



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:21 AM UTC

PDB ID : 2QM4 / pdb_00002qm4
Title : Crystal structure of human XLF/Cernunnos, a non-homologous end-joining factor
Authors : Li, Y.; Chirgadze, D.Y.; Sibanda, B.L.; Bolanos-Garcia, V.M.; Davies, O.R.; Blundell, T.L.
Deposited on : 2007-07-14
Resolution : 2.30 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
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

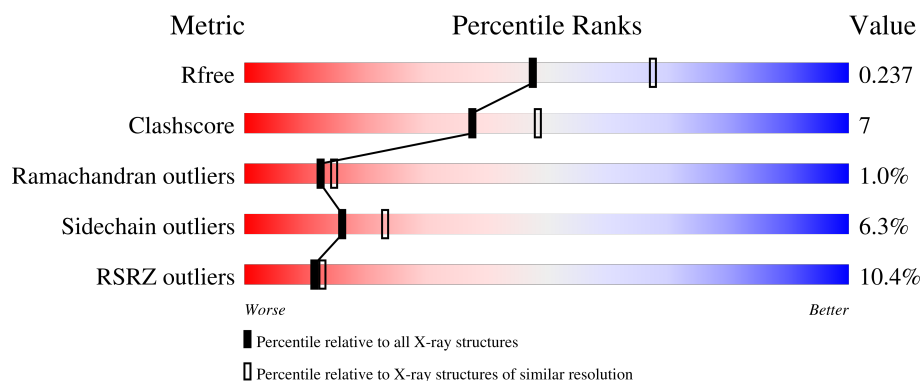
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	235	7% 79% 17% ..
1	B	235	16% 78% 16% ..
1	C	235	11% 76% 20% ..
1	D	235	5% 75% 19% ..

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 7510 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	232	Total	C	N	O	S	Se	0	7	0
			1864	1191	312	344	6	11			
1	B	225	Total	C	N	O	S	Se	0	10	0
			1821	1167	304	334	6	10			
1	C	229	Total	C	N	O	S	Se	0	4	0
			1818	1160	305	337	6	10			
1	D	225	Total	C	N	O	S	Se	0	7	0
			1775	1137	292	330	6	10			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-1	SER	-	expression tag	UNP Q9H9Q4
A	0	GLY	-	expression tag	UNP Q9H9Q4
A	1	MSE	MET	modified residue	UNP Q9H9Q4
A	10	MSE	MET	modified residue	UNP Q9H9Q4
A	124	MSE	MET	modified residue	UNP Q9H9Q4
A	140	MSE	MET	modified residue	UNP Q9H9Q4
A	142	MSE	MET	modified residue	UNP Q9H9Q4
A	159	MSE	MET	modified residue	UNP Q9H9Q4
A	194	MSE	MET	modified residue	UNP Q9H9Q4
A	212	MSE	MET	modified residue	UNP Q9H9Q4
A	219	MSE	MET	modified residue	UNP Q9H9Q4
B	-1	SER	-	expression tag	UNP Q9H9Q4
B	0	GLY	-	expression tag	UNP Q9H9Q4
B	1	MSE	MET	modified residue	UNP Q9H9Q4
B	10	MSE	MET	modified residue	UNP Q9H9Q4
B	124	MSE	MET	modified residue	UNP Q9H9Q4
B	140	MSE	MET	modified residue	UNP Q9H9Q4
B	142	MSE	MET	modified residue	UNP Q9H9Q4
B	159	MSE	MET	modified residue	UNP Q9H9Q4
B	194	MSE	MET	modified residue	UNP Q9H9Q4
B	212	MSE	MET	modified residue	UNP Q9H9Q4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	219	MSE	MET	modified residue	UNP Q9H9Q4
C	-1	SER	-	expression tag	UNP Q9H9Q4
C	0	GLY	-	expression tag	UNP Q9H9Q4
C	1	MSE	MET	modified residue	UNP Q9H9Q4
C	10	MSE	MET	modified residue	UNP Q9H9Q4
C	124	MSE	MET	modified residue	UNP Q9H9Q4
C	140	MSE	MET	modified residue	UNP Q9H9Q4
C	142	MSE	MET	modified residue	UNP Q9H9Q4
C	159	MSE	MET	modified residue	UNP Q9H9Q4
C	194	MSE	MET	modified residue	UNP Q9H9Q4
C	212	MSE	MET	modified residue	UNP Q9H9Q4
C	219	MSE	MET	modified residue	UNP Q9H9Q4
D	-1	SER	-	expression tag	UNP Q9H9Q4
D	0	GLY	-	expression tag	UNP Q9H9Q4
D	1	MSE	MET	modified residue	UNP Q9H9Q4
D	10	MSE	MET	modified residue	UNP Q9H9Q4
D	124	MSE	MET	modified residue	UNP Q9H9Q4
D	140	MSE	MET	modified residue	UNP Q9H9Q4
D	142	MSE	MET	modified residue	UNP Q9H9Q4
D	159	MSE	MET	modified residue	UNP Q9H9Q4
D	194	MSE	MET	modified residue	UNP Q9H9Q4
D	212	MSE	MET	modified residue	UNP Q9H9Q4
D	219	MSE	MET	modified residue	UNP Q9H9Q4

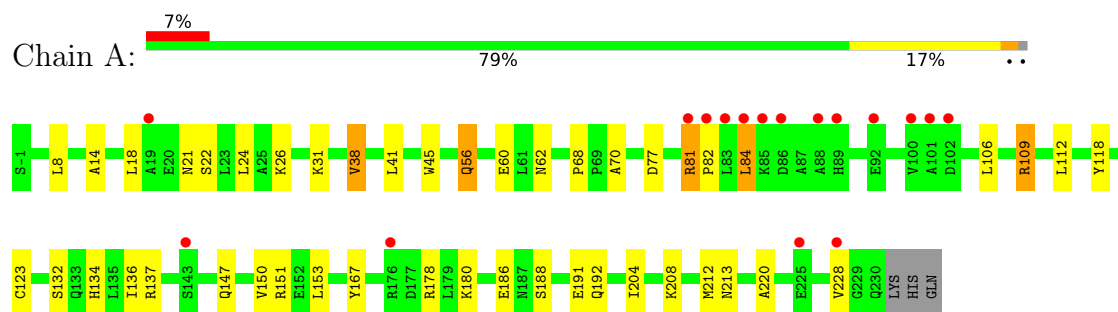
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	73	Total O 73 73	0	0
2	B	49	Total O 49 49	0	0
2	C	47	Total O 47 47	0	0
2	D	63	Total O 63 63	0	0

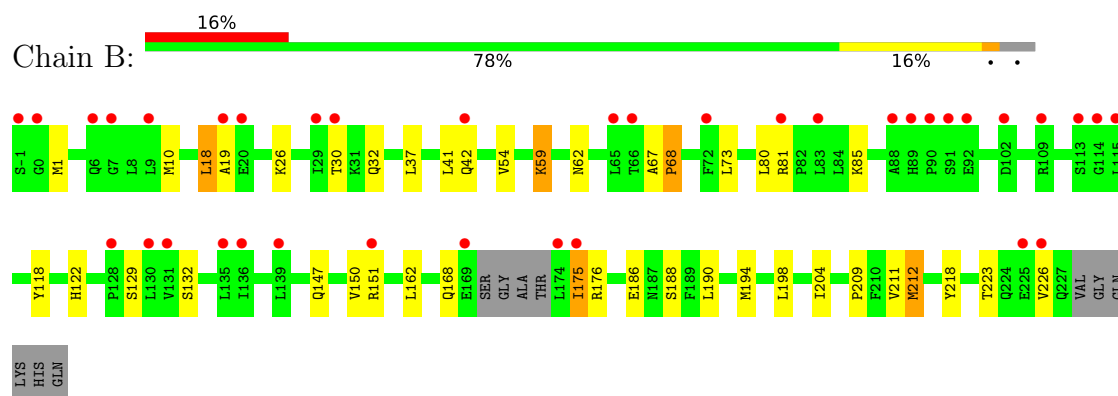
3 Residue-property plots [i](#)

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.

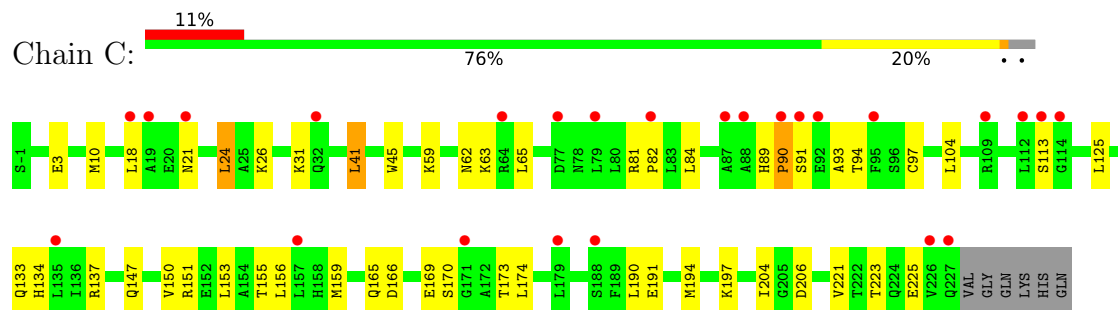
- Molecule 1: Non-homologous end-joining factor 1



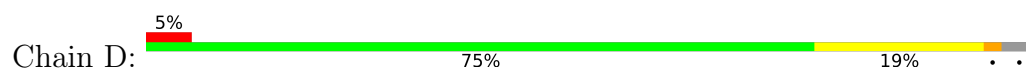
- Molecule 1: Non-homologous end-joining factor 1

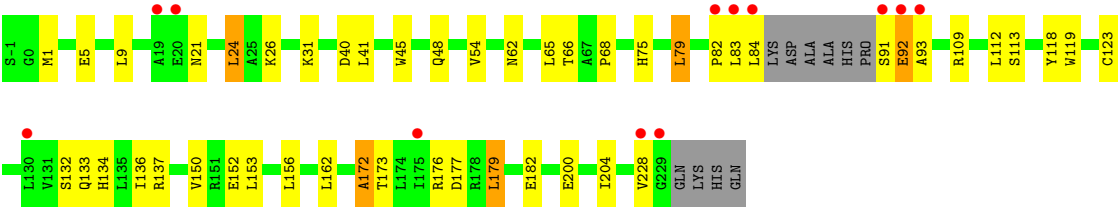


- Molecule 1: Non-homologous end-joining factor 1



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





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.74Å 92.91Å 103.69Å 90.00° 106.22° 90.00°	Depositor
Resolution (Å)	37.01 – 2.30 37.01 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (37.01-2.30) 92.7 (37.01-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.75 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.182 , 0.239 0.180 , 0.237	Depositor DCC
R_{free} test set	2000 reflections (4.17%)	wwPDB-VP
Wilson B-factor (Å ²)	38.4	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 45.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.035 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7510	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.62% 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 [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	0.79	1/1909 (0.1%)	0.99	2/2570 (0.1%)
1	B	0.75	0/1875	0.93	3/2526 (0.1%)
1	C	0.71	0/1856	0.95	3/2500 (0.1%)
1	D	0.81	2/1821 (0.1%)	0.97	1/2454 (0.0%)
All	All	0.77	3/7461 (0.0%)	0.96	9/10050 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	68	PRO	CA-C	7.20	1.55	1.51
1	D	68	PRO	CA-C	7.13	1.55	1.51
1	D	92	GLU	CA-CB	6.44	1.64	1.53

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	90	PRO	N-CA-CB	7.36	110.98	103.25
1	B	68	PRO	CA-C-N	6.05	125.53	119.24
1	B	68	PRO	C-N-CA	6.05	125.53	119.24
1	A	136	ILE	N-CA-C	5.79	115.92	110.30
1	D	172	ALA	N-CA-C	5.70	122.95	110.80
1	A	150	VAL	N-CA-C	5.56	116.19	110.36
1	B	186	GLU	N-CA-C	5.34	116.79	111.07
1	C	133	GLN	N-CA-C	5.16	116.98	111.36
1	C	93	ALA	N-CA-C	5.08	117.26	110.35

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	1864	0	1886	35	0
1	B	1821	0	1825	29	0
1	C	1818	0	1821	20	0
1	D	1775	0	1765	33	0
2	A	73	0	0	1	0
2	B	49	0	0	2	0
2	C	47	0	0	1	0
2	D	63	0	0	4	0
All	All	7510	0	7297	102	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 7.

All (102) 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:109:ARG:HH11	1:A:109:ARG:HG3	1.24	1.00
1:A:62:ASN:HD21	1:A:118:TYR:H	1.20	0.89
1:A:208:LYS:NZ	1:A:212[A]:MSE:HE3	1.87	0.89
1:B:211:VAL:HG12	1:B:212[A]:MSE:HE3	1.55	0.87
1:D:62:ASN:HD21	1:D:118:TYR:H	1.28	0.82
1:A:188:SER:O	1:A:192:GLN:HG2	1.81	0.81
1:A:208:LYS:HZ2	1:A:212[A]:MSE:HE3	1.47	0.78
1:D:1:MSE:HE2	2:D:283:HOH:O	1.84	0.77
1:A:109:ARG:HH11	1:A:109:ARG:CG	1.98	0.75
1:B:212[A]:MSE:HA	1:B:212[A]:MSE:HE2	1.69	0.74
1:D:75:HIS:CE1	1:D:79:LEU:HD21	2.26	0.71
1:B:30:THR:OG1	1:B:32:GLN:HG2	1.91	0.71
1:C:155:THR:O	1:C:159[B]:MSE:HG2	1.91	0.69
1:A:81:ARG:HB3	1:A:82:PRO:HD3	1.77	0.67
1:C:206:ASP:HB2	2:C:266:HOH:O	1.93	0.67
1:C:165:GLN:O	1:C:169:GLU:HG2	1.97	0.65
1:C:18:LEU:HD22	1:C:97:CYS:HB2	1.77	0.64
1:D:228:VAL:HG12	1:D:228:VAL:O	1.98	0.63
1:B:62:ASN:HD21	1:B:118:TYR:H	1.48	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:109:ARG:HG3	1:A:109:ARG:NH1	2.01	0.62
1:A:21:ASN:CG	1:A:22:SER:H	2.08	0.62
1:B:223:THR:O	1:B:226:VAL:HG22	2.00	0.62
1:C:26:LYS:NZ	1:C:134:HIS:HD2	1.98	0.61
1:A:26:LYS:NZ	1:A:134:HIS:HD2	2.01	0.59
1:A:208:LYS:HZ3	1:A:212[A]:MSE:HE3	1.67	0.57
1:C:166:ASP:HB2	1:D:179:LEU:HD23	1.85	0.56
1:B:54:VAL:HG21	1:B:73:LEU:HD21	1.88	0.56
1:B:175:ILE:HG23	1:B:176:ARG:H	1.70	0.56
1:A:153:LEU:HD11	1:B:150:VAL:HG13	1.87	0.56
1:B:175:ILE:HG23	1:B:176:ARG:N	2.20	0.56
1:B:10:MSE:HA	1:B:10:MSE:HE2	1.87	0.56
1:C:197:LYS:HG2	1:D:152:GLU:OE2	2.06	0.56
1:C:10:MSE:HA	1:C:10:MSE:HE2	1.88	0.55
1:A:14:ALA:HB2	1:A:84:LEU:HD12	1.89	0.55
1:C:190:LEU:O	1:C:194:MSE:HG2	2.06	0.55
1:A:31[B]:LYS:HE3	1:A:70:ALA:HB1	1.88	0.55
1:B:26:LYS:NZ	1:B:218:TYR:OH	2.41	0.54
1:C:153:LEU:HD11	1:D:150:VAL:HG13	1.90	0.53
1:D:1:MSE:HE3	1:D:48:GLN:OE1	2.09	0.53
1:C:147:GLN:HE21	1:C:151:ARG:NH1	2.06	0.53
1:B:212[A]:MSE:HA	1:B:212[A]:MSE:CE	2.39	0.53
1:A:8:LEU:HD11	1:A:26:LYS:HG3	1.92	0.51
1:B:59:LYS:HG3	2:B:247:HOH:O	2.10	0.50
1:C:137:ARG:O	1:D:204:ILE:HD11	2.11	0.50
1:D:1:MSE:CE	2:D:283:HOH:O	2.50	0.49
1:A:109:ARG:CG	1:A:109:ARG:NH1	2.68	0.48
1:A:137[B]:ARG:NH1	1:B:42[B]:GLN:NE2	2.60	0.48
1:D:176:ARG:HB2	1:D:179:LEU:HD22	1.94	0.48
1:A:31[B]:LYS:HG3	1:A:70:ALA:HB1	1.95	0.48
1:A:41:LEU:HD13	1:A:204:ILE:CG2	2.44	0.48
1:B:209:PRO:HA	1:B:212[B]:MSE:HE2	1.95	0.48
1:B:1:MSE:CE	1:D:66:THR:HG21	2.45	0.47
1:D:109:ARG:NH2	2:D:287:HOH:O	2.47	0.47
1:B:10:MSE:HB2	2:B:245:HOH:O	2.14	0.46
1:C:26:LYS:HZ2	1:C:134:HIS:HD2	1.63	0.46
1:D:54:VAL:HA	1:D:119:TRP:HH2	1.80	0.46
1:A:147:GLN:O	1:A:151:ARG:HG2	2.14	0.46
1:C:204:ILE:HD11	1:D:137[B]:ARG:CZ	2.45	0.46
1:D:9:LEU:HD21	1:D:133:GLN:HG2	1.97	0.46
1:A:167:TYR:OH	1:B:168:GLN:HG2	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:26:LYS:HE2	1:D:134:HIS:HD2	1.80	0.45
1:D:31:LYS:NZ	2:D:281:HOH:O	2.38	0.45
1:B:190:LEU:O	1:B:194:MSE:HG2	2.16	0.45
1:B:41:LEU:HD13	1:B:204:ILE:HD12	1.99	0.45
1:D:45:TRP:HB3	1:D:123:CYS:HB3	1.99	0.44
1:A:134:HIS:HE1	2:A:234:HOH:O	2.01	0.44
1:B:122:HIS:HB3	1:D:113:SER:O	2.16	0.44
1:D:5:GLU:HG2	1:D:48:GLN:NE2	2.32	0.44
1:D:75:HIS:CD2	1:D:112:LEU:HD13	2.52	0.44
1:A:77:ASP:O	1:A:81:ARG:HB2	2.17	0.44
1:D:83:LEU:O	1:D:84:LEU:CB	2.65	0.43
1:A:26:LYS:HZ3	1:A:134:HIS:HD2	1.63	0.43
1:D:132:SER:HA	1:D:136:ILE:HB	2.00	0.43
1:B:175:ILE:CG2	1:B:176:ARG:N	2.82	0.43
1:D:62:ASN:HD21	1:D:118:TYR:N	2.06	0.43
1:A:56:GLN:NE2	1:A:60:GLU:OE2	2.52	0.43
1:D:228:VAL:O	1:D:228:VAL:CG1	2.65	0.43
1:A:18:LEU:HD21	1:A:106:LEU:CD1	2.49	0.42
1:A:45:TRP:HB3	1:A:123:CYS:HB3	2.01	0.42
1:C:10:MSE:HE1	1:C:223:THR:HA	2.01	0.42
1:A:178:ARG:HD2	1:A:180:LYS:HE3	2.01	0.42
1:B:18:LEU:O	1:B:19:ALA:C	2.61	0.42
1:A:38[A]:VAL:HG12	1:A:45:TRP:HB2	2.01	0.42
1:C:24:LEU:HD22	1:C:41:LEU:CD1	2.48	0.42
1:C:81:ARG:N	1:C:82:PRO:HD2	2.34	0.42
1:A:213:ASN:HB3	1:B:147[A]:GLN:OE1	2.19	0.42
1:D:1:MSE:CE	1:D:48:GLN:HG2	2.49	0.42
1:D:24:LEU:HD22	1:D:41:LEU:HD21	2.02	0.42
1:C:45:TRP:CE3	1:C:104:LEU:HD22	2.55	0.42
1:D:91:SER:C	1:D:93:ALA:H	2.27	0.42
1:D:176:ARG:CB	1:D:179:LEU:HD22	2.50	0.41
1:A:31[B]:LYS:HG3	1:A:70:ALA:CB	2.51	0.41
1:A:137[A]:ARG:CZ	1:B:204:ILE:HD11	2.51	0.41
1:A:153:LEU:CD1	1:B:150:VAL:HG13	2.50	0.41
1:D:40:ASP:O	1:D:41:LEU:HB2	2.20	0.41
1:A:220:ALA:HB1	1:B:198:LEU:CD2	2.51	0.41
1:C:221:VAL:O	1:C:225:GLU:HG3	2.20	0.41
1:A:109:ARG:HD2	1:A:118:TYR:CE1	2.56	0.40
1:B:67:ALA:HB1	1:B:68:PRO:HD2	2.04	0.40
1:B:80:LEU:HD12	1:B:80:LEU:HA	1.95	0.40
1:C:150:VAL:HG13	1:D:153:LEU:HD11	2.04	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	237/235 (101%)	228 (96%)	9 (4%)	0	100	100
1	B	231/235 (98%)	219 (95%)	12 (5%)	0	100	100
1	C	231/235 (98%)	218 (94%)	8 (4%)	5 (2%)	5	4
1	D	228/235 (97%)	214 (94%)	10 (4%)	4 (2%)	6	6
All	All	927/940 (99%)	879 (95%)	39 (4%)	9 (1%)	12	15

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	89	HIS
1	C	90	PRO
1	C	91	SER
1	D	92	GLU
1	D	172	ALA
1	C	113	SER
1	D	82	PRO
1	D	182	GLU
1	C	170	SER

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	208/198 (105%)	196 (94%)	12 (6%)	18	26
1	B	201/198 (102%)	187 (93%)	14 (7%)	14	19
1	C	200/198 (101%)	184 (92%)	16 (8%)	11	15
1	D	195/198 (98%)	184 (94%)	11 (6%)	19	28
All	All	804/792 (102%)	751 (93%)	53 (7%)	16	21

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LEU
1	A	38[A]	VAL
1	A	38[B]	VAL
1	A	56	GLN
1	A	81	ARG
1	A	84	LEU
1	A	109	ARG
1	A	112	LEU
1	A	132	SER
1	A	186	GLU
1	A	191	GLU
1	A	228	VAL
1	B	18	LEU
1	B	37	LEU
1	B	59	LYS
1	B	81	ARG
1	B	85	LYS
1	B	129	SER
1	B	132	SER
1	B	151[A]	ARG
1	B	151[B]	ARG
1	B	162	LEU
1	B	175	ILE
1	B	188	SER
1	B	212[A]	MSE
1	B	212[B]	MSE
1	C	3	GLU
1	C	21	ASN
1	C	24	LEU
1	C	31	LYS
1	C	41	LEU
1	C	59	LYS
1	C	62	ASN

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Mol	Chain	Res	Type
1	C	63	LYS
1	C	65	LEU
1	C	84	LEU
1	C	94	THR
1	C	125	LEU
1	C	156	LEU
1	C	173	THR
1	C	174	LEU
1	C	191	GLU
1	D	21	ASN
1	D	24	LEU
1	D	65	LEU
1	D	79	LEU
1	D	156	LEU
1	D	162	LEU
1	D	173	THR
1	D	177	ASP
1	D	179	LEU
1	D	200[A]	GLU
1	D	200[B]	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	GLN
1	A	32	GLN
1	A	62	ASN
1	A	120	ASN
1	A	134	HIS
1	A	192	GLN
1	A	227	GLN
1	B	11	GLN
1	B	48	GLN
1	B	62	ASN
1	B	120	ASN
1	C	32	GLN
1	C	120	ASN
1	C	122	HIS
1	C	134	HIS
1	C	147	GLN
1	C	187	ASN
1	D	56	GLN

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Mol	Chain	Res	Type
1	D	62	ASN
1	D	75	HIS
1	D	134	HIS
1	D	165	GLN
1	D	192	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 [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	223/235 (94%)	0.69	17 (7%) 20 21	25, 54, 85, 112	7 (3%)
1	B	216/235 (91%)	1.20	37 (17%) 4 5	28, 56, 89, 106	10 (4%)
1	C	220/235 (93%)	1.00	25 (11%) 10 11	30, 58, 82, 94	6 (2%)
1	D	216/235 (91%)	0.74	12 (5%) 30 32	29, 52, 74, 96	6 (2%)
All	All	875/940 (93%)	0.91	91 (10%) 11 13	25, 55, 85, 112	29 (3%)

All (91) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	83	LEU	6.6
1	B	174	LEU	6.0
1	A	84	LEU	4.7
1	B	90	PRO	4.0
1	A	82	PRO	3.7
1	B	175	ILE	3.7
1	B	225[A]	GLU	3.6
1	B	66	THR	3.5
1	B	113	SER	3.5
1	D	92	GLU	3.4
1	A	86	ASP	3.4
1	A	100	VAL	3.3
1	B	88	ALA	3.3
1	D	84	LEU	3.3
1	C	91	SER	3.3
1	C	113	SER	3.3
1	C	112	LEU	3.2
1	C	109[A]	ARG	3.1
1	C	18	LEU	3.1
1	D	175	ILE	3.0
1	D	91	SER	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	6	GLN	3.0
1	B	42[A]	GLN	3.0
1	D	229	GLY	2.9
1	C	88	ALA	2.9
1	B	135	LEU	2.9
1	D	19	ALA	2.9
1	D	82	PRO	2.8
1	B	89	HIS	2.8
1	C	90	PRO	2.8
1	A	83	LEU	2.7
1	B	91	SER	2.7
1	A	89	HIS	2.7
1	D	228	VAL	2.7
1	B	130	LEU	2.7
1	A	101	ALA	2.6
1	B	131	VAL	2.6
1	C	21	ASN	2.6
1	C	114	GLY	2.6
1	C	135	LEU	2.5
1	C	227	GLN	2.5
1	A	81	ARG	2.5
1	C	32	GLN	2.5
1	B	-1	SER	2.5
1	C	92	GLU	2.5
1	C	95	PHE	2.5
1	B	109[A]	ARG	2.5
1	A	19	ALA	2.5
1	B	136	ILE	2.4
1	B	226	VAL	2.4
1	B	169	GLU	2.4
1	B	9	LEU	2.4
1	B	30	THR	2.4
1	B	151[A]	ARG	2.4
1	D	20	GLU	2.4
1	C	171	GLY	2.3
1	B	19	ALA	2.3
1	C	188	SER	2.3
1	B	128	PRO	2.3
1	B	29	ILE	2.3
1	B	83	LEU	2.3
1	B	115	LEU	2.3
1	C	19	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	228	VAL	2.3
1	B	72	PHE	2.2
1	C	179	LEU	2.2
1	B	102	ASP	2.2
1	C	77	ASP	2.2
1	B	7	GLY	2.2
1	B	0	GLY	2.2
1	A	225	GLU	2.1
1	C	82	PRO	2.1
1	B	114	GLY	2.1
1	C	87	ALA	2.1
1	B	65	LEU	2.1
1	C	79	LEU	2.1
1	B	139	LEU	2.1
1	A	88	ALA	2.1
1	D	93	ALA	2.1
1	A	92	GLU	2.1
1	D	130	LEU	2.1
1	A	102	ASP	2.1
1	C	226	VAL	2.1
1	A	143	SER	2.0
1	B	20	GLU	2.0
1	C	157	LEU	2.0
1	B	81	ARG	2.0
1	A	85	LYS	2.0
1	B	92	GLU	2.0
1	A	176	ARG	2.0
1	C	64	ARG	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.