



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:57 AM UTC

PDB ID : 4CMY / pdb_00004cmy
Title : Chlorobium tepidum Ferritin
Authors : Pohl, E.; Arenas, M.; Townsend, P.D.; Yevenes, A.
Deposited on : 2014-01-18
Resolution : 2.59 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

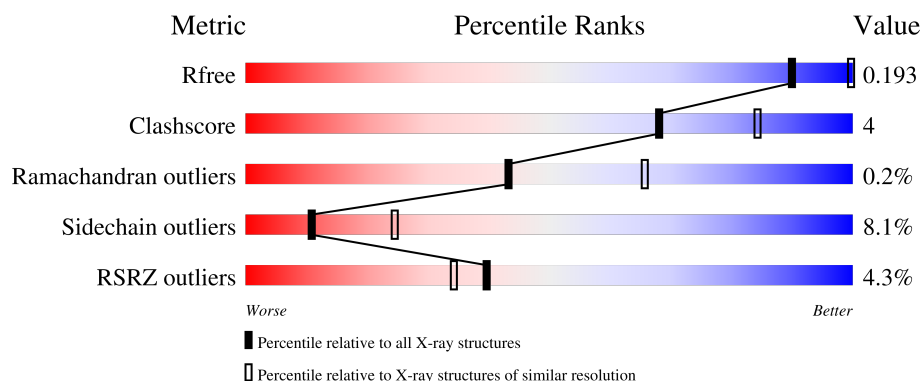
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.59 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.



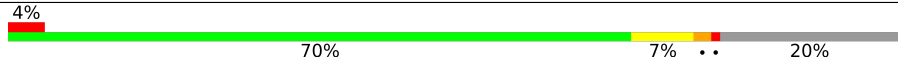
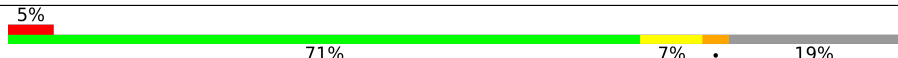
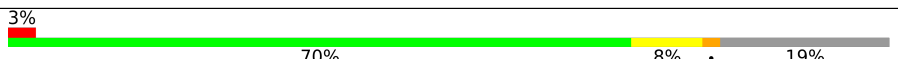
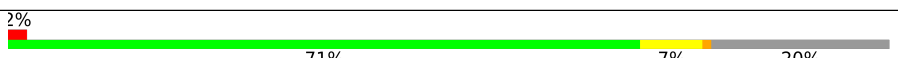
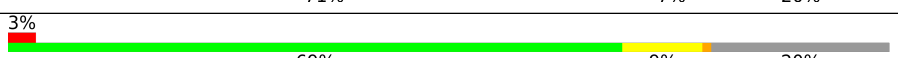
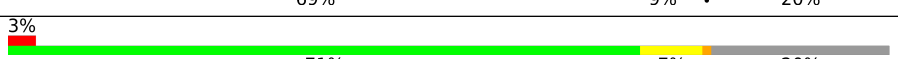
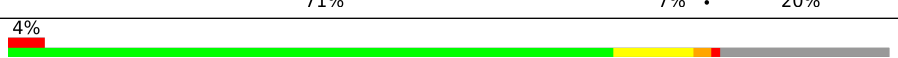
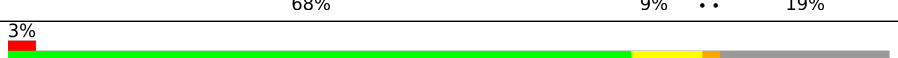
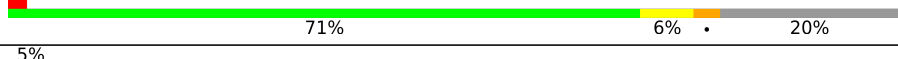
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	203	
1	B	203	
1	C	203	
1	D	203	
1	E	203	

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Mol	Chain	Length	Quality of chain
1	F	203	
1	G	203	
1	H	203	
1	I	203	
1	J	203	
1	K	203	
1	L	203	
1	M	203	
1	N	203	
1	O	203	
1	P	203	
1	Q	203	
1	R	203	
1	S	203	
1	T	203	
1	U	203	
1	W	203	
1	X	203	
2	V	203	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 32293 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 FERRITIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	163	Total	C	N	O	S	0	0	0
			1318	842	220	250	6			
1	B	164	Total	C	N	O	S	0	2	0
			1335	854	223	251	7			
1	C	163	Total	C	N	O	S	0	0	0
			1318	842	220	250	6			
1	D	163	Total	C	N	O	S	0	0	0
			1318	842	220	250	6			
1	E	163	Total	C	N	O	S	0	0	0
			1318	842	220	250	6			
1	F	163	Total	C	N	O	S	0	1	0
			1325	847	222	250	6			
1	G	163	Total	C	N	O	S	0	0	0
			1318	842	220	250	6			
1	H	163	Total	C	N	O	S	0	0	0
			1318	842	220	250	6			
1	I	164	Total	C	N	O	S	0	0	0
			1323	845	221	251	6			
1	J	164	Total	C	N	O	S	0	1	0
			1330	850	223	251	6			
1	K	163	Total	C	N	O	S	0	1	0
			1325	847	222	250	6			
1	L	163	Total	C	N	O	S	0	1	0
			1325	847	222	250	6			
1	M	163	Total	C	N	O	S	0	1	0
			1325	847	222	250	6			
1	N	164	Total	C	N	O	S	0	0	0
			1323	845	221	251	6			
1	O	164	Total	C	N	O	S	0	1	0
			1330	850	223	251	6			
1	P	163	Total	C	N	O	S	0	1	0
			1322	845	220	251	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	163	Total	C	N	O	S	0	1	0
			1325	847	222	250	6			
1	R	163	Total	C	N	O	S	0	2	0
			1329	850	222	251	6			
1	S	162	Total	C	N	O	S	0	0	0
			1313	839	219	249	6			
1	T	163	Total	C	N	O	S	0	2	0
			1333	852	223	251	7			
1	U	164	Total	C	N	O	S	0	1	0
			1330	850	223	251	6			
1	W	164	Total	C	N	O	S	0	1	0
			1331	850	222	252	7			
1	X	163	Total	C	N	O	S	0	1	0
			1325	847	222	250	6			

- Molecule 2 is a protein called FERRITIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	V	163	Total	C	N	O	S	0	1	0
			1329	850	222	251	6			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
V	145	ASP	ASN	conflict	UNP Q8KBP5

- Molecule 3 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Fe	0	0
			2	2		
3	B	2	Total	Fe	0	0
			2	2		
3	C	2	Total	Fe	0	0
			2	2		
3	D	2	Total	Fe	0	0
			2	2		
3	E	2	Total	Fe	0	0
			2	2		
3	F	2	Total	Fe	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	2	Total 2	Fe 2	0	0
3	H	2	Total 2	Fe 2	0	0
3	I	2	Total 2	Fe 2	0	0
3	J	2	Total 2	Fe 2	0	0
3	K	2	Total 2	Fe 2	0	0
3	L	2	Total 2	Fe 2	0	0
3	M	2	Total 2	Fe 2	0	0
3	N	2	Total 2	Fe 2	0	0
3	O	2	Total 2	Fe 2	0	0
3	P	2	Total 2	Fe 2	0	0
3	Q	2	Total 2	Fe 2	0	0
3	R	2	Total 2	Fe 2	0	0
3	S	2	Total 2	Fe 2	0	0
3	T	2	Total 2	Fe 2	0	0
3	U	2	Total 2	Fe 2	0	0
3	V	2	Total 2	Fe 2	0	0
3	W	2	Total 2	Fe 2	0	0
3	X	2	Total 2	Fe 2	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	31	Total 31	O 31	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	16	Total O 16 16	0	0
4	C	30	Total O 30 30	0	0
4	D	17	Total O 17 17	0	0
4	E	20	Total O 20 20	0	0
4	F	6	Total O 6 6	0	0
4	G	14	Total O 14 14	0	0
4	H	16	Total O 16 16	0	0
4	I	18	Total O 18 18	0	0
4	J	29	Total O 29 29	0	0
4	K	33	Total O 33 33	0	0
4	L	23	Total O 23 23	0	0
4	M	10	Total O 10 10	0	0
4	N	20	Total O 20 20	0	0
4	O	10	Total O 10 10	0	0
4	P	10	Total O 10 10	0	0
4	Q	11	Total O 11 11	0	0
4	R	13	Total O 13 13	0	0
4	S	17	Total O 17 17	0	0
4	T	13	Total O 13 13	0	0
4	U	30	Total O 30 30	0	0
4	V	26	Total O 26 26	0	0

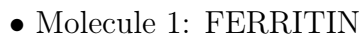
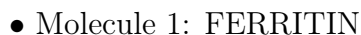
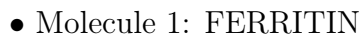
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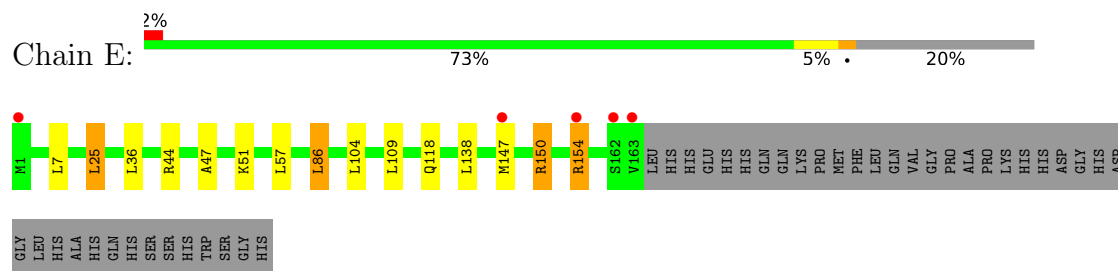
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	W	25	Total	O	0	0
			25	25		
4	X	21	Total	O	0	0
			21	21		

i

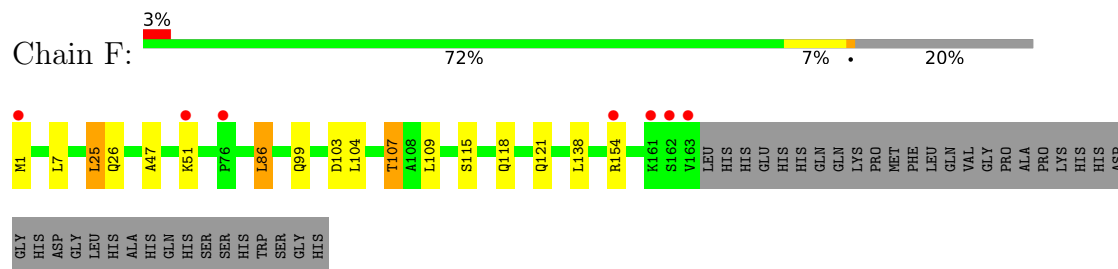
- Molecule 1: FERRITIN



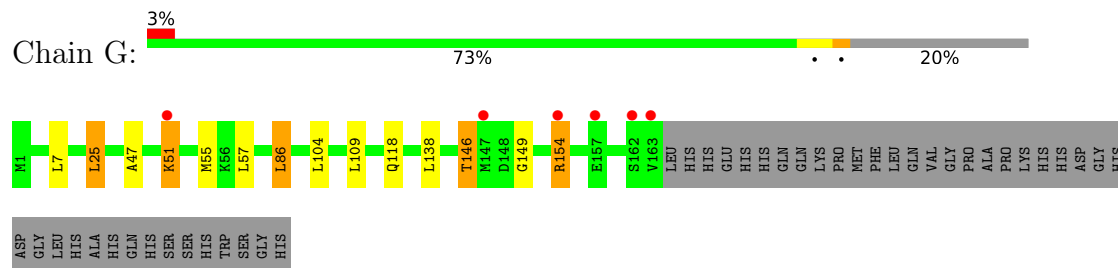
- Molecule 1: FERRITIN



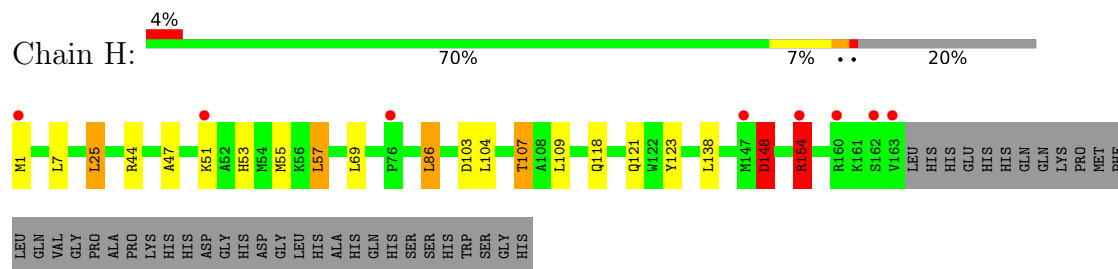
- Molecule 1: FERRITIN



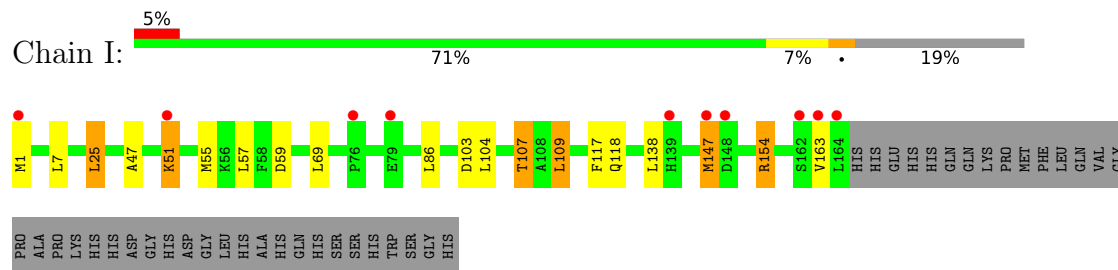
- Molecule 1: FERRITIN



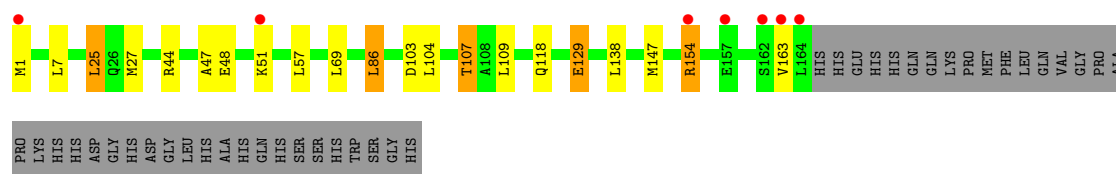
- Molecule 1: FERRITIN



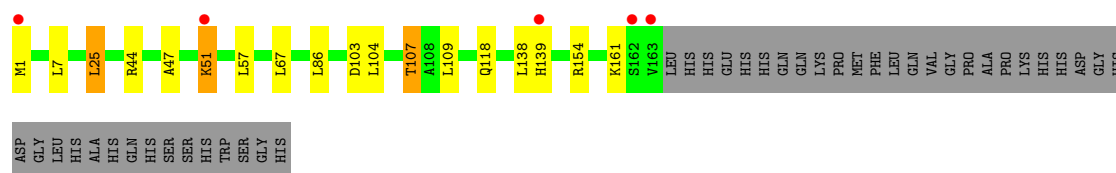
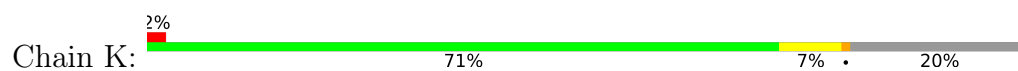
- Molecule 1: FERRITIN



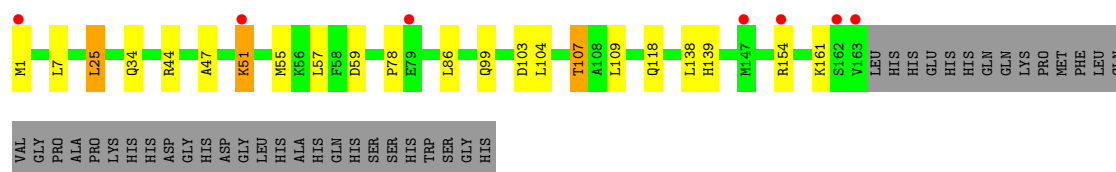
- Molecule 1: FERRITIN



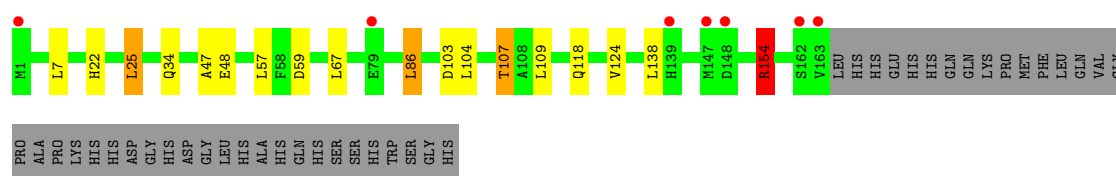
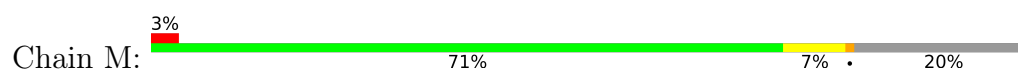
• Molecule 1: FERRITIN



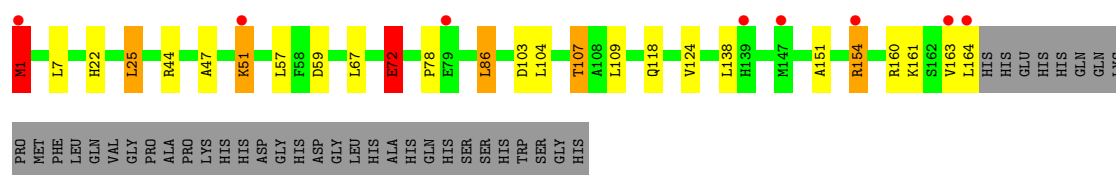
• Molecule 1: FERRITIN



• Molecule 1: FERRITIN



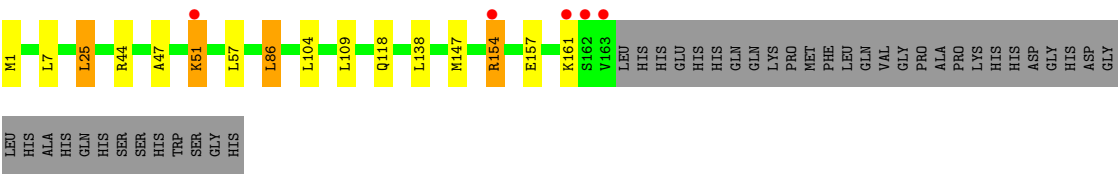
• Molecule 1: FERRITIN



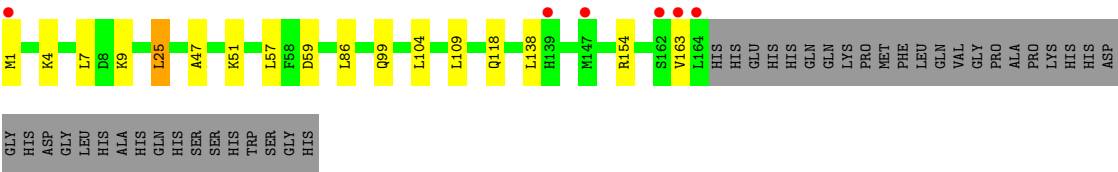
• Molecule 1: FERRITIN



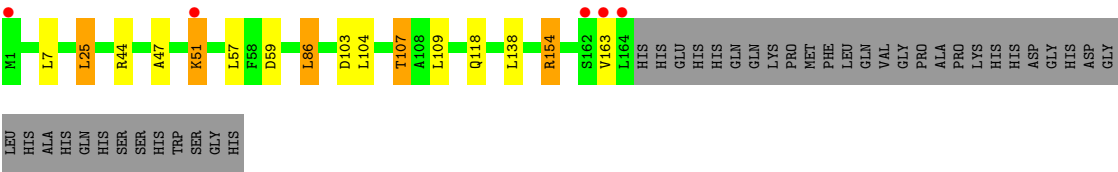




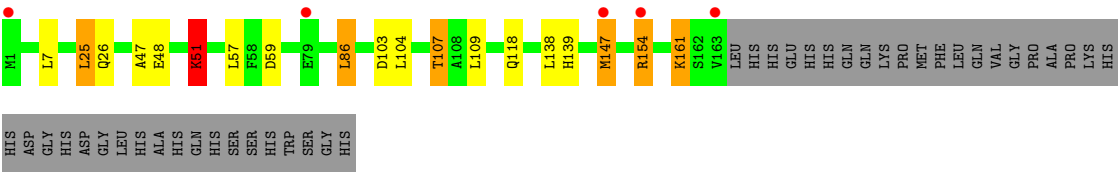
● Molecule 1: FERRITIN



● Molecule 1: FERRITIN



● Molecule 1: FERRITIN



● Molecule 2: FERRITIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	123.31Å 190.37Å 242.07Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	103.71 – 2.59 103.71 – 2.59	Depositor EDS
% Data completeness (in resolution range)	100.0 (103.71-2.59) 100.0 (103.71-2.59)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.58Å)	Xtriage
Refinement program	REFMAC 5.7.0029	Depositor
R, R_{free}	0.193 , 0.218 0.195 , 0.193	Depositor DCC
R_{free} test set	8902 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	37.5	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 26.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	32293	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% 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

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

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	0.94	0/1348	0.95	3/1825 (0.2%)
1	B	0.93	1/1372 (0.1%)	0.95	3/1857 (0.2%)
1	C	0.93	1/1348 (0.1%)	0.95	4/1825 (0.2%)
1	D	0.98	1/1348 (0.1%)	1.00	6/1825 (0.3%)
1	E	0.91	0/1348	0.88	0/1825
1	F	0.86	0/1359	0.87	3/1840 (0.2%)
1	G	0.91	1/1348 (0.1%)	0.90	2/1825 (0.1%)
1	H	1.01	3/1348 (0.2%)	1.02	8/1825 (0.4%)
1	I	0.90	0/1353	0.89	1/1832 (0.1%)
1	J	0.99	0/1364	0.94	3/1847 (0.2%)
1	K	0.96	0/1359	0.91	1/1840 (0.1%)
1	L	0.94	0/1359	0.92	3/1840 (0.2%)
1	M	0.93	0/1359	0.91	3/1840 (0.2%)
1	N	0.91	0/1353	0.96	4/1832 (0.2%)
1	O	0.91	2/1364 (0.1%)	0.99	6/1847 (0.3%)
1	P	0.97	5/1355 (0.4%)	0.94	2/1835 (0.1%)
1	Q	0.92	0/1359	0.91	3/1840 (0.2%)
1	R	0.95	0/1366	0.99	4/1850 (0.2%)
1	S	0.91	0/1343	0.91	3/1818 (0.2%)
1	T	0.93	0/1367	0.90	2/1850 (0.1%)
1	U	0.98	0/1364	0.92	2/1847 (0.1%)
1	W	0.94	0/1361	0.89	0/1842
1	X	1.01	2/1359 (0.1%)	0.98	6/1840 (0.3%)
2	V	0.89	0/1363	0.87	0/1844
All	All	0.94	16/32567 (0.0%)	0.93	72/44091 (0.2%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	X	51	LYS	CA-CB	9.58	1.68	1.53
1	D	146	THR	N-CA	9.47	1.58	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	148	ASP	CA-CB	9.39	1.68	1.53
1	O	110	ARG	CA-C	-7.14	1.43	1.52
1	P	162	SER	CA-C	6.96	1.61	1.52
1	X	51	LYS	CA-C	6.51	1.61	1.52
1	P	147	MET	CA-CB	6.10	1.62	1.53
1	O	110	ARG	C-O	-5.95	1.17	1.24
1	B	78	PRO	CA-C	5.86	1.60	1.52
1	H	148	ASP	N-CA	5.57	1.53	1.46
1	P	1	MET	C-N	-5.55	1.25	1.33
1	H	148	ASP	CA-C	5.39	1.59	1.52
1	P	139	HIS	CA-CB	5.39	1.61	1.53
1	G	146	THR	CA-C	-5.26	1.46	1.52
1	C	139	HIS	CA-CB	5.18	1.61	1.53
1	P	2	LEU	CA-C	-5.18	1.45	1.52

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	R	144	ILE	CA-CB-CG1	11.70	130.29	110.40
1	P	2	LEU	N-CA-CB	9.95	127.30	110.49
1	M	154	ARG	CG-CD-NE	9.79	133.55	112.00
1	O	110	ARG	CB-CA-C	-9.51	95.95	110.88
1	H	57	LEU	CB-CG-CD1	-9.36	82.61	110.70
1	C	1	MET	CB-CG-SD	9.20	140.31	112.70
1	N	1	MET	CB-CG-SD	8.84	139.21	112.70
1	D	145	ASN	N-CA-C	-8.57	92.91	107.99
1	X	51	LYS	CB-CA-C	8.40	125.13	110.85
1	N	161	LYS	CB-CG-CD	8.08	129.89	111.30
1	D	146	THR	N-CA-CB	8.05	124.10	110.49
1	X	147	MET	CG-SD-CE	8.05	118.60	100.90
1	I	147	MET	CB-CG-SD	8.00	136.71	112.70
1	Q	147	MET	CG-SD-CE	7.92	118.34	100.90
1	H	154	ARG	CG-CD-NE	7.88	129.33	112.00
1	A	154	ARG	CG-CD-NE	7.78	129.12	112.00
1	J	129	GLU	CG-CD-OE2	7.77	136.28	118.40
1	G	146	THR	N-CA-CB	-7.64	98.02	111.08
1	C	146	THR	N-CA-CB	-7.62	98.05	111.08
1	X	51	LYS	N-CA-C	-7.58	103.10	111.36
1	D	161	LYS	CB-CG-CD	7.55	128.67	111.30
1	O	110	ARG	N-CA-CB	7.29	120.57	110.01
1	B	80	TRP	N-CA-C	7.19	121.20	109.76
1	Q	146	THR	CA-CB-OG1	7.16	120.34	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	O	161	LYS	CG-CD-CE	7.12	127.68	111.30
1	R	49	GLU	CG-CD-OE2	7.08	134.69	118.40
1	F	121	GLN	CG-CD-NE2	6.99	126.89	116.40
1	T	147	MET	CG-SD-CE	6.98	116.26	100.90
1	O	110	ARG	CA-CB-CG	6.94	127.98	114.10
1	H	121	GLN	CG-CD-NE2	6.92	126.77	116.40
1	L	161	LYS	CB-CG-CD	6.89	127.15	111.30
1	P	1	MET	CA-C-O	6.79	132.35	120.80
1	R	51	LYS	CG-CD-CE	6.76	126.84	111.30
1	J	129	GLU	CG-CD-OE1	-6.74	102.89	118.40
1	N	72	GLU	CA-CB-CG	6.56	127.23	114.10
1	R	49	GLU	CG-CD-OE1	-6.42	103.63	118.40
1	H	148	ASP	CB-CA-C	6.35	121.33	110.79
1	S	34	GLN	CG-CD-NE2	6.35	125.92	116.40
1	M	34	GLN	CG-CD-NE2	6.28	125.81	116.40
1	F	1	MET	CG-SD-CE	6.25	114.66	100.90
1	J	147	MET	CG-SD-CE	-6.23	87.19	100.90
1	Q	1	MET	CG-SD-CE	6.23	114.60	100.90
1	B	78	PRO	CB-CA-C	6.19	121.77	111.56
1	K	161	LYS	CB-CG-CD	6.18	125.51	111.30
1	X	161	LYS	CA-CB-CG	6.17	126.44	114.10
1	L	34	GLN	CG-CD-NE2	6.01	125.41	116.40
1	O	161	LYS	CD-CE-NZ	5.99	131.07	111.90
1	G	146	THR	CA-CB-CG2	5.96	120.63	110.50
1	X	51	LYS	CB-CG-CD	5.90	124.87	111.30
1	A	154	ARG	CB-CG-CD	5.87	124.79	111.30
1	S	147	MET	N-CA-CB	5.87	118.84	110.16
1	D	51	LYS	CB-CG-CD	5.84	124.72	111.30
1	C	146	THR	CA-CB-CG2	5.83	120.41	110.50
1	H	154	ARG	CB-CG-CD	5.76	124.55	111.30
1	T	147	MET	N-CA-CB	5.72	118.62	110.16
1	D	146	THR	N-CA-C	-5.71	98.63	110.80
1	L	44	ARG	CG-CD-NE	5.69	124.52	112.00
1	F	121	GLN	CG-CD-OE1	-5.66	109.49	120.80
1	H	57	LEU	CB-CG-CD2	5.61	127.53	110.70
1	A	146	THR	CA-CB-CG2	5.60	120.02	110.50
1	D	146	THR	CA-CB-OG1	5.58	117.97	109.60
1	B	146	THR	N-CA-CB	5.57	119.76	110.85
1	U	4	LYS	N-CA-CB	5.49	118.79	110.28
1	H	121	GLN	CG-CD-OE1	-5.45	109.90	120.80
1	X	51	LYS	CA-CB-CG	-5.33	103.44	114.10
1	H	148	ASP	N-CA-C	-5.19	105.62	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	44	ARG	NE-CZ-NH1	5.16	126.66	121.50
1	N	160	ARG	CD-NE-CZ	5.16	131.62	124.40
1	O	154	ARG	CB-CG-CD	5.12	123.07	111.30
1	S	34	GLN	CG-CD-OE1	-5.08	110.65	120.80
1	U	4	LYS	CD-CE-NZ	-5.02	95.84	111.90
1	M	34	GLN	CG-CD-OE1	-5.02	110.76	120.80

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	1318	0	1263	11	0
1	B	1335	0	1283	12	0
1	C	1318	0	1263	11	0
1	D	1318	0	1263	16	0
1	E	1318	0	1263	10	0
1	F	1325	0	1270	6	0
1	G	1318	0	1263	10	0
1	H	1318	0	1263	21	0
1	I	1323	0	1265	13	0
1	J	1330	0	1272	16	0
1	K	1325	0	1270	11	0
1	L	1325	0	1270	12	0
1	M	1325	0	1270	10	0
1	N	1323	0	1265	20	0
1	O	1330	0	1272	15	0
1	P	1322	0	1270	16	0
1	Q	1325	0	1270	13	0
1	R	1329	0	1277	10	0
1	S	1313	0	1261	10	0
1	T	1333	0	1278	10	0
1	U	1330	0	1272	5	0
1	W	1331	0	1275	12	0
1	X	1325	0	1269	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	V	1329	0	1279	12	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
3	G	2	0	0	0	0
3	H	2	0	0	0	0
3	I	2	0	0	0	0
3	J	2	0	0	0	0
3	K	2	0	0	0	0
3	L	2	0	0	0	0
3	M	2	0	0	0	0
3	N	2	0	0	0	0
3	O	2	0	0	0	0
3	P	2	0	0	0	0
3	Q	2	0	0	0	0
3	R	2	0	0	0	0
3	S	2	0	0	0	0
3	T	2	0	0	0	0
3	U	2	0	0	0	0
3	V	2	0	0	0	0
3	W	2	0	0	0	0
3	X	2	0	0	0	0
4	A	31	0	0	4	0
4	B	16	0	0	1	0
4	C	30	0	0	0	0
4	D	17	0	0	3	0
4	E	20	0	0	0	0
4	F	6	0	0	0	0
4	G	14	0	0	0	0
4	H	16	0	0	2	0
4	I	18	0	0	0	0
4	J	29	0	0	1	0
4	K	33	0	0	3	0
4	L	23	0	0	0	0
4	M	10	0	0	0	0
4	N	20	0	0	3	0
4	O	10	0	0	2	0
4	P	10	0	0	1	0
4	Q	11	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	R	13	0	0	0	0
4	S	17	0	0	3	0
4	T	13	0	0	0	0
4	U	30	0	0	1	0
4	V	26	0	0	2	0
4	W	25	0	0	0	0
4	X	21	0	0	3	0
All	All	32293	0	30466	227	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 (227) 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:J:48:GLU:OE2	1:X:51:LYS:NZ	1.59	1.33
1:K:51:LYS:HE2	1:L:51:LYS:CE	1.58	1.32
1:K:51:LYS:CE	1:L:51:LYS:HE2	1.62	1.28
1:I:51:LYS:HE2	1:Q:51:LYS:CE	1.70	1.20
1:I:51:LYS:CE	1:Q:51:LYS:HE2	1.70	1.20
2:V:51:LYS:HE2	1:W:51:LYS:CE	1.70	1.19
2:V:51:LYS:CE	1:W:51:LYS:HE2	1.71	1.19
1:D:146:THR:HG22	1:D:149:GLY:H	1.05	1.13
1:X:48:GLU:OE2	1:X:51:LYS:HE3	1.52	1.09
1:C:48:GLU:OE1	1:E:51:LYS:NZ	1.84	1.08
1:J:48:GLU:CD	1:X:51:LYS:NZ	2.20	0.99
1:H:57:LEU:N	1:H:57:LEU:HD12	1.76	0.98
1:H:57:LEU:HD12	1:H:57:LEU:H	1.30	0.95
2:V:51:LYS:HE2	1:W:51:LYS:HE2	0.89	0.89
1:D:146:THR:HG22	1:D:149:GLY:N	1.90	0.85
1:G:51:LYS:HE3	1:T:51:LYS:HE2	1.60	0.83
1:N:51:LYS:HE3	1:O:51:LYS:HE2	1.60	0.83
1:I:51:LYS:HE2	1:Q:51:LYS:HE2	0.83	0.78
1:D:146:THR:CG2	1:D:149:GLY:H	1.93	0.76
1:N:72:GLU:OE1	1:N:72:GLU:N	2.19	0.76
1:R:5[B]:THR:HG21	1:R:111:GLU:OE2	1.87	0.74
1:G:51:LYS:CE	1:T:51:LYS:HE2	2.22	0.70
1:R:5[B]:THR:CG2	1:R:111:GLU:OE2	2.40	0.70
1:J:129:GLU:HA	1:J:129:GLU:OE1	1.92	0.70
1:P:159:VAL:O	1:P:162:SER:O	2.09	0.69
1:H:53:HIS:O	1:H:57:LEU:CD1	2.41	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:48:GLU:CD	1:X:51:LYS:HZ3	1.97	0.68
1:A:153:PHE:CD1	4:A:2031:HOH:O	2.46	0.68
1:S:22:HIS:HD2	4:S:2011:HOH:O	1.76	0.67
1:P:1:MET:O	4:P:2002:HOH:O	2.11	0.67
1:K:51:LYS:CE	1:L:51:LYS:CE	2.44	0.67
1:N:51:LYS:CE	1:O:51:LYS:HE2	2.24	0.67
2:V:81:LYS:HG3	2:V:85:GLU:OE1	1.94	0.67
1:P:146[A]:THR:HG22	1:P:149:GLY:H	1.60	0.66
1:T:1[A]:MET:HE1	4:V:2020:HOH:O	1.96	0.65
1:D:69:LEU:HA	1:F:26:GLN:HE22	1.62	0.64
1:H:57:LEU:CD1	1:H:123:TYR:OH	2.46	0.64
1:G:146:THR:HG22	1:G:149:GLY:H	1.62	0.64
1:A:153:PHE:CE1	4:A:2031:HOH:O	2.50	0.64
1:C:146:THR:HG22	1:C:149:GLY:H	1.62	0.64
1:L:103:ASP:O	1:L:107:THR:HG23	1.98	0.64
1:D:143:ILE:O	1:D:146:THR:HB	1.98	0.64
1:C:103:ASP:O	1:C:107:THR:HG23	1.98	0.63
1:H:57:LEU:HD11	1:H:123:TYR:OH	1.98	0.63
1:K:51:LYS:HE2	1:L:51:LYS:HE2	0.73	0.63
1:I:103:ASP:O	1:I:107:THR:HG23	1.99	0.63
1:N:154:ARG:HG2	1:P:154:ARG:HH22	1.64	0.62
1:H:103:ASP:O	1:H:107:THR:HG23	2.00	0.62
1:M:103:ASP:O	1:M:107:THR:HG23	1.99	0.62
1:A:103:ASP:O	1:A:107:THR:HG23	2.00	0.62
1:J:103:ASP:O	1:J:107:THR:HG23	2.00	0.62
1:K:103:ASP:O	1:K:107:THR:HG23	2.00	0.61
1:O:139[B]:HIS:HE1	4:O:2010:HOH:O	1.82	0.61
1:N:103:ASP:O	1:N:107:THR:HG23	2.00	0.61
1:C:48:GLU:CD	1:E:51:LYS:NZ	2.59	0.61
1:O:103:ASP:O	1:O:107:THR:HG23	2.01	0.60
1:B:103:ASP:O	1:B:107:THR:HG23	2.00	0.60
4:N:2006:HOH:O	1:P:139:HIS:HB3	2.00	0.60
1:F:103:ASP:O	1:F:107:THR:HG23	2.01	0.60
1:Q:103:ASP:O	1:Q:107:THR:HG23	2.01	0.60
1:W:103:ASP:O	1:W:107:THR:HG23	2.01	0.60
1:X:103:ASP:O	1:X:107:THR:HG23	2.01	0.60
1:S:103:ASP:O	1:S:107:THR:HG23	2.01	0.60
1:H:57:LEU:H	1:H:57:LEU:CD1	2.11	0.59
2:V:44:ARG:NH1	1:W:59:ASP:OD1	2.28	0.59
1:H:154:ARG:HD3	4:K:2032:HOH:O	2.01	0.59
1:W:154:ARG:NH2	1:X:154:ARG:HG2	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:69:LEU:HA	1:Q:26:GLN:HE22	1.66	0.59
1:H:69:LEU:HA	1:P:26:GLN:HE22	1.67	0.58
1:J:48:GLU:OE1	1:X:51:LYS:NZ	2.27	0.58
2:V:103:ASP:O	2:V:107:THR:HG23	2.03	0.58
1:H:51:LYS:NZ	1:P:51:LYS:HD3	2.19	0.57
4:H:2013:HOH:O	1:L:1:MET:HE1	2.03	0.57
1:J:69:LEU:HA	1:X:26:GLN:HE22	1.68	0.57
1:A:51:LYS:NZ	1:M:48:GLU:OE1	2.33	0.56
1:B:163:VAL:O	1:B:164:LEU:CB	2.53	0.55
1:J:1:MET:N	4:J:2001:HOH:O	2.39	0.55
2:V:81:LYS:CG	2:V:85:GLU:OE1	2.56	0.54
1:N:59:ASP:OD1	1:O:44:ARG:NH1	2.29	0.54
1:D:152:LEU:HB3	1:E:36:LEU:HD21	1.90	0.53
1:H:148:ASP:OD1	1:H:148:ASP:N	2.41	0.53
1:J:129:GLU:OE1	1:J:129:GLU:CA	2.56	0.52
1:N:163:VAL:O	1:N:164:LEU:CB	2.56	0.52
1:U:1:MET:HA	1:U:1:MET:HE2	1.91	0.52
1:N:22:HIS:HD2	4:N:2015:HOH:O	1.92	0.52
1:G:154:ARG:HH22	1:O:154:ARG:HG2	1.75	0.51
1:P:1:MET:O	1:P:2:LEU:CB	2.59	0.51
1:S:48:GLU:OE1	1:U:51:LYS:NZ	2.35	0.51
1:A:22:HIS:HD2	4:A:2018:HOH:O	1.93	0.51
4:H:2010:HOH:O	1:P:22:HIS:HD2	1.94	0.50
1:I:1:MET:HA	1:I:1:MET:HE2	1.93	0.50
1:D:150:ARG:NH1	1:N:151:ALA:HA	2.24	0.50
1:E:147:MET:HA	1:E:150:ARG:NH1	2.27	0.50
1:G:51:LYS:NZ	1:T:51:LYS:HE2	2.26	0.49
1:J:154:ARG:HH22	1:M:154:ARG:HG2	1.77	0.49
1:G:55:MET:HE3	1:T:44:ARG:HH21	1.76	0.49
1:I:51:LYS:CE	1:Q:51:LYS:CE	2.54	0.49
1:P:1:MET:O	1:P:2:LEU:HB2	2.12	0.49
2:V:51:LYS:CE	1:W:51:LYS:CE	2.53	0.49
1:X:147:MET:HG3	4:X:2020:HOH:O	2.13	0.49
1:H:44:ARG:HH21	1:P:55:MET:HE3	1.79	0.48
1:G:51:LYS:NZ	1:T:51:LYS:CE	2.76	0.48
1:N:103:ASP:O	1:N:107:THR:CG2	2.62	0.48
1:D:142:ARG:O	1:D:145:ASN:O	2.32	0.48
1:C:103:ASP:O	1:C:107:THR:CG2	2.62	0.48
1:H:51:LYS:HZ3	1:P:48:GLU:CD	2.22	0.48
1:T:157:GLU:O	1:T:161:LYS:HG3	2.14	0.47
1:F:103:ASP:O	1:F:107:THR:CG2	2.63	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:103:ASP:O	1:L:107:THR:CG2	2.62	0.47
1:I:103:ASP:O	1:I:107:THR:CG2	2.63	0.47
1:K:103:ASP:O	1:K:107:THR:CG2	2.63	0.47
1:S:154:ARG:HG2	1:X:154:ARG:HH22	1.80	0.47
1:W:103:ASP:O	1:W:107:THR:CG2	2.62	0.47
1:B:146:THR:HG23	1:B:149:GLY:H	1.79	0.47
1:H:53:HIS:O	1:H:57:LEU:HD13	2.15	0.47
1:I:154:ARG:HG2	1:T:154:ARG:NH2	2.30	0.47
1:J:103:ASP:O	1:J:107:THR:CG2	2.63	0.47
1:N:44:ARG:HH21	1:O:55:MET:HE3	1.80	0.46
1:X:139[B]:HIS:CE1	4:X:2017:HOH:O	2.68	0.46
1:H:103:ASP:O	1:H:107:THR:CG2	2.62	0.46
1:S:103:ASP:O	1:S:107:THR:CG2	2.63	0.46
1:H:57:LEU:HD11	1:H:123:TYR:CZ	2.49	0.46
1:Q:103:ASP:O	1:Q:107:THR:CG2	2.64	0.46
1:N:25:LEU:HD13	1:N:47:ALA:CB	2.46	0.46
1:X:103:ASP:O	1:X:107:THR:CG2	2.64	0.46
2:V:25:LEU:HD13	2:V:47:ALA:CB	2.45	0.46
2:V:103:ASP:O	2:V:107:THR:CG2	2.64	0.46
1:D:49:GLU:OE1	4:D:2006:HOH:O	2.21	0.46
1:H:57:LEU:HD12	1:H:123:TYR:OH	2.13	0.46
1:O:25:LEU:HD13	1:O:47:ALA:CB	2.45	0.46
1:L:25:LEU:HD13	1:L:47:ALA:CB	2.46	0.46
1:O:103:ASP:O	1:O:107:THR:CG2	2.63	0.45
1:S:110:ARG:NH2	4:S:2015:HOH:O	2.49	0.45
1:C:48:GLU:CD	1:E:51:LYS:HZ1	2.15	0.45
1:J:25:LEU:HD13	1:J:47:ALA:CB	2.46	0.45
1:R:25:LEU:HD13	1:R:47:ALA:CB	2.46	0.45
1:S:109:LEU:HD22	1:S:117:PHE:CE1	2.51	0.45
1:D:25:LEU:HD13	1:D:47:ALA:CB	2.47	0.45
1:B:78:PRO:HG2	1:R:67:LEU:HD21	1.99	0.45
1:B:103:ASP:O	1:B:107:THR:CG2	2.63	0.45
1:I:55:MET:HE3	1:Q:44:ARG:HH21	1.81	0.45
1:B:25:LEU:HD13	1:B:47:ALA:CB	2.46	0.45
1:I:25:LEU:HD13	1:I:47:ALA:CB	2.47	0.45
1:M:103:ASP:O	1:M:107:THR:CG2	2.63	0.45
1:A:44:ARG:NH2	1:M:59:ASP:OD2	2.48	0.45
1:M:25:LEU:HD13	1:M:47:ALA:CB	2.47	0.45
1:O:139[B]:HIS:CE1	4:O:2010:HOH:O	2.64	0.45
1:S:25:LEU:HD13	1:S:47:ALA:CB	2.47	0.45
1:A:25:LEU:HD13	1:A:47:ALA:CB	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:ASP:O	1:A:107:THR:CG2	2.63	0.45
1:B:21:ALA:CB	1:B:51:LYS:HE2	2.47	0.45
1:K:139[A]:HIS:CD2	4:K:2031:HOH:O	2.70	0.45
1:Q:86:LEU:C	1:Q:86:LEU:HD12	2.42	0.45
1:E:86:LEU:HD12	1:E:86:LEU:C	2.42	0.44
1:G:86:LEU:C	1:G:86:LEU:HD12	2.42	0.44
1:D:126:GLU:OE2	4:D:2006:HOH:O	2.21	0.44
1:H:55:MET:HE3	1:P:44:ARG:HH21	1.82	0.44
1:J:86:LEU:HD12	1:J:86:LEU:C	2.42	0.44
1:B:22:HIS:HD2	4:B:2010:HOH:O	2.01	0.44
1:C:25:LEU:HD13	1:C:47:ALA:CB	2.48	0.44
1:Q:25:LEU:HD13	1:Q:47:ALA:CB	2.48	0.44
1:P:25:LEU:HD13	1:P:47:ALA:CB	2.47	0.44
1:U:25:LEU:HD13	1:U:47:ALA:CB	2.47	0.44
1:T:25:LEU:HD13	1:T:47:ALA:CB	2.48	0.44
1:C:35:SER:OG	1:L:139[B]:HIS:CD2	2.71	0.44
1:X:139[B]:HIS:HE1	4:X:2017:HOH:O	1.99	0.44
1:E:25:LEU:HD13	1:E:47:ALA:CB	2.48	0.43
1:A:153:PHE:HB3	1:O:154:ARG:NH1	2.33	0.43
1:D:86:LEU:C	1:D:86:LEU:HD12	2.43	0.43
1:G:25:LEU:HD13	1:G:47:ALA:CB	2.47	0.43
1:N:67:LEU:HD21	1:O:78:PRO:HG2	1.99	0.43
1:S:44:ARG:NH2	1:U:59:ASP:OD2	2.51	0.43
1:B:86:LEU:C	1:B:86:LEU:HD12	2.42	0.43
1:K:25:LEU:HD13	1:K:47:ALA:CB	2.47	0.43
1:X:25:LEU:HD13	1:X:47:ALA:CB	2.48	0.43
1:R:5[B]:THR:HG22	1:R:111:GLU:OE2	2.16	0.43
1:X:161:LYS:HB2	1:X:161:LYS:HE2	1.80	0.43
1:W:25:LEU:HD13	1:W:47:ALA:CB	2.48	0.43
1:F:25:LEU:HD13	1:F:47:ALA:CB	2.49	0.42
1:N:86:LEU:C	1:N:86:LEU:HD12	2.44	0.42
1:N:1:MET:N	4:N:2001:HOH:O	2.43	0.42
1:H:25:LEU:HD13	1:H:47:ALA:CB	2.49	0.42
1:I:109:LEU:HD22	1:I:117:PHE:CE1	2.54	0.42
1:A:86:LEU:HD12	1:A:86:LEU:C	2.44	0.42
1:H:86:LEU:HD12	1:H:86:LEU:C	2.45	0.42
1:K:44:ARG:HH21	1:L:55:MET:HE3	1.83	0.42
1:R:140:LYS:O	1:R:144:ILE:HG12	2.20	0.42
1:S:139:HIS:HB2	4:S:2017:HOH:O	2.19	0.42
1:C:59:ASP:OD2	1:E:44:ARG:NH2	2.53	0.42
1:K:44:ARG:NH2	1:L:59:ASP:OD2	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:51:LYS:NZ	1:O:51:LYS:CE	2.83	0.42
1:Q:86:LEU:C	1:Q:86:LEU:CD1	2.93	0.42
1:K:67:LEU:HD21	1:L:78:PRO:HG2	2.02	0.42
1:W:154:ARG:HH21	1:X:154:ARG:HG2	1.83	0.42
1:M:86:LEU:HD12	1:M:86:LEU:C	2.44	0.42
1:T:86:LEU:HD12	1:T:86:LEU:C	2.44	0.42
1:C:124:VAL:CG1	1:R:115:SER:HA	2.50	0.41
1:A:78:PRO:HG2	1:M:67:LEU:HD21	2.02	0.41
1:J:86:LEU:C	1:J:86:LEU:CD1	2.94	0.41
1:R:86:LEU:HD12	1:R:86:LEU:C	2.45	0.41
1:D:99:GLN:HG2	4:D:2010:HOH:O	2.20	0.41
4:A:2016:HOH:O	1:M:22:HIS:HD2	2.03	0.41
1:H:154:ARG:CD	4:K:2032:HOH:O	2.65	0.41
1:D:86:LEU:C	1:D:86:LEU:CD1	2.94	0.41
1:E:154:ARG:HH22	1:P:154:ARG:HG2	1.85	0.41
1:B:86:LEU:C	1:B:86:LEU:CD1	2.94	0.41
1:D:154:ARG:HG2	1:N:154:ARG:HH22	1.86	0.41
1:F:86:LEU:HD12	1:F:86:LEU:C	2.44	0.41
1:N:78:PRO:HG2	1:O:67:LEU:HD21	2.03	0.41
1:Q:27:MET:SD	1:Q:86:LEU:HD13	2.61	0.41
2:V:1:MET:N	4:V:2001:HOH:O	2.46	0.41
1:F:115:SER:HA	1:N:124:VAL:CG1	2.51	0.40
1:J:27:MET:SD	1:J:86:LEU:HD13	2.61	0.40
1:W:86:LEU:C	1:W:86:LEU:HD12	2.46	0.40
1:B:55:MET:HE3	1:R:44:ARG:HH21	1.85	0.40
2:V:59:ASP:OD1	1:W:44:ARG:NH1	2.38	0.40
1:E:86:LEU:C	1:E:86:LEU:CD1	2.94	0.40
1:I:59:ASP:OD2	1:Q:44:ARG:NH2	2.54	0.40
1:J:44:ARG:NH2	1:X:59:ASP:OD2	2.55	0.40
1:N:51:LYS:HE3	1:O:51:LYS:CE	2.40	0.40
1:X:86:LEU:C	1:X:86:LEU:HD12	2.46	0.40
1:B:154:ARG:HG2	1:C:154:ARG:HH22	1.86	0.40
1:D:27:MET:SD	1:D:86:LEU:HD13	2.62	0.40
1:M:124:VAL:CG1	1:P:115:SER:HA	2.52	0.40
1:G:86:LEU:C	1:G:86:LEU:CD1	2.95	0.40
1:R:5[B]:THR:CG2	1:R:6:ILE:N	2.83	0.40
1:U:9:LYS:HE2	4:U:2005:HOH:O	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	161/203 (79%)	160 (99%)	1 (1%)	0	100	100
1	B	163/203 (80%)	161 (99%)	2 (1%)	0	100	100
1	C	161/203 (79%)	161 (100%)	0	0	100	100
1	D	161/203 (79%)	160 (99%)	0	1 (1%)	21	42
1	E	161/203 (79%)	161 (100%)	0	0	100	100
1	F	162/203 (80%)	162 (100%)	0	0	100	100
1	G	161/203 (79%)	161 (100%)	0	0	100	100
1	H	161/203 (79%)	161 (100%)	0	0	100	100
1	I	162/203 (80%)	160 (99%)	1 (1%)	1 (1%)	21	42
1	J	163/203 (80%)	161 (99%)	1 (1%)	1 (1%)	21	42
1	K	162/203 (80%)	162 (100%)	0	0	100	100
1	L	162/203 (80%)	161 (99%)	1 (1%)	0	100	100
1	M	162/203 (80%)	162 (100%)	0	0	100	100
1	N	162/203 (80%)	161 (99%)	1 (1%)	0	100	100
1	O	163/203 (80%)	162 (99%)	0	1 (1%)	21	42
1	P	162/203 (80%)	161 (99%)	0	1 (1%)	21	42
1	Q	162/203 (80%)	162 (100%)	0	0	100	100
1	R	163/203 (80%)	163 (100%)	0	0	100	100
1	S	160/203 (79%)	160 (100%)	0	0	100	100
1	T	162/203 (80%)	162 (100%)	0	0	100	100
1	U	163/203 (80%)	161 (99%)	1 (1%)	1 (1%)	21	42
1	W	162/203 (80%)	160 (99%)	1 (1%)	1 (1%)	21	42
1	X	162/203 (80%)	162 (100%)	0	0	100	100
2	V	162/203 (80%)	161 (99%)	1 (1%)	0	100	100
All	All	3885/4872 (80%)	3868 (100%)	10 (0%)	7 (0%)	43	66

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	146	THR
1	P	2	LEU
1	I	163	VAL
1	J	163	VAL
1	O	163	VAL
1	W	163	VAL
1	U	163	VAL

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	137/175 (78%)	126 (92%)	11 (8%)	11	25
1	B	139/175 (79%)	129 (93%)	10 (7%)	13	30
1	C	137/175 (78%)	125 (91%)	12 (9%)	9	21
1	D	137/175 (78%)	126 (92%)	11 (8%)	11	25
1	E	137/175 (78%)	127 (93%)	10 (7%)	13	29
1	F	138/175 (79%)	127 (92%)	11 (8%)	11	25
1	G	137/175 (78%)	127 (93%)	10 (7%)	13	29
1	H	137/175 (78%)	126 (92%)	11 (8%)	11	25
1	I	137/175 (78%)	125 (91%)	12 (9%)	9	21
1	J	138/175 (79%)	127 (92%)	11 (8%)	11	25
1	K	138/175 (79%)	126 (91%)	12 (9%)	9	21
1	L	138/175 (79%)	126 (91%)	12 (9%)	9	21
1	M	138/175 (79%)	128 (93%)	10 (7%)	13	30
1	N	137/175 (78%)	124 (90%)	13 (10%)	8	18
1	O	138/175 (79%)	126 (91%)	12 (9%)	9	21
1	P	138/175 (79%)	128 (93%)	10 (7%)	13	30
1	Q	138/175 (79%)	126 (91%)	12 (9%)	9	21

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	R	139/175 (79%)	129 (93%)	10 (7%)	13	30
1	S	137/175 (78%)	125 (91%)	12 (9%)	9	21
1	T	139/175 (79%)	129 (93%)	10 (7%)	13	30
1	U	138/175 (79%)	128 (93%)	10 (7%)	13	30
1	W	138/175 (79%)	127 (92%)	11 (8%)	11	25
1	X	138/175 (79%)	127 (92%)	11 (8%)	11	25
2	V	139/175 (79%)	126 (91%)	13 (9%)	8	18
All	All	3307/4200 (79%)	3040 (92%)	267 (8%)	11	24

All (267) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	LEU
1	A	25	LEU
1	A	57	LEU
1	A	86	LEU
1	A	104	LEU
1	A	107	THR
1	A	109	LEU
1	A	118	GLN
1	A	138	LEU
1	A	146	THR
1	A	154	ARG
1	B	7	LEU
1	B	25	LEU
1	B	57	LEU
1	B	86	LEU
1	B	104	LEU
1	B	107	THR
1	B	109	LEU
1	B	118	GLN
1	B	138	LEU
1	B	154	ARG
1	C	1	MET
1	C	7	LEU
1	C	25	LEU
1	C	51	LYS
1	C	57	LEU
1	C	86	LEU
1	C	104	LEU

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Mol	Chain	Res	Type
1	C	107	THR
1	C	109	LEU
1	C	118	GLN
1	C	138	LEU
1	C	154	ARG
1	D	7	LEU
1	D	25	LEU
1	D	51	LYS
1	D	57	LEU
1	D	86	LEU
1	D	104	LEU
1	D	109	LEU
1	D	118	GLN
1	D	138	LEU
1	D	145	ASN
1	D	154	ARG
1	E	7	LEU
1	E	25	LEU
1	E	57	LEU
1	E	86	LEU
1	E	104	LEU
1	E	109	LEU
1	E	118	GLN
1	E	138	LEU
1	E	150	ARG
1	E	154	ARG
1	F	7	LEU
1	F	25	LEU
1	F	51	LYS
1	F	86	LEU
1	F	99	GLN
1	F	104	LEU
1	F	107	THR
1	F	109	LEU
1	F	118	GLN
1	F	138	LEU
1	F	154	ARG
1	G	7	LEU
1	G	25	LEU
1	G	51	LYS
1	G	57	LEU
1	G	86	LEU

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Mol	Chain	Res	Type
1	G	104	LEU
1	G	109	LEU
1	G	118	GLN
1	G	138	LEU
1	G	154	ARG
1	H	1	MET
1	H	7	LEU
1	H	25	LEU
1	H	86	LEU
1	H	104	LEU
1	H	107	THR
1	H	109	LEU
1	H	118	GLN
1	H	138	LEU
1	H	148	ASP
1	H	154	ARG
1	I	7	LEU
1	I	25	LEU
1	I	51	LYS
1	I	57	LEU
1	I	86	LEU
1	I	104	LEU
1	I	107	THR
1	I	109	LEU
1	I	118	GLN
1	I	138	LEU
1	I	147	MET
1	I	154	ARG
1	J	7	LEU
1	J	25	LEU
1	J	51	LYS
1	J	57	LEU
1	J	86	LEU
1	J	104	LEU
1	J	107	THR
1	J	109	LEU
1	J	118	GLN
1	J	138	LEU
1	J	154	ARG
1	K	1	MET
1	K	7	LEU
1	K	25	LEU

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Mol	Chain	Res	Type
1	K	51	LYS
1	K	57	LEU
1	K	86	LEU
1	K	104	LEU
1	K	107	THR
1	K	109	LEU
1	K	118	GLN
1	K	138	LEU
1	K	154	ARG
1	L	7	LEU
1	L	25	LEU
1	L	51	LYS
1	L	57	LEU
1	L	86	LEU
1	L	99	GLN
1	L	104	LEU
1	L	107	THR
1	L	109	LEU
1	L	118	GLN
1	L	138	LEU
1	L	154	ARG
1	M	7	LEU
1	M	25	LEU
1	M	57	LEU
1	M	86	LEU
1	M	104	LEU
1	M	107	THR
1	M	109	LEU
1	M	118	GLN
1	M	138	LEU
1	M	154	ARG
1	N	1	MET
1	N	7	LEU
1	N	25	LEU
1	N	51	LYS
1	N	57	LEU
1	N	72	GLU
1	N	86	LEU
1	N	104	LEU
1	N	107	THR
1	N	109	LEU
1	N	118	GLN

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Mol	Chain	Res	Type
1	N	138	LEU
1	N	154	ARG
1	O	7	LEU
1	O	25	LEU
1	O	51	LYS
1	O	57	LEU
1	O	86	LEU
1	O	104	LEU
1	O	107	THR
1	O	109	LEU
1	O	110	ARG
1	O	118	GLN
1	O	138	LEU
1	O	154	ARG
1	P	7	LEU
1	P	25	LEU
1	P	51	LYS
1	P	57	LEU
1	P	86	LEU
1	P	104	LEU
1	P	109	LEU
1	P	118	GLN
1	P	138	LEU
1	P	154	ARG
1	Q	7	LEU
1	Q	25	LEU
1	Q	51	LYS
1	Q	57	LEU
1	Q	86	LEU
1	Q	104	LEU
1	Q	107	THR
1	Q	109	LEU
1	Q	118	GLN
1	Q	138	LEU
1	Q	146	THR
1	Q	154	ARG
1	R	7	LEU
1	R	25	LEU
1	R	51	LYS
1	R	57	LEU
1	R	86	LEU
1	R	104	LEU

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Mol	Chain	Res	Type
1	R	109	LEU
1	R	118	GLN
1	R	138	LEU
1	R	154	ARG
1	S	7	LEU
1	S	25	LEU
1	S	51	LYS
1	S	57	LEU
1	S	86	LEU
1	S	99	GLN
1	S	104	LEU
1	S	107	THR
1	S	109	LEU
1	S	118	GLN
1	S	138	LEU
1	S	154	ARG
1	T	7	LEU
1	T	25	LEU
1	T	51	LYS
1	T	57	LEU
1	T	86	LEU
1	T	104	LEU
1	T	109	LEU
1	T	118	GLN
1	T	138	LEU
1	T	154	ARG
1	U	7	LEU
1	U	25	LEU
1	U	57	LEU
1	U	86	LEU
1	U	99	GLN
1	U	104	LEU
1	U	109	LEU
1	U	118	GLN
1	U	138	LEU
1	U	154	ARG
2	V	1	MET
2	V	7	LEU
2	V	25	LEU
2	V	51	LYS
2	V	57	LEU
2	V	81	LYS

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Mol	Chain	Res	Type
2	V	86	LEU
2	V	104	LEU
2	V	107	THR
2	V	109	LEU
2	V	118	GLN
2	V	138	LEU
2	V	154	ARG
1	W	7	LEU
1	W	25	LEU
1	W	51	LYS
1	W	57	LEU
1	W	86	LEU
1	W	104	LEU
1	W	107	THR
1	W	109	LEU
1	W	118	GLN
1	W	138	LEU
1	W	154	ARG
1	X	7	LEU
1	X	25	LEU
1	X	51	LYS
1	X	57	LEU
1	X	86	LEU
1	X	104	LEU
1	X	107	THR
1	X	109	LEU
1	X	118	GLN
1	X	138	LEU
1	X	154	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (66) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	HIS
1	A	102	ASN
1	A	118	GLN
1	A	139	HIS
1	B	22	HIS
1	B	102	ASN
1	C	99	GLN
1	C	102	ASN
1	C	118	GLN

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Mol	Chain	Res	Type
1	D	22	HIS
1	D	102	ASN
1	D	118	GLN
1	E	102	ASN
1	E	118	GLN
1	E	139	HIS
1	F	22	HIS
1	F	26	GLN
1	F	102	ASN
1	F	118	GLN
1	F	121	GLN
1	G	102	ASN
1	G	118	GLN
1	G	139	HIS
1	H	102	ASN
1	H	118	GLN
1	H	121	GLN
1	I	102	ASN
1	I	118	GLN
1	J	102	ASN
1	K	102	ASN
1	K	118	GLN
1	L	102	ASN
1	L	118	GLN
1	M	22	HIS
1	M	102	ASN
1	M	118	GLN
1	N	22	HIS
1	N	102	ASN
1	N	118	GLN
1	O	102	ASN
1	O	118	GLN
1	P	22	HIS
1	P	26	GLN
1	P	102	ASN
1	P	118	GLN
1	Q	26	GLN
1	Q	102	ASN
1	Q	118	GLN
1	R	102	ASN
1	R	118	GLN
1	S	22	HIS

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Mol	Chain	Res	Type
1	S	102	ASN
1	S	139	HIS
1	T	102	ASN
1	T	121	GLN
1	U	99	GLN
1	U	102	ASN
1	U	118	GLN
2	V	102	ASN
2	V	118	GLN
1	W	102	ASN
1	W	118	GLN
1	W	139	HIS
1	X	26	GLN
1	X	102	ASN
1	X	118	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 48 ligands modelled in this entry, 48 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	163/203 (80%)	-0.21	4 (2%) 58 52	24, 33, 61, 79	0
1	B	164/203 (80%)	0.10	10 (6%) 27 22	23, 40, 69, 84	2 (1%)
1	C	163/203 (80%)	0.01	6 (3%) 45 39	26, 40, 67, 84	0
1	D	163/203 (80%)	0.19	8 (4%) 35 29	30, 41, 70, 96	0
1	E	163/203 (80%)	-0.09	5 (3%) 51 45	25, 38, 70, 94	0
1	F	163/203 (80%)	0.09	7 (4%) 40 34	28, 43, 68, 84	1 (0%)
1	G	163/203 (80%)	0.08	6 (3%) 45 39	31, 42, 70, 86	0
1	H	163/203 (80%)	0.01	8 (4%) 35 29	27, 40, 71, 94	0
1	I	164/203 (80%)	-0.02	10 (6%) 27 22	28, 41, 71, 98	0
1	J	164/203 (80%)	-0.14	7 (4%) 40 34	22, 33, 58, 79	2 (1%)
1	K	163/203 (80%)	-0.27	5 (3%) 51 45	19, 32, 59, 87	1 (0%)
1	L	163/203 (80%)	-0.19	7 (4%) 40 34	22, 33, 65, 82	1 (0%)
1	M	163/203 (80%)	-0.04	7 (4%) 40 34	23, 38, 66, 99	1 (0%)
1	N	164/203 (80%)	0.17	8 (4%) 35 29	29, 43, 74, 91	0
1	O	164/203 (80%)	0.05	7 (4%) 40 34	24, 42, 71, 85	1 (0%)
1	P	163/203 (80%)	0.15	9 (5%) 30 25	29, 43, 77, 96	1 (0%)
1	Q	163/203 (80%)	-0.04	8 (4%) 35 29	23, 37, 66, 78	1 (0%)
1	R	163/203 (80%)	0.09	9 (5%) 30 25	25, 38, 71, 94	2 (1%)
1	S	162/203 (79%)	-0.12	6 (3%) 45 39	24, 38, 65, 88	0
1	T	163/203 (80%)	-0.13	5 (3%) 51 45	21, 37, 62, 93	2 (1%)
1	U	164/203 (80%)	-0.19	6 (3%) 45 39	20, 32, 65, 82	1 (0%)
1	W	164/203 (80%)	-0.17	5 (3%) 52 47	18, 32, 62, 77	1 (0%)
1	X	163/203 (80%)	-0.27	5 (3%) 51 45	22, 32, 65, 87	1 (0%)
2	V	163/203 (80%)	-0.06	11 (6%) 24 19	21, 36, 69, 92	1 (0%)

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
All	All	3918/4872 (80%)	-0.04	169 (4%)	40	34	18, 38, 69, 99	19 (0%)

All (169) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	W	1[A]	MET	7.5
1	B	1[A]	MET	6.7
1	J	1	MET	5.7
1	P	1	MET	5.3
1	A	163	VAL	5.1
1	L	163	VAL	5.1
1	R	163	VAL	5.1
1	H	147	MET	4.9
2	V	76	PRO	4.8
1	Q	163	VAL	4.8
1	K	1	MET	4.7
1	C	1	MET	4.6
1	D	163	VAL	4.5
1	G	163	VAL	4.4
2	V	163	VAL	4.4
2	V	79	GLU	4.4
1	C	163	VAL	4.3
1	N	164	LEU	4.2
1	K	163	VAL	4.2
1	I	148	ASP	4.1
1	P	162	SER	4.1
1	M	163	VAL	4.0
1	A	1	MET	4.0
2	V	1	MET	3.9
1	D	1	MET	3.9
1	N	1	MET	3.9
1	W	162	SER	3.9
1	T	163	VAL	3.9
1	J	164	LEU	3.8
1	S	1	MET	3.7
1	H	1	MET	3.7
1	T	162	SER	3.7
1	X	163	VAL	3.6
2	V	74	ALA	3.6
1	P	51	LYS	3.6
1	M	148	ASP	3.6
1	H	162	SER	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	51	LYS	3.6
1	W	164	LEU	3.6
1	F	1	MET	3.5
1	O	164	LEU	3.5
2	V	51	LYS	3.5
1	S	147	MET	3.4
1	I	147	MET	3.4
1	C	162	SER	3.4
1	U	162	SER	3.4
1	E	163	VAL	3.4
1	B	79	GLU	3.4
1	M	147	MET	3.3
1	O	51	LYS	3.3
1	S	51	LYS	3.3
1	U	1	MET	3.3
1	B	162	SER	3.3
1	O	1	MET	3.3
1	J	162	SER	3.2
1	M	162	SER	3.2
1	C	51	LYS	3.2
1	D	147	MET	3.2
1	N	154	ARG	3.2
1	R	139[A]	HIS	3.2
1	G	157	GLU	3.1
1	L	51	LYS	3.1
1	F	163	VAL	3.1
1	P	79	GLU	3.1
1	C	147	MET	3.1
1	U	164	LEU	3.1
1	P	163	VAL	3.1
1	E	147	MET	3.1
1	J	154	ARG	3.1
1	N	163	VAL	3.1
2	V	75	THR	3.0
1	L	1	MET	3.0
1	Q	1	MET	3.0
1	Q	162	SER	3.0
1	D	81	LYS	3.0
1	J	51	LYS	3.0
2	V	147	MET	2.9
1	L	147	MET	2.9
1	I	76	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
1	G	51	LYS	2.9
1	I	1	MET	2.9
1	D	145	ASN	2.9
1	S	162	SER	2.9
1	G	147	MET	2.9
1	R	147	MET	2.9
1	F	162	SER	2.8
1	D	148	ASP	2.8
2	V	162	SER	2.8
1	H	76	PRO	2.8
1	M	79	GLU	2.8
1	S	79	GLU	2.8
1	G	162	SER	2.7
1	I	51	LYS	2.7
1	W	51	LYS	2.7
1	E	1	MET	2.7
1	S	154	ARG	2.7
1	B	76	PRO	2.7
1	A	162	SER	2.7
1	I	163	VAL	2.7
1	F	51	LYS	2.7
1	F	76	PRO	2.6
1	R	4	LYS	2.6
1	P	77	ALA	2.6
1	L	79	GLU	2.6
1	D	162	SER	2.6
1	E	154	ARG	2.6
1	X	1	MET	2.6
1	B	163	VAL	2.5
1	G	154	ARG	2.5
1	R	51	LYS	2.5
1	Q	146	THR	2.5
1	B	161	LYS	2.5
1	I	162	SER	2.5
1	O	162	SER	2.5
1	A	147	MET	2.5
1	R	148	ASP	2.5
1	B	147	MET	2.5
1	H	163	VAL	2.4
1	U	163	VAL	2.4
1	P	147	MET	2.4
2	V	77	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	N	79	GLU	2.4
1	M	139[A]	HIS	2.4
1	C	154	ARG	2.4
1	L	154	ARG	2.4
1	K	51	LYS	2.4
1	Q	4	LYS	2.4
1	T	154	ARG	2.3
1	J	163	VAL	2.3
1	B	157	GLU	2.3
1	Q	51	LYS	2.3
1	K	162	SER	2.3
1	B	164	LEU	2.3
1	T	161	LYS	2.3
1	R	79	GLU	2.3
1	H	160	ARG	2.3
1	P	154	ARG	2.3
1	P	146[A]	THR	2.3
1	I	79	GLU	2.3
1	D	154	ARG	2.3
1	N	139	HIS	2.2
1	W	163	VAL	2.2
1	X	154	ARG	2.2
1	F	154	ARG	2.2
1	H	51	LYS	2.2
1	I	139	HIS	2.2
1	Q	147	MET	2.2
1	X	147	MET	2.2
1	I	164	LEU	2.2
1	O	79	GLU	2.2
1	N	147	MET	2.1
1	R	149	GLY	2.1
1	L	162	SER	2.1
1	N	51	LYS	2.1
1	O	163	VAL	2.1
1	X	79	GLU	2.1
1	M	1	MET	2.1
1	E	162	SER	2.1
1	U	147	MET	2.1
1	H	154	ARG	2.0
2	V	154	ARG	2.0
1	R	162	SER	2.0
1	K	139[A]	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
1	U	139[A]	HIS	2.0
1	O	110	ARG	2.0
1	Q	154	ARG	2.0
1	F	161	LYS	2.0
1	J	157	GLU	2.0
1	T	51	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](#)

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(Å ²)	Q<0.9
3	FE	O	1166	1/1	0.98	0.04	34,34,34,34	0
3	FE	P	1165	1/1	0.98	0.04	45,45,45,45	0
3	FE	V	1164	1/1	0.98	0.04	37,37,37,37	0
3	FE	B	1166	1/1	0.99	0.01	27,27,27,27	0
3	FE	C	1164	1/1	0.99	0.03	30,30,30,30	0
3	FE	D	1164	1/1	0.99	0.03	29,29,29,29	0
3	FE	D	1165	1/1	0.99	0.02	38,38,38,38	0
3	FE	E	1164	1/1	0.99	0.02	35,35,35,35	0
3	FE	F	1164	1/1	0.99	0.05	36,36,36,36	0
3	FE	F	1165	1/1	0.99	0.03	42,42,42,42	0
3	FE	G	1164	1/1	0.99	0.02	34,34,34,34	0
3	FE	G	1165	1/1	0.99	0.02	48,48,48,48	0
3	FE	H	1164	1/1	0.99	0.03	34,34,34,34	0
3	FE	H	1165	1/1	0.99	0.02	39,39,39,39	0
3	FE	I	1165	1/1	0.99	0.02	38,38,38,38	0
3	FE	J	1165	1/1	0.99	0.03	31,31,31,31	0
3	FE	K	1164	1/1	0.99	0.03	31,31,31,31	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	FE	L	1164	1/1	0.99	0.03	34,34,34,34	0
3	FE	M	1164	1/1	0.99	0.02	38,38,38,38	0
3	FE	M	1165	1/1	0.99	0.01	30,30,30,30	0
3	FE	N	1165	1/1	0.99	0.02	36,36,36,36	0
3	FE	N	1166	1/1	0.99	0.02	45,45,45,45	0
3	FE	O	1165	1/1	0.99	0.03	41,41,41,41	0
3	FE	A	1164	1/1	0.99	0.02	27,27,27,27	0
3	FE	A	1165	1/1	0.99	0.03	35,35,35,35	0
3	FE	Q	1164	1/1	0.99	0.03	39,39,39,39	0
3	FE	R	1164	1/1	0.99	0.03	32,32,32,32	0
3	FE	R	1165	1/1	0.99	0.02	35,35,35,35	0
3	FE	S	1163	1/1	0.99	0.03	36,36,36,36	0
3	FE	T	1165	1/1	0.99	0.03	39,39,39,39	0
3	FE	U	1166	1/1	0.99	0.03	31,31,31,31	0
3	FE	B	1165	1/1	0.99	0.02	39,39,39,39	0
3	FE	V	1165	1/1	0.99	0.03	32,32,32,32	0
3	FE	X	1164	1/1	0.99	0.03	21,21,21,21	0
3	FE	X	1165	1/1	0.99	0.02	31,31,31,31	0
3	FE	K	1165	1/1	1.00	0.03	24,24,24,24	0
3	FE	I	1166	1/1	1.00	0.01	32,32,32,32	0
3	FE	S	1164	1/1	1.00	0.03	27,27,27,27	0
3	FE	T	1164	1/1	1.00	0.03	27,27,27,27	0
3	FE	L	1165	1/1	1.00	0.03	25,25,25,25	0
3	FE	U	1165	1/1	1.00	0.03	26,26,26,26	0
3	FE	P	1164	1/1	1.00	0.02	32,32,32,32	0
3	FE	E	1165	1/1	1.00	0.01	30,30,30,30	0
3	FE	J	1166	1/1	1.00	0.02	26,26,26,26	0
3	FE	W	1165	1/1	1.00	0.02	22,22,22,22	0
3	FE	W	1166	1/1	1.00	0.02	33,33,33,33	0
3	FE	Q	1165	1/1	1.00	0.02	28,28,28,28	0
3	FE	C	1165	1/1	1.00	0.02	38,38,38,38	0

6.5 Other polymers [i](#)

There are no such residues in this entry.