



Full wwPDB X-ray Structure Validation Report ⓘ

May 29, 2020 – 03:17 am BST

PDB ID : 3MWP
Title : Nucleoprotein structure of lassa fever virus
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Deposited on : 2010-05-06
Resolution : 1.79 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.11
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

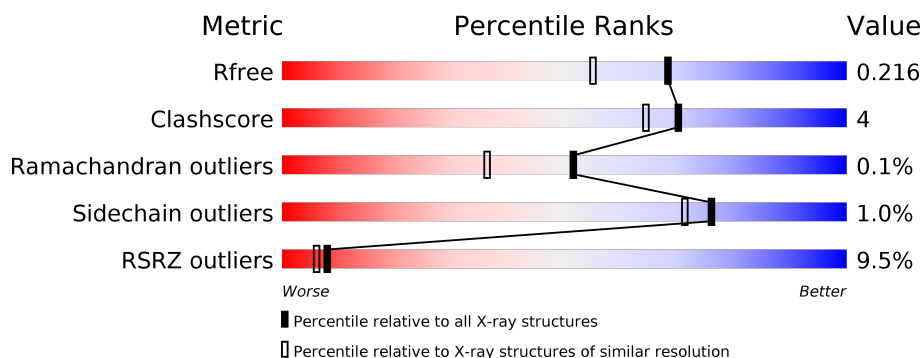
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.79 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}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 on 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	577	
1	B	577	
1	C	577	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZN	B	570	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12477 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 Nucleoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	512	Total	C	N	O	S	0	1	0
			4011	2521	697	766	27			
1	B	515	Total	C	N	O	S	0	0	0
			4027	2528	701	771	27			
1	C	506	Total	C	N	O	S	0	0	0
			3961	2488	688	758	27			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	GLY	-	EXPRESSION TAG	UNP P13699
A	-6	ALA	-	EXPRESSION TAG	UNP P13699
A	-5	MET	-	EXPRESSION TAG	UNP P13699
A	-4	ASP	-	EXPRESSION TAG	UNP P13699
A	-3	HIS	-	EXPRESSION TAG	UNP P13699
A	-2	VAL	-	EXPRESSION TAG	UNP P13699
A	-1	GLU	-	EXPRESSION TAG	UNP P13699
A	0	PHE	-	EXPRESSION TAG	UNP P13699
B	-7	GLY	-	EXPRESSION TAG	UNP P13699
B	-6	ALA	-	EXPRESSION TAG	UNP P13699
B	-5	MET	-	EXPRESSION TAG	UNP P13699
B	-4	ASP	-	EXPRESSION TAG	UNP P13699
B	-3	HIS	-	EXPRESSION TAG	UNP P13699
B	-2	VAL	-	EXPRESSION TAG	UNP P13699
B	-1	GLU	-	EXPRESSION TAG	UNP P13699
B	0	PHE	-	EXPRESSION TAG	UNP P13699
C	-7	GLY	-	EXPRESSION TAG	UNP P13699
C	-6	ALA	-	EXPRESSION TAG	UNP P13699
C	-5	MET	-	EXPRESSION TAG	UNP P13699
C	-4	ASP	-	EXPRESSION TAG	UNP P13699
C	-3	HIS	-	EXPRESSION TAG	UNP P13699
C	-2	VAL	-	EXPRESSION TAG	UNP P13699
C	-1	GLU	-	EXPRESSION TAG	UNP P13699

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Chain	Residue	Modelled	Actual	Comment	Reference
C	0	PHE	-	EXPRESSION TAG	UNP P13699

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Zn	0	0
			1	1		
2	A	1	Total	Zn	0	0
			1	1		
2	C	1	Total	Zn	0	0
			1	1		

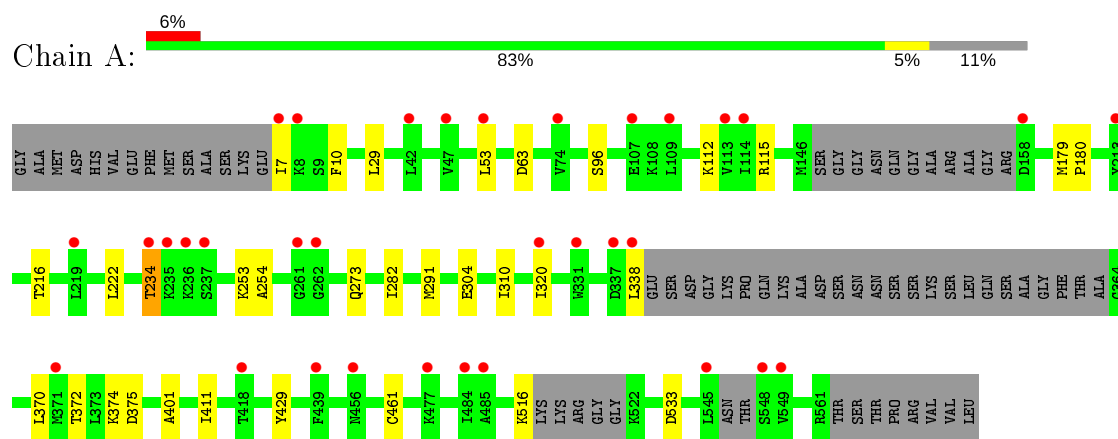
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	175	Total	O	0	0
			175	175		
3	B	192	Total	O	0	0
			192	192		
3	C	108	Total	O	0	0
			108	108		

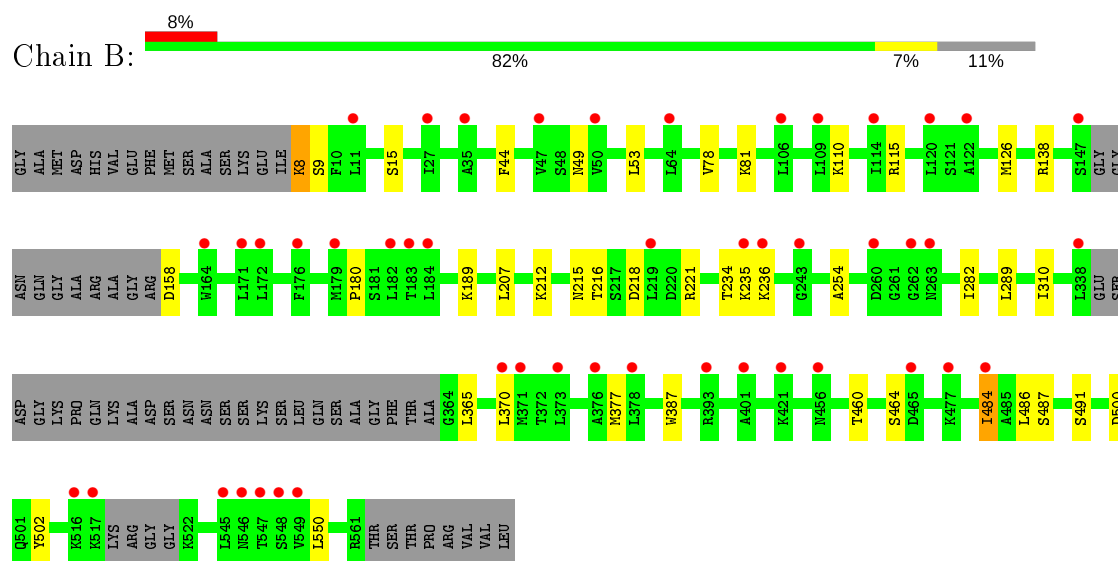
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

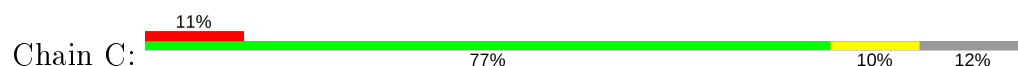
• Molecule 1: Nucleoprotein

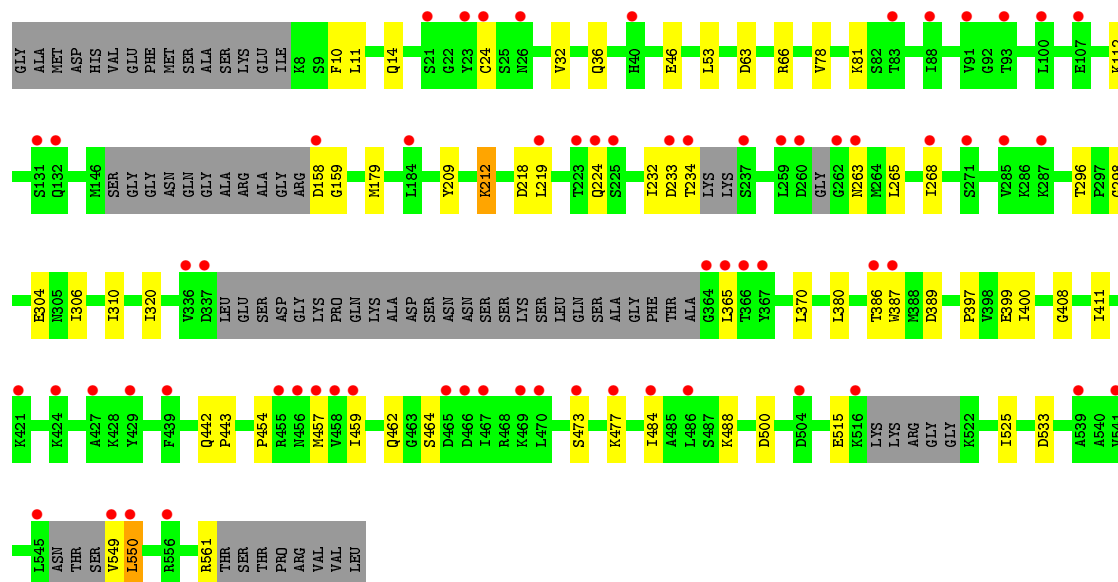


• Molecule 1: Nucleoprotein



• Molecule 1: Nucleoprotein





4 Data and refinement statistics

Property	Value	Source
Space group	P 3	Depositor
Cell constants a, b, c, α , β , γ	176.82Å 176.82Å 56.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	153.18 – 1.79 47.59 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.7 (153.18-1.79) 99.9 (47.59-1.80)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.70 (at 1.79Å)	Xtriage
Refinement program	REFMAC 5.5.0070	Depositor
R, R_{free}	0.178 , 0.196 0.199 , 0.216	Depositor DCC
R_{free} test set	9260 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtriage
Anisotropy	0.560	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	0.008 for -h,-k,l 0.021 for h,-h-k,-l 0.000 for -k,-h,-l	Xtriage
Reported twinning fraction	0.348 for H, K, L 0.652 for K, H, -L	Depositor
Outliers	0 of 184177 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12477	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.53% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.61	1/4070 (0.0%)	0.71	1/5493 (0.0%)
1	B	0.64	0/4084	0.76	4/5513 (0.1%)
1	C	0.57	0/4015	0.68	1/5418 (0.0%)
All	All	0.61	1/12169 (0.0%)	0.72	6/16424 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	461	CYS	CB-SG	-8.33	1.68	1.82

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	377	MET	CG-SD-CE	-13.53	78.55	100.20
1	B	115	ARG	NE-CZ-NH1	6.53	123.56	120.30
1	B	115	ARG	NE-CZ-NH2	-6.32	117.14	120.30
1	A	461	CYS	CB-CA-C	-5.73	98.95	110.40
1	B	500	ASP	CB-CG-OD1	5.07	122.87	118.30
1	C	500	ASP	CB-CG-OD1	5.00	122.80	118.30

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	4011	0	4085	22	0
1	B	4027	0	4098	28	0
1	C	3961	0	4019	38	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	175	0	0	3	0
3	B	192	0	0	2	0
3	C	108	0	0	2	0
All	All	12477	0	12202	86	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 4.

All (86) 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:C:219:LEU:HD23	1:C:234:THR:HG22	1.37	1.01
1:C:24:CYS:SG	1:C:268:ILE:HG12	2.17	0.85
1:B:502:TYR:CE1	1:B:550:LEU:HD13	2.16	0.80
1:C:265:LEU:O	1:C:268:ILE:HG13	1.82	0.78
1:C:408:GLY:O	1:C:550:LEU:HA	1.83	0.78
1:C:533:ASP:OD2	3:C:655:HOH:O	2.01	0.78
1:C:397:PRO:HB2	1:C:400:ILE:HD11	1.75	0.69
1:C:219:LEU:HD23	1:C:234:THR:CG2	2.19	0.66
1:B:502:TYR:CE1	1:B:550:LEU:CD1	2.79	0.66
1:A:533:ASP:OD2	3:A:728:HOH:O	2.13	0.65
1:C:219:LEU:CD2	1:C:234:THR:HG22	2.22	0.63
1:A:7:ILE:HG23	1:A:10:PHE:H	1.65	0.62
1:C:112:LYS:HE3	1:C:304:GLU:OE1	2.00	0.61
1:C:515:GLU:HG3	1:C:525:ILE:HD13	1.82	0.61
1:A:216:THR:HG23	1:A:234:THR:HG21	1.86	0.57
1:C:365:LEU:HD22	1:C:370:LEU:HD23	1.86	0.57
1:C:387:TRP:CH2	1:C:484:ILE:HG12	2.41	0.56
1:C:389:ASP:OD1	3:C:655:HOH:O	2.18	0.56
1:B:212:LYS:HE3	1:B:218:ASP:OD2	2.07	0.54
1:A:115:ARG:HD2	1:A:375:ASP:OD2	2.08	0.54
1:B:49:ASN:O	1:B:53:LEU:HG	2.08	0.53
1:B:365:LEU:HD22	1:B:370:LEU:HD23	1.88	0.53
1:B:78:VAL:O	1:B:81:LYS:HE2	2.09	0.52
1:A:179:MET:HG3	1:A:320:ILE:O	2.10	0.52
1:B:215:ASN:HB2	3:B:681:HOH:O	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:112:LYS:NZ	1:C:298:GLY:O	2.44	0.50
1:C:442:GLN:NE2	1:C:443:PRO:HD2	2.25	0.50
1:C:306:ILE:HG12	1:C:310:ILE:HD12	1.94	0.50
1:C:179:MET:HG3	1:C:320:ILE:O	2.10	0.50
1:C:386:THR:HB	1:C:459:ILE:HD13	1.93	0.50
1:A:216:THR:HA	1:A:234:THR:HG21	1.94	0.50
1:C:265:LEU:HD22	1:C:268:ILE:HD11	1.92	0.50
1:A:370:LEU:HG	1:A:374:LYS:HE3	1.92	0.49
1:B:484:ILE:CG2	1:B:486:LEU:HG	2.42	0.49
1:A:273:GLN:HG2	3:A:646:HOH:O	2.12	0.49
1:B:502:TYR:HE1	1:B:550:LEU:HD13	1.71	0.49
1:C:219:LEU:HD21	1:C:232:ILE:HB	1.96	0.48
1:B:216:THR:HG22	1:B:234:THR:CB	2.43	0.48
1:B:126:MET:CE	1:B:138:ARG:HG3	2.44	0.47
1:C:10:PHE:CE2	1:C:14:GLN:HG3	2.49	0.47
1:B:216:THR:HG22	1:B:234:THR:HB	1.95	0.47
1:C:399:GLU:O	1:C:400:ILE:HD13	2.15	0.46
1:C:454:PRO:HD2	1:C:457:MET:SD	2.55	0.46
1:C:212:LYS:HE2	1:C:218:ASP:OD2	2.16	0.46
1:C:112:LYS:HE2	1:C:296:THR:O	2.16	0.46
1:A:222:LEU:HD23	1:B:207:LEU:HD13	1.98	0.46
1:A:96:SER:OG	1:A:338:LEU:HD21	2.16	0.45
1:C:380:LEU:HD11	1:C:411:ILE:HD11	1.98	0.45
1:A:429:TYR:CZ	1:A:516:LYS:HE2	2.52	0.45
1:C:462:GLN:HB2	1:C:484:ILE:HD11	1.98	0.45
1:C:484:ILE:HG13	1:C:484:ILE:H	1.71	0.44
1:C:209:TYR:HA	1:C:212:LYS:O	2.18	0.44
1:B:180:PRO:HB3	1:B:254:ALA:HA	1.99	0.44
1:A:53:LEU:HD22	1:A:63:ASP:OD1	2.18	0.44
1:A:115:ARG:HD3	1:A:372:THR:HA	2.00	0.44
1:A:338:LEU:HD11	3:A:654:HOH:O	2.18	0.44
1:B:484:ILE:HG22	1:B:486:LEU:HG	2.00	0.44
1:B:282:ILE:HD11	1:B:310:ILE:CD1	2.48	0.43
1:B:110:LYS:NZ	3:B:695:HOH:O	2.51	0.43
1:C:443:PRO:HG3	1:C:561:ARG:HH21	1.83	0.43
1:C:387:TRP:HH2	1:C:484:ILE:HG12	1.84	0.42
1:B:387:TRP:CH2	1:B:484:ILE:HG13	2.54	0.42
1:A:180:PRO:HG3	1:A:253:LYS:HG2	2.01	0.42
1:B:502:TYR:CZ	1:B:550:LEU:HD11	2.54	0.42
1:A:216:THR:CB	1:A:234:THR:HG21	2.50	0.42
1:A:180:PRO:HB3	1:A:254:ALA:HA	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:460:THR:CG2	1:B:484:ILE:HG12	2.50	0.42
1:A:112:LYS:HD2	1:A:304:GLU:OE2	2.19	0.42
1:B:484:ILE:HD13	1:B:484:ILE:HA	1.80	0.42
1:A:401:ALA:HA	1:A:411:ILE:O	2.20	0.41
1:A:29:LEU:HD23	1:A:29:LEU:C	2.41	0.41
1:B:487:SER:O	1:B:491:SER:HB2	2.20	0.41
1:C:32:VAL:O	1:C:36:GLN:HG3	2.21	0.41
1:A:282:ILE:HD11	1:A:310:ILE:CD1	2.51	0.41
1:A:291:MET:O	1:C:477:LYS:NZ	2.53	0.41
1:C:488:LYS:HD3	1:C:488:LYS:HA	1.80	0.41
1:C:53:LEU:HD22	1:C:63:ASP:OD1	2.20	0.41
1:B:221:ARG:HD3	1:B:221:ARG:HA	1.91	0.41
1:B:44:PHE:CG	1:B:189:LYS:HG2	2.56	0.41
1:B:8:LYS:HB2	1:B:9:SER:H	1.68	0.41
1:C:46:GLU:OE2	1:C:66:ARG:HD2	2.21	0.41
1:B:235:LYS:HG3	1:B:236:LYS:N	2.36	0.41
1:B:15:SER:CB	1:B:289:LEU:HD11	2.51	0.40
1:B:550:LEU:N	1:B:550:LEU:HD12	2.35	0.40
1:C:158:ASP:HB3	1:C:159:GLY:H	1.67	0.40
1:C:78:VAL:O	1:C:81:LYS:HE3	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	503/577 (87%)	496 (99%)	7 (1%)	0	100	100
1	B	507/577 (88%)	502 (99%)	5 (1%)	0	100	100
1	C	492/577 (85%)	483 (98%)	8 (2%)	1 (0%)	47	33
All	All	1502/1731 (87%)	1481 (99%)	20 (1%)	1 (0%)	51	36

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	550	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	450/498 (90%)	449 (100%)	1 (0%)	93	92
1	B	452/498 (91%)	448 (99%)	4 (1%)	78	75
1	C	444/498 (89%)	436 (98%)	8 (2%)	59	48
All	All	1346/1494 (90%)	1333 (99%)	13 (1%)	76	71

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	234	THR
1	B	8	LYS
1	B	158	ASP
1	B	464	SER
1	B	484	ILE
1	C	11	LEU
1	C	212	LYS
1	C	224	GLN
1	C	233	ASP
1	C	263	ASN
1	C	464	SER
1	C	473	SER
1	C	549	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	501	GLN
1	B	132	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 3 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

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	512/577 (88%)	0.75	33 (6%) 19 15	31, 33, 45, 66	0
1	B	515/577 (89%)	0.89	47 (9%) 9 7	31, 33, 45, 58	0
1	C	506/577 (87%)	0.97	65 (12%) 3 2	32, 36, 49, 65	0
All	All	1533/1731 (88%)	0.87	145 (9%) 8 6	31, 34, 47, 66	0

All (145) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	545	LEU	6.4
1	B	262	GLY	6.0
1	C	549	VAL	5.8
1	A	7	ILE	5.6
1	B	547	THR	5.3
1	B	338	LEU	5.2
1	C	439	PHE	5.1
1	C	260	ASP	5.1
1	C	556	ARG	5.0
1	A	549	VAL	4.8
1	C	365	LEU	4.7
1	B	549	VAL	4.6
1	B	147	SER	4.5
1	A	548	SER	4.3
1	B	546	ASN	4.2
1	A	234	THR	4.1
1	C	219	LEU	4.0
1	C	367	TYR	3.8
1	B	456	ASN	3.8
1	C	364	GLY	3.8
1	C	516	LYS	3.7
1	A	338	LEU	3.7
1	C	132	GLN	3.7

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Mol	Chain	Res	Type	RSRZ
1	C	158	ASP	3.6
1	B	548	SER	3.5
1	A	236	LYS	3.5
1	C	88	ILE	3.5
1	C	131	SER	3.4
1	B	260	ASP	3.4
1	A	484	ILE	3.4
1	A	545	LEU	3.3
1	A	213	TYR	3.2
1	C	465	ASP	3.2
1	C	224	GLN	3.2
1	B	484	ILE	3.2
1	C	458	VAL	3.2
1	C	23	TYR	3.2
1	C	473	SER	3.1
1	A	113	VAL	3.1
1	C	287	LYS	3.1
1	B	477	LYS	3.0
1	C	259	LEU	2.9
1	C	469	LYS	2.9
1	A	439	PHE	2.9
1	A	456	ASN	2.9
1	B	179	MET	2.9
1	C	539	ALA	2.8
1	C	336	VAL	2.7
1	B	64	LEU	2.7
1	A	8	LYS	2.7
1	B	401	ALA	2.7
1	C	184	LEU	2.7
1	B	263	ASN	2.7
1	C	223	THR	2.7
1	C	83	THR	2.7
1	B	47	VAL	2.6
1	C	268	ILE	2.6
1	C	429	TYR	2.6
1	B	176	PHE	2.6
1	C	424	LYS	2.6
1	C	40	HIS	2.6
1	C	234	THR	2.6
1	B	421	LYS	2.6
1	C	477	LYS	2.6
1	C	271	SER	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	427	ALA	2.5
1	C	467	ILE	2.5
1	B	172	LEU	2.5
1	B	27	ILE	2.5
1	C	459	ILE	2.5
1	C	93	THR	2.5
1	C	26	ASN	2.5
1	C	486	LEU	2.5
1	A	114	ILE	2.5
1	A	320	ILE	2.5
1	C	484	ILE	2.5
1	C	541	VAL	2.5
1	B	11	LEU	2.5
1	C	386	THR	2.4
1	B	182	LEU	2.4
1	B	106	LEU	2.4
1	B	376	ALA	2.4
1	C	100	LEU	2.4
1	C	466	ASP	2.4
1	C	457	MET	2.4
1	A	47	VAL	2.4
1	A	418	THR	2.4
1	C	504	ASP	2.4
1	C	237	SER	2.4
1	B	50	VAL	2.3
1	C	366	THR	2.3
1	C	21	SER	2.3
1	B	164	TRP	2.3
1	A	371	MET	2.3
1	A	485	ALA	2.3
1	C	107	GLU	2.3
1	C	233	ASP	2.3
1	C	91	VAL	2.3
1	C	387	TRP	2.3
1	B	373	LEU	2.3
1	B	122	ALA	2.3
1	B	393	ARG	2.3
1	C	455	ARG	2.2
1	A	237	SER	2.2
1	B	109	LEU	2.2
1	B	219	LEU	2.2
1	B	545	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	421	LYS	2.2
1	A	53	LEU	2.2
1	B	171	LEU	2.2
1	B	371	MET	2.2
1	B	243	GLY	2.2
1	C	262	GLY	2.2
1	C	263	ASN	2.2
1	A	109	LEU	2.2
1	A	219	LEU	2.2
1	B	370	LEU	2.2
1	C	337	ASP	2.2
1	B	516	LYS	2.2
1	A	158	ASP	2.2
1	A	235	LYS	2.2
1	C	24	CYS	2.2
1	A	74	VAL	2.2
1	B	378	LEU	2.2
1	B	35	ALA	2.2
1	B	517	LYS	2.1
1	B	465	ASP	2.1
1	A	331	TRP	2.1
1	A	337	ASP	2.1
1	B	114	ILE	2.1
1	A	261	GLY	2.1
1	A	477	LYS	2.1
1	B	236	LYS	2.1
1	C	456	ASN	2.1
1	B	184	LEU	2.1
1	B	235	LYS	2.1
1	C	470	LEU	2.1
1	A	42	LEU	2.0
1	A	262	GLY	2.0
1	C	225	SER	2.0
1	A	107	GLU	2.0
1	C	285	VAL	2.0
1	B	120	LEU	2.0
1	C	550	LEU	2.0
1	B	183	THR	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZN	B	570	1/1	0.77	0.52	103,103,103,103	0
2	ZN	C	570	1/1	0.98	0.10	44,44,44,44	0
2	ZN	A	570	1/1	0.99	0.10	33,33,33,33	0

6.5 Other polymers [i](#)

There are no such residues in this entry.