



# wwPDB X-ray Structure Validation Summary Report ⓘ

Feb 15, 2017 – 06:09 am GMT

PDB ID : 1MFR  
Title : CRYSTAL STRUCTURE OF M FERRITIN  
Authors : Ha, Y.; Shi, D.; Allewell, N.M.  
Deposited on : 1998-06-18  
Resolution : 2.80 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Xtriage (Phenix) : NOT EXECUTED  
EDS : NOT EXECUTED  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

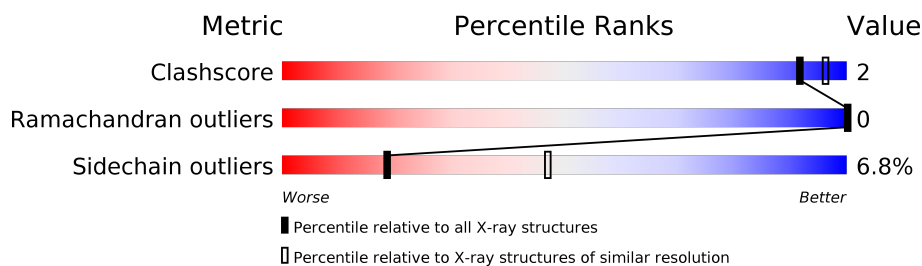
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 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))
Clashscore	112137	3033 (2.80-2.80)
Ramachandran outliers	110173	2983 (2.80-2.80)
Sidechain outliers	110143	2985 (2.80-2.80)


















The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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  $\leq 5\%$ . 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	176	
1	B	176	
1	C	176	
1	D	176	
1	E	176	
1	F	176	
1	G	176	

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Mol	Chain	Length	Quality of chain
1	H	176	 87% 7% . .
1	I	176	 86% 9% . .
1	J	176	 86% 8% . .
1	K	176	 88% 7% . .
1	L	176	 86% 7% . .
1	M	176	 88% 7% . .
1	N	176	 87% 7% . .
1	O	176	 87% 7% . .
1	P	176	 87% 7% . .
1	Q	176	 86% 8% . .
1	R	176	 88% 7% . .
1	S	176	 87% 7% . .
1	T	176	 88% 7% . .
1	U	176	 88% 7% . .
1	V	176	 88% 7% . .
1	W	176	 87% 7% . .
1	X	176	 86% 8% . .

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 34320 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 M FERRITIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	B	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	C	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	D	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	E	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	F	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	G	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	H	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	I	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	J	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	K	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	L	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	M	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	N	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	O	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	P	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	R	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	S	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	T	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	U	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	V	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	W	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			
1	X	171	Total	C	N	O	S	0	0	0
			1373	868	238	260	7			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	P	2	Total	Mg	0	0
			2	2		
2	K	2	Total	Mg	0	0
			2	2		
2	B	2	Total	Mg	0	0
			2	2		
2	W	2	Total	Mg	0	0
			2	2		
2	N	2	Total	Mg	0	0
			2	2		
2	X	2	Total	Mg	0	0
			2	2		
2	S	2	Total	Mg	0	0
			2	2		
2	J	2	Total	Mg	0	0
			2	2		
2	E	2	Total	Mg	0	0
			2	2		
2	V	2	Total	Mg	0	0
			2	2		
2	A	2	Total	Mg	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	R	2	Total 2	Mg 2	0	0
2	M	2	Total 2	Mg 2	0	0
2	D	2	Total 2	Mg 2	0	0
2	I	2	Total 2	Mg 2	0	0
2	U	2	Total 2	Mg 2	0	0
2	L	2	Total 2	Mg 2	0	0
2	G	2	Total 2	Mg 2	0	0
2	Q	2	Total 2	Mg 2	0	0
2	H	2	Total 2	Mg 2	0	0
2	C	2	Total 2	Mg 2	0	0
2	T	2	Total 2	Mg 2	0	0
2	O	2	Total 2	Mg 2	0	0
2	F	2	Total 2	Mg 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	54	Total 54	O 54	0	0
3	B	56	Total 56	O 56	0	0
3	C	57	Total 57	O 57	0	0
3	D	55	Total 55	O 55	0	0
3	E	55	Total 55	O 55	0	0
3	F	54	Total 54	O 54	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	G	55	Total 55	O 55	0	0
3	H	57	Total 57	O 57	0	0
3	I	55	Total 55	O 55	0	0
3	J	55	Total 55	O 55	0	0
3	K	56	Total 56	O 56	0	0
3	L	53	Total 53	O 53	0	0
3	M	55	Total 55	O 55	0	0
3	N	57	Total 57	O 57	0	0
3	O	55	Total 55	O 55	0	0
3	P	55	Total 55	O 55	0	0
3	Q	54	Total 54	O 54	0	0
3	R	54	Total 54	O 54	0	0
3	S	54	Total 54	O 54	0	0
3	T	58	Total 58	O 58	0	0
3	U	54	Total 54	O 54	0	0
3	V	54	Total 54	O 54	0	0
3	W	53	Total 53	O 53	0	0
3	X	55	Total 55	O 55	0	0

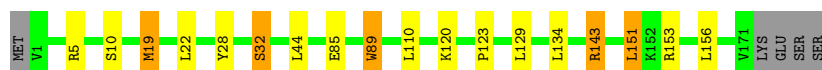
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.


Note EDS was not executed.

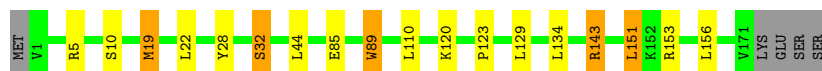
#### • Molecule 1: M FERRITIN

Chain A: 



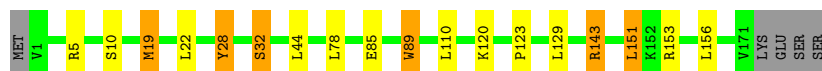
#### • Molecule 1: M FERRITIN

Chain B: 




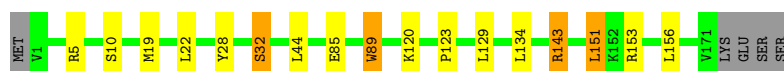
#### • Molecule 1: M FERRITIN

Chain C: 




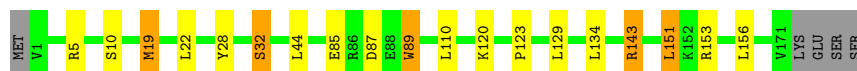
#### • Molecule 1: M FERRITIN

Chain D: 




#### • Molecule 1: M FERRITIN

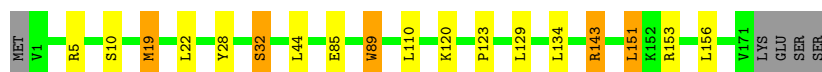
Chain E: 



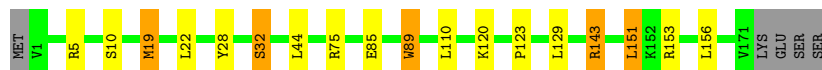
#### • Molecule 1: M FERRITIN

Chain F: 

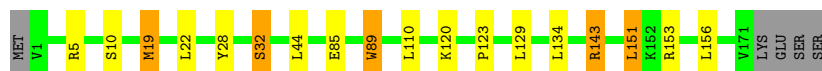
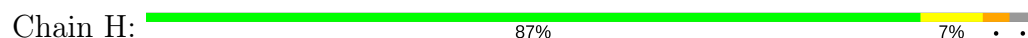




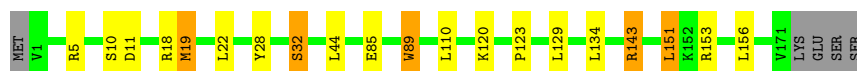
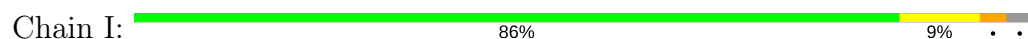
- Molecule 1: M FERRITIN



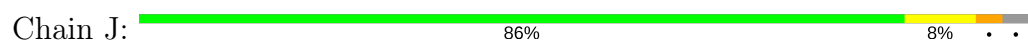
- Molecule 1: M FERRITIN



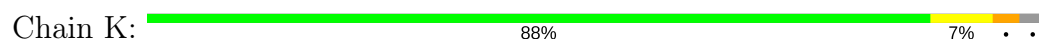
- Molecule 1: M FERRITIN



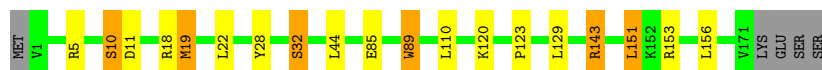
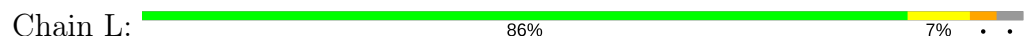
- Molecule 1: M FERRITIN



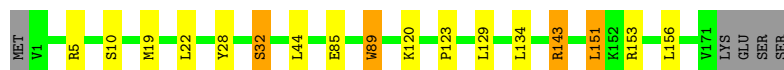
- Molecule 1: M FERRITIN




- Molecule 1: M FERRITIN



- Molecule 1: M FERRITIN




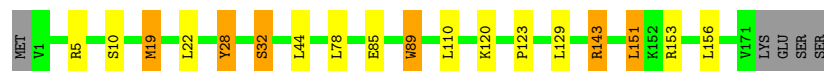
- Molecule 1: M FERRITIN

Chain N:  87% 7% . .




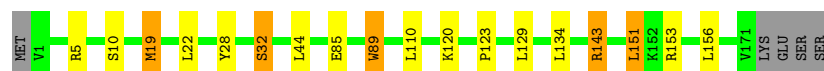
• Molecule 1: M FERRITIN

Chain O:  87% 7% . .




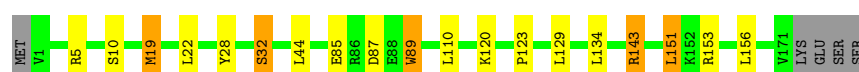
• Molecule 1: M FERRITIN

Chain P:  87% 7% . .




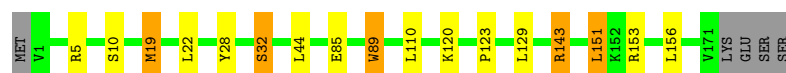
• Molecule 1: M FERRITIN

Chain Q:  86% 8% . .




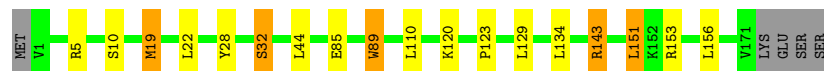
• Molecule 1: M FERRITIN

Chain R:  88% 7% . .




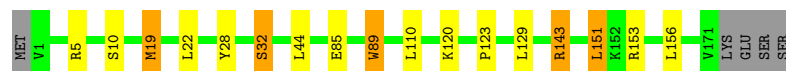
• Molecule 1: M FERRITIN

Chain S:  87% 7% . .




• Molecule 1: M FERRITIN

Chain T:  88% 7% . .



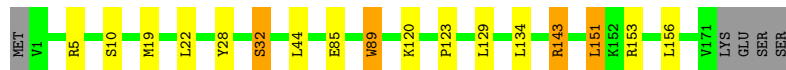
• Molecule 1: M FERRITIN

Chain U:  88% 7% . .



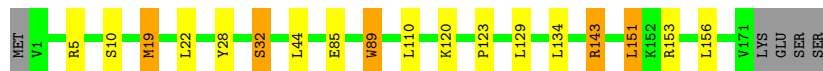
- Molecule 1: M FERRITIN

Chain V: 88% 7% . .



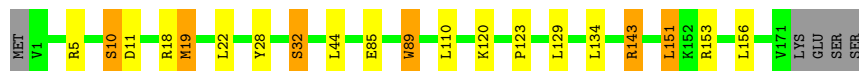
- Molecule 1: M FERRITIN

Chain W: 87% 7% . .



- Molecule 1: M FERRITIN

Chain X: 86% 8% . .



## 4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	210.80Å 210.80Å 328.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.80	Depositor
% Data completeness (in resolution range)	70.0 (10.00-2.80)	Depositor
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.08	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.190 , 0.194	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	34320	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	30.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

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

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	0.75	0/1401	1.21	11/1890 (0.6%)
1	B	0.75	0/1401	1.21	11/1890 (0.6%)
1	C	0.75	0/1401	1.21	11/1890 (0.6%)
1	D	0.75	0/1401	1.21	11/1890 (0.6%)
1	E	0.76	0/1401	1.21	11/1890 (0.6%)
1	F	0.75	0/1401	1.21	11/1890 (0.6%)
1	G	0.75	0/1401	1.21	11/1890 (0.6%)
1	H	0.75	0/1401	1.21	11/1890 (0.6%)
1	I	0.75	0/1401	1.21	11/1890 (0.6%)
1	J	0.75	0/1401	1.21	11/1890 (0.6%)
1	K	0.75	0/1401	1.21	11/1890 (0.6%)
1	L	0.76	0/1401	1.21	11/1890 (0.6%)
1	M	0.76	0/1401	1.21	11/1890 (0.6%)
1	N	0.75	0/1401	1.21	11/1890 (0.6%)
1	O	0.75	0/1401	1.21	11/1890 (0.6%)
1	P	0.76	0/1401	1.21	11/1890 (0.6%)
1	Q	0.76	0/1401	1.21	11/1890 (0.6%)
1	R	0.76	0/1401	1.21	11/1890 (0.6%)
1	S	0.75	0/1401	1.20	11/1890 (0.6%)
1	T	0.75	0/1401	1.21	11/1890 (0.6%)
1	U	0.76	0/1401	1.21	11/1890 (0.6%)
1	V	0.75	0/1401	1.21	11/1890 (0.6%)
1	W	0.75	0/1401	1.21	11/1890 (0.6%)
1	X	0.76	0/1401	1.21	11/1890 (0.6%)
All	All	0.75	0/33624	1.21	264/45360 (0.6%)

There are no bond length outliers.

The worst 5 of 264 bond angle outliers are listed below:

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	T	153	ARG	NE-CZ-NH2	-9.34	115.63	120.30
1	N	153	ARG	NE-CZ-NH2	-9.31	115.65	120.30
1	L	153	ARG	NE-CZ-NH2	-9.30	115.65	120.30
1	M	153	ARG	NE-CZ-NH2	-9.30	115.65	120.30
1	X	153	ARG	NE-CZ-NH2	-9.29	115.66	120.30

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	1373	0	1302	5	0
1	B	1373	0	1302	5	0
1	C	1373	0	1302	7	0
1	D	1373	0	1302	4	0
1	E	1373	0	1302	5	2
1	F	1373	0	1302	5	0
1	G	1373	0	1302	6	0
1	H	1373	0	1302	5	0
1	I	1373	0	1302	5	3
1	J	1373	0	1302	5	1
1	K	1373	0	1302	5	0
1	L	1373	0	1302	5	7
1	M	1373	0	1302	4	0
1	N	1373	0	1302	5	0
1	O	1373	0	1302	7	0
1	P	1373	0	1302	5	0
1	Q	1373	0	1302	5	2
1	R	1373	0	1302	5	0
1	S	1373	0	1302	5	0
1	T	1373	0	1302	5	0
1	U	1373	0	1302	5	0
1	V	1373	0	1302	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	W	1373	0	1302	5	0
1	X	1373	0	1302	5	5
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0
2	K	2	0	0	0	0
2	L	2	0	0	0	0
2	M	2	0	0	0	0
2	N	2	0	0	0	0
2	O	2	0	0	0	0
2	P	2	0	0	0	0
2	Q	2	0	0	0	0
2	R	2	0	0	0	0
2	S	2	0	0	0	0
2	T	2	0	0	0	0
2	U	2	0	0	0	0
2	V	2	0	0	0	0
2	W	2	0	0	0	0
2	X	2	0	0	0	0
3	A	54	0	0	2	0
3	B	56	0	0	2	0
3	C	57	0	0	2	0
3	D	55	0	0	2	0
3	E	55	0	0	2	2
3	F	54	0	0	2	0
3	G	55	0	0	2	0
3	H	57	0	0	2	0
3	I	55	0	0	2	3
3	J	55	0	0	2	2
3	K	56	0	0	2	0
3	L	53	0	0	2	1
3	M	55	0	0	2	0
3	N	57	0	0	2	0
3	O	55	0	0	2	0
3	P	55	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	Q	54	0	0	2	4
3	R	54	0	0	2	0
3	S	54	0	0	2	0
3	T	58	0	0	2	0
3	U	54	0	0	2	2
3	V	54	0	0	2	0
3	W	53	0	0	2	0
3	X	55	0	0	2	2
All	All	34320	0	31248	120	18

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 2.

The worst 5 of 120 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:K:151:LEU:HD12	1:K:156:LEU:HD13	1.82	0.62
1:N:151:LEU:HD12	1:N:156:LEU:HD13	1.82	0.62
1:W:151:LEU:HD12	1:W:156:LEU:HD13	1.82	0.62
1:D:151:LEU:HD12	1:D:156:LEU:HD13	1.82	0.61
1:E:151:LEU:HD12	1:E:156:LEU:HD13	1.82	0.61

The worst 5 of 18 symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:18:ARG:NH1	1:Q:87:ASP:OD2[4_566]	0.93	1.27
1:L:11:ASP:CG	3:Q:224:HOH:O[4_566]	1.31	0.89
1:L:11:ASP:OD1	3:Q:224:HOH:O[4_566]	1.50	0.70
1:E:87:ASP:OD2	1:X:18:ARG:NH1[5_665]	1.53	0.67
1:X:11:ASP:OD1	3:E:221:HOH:O[5_565]	1.56	0.64

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.



The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	B	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	C	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	D	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	E	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	F	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	G	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	H	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	I	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	J	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	K	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	L	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	M	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	N	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	O	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	P	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	Q	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	R	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	S	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	T	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	U	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	V	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	W	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
1	X	169/176 (96%)	165 (98%)	4 (2%)	0	100	100
All	All	4056/4224 (96%)	3960 (98%)	96 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar

resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	B	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	C	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	D	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	E	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	F	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	G	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	H	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	I	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	J	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	K	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	L	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	M	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	N	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	O	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	P	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	Q	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	R	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	S	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	T	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	U	142/157 (90%)	133 (94%)	9 (6%)	21	51
1	V	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	W	142/157 (90%)	132 (93%)	10 (7%)	18	45
1	X	142/157 (90%)	132 (93%)	10 (7%)	18	45
All	All	3408/3768 (90%)	3176 (93%)	232 (7%)	18	47

5 of 232 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	120	LYS
1	N	32	SER

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Mol	Chain	Res	Type
1	W	28	TYR
1	L	22	LEU
1	M	22	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

### 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 carbohydrates 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 ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data

### 6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

### 6.4 Ligands

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers

EDS was not executed - this section is therefore empty.