20.00 – 2.15	Depositor EDS
% Data completeness (in resolution range)	83.6 (20.00-2.15) 86.6 (20.00-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.65 (at 2.16Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.189 , 0.252 0.199 , 0.254	Depositor DCC
R_{free} test set	1796 reflections (2.89%)	wwPDB-VP
Wilson B-factor (Å ²)	16.8	Xtriage
Anisotropy	0.285	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 53.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6234	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.62% 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	0.51	0/2825	0.99	9/3821 (0.2%)
1	B	0.53	0/2877	1.04	18/3889 (0.5%)
All	All	0.52	0/5702	1.02	27/7710 (0.4%)

There are no bond length outliers.

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	100	HIS	N-CA-C	-9.49	98.16	109.93
1	B	100	HIS	N-CA-C	-8.30	99.34	109.72
1	B	215	THR	N-CA-C	7.99	119.98	111.28
1	A	241	MSE	N-CA-C	7.57	122.35	110.32
1	B	10	GLU	N-CA-C	-6.72	99.48	109.42
1	A	236	VAL	N-CA-C	-6.32	99.26	108.11
1	A	271	LYS	N-CA-C	-6.26	96.90	107.99
1	B	268	ARG	CA-C-N	6.13	125.75	119.56
1	B	268	ARG	C-N-CA	6.13	125.75	119.56
1	B	179	ALA	N-CA-C	-6.09	104.19	111.69
1	B	241	MSE	N-CA-C	6.06	119.45	110.48
1	A	187	SER	N-CA-C	5.95	120.07	112.34
1	B	217	ILE	N-CA-C	5.89	116.49	107.77
1	B	101	PRO	N-CA-C	5.84	122.46	113.75
1	A	179	ALA	N-CA-C	-5.80	104.55	111.69
1	B	237	GLU	N-CA-C	5.77	118.21	107.99
1	B	298	ASP	CA-C-N	5.59	125.21	119.56
1	B	298	ASP	C-N-CA	5.59	125.21	119.56
1	B	64	ASP	N-CA-C	-5.53	101.30	109.86
1	B	123	PHE	N-CA-C	5.51	120.06	113.23
1	B	5	ASP	N-CA-C	-5.44	102.83	110.50
1	B	197	TRP	N-CA-C	5.43	118.10	109.96
1	B	142	ALA	N-CA-C	5.35	118.97	112.23
1	B	58	MSE	N-CA-C	5.17	117.82	109.40
1	A	189	PHE	CA-C-N	5.14	125.17	119.32

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	189	PHE	C-N-CA	5.14	125.17	119.32
1	A	198	VAL	N-CA-C	-5.08	101.16	108.42

There are no chirality outliers.

There are no planarity outliers.

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	2763	0	2691	90	0
1	B	2815	0	2747	80	0
2	A	324	0	0	13	0
2	B	332	0	0	10	0
All	All	6234	0	5438	165	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 15.

All (165) 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:38:ILE:HD11	1:A:52:HIS:HB3	1.47	0.96
1:B:117:GLN:HE21	1:B:118:ARG:HH21	1.10	0.93
1:B:348:MSE:HE3	1:B:365:HIS:HE1	1.33	0.91
1:B:348:MSE:HE2	1:B:361:LEU:HD21	1.56	0.88
2:A:480:HOH:O	1:B:210:CYS:HB2	1.74	0.86
1:A:29:ASP:HB2	1:A:103:GLN:OE1	1.79	0.82
1:A:275:LYS:HG3	1:A:278:LEU:HD12	1.60	0.81
1:B:312:LEU:O	1:B:316:GLU:HG3	1.81	0.80
1:B:37:LEU:HD11	1:B:87:ALA:HA	1.62	0.80
1:B:117:GLN:HG3	1:B:118:ARG:CZ	2.11	0.79
1:B:348:MSE:HE3	1:B:365:HIS:CE1	2.16	0.79
1:A:305:ARG:HG2	1:A:313:ARG:HH12	1.46	0.78
1:B:117:GLN:CG	1:B:118:ARG:H	1.96	0.78
1:A:44:LYS:O	1:A:82:VAL:HG21	1.84	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:ASP:OD2	1:A:71:GLN:HG3	1.86	0.75
1:A:11:PRO:HG2	1:A:67:ARG:HH22	1.53	0.74
1:A:53:ASP:OD1	1:A:55:THR:HB	1.87	0.73
1:B:42:LYS:O	1:B:43:ASN:HB2	1.90	0.71
1:A:41:VAL:HG12	1:A:82:VAL:HG12	1.71	0.70
1:A:36:MSE:HG3	1:A:90:HIS:ND1	2.07	0.70
1:A:39:ASP:HA	1:A:42:LYS:HD2	1.73	0.69
1:A:217:ILE:O	2:A:399:HOH:O	2.08	0.69
1:B:117:GLN:HG2	1:B:118:ARG:H	1.57	0.69
1:B:117:GLN:HG3	1:B:118:ARG:NH2	2.07	0.69
1:B:117:GLN:NE2	1:B:118:ARG:HH21	1.90	0.67
1:B:33:ASP:C	1:B:35:SER:H	2.04	0.66
1:B:195:MSE:HE3	1:B:198:VAL:HG22	1.78	0.66
1:A:51:LYS:HB3	1:A:51:LYS:NZ	2.11	0.65
1:A:303:ASP:OD2	1:A:305:ARG:HD2	1.95	0.65
1:A:363:LYS:O	1:A:367:GLU:HG3	1.97	0.64
1:B:117:GLN:HG3	1:B:118:ARG:NE	2.11	0.64
1:A:55:THR:O	1:A:56:GLU:HB2	1.97	0.64
1:B:309:GLU:HG3	1:B:313:ARG:HH11	1.63	0.64
1:B:117:GLN:HG2	1:B:118:ARG:N	2.14	0.63
1:B:117:GLN:CG	1:B:118:ARG:N	2.60	0.62
1:A:195:MSE:HE3	1:A:198:VAL:HG22	1.81	0.62
1:B:241:MSE:HE3	2:B:408:HOH:O	2.01	0.61
1:A:38:ILE:HD11	1:A:52:HIS:CB	2.25	0.61
1:B:176:THR:O	1:B:177:ARG:HB2	2.02	0.59
1:A:21:GLN:HG2	1:A:133:PHE:CE2	2.39	0.57
1:B:313:ARG:O	1:B:317:LYS:HE2	2.05	0.57
1:A:215:THR:HG23	1:B:304:ARG:CZ	2.35	0.56
1:B:11:PRO:HG2	2:B:548:HOH:O	2.04	0.56
1:B:30:LEU:HD12	1:B:103:GLN:HG2	1.87	0.56
1:A:227:ILE:HD12	1:A:238:VAL:HG22	1.87	0.56
1:A:10:GLU:HB3	1:A:11:PRO:HD2	1.87	0.56
1:A:319:ALA:HB3	1:A:320:PRO:HD3	1.88	0.55
1:A:99:THR:HG21	1:A:171:MSE:HE3	1.88	0.55
1:B:217:ILE:HG13	1:B:217:ILE:O	2.02	0.55
1:B:348:MSE:CE	1:B:361:LEU:HD21	2.34	0.55
1:A:176:THR:O	1:A:177:ARG:HB2	2.07	0.55
1:B:33:ASP:C	1:B:35:SER:N	2.64	0.55
1:B:165:SER:CB	1:B:241:MSE:HG2	2.37	0.54
1:A:39:ASP:CA	1:A:42:LYS:HD2	2.38	0.54
1:A:343:ASN:HD22	1:A:346:GLN:HB2	1.72	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:364:LYS:HE3	1:A:368:ILE:HD11	1.90	0.53
1:B:151:LEU:HD23	1:B:151:LEU:C	2.33	0.53
1:A:38:ILE:C	1:A:40:ALA:H	2.17	0.53
1:A:227:ILE:CD1	1:A:238:VAL:HG13	2.38	0.53
1:A:10:GLU:OE1	1:A:67:ARG:NH1	2.42	0.53
1:B:117:GLN:HE21	1:B:118:ARG:NH2	1.93	0.52
1:A:226:ASP:O	1:A:227:ILE:HD13	2.09	0.52
1:A:51:LYS:HG2	2:A:429:HOH:O	2.09	0.52
1:A:339:VAL:HG13	2:A:679:HOH:O	2.10	0.52
1:B:348:MSE:HE1	2:B:447:HOH:O	2.10	0.51
1:A:39:ASP:HA	1:A:42:LYS:CD	2.40	0.51
1:A:209:ARG:HH11	1:A:209:ARG:HG2	1.75	0.51
1:B:296:ILE:HB	1:B:307:LEU:CD1	2.40	0.51
1:A:125:TYR:CD1	1:B:176:THR:HB	2.45	0.51
1:B:223:LEU:HD13	1:B:225:TRP:CH2	2.46	0.51
1:A:215:THR:HG23	1:B:304:ARG:NH2	2.26	0.51
1:A:98:GLY:HA3	1:A:168:SER:HB3	1.93	0.51
1:B:12:PHE:CE1	1:B:262:HIS:HA	2.46	0.51
1:A:305:ARG:HG2	1:A:313:ARG:NH1	2.20	0.50
1:A:105:TRP:CZ2	1:A:128:GLN:HG2	2.46	0.50
1:A:347:LEU:HD13	2:A:690:HOH:O	2.11	0.50
1:A:328:SER:O	1:A:332:GLU:HG3	2.11	0.50
1:B:348:MSE:HE2	1:B:361:LEU:CD2	2.37	0.50
1:A:343:ASN:ND2	1:A:346:GLN:H	2.09	0.50
1:A:39:ASP:HA	1:A:42:LYS:CE	2.42	0.50
1:B:35:SER:HB3	1:B:56:GLU:OE1	2.12	0.50
1:B:117:GLN:C	1:B:119:THR:H	2.18	0.49
1:B:33:ASP:O	1:B:35:SER:N	2.45	0.49
1:B:36:MSE:HE3	1:B:90:HIS:HB3	1.94	0.49
1:A:267:GLU:O	1:A:268:ARG:C	2.55	0.49
1:A:195:MSE:HG2	1:A:196:PRO:O	2.13	0.48
1:B:244:PRO:HG3	2:B:436:HOH:O	2.12	0.48
1:A:41:VAL:C	1:A:43:ASN:H	2.22	0.48
1:B:319:ALA:HB3	1:B:320:PRO:HD3	1.96	0.48
1:A:41:VAL:HG12	1:A:82:VAL:CG1	2.42	0.48
1:A:12:PHE:CE1	1:A:262:HIS:HA	2.49	0.47
1:B:45:ILE:HD11	1:B:50:VAL:HB	1.96	0.47
1:A:273:GLN:NE2	2:A:495:HOH:O	2.46	0.47
1:B:117:GLN:HG3	1:B:118:ARG:H	1.76	0.47
1:B:158:VAL:N	1:B:159:PRO:CD	2.77	0.47
1:A:341:GLY:O	1:A:342:LEU:C	2.58	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:165:SER:OG	1:B:241:MSE:HG2	2.14	0.47
1:B:296:ILE:HG22	1:B:305:ARG:O	2.14	0.47
1:A:39:ASP:C	1:A:42:LYS:HD2	2.40	0.47
1:B:165:SER:HB2	1:B:241:MSE:HG2	1.96	0.47
1:B:119:THR:HG22	1:B:127:ILE:HD13	1.95	0.47
1:A:21:GLN:HG2	1:A:133:PHE:CZ	2.49	0.47
1:A:185:ILE:HD12	2:B:466:HOH:O	2.15	0.46
1:A:308:THR:HG23	1:A:338:VAL:CG1	2.45	0.46
1:A:10:GLU:HB2	1:A:13:THR:CG2	2.45	0.46
1:A:245:LEU:HD12	1:A:344:GLU:OE1	2.15	0.46
1:B:209:ARG:HE	1:B:209:ARG:CA	2.29	0.46
1:B:18:LEU:HD21	1:B:250:ALA:HB1	1.99	0.45
1:B:326:GLY:HA2	2:B:416:HOH:O	2.16	0.45
1:A:207:LEU:CD1	1:B:278:LEU:HD11	2.45	0.45
1:A:182:ARG:HB3	1:A:183:PRO:HD3	1.98	0.45
1:A:51:LYS:HB3	1:A:51:LYS:HZ2	1.81	0.45
1:A:132:VAL:HB	1:A:182:ARG:HG3	1.98	0.45
1:A:308:THR:HG23	1:A:338:VAL:HG12	1.98	0.45
1:B:34:SER:O	1:B:38:ILE:HG13	2.17	0.45
1:A:123:PHE:CD2	1:A:189:PHE:HE1	2.36	0.44
1:B:267:GLU:O	1:B:268:ARG:C	2.59	0.44
1:A:268:ARG:NH1	2:A:389:HOH:O	2.48	0.44
1:A:35:SER:HB3	1:A:56:GLU:OE2	2.18	0.44
1:A:278:LEU:HD13	1:B:210:CYS:SG	2.58	0.44
1:A:25:PRO:CG	1:A:93:GLU:HG3	2.48	0.44
1:A:100:HIS:CG	1:A:103:GLN:HG2	2.53	0.44
1:A:322:ALA:HB3	2:A:687:HOH:O	2.17	0.44
1:B:201:TRP:CZ2	1:B:231:PRO:HG3	2.52	0.44
1:A:170:TYR:HE1	2:A:555:HOH:O	2.01	0.43
1:B:256:LEU:HB2	1:B:334:LEU:CD1	2.48	0.43
1:B:209:ARG:HE	1:B:209:ARG:C	2.25	0.43
1:A:55:THR:O	1:A:56:GLU:CB	2.64	0.43
1:A:37:LEU:O	1:A:40:ALA:HB3	2.19	0.43
1:B:67:ARG:HG3	1:B:68:ASP:OD1	2.19	0.43
1:B:191:ASP:OD1	1:B:223:LEU:HA	2.19	0.43
1:A:51:LYS:NZ	1:A:51:LYS:CB	2.80	0.43
1:B:118:ARG:N	1:B:118:ARG:CD	2.82	0.43
1:B:268:ARG:N	1:B:269:PRO:CD	2.82	0.42
1:A:182:ARG:HD3	2:A:385:HOH:O	2.19	0.42
1:B:125:TYR:HB3	2:B:624:HOH:O	2.18	0.42
1:A:189:PHE:HA	1:A:190:PRO:HD3	1.86	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:217:ILE:O	1:B:217:ILE:CG1	2.65	0.42
1:B:221:LYS:HD3	2:B:532:HOH:O	2.19	0.42
1:A:38:ILE:CD1	1:A:52:HIS:HB3	2.33	0.42
1:A:51:LYS:HB3	1:A:51:LYS:HZ3	1.81	0.42
1:A:226:ASP:HB3	1:A:240:VAL:CG1	2.50	0.42
1:B:201:TRP:CH2	1:B:231:PRO:HG3	2.55	0.42
1:A:30:LEU:HG	1:A:103:GLN:HG3	2.02	0.41
1:B:103:GLN:NE2	2:B:627:HOH:O	2.53	0.41
1:B:195:MSE:HG2	1:B:196:PRO:O	2.20	0.41
1:B:195:MSE:CE	1:B:198:VAL:HG22	2.48	0.41
1:A:25:PRO:HG2	2:A:475:HOH:O	2.20	0.41
1:A:197:TRP:HA	2:A:471:HOH:O	2.20	0.41
1:B:102:PHE:HA	1:B:173:GLY:N	2.35	0.41
1:B:45:ILE:HG23	1:B:79:MSE:HG2	2.03	0.41
1:B:118:ARG:N	1:B:118:ARG:HD3	2.35	0.41
1:A:30:LEU:CG	1:A:103:GLN:HG3	2.51	0.41
1:A:157:PHE:HA	1:A:160:HIS:ND1	2.36	0.41
1:B:330:ALA:O	1:B:334:LEU:HG	2.21	0.41
1:B:348:MSE:CE	1:B:365:HIS:HE1	2.18	0.41
1:A:85:GLN:O	1:A:88:THR:OG1	2.40	0.40
1:A:272:HIS:HB3	2:A:536:HOH:O	2.20	0.40
1:A:325:ILE:H	1:A:325:ILE:HG13	1.77	0.40
1:B:38:ILE:O	1:B:41:VAL:HG22	2.20	0.40
1:A:158:VAL:N	1:A:159:PRO:CD	2.84	0.40
1:A:171:MSE:HE1	1:A:185:ILE:HD13	2.03	0.40
1:A:343:ASN:HD21	1:A:345:ALA:HB3	1.87	0.40
1:B:37:LEU:HD11	1:B:87:ALA:CA	2.41	0.40
1:B:21:GLN:HG3	2:B:455:HOH:O	2.21	0.40

There are no symmetry-related clashes.

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	348/372 (94%)	334 (96%)	11 (3%)	3 (1%)	14	9
1	B	352/372 (95%)	344 (98%)	6 (2%)	2 (1%)	21	15
All	All	700/744 (94%)	678 (97%)	17 (2%)	5 (1%)	18	13

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	B	34	SER
1	B	55	THR
1	A	42	LYS
1	A	342	LEU

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	298/306 (97%)	285 (96%)	13 (4%)	25	23
1	B	304/306 (99%)	293 (96%)	11 (4%)	31	31
All	All	602/612 (98%)	578 (96%)	24 (4%)	28	27

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

Mol	Chain	Res	Type
1	A	36	MSE
1	A	37	LEU
1	A	43	ASN
1	A	46	THR
1	A	51	LYS
1	A	60	GLU
1	A	85	GLN
1	A	127	ILE
1	A	158	VAL
1	A	185	ILE
1	A	278	LEU

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Mol	Chain	Res	Type
1	A	316	GLU
1	A	347	LEU
1	B	29	ASP
1	B	32	GLN
1	B	43	ASN
1	B	70	ASN
1	B	106	GLN
1	B	129	GLN
1	B	192	ASN
1	B	211	LEU
1	B	217	ILE
1	B	307	LEU
1	B	315	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	43	ASN
1	A	80	GLN
1	A	128	GLN
1	A	203	GLN
1	A	273	GLN
1	A	337	GLN
1	A	343	ASN
1	B	7	HIS
1	B	32	GLN
1	B	43	ASN
1	B	80	GLN
1	B	117	GLN
1	B	128	GLN
1	B	136	HIS
1	B	172	GLN
1	B	203	GLN
1	B	232	HIS

5.3.3 RNA ⓘ

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	342/372 (91%)	-0.10	11 (3%) 50 54	7, 19, 46, 64	0
1	B	348/372 (93%)	-0.22	4 (1%) 78 80	8, 19, 44, 62	0
All	All	690/744 (92%)	-0.16	15 (2%) 62 66	7, 19, 45, 64	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	343	ASN	3.3
1	A	342	LEU	3.2
1	A	11	PRO	3.2
1	A	41	VAL	3.0
1	A	38	ILE	2.8
1	A	40	ALA	2.6
1	B	54	ILE	2.6
1	A	39	ASP	2.6
1	A	37	LEU	2.5
1	B	43	ASN	2.4
1	A	45	ILE	2.1
1	B	52	HIS	2.1
1	A	54	ILE	2.1
1	A	341	GLY	2.0
1	B	45	ILE	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.