wwPDB EM Map/Model Validation Report (i)

Apr 10, 2016 – 02:08 PM BST

PDB ID	:	4V1W
EMDB ID:	:	EMD-2788
Title	:	3D structure of horse spleen apoferritin determined by electron cryomicroscopy
Authors	:	Russo, C.J.; Passmore, L.A.
Deposited on	:	2014-10-02
Resolution	:	4.70 Å(reported)
Based on PDB ID	:	2W0O

This is a wwPDB EM Map/Model Validation Report for a publicly released PDB/EMDB entry.
For rigid body fitted models, validation errors reported here could
stem from errors in the original structure(s) used in the fitting.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at
http://wwpdb.org/validation/2016/EMValidationReportHelp

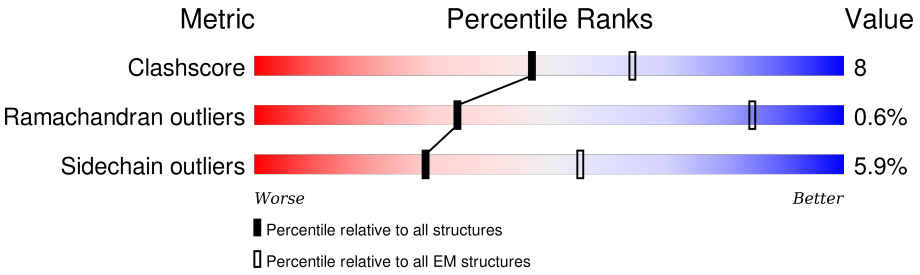
MolProbity	:	4.02b-467
Mogul	:	unknown
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk27241

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 4.70 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)	EM structures (#Entries)
Clashscore	114402	924
Ramachandran outliers	111179	726
Sidechain outliers	111093	686

The table below summarises the geometric issues observed across the polymeric chains. The red, orange, yellow and green segments on the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. 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 <=5%

Mol	Chain	Length	Quality of chain
1	A	174	75% 20% • •
1	B	174	74% 18% 5% •
1	C	174	73% 20% 5% •
1	D	174	74% 19% 5% •
1	E	174	75% 18% 5% •
1	F	174	72% 21% 5% •
1	G	174	74% 20% 5% •
1	H	174	72% 22% • •
1	I	174	74% 19% 5% •

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Mol	Chain	Length	Quality of chain
1	J	174	75% 20%
1	K	174	75% 17% 5%
1	L	174	73% 20% 5%
1	M	174	74% 20% 5%
1	N	174	74% 19% 5%
1	O	174	75% 18% 5%
1	P	174	74% 20%
1	Q	174	74% 20% 5%
1	R	174	72% 22%
1	S	174	73% 21%
1	T	174	74% 20% 5%
1	U	174	72% 21%
1	V	174	75% 19%
1	W	174	76% 17% 5%
1	X	174	74% 20% 5%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 32736 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 LIGHT CHAIN.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	B	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	C	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	D	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	E	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	F	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	G	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	H	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	I	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	J	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	K	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	L	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	M	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	N	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	O	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	P	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	Q	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		

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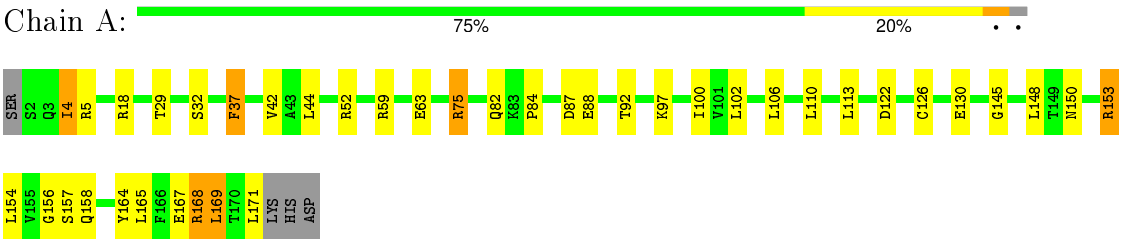
Mol	Chain	Residues	Atoms					AltConf	Trace
1	R	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	S	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	T	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	U	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	V	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	W	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		
1	X	170	Total	C	N	O	S	0	0
			1364	860	239	260	5		

3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

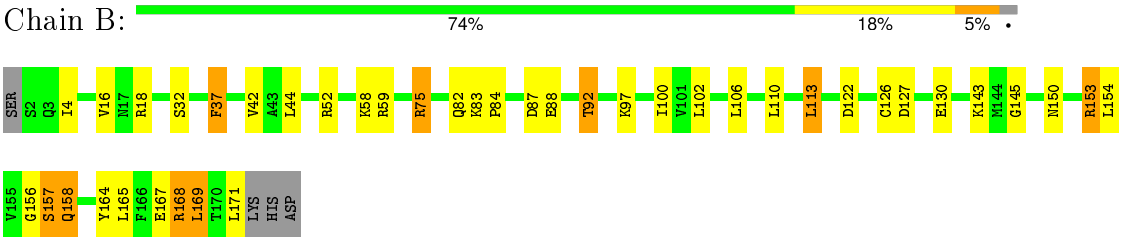
- Molecule 1: FERRITIN LIGHT CHAIN

Chain A:



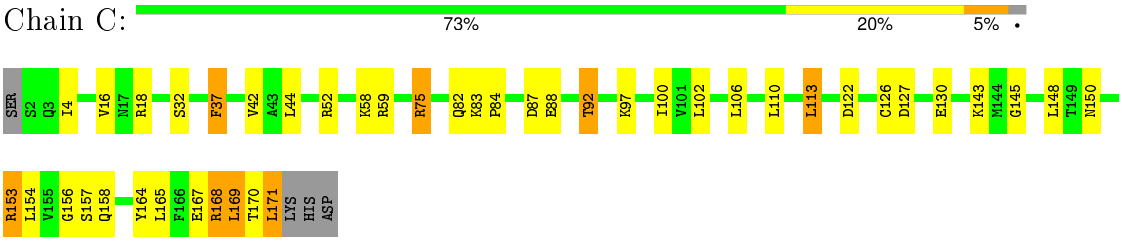
- Molecule 1: FERRITIN LIGHT CHAIN

Chain B:



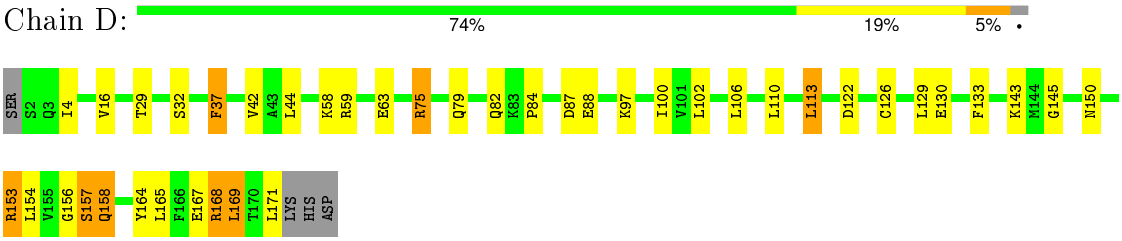
- Molecule 1: FERRITIN LIGHT CHAIN

Chain C:



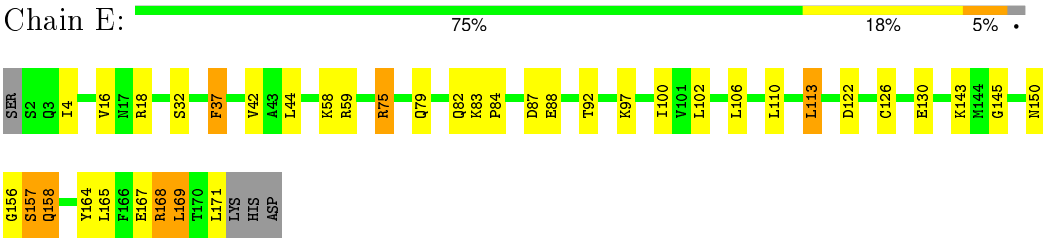
- Molecule 1: FERRITIN LIGHT CHAIN

Chain D:



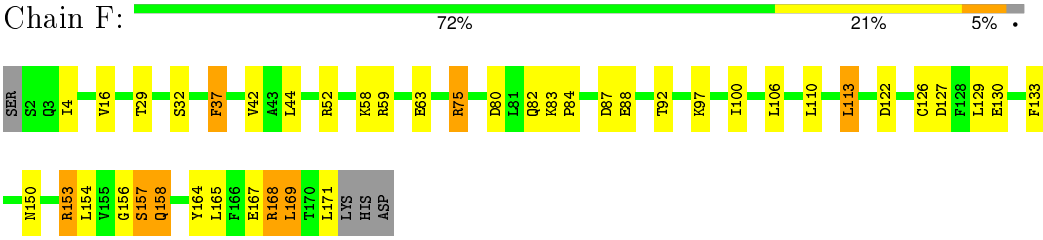
- Molecule 1: FERRITIN LIGHT CHAIN

Chain E:



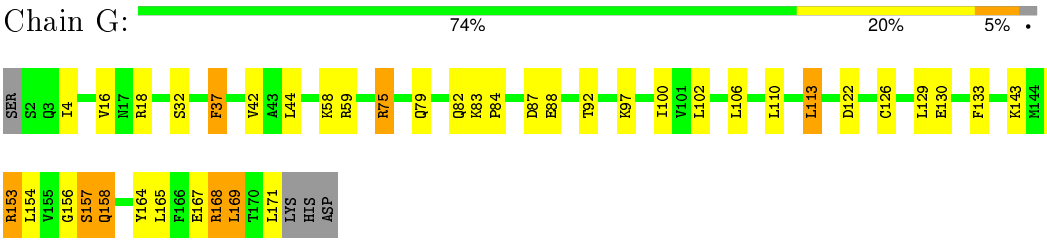
• Molecule 1: FERRITIN LIGHT CHAIN

Chain F:



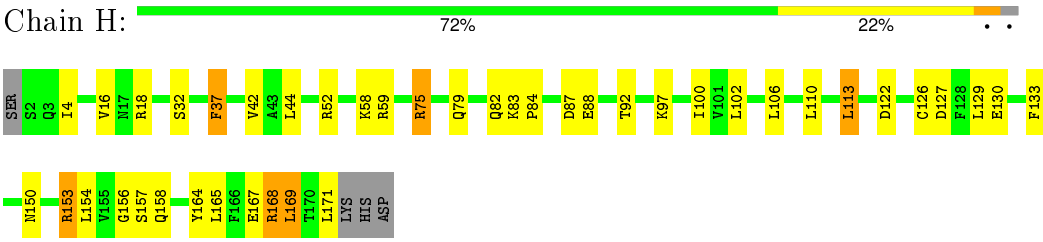
• Molecule 1: FERRITIN LIGHT CHAIN

Chain G:



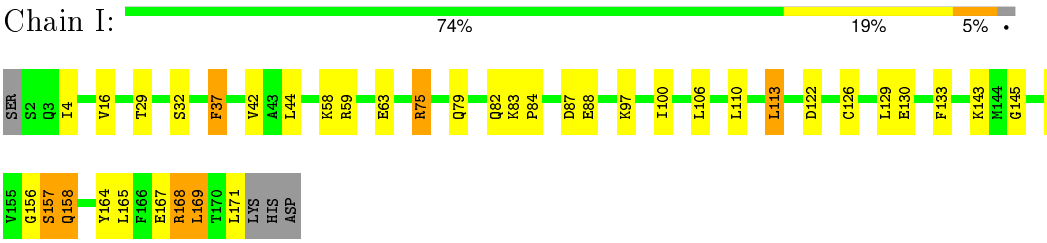
• Molecule 1: FERRITIN LIGHT CHAIN

Chain H:



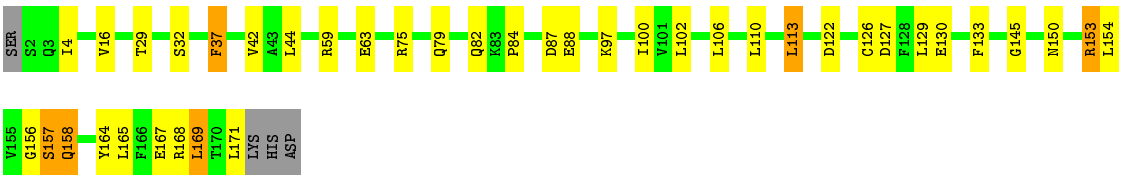
• Molecule 1: FERRITIN LIGHT CHAIN

Chain I:



• Molecule 1: FERRITIN LIGHT CHAIN

Chain J:

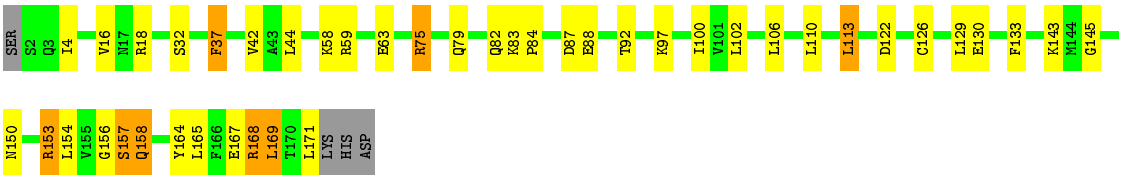


- Molecule 1: FERRITIN LIGHT CHAIN

Chain K:

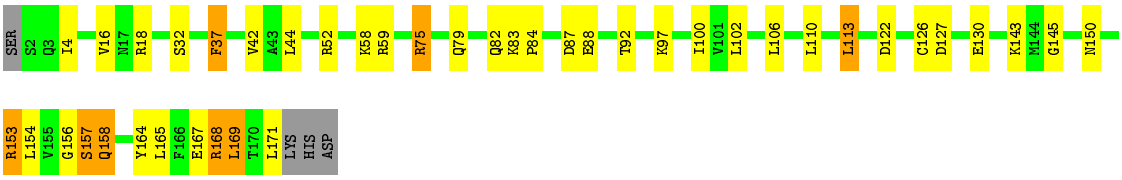
- Molecule 1: FERRITIN LIGHT CHAIN

Chain L:



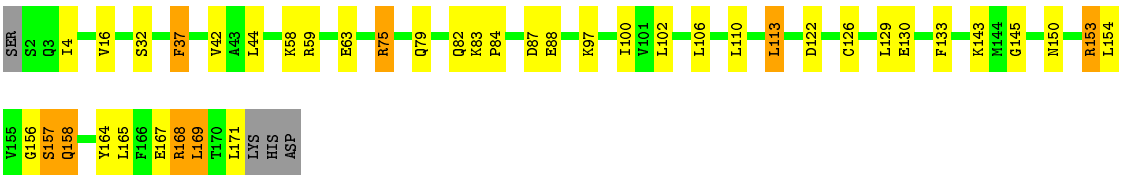
- Molecule 1: FERRITIN LIGHT CHAIN

Chain M:



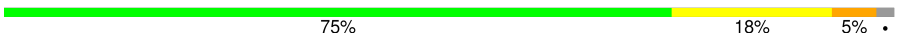
- Molecule 1: FERRITIN LIGHT CHAIN

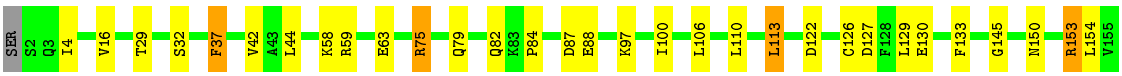
Chain N:



- Molecule 1: FERRITIN LIGHT CHAIN

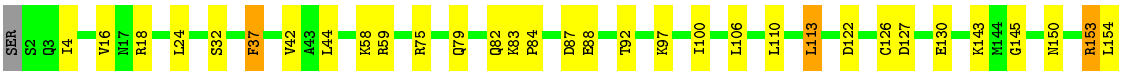
Chain O:





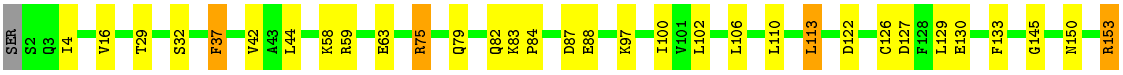
- Molecule 1: FERRITIN LIGHT CHAIN

Chain P:



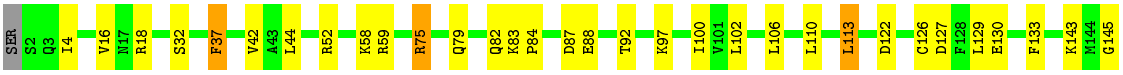
- Molecule 1: FERRITIN LIGHT CHAIN

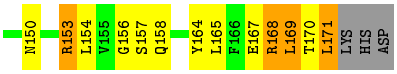
Chain Q:



- Molecule 1: FERRITIN LIGHT CHAIN

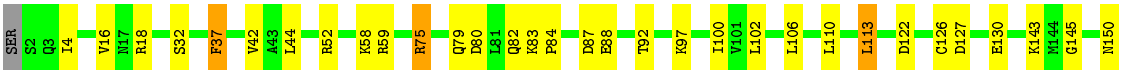
Chain R:





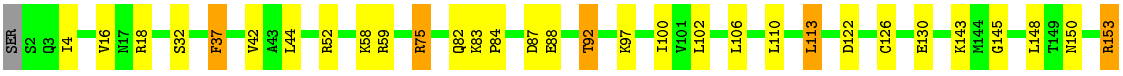
- Molecule 1: FERRITIN LIGHT CHAIN

Chain S:



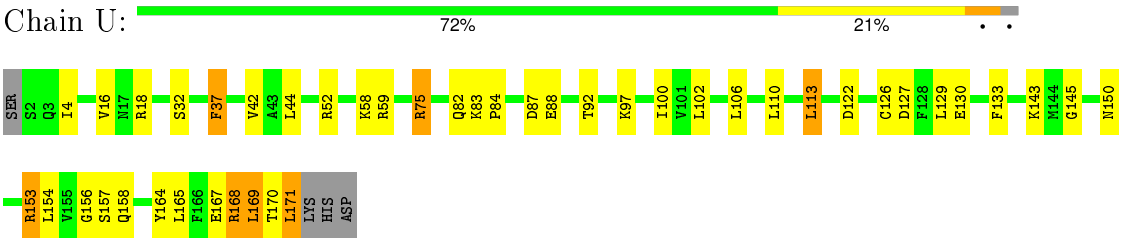
- Molecule 1: FERRITIN LIGHT CHAIN

Chain T:



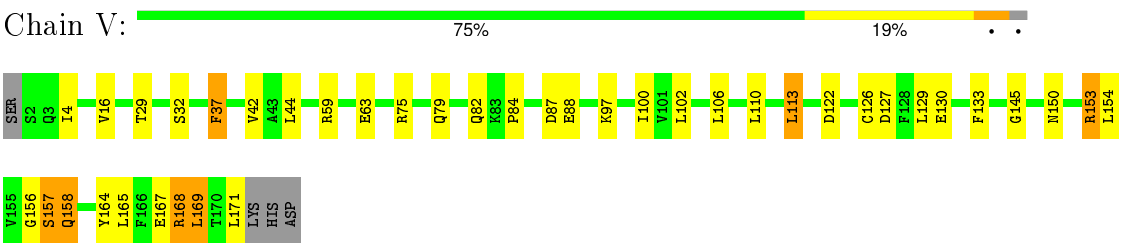
• Molecule 1: FERRITIN LIGHT CHAIN

Chain U:



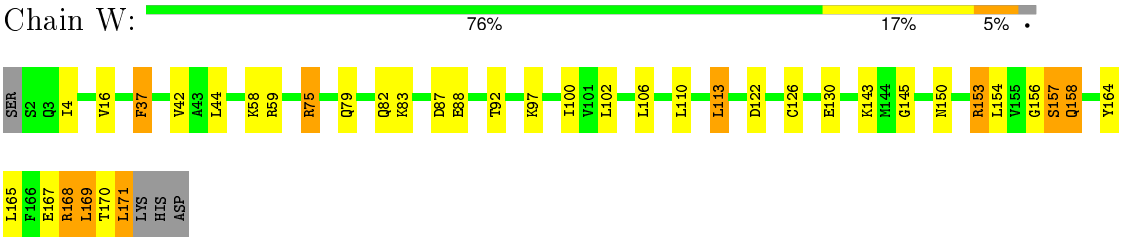
• Molecule 1: FERRITIN LIGHT CHAIN

Chain V:



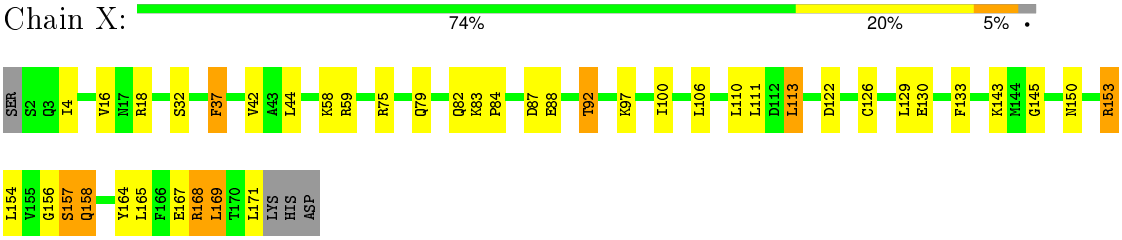
• Molecule 1: FERRITIN LIGHT CHAIN

Chain W:



• Molecule 1: FERRITIN LIGHT CHAIN

Chain X:



4 Experimental information (i)

Property	Value	Source
Reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	Depositor
Number of images	Not provided	Depositor
Resolution determination method	Not provided	Depositor
CTF correction method	PER PARTICLE	Depositor
Microscope	OTHER	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	Not provided	Depositor
Minimum defocus (nm)	1573	Depositor
Maximum defocus (nm)	3639	Depositor
Magnification	59000	Depositor
Image detector	FEI FALCON II (4K X 4K)	Depositor

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	#$\|Z\|$ >2	RMSZ	#$\|Z\|$ >2
1	A	0.98	1/1386 (0.1%)	1.26	11/1866 (0.6%)
1	B	0.98	1/1386 (0.1%)	1.25	10/1866 (0.5%)
1	C	0.98	1/1386 (0.1%)	1.26	11/1866 (0.6%)
1	D	0.98	1/1386 (0.1%)	1.26	9/1866 (0.5%)
1	E	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	F	0.98	2/1386 (0.1%)	1.27	11/1866 (0.6%)
1	G	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	H	0.98	1/1386 (0.1%)	1.26	11/1866 (0.6%)
1	I	0.98	1/1386 (0.1%)	1.27	9/1866 (0.5%)
1	J	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	K	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	L	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	M	0.98	1/1386 (0.1%)	1.26	11/1866 (0.6%)
1	N	0.97	1/1386 (0.1%)	1.26	9/1866 (0.5%)
1	O	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	P	0.98	1/1386 (0.1%)	1.26	11/1866 (0.6%)
1	Q	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	R	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	S	0.98	2/1386 (0.1%)	1.26	10/1866 (0.5%)
1	T	0.98	1/1386 (0.1%)	1.25	10/1866 (0.5%)
1	U	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	V	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	W	0.98	1/1386 (0.1%)	1.26	10/1866 (0.5%)
1	X	0.98	1/1386 (0.1%)	1.26	11/1866 (0.6%)
All	All	0.98	26/33264 (0.1%)	1.26	244/44784 (0.5%)

All (26) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	I	37	PHE	CB-CG	-5.21	1.42	1.51
1	V	37	PHE	CB-CG	-5.21	1.42	1.51
1	G	37	PHE	CB-CG	-5.17	1.42	1.51
1	P	37	PHE	CB-CG	-5.17	1.42	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	M	37	PHE	CB-CG	-5.17	1.42	1.51
1	Q	37	PHE	CB-CG	-5.16	1.42	1.51
1	E	37	PHE	CB-CG	-5.16	1.42	1.51
1	K	37	PHE	CB-CG	-5.16	1.42	1.51
1	W	37	PHE	CB-CG	-5.16	1.42	1.51
1	R	37	PHE	CB-CG	-5.16	1.42	1.51
1	X	37	PHE	CB-CG	-5.16	1.42	1.51
1	U	37	PHE	CB-CG	-5.16	1.42	1.51
1	F	80	ASP	CB-CG	5.15	1.62	1.51
1	J	37	PHE	CB-CG	-5.13	1.42	1.51
1	F	37	PHE	CB-CG	-5.13	1.42	1.51
1	D	37	PHE	CB-CG	-5.12	1.42	1.51
1	C	37	PHE	CB-CG	-5.12	1.42	1.51
1	L	37	PHE	CB-CG	-5.12	1.42	1.51
1	O	37	PHE	CB-CG	-5.12	1.42	1.51
1	N	37	PHE	CB-CG	-5.12	1.42	1.51
1	T	37	PHE	CB-CG	-5.12	1.42	1.51
1	H	37	PHE	CB-CG	-5.09	1.42	1.51
1	B	37	PHE	CB-CG	-5.08	1.42	1.51
1	S	37	PHE	CB-CG	-5.08	1.42	1.51
1	A	37	PHE	CB-CG	-5.04	1.42	1.51
1	S	80	ASP	CB-CG	5.01	1.62	1.51

All (244) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	I	59	ARG	NE-CZ-NH1	9.94	125.27	120.30
1	V	59	ARG	NE-CZ-NH1	9.73	125.17	120.30
1	Q	59	ARG	NE-CZ-NH1	9.60	125.10	120.30
1	J	59	ARG	NE-CZ-NH1	9.52	125.06	120.30
1	D	59	ARG	NE-CZ-NH1	9.51	125.06	120.30
1	F	59	ARG	NE-CZ-NH1	9.49	125.05	120.30
1	N	59	ARG	NE-CZ-NH1	9.46	125.03	120.30
1	X	59	ARG	NE-CZ-NH1	9.39	125.00	120.30
1	L	59	ARG	NE-CZ-NH1	9.38	124.99	120.30
1	O	59	ARG	NE-CZ-NH1	9.34	124.97	120.30
1	E	59	ARG	NE-CZ-NH1	9.26	124.93	120.30
1	M	59	ARG	NE-CZ-NH1	9.16	124.88	120.30
1	W	59	ARG	NE-CZ-NH1	9.12	124.86	120.30
1	G	59	ARG	NE-CZ-NH1	9.05	124.83	120.30
1	P	59	ARG	NE-CZ-NH1	8.97	124.79	120.30
1	K	59	ARG	NE-CZ-NH1	8.91	124.76	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	S	59	ARG	NE-CZ-NH1	8.91	124.75	120.30
1	U	59	ARG	NE-CZ-NH1	8.83	124.72	120.30
1	T	59	ARG	NE-CZ-NH1	8.77	124.69	120.30
1	B	59	ARG	NE-CZ-NH1	8.74	124.67	120.30
1	H	59	ARG	NE-CZ-NH1	8.74	124.67	120.30
1	C	59	ARG	NE-CZ-NH1	8.68	124.64	120.30
1	R	59	ARG	NE-CZ-NH1	8.52	124.56	120.30
1	A	59	ARG	NE-CZ-NH1	8.51	124.56	120.30
1	K	169	LEU	CB-CG-CD1	-7.10	98.94	111.00
1	V	169	LEU	CB-CG-CD1	-6.68	99.64	111.00
1	U	169	LEU	CB-CG-CD2	6.63	122.27	111.00
1	H	169	LEU	CB-CG-CD2	6.58	122.18	111.00
1	D	169	LEU	CB-CG-CD1	-6.50	99.94	111.00
1	O	153	ARG	NE-CZ-NH1	6.49	123.55	120.30
1	F	169	LEU	CB-CG-CD2	6.47	122.00	111.00
1	F	169	LEU	CB-CG-CD1	-6.45	100.03	111.00
1	I	153	ARG	NE-CZ-NH1	6.45	123.52	120.30
1	J	169	LEU	CB-CG-CD1	-6.38	100.15	111.00
1	O	169	LEU	CB-CG-CD2	6.38	121.84	111.00
1	J	169	LEU	CB-CG-CD2	6.37	121.83	111.00
1	B	169	LEU	CB-CG-CD2	6.34	121.78	111.00
1	J	153	ARG	NE-CZ-NH1	6.29	123.44	120.30
1	E	169	LEU	CB-CG-CD2	6.28	121.67	111.00
1	S	169	LEU	CB-CG-CD2	6.27	121.66	111.00
1	X	153	ARG	NE-CZ-NH1	6.26	123.43	120.30
1	D	153	ARG	NE-CZ-NH1	6.23	123.42	120.30
1	P	153	ARG	NE-CZ-NH1	6.23	123.42	120.30
1	A	169	LEU	CB-CG-CD1	-6.22	100.43	111.00
1	N	153	ARG	NE-CZ-NH1	6.21	123.41	120.30
1	L	169	LEU	CB-CG-CD2	6.20	121.54	111.00
1	V	153	ARG	NE-CZ-NH1	6.19	123.39	120.30
1	F	153	ARG	NE-CZ-NH1	6.17	123.38	120.30
1	Q	153	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	M	153	ARG	NE-CZ-NH1	6.15	123.38	120.30
1	X	169	LEU	CB-CG-CD2	6.14	121.44	111.00
1	G	169	LEU	CB-CG-CD1	-6.13	100.57	111.00
1	N	169	LEU	CB-CG-CD1	-6.13	100.58	111.00
1	M	169	LEU	CB-CG-CD2	6.13	121.42	111.00
1	Q	169	LEU	CB-CG-CD1	-6.11	100.62	111.00
1	I	169	LEU	CB-CG-CD2	6.10	121.37	111.00
1	N	169	LEU	CB-CG-CD2	6.10	121.37	111.00
1	U	153	ARG	NE-CZ-NH1	6.10	123.35	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	R	169	LEU	CB-CG-CD2	6.09	121.36	111.00
1	W	169	LEU	CB-CG-CD1	-6.08	100.67	111.00
1	C	153	ARG	NE-CZ-NH1	6.07	123.33	120.30
1	G	169	LEU	CB-CG-CD2	6.06	121.30	111.00
1	W	153	ARG	NE-CZ-NH1	6.04	123.32	120.30
1	G	153	ARG	NE-CZ-NH1	6.03	123.32	120.30
1	R	153	ARG	NE-CZ-NH1	6.03	123.31	120.30
1	T	153	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	V	169	LEU	CB-CG-CD2	6.02	121.24	111.00
1	P	169	LEU	CB-CG-CD2	6.01	121.21	111.00
1	W	169	LEU	CB-CG-CD2	5.94	121.10	111.00
1	O	169	LEU	CB-CG-CD1	-5.94	100.91	111.00
1	S	122	ASP	CB-CG-OD1	5.93	123.64	118.30
1	S	153	ARG	NE-CZ-NH1	5.93	123.27	120.30
1	U	122	ASP	CB-CG-OD1	5.93	123.64	118.30
1	I	122	ASP	CB-CG-OD1	5.93	123.64	118.30
1	A	122	ASP	CB-CG-OD1	5.93	123.64	118.30
1	L	153	ARG	NE-CZ-NH1	5.92	123.26	120.30
1	C	122	ASP	CB-CG-OD1	5.92	123.63	118.30
1	R	122	ASP	CB-CG-OD1	5.91	123.62	118.30
1	Q	169	LEU	CB-CG-CD2	5.90	121.03	111.00
1	K	169	LEU	CB-CG-CD2	5.89	121.02	111.00
1	B	122	ASP	CB-CG-OD1	5.89	123.60	118.30
1	F	122	ASP	CB-CG-OD1	5.89	123.60	118.30
1	H	122	ASP	CB-CG-OD1	5.88	123.59	118.30
1	T	122	ASP	CB-CG-OD1	5.88	123.59	118.30
1	K	153	ARG	NE-CZ-NH1	5.87	123.24	120.30
1	H	153	ARG	NE-CZ-NH1	5.87	123.23	120.30
1	O	122	ASP	CB-CG-OD1	5.86	123.57	118.30
1	G	122	ASP	CB-CG-OD1	5.85	123.57	118.30
1	D	122	ASP	CB-CG-OD1	5.85	123.56	118.30
1	T	169	LEU	CB-CG-CD1	-5.85	101.06	111.00
1	K	122	ASP	CB-CG-OD1	5.85	123.56	118.30
1	M	122	ASP	CB-CG-OD1	5.84	123.56	118.30
1	J	122	ASP	CB-CG-OD1	5.84	123.56	118.30
1	L	122	ASP	CB-CG-OD1	5.84	123.56	118.30
1	X	122	ASP	CB-CG-OD1	5.84	123.56	118.30
1	E	122	ASP	CB-CG-OD1	5.83	123.55	118.30
1	V	122	ASP	CB-CG-OD1	5.82	123.54	118.30
1	D	169	LEU	CB-CG-CD2	5.81	120.88	111.00
1	N	122	ASP	CB-CG-OD1	5.81	123.53	118.30
1	E	153	ARG	NE-CZ-NH1	5.80	123.20	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	P	122	ASP	CB-CG-OD1	5.79	123.51	118.30
1	I	169	LEU	CB-CG-CD1	-5.79	101.16	111.00
1	Q	122	ASP	CB-CG-OD1	5.79	123.51	118.30
1	E	169	LEU	CB-CG-CD1	-5.78	101.17	111.00
1	V	113	LEU	CB-CG-CD1	5.77	120.81	111.00
1	Q	113	LEU	CB-CG-CD1	5.75	120.77	111.00
1	W	122	ASP	CB-CG-OD1	5.74	123.47	118.30
1	A	153	ARG	NE-CZ-NH1	5.73	123.17	120.30
1	B	113	LEU	CB-CG-CD1	5.71	120.70	111.00
1	R	113	LEU	CB-CG-CD1	5.70	120.70	111.00
1	T	113	LEU	CB-CG-CD1	5.70	120.69	111.00
1	G	113	LEU	CB-CG-CD1	5.69	120.67	111.00
1	K	113	LEU	CB-CG-CD1	5.68	120.65	111.00
1	A	169	LEU	CB-CG-CD2	5.67	120.64	111.00
1	D	113	LEU	CB-CG-CD1	5.66	120.63	111.00
1	A	110	LEU	CA-CB-CG	5.66	128.32	115.30
1	J	113	LEU	CB-CG-CD1	5.65	120.61	111.00
1	C	113	LEU	CB-CG-CD1	5.65	120.61	111.00
1	F	113	LEU	CB-CG-CD1	5.64	120.59	111.00
1	S	113	LEU	CB-CG-CD1	5.64	120.59	111.00
1	K	75	ARG	NE-CZ-NH1	5.63	123.11	120.30
1	N	113	LEU	CB-CG-CD1	5.63	120.57	111.00
1	B	153	ARG	NE-CZ-NH1	5.62	123.11	120.30
1	W	113	LEU	CB-CG-CD1	5.62	120.55	111.00
1	V	75	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	O	75	ARG	NE-CZ-NH1	5.61	123.11	120.30
1	U	113	LEU	CB-CG-CD1	5.61	120.54	111.00
1	I	75	ARG	NE-CZ-NH1	5.61	123.10	120.30
1	I	113	LEU	CB-CG-CD1	5.61	120.53	111.00
1	P	113	LEU	CB-CG-CD1	5.61	120.53	111.00
1	H	113	LEU	CB-CG-CD1	5.61	120.53	111.00
1	E	113	LEU	CB-CG-CD1	5.60	120.52	111.00
1	X	75	ARG	NE-CZ-NH1	5.60	123.10	120.30
1	M	113	LEU	CB-CG-CD1	5.60	120.51	111.00
1	L	113	LEU	CB-CG-CD1	5.59	120.50	111.00
1	O	113	LEU	CB-CG-CD1	5.58	120.49	111.00
1	G	75	ARG	NE-CZ-NH1	5.58	123.09	120.30
1	M	75	ARG	NE-CZ-NH1	5.58	123.09	120.30
1	T	169	LEU	CB-CG-CD2	5.58	120.48	111.00
1	X	106	LEU	CB-CG-CD1	-5.57	101.53	111.00
1	Q	75	ARG	NE-CZ-NH1	5.57	123.08	120.30
1	C	169	LEU	CB-CG-CD2	5.57	120.47	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	L	75	ARG	NE-CZ-NH1	5.57	123.08	120.30
1	W	75	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	U	110	LEU	CA-CB-CG	5.56	128.09	115.30
1	P	75	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	R	110	LEU	CA-CB-CG	5.55	128.06	115.30
1	B	110	LEU	CA-CB-CG	5.55	128.06	115.30
1	D	75	ARG	NE-CZ-NH1	5.54	123.07	120.30
1	S	110	LEU	CA-CB-CG	5.53	128.03	115.30
1	I	110	LEU	CA-CB-CG	5.53	128.02	115.30
1	H	110	LEU	CA-CB-CG	5.53	128.01	115.30
1	O	110	LEU	CA-CB-CG	5.52	127.99	115.30
1	L	169	LEU	CB-CG-CD1	-5.51	101.64	111.00
1	L	106	LEU	CB-CG-CD1	-5.51	101.64	111.00
1	C	110	LEU	CA-CB-CG	5.50	127.96	115.30
1	D	110	LEU	CA-CB-CG	5.50	127.95	115.30
1	T	110	LEU	CA-CB-CG	5.49	127.92	115.30
1	R	75	ARG	NE-CZ-NH1	5.48	123.04	120.30
1	K	110	LEU	CA-CB-CG	5.47	127.89	115.30
1	N	75	ARG	NE-CZ-NH1	5.47	123.04	120.30
1	J	75	ARG	NE-CZ-NH1	5.47	123.04	120.30
1	Q	110	LEU	CA-CB-CG	5.47	127.88	115.30
1	F	75	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	A	75	ARG	NE-CZ-NH1	5.46	123.03	120.30
1	I	106	LEU	CB-CG-CD1	-5.46	101.72	111.00
1	P	110	LEU	CA-CB-CG	5.46	127.85	115.30
1	M	106	LEU	CB-CG-CD1	-5.45	101.73	111.00
1	X	110	LEU	CA-CB-CG	5.45	127.83	115.30
1	J	110	LEU	CA-CB-CG	5.44	127.82	115.30
1	N	110	LEU	CA-CB-CG	5.44	127.81	115.30
1	R	106	LEU	CB-CG-CD1	-5.44	101.75	111.00
1	S	75	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	B	75	ARG	NE-CZ-NH1	5.43	123.02	120.30
1	E	110	LEU	CA-CB-CG	5.43	127.80	115.30
1	W	110	LEU	CA-CB-CG	5.43	127.78	115.30
1	V	110	LEU	CA-CB-CG	5.42	127.78	115.30
1	A	106	LEU	CB-CG-CD1	-5.42	101.78	111.00
1	C	169	LEU	CB-CG-CD1	-5.42	101.78	111.00
1	E	75	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	G	110	LEU	CA-CB-CG	5.42	127.76	115.30
1	K	106	LEU	CB-CG-CD1	-5.42	101.79	111.00
1	L	110	LEU	CA-CB-CG	5.41	127.75	115.30
1	V	106	LEU	CB-CG-CD1	-5.41	101.80	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	W	106	LEU	CB-CG-CD1	-5.41	101.80	111.00
1	X	169	LEU	CB-CG-CD1	-5.41	101.80	111.00
1	F	106	LEU	CB-CG-CD1	-5.41	101.80	111.00
1	T	75	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	C	75	ARG	NE-CZ-NH1	5.41	123.00	120.30
1	G	106	LEU	CB-CG-CD1	-5.41	101.81	111.00
1	X	113	LEU	CB-CG-CD1	5.41	120.19	111.00
1	E	106	LEU	CB-CG-CD1	-5.40	101.82	111.00
1	P	106	LEU	CB-CG-CD1	-5.40	101.82	111.00
1	T	106	LEU	CB-CG-CD1	-5.39	101.83	111.00
1	A	113	LEU	CB-CG-CD1	5.39	120.16	111.00
1	F	110	LEU	CA-CB-CG	5.39	127.69	115.30
1	B	106	LEU	CB-CG-CD1	-5.38	101.85	111.00
1	Q	106	LEU	CB-CG-CD1	-5.38	101.85	111.00
1	M	110	LEU	CA-CB-CG	5.37	127.66	115.30
1	O	106	LEU	CB-CG-CD1	-5.37	101.87	111.00
1	U	75	ARG	NE-CZ-NH1	5.37	122.98	120.30
1	D	106	LEU	CB-CG-CD1	-5.36	101.89	111.00
1	C	92	THR	N-CA-CB	5.36	120.48	110.30
1	C	106	LEU	CB-CG-CD1	-5.36	101.89	111.00
1	H	106	LEU	CB-CG-CD1	-5.36	101.89	111.00
1	S	106	LEU	CB-CG-CD1	-5.35	101.91	111.00
1	M	169	LEU	CB-CG-CD1	-5.33	101.94	111.00
1	J	106	LEU	CB-CG-CD1	-5.33	101.95	111.00
1	U	106	LEU	CB-CG-CD1	-5.32	101.96	111.00
1	N	106	LEU	CB-CG-CD1	-5.31	101.98	111.00
1	A	92	THR	N-CA-CB	5.27	120.31	110.30
1	W	92	THR	N-CA-CB	5.23	120.24	110.30
1	H	75	ARG	NE-CZ-NH1	5.17	122.89	120.30
1	P	169	LEU	CB-CG-CD1	-5.16	102.23	111.00
1	S	127	ASP	CB-CG-OD1	5.16	122.94	118.30
1	B	127	ASP	CB-CG-OD1	5.15	122.94	118.30
1	X	111	LEU	CB-CG-CD1	-5.14	102.27	111.00
1	A	5	ARG	NE-CZ-NH2	-5.13	117.73	120.30
1	U	92	THR	N-CA-CB	5.13	120.05	110.30
1	F	127	ASP	CB-CG-OD1	5.12	122.91	118.30
1	M	92	THR	N-CA-CB	5.12	120.03	110.30
1	T	92	THR	N-CA-CB	5.12	120.03	110.30
1	U	127	ASP	CB-CG-OD1	5.11	122.90	118.30
1	H	127	ASP	CB-CG-OD1	5.10	122.89	118.30
1	B	92	THR	N-CA-CB	5.09	119.98	110.30
1	R	92	THR	N-CA-CB	5.09	119.97	110.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	J	127	ASP	CB-CG-OD1	5.08	122.88	118.30
1	X	92	THR	N-CA-CB	5.08	119.94	110.30
1	P	92	THR	N-CA-CB	5.07	119.93	110.30
1	Q	127	ASP	CB-CG-OD1	5.07	122.86	118.30
1	S	92	THR	N-CA-CB	5.07	119.93	110.30
1	G	92	THR	N-CA-CB	5.06	119.92	110.30
1	H	92	THR	N-CA-CB	5.06	119.92	110.30
1	K	92	THR	N-CA-CB	5.06	119.91	110.30
1	V	127	ASP	CB-CG-OD1	5.06	122.85	118.30
1	R	127	ASP	CB-CG-OD1	5.04	122.84	118.30
1	E	92	THR	N-CA-CB	5.03	119.86	110.30
1	C	127	ASP	CB-CG-OD1	5.02	122.82	118.30
1	M	127	ASP	CB-CG-OD1	5.01	122.81	118.30
1	O	127	ASP	CB-CG-OD1	5.01	122.81	118.30
1	H	169	LEU	CB-CG-CD1	-5.01	102.48	111.00
1	L	92	THR	N-CA-CB	5.01	119.81	110.30
1	F	92	THR	N-CA-CB	5.00	119.81	110.30
1	P	127	ASP	CB-CG-OD1	5.00	122.80	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1364	0	1345	27	0
1	B	1364	0	1345	29	0
1	C	1364	0	1345	26	0
1	D	1364	0	1345	28	0
1	E	1364	0	1345	24	0
1	F	1364	0	1345	29	0
1	G	1364	0	1345	25	0
1	H	1364	0	1345	26	0
1	I	1364	0	1345	27	0
1	J	1364	0	1345	27	0
1	K	1364	0	1345	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	L	1364	0	1345	24	0
1	M	1364	0	1345	25	0
1	N	1364	0	1345	29	0
1	O	1364	0	1345	27	0
1	P	1364	0	1345	23	0
1	Q	1364	0	1345	30	0
1	R	1364	0	1345	28	0
1	S	1364	0	1345	27	0
1	T	1364	0	1345	25	0
1	U	1364	0	1345	26	0
1	V	1364	0	1345	24	0
1	W	1364	0	1345	25	0
1	X	1364	0	1345	22	0
All	All	32736	0	32280	506	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 8.

All (506) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:154:LEU:HD21	1:H:164:TYR:HB3	1.77	0.67
1:S:164:TYR:HB3	1:T:154:LEU:HD21	1.76	0.66
1:R:154:LEU:HD21	1:U:164:TYR:HB3	1.78	0.66
1:O:164:TYR:HB3	1:W:154:LEU:HD21	1.79	0.65
1:L:164:TYR:HB3	1:Q:154:LEU:HD21	1.79	0.65
1:T:164:TYR:HB3	1:U:154:LEU:HD21	1.79	0.64
1:F:154:LEU:HD21	1:G:164:TYR:HB3	1.80	0.64
1:R:164:TYR:HB3	1:S:154:LEU:HD21	1.78	0.64
1:C:164:TYR:HB3	1:H:154:LEU:HD21	1.79	0.64
1:S:164:TYR:CZ	1:S:168:ARG:HG3	2.33	0.64
1:G:154:LEU:HD21	1:V:164:TYR:HB3	1.80	0.64
1:F:164:TYR:HB3	1:X:154:LEU:HD21	1.80	0.63
1:E:164:TYR:HB3	1:N:154:LEU:HD21	1.80	0.63
1:B:164:TYR:HB3	1:C:154:LEU:HD21	1.78	0.63
1:V:154:LEU:HD21	1:X:164:TYR:HB3	1.79	0.63
1:J:154:LEU:HD21	1:P:164:TYR:HB3	1.80	0.63
1:B:164:TYR:CZ	1:B:168:ARG:HG3	2.34	0.63
1:P:154:LEU:HD21	1:Q:164:TYR:HB3	1.81	0.62
1:I:164:TYR:HB3	1:K:154:LEU:HD21	1.81	0.62
1:K:164:TYR:HB3	1:O:154:LEU:HD21	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:164:TYR:CZ	1:H:168:ARG:HG3	2.34	0.62
1:R:164:TYR:CZ	1:R:168:ARG:HG3	2.35	0.62
1:J:164:TYR:HB3	1:L:154:LEU:HD21	1.82	0.62
1:W:164:TYR:CZ	1:W:168:ARG:HG3	2.35	0.62
1:I:164:TYR:CZ	1:I:168:ARG:HG3	2.34	0.62
1:D:154:LEU:HD21	1:M:164:TYR:HB3	1.81	0.62
1:A:164:TYR:HB3	1:B:154:LEU:HD21	1.80	0.61
1:O:164:TYR:CZ	1:O:168:ARG:HG3	2.35	0.61
1:U:164:TYR:CZ	1:U:168:ARG:HG3	2.35	0.61
1:C:164:TYR:CZ	1:C:168:ARG:HG3	2.36	0.61
1:F:164:TYR:CZ	1:F:168:ARG:HG3	2.36	0.61
1:K:164:TYR:CZ	1:K:168:ARG:HG3	2.36	0.61
1:D:164:TYR:HB3	1:E:154:LEU:HD21	1.82	0.61
1:M:154:LEU:HD21	1:N:164:TYR:HB3	1.82	0.61
1:R:154:LEU:HD21	1:U:164:TYR:CB	2.31	0.60
1:L:164:TYR:CZ	1:L:168:ARG:HG3	2.36	0.60
1:E:164:TYR:CZ	1:E:168:ARG:HG3	2.37	0.60
1:X:164:TYR:CZ	1:X:168:ARG:HG3	2.36	0.60
1:I:154:LEU:HD21	1:W:164:TYR:HB3	1.83	0.60
1:Q:37:PHE:CD1	1:Q:42:VAL:HG11	2.37	0.60
1:X:37:PHE:CD1	1:X:42:VAL:HG11	2.36	0.60
1:O:37:PHE:CD1	1:O:42:VAL:HG11	2.37	0.60
1:H:52:ARG:HH22	1:I:63:GLU:CG	2.13	0.60
1:M:37:PHE:CD1	1:M:42:VAL:HG11	2.37	0.60
1:T:164:TYR:CZ	1:T:168:ARG:HG3	2.37	0.60
1:E:37:PHE:CD1	1:E:42:VAL:HG11	2.37	0.60
1:N:37:PHE:CD1	1:N:42:VAL:HG11	2.37	0.60
1:M:164:TYR:CZ	1:M:168:ARG:HG3	2.37	0.60
1:W:37:PHE:CD1	1:W:42:VAL:HG11	2.36	0.60
1:P:164:TYR:CZ	1:P:168:ARG:HG3	2.36	0.60
1:S:37:PHE:CD1	1:S:42:VAL:HG11	2.37	0.60
1:S:164:TYR:CB	1:T:154:LEU:HD21	2.31	0.59
1:T:164:TYR:CB	1:U:154:LEU:HD21	2.32	0.59
1:R:164:TYR:CB	1:S:154:LEU:HD21	2.31	0.59
1:G:164:TYR:CZ	1:G:168:ARG:HG3	2.37	0.59
1:J:164:TYR:CZ	1:J:168:ARG:HG3	2.36	0.59
1:A:164:TYR:CZ	1:A:168:ARG:HG3	2.37	0.59
1:L:37:PHE:CD1	1:L:42:VAL:HG11	2.36	0.59
1:C:37:PHE:CD1	1:C:42:VAL:HG11	2.37	0.59
1:A:37:PHE:CD1	1:A:42:VAL:HG11	2.37	0.59
1:P:37:PHE:CD1	1:P:42:VAL:HG11	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:V:37:PHE:CD1	1:V:42:VAL:HG11	2.37	0.59
1:Q:164:TYR:CZ	1:Q:168:ARG:HG3	2.37	0.59
1:J:37:PHE:CD1	1:J:42:VAL:HG11	2.37	0.59
1:G:154:LEU:HD21	1:V:164:TYR:CB	2.33	0.59
1:F:37:PHE:CD1	1:F:42:VAL:HG11	2.38	0.59
1:B:37:PHE:CD1	1:B:42:VAL:HG11	2.38	0.59
1:Q:63:GLU:CG	1:R:52:ARG:HH22	2.16	0.59
1:I:37:PHE:CD1	1:I:42:VAL:HG11	2.37	0.59
1:D:37:PHE:CD1	1:D:42:VAL:HG11	2.37	0.59
1:N:164:TYR:CZ	1:N:168:ARG:HG3	2.37	0.59
1:A:154:LEU:HD21	1:H:164:TYR:CB	2.33	0.59
1:O:164:TYR:CB	1:W:154:LEU:HD21	2.33	0.59
1:D:164:TYR:CZ	1:D:168:ARG:HG3	2.37	0.59
1:R:37:PHE:CD1	1:R:42:VAL:HG11	2.38	0.59
1:T:37:PHE:CD1	1:T:42:VAL:HG11	2.38	0.59
1:J:154:LEU:HD21	1:P:164:TYR:CB	2.33	0.59
1:F:164:TYR:CB	1:X:154:LEU:HD21	2.32	0.58
1:G:37:PHE:CD1	1:G:42:VAL:HG11	2.37	0.58
1:H:37:PHE:CD1	1:H:42:VAL:HG11	2.37	0.58
1:U:37:PHE:CD1	1:U:42:VAL:HG11	2.37	0.58
1:K:37:PHE:CD1	1:K:42:VAL:HG11	2.38	0.58
1:B:164:TYR:CB	1:C:154:LEU:HD21	2.32	0.58
1:F:154:LEU:HD21	1:G:164:TYR:CB	2.33	0.58
1:V:164:TYR:CZ	1:V:168:ARG:HG3	2.37	0.58
1:P:154:LEU:HD21	1:Q:164:TYR:CB	2.33	0.58
1:H:52:ARG:NH2	1:I:63:GLU:HG2	2.19	0.58
1:V:154:LEU:HD21	1:X:164:TYR:CB	2.33	0.58
1:K:164:TYR:CB	1:O:154:LEU:HD21	2.33	0.58
1:L:164:TYR:CB	1:Q:154:LEU:HD21	2.34	0.57
1:N:79:GLN:HA	1:S:83:LYS:HD2	1.86	0.57
1:J:164:TYR:CB	1:L:154:LEU:HD21	2.35	0.57
1:E:164:TYR:CB	1:N:154:LEU:HD21	2.35	0.57
1:D:164:TYR:CB	1:E:154:LEU:HD21	2.33	0.57
1:Q:79:GLN:HA	1:R:83:LYS:HD2	1.87	0.57
1:C:83:LYS:HD2	1:J:79:GLN:HA	1.87	0.57
1:O:79:GLN:HA	1:U:83:LYS:HD2	1.85	0.57
1:C:164:TYR:CB	1:H:154:LEU:HD21	2.35	0.56
1:I:164:TYR:CB	1:K:154:LEU:HD21	2.35	0.56
1:V:44:LEU:HB3	1:V:167:GLU:HG3	1.87	0.56
1:M:154:LEU:HD21	1:N:164:TYR:CB	2.34	0.56
1:I:44:LEU:HB3	1:I:167:GLU:HG3	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:79:GLN:HA	1:M:83:LYS:HD2	1.88	0.56
1:I:154:LEU:HD21	1:W:164:TYR:CB	2.35	0.56
1:O:44:LEU:HB3	1:O:167:GLU:HG3	1.88	0.56
1:J:44:LEU:HB3	1:J:167:GLU:HG3	1.88	0.56
1:H:52:ARG:HH22	1:I:63:GLU:HG2	1.70	0.56
1:D:97:LYS:O	1:D:100:ILE:HG22	2.06	0.56
1:S:97:LYS:O	1:S:100:ILE:HG22	2.06	0.56
1:B:83:LYS:HD2	1:D:79:GLN:HA	1.88	0.56
1:T:97:LYS:O	1:T:100:ILE:HG22	2.06	0.56
1:C:97:LYS:O	1:C:100:ILE:HG22	2.06	0.56
1:D:154:LEU:HD21	1:M:164:TYR:CB	2.36	0.55
1:W:97:LYS:O	1:W:100:ILE:HG22	2.07	0.55
1:B:97:LYS:O	1:B:100:ILE:HG22	2.07	0.55
1:G:44:LEU:HB3	1:G:167:GLU:HG3	1.88	0.55
1:J:97:LYS:O	1:J:100:ILE:HG22	2.06	0.55
1:M:97:LYS:O	1:M:100:ILE:HG22	2.06	0.55
1:O:97:LYS:O	1:O:100:ILE:HG22	2.07	0.55
1:G:97:LYS:O	1:G:100:ILE:HG22	2.06	0.55
1:A:164:TYR:CB	1:B:154:LEU:HD21	2.35	0.55
1:N:97:LYS:O	1:N:100:ILE:HG22	2.07	0.55
1:E:44:LEU:HB3	1:E:167:GLU:HG3	1.88	0.55
1:I:97:LYS:O	1:I:100:ILE:HG22	2.07	0.55
1:Q:63:GLU:HG2	1:R:52:ARG:NH2	2.21	0.55
1:E:97:LYS:O	1:E:100:ILE:HG22	2.07	0.55
1:P:97:LYS:O	1:P:100:ILE:HG22	2.06	0.55
1:V:97:LYS:O	1:V:100:ILE:HG22	2.06	0.55
1:K:97:LYS:O	1:K:100:ILE:HG22	2.07	0.55
1:Q:44:LEU:HB3	1:Q:167:GLU:HG3	1.87	0.55
1:N:44:LEU:HB3	1:N:167:GLU:HG3	1.87	0.55
1:F:97:LYS:O	1:F:100:ILE:HG22	2.06	0.55
1:R:44:LEU:HB3	1:R:167:GLU:HG3	1.89	0.55
1:L:97:LYS:O	1:L:100:ILE:HG22	2.06	0.55
1:P:44:LEU:HB3	1:P:167:GLU:HG3	1.89	0.55
1:M:44:LEU:HB3	1:M:167:GLU:HG3	1.89	0.55
1:Q:97:LYS:O	1:Q:100:ILE:HG22	2.06	0.55
1:H:83:LYS:HD2	1:I:79:GLN:HA	1.88	0.55
1:K:79:GLN:HA	1:P:83:LYS:HD2	1.89	0.55
1:X:97:LYS:O	1:X:100:ILE:HG22	2.07	0.55
1:F:168:ARG:HA	1:F:168:ARG:HE	1.72	0.54
1:D:44:LEU:HB3	1:D:167:GLU:HG3	1.88	0.54
1:A:97:LYS:O	1:A:100:ILE:HG22	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:97:LYS:O	1:H:100:ILE:HG22	2.07	0.54
1:U:97:LYS:O	1:U:100:ILE:HG22	2.06	0.54
1:L:44:LEU:HB3	1:L:167:GLU:HG3	1.89	0.54
1:A:44:LEU:HB3	1:A:167:GLU:HG3	1.90	0.54
1:W:79:GLN:HA	1:X:83:LYS:HD2	1.90	0.54
1:U:44:LEU:HB3	1:U:167:GLU:HG3	1.89	0.54
1:R:97:LYS:O	1:R:100:ILE:HG22	2.07	0.54
1:N:83:LYS:HD3	1:S:79:GLN:HA	1.89	0.54
1:K:44:LEU:HB3	1:K:167:GLU:HG3	1.89	0.54
1:W:44:LEU:HB3	1:W:167:GLU:HG3	1.89	0.54
1:H:126:CYS:O	1:H:130:GLU:HG3	2.08	0.54
1:U:126:CYS:O	1:U:130:GLU:HG3	2.08	0.54
1:X:44:LEU:HB3	1:X:167:GLU:HG3	1.90	0.54
1:T:83:LYS:HD2	1:V:79:GLN:HA	1.90	0.54
1:H:44:LEU:HB3	1:H:167:GLU:HG3	1.90	0.54
1:L:83:LYS:HD2	1:M:79:GLN:HA	1.90	0.54
1:Q:126:CYS:O	1:Q:130:GLU:HG3	2.08	0.53
1:T:126:CYS:O	1:T:130:GLU:HG3	2.08	0.53
1:H:52:ARG:HH22	1:I:63:GLU:CD	2.12	0.53
1:F:126:CYS:O	1:F:130:GLU:HG3	2.09	0.53
1:E:79:GLN:HA	1:G:83:LYS:HD2	1.91	0.53
1:B:52:ARG:HH22	1:D:63:GLU:CG	2.21	0.53
1:C:44:LEU:HB3	1:C:167:GLU:HG3	1.89	0.53
1:T:44:LEU:HB3	1:T:167:GLU:HG3	1.89	0.53
1:B:126:CYS:O	1:B:130:GLU:HG3	2.08	0.53
1:C:126:CYS:O	1:C:130:GLU:HG3	2.09	0.53
1:E:83:LYS:HD2	1:G:79:GLN:HA	1.89	0.53
1:R:126:CYS:O	1:R:130:GLU:HG3	2.08	0.53
1:S:126:CYS:O	1:S:130:GLU:HG3	2.09	0.53
1:X:126:CYS:O	1:X:130:GLU:HG3	2.09	0.53
1:N:126:CYS:O	1:N:130:GLU:HG3	2.08	0.53
1:B:52:ARG:NH2	1:D:63:GLU:HG2	2.24	0.53
1:P:126:CYS:O	1:P:130:GLU:HG3	2.09	0.53
1:F:44:LEU:HB3	1:F:167:GLU:HG3	1.89	0.53
1:A:126:CYS:O	1:A:130:GLU:HG3	2.09	0.53
1:K:126:CYS:O	1:K:130:GLU:HG3	2.09	0.53
1:E:126:CYS:O	1:E:130:GLU:HG3	2.09	0.53
1:O:126:CYS:O	1:O:130:GLU:HG3	2.08	0.53
1:D:126:CYS:O	1:D:130:GLU:HG3	2.09	0.53
1:V:126:CYS:O	1:V:130:GLU:HG3	2.08	0.53
1:W:126:CYS:O	1:W:130:GLU:HG3	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:S:44:LEU:HB3	1:S:167:GLU:HG3	1.89	0.52
1:M:126:CYS:O	1:M:130:GLU:HG3	2.09	0.52
1:L:126:CYS:O	1:L:130:GLU:HG3	2.09	0.52
1:I:126:CYS:O	1:I:130:GLU:HG3	2.09	0.52
1:J:126:CYS:O	1:J:130:GLU:HG3	2.09	0.52
1:G:126:CYS:O	1:G:130:GLU:HG3	2.09	0.52
1:B:44:LEU:HB3	1:B:167:GLU:HG3	1.90	0.51
1:C:150:ASN:HA	1:C:153:ARG:HG2	1.93	0.51
1:N:63:GLU:HG2	1:S:52:ARG:NH2	2.25	0.51
1:T:150:ASN:HA	1:T:153:ARG:HG2	1.93	0.51
1:Q:63:GLU:CD	1:R:52:ARG:HH22	2.14	0.51
1:A:150:ASN:HA	1:A:153:ARG:HG2	1.93	0.51
1:N:63:GLU:CG	1:S:52:ARG:HH22	2.23	0.50
1:H:79:GLN:HA	1:I:83:LYS:HD3	1.92	0.50
1:Q:63:GLU:HG2	1:R:52:ARG:HH22	1.74	0.50
1:U:150:ASN:HA	1:U:153:ARG:HG2	1.94	0.50
1:W:83:LYS:HD2	1:X:79:GLN:HA	1.94	0.50
1:K:83:LYS:HD2	1:P:79:GLN:HA	1.92	0.50
1:H:150:ASN:HA	1:H:153:ARG:HG2	1.94	0.50
1:B:150:ASN:HA	1:B:153:ARG:HG2	1.94	0.50
1:Q:150:ASN:HA	1:Q:153:ARG:HG2	1.94	0.50
1:C:52:ARG:HH22	1:J:63:GLU:CG	2.24	0.50
1:G:150:ASN:HA	1:G:153:ARG:HG2	1.93	0.50
1:F:150:ASN:HA	1:F:153:ARG:HG2	1.94	0.50
1:O:63:GLU:CG	1:U:52:ARG:HH22	2.25	0.49
1:R:150:ASN:HA	1:R:153:ARG:HG2	1.94	0.49
1:S:165:LEU:O	1:S:169:LEU:HB2	2.12	0.49
1:U:165:LEU:O	1:U:169:LEU:HB2	2.13	0.49
1:R:165:LEU:O	1:R:169:LEU:HB2	2.13	0.49
1:A:52:ARG:NH2	1:F:63:GLU:HG2	2.27	0.49
1:S:150:ASN:HA	1:S:153:ARG:HG2	1.94	0.49
1:H:165:LEU:O	1:H:169:LEU:HB2	2.13	0.49
1:C:52:ARG:NH2	1:J:63:GLU:HG2	2.27	0.49
1:O:63:GLU:HG2	1:U:52:ARG:NH2	2.27	0.49
1:T:165:LEU:O	1:T:169:LEU:HB2	2.13	0.49
1:P:150:ASN:HA	1:P:153:ARG:HG2	1.94	0.48
1:A:165:LEU:O	1:A:169:LEU:HB2	2.13	0.48
1:V:150:ASN:HA	1:V:153:ARG:HG2	1.93	0.48
1:A:52:ARG:HH22	1:F:63:GLU:CG	2.26	0.48
1:I:150:ASN:HA	1:I:153:ARG:HG2	1.95	0.48
1:D:150:ASN:HA	1:D:153:ARG:HG2	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:150:ASN:HA	1:E:153:ARG:HG2	1.95	0.48
1:N:150:ASN:HA	1:N:153:ARG:HG2	1.95	0.48
1:K:165:LEU:O	1:K:169:LEU:HB2	2.14	0.48
1:L:150:ASN:HA	1:L:153:ARG:HG2	1.95	0.48
1:H:58:LYS:HA	1:H:58:LYS:HD3	1.66	0.47
1:M:150:ASN:HA	1:M:153:ARG:HG2	1.95	0.47
1:G:165:LEU:O	1:G:169:LEU:HB2	2.14	0.47
1:B:165:LEU:O	1:B:169:LEU:HB2	2.14	0.47
1:O:150:ASN:HA	1:O:153:ARG:HG2	1.96	0.47
1:C:165:LEU:O	1:C:169:LEU:HB2	2.14	0.47
1:J:150:ASN:HA	1:J:153:ARG:HG2	1.94	0.47
1:W:150:ASN:HA	1:W:153:ARG:HG2	1.95	0.47
1:U:58:LYS:HA	1:U:58:LYS:HD3	1.66	0.47
1:W:165:LEU:O	1:W:169:LEU:HB2	2.14	0.47
1:J:165:LEU:O	1:J:169:LEU:HB2	2.15	0.47
1:F:75:ARG:HD3	1:F:75:ARG:HA	1.70	0.47
1:A:145:GLY:HA3	1:I:4:ILE:HG21	1.96	0.47
1:F:165:LEU:O	1:F:169:LEU:HB2	2.14	0.47
1:X:150:ASN:HA	1:X:153:ARG:HG2	1.96	0.47
1:K:150:ASN:HA	1:K:153:ARG:HG2	1.96	0.47
1:Q:75:ARG:HD3	1:Q:75:ARG:HA	1.72	0.47
1:O:165:LEU:O	1:O:169:LEU:HB2	2.14	0.47
1:A:32:SER:CB	1:A:84:PRO:HG2	2.45	0.47
1:D:75:ARG:HA	1:D:75:ARG:HD3	1.72	0.46
1:X:16:VAL:HG22	1:X:113:LEU:HD21	1.97	0.46
1:N:75:ARG:HA	1:N:75:ARG:HD3	1.72	0.46
1:M:165:LEU:O	1:M:169:LEU:HB2	2.16	0.46
1:K:58:LYS:HA	1:K:58:LYS:HD3	1.66	0.46
1:X:165:LEU:O	1:X:169:LEU:HB2	2.16	0.46
1:I:58:LYS:HD3	1:I:58:LYS:HA	1.65	0.46
1:B:52:ARG:HH22	1:D:63:GLU:CD	2.19	0.46
1:P:165:LEU:O	1:P:169:LEU:HB2	2.16	0.46
1:R:58:LYS:HA	1:R:58:LYS:HD3	1.66	0.46
1:E:4:ILE:HG21	1:F:145:GLY:HA3	1.98	0.46
1:B:16:VAL:HG22	1:B:113:LEU:HD21	1.98	0.46
1:Q:165:LEU:HD23	1:Q:165:LEU:HA	1.80	0.46
1:I:145:GLY:HA3	1:X:4:ILE:HG21	1.98	0.46
1:E:58:LYS:HA	1:E:58:LYS:HD3	1.66	0.46
1:M:4:ILE:HG21	1:Q:145:GLY:HA3	1.98	0.45
1:O:58:LYS:HD3	1:O:58:LYS:HA	1.66	0.45
1:W:58:LYS:HA	1:W:58:LYS:HD3	1.66	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:W:75:ARG:HD3	1:W:75:ARG:HA	1.73	0.45
1:Q:165:LEU:O	1:Q:169:LEU:HB2	2.16	0.45
1:E:165:LEU:O	1:E:169:LEU:HB2	2.17	0.45
1:R:75:ARG:HA	1:R:75:ARG:HD3	1.72	0.45
1:J:165:LEU:HD23	1:J:165:LEU:HA	1.77	0.45
1:S:16:VAL:HG22	1:S:113:LEU:HD21	1.98	0.45
1:V:145:GLY:HA3	1:W:4:ILE:HG21	1.98	0.45
1:D:145:GLY:HA3	1:L:4:ILE:HG21	1.98	0.45
1:L:16:VAL:HG22	1:L:113:LEU:HD21	1.98	0.45
1:T:16:VAL:HG22	1:T:113:LEU:HD21	1.99	0.45
1:D:58:LYS:HA	1:D:58:LYS:HD3	1.66	0.45
1:M:157:SER:HB2	1:M:158:GLN:HE21	1.82	0.45
1:O:32:SER:CB	1:O:84:PRO:HG2	2.47	0.45
1:A:75:ARG:HD3	1:A:75:ARG:HA	1.70	0.45
1:P:58:LYS:HA	1:P:58:LYS:HD3	1.66	0.45
1:E:16:VAL:HG22	1:E:113:LEU:HD21	1.99	0.45
1:J:32:SER:CB	1:J:84:PRO:HG2	2.47	0.45
1:N:165:LEU:O	1:N:169:LEU:HB2	2.16	0.45
1:X:58:LYS:HD3	1:X:58:LYS:HA	1.66	0.45
1:K:75:ARG:HD3	1:K:75:ARG:HA	1.73	0.45
1:M:58:LYS:HA	1:M:58:LYS:HD3	1.66	0.45
1:M:102:LEU:HD12	1:M:102:LEU:HA	1.81	0.45
1:G:102:LEU:HA	1:G:102:LEU:HD12	1.81	0.45
1:H:16:VAL:HG22	1:H:113:LEU:HD21	1.98	0.45
1:U:16:VAL:HG22	1:U:113:LEU:HD21	1.99	0.45
1:O:145:GLY:HA3	1:P:4:ILE:HG21	1.99	0.45
1:V:165:LEU:O	1:V:169:LEU:HB2	2.17	0.45
1:C:164:TYR:CE1	1:C:168:ARG:HG3	2.52	0.44
1:H:165:LEU:HD23	1:H:165:LEU:HA	1.77	0.44
1:J:145:GLY:HA3	1:K:4:ILE:HG21	1.99	0.44
1:Q:157:SER:HB2	1:Q:158:GLN:HE21	1.83	0.44
1:V:32:SER:CB	1:V:84:PRO:HG2	2.47	0.44
1:H:145:GLY:HA3	1:J:4:ILE:HG21	1.99	0.44
1:V:157:SER:HB2	1:V:158:GLN:HE21	1.82	0.44
1:G:4:ILE:HG21	1:N:145:GLY:HA3	1.98	0.44
1:O:75:ARG:HD3	1:O:75:ARG:HA	1.72	0.44
1:R:16:VAL:HG22	1:R:113:LEU:HD21	1.99	0.44
1:U:164:TYR:CE1	1:U:168:ARG:HG3	2.52	0.44
1:T:164:TYR:CE1	1:T:168:ARG:HG3	2.52	0.44
1:C:16:VAL:HG22	1:C:113:LEU:HD21	1.99	0.44
1:M:16:VAL:HG22	1:M:113:LEU:HD21	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:58:LYS:HA	1:L:58:LYS:HD3	1.66	0.44
1:L:102:LEU:HD12	1:L:102:LEU:HA	1.81	0.44
1:I:157:SER:HB2	1:I:158:GLN:HE21	1.82	0.44
1:P:16:VAL:HG22	1:P:113:LEU:HD21	1.99	0.44
1:T:52:ARG:NH2	1:V:63:GLU:HG2	2.32	0.44
1:I:75:ARG:HA	1:I:75:ARG:HD3	1.73	0.44
1:N:58:LYS:HA	1:N:58:LYS:HD3	1.66	0.44
1:R:170:THR:HG22	1:R:171:LEU:HD23	1.99	0.44
1:R:164:TYR:CE1	1:R:168:ARG:HG3	2.52	0.44
1:U:145:GLY:HA3	1:V:4:ILE:HG21	2.00	0.44
1:H:4:ILE:HG21	1:K:145:GLY:HA3	1.99	0.44
1:W:157:SER:HB2	1:W:158:GLN:HE21	1.82	0.44
1:D:165:LEU:O	1:D:169:LEU:HB2	2.17	0.44
1:E:102:LEU:HA	1:E:102:LEU:HD12	1.81	0.44
1:F:16:VAL:HG22	1:F:113:LEU:HD21	2.00	0.44
1:M:165:LEU:HA	1:M:165:LEU:HD23	1.75	0.44
1:N:4:ILE:HG21	1:T:145:GLY:HA3	1.99	0.44
1:D:32:SER:CB	1:D:84:PRO:HG2	2.47	0.44
1:O:157:SER:HB2	1:O:158:GLN:HE21	1.83	0.44
1:I:32:SER:CB	1:I:84:PRO:HG2	2.48	0.44
1:B:102:LEU:HD12	1:B:102:LEU:HA	1.81	0.44
1:T:75:ARG:HA	1:T:75:ARG:HD3	1.73	0.44
1:R:165:LEU:HD23	1:R:165:LEU:HA	1.79	0.44
1:E:165:LEU:HA	1:E:165:LEU:HD23	1.72	0.44
1:C:145:GLY:HA3	1:D:4:ILE:HG21	1.99	0.44
1:G:16:VAL:HG22	1:G:113:LEU:HD21	1.99	0.44
1:T:102:LEU:HD12	1:T:102:LEU:HA	1.82	0.44
1:C:4:ILE:HG21	1:L:145:GLY:HA3	1.99	0.44
1:V:157:SER:HB2	1:V:158:GLN:NE2	2.33	0.44
1:U:4:ILE:HG21	1:W:145:GLY:HA3	1.99	0.44
1:C:102:LEU:HD12	1:C:102:LEU:HA	1.81	0.44
1:Q:32:SER:CB	1:Q:84:PRO:HG2	2.48	0.44
1:J:157:SER:HB2	1:J:158:GLN:HE21	1.83	0.44
1:L:165:LEU:O	1:L:169:LEU:HB2	2.18	0.44
1:C:52:ARG:HH22	1:J:63:GLU:CD	2.21	0.43
1:N:157:SER:HB2	1:N:158:GLN:NE2	2.34	0.43
1:C:75:ARG:HD3	1:C:75:ARG:HA	1.73	0.43
1:T:58:LYS:HA	1:T:58:LYS:HD3	1.66	0.43
1:G:145:GLY:HA3	1:T:4:ILE:HG21	1.99	0.43
1:G:165:LEU:HA	1:G:165:LEU:HD23	1.77	0.43
1:Q:16:VAL:HG22	1:Q:113:LEU:HD21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:16:VAL:HG22	1:I:113:LEU:HD21	2.00	0.43
1:D:157:SER:HB2	1:D:158:GLN:HE21	1.83	0.43
1:D:157:SER:HB2	1:D:158:GLN:NE2	2.34	0.43
1:B:75:ARG:HD3	1:B:75:ARG:HA	1.72	0.43
1:M:157:SER:HB2	1:M:158:GLN:NE2	2.33	0.43
1:Q:157:SER:HB2	1:Q:158:GLN:NE2	2.33	0.43
1:J:157:SER:HB2	1:J:158:GLN:NE2	2.33	0.43
1:M:145:GLY:HA3	1:S:4:ILE:HG21	1.99	0.43
1:N:63:GLU:CD	1:S:52:ARG:HH22	2.22	0.43
1:W:157:SER:HB2	1:W:158:GLN:NE2	2.33	0.43
1:N:157:SER:HB2	1:N:158:GLN:HE21	1.84	0.43
1:U:102:LEU:HA	1:U:102:LEU:HD12	1.82	0.43
1:Q:58:LYS:HA	1:Q:58:LYS:HD3	1.66	0.43
1:W:165:LEU:HD23	1:W:165:LEU:HA	1.75	0.43
1:I:157:SER:HB2	1:I:158:GLN:NE2	2.33	0.43
1:B:4:ILE:HG21	1:E:145:GLY:HA3	1.99	0.43
1:L:75:ARG:HA	1:L:75:ARG:HD3	1.74	0.43
1:B:164:TYR:CE1	1:B:168:ARG:HG3	2.53	0.43
1:U:165:LEU:HD23	1:U:165:LEU:HA	1.79	0.43
1:B:165:LEU:HA	1:B:165:LEU:HD23	1.78	0.43
1:A:32:SER:HB3	1:A:84:PRO:HG2	2.01	0.43
1:O:157:SER:HB2	1:O:158:GLN:NE2	2.33	0.43
1:I:165:LEU:O	1:I:169:LEU:HB2	2.18	0.43
1:A:4:ILE:HG21	1:X:145:GLY:HA3	2.00	0.43
1:A:29:THR:HA	1:A:84:PRO:HG3	2.01	0.43
1:A:102:LEU:HA	1:A:102:LEU:HD12	1.81	0.43
1:N:32:SER:CB	1:N:84:PRO:HG2	2.48	0.43
1:O:16:VAL:HG22	1:O:113:LEU:HD21	2.01	0.43
1:G:157:SER:HB2	1:G:158:GLN:NE2	2.33	0.43
1:P:32:SER:CB	1:P:84:PRO:HG2	2.49	0.43
1:X:165:LEU:HA	1:X:165:LEU:HD23	1.75	0.42
1:P:145:GLY:HA3	1:R:4:ILE:HG21	1.99	0.42
1:N:16:VAL:HG22	1:N:113:LEU:HD21	2.00	0.42
1:E:157:SER:HB2	1:E:158:GLN:NE2	2.34	0.42
1:D:16:VAL:HG22	1:D:113:LEU:HD21	2.01	0.42
1:C:58:LYS:HA	1:C:58:LYS:HD3	1.67	0.42
1:G:75:ARG:HD3	1:G:75:ARG:HA	1.74	0.42
1:A:164:TYR:CE1	1:A:168:ARG:HG3	2.53	0.42
1:F:32:SER:CB	1:F:84:PRO:HG2	2.48	0.42
1:T:170:THR:HG22	1:T:171:LEU:HD23	2.00	0.42
1:H:102:LEU:HA	1:H:102:LEU:HD12	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:P:165:LEU:HD23	1:P:165:LEU:HA	1.76	0.42
1:G:157:SER:HB2	1:G:158:GLN:HE21	1.84	0.42
1:C:170:THR:HG22	1:C:171:LEU:HD23	2.00	0.42
1:S:102:LEU:HD12	1:S:102:LEU:HA	1.82	0.42
1:N:102:LEU:HA	1:N:102:LEU:HD12	1.81	0.42
1:E:32:SER:CB	1:E:84:PRO:HG2	2.49	0.42
1:A:52:ARG:HH22	1:F:63:GLU:CD	2.23	0.42
1:V:165:LEU:HD23	1:V:165:LEU:HA	1.80	0.42
1:B:145:GLY:HA3	1:F:4:ILE:HG21	2.00	0.42
1:O:4:ILE:HG21	1:R:145:GLY:HA3	2.00	0.42
1:M:32:SER:CB	1:M:84:PRO:HG2	2.50	0.42
1:G:32:SER:CB	1:G:84:PRO:HG2	2.49	0.42
1:A:52:ARG:NH2	1:F:63:GLU:CD	2.73	0.42
1:J:29:THR:HA	1:J:84:PRO:HG3	2.02	0.42
1:V:29:THR:HA	1:V:84:PRO:HG3	2.02	0.42
1:S:32:SER:CB	1:S:84:PRO:HG2	2.50	0.42
1:F:58:LYS:HD3	1:F:58:LYS:HA	1.66	0.42
1:W:16:VAL:HG22	1:W:113:LEU:HD21	2.01	0.42
1:O:63:GLU:CD	1:U:52:ARG:HH22	2.22	0.42
1:O:29:THR:HA	1:O:84:PRO:HG3	2.02	0.42
1:B:157:SER:HB2	1:B:158:GLN:NE2	2.34	0.42
1:P:157:SER:HB2	1:P:158:GLN:NE2	2.34	0.42
1:W:102:LEU:HA	1:W:102:LEU:HD12	1.80	0.42
1:S:164:TYR:CE1	1:S:168:ARG:HG3	2.54	0.42
1:A:37:PHE:CD1	1:A:42:VAL:CG1	3.03	0.42
1:K:16:VAL:HG22	1:K:113:LEU:HD21	2.02	0.42
1:Q:83:LYS:HE3	1:R:79:GLN:HA	2.01	0.42
1:V:16:VAL:HG22	1:V:113:LEU:HD21	2.01	0.42
1:B:58:LYS:HA	1:B:58:LYS:HD3	1.66	0.42
1:Q:4:ILE:HG21	1:S:145:GLY:HA3	2.01	0.42
1:Q:29:THR:HA	1:Q:84:PRO:HG3	2.02	0.41
1:J:16:VAL:HG22	1:J:113:LEU:HD21	2.01	0.41
1:K:102:LEU:HA	1:K:102:LEU:HD12	1.81	0.41
1:L:157:SER:HB2	1:L:158:GLN:NE2	2.34	0.41
1:F:157:SER:HB2	1:F:158:GLN:NE2	2.35	0.41
1:X:32:SER:CB	1:X:84:PRO:HG2	2.51	0.41
1:H:32:SER:CB	1:H:84:PRO:HG2	2.50	0.41
1:B:32:SER:CB	1:B:84:PRO:HG2	2.50	0.41
1:U:32:SER:CB	1:U:84:PRO:HG2	2.49	0.41
1:C:32:SER:CB	1:C:84:PRO:HG2	2.50	0.41
1:F:164:TYR:CE1	1:F:168:ARG:HG3	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:52:ARG:NH2	1:D:63:GLU:CG	2.83	0.41
1:R:32:SER:CB	1:R:84:PRO:HG2	2.50	0.41
1:G:164:TYR:CE1	1:G:168:ARG:HG3	2.56	0.41
1:T:32:SER:CB	1:T:84:PRO:HG2	2.50	0.41
1:K:157:SER:HB2	1:K:158:GLN:NE2	2.35	0.41
1:N:165:LEU:HD23	1:N:165:LEU:HA	1.78	0.41
1:E:157:SER:HB2	1:E:158:GLN:HE21	1.86	0.41
1:V:129:LEU:O	1:V:133:PHE:HB2	2.21	0.41
1:S:58:LYS:HA	1:S:58:LYS:HD3	1.66	0.41
1:S:75:ARG:HA	1:S:75:ARG:HD3	1.72	0.41
1:K:165:LEU:HD23	1:K:165:LEU:HA	1.73	0.41
1:P:157:SER:HB2	1:P:158:GLN:HE21	1.86	0.41
1:V:102:LEU:HD12	1:V:102:LEU:HA	1.80	0.41
1:D:102:LEU:HD12	1:D:102:LEU:HA	1.82	0.41
1:G:58:LYS:HA	1:G:58:LYS:HD3	1.66	0.41
1:O:164:TYR:CE1	1:O:168:ARG:HG3	2.56	0.41
1:V:37:PHE:CD1	1:V:42:VAL:CG1	3.04	0.41
1:J:37:PHE:CD1	1:J:42:VAL:CG1	3.04	0.41
1:T:165:LEU:HD23	1:T:165:LEU:HA	1.81	0.41
1:C:165:LEU:HD23	1:C:165:LEU:HA	1.81	0.41
1:L:165:LEU:HD23	1:L:165:LEU:HA	1.75	0.41
1:F:157:SER:HB2	1:F:158:GLN:HE21	1.86	0.41
1:L:32:SER:CB	1:L:84:PRO:HG2	2.50	0.41
1:J:102:LEU:HA	1:J:102:LEU:HD12	1.81	0.41
1:Q:164:TYR:CE1	1:Q:168:ARG:HG3	2.55	0.41
1:D:164:TYR:CE1	1:D:168:ARG:HG3	2.54	0.41
1:N:164:TYR:CE1	1:N:168:ARG:HG3	2.54	0.41
1:O:37:PHE:CD1	1:O:42:VAL:CG1	3.04	0.41
1:N:63:GLU:CG	1:S:52:ARG:NH2	2.84	0.41
1:J:129:LEU:O	1:J:133:PHE:HB2	2.21	0.41
1:A:63:GLU:HG2	1:F:52:ARG:NH2	2.36	0.41
1:J:164:TYR:CE1	1:J:168:ARG:HG3	2.56	0.41
1:W:164:TYR:CE1	1:W:168:ARG:HG3	2.55	0.41
1:A:165:LEU:HD23	1:A:165:LEU:HA	1.83	0.41
1:X:157:SER:HB2	1:X:158:GLN:NE2	2.35	0.41
1:L:129:LEU:O	1:L:133:PHE:HB2	2.21	0.41
1:E:75:ARG:HD3	1:E:75:ARG:HA	1.74	0.41
1:O:129:LEU:O	1:O:133:PHE:HB2	2.21	0.41
1:G:129:LEU:O	1:G:133:PHE:HB2	2.21	0.41
1:W:170:THR:HG22	1:W:171:LEU:HD23	2.03	0.41
1:S:37:PHE:CD1	1:S:42:VAL:CG1	3.04	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:37:PHE:CD1	1:B:42:VAL:CG1	3.04	0.41
1:R:102:LEU:HA	1:R:102:LEU:HD12	1.82	0.41
1:N:129:LEU:O	1:N:133:PHE:HB2	2.21	0.41
1:M:37:PHE:CD1	1:M:42:VAL:CG1	3.04	0.40
1:S:165:LEU:HD23	1:S:165:LEU:HA	1.78	0.40
1:B:157:SER:HB2	1:B:158:GLN:HE21	1.85	0.40
1:B:92:THR:O	1:B:92:THR:HG22	2.22	0.40
1:Q:102:LEU:HA	1:Q:102:LEU:HD12	1.81	0.40
1:C:92:THR:HG22	1:C:92:THR:O	2.21	0.40
1:P:164:TYR:CE1	1:P:168:ARG:HG3	2.56	0.40
1:D:29:THR:HA	1:D:84:PRO:HG3	2.02	0.40
1:F:29:THR:HA	1:F:84:PRO:HG3	2.03	0.40
1:Q:129:LEU:O	1:Q:133:PHE:HB2	2.21	0.40
1:T:37:PHE:CD1	1:T:42:VAL:CG1	3.05	0.40
1:A:52:ARG:NH2	1:F:63:GLU:CG	2.83	0.40
1:I:29:THR:HA	1:I:84:PRO:HG3	2.04	0.40
1:I:129:LEU:O	1:I:133:PHE:HB2	2.22	0.40
1:K:32:SER:CB	1:K:84:PRO:HG2	2.51	0.40
1:T:92:THR:HG22	1:T:92:THR:O	2.21	0.40
1:W:37:PHE:CD1	1:W:42:VAL:CG1	3.04	0.40
1:D:129:LEU:O	1:D:133:PHE:HB2	2.21	0.40
1:L:63:GLU:HG2	1:M:52:ARG:NH2	2.36	0.40
1:X:129:LEU:O	1:X:133:PHE:HB2	2.21	0.40
1:R:129:LEU:O	1:R:133:PHE:HB2	2.22	0.40
1:K:24:LEU:HA	1:K:24:LEU:HD23	1.95	0.40
1:U:75:ARG:HD3	1:U:75:ARG:HA	1.73	0.40
1:X:92:THR:O	1:X:92:THR:HG22	2.22	0.40
1:M:75:ARG:HD3	1:M:75:ARG:HA	1.74	0.40
1:P:24:LEU:HD23	1:P:24:LEU:HA	1.97	0.40
1:E:37:PHE:CD1	1:E:42:VAL:CG1	3.04	0.40
1:U:170:THR:HG22	1:U:171:LEU:HD23	2.02	0.40
1:F:129:LEU:O	1:F:133:PHE:HB2	2.21	0.40
1:H:75:ARG:HA	1:H:75:ARG:HD3	1.73	0.40
1:H:129:LEU:O	1:H:133:PHE:HB2	2.21	0.40
1:U:129:LEU:O	1:U:133:PHE:HB2	2.21	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 PDB entries followed by that with respect to all EM entries.

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	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	B	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	C	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	D	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	E	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	F	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	G	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	H	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	I	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	J	168/174 (97%)	163 (97%)	4 (2%)	1 (1%)	30	74
1	K	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	L	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	M	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	N	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	O	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	P	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	Q	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	R	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	S	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	T	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	U	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	V	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	W	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
1	X	168/174 (97%)	162 (96%)	5 (3%)	1 (1%)	30	74
All	All	4032/4176 (97%)	3889 (96%)	119 (3%)	24 (1%)	34	74

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	156	GLY
1	F	156	GLY
1	G	156	GLY
1	J	156	GLY
1	L	156	GLY
1	O	156	GLY
1	B	156	GLY
1	D	156	GLY
1	E	156	GLY
1	H	156	GLY
1	I	156	GLY
1	K	156	GLY
1	N	156	GLY
1	P	156	GLY
1	Q	156	GLY
1	S	156	GLY
1	V	156	GLY
1	X	156	GLY
1	C	156	GLY
1	M	156	GLY
1	R	156	GLY
1	T	156	GLY
1	U	156	GLY
1	W	156	GLY

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

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	144/148 (97%)	134 (93%)	10 (7%)	19 58
1	B	144/148 (97%)	135 (94%)	9 (6%)	22 61
1	C	144/148 (97%)	134 (93%)	10 (7%)	19 58
1	D	144/148 (97%)	136 (94%)	8 (6%)	26 65
1	E	144/148 (97%)	135 (94%)	9 (6%)	22 61

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	F	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	G	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	H	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	I	144/148 (97%)	136 (94%)	8 (6%)	26	65
1	J	144/148 (97%)	138 (96%)	6 (4%)	36	71
1	K	144/148 (97%)	136 (94%)	8 (6%)	26	65
1	L	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	M	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	N	144/148 (97%)	136 (94%)	8 (6%)	26	65
1	O	144/148 (97%)	137 (95%)	7 (5%)	31	68
1	P	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	Q	144/148 (97%)	137 (95%)	7 (5%)	31	68
1	R	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	S	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	T	144/148 (97%)	134 (93%)	10 (7%)	19	58
1	U	144/148 (97%)	135 (94%)	9 (6%)	22	61
1	V	144/148 (97%)	137 (95%)	7 (5%)	31	68
1	W	144/148 (97%)	136 (94%)	8 (6%)	26	65
1	X	144/148 (97%)	135 (94%)	9 (6%)	22	61
All	All	3456/3552 (97%)	3251 (94%)	205 (6%)	29	63

All (205) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	ILE
1	A	18	ARG
1	A	82	GLN
1	A	87	ASP
1	A	88	GLU
1	A	148	LEU
1	A	157	SER
1	A	158	GLN
1	A	168	ARG
1	A	171	LEU
1	B	18	ARG

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Mol	Chain	Res	Type
1	B	82	GLN
1	B	87	ASP
1	B	88	GLU
1	B	143	LYS
1	B	157	SER
1	B	158	GLN
1	B	168	ARG
1	B	171	LEU
1	C	18	ARG
1	C	82	GLN
1	C	87	ASP
1	C	88	GLU
1	C	143	LYS
1	C	148	LEU
1	C	157	SER
1	C	158	GLN
1	C	168	ARG
1	C	171	LEU
1	D	82	GLN
1	D	87	ASP
1	D	88	GLU
1	D	143	LYS
1	D	157	SER
1	D	158	GLN
1	D	168	ARG
1	D	171	LEU
1	E	18	ARG
1	E	82	GLN
1	E	87	ASP
1	E	88	GLU
1	E	143	LYS
1	E	157	SER
1	E	158	GLN
1	E	168	ARG
1	E	171	LEU
1	F	82	GLN
1	F	83	LYS
1	F	87	ASP
1	F	88	GLU
1	F	143	LYS
1	F	157	SER
1	F	158	GLN

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Mol	Chain	Res	Type
1	F	168	ARG
1	F	171	LEU
1	G	18	ARG
1	G	82	GLN
1	G	87	ASP
1	G	88	GLU
1	G	143	LYS
1	G	157	SER
1	G	158	GLN
1	G	168	ARG
1	G	171	LEU
1	H	18	ARG
1	H	82	GLN
1	H	87	ASP
1	H	88	GLU
1	H	143	LYS
1	H	157	SER
1	H	158	GLN
1	H	168	ARG
1	H	171	LEU
1	I	82	GLN
1	I	87	ASP
1	I	88	GLU
1	I	143	LYS
1	I	157	SER
1	I	158	GLN
1	I	168	ARG
1	I	171	LEU
1	J	82	GLN
1	J	87	ASP
1	J	88	GLU
1	J	157	SER
1	J	158	GLN
1	J	171	LEU
1	K	4	ILE
1	K	82	GLN
1	K	87	ASP
1	K	88	GLU
1	K	157	SER
1	K	158	GLN
1	K	168	ARG
1	K	171	LEU

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Mol	Chain	Res	Type
1	L	18	ARG
1	L	82	GLN
1	L	87	ASP
1	L	88	GLU
1	L	143	LYS
1	L	157	SER
1	L	158	GLN
1	L	168	ARG
1	L	171	LEU
1	M	18	ARG
1	M	82	GLN
1	M	87	ASP
1	M	88	GLU
1	M	143	LYS
1	M	157	SER
1	M	158	GLN
1	M	168	ARG
1	M	171	LEU
1	N	82	GLN
1	N	87	ASP
1	N	88	GLU
1	N	143	LYS
1	N	157	SER
1	N	158	GLN
1	N	168	ARG
1	N	171	LEU
1	O	82	GLN
1	O	87	ASP
1	O	88	GLU
1	O	157	SER
1	O	158	GLN
1	O	168	ARG
1	O	171	LEU
1	P	18	ARG
1	P	82	GLN
1	P	87	ASP
1	P	88	GLU
1	P	143	LYS
1	P	157	SER
1	P	158	GLN
1	P	168	ARG
1	P	171	LEU

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Mol	Chain	Res	Type
1	Q	82	GLN
1	Q	87	ASP
1	Q	88	GLU
1	Q	157	SER
1	Q	158	GLN
1	Q	168	ARG
1	Q	171	LEU
1	R	18	ARG
1	R	82	GLN
1	R	87	ASP
1	R	88	GLU
1	R	143	LYS
1	R	157	SER
1	R	158	GLN
1	R	168	ARG
1	R	171	LEU
1	S	18	ARG
1	S	82	GLN
1	S	87	ASP
1	S	88	GLU
1	S	143	LYS
1	S	157	SER
1	S	158	GLN
1	S	168	ARG
1	S	171	LEU
1	T	18	ARG
1	T	82	GLN
1	T	87	ASP
1	T	88	GLU
1	T	143	LYS
1	T	148	LEU
1	T	157	SER
1	T	158	GLN
1	T	168	ARG
1	T	171	LEU
1	U	18	ARG
1	U	82	GLN
1	U	87	ASP
1	U	88	GLU
1	U	143	LYS
1	U	157	SER
1	U	158	GLN

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Mol	Chain	Res	Type
1	U	168	ARG
1	U	171	LEU
1	V	82	GLN
1	V	87	ASP
1	V	88	GLU
1	V	157	SER
1	V	158	GLN
1	V	168	ARG
1	V	171	LEU
1	W	82	GLN
1	W	87	ASP
1	W	88	GLU
1	W	143	LYS
1	W	157	SER
1	W	158	GLN
1	W	168	ARG
1	W	171	LEU
1	X	18	ARG
1	X	82	GLN
1	X	87	ASP
1	X	88	GLU
1	X	143	LYS
1	X	157	SER
1	X	158	GLN
1	X	168	ARG
1	X	171	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	GLN
1	B	71	GLN
1	C	71	GLN
1	D	71	GLN
1	E	71	GLN
1	F	71	GLN
1	G	71	GLN
1	H	71	GLN
1	I	71	GLN
1	J	71	GLN
1	K	71	GLN
1	L	71	GLN

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Mol	Chain	Res	Type
1	M	71	GLN
1	N	71	GLN
1	O	71	GLN
1	P	71	GLN
1	Q	71	GLN
1	R	71	GLN
1	S	71	GLN
1	T	71	GLN
1	U	71	GLN
1	V	71	GLN
1	W	71	GLN
1	X	71	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no carbohydrates in this entry.

5.6 Ligand geometry (i)

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.