



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 07:49 AM UTC

PDB ID : 2BKC / pdb_00002bkc
Title : The X-ray structure of the H43G *Listeria innocua* Dps mutant
Authors : Ilari, A.; Stefanini, S.; Chiancone, E.
Deposited on : 2005-02-15
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
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

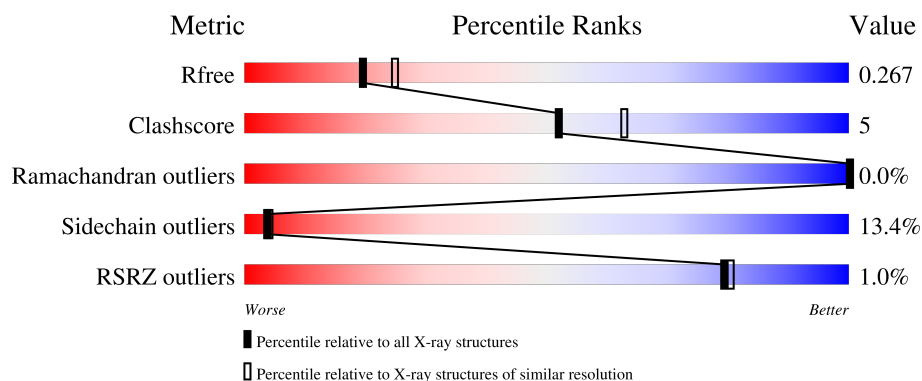
1 Overall quality at a glance

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.



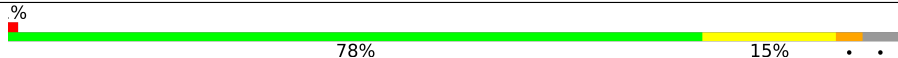
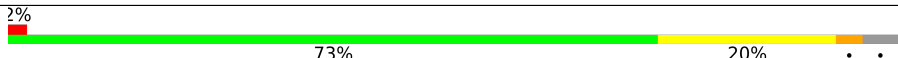
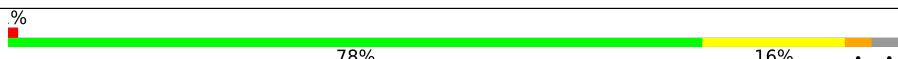
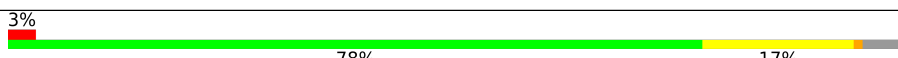
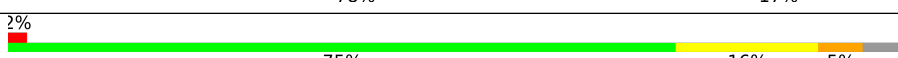
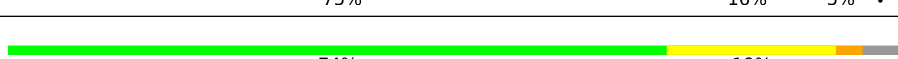
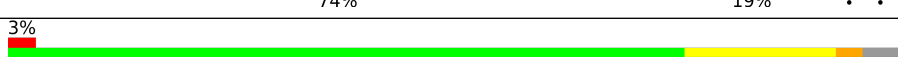
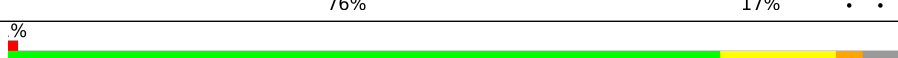
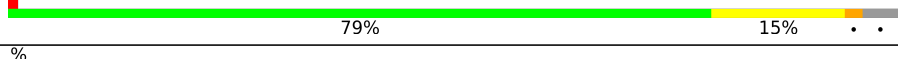
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	156	 80% 13% . .
1	B	156	 79% 13% . . .
1	C	156	 80% 12% . . .
1	D	156	 76% 19% . .
1	E	156	 77% 16% . .

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Mol	Chain	Length	Quality of chain
1	F	156	
1	G	156	
1	H	156	
1	I	156	
1	J	156	
1	K	156	
1	L	156	
1	M	156	
1	N	156	
1	O	156	
1	P	156	
1	Q	156	
1	R	156	
1	S	156	
1	T	156	
1	U	156	
1	V	156	
1	X	156	
1	Y	156	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 29380 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-HEME IRON-CONTAINING FERRITIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	B	150	Total	C	N	O	S	0	1	0
			1220	782	194	237	7			
1	C	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	D	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	E	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	F	150	Total	C	N	O	S	0	1	0
			1220	782	194	237	7			
1	G	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	H	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	I	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	J	150	Total	C	N	O	S	0	1	0
			1220	782	194	237	7			
1	K	150	Total	C	N	O	S	0	1	0
			1220	782	194	237	7			
1	L	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	M	150	Total	C	N	O	S	0	1	0
			1220	782	194	237	7			
1	N	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	O	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	P	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	R	149	Total	C	N	O	S	0	1	0
			1213	777	193	236	7			
1	S	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	T	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	U	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	V	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	X	150	Total	C	N	O	S	0	0	0
			1216	780	194	235	7			
1	Y	150	Total	C	N	O	S	0	0	0
			1213	778	193	235	7			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	43	GLY	HIS	engineered mutation	UNP P80725
B	43	GLY	HIS	engineered mutation	UNP P80725
C	43	GLY	HIS	engineered mutation	UNP P80725
D	43	GLY	HIS	engineered mutation	UNP P80725
E	43	GLY	HIS	engineered mutation	UNP P80725
F	43	GLY	HIS	engineered mutation	UNP P80725
G	43	GLY	HIS	engineered mutation	UNP P80725
H	43	GLY	HIS	engineered mutation	UNP P80725
I	43	GLY	HIS	engineered mutation	UNP P80725
J	43	GLY	HIS	engineered mutation	UNP P80725
K	43	GLY	HIS	engineered mutation	UNP P80725
L	43	GLY	HIS	engineered mutation	UNP P80725
M	43	GLY	HIS	engineered mutation	UNP P80725
N	43	GLY	HIS	engineered mutation	UNP P80725
O	43	GLY	HIS	engineered mutation	UNP P80725
P	43	GLY	HIS	engineered mutation	UNP P80725
Q	43	GLY	HIS	engineered mutation	UNP P80725
R	43	GLY	HIS	engineered mutation	UNP P80725
S	43	GLY	HIS	engineered mutation	UNP P80725
T	43	GLY	HIS	engineered mutation	UNP P80725
U	43	GLY	HIS	engineered mutation	UNP P80725
V	43	GLY	HIS	engineered mutation	UNP P80725
X	43	GLY	HIS	engineered mutation	UNP P80725

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Chain	Residue	Modelled	Actual	Comment	Reference
Y	43	GLY	HIS	engineered mutation	UNP P80725

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	12	Total O 12 12	0	0
2	B	8	Total O 8 8	0	0
2	C	14	Total O 14 14	0	0
2	D	9	Total O 9 9	0	0
2	E	10	Total O 10 10	0	0
2	F	7	Total O 7 7	0	0
2	G	6	Total O 6 6	0	0
2	H	10	Total O 10 10	0	0
2	I	7	Total O 7 7	0	0
2	J	7	Total O 7 7	0	0
2	K	16	Total O 16 16	0	0
2	L	7	Total O 7 7	0	0
2	M	9	Total O 9 9	0	0
2	N	5	Total O 5 5	0	0
2	O	5	Total O 5 5	0	0
2	P	3	Total O 3 3	0	0
2	Q	3	Total O 3 3	0	0
2	R	13	Total O 13 13	0	0
2	S	4	Total O 4 4	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	T	9	Total	O	0	0
			9	9		
2	U	4	Total	O	0	0
			4	4		
2	V	5	Total	O	0	0
			5	5		
2	X	5	Total	O	0	0
			5	5		
2	Y	4	Total	O	0	0
			4	4		

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-HEME IRON-CONTAINING FERRITIN

Chain A: 



• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain B: 



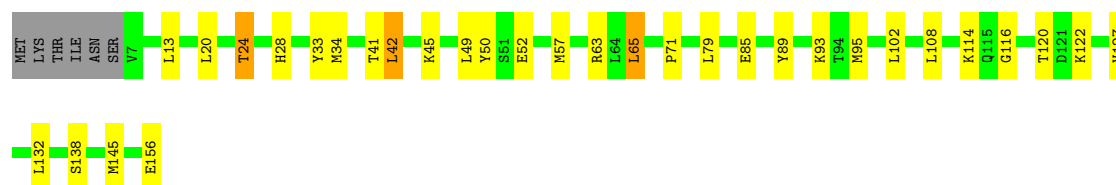
• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain C: 



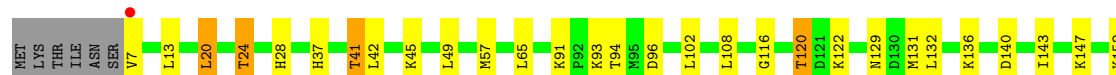
• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain D: 



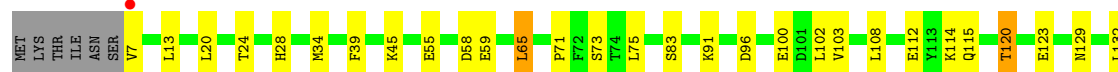
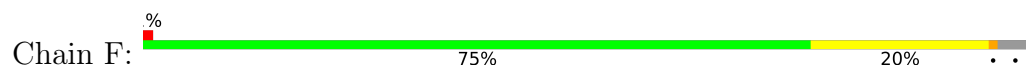
• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain E: 

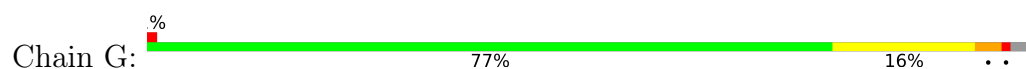




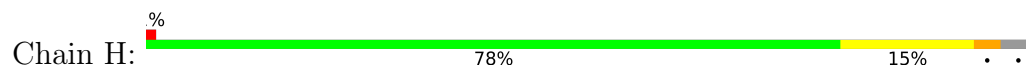
• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



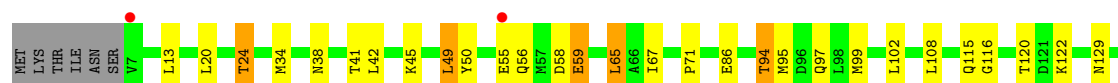
• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



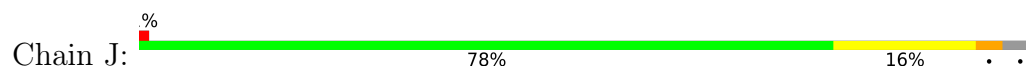
• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

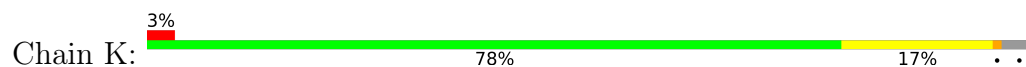


• Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

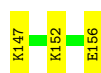
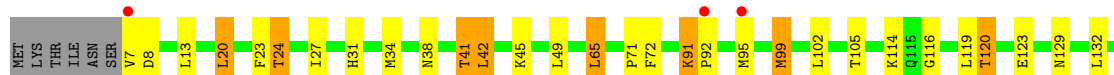
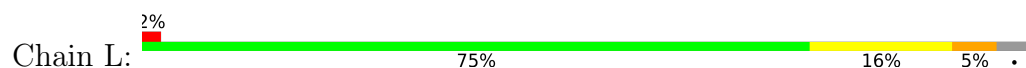




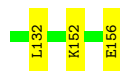
- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



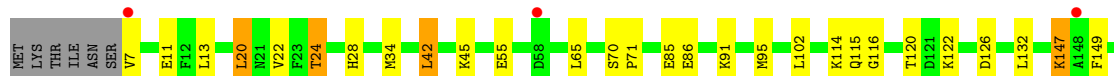
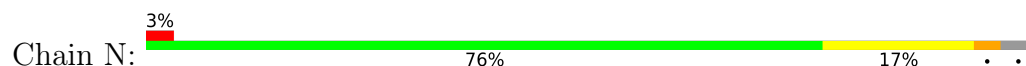
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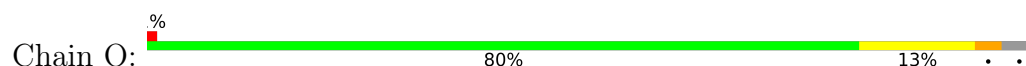
- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



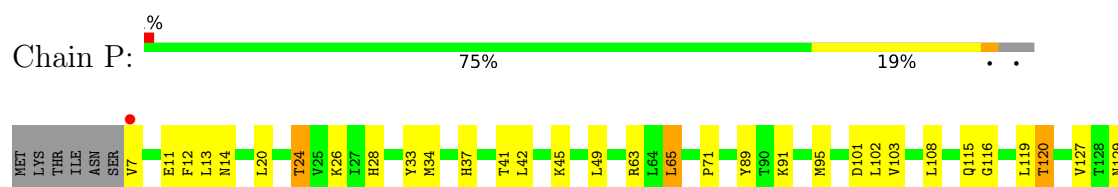
- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



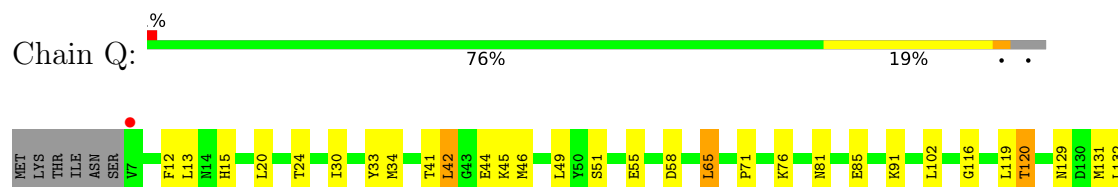
- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



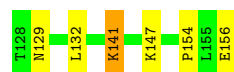
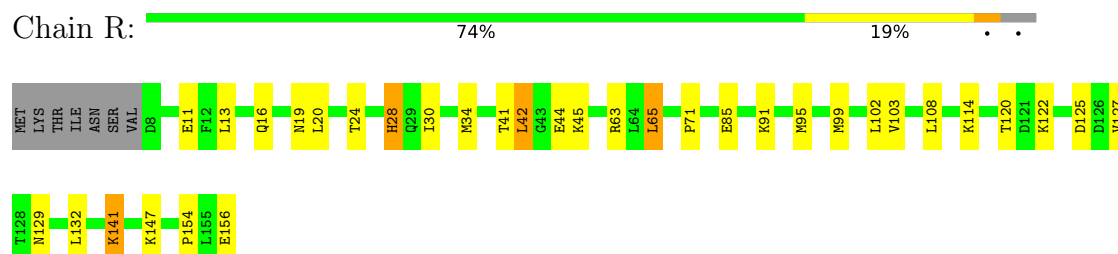
- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



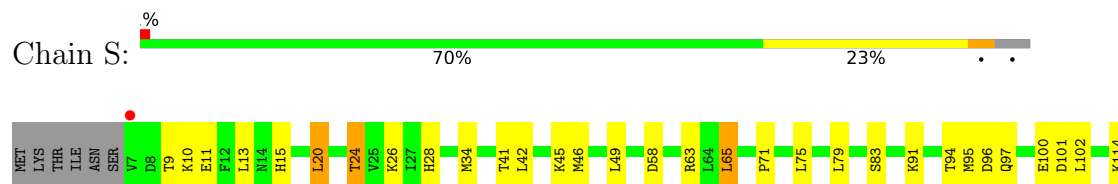
- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



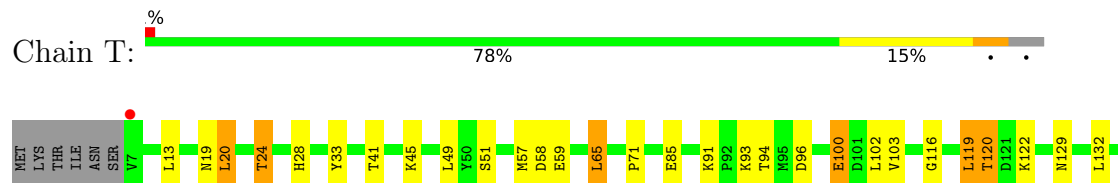
- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

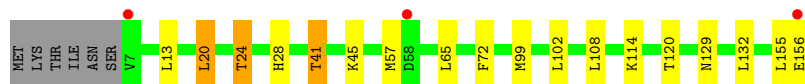


- Molecule 1: NON-HEME IRON-CONTAINING FERRITIN



● Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain U:  2% 85% 10% . .



● Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain V:  % 78% 15% . .



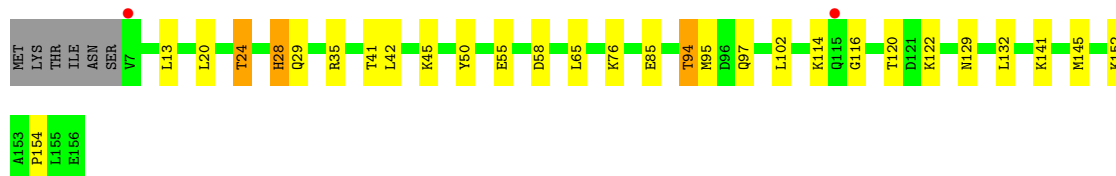
● Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain X:  % 79% 15% . .



● Molecule 1: NON-HEME IRON-CONTAINING FERRITIN

Chain Y:  % 78% 17% . .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	90.73Å 175.65Å 135.96Å 90.00° 92.53° 90.00°	Depositor
Resolution (Å)	25.00 – 2.30 25.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	91.5 (25.00-2.30) 91.4 (25.00-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.209 , 0.266 0.212 , 0.267	Depositor DCC
R_{free} test set	8694 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	27.6	Xtriage
Anisotropy	0.406	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 22.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.025 for h,-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	29380	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.63% 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	1.07	2/1242 (0.2%)	1.12	0/1674
1	B	1.04	1/1250 (0.1%)	1.11	1/1685 (0.1%)
1	C	1.04	0/1242	1.10	1/1674 (0.1%)
1	D	1.08	1/1242 (0.1%)	1.05	0/1674
1	E	1.03	1/1242 (0.1%)	1.08	0/1674
1	F	1.06	1/1250 (0.1%)	1.06	1/1685 (0.1%)
1	G	1.00	1/1242 (0.1%)	1.05	0/1674
1	H	1.05	0/1242	1.11	1/1674 (0.1%)
1	I	1.10	2/1242 (0.2%)	1.10	1/1674 (0.1%)
1	J	0.99	0/1250	1.10	1/1685 (0.1%)
1	K	1.06	0/1250	1.15	3/1685 (0.2%)
1	L	1.03	1/1242 (0.1%)	1.12	1/1674 (0.1%)
1	M	1.00	0/1250	1.11	0/1685
1	N	1.00	0/1242	1.07	2/1674 (0.1%)
1	O	1.04	2/1242 (0.2%)	1.09	0/1674
1	P	0.96	1/1242 (0.1%)	1.08	0/1674
1	Q	1.01	1/1242 (0.1%)	1.09	0/1674
1	R	1.01	1/1243 (0.1%)	1.08	0/1675
1	S	1.02	0/1242	1.08	0/1674
1	T	0.99	0/1242	1.09	1/1674 (0.1%)
1	U	1.05	2/1242 (0.2%)	1.08	1/1674 (0.1%)
1	V	1.00	2/1242 (0.2%)	1.07	0/1674
1	X	0.97	0/1242	1.08	0/1674
1	Y	0.99	0/1239	1.05	1/1671 (0.1%)
All	All	1.03	19/29846 (0.1%)	1.09	15/40229 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	J	0	1

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	99	MET	SD-CE	-7.93	1.59	1.79
1	I	145	MET	SD-CE	-7.06	1.61	1.79
1	O	46	MET	SD-CE	-6.50	1.63	1.79
1	F	134	ALA	CA-CB	6.14	1.62	1.53
1	I	99	MET	SD-CE	-5.96	1.64	1.79
1	Q	131	MET	SD-CE	-5.78	1.65	1.79
1	V	145	MET	SD-CE	5.71	1.93	1.79
1	U	99	MET	SD-CE	-5.71	1.65	1.79
1	U	41	THR	CA-CB	5.67	1.62	1.53
1	B	99	MET	SD-CE	-5.39	1.66	1.79
1	R	141	LYS	CD-CE	5.30	1.68	1.52
1	A	70	SER	C-O	-5.29	1.18	1.24
1	L	99	MET	SD-CE	-5.17	1.66	1.79
1	O	95	MET	CG-SD	5.15	1.93	1.80
1	E	131	MET	CA-C	5.15	1.59	1.52
1	P	24	THR	CA-CB	5.15	1.61	1.53
1	G	24	THR	CA-CB	5.14	1.61	1.53
1	D	145	MET	SD-CE	-5.07	1.66	1.79
1	V	46	MET	SD-CE	5.05	1.92	1.79

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	91	LYS	CB-CA-C	6.52	118.25	108.86
1	H	55	GLU	N-CA-CB	6.20	119.00	110.01
1	L	72	PHE	N-CA-C	-6.13	102.01	110.35
1	N	147	LYS	N-CA-C	6.00	117.90	111.36
1	U	72	PHE	N-CA-C	-5.98	101.55	110.23
1	B	120	THR	N-CA-CB	-5.75	101.47	110.44
1	F	73	SER	N-CA-C	5.68	117.72	110.61
1	N	22	VAL	CB-CA-C	-5.59	104.74	112.24
1	C	120	THR	N-CA-CB	-5.53	101.79	110.30
1	I	67	ILE	N-CA-C	5.28	117.25	112.29
1	K	91	LYS	CA-C-N	5.24	125.00	119.76
1	K	91	LYS	C-N-CA	5.24	125.00	119.76
1	Y	122	LYS	CB-CA-C	-5.15	102.74	110.92
1	T	19	ASN	N-CA-C	5.11	117.24	111.11
1	J	105	THR	OG1-CB-CG2	-5.02	99.26	109.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	J	7	VAL	Peptide

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	1216	0	1182	10	0
1	B	1220	0	1182	10	0
1	C	1216	0	1182	12	0
1	D	1216	0	1182	12	0
1	E	1216	0	1182	17	0
1	F	1220	0	1182	11	0
1	G	1216	0	1182	14	0
1	H	1216	0	1182	18	0
1	I	1216	0	1182	18	0
1	J	1220	0	1182	8	0
1	K	1220	0	1182	18	0
1	L	1216	0	1182	22	0
1	M	1220	0	1182	14	0
1	N	1216	0	1182	8	0
1	O	1216	0	1182	14	0
1	P	1216	0	1182	14	0
1	Q	1216	0	1182	12	0
1	R	1213	0	1173	15	0
1	S	1216	0	1182	19	0
1	T	1216	0	1182	12	0
1	U	1216	0	1182	9	0
1	V	1216	0	1182	12	0
1	X	1216	0	1182	8	0
1	Y	1213	0	1173	13	0
2	A	12	0	0	2	0
2	B	8	0	0	0	0
2	C	14	0	0	0	0
2	D	9	0	0	0	0
2	E	10	0	0	0	0
2	F	7	0	0	0	0
2	G	6	0	0	0	0
2	H	10	0	0	0	0
2	I	7	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	J	7	0	0	0	0
2	K	16	0	0	0	0
2	L	7	0	0	0	0
2	M	9	0	0	0	0
2	N	5	0	0	0	0
2	O	5	0	0	0	0
2	P	3	0	0	0	0
2	Q	3	0	0	0	0
2	R	13	0	0	0	0
2	S	4	0	0	0	0
2	T	9	0	0	0	0
2	U	4	0	0	0	0
2	V	5	0	0	0	0
2	X	5	0	0	0	0
2	Y	4	0	0	0	0
All	All	29380	0	28350	295	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 5.

All (295) 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:F:120:THR:HG22	1:F:129:ASN:HB2	1.55	0.87
1:A:37:HIS:NE2	1:K:37:HIS:ND1	2.21	0.86
1:D:42:LEU:HD23	1:D:95:MET:SD	2.17	0.85
1:C:116:GLY:O	1:C:120:THR:HB	1.80	0.81
1:L:34:MET:HB2	1:L:95:MET:SD	2.21	0.80
1:Q:116:GLY:O	1:Q:120:THR:HB	1.81	0.80
1:I:94:THR:H	1:I:97:GLN:HE21	1.30	0.80
1:A:37:HIS:HE2	1:K:37:HIS:HD1	0.83	0.80
1:I:38:ASN:HB3	1:I:42:LEU:HD23	1.63	0.80
1:T:116:GLY:O	1:T:120:THR:HG22	1.82	0.78
1:G:38:ASN:HB3	1:G:42:LEU:HD23	1.65	0.77
1:A:37:HIS:CE1	1:K:37:HIS:HD1	2.05	0.75
1:V:116:GLY:O	1:V:120:THR:HB	1.88	0.74
1:L:120:THR:CG2	1:L:129:ASN:HB2	2.16	0.74
1:C:100:GLU:OE1	1:C:152:LYS:NZ	2.21	0.73
1:H:28:HIS:CD2	1:H:50:TYR:CE2	2.76	0.73
1:T:119:LEU:C	1:T:119:LEU:HD23	2.14	0.72
2:A:2003:HOH:O	1:K:37:HIS:HE1	1.72	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:R:120:THR:HG22	1:R:129:ASN:HB2	1.72	0.71
1:P:116:GLY:O	1:P:120:THR:HB	1.92	0.70
1:S:118:GLU:O	1:S:122:LYS:HG2	1.90	0.70
1:I:94:THR:HG22	1:I:97:GLN:H	1.57	0.70
1:S:26:LYS:NZ	1:S:101:ASP:OD2	2.22	0.69
1:Q:120:THR:HG23	1:Q:129:ASN:HB2	1.75	0.69
1:O:116:GLY:O	1:O:120:THR:HB	1.92	0.69
1:L:31:HIS:HA	1:L:34:MET:HE3	1.76	0.68
1:T:119:LEU:HD23	1:T:120:THR:N	2.09	0.68
1:L:120:THR:HG21	1:L:129:ASN:CA	2.24	0.67
1:R:147:LYS:NZ	1:R:154:PRO:O	2.26	0.67
1:Y:116:GLY:O	1:Y:120:THR:HG22	1.95	0.67
1:N:147:LYS:NZ	1:N:156:GLU:HB2	2.10	0.67
1:F:120:THR:CG2	1:F:129:ASN:HB2	2.25	0.67
1:H:116:GLY:O	1:H:120:THR:HB	1.95	0.66
1:E:116:GLY:O	1:E:120:THR:HB	1.96	0.65
1:K:94:THR:H	1:K:97:GLN:HE21	1.44	0.65
1:G:116:GLY:O	1:G:120:THR:HB	1.95	0.65
1:O:120:THR:HG23	1:O:129:ASN:HB2	1.79	0.65
1:P:120:THR:HG22	1:P:129:ASN:HD22	1.60	0.65
1:C:42:LEU:HD23	1:C:95:MET:SD	2.37	0.64
1:E:37:HIS:HB3	1:L:38:ASN:HD21	1.63	0.64
1:X:120:THR:HG22	1:X:129:ASN:HB2	1.79	0.64
1:A:141:LYS:HG2	1:A:145:MET:HE2	1.78	0.64
1:B:116:GLY:O	1:B:120:THR:HB	1.97	0.63
1:X:120:THR:HG21	1:X:129:ASN:N	2.13	0.63
1:V:34:MET:HB2	1:V:95:MET:HE1	1.80	0.63
1:T:20:LEU:O	1:T:24:THR:HB	1.98	0.63
1:H:34:MET:HG3	1:H:95:MET:HE1	1.82	0.62
1:H:120:THR:HG23	1:H:129:ASN:HB2	1.83	0.60
1:G:120:THR:HG21	1:G:129:ASN:HA	1.84	0.60
1:U:120:THR:HG21	1:U:129:ASN:HA	1.84	0.60
1:Y:120:THR:HG23	1:Y:129:ASN:HD22	1.67	0.59
1:C:120:THR:HG23	1:C:129:ASN:HB2	1.84	0.59
1:O:34:MET:HB2	1:O:95:MET:SD	2.43	0.59
1:E:120:THR:HG22	1:E:129:ASN:HD22	1.69	0.58
1:C:37:HIS:HE1	1:E:96:ASP:OD1	1.86	0.58
1:R:65:LEU:HD13	1:R:71:PRO:HD3	1.85	0.58
1:I:24:THR:HG23	1:I:50:TYR:CE2	2.39	0.58
1:C:120:THR:HG22	1:C:129:ASN:HD22	1.68	0.58
1:E:120:THR:HG21	1:E:129:ASN:HA	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:83:SER:OG	1:J:112:GLU:OE2	2.22	0.57
1:S:94:THR:HG23	1:S:97:GLN:H	1.67	0.57
1:F:96:ASP:O	1:F:100:GLU:HG3	2.05	0.57
1:O:120:THR:CG2	1:O:129:ASN:HD22	2.18	0.57
1:H:120:THR:HG21	1:H:129:ASN:HA	1.85	0.57
1:I:116:GLY:O	1:I:120:THR:HG22	2.05	0.57
1:L:120:THR:HG22	1:L:129:ASN:HB2	1.84	0.57
1:S:65:LEU:HD13	1:S:71:PRO:HD3	1.86	0.57
1:E:147:LYS:NZ	1:E:156:GLU:O	2.37	0.56
1:Y:120:THR:HG21	1:Y:129:ASN:HA	1.87	0.56
1:N:147:LYS:HZ3	1:N:156:GLU:HB2	1.69	0.56
1:H:34:MET:CG	1:H:95:MET:HE1	2.36	0.56
1:H:30:ILE:O	1:H:34:MET:HB3	2.05	0.56
1:L:116:GLY:O	1:L:120:THR:HB	2.06	0.56
1:F:65:LEU:HD13	1:F:71:PRO:HD3	1.88	0.55
1:T:65:LEU:HD13	1:T:71:PRO:HD3	1.87	0.55
1:P:34:MET:HB2	1:P:95:MET:HE1	1.89	0.55
1:S:63:ARG:NH2	1:S:127:VAL:HB	2.21	0.55
1:Y:94:THR:HG22	1:Y:97:GLN:HG3	1.88	0.55
1:T:96:ASP:O	1:T:100:GLU:HG3	2.08	0.54
1:N:116:GLY:O	1:N:120:THR:HG23	2.07	0.54
1:A:38:ASN:HD21	1:K:37:HIS:HB3	1.72	0.54
1:H:28:HIS:NE2	1:H:50:TYR:CZ	2.75	0.54
1:L:7:VAL:HG23	1:L:119:LEU:HD22	1.90	0.53
1:L:65:LEU:HD13	1:L:71:PRO:HD3	1.89	0.53
1:C:20:LEU:O	1:C:24:THR:HB	2.08	0.53
1:H:94:THR:HB	1:H:97:GLN:H	1.73	0.53
1:H:74:THR:OG1	1:H:77:GLU:OE1	2.19	0.53
1:J:108:LEU:C	1:J:108:LEU:HD23	2.34	0.53
1:M:76:LYS:O	1:M:80:GLU:HG3	2.08	0.53
1:G:120:THR:HG23	1:G:129:ASN:HB2	1.91	0.53
1:L:120:THR:HG21	1:L:129:ASN:HA	1.89	0.53
1:C:120:THR:HG21	1:C:129:ASN:HA	1.89	0.53
1:M:127:VAL:HG13	1:Q:137:ALA:HB2	1.89	0.53
1:P:120:THR:HG21	1:P:129:ASN:HA	1.91	0.53
1:I:155:LEU:O	1:I:156:GLU:C	2.51	0.53
1:E:41:THR:CG2	1:L:41:THR:HG21	2.39	0.52
1:H:120:THR:HG22	1:H:129:ASN:HD22	1.74	0.52
1:I:120:THR:HG21	1:I:129:ASN:HA	1.92	0.52
1:O:34:MET:HB2	1:O:95:MET:HE1	1.91	0.52
1:S:120:THR:HG22	1:S:129:ASN:HB2	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:147:LYS:HG2	1:S:153:ALA:O	2.10	0.52
1:J:20:LEU:O	1:J:24:THR:HB	2.10	0.51
1:S:120:THR:HG23	1:S:125:ASP:HB3	1.93	0.51
1:F:108:LEU:C	1:F:108:LEU:HD23	2.35	0.51
1:J:120:THR:HG23	1:J:129:ASN:HB2	1.93	0.51
1:V:120:THR:HG21	1:V:129:ASN:HA	1.92	0.51
1:E:41:THR:HG21	1:L:41:THR:HG21	1.93	0.51
1:L:120:THR:HG21	1:L:129:ASN:N	2.26	0.51
1:M:42:LEU:HD23	1:M:95:MET:SD	2.51	0.51
1:O:120:THR:HG21	1:O:129:ASN:HA	1.93	0.51
1:U:20:LEU:HD23	1:U:57:MET:HA	1.93	0.51
1:X:42:LEU:HD23	1:X:95:MET:SD	2.51	0.51
1:D:108:LEU:C	1:D:108:LEU:HD23	2.36	0.50
1:Y:94:THR:HG22	1:Y:97:GLN:H	1.76	0.50
1:M:15:HIS:HE1	1:M:112:GLU:OE1	1.93	0.50
1:P:14:ASN:OD1	1:P:71:PRO:HA	2.12	0.50
1:O:38:ASN:HB3	1:O:42:LEU:HD23	1.93	0.50
1:Q:42:LEU:HD21	1:Q:149:PHE:CZ	2.46	0.50
1:E:108:LEU:HD23	1:E:108:LEU:C	2.37	0.50
1:F:58[A]:ASP:OD2	1:H:28:HIS:HE1	1.95	0.50
1:K:120:THR:HG23	1:K:129:ASN:HB2	1.94	0.50
1:P:37:HIS:HE1	1:S:96:ASP:OD1	1.94	0.50
1:X:15:HIS:HE1	1:X:83:SER:OG	1.95	0.50
1:L:20:LEU:O	1:L:24:THR:HB	2.12	0.49
1:C:42:LEU:CD2	1:C:95:MET:SD	3.01	0.49
1:F:83:SER:OG	1:F:112:GLU:OE2	2.24	0.49
1:L:120:THR:CG2	1:L:129:ASN:CB	2.90	0.49
1:S:120:THR:HG21	1:S:129:ASN:N	2.27	0.49
1:K:155:LEU:O	1:K:156:GLU:HB2	2.12	0.49
1:Q:30:ILE:HG22	1:Q:34:MET:HE3	1.95	0.49
2:A:2003:HOH:O	1:K:37:HIS:CE1	2.55	0.49
1:L:99:MET:HE2	1:L:99:MET:HA	1.95	0.49
1:V:120:THR:HG23	1:V:129:ASN:HB2	1.95	0.48
1:V:57:MET:HE1	1:Y:28:HIS:CD2	2.49	0.48
1:M:94:THR:HG22	1:M:96:ASP:H	1.78	0.48
1:R:63:ARG:NH2	1:R:127:VAL:HB	2.28	0.48
1:U:108:LEU:C	1:U:108:LEU:HD23	2.37	0.48
1:O:20:LEU:O	1:O:24:THR:HB	2.14	0.48
1:D:28:HIS:CD2	1:D:50:TYR:CE2	3.01	0.48
1:D:63:ARG:NH2	1:D:127:VAL:HB	2.29	0.48
1:F:120:THR:HG21	1:F:129:ASN:N	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:Y:141:LYS:HG2	1:Y:145:MET:HE2	1.95	0.48
1:D:34:MET:HB2	1:D:95:MET:HE1	1.96	0.48
1:A:95:MET:SD	1:A:95:MET:C	2.96	0.48
1:E:120:THR:CG2	1:E:129:ASN:HD22	2.27	0.48
1:H:65:LEU:HD13	1:H:71:PRO:HD3	1.95	0.48
1:T:20:LEU:HD23	1:T:57:MET:HA	1.96	0.47
1:U:120:THR:HG23	1:U:129:ASN:HD22	1.78	0.47
1:K:94:THR:H	1:K:97:GLN:NE2	2.11	0.47
1:T:65:LEU:HD13	1:T:71:PRO:CD	2.44	0.47
1:E:37:HIS:HB3	1:L:38:ASN:ND2	2.28	0.47
1:S:143:ILE:O	1:S:147:LYS:HB2	2.15	0.47
1:V:73:SER:HB2	1:Y:29:GLN:HB2	1.97	0.47
1:D:65:LEU:HD13	1:D:71:PRO:HD3	1.97	0.47
1:H:52:GLU:O	1:H:55:GLU:HG3	2.15	0.47
1:L:42:LEU:HD12	1:L:42:LEU:HA	1.81	0.47
1:U:57:MET:HE1	1:X:28:HIS:CD2	2.50	0.46
1:X:120:THR:HG23	1:X:125:ASP:HB3	1.97	0.46
1:P:108:LEU:C	1:P:108:LEU:HD23	2.40	0.46
1:M:76:LYS:O	1:M:80:GLU:CG	2.63	0.46
1:J:52:GLU:OE2	1:J:138:SER:OG	2.20	0.46
1:P:33:TYR:CD2	1:P:89:TYR:CD2	3.03	0.46
1:U:120:THR:CG2	1:U:129:ASN:HD22	2.29	0.46
1:L:91:LYS:O	1:L:92:PRO:C	2.58	0.46
1:O:42:LEU:O	1:O:46:MET:HG2	2.16	0.46
1:R:120:THR:HG21	1:R:129:ASN:N	2.29	0.46
1:E:120:THR:HG23	1:E:129:ASN:HB2	1.98	0.46
1:K:15:HIS:HE1	1:K:83:SER:OG	1.99	0.46
1:M:63:ARG:NH2	1:M:127:VAL:HB	2.30	0.46
1:N:42:LEU:HD21	1:N:149:PHE:CZ	2.51	0.46
1:N:71:PRO:O	1:P:33:TYR:OH	2.30	0.46
1:I:55:GLU:OE1	1:I:59:GLU:OE1	2.33	0.46
1:L:95:MET:O	1:L:99:MET:HG2	2.16	0.46
1:Q:120:THR:CG2	1:Q:129:ASN:HB2	2.45	0.46
1:B:83:SER:OG	1:B:112:GLU:OE2	2.25	0.46
1:G:120:THR:HG22	1:G:129:ASN:HD22	1.81	0.46
1:R:34:MET:HE1	1:R:42:LEU:HB3	1.97	0.46
1:S:42:LEU:N	1:S:42:LEU:HD12	2.30	0.46
1:I:56:GLN:HG3	1:I:135:PHE:CE2	2.51	0.45
1:A:65:LEU:HD13	1:A:71:PRO:CD	2.47	0.45
1:B:24:THR:HG23	1:B:50:TYR:CE2	2.51	0.45
1:I:49:LEU:HD11	1:I:142:HIS:ND1	2.30	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:138:SER:O	1:I:142:HIS:HD2	1.99	0.45
1:O:120:THR:HG22	1:O:129:ASN:HD22	1.81	0.45
1:M:15:HIS:CE1	1:M:112:GLU:OE1	2.68	0.45
1:B:120:THR:HG23	1:B:129:ASN:HB2	1.97	0.45
1:H:12:PHE:CD2	1:H:12:PHE:C	2.94	0.45
1:I:156:GLU:OE1	1:I:156:GLU:HA	2.14	0.45
1:P:26:LYS:NZ	1:P:101:ASP:OD2	2.33	0.45
1:S:34:MET:HB2	1:S:95:MET:HE1	1.97	0.45
1:O:34:MET:HB2	1:O:95:MET:CE	2.46	0.45
1:K:94:THR:N	1:K:97:GLN:HE21	2.10	0.45
1:Q:120:THR:HG21	1:Q:129:ASN:HA	1.99	0.45
1:R:71:PRO:O	1:T:33:TYR:OH	2.34	0.45
1:D:79:LEU:HD23	1:D:79:LEU:HA	1.89	0.45
1:L:23:PHE:CD1	1:L:105:THR:HG21	2.51	0.45
1:P:65:LEU:O	1:R:154:PRO:HG2	2.17	0.45
1:J:33:TYR:OH	1:L:71:PRO:O	2.31	0.44
1:Q:12:PHE:CD2	1:Q:12:PHE:C	2.95	0.44
1:A:24:THR:HG23	1:A:50:TYR:CE2	2.53	0.44
1:O:120:THR:HG23	1:O:129:ASN:CB	2.46	0.44
1:G:65:LEU:HD13	1:G:71:PRO:HD3	1.99	0.44
1:X:155:LEU:O	1:X:156:GLU:HB2	2.18	0.44
1:M:20:LEU:O	1:M:24:THR:HB	2.17	0.44
1:G:120:THR:CG2	1:G:129:ASN:HA	2.48	0.44
1:K:21:ASN:CG	1:K:57:MET:HE2	2.42	0.44
1:E:20:LEU:HD23	1:E:57:MET:HA	1.99	0.44
1:V:46:MET:HE1	1:V:102:LEU:HD11	2.00	0.44
1:D:116:GLY:O	1:D:120:THR:HG22	2.18	0.43
1:M:108:LEU:C	1:M:108:LEU:HD23	2.43	0.43
1:R:30:ILE:HG22	1:R:34:MET:HE3	2.00	0.43
1:U:155:LEU:O	1:U:156:GLU:C	2.60	0.43
1:V:120:THR:CG2	1:V:129:ASN:HA	2.47	0.43
1:I:108:LEU:C	1:I:108:LEU:HD23	2.44	0.43
1:P:63:ARG:NH2	1:P:127:VAL:HB	2.33	0.43
1:S:75:LEU:N	1:S:75:LEU:HD12	2.33	0.43
1:M:42:LEU:HD12	1:M:42:LEU:HA	1.85	0.43
1:I:65:LEU:HD13	1:I:71:PRO:HD3	2.00	0.43
1:V:120:THR:HG22	1:V:129:ASN:HD22	1.83	0.43
1:Y:24:THR:HG23	1:Y:50:TYR:CE2	2.53	0.43
1:B:108:LEU:HD23	1:B:108:LEU:C	2.44	0.43
1:P:12:PHE:CZ	1:P:116:GLY:HA3	2.54	0.43
1:S:42:LEU:O	1:S:46:MET:HG2	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:120:THR:HG22	1:B:129:ASN:HD22	1.84	0.43
1:E:120:THR:CG2	1:E:129:ASN:HA	2.47	0.43
1:U:20:LEU:O	1:U:24:THR:HB	2.19	0.43
1:R:16:GLN:HA	1:R:19:ASN:HB2	2.01	0.43
1:Y:35:ARG:O	1:Y:95:MET:HE2	2.19	0.43
1:D:33:TYR:CD2	1:D:89:TYR:CD2	3.07	0.42
1:J:112:GLU:HA	1:J:115:GLN:HE21	1.84	0.42
1:O:56:GLN:O	1:O:60:VAL:HG23	2.19	0.42
1:I:120:THR:HG23	1:I:129:ASN:HB2	2.01	0.42
1:T:96:ASP:O	1:T:100:GLU:CG	2.67	0.42
1:F:141:LYS:O	1:F:145:MET:HG3	2.19	0.42
1:K:94:THR:HB	1:K:97:GLN:HE21	1.84	0.42
1:M:94:THR:HG22	1:M:96:ASP:N	2.33	0.42
1:I:34:MET:HB2	1:I:95:MET:HE1	2.01	0.42
1:P:120:THR:HG23	1:P:129:ASN:HB2	2.01	0.42
1:X:42:LEU:O	1:X:46:MET:HG2	2.20	0.42
1:A:20:LEU:HD23	1:A:57:MET:HA	2.01	0.42
1:G:65:LEU:HD13	1:G:71:PRO:CD	2.49	0.42
1:H:120:THR:HG21	1:H:129:ASN:CA	2.50	0.42
1:J:28:HIS:CD2	1:J:50:TYR:CE2	3.08	0.42
1:Y:120:THR:CG2	1:Y:129:ASN:HD22	2.32	0.42
1:G:42:LEU:HD21	1:G:149:PHE:CE1	2.54	0.42
1:S:20:LEU:O	1:S:24:THR:HB	2.19	0.42
1:D:52:GLU:OE2	1:D:138:SER:OG	2.23	0.42
1:F:39:PHE:CD2	1:F:39:PHE:C	2.97	0.42
1:K:42:LEU:HD12	1:K:42:LEU:HA	1.82	0.42
1:N:20:LEU:O	1:N:24:THR:HB	2.20	0.42
1:N:155:LEU:O	1:N:156:GLU:C	2.62	0.42
1:O:120:THR:HG23	1:O:129:ASN:HD22	1.85	0.42
1:D:24:THR:HG23	1:D:50:TYR:CE2	2.55	0.42
1:E:20:LEU:O	1:E:24:THR:HB	2.19	0.42
1:G:75:LEU:N	1:G:75:LEU:HD12	2.34	0.42
1:H:28:HIS:NE2	1:H:50:TYR:CE2	2.87	0.42
1:R:120:THR:HG23	1:R:125:ASP:HB3	2.01	0.42
1:G:20:LEU:O	1:G:24:THR:HB	2.20	0.41
1:U:120:THR:HG23	1:U:129:ASN:HB2	2.02	0.41
1:B:42:LEU:HA	1:B:45:LYS:HB2	2.02	0.41
1:C:37:HIS:CE1	1:E:96:ASP:OD1	2.69	0.41
1:I:108:LEU:HD23	1:I:108:LEU:O	2.20	0.41
1:R:34:MET:HE2	1:R:95:MET:SD	2.59	0.41
1:S:65:LEU:O	1:Y:154:PRO:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:86:GLU:CD	1:K:74:THR:HB	2.45	0.41
1:Q:42:LEU:O	1:Q:46:MET:HG2	2.20	0.41
1:M:23:PHE:CD1	1:M:105:THR:HG21	2.56	0.41
1:R:108:LEU:C	1:R:108:LEU:HD23	2.45	0.41
1:G:42:LEU:N	1:G:42:LEU:HD22	2.35	0.41
1:N:34:MET:HB2	1:N:95:MET:SD	2.61	0.41
1:Q:15:HIS:CD2	1:Q:81:ASN:O	2.73	0.41
1:R:99:MET:O	1:R:103:VAL:HG23	2.21	0.41
1:V:38:ASN:HB3	1:V:42:LEU:HD23	2.03	0.41
1:V:65:LEU:HD13	1:V:71:PRO:HD3	2.03	0.41
1:Y:24:THR:HG23	1:Y:50:TYR:HE2	1.85	0.41
1:G:15:HIS:HE1	1:G:83:SER:OG	2.03	0.41
1:K:94:THR:HB	1:K:97:GLN:NE2	2.35	0.41
1:M:45:LYS:HZ3	1:M:45:LYS:HG3	1.74	0.41
1:E:136:LYS:NZ	1:E:140:ASP:OD2	2.45	0.41
1:Q:65:LEU:HD13	1:Q:71:PRO:HD3	2.03	0.41
1:R:28:HIS:CD2	1:T:57:MET:HE1	2.55	0.41
1:B:49:LEU:HD12	1:B:49:LEU:HA	1.91	0.41
1:T:120:THR:HG23	1:T:129:ASN:HD22	1.84	0.40
1:B:52:GLU:OE2	1:B:138:SER:OG	2.24	0.40
1:C:155:LEU:O	1:C:156:GLU:C	2.64	0.40
1:G:34:MET:HE2	1:G:95:MET:SD	2.61	0.40
1:S:15:HIS:HE1	1:S:83:SER:OG	2.04	0.40
1:V:42:LEU:HD12	1:V:99:MET:SD	2.60	0.40
1:B:28:HIS:CD2	1:D:57:MET:HE1	2.56	0.40
1:Q:33:TYR:OH	1:S:71:PRO:O	2.28	0.40
1:C:15:HIS:HE1	1:C:83:SER:OG	2.04	0.40
1:A:38:ASN:HB3	1:A:42:LEU:HD22	2.03	0.40
1:F:75:LEU:N	1:F:75:LEU:HD12	2.37	0.40
1:H:120:THR:CG2	1:H:129:ASN:HD22	2.35	0.40
1:K:116:GLY:O	1:K:120:THR:HB	2.22	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	148/156 (95%)	147 (99%)	1 (1%)	0	100	100
1	B	149/156 (96%)	147 (99%)	2 (1%)	0	100	100
1	C	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	D	148/156 (95%)	147 (99%)	1 (1%)	0	100	100
1	E	148/156 (95%)	145 (98%)	3 (2%)	0	100	100
1	F	149/156 (96%)	147 (99%)	2 (1%)	0	100	100
1	G	148/156 (95%)	145 (98%)	2 (1%)	1 (1%)	18	23
1	H	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	I	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	J	149/156 (96%)	145 (97%)	4 (3%)	0	100	100
1	K	149/156 (96%)	146 (98%)	3 (2%)	0	100	100
1	L	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	M	149/156 (96%)	147 (99%)	2 (1%)	0	100	100
1	N	148/156 (95%)	144 (97%)	4 (3%)	0	100	100
1	O	148/156 (95%)	145 (98%)	3 (2%)	0	100	100
1	P	148/156 (95%)	144 (97%)	4 (3%)	0	100	100
1	Q	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	R	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	S	148/156 (95%)	144 (97%)	4 (3%)	0	100	100
1	T	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	U	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	V	148/156 (95%)	144 (97%)	4 (3%)	0	100	100
1	X	148/156 (95%)	146 (99%)	2 (1%)	0	100	100
1	Y	148/156 (95%)	142 (96%)	6 (4%)	0	100	100
All	All	3557/3744 (95%)	3493 (98%)	63 (2%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	92	PRO

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	132/138 (96%)	117 (89%)	15 (11%)	5	6
1	B	133/138 (96%)	117 (88%)	16 (12%)	5	5
1	C	132/138 (96%)	115 (87%)	17 (13%)	4	4
1	D	132/138 (96%)	117 (89%)	15 (11%)	5	6
1	E	132/138 (96%)	112 (85%)	20 (15%)	3	3
1	F	133/138 (96%)	114 (86%)	19 (14%)	3	3
1	G	132/138 (96%)	113 (86%)	19 (14%)	3	3
1	H	132/138 (96%)	116 (88%)	16 (12%)	5	5
1	I	132/138 (96%)	118 (89%)	14 (11%)	6	8
1	J	133/138 (96%)	115 (86%)	18 (14%)	4	4
1	K	133/138 (96%)	118 (89%)	15 (11%)	5	7
1	L	132/138 (96%)	113 (86%)	19 (14%)	3	3
1	M	133/138 (96%)	110 (83%)	23 (17%)	2	2
1	N	132/138 (96%)	110 (83%)	22 (17%)	2	2
1	O	132/138 (96%)	116 (88%)	16 (12%)	5	5
1	P	132/138 (96%)	114 (86%)	18 (14%)	3	4
1	Q	132/138 (96%)	113 (86%)	19 (14%)	3	3
1	R	132/138 (96%)	114 (86%)	18 (14%)	3	4
1	S	132/138 (96%)	112 (85%)	20 (15%)	3	3
1	T	132/138 (96%)	110 (83%)	22 (17%)	2	2
1	U	132/138 (96%)	122 (92%)	10 (8%)	12	16
1	V	132/138 (96%)	116 (88%)	16 (12%)	5	5
1	X	132/138 (96%)	113 (86%)	19 (14%)	3	3
1	Y	131/138 (95%)	114 (87%)	17 (13%)	4	4
All	All	3172/3312 (96%)	2749 (87%)	423 (13%)	4	4

All (423) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	20	LEU
1	A	24	THR
1	A	28	HIS
1	A	42	LEU
1	A	44	GLU
1	A	45	LYS
1	A	55	GLU
1	A	65	LEU
1	A	96	ASP
1	A	102	LEU
1	A	114	LYS
1	A	132	LEU
1	A	152	LYS
1	A	156	GLU
1	B	7	VAL
1	B	13	LEU
1	B	20	LEU
1	B	24	THR
1	B	41	THR
1	B	42	LEU
1	B	44	GLU
1	B	45	LYS
1	B	49	LEU
1	B	65	LEU
1	B	91	LYS
1	B	102	LEU
1	B	114	LYS
1	B	119	LEU
1	B	120	THR
1	B	132	LEU
1	C	13	LEU
1	C	20	LEU
1	C	24	THR
1	C	34	MET
1	C	42	LEU
1	C	45	LYS
1	C	55	GLU
1	C	60	VAL
1	C	65	LEU
1	C	76	LYS
1	C	96	ASP
1	C	100	GLU

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Mol	Chain	Res	Type
1	C	102	LEU
1	C	120	THR
1	C	122	LYS
1	C	132	LEU
1	C	156	GLU
1	D	13	LEU
1	D	20	LEU
1	D	24	THR
1	D	41	THR
1	D	42	LEU
1	D	45	LYS
1	D	49	LEU
1	D	65	LEU
1	D	85	GLU
1	D	93	LYS
1	D	102	LEU
1	D	114	LYS
1	D	122	LYS
1	D	132	LEU
1	D	156	GLU
1	E	7	VAL
1	E	13	LEU
1	E	20	LEU
1	E	24	THR
1	E	28	HIS
1	E	41	THR
1	E	42	LEU
1	E	45	LYS
1	E	49	LEU
1	E	65	LEU
1	E	91	LYS
1	E	93	LYS
1	E	94	THR
1	E	102	LEU
1	E	120	THR
1	E	122	LYS
1	E	132	LEU
1	E	143	ILE
1	E	152	LYS
1	E	156	GLU
1	F	7	VAL
1	F	13	LEU

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Mol	Chain	Res	Type
1	F	20	LEU
1	F	24	THR
1	F	28	HIS
1	F	34	MET
1	F	45	LYS
1	F	55	GLU
1	F	59	GLU
1	F	65	LEU
1	F	91	LYS
1	F	102	LEU
1	F	103	VAL
1	F	114	LYS
1	F	115	GLN
1	F	120	THR
1	F	123	GLU
1	F	132	LEU
1	F	156	GLU
1	G	13	LEU
1	G	20	LEU
1	G	24	THR
1	G	28	HIS
1	G	34	MET
1	G	41	THR
1	G	45	LYS
1	G	49	LEU
1	G	58	ASP
1	G	65	LEU
1	G	76	LYS
1	G	93	LYS
1	G	100	GLU
1	G	102	LEU
1	G	103	VAL
1	G	114	LYS
1	G	120	THR
1	G	132	LEU
1	G	156	GLU
1	H	7	VAL
1	H	13	LEU
1	H	20	LEU
1	H	24	THR
1	H	34	MET
1	H	45	LYS

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Mol	Chain	Res	Type
1	H	51	SER
1	H	65	LEU
1	H	93	LYS
1	H	94	THR
1	H	102	LEU
1	H	114	LYS
1	H	120	THR
1	H	132	LEU
1	H	141	LYS
1	H	152	LYS
1	I	13	LEU
1	I	20	LEU
1	I	24	THR
1	I	41	THR
1	I	45	LYS
1	I	49	LEU
1	I	58	ASP
1	I	59	GLU
1	I	65	LEU
1	I	94	THR
1	I	102	LEU
1	I	115	GLN
1	I	122	LYS
1	I	132	LEU
1	J	7	VAL
1	J	11	GLU
1	J	13	LEU
1	J	20	LEU
1	J	24	THR
1	J	28	HIS
1	J	41	THR
1	J	42	LEU
1	J	44	GLU
1	J	45	LYS
1	J	65	LEU
1	J	76	LYS
1	J	90	THR
1	J	94	THR
1	J	102	LEU
1	J	132	LEU
1	J	141	LYS
1	J	147	LYS

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Mol	Chain	Res	Type
1	K	13	LEU
1	K	20	LEU
1	K	24	THR
1	K	41	THR
1	K	42	LEU
1	K	45	LYS
1	K	49	LEU
1	K	65	LEU
1	K	93	LYS
1	K	94	THR
1	K	100	GLU
1	K	102	LEU
1	K	118	GLU
1	K	119	LEU
1	K	132	LEU
1	L	8	ASP
1	L	13	LEU
1	L	20	LEU
1	L	24	THR
1	L	27	ILE
1	L	41	THR
1	L	42	LEU
1	L	45	LYS
1	L	49	LEU
1	L	65	LEU
1	L	91	LYS
1	L	102	LEU
1	L	114	LYS
1	L	120	THR
1	L	123	GLU
1	L	132	LEU
1	L	147	LYS
1	L	152	LYS
1	L	156	GLU
1	M	7	VAL
1	M	10	LYS
1	M	13	LEU
1	M	20	LEU
1	M	24	THR
1	M	41	THR
1	M	42	LEU
1	M	45	LYS

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Mol	Chain	Res	Type
1	M	49	LEU
1	M	59	GLU
1	M	65	LEU
1	M	80	GLU
1	M	91	LYS
1	M	93	LYS
1	M	100	GLU
1	M	102	LEU
1	M	103	VAL
1	M	114	LYS
1	M	118	GLU
1	M	119	LEU
1	M	132	LEU
1	M	152	LYS
1	M	156	GLU
1	N	7	VAL
1	N	11	GLU
1	N	13	LEU
1	N	20	LEU
1	N	24	THR
1	N	28	HIS
1	N	42	LEU
1	N	45	LYS
1	N	55	GLU
1	N	65	LEU
1	N	70	SER
1	N	85	GLU
1	N	86	GLU
1	N	91	LYS
1	N	102	LEU
1	N	114	LYS
1	N	115	GLN
1	N	122	LYS
1	N	126	ASP
1	N	132	LEU
1	N	152	LYS
1	N	155	LEU
1	O	7	VAL
1	O	11	GLU
1	O	13	LEU
1	O	20	LEU
1	O	24	THR

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Mol	Chain	Res	Type
1	O	41	THR
1	O	49	LEU
1	O	65	LEU
1	O	85	GLU
1	O	93	LYS
1	O	102	LEU
1	O	114	LYS
1	O	119	LEU
1	O	120	THR
1	O	132	LEU
1	O	141	LYS
1	P	7	VAL
1	P	11	GLU
1	P	13	LEU
1	P	20	LEU
1	P	24	THR
1	P	28	HIS
1	P	41	THR
1	P	42	LEU
1	P	45	LYS
1	P	49	LEU
1	P	65	LEU
1	P	91	LYS
1	P	102	LEU
1	P	103	VAL
1	P	115	GLN
1	P	119	LEU
1	P	120	THR
1	P	132	LEU
1	Q	13	LEU
1	Q	20	LEU
1	Q	24	THR
1	Q	41	THR
1	Q	42	LEU
1	Q	44	GLU
1	Q	45	LYS
1	Q	49	LEU
1	Q	51	SER
1	Q	55	GLU
1	Q	58	ASP
1	Q	65	LEU
1	Q	76	LYS

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Mol	Chain	Res	Type
1	Q	85	GLU
1	Q	91	LYS
1	Q	102	LEU
1	Q	119	LEU
1	Q	120	THR
1	Q	132	LEU
1	R	11	GLU
1	R	13	LEU
1	R	20	LEU
1	R	24	THR
1	R	28	HIS
1	R	41	THR
1	R	42	LEU
1	R	44	GLU
1	R	45	LYS
1	R	65	LEU
1	R	85	GLU
1	R	91	LYS
1	R	102	LEU
1	R	114	LYS
1	R	122	LYS
1	R	132	LEU
1	R	141	LYS
1	R	156	GLU
1	S	9	THR
1	S	10	LYS
1	S	11	GLU
1	S	13	LEU
1	S	20	LEU
1	S	24	THR
1	S	28	HIS
1	S	41	THR
1	S	45	LYS
1	S	49	LEU
1	S	58	ASP
1	S	65	LEU
1	S	79	LEU
1	S	91	LYS
1	S	100	GLU
1	S	102	LEU
1	S	114	LYS
1	S	122	LYS

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Mol	Chain	Res	Type
1	S	132	LEU
1	S	147	LYS
1	T	13	LEU
1	T	20	LEU
1	T	24	THR
1	T	28	HIS
1	T	41	THR
1	T	45	LYS
1	T	49	LEU
1	T	51	SER
1	T	58	ASP
1	T	59	GLU
1	T	65	LEU
1	T	85	GLU
1	T	91	LYS
1	T	93	LYS
1	T	94	THR
1	T	100	GLU
1	T	102	LEU
1	T	103	VAL
1	T	119	LEU
1	T	120	THR
1	T	122	LYS
1	T	132	LEU
1	U	13	LEU
1	U	20	LEU
1	U	24	THR
1	U	28	HIS
1	U	41	THR
1	U	45	LYS
1	U	65	LEU
1	U	102	LEU
1	U	114	LYS
1	U	132	LEU
1	V	7	VAL
1	V	11	GLU
1	V	20	LEU
1	V	24	THR
1	V	28	HIS
1	V	41	THR
1	V	44	GLU
1	V	45	LYS

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Mol	Chain	Res	Type
1	V	49	LEU
1	V	58	ASP
1	V	65	LEU
1	V	85	GLU
1	V	91	LYS
1	V	102	LEU
1	V	120	THR
1	V	132	LEU
1	X	7	VAL
1	X	13	LEU
1	X	20	LEU
1	X	24	THR
1	X	28	HIS
1	X	34	MET
1	X	41	THR
1	X	42	LEU
1	X	45	LYS
1	X	65	LEU
1	X	100	GLU
1	X	102	LEU
1	X	103	VAL
1	X	108	LEU
1	X	114	LYS
1	X	119	LEU
1	X	120	THR
1	X	122	LYS
1	X	132	LEU
1	Y	13	LEU
1	Y	20	LEU
1	Y	24	THR
1	Y	28	HIS
1	Y	41	THR
1	Y	42	LEU
1	Y	45	LYS
1	Y	55	GLU
1	Y	58	ASP
1	Y	65	LEU
1	Y	76	LYS
1	Y	85	GLU
1	Y	94	THR
1	Y	102	LEU
1	Y	114	LYS

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Mol	Chain	Res	Type
1	Y	132	LEU
1	Y	152	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (83) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	38	ASN
1	A	115	GLN
1	B	15	HIS
1	B	21	ASN
1	B	28	HIS
1	B	115	GLN
1	B	129	ASN
1	C	15	HIS
1	C	38	ASN
1	C	81	ASN
1	C	129	ASN
1	D	15	HIS
1	E	15	HIS
1	E	81	ASN
1	E	115	GLN
1	E	129	ASN
1	F	15	HIS
1	F	29	GLN
1	G	15	HIS
1	G	28	HIS
1	G	29	GLN
1	G	56	GLN
1	G	81	ASN
1	G	97	GLN
1	G	115	GLN
1	G	129	ASN
1	H	15	HIS
1	H	81	ASN
1	H	129	ASN
1	I	15	HIS
1	I	97	GLN
1	I	142	HIS
1	J	38	ASN
1	J	56	GLN
1	J	81	ASN
1	J	115	GLN

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Mol	Chain	Res	Type
1	K	15	HIS
1	K	81	ASN
1	K	97	GLN
1	K	115	GLN
1	K	129	ASN
1	L	15	HIS
1	L	29	GLN
1	L	38	ASN
1	M	81	ASN
1	N	81	ASN
1	O	29	GLN
1	O	115	GLN
1	O	129	ASN
1	P	15	HIS
1	P	31	HIS
1	P	37	HIS
1	P	129	ASN
1	Q	15	HIS
1	Q	81	ASN
1	R	15	HIS
1	R	29	GLN
1	R	129	ASN
1	S	15	HIS
1	S	29	GLN
1	S	81	ASN
1	T	15	HIS
1	T	29	GLN
1	T	115	GLN
1	T	129	ASN
1	U	15	HIS
1	U	37	HIS
1	U	38	ASN
1	U	81	ASN
1	U	97	GLN
1	U	129	ASN
1	V	15	HIS
1	V	29	GLN
1	V	129	ASN
1	X	15	HIS
1	X	29	GLN
1	X	56	GLN
1	Y	15	HIS

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Mol	Chain	Res	Type
1	Y	28	HIS
1	Y	29	GLN
1	Y	38	ASN
1	Y	81	ASN
1	Y	129	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	150/156 (96%)	-0.11	1 (0%) 84 85	20, 27, 38, 49	0
1	B	150/156 (96%)	-0.07	1 (0%) 84 85	18, 28, 39, 53	1 (0%)
1	C	150/156 (96%)	-0.10	1 (0%) 84 85	22, 28, 40, 52	0
1	D	150/156 (96%)	-0.16	0 100 100	22, 28, 40, 50	0
1	E	150/156 (96%)	-0.12	1 (0%) 84 85	22, 28, 39, 49	0
1	F	150/156 (96%)	-0.03	2 (1%) 75 76	18, 28, 40, 52	1 (0%)
1	G	150/156 (96%)	-0.09	1 (0%) 84 85	21, 29, 40, 54	0
1	H	150/156 (96%)	-0.05	1 (0%) 84 85	22, 28, 41, 50	0
1	I	150/156 (96%)	-0.08	3 (2%) 65 66	21, 28, 40, 52	0
1	J	150/156 (96%)	-0.07	1 (0%) 84 85	18, 29, 42, 52	1 (0%)
1	K	150/156 (96%)	-0.13	4 (2%) 56 58	17, 27, 37, 52	1 (0%)
1	L	150/156 (96%)	0.04	3 (2%) 65 66	23, 29, 43, 50	0
1	M	150/156 (96%)	-0.07	0 100 100	17, 28, 41, 52	1 (0%)
1	N	150/156 (96%)	0.17	4 (2%) 56 58	25, 31, 41, 53	0
1	O	150/156 (96%)	-0.04	1 (0%) 84 85	23, 29, 42, 53	0
1	P	150/156 (96%)	0.10	1 (0%) 84 85	24, 31, 42, 52	0
1	Q	150/156 (96%)	-0.05	1 (0%) 84 85	23, 29, 41, 51	0
1	R	149/156 (95%)	0.04	0 100 100	18, 29, 41, 51	1 (0%)
1	S	150/156 (96%)	0.08	1 (0%) 84 85	25, 31, 42, 50	0
1	T	150/156 (96%)	-0.07	1 (0%) 84 85	24, 29, 40, 50	0
1	U	150/156 (96%)	0.01	3 (2%) 65 66	23, 30, 41, 51	0
1	V	150/156 (96%)	0.03	1 (0%) 84 85	24, 30, 42, 51	0
1	X	150/156 (96%)	0.01	2 (1%) 75 76	23, 29, 41, 54	0
1	Y	150/156 (96%)	0.17	2 (1%) 75 76	25, 31, 42, 51	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
All	All	3599/3744 (96%)	-0.03	36 (1%) 79 80	17, 29, 42, 54	6 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	L	7	VAL	7.1
1	U	7	VAL	4.6
1	F	7	VAL	4.1
1	Y	7	VAL	4.0
1	C	7	VAL	3.9
1	J	7	VAL	3.8
1	E	7	VAL	3.6
1	H	7	VAL	3.1
1	V	7	VAL	2.8
1	N	156	GLU	2.7
1	I	7	VAL	2.7
1	Q	7	VAL	2.6
1	I	156	GLU	2.6
1	N	58	ASP	2.6
1	S	7	VAL	2.5
1	N	7	VAL	2.5
1	N	148	ALA	2.5
1	P	7	VAL	2.4
1	B	156	GLU	2.4
1	K	156	GLU	2.4
1	X	92	PRO	2.4
1	T	7	VAL	2.4
1	U	58	ASP	2.3
1	O	95	MET	2.3
1	F	156	GLU	2.3
1	K	93	LYS	2.3
1	G	156	GLU	2.2
1	I	55	GLU	2.2
1	K	118	GLU	2.2
1	Y	115	GLN	2.2
1	L	92	PRO	2.2
1	L	95	MET	2.1
1	A	91	LYS	2.1
1	U	156	GLU	2.0
1	K	91	LYS	2.0
1	X	122	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.