



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 06:53 PM UTC

PDB ID : 2R9A / pdb_00002r9a
Title : Crystal structure of human XLF
Authors : Andres, S.N.; Junop, M.S.
Deposited on : 2007-09-12
Resolution : 2.50 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

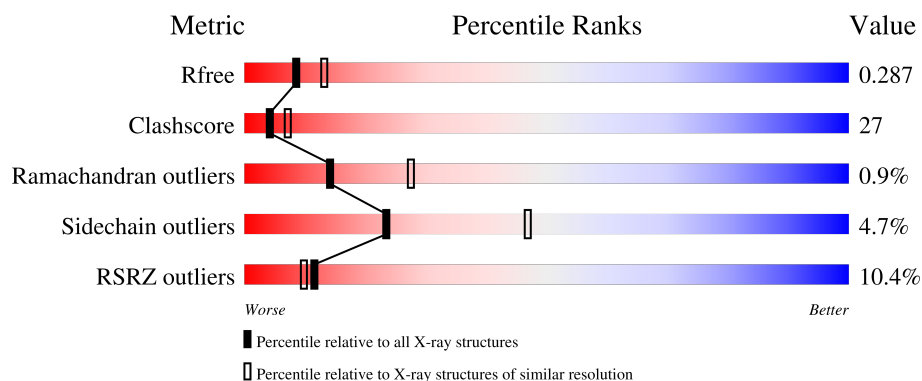
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	230	10% 53% 36% 6% 5%
1	B	230	9% 53% 36% • 7%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3744 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 Non-homologous end-joining factor 1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	219	Total	C	N	O	S	Se	0	0	0
			1752	1121	296	320	6	9			
1	B	214	Total	C	N	O	S	Se	0	0	0
			1710	1092	288	315	6	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	HIS	-	expression tag	UNP Q9H9Q4
A	226	HIS	-	expression tag	UNP Q9H9Q4
A	227	HIS	-	expression tag	UNP Q9H9Q4
A	228	HIS	-	expression tag	UNP Q9H9Q4
A	229	HIS	-	expression tag	UNP Q9H9Q4
A	230	HIS	-	expression tag	UNP Q9H9Q4
B	225	HIS	-	expression tag	UNP Q9H9Q4
B	226	HIS	-	expression tag	UNP Q9H9Q4
B	227	HIS	-	expression tag	UNP Q9H9Q4
B	228	HIS	-	expression tag	UNP Q9H9Q4
B	229	HIS	-	expression tag	UNP Q9H9Q4
B	230	HIS	-	expression tag	UNP Q9H9Q4

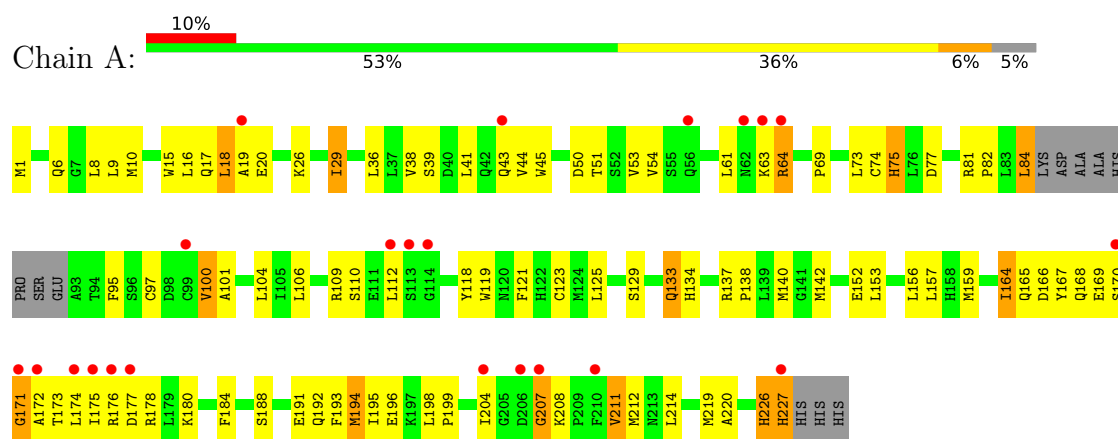
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	122	Total	O	0	0
			122	122		
2	B	160	Total	O	0	0
			160	160		

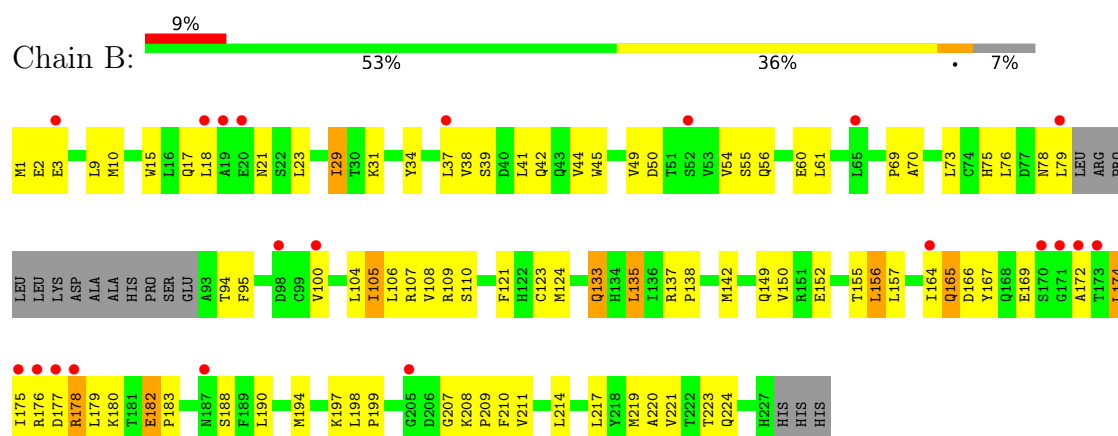
3 Residue-property plots [i](#)

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: Non-homologous end-joining factor 1



• Molecule 1: Non-homologous end-joining factor 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	63.46Å 86.95Å 91.91Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	25.00 – 2.50 25.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (25.00-2.50) 83.2 (25.00-2.50)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.78 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.251 , 0.287 0.253 , 0.287	Depositor DCC
R_{free} test set	1164 reflections (6.83%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.391	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.3	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	0.89	EDS
Total number of atoms	3744	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 5.27% 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

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	1.04	5/1781 (0.3%)	1.10	12/2401 (0.5%)
1	B	0.97	3/1738 (0.2%)	1.09	2/2342 (0.1%)
All	All	1.01	8/3519 (0.2%)	1.10	14/4743 (0.3%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	133	GLN	CD-OE1	6.91	1.36	1.23
1	B	133	GLN	CD-OE1	5.95	1.34	1.23
1	A	227	HIS	ND1-CE1	5.87	1.38	1.32
1	A	75	HIS	ND1-CE1	5.55	1.38	1.32
1	A	134	HIS	ND1-CE1	5.47	1.38	1.32
1	B	42	GLN	CD-OE1	5.23	1.33	1.23
1	A	226	HIS	ND1-CE1	5.19	1.37	1.32
1	B	165	GLN	CD-NE2	-5.00	1.22	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	182	GLU	CA-C-N	6.88	128.44	119.84
1	B	182	GLU	C-N-CA	6.88	128.44	119.84
1	A	226	HIS	CB-CG-CD2	-6.70	122.49	131.20
1	A	134	HIS	CB-CG-CD2	-6.60	122.63	131.20
1	A	75	HIS	CB-CG-CD2	-6.54	122.70	131.20
1	A	227	HIS	CB-CG-CD2	-6.43	122.85	131.20
1	A	195	ILE	N-CA-C	-6.00	107.04	111.90
1	A	194	MSE	CB-CG-SE	-5.77	95.38	112.70
1	A	75	HIS	CB-CG-ND1	5.77	131.35	122.70
1	A	227	HIS	CB-CG-ND1	5.73	131.30	122.70
1	A	134	HIS	CB-CG-ND1	5.69	131.23	122.70
1	A	226	HIS	CB-CG-ND1	5.64	131.16	122.70
1	A	29	ILE	CB-CA-C	-5.29	102.94	110.83
1	A	207	GLY	N-CA-C	-5.13	108.75	115.47

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	1752	0	1758	111	0
1	B	1710	0	1705	105	0
2	A	122	0	0	26	0
2	B	160	0	0	15	0
All	All	3744	0	3463	189	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 27.

All (189) 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:140:MSE:HE3	1:B:210:PHE:HB2	1.18	1.13
1:B:219:MSE:HG2	2:B:390:HOH:O	1.53	1.08
1:A:6:GLN:HG3	2:A:293:HOH:O	1.57	1.05
1:B:10:MSE:HE3	1:B:223:THR:HG22	1.36	1.04
1:A:194:MSE:HE2	1:B:217:LEU:HD13	1.39	1.02
1:A:140:MSE:HE3	1:B:210:PHE:CB	1.89	1.02
1:A:165:GLN:O	1:A:169:GLU:HG2	1.60	0.99
1:B:10:MSE:HE3	1:B:223:THR:CG2	1.94	0.97
1:A:164:ILE:HG22	1:B:164:ILE:HG12	1.46	0.96
1:A:173:THR:O	1:A:174:LEU:HD12	1.68	0.92
1:B:10:MSE:CE	1:B:223:THR:HG22	2.01	0.90
1:B:177:ASP:O	1:B:178:ARG:HG3	1.73	0.88
1:A:75:HIS:HD2	2:A:260:HOH:O	1.56	0.88
1:B:2:GLU:HG3	2:B:362:HOH:O	1.75	0.87
1:A:63:LYS:O	1:A:64:ARG:HG3	1.77	0.84
1:B:15:TRP:CZ3	1:B:17:GLN:HG2	2.14	0.82
1:B:165:GLN:O	1:B:169:GLU:HG2	1.81	0.80
1:B:15:TRP:HZ3	1:B:17:GLN:HG2	1.44	0.80
1:A:220:ALA:HB1	1:B:194:MSE:CE	2.13	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:MSE:HE3	2:A:280:HOH:O	1.83	0.78
1:A:152:GLU:OE2	1:B:197:LYS:HD2	1.85	0.76
1:B:219:MSE:HA	2:B:390:HOH:O	1.84	0.76
1:B:133:GLN:HG3	2:B:357:HOH:O	1.85	0.75
1:A:110:SER:HB2	2:A:253:HOH:O	1.85	0.75
1:A:164:ILE:HG22	1:B:164:ILE:CG1	2.16	0.74
1:A:75:HIS:CD2	2:A:260:HOH:O	2.36	0.73
1:B:78:ASN:O	1:B:79:LEU:O	2.06	0.73
1:B:18:LEU:HB2	1:B:21:ASN:O	1.90	0.72
1:B:37:LEU:HD12	1:B:45:TRP:O	1.90	0.72
1:B:37:LEU:HD11	1:B:44:VAL:HG12	1.69	0.72
2:A:310:HOH:O	1:B:197:LYS:HE2	1.88	0.72
1:A:165:GLN:O	1:A:169:GLU:CG	2.37	0.71
1:A:8:LEU:HD11	1:A:26:LYS:HG3	1.74	0.69
1:A:140:MSE:CE	1:B:210:PHE:HB2	2.10	0.69
1:B:106:LEU:HB3	1:B:121:PHE:HB2	1.73	0.69
1:B:107:ARG:NH1	2:B:332:HOH:O	2.25	0.68
1:A:51:THR:HG23	1:A:69:PRO:CG	2.23	0.68
1:A:219:MSE:HG2	2:A:325:HOH:O	1.93	0.67
1:B:177:ASP:O	1:B:178:ARG:CG	2.42	0.67
1:B:15:TRP:HB2	1:B:211:VAL:HG21	1.75	0.67
1:A:165:GLN:C	1:A:169:GLU:HG2	2.21	0.66
1:A:220:ALA:CB	1:B:194:MSE:CE	2.74	0.65
1:A:54:VAL:HG21	1:A:73:LEU:HD11	1.79	0.65
1:B:10:MSE:HG2	1:B:223:THR:HG22	1.78	0.65
1:A:178:ARG:HG2	2:A:342:HOH:O	1.98	0.64
1:B:79:LEU:HD11	1:B:95:PHE:HZ	1.63	0.64
1:A:43:GLN:OE1	1:A:125:LEU:HD21	1.98	0.64
1:A:170:SER:HB3	2:A:333:HOH:O	1.97	0.64
1:A:220:ALA:CB	1:B:194:MSE:HE1	2.28	0.63
1:A:220:ALA:HB1	1:B:194:MSE:HE1	1.79	0.63
1:A:194:MSE:C	2:A:331:HOH:O	2.41	0.62
1:B:37:LEU:HD11	1:B:44:VAL:CG1	2.28	0.62
1:B:1:MSE:HE1	1:B:50:ASP:HB3	1.82	0.62
1:A:41:LEU:HD13	1:A:204:ILE:HG22	1.81	0.62
1:A:173:THR:C	1:A:174:LEU:HD12	2.24	0.62
1:A:208:LYS:HG3	1:A:212:MSE:HE2	1.81	0.62
1:B:142:MSE:HE1	1:B:214:LEU:HB3	1.82	0.61
1:A:227:HIS:HE1	2:A:242:HOH:O	1.82	0.61
1:A:175:ILE:HG22	1:B:166:ASP:OD1	2.00	0.60
1:B:194:MSE:CE	1:B:198:LEU:HD22	2.32	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:61:LEU:HD12	1:A:119:TRP:HA	1.83	0.59
1:A:169:GLU:HG3	1:A:170:SER:N	2.17	0.59
1:A:226:HIS:CE1	2:A:233:HOH:O	2.55	0.59
1:A:106:LEU:HB3	1:A:121:PHE:HB2	1.84	0.58
1:A:194:MSE:HE2	1:B:217:LEU:CD1	2.24	0.58
1:A:129:SER:O	1:A:133:GLN:HG2	2.04	0.57
1:A:212:MSE:HE3	2:A:347:HOH:O	2.04	0.57
1:A:156:LEU:HA	1:A:159:MSE:HE3	1.85	0.57
2:A:310:HOH:O	1:B:197:LYS:CE	2.49	0.57
1:A:164:ILE:CG2	1:B:164:ILE:CG1	2.83	0.57
1:A:15:TRP:CD2	1:A:208:LYS:HE2	2.40	0.56
1:A:109:ARG:CZ	1:A:109:ARG:HB3	2.34	0.56
1:A:178:ARG:CG	2:A:342:HOH:O	2.53	0.56
1:A:167:TYR:CZ	1:B:174:LEU:HD13	2.40	0.56
1:A:226:HIS:HE1	2:A:233:HOH:O	1.87	0.56
1:A:51:THR:HG23	1:A:69:PRO:HG2	1.86	0.56
1:B:104:LEU:O	1:B:123:CYS:HB2	2.06	0.56
1:A:43:GLN:HB2	1:A:45:TRP:CH2	2.41	0.56
1:A:192:GLN:O	1:A:196:GLU:HB2	2.05	0.55
1:A:129:SER:HB2	2:A:285:HOH:O	2.05	0.55
1:A:138:PRO:O	1:A:142:MSE:HG3	2.08	0.54
1:B:219:MSE:CG	2:B:390:HOH:O	2.31	0.54
1:A:188:SER:HA	1:A:191:GLU:HG2	1.90	0.54
1:A:137:ARG:NH1	1:B:41:LEU:O	2.36	0.53
1:B:152:GLU:HA	1:B:155:THR:HG22	1.91	0.53
1:A:9:LEU:HD22	1:A:226:HIS:CE1	2.43	0.53
1:A:193:PHE:HD2	1:A:194:MSE:HG3	1.73	0.53
1:A:175:ILE:HG23	1:B:167:TYR:CE2	2.44	0.53
1:B:10:MSE:HG2	1:B:223:THR:CG2	2.39	0.53
1:B:190:LEU:O	1:B:194:MSE:HG2	2.09	0.53
1:A:18:LEU:HB3	1:A:95:PHE:HB2	1.91	0.52
1:A:77:ASP:O	1:A:81:ARG:HG2	2.08	0.52
1:A:157:LEU:HD13	1:A:184:PHE:CD1	2.44	0.52
1:B:188:SER:HB2	2:B:328:HOH:O	2.09	0.52
1:A:180:LYS:HB3	2:A:345:HOH:O	2.09	0.52
1:A:219:MSE:CE	2:A:280:HOH:O	2.51	0.52
1:A:39:SER:HB2	1:A:44:VAL:HG22	1.90	0.52
1:A:170:SER:CA	2:A:344:HOH:O	2.58	0.52
1:B:31:LYS:HD3	2:B:343:HOH:O	2.10	0.52
1:A:15:TRP:CH2	1:A:17:GLN:HB3	2.45	0.52
1:A:194:MSE:CE	1:B:217:LEU:HD13	2.27	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:GLU:HA	1:A:20:GLU:OE1	2.10	0.51
1:A:157:LEU:HG	1:B:157:LEU:HG	1.93	0.51
1:A:170:SER:O	1:A:171:GLY:C	2.53	0.51
1:B:75:HIS:O	1:B:78:ASN:HB2	2.11	0.51
1:B:9:LEU:HB3	2:B:279:HOH:O	2.10	0.50
1:B:177:ASP:C	1:B:178:ARG:HG3	2.34	0.50
1:A:36:LEU:HD22	1:A:121:PHE:CD2	2.47	0.50
1:B:10:MSE:SE	1:B:223:THR:HG22	2.61	0.50
1:B:3:GLU:HB3	2:B:356:HOH:O	2.11	0.50
1:A:176:ARG:O	1:A:177:ASP:HB2	2.11	0.50
1:B:10:MSE:CG	1:B:223:THR:HG22	2.42	0.50
1:A:51:THR:HG23	1:A:69:PRO:HG3	1.94	0.49
1:A:75:HIS:HB2	1:A:112:LEU:HD12	1.94	0.49
1:A:100:VAL:O	1:A:101:ALA:C	2.55	0.49
1:A:170:SER:N	2:A:344:HOH:O	2.45	0.49
1:B:78:ASN:O	1:B:79:LEU:C	2.55	0.49
1:A:97:CYS:SG	1:A:104:LEU:HD11	2.52	0.49
1:B:38:VAL:HG12	1:B:39:SER:N	2.26	0.49
1:A:142:MSE:HE1	1:A:214:LEU:HB3	1.94	0.49
1:A:227:HIS:CE1	2:A:242:HOH:O	2.62	0.49
1:A:109:ARG:HE	1:A:118:TYR:HE2	1.61	0.49
1:A:170:SER:C	1:A:172:ALA:N	2.70	0.48
1:B:219:MSE:O	1:B:223:THR:HG23	2.13	0.48
1:A:175:ILE:HG23	1:B:167:TYR:CZ	2.47	0.48
1:A:198:LEU:N	1:A:199:PRO:CD	2.76	0.48
1:A:15:TRP:CE2	1:A:208:LYS:HE2	2.49	0.48
1:A:157:LEU:HD21	1:B:156:LEU:HB3	1.96	0.48
1:B:34:TYR:C	1:B:34:TYR:CD1	2.91	0.48
1:B:54:VAL:HG21	1:B:73:LEU:HD21	1.96	0.48
1:B:124:MSE:HG2	2:B:367:HOH:O	2.13	0.48
1:B:45:TRP:HB3	1:B:123:CYS:HB3	1.96	0.48
1:B:1:MSE:CE	1:B:50:ASP:HB3	2.43	0.47
1:B:56:GLN:HG2	1:B:60:GLU:OE2	2.14	0.47
1:B:194:MSE:HE3	1:B:198:LEU:HD22	1.96	0.47
1:A:16:LEU:HD22	1:A:84:LEU:HD11	1.96	0.47
1:B:76:LEU:O	1:B:79:LEU:HB2	2.14	0.47
1:A:81:ARG:HB2	1:A:82:PRO:HD3	1.96	0.47
1:B:70:ALA:HA	1:B:73:LEU:HD12	1.96	0.47
1:B:182:GLU:HA	1:B:183:PRO:HD2	1.68	0.47
1:A:220:ALA:CB	1:B:194:MSE:HE2	2.44	0.47
1:A:170:SER:O	1:A:172:ALA:N	2.48	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:194:MSE:HE1	1:B:198:LEU:HD22	1.97	0.46
1:B:2:GLU:HB3	2:B:363:HOH:O	2.15	0.46
1:B:56:GLN:NE2	2:B:344:HOH:O	2.29	0.46
1:B:198:LEU:HB3	1:B:199:PRO:HD3	1.98	0.46
1:A:50:ASP:OD1	1:A:53:VAL:HG23	2.16	0.46
1:B:29:ILE:HD13	1:B:34:TYR:HB2	1.98	0.46
1:B:94:THR:HB	1:B:109:ARG:HB3	1.97	0.45
1:A:45:TRP:HB3	1:A:123:CYS:HB2	1.98	0.45
1:A:170:SER:HB2	1:B:175:ILE:HD12	1.97	0.45
1:A:19:ALA:O	1:A:20:GLU:HB3	2.16	0.45
1:B:176:ARG:CB	1:B:179:LEU:HD13	2.47	0.45
1:B:190:LEU:HD23	1:B:190:LEU:HA	1.87	0.45
1:A:208:LYS:HG3	1:A:212:MSE:CE	2.47	0.45
1:B:79:LEU:HA	2:B:270:HOH:O	2.16	0.45
1:A:106:LEU:O	1:A:121:PHE:N	2.39	0.45
1:A:207:GLY:O	1:A:211:VAL:HG13	2.17	0.44
1:B:23:LEU:HD13	1:B:38:VAL:HG11	1.98	0.44
1:B:149:GLN:O	1:B:150:VAL:C	2.60	0.44
1:B:179:LEU:O	1:B:180:LYS:C	2.60	0.44
1:A:153:LEU:HD23	1:A:156:LEU:HD12	2.00	0.43
1:A:169:GLU:C	2:A:344:HOH:O	2.61	0.43
1:A:61:LEU:CD1	1:A:119:TRP:HA	2.48	0.43
1:A:165:GLN:HA	1:A:168:GLN:HG2	2.00	0.43
2:A:268:HOH:O	1:B:156:LEU:HG	2.17	0.43
1:A:170:SER:C	2:A:344:HOH:O	2.61	0.42
1:A:45:TRP:HB3	1:A:123:CYS:CB	2.50	0.42
1:B:217:LEU:O	1:B:221:VAL:HG23	2.19	0.42
1:B:105:ILE:HA	1:B:121:PHE:O	2.19	0.42
1:A:17:GLN:NE2	2:A:346:HOH:O	2.52	0.42
1:B:176:ARG:HB2	1:B:179:LEU:HD13	2.02	0.42
1:A:1:MSE:HA	1:A:1:MSE:HE2	2.01	0.42
1:A:166:ASP:OD2	1:B:174:LEU:HD12	2.19	0.42
1:B:208:LYS:HB3	1:B:209:PRO:HD3	2.02	0.42
1:B:37:LEU:CD1	1:B:44:VAL:HG12	2.43	0.41
1:B:135:LEU:C	1:B:138:PRO:HD2	2.45	0.41
1:A:10:MSE:HE3	1:A:227:HIS:NE2	2.35	0.41
1:A:198:LEU:CD2	1:B:220:ALA:HB1	2.49	0.41
1:B:108:VAL:HG21	1:B:121:PHE:HE2	1.86	0.41
1:A:8:LEU:HD12	1:A:8:LEU:HA	1.95	0.41
1:A:176:ARG:O	1:A:176:ARG:HG2	2.21	0.41
1:B:208:LYS:HA	1:B:211:VAL:HG22	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:137:ARG:HB2	1:B:138:PRO:HD3	2.03	0.41
1:A:198:LEU:HD21	1:B:220:ALA:HB1	2.02	0.40
1:B:55:SER:HB3	1:B:69:PRO:HG3	2.03	0.40
1:A:140:MSE:HE1	1:B:207:GLY:HA2	2.03	0.40
1:B:50:ASP:HB2	2:B:335:HOH:O	2.21	0.40
1:B:17:GLN:C	1:B:18:LEU:HD23	2.46	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	215/230 (94%)	198 (92%)	15 (7%)	2 (1%)	14	27
1	B	210/230 (91%)	201 (96%)	7 (3%)	2 (1%)	12	24
All	All	425/460 (92%)	399 (94%)	22 (5%)	4 (1%)	14	27

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	172	ALA
1	B	178	ARG
1	A	64	ARG
1	A	171	GLY

5.3.2 Protein sidechains [i](#)

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	195/195 (100%)	187 (96%)	8 (4%)	27	53
1	B	190/195 (97%)	180 (95%)	10 (5%)	20	42
All	All	385/390 (99%)	367 (95%)	18 (5%)	23	47

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	LEU
1	A	29	ILE
1	A	38	VAL
1	A	74	CYS
1	A	84	LEU
1	A	100	VAL
1	A	164	ILE
1	A	211	VAL
1	B	29	ILE
1	B	49	VAL
1	B	61	LEU
1	B	100	VAL
1	B	105	ILE
1	B	110	SER
1	B	135	LEU
1	B	156	LEU
1	B	174	LEU
1	B	224	GLN

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	11	GLN
1	A	226	HIS
1	B	21	ASN
1	B	42	GLN
1	B	43	GLN
1	B	120	ASN
1	B	187	ASN
1	B	213	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	210/230 (91%)	0.96	22 (10%)	11 10	31, 43, 62, 71	0
1	B	205/230 (89%)	0.88	21 (10%)	12 10	31, 42, 60, 71	0
All	All	415/460 (90%)	0.92	43 (10%)	11 10	31, 42, 61, 71	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	63	LYS	7.0
1	A	176	ARG	4.3
1	B	177	ASP	4.2
1	B	178	ARG	4.0
1	B	79	LEU	4.0
1	A	170	SER	3.7
1	B	170	SER	3.6
1	B	176	ARG	3.5
1	A	64	ARG	3.5
1	A	175	ILE	3.3
1	B	100	VAL	3.2
1	B	98	ASP	3.1
1	B	65	LEU	3.1
1	B	171	GLY	3.1
1	A	56	GLN	3.1
1	A	227	HIS	3.0
1	A	171	GLY	3.0
1	B	18	LEU	2.9
1	A	112	LEU	2.8
1	A	174	LEU	2.7
1	B	3	GLU	2.6
1	A	206	ASP	2.6
1	A	177	ASP	2.5
1	A	172	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	113	SER	2.5
1	B	164	ILE	2.4
1	B	37	LEU	2.4
1	A	207	GLY	2.4
1	B	20	GLU	2.4
1	B	19	ALA	2.3
1	B	205	GLY	2.3
1	A	204	ILE	2.3
1	A	62	ASN	2.3
1	A	43	GLN	2.2
1	B	172	ALA	2.1
1	B	173	THR	2.1
1	A	99	CYS	2.1
1	A	114	GLY	2.1
1	B	52	SER	2.1
1	A	19	ALA	2.1
1	B	187	ASN	2.1
1	B	175	ILE	2.1
1	A	210	PHE	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.