



# Full wwPDB X-ray Structure Validation Report ⓘ

Nov 16, 2023 – 07:55 AM JST

PDB ID : 6LBC  
Title : shrimp ferritin-T158R  
Authors : Zhao, G.; Chen, H.; Zhang, T.  
Deposited on : 2019-11-14  
Resolution : 1.80 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](mailto: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 : 4.02b-467  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.36

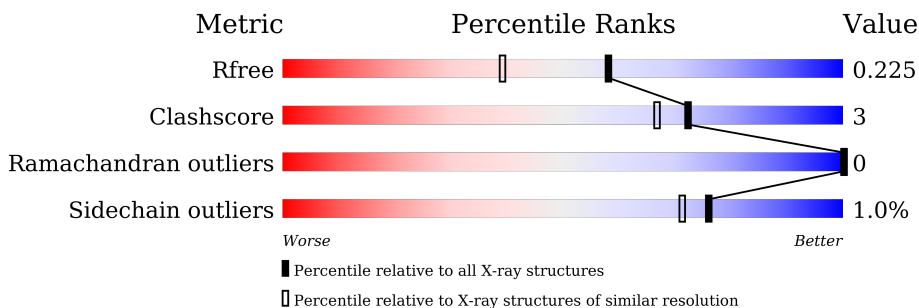
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.80 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))
$R_{free}$	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)

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  $\geq 3$ , 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  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	170	
1	B	170	
1	C	170	
1	D	170	
1	E	170	
1	F	170	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9187 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
			Total	C	N	O	S			
1	A	169	1371	858	232	274	7	0	1	0
1	B	169	1363	854	231	271	7	0	0	0
1	C	169	1371	858	232	274	7	0	1	0
1	D	169	1363	854	231	271	7	0	0	0
1	E	169	1363	854	231	271	7	0	0	0
1	F	169	1372	859	233	273	7	0	1	0

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	158	ARG	THR	engineered mutation	UNP T2B7E1
B	158	ARG	THR	engineered mutation	UNP T2B7E1
C	158	ARG	THR	engineered mutation	UNP T2B7E1
D	158	ARG	THR	engineered mutation	UNP T2B7E1
E	158	ARG	THR	engineered mutation	UNP T2B7E1
F	158	ARG	THR	engineered mutation	UNP T2B7E1

- Molecule 2 is FE (III) ION (three-letter code: FE) (formula: Fe).

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

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Fe 1 1	0	0
2	E	1	Total Fe 1 1	0	0
2	F	1	Total Fe 1 1	0	0


- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	157	Total O 157 157	0	0
3	B	157	Total O 157 157	0	0
3	C	155	Total O 155 155	0	0
3	D	165	Total O 165 165	0	0
3	E	171	Total O 171 171	0	0
3	F	173	Total O 173 173	0	0

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

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

Chain A:  91% 9%



- Molecule 1: Ferritin

Chain B:  93% 6%




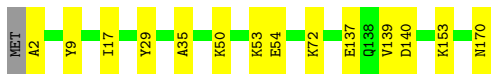
- Molecule 1: Ferritin

Chain C:  92% 8%



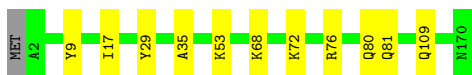
- Molecule 1: Ferritin

Chain D:  91% 8%




- Molecule 1: Ferritin

Chain E:  93% 6%



- Molecule 1: Ferritin

Chain F:  90% 9%



## 4 Data and refinement statistics

Property	Value	Source
Space group	I 4	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	124.78Å 124.78Å 175.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.46 – 1.80 39.46 – 1.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (39.46-1.80) 99.9 (39.46-1.80)	Depositor EDS
$R_{merge}$	0.15	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.55 (at 1.81Å)	Xtrriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, $R_{free}$	0.195 , 0.225 0.195 , 0.225	Depositor DCC
$R_{free}$ test set	1970 reflections (1.60%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	18.3	Xtrriage
Anisotropy	0.046	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 24.7	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.40$ , $\langle L^2 \rangle = 0.22$	Xtrriage
Estimated twinning fraction	0.095 for $-1/2^*h+1/2^*k-1/2^*l, 1/2^*h-1/2^*k-1/2^*l, -h-k$ 0.088 for $-1/2^*h+1/2^*k+1/2^*l, 1/2^*h-1/2^*k+1/2^*l, h+k$ 0.084 for $-1/2^*h-1/2^*k+1/2^*l, -1/2^*h-1/2^*k-1/2^*l, h-k$ 0.084 for $-1/2^*h-1/2^*k-1/2^*l, -1/2^*h-1/2^*k+1/2^*l, -h+k$ 0.267 for $-h, k, -l$	Xtrriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	9187	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.46% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.34	0/1395	0.49	1/1874 (0.1%)
1	B	0.36	0/1387	0.48	0/1863
1	C	0.34	0/1395	0.50	0/1874
1	D	0.35	0/1387	0.49	0/1863
1	E	0.34	0/1387	0.49	0/1863
1	F	0.35	0/1396	0.49	0/1875
All	All	0.35	0/8347	0.49	1/11212 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	114	LEU	CA-CB-CG	5.62	128.22	115.30

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	1371	0	1320	14	0
1	B	1363	0	1317	7	0
1	C	1371	0	1320	8	0
1	D	1363	0	1317	11	0

*Continued on next page...*



Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1363	0	1317	8	0
1	F	1372	0	1324	13	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
2	E	1	0	0	0	0
2	F	1	0	0	0	0
3	A	157	0	0	7	1
3	B	157	0	0	1	0
3	C	155	0	0	3	0
3	D	165	0	0	5	0
3	E	171	0	0	4	0
3	F	173	0	0	2	0
All	All	9187	0	7915	54	1

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 3.

All (54) 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:106:GLN:NE2	3:A:301:HOH:O	1.87	1.05
1:C:60:ARG:NE	3:C:301:HOH:O	1.95	0.99
1:A:102[A]:ASP:OD2	3:A:301:HOH:O	1.98	0.79
1:A:102[A]:ASP:OD2	1:A:106:GLN:NE2	2.18	0.76
1:E:68:LYS:NZ	3:E:302:HOH:O	2.24	0.70
1:A:19:LYS:NZ	3:A:303:HOH:O	2.24	0.70
1:C:170:ASN:ND2	3:C:304:HOH:O	2.27	0.67
1:A:109:GLN:NE2	3:A:304:HOH:O	2.27	0.67
1:D:54:GLU:OE2	3:D:301:HOH:O	2.12	0.67
1:C:133:GLU:OE2	3:C:302:HOH:O	2.13	0.65
1:F:75:GLY:C	1:F:76:ARG:HE	2.00	0.65
1:D:137:GLU:OE2	3:D:302:HOH:O	2.15	0.64
1:E:109:GLN:OE1	3:E:301:HOH:O	2.16	0.62
1:A:137:GLU:OE2	3:A:302:HOH:O	2.16	0.59
1:B:35:ALA:HB1	1:B:53:LYS:HG3	1.85	0.58
1:A:35:ALA:HB1	1:A:53:LYS:HG3	1.87	0.57
1:F:3:SER:HB3	1:F:6:ARG:HB2	1.88	0.56
1:F:35:ALA:HB1	1:F:53:LYS:HG3	1.87	0.55
1:E:9:TYR:OH	1:E:17:ILE:HD12	2.07	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:9:TYR:OH	1:B:17:ILE:HD12	2.08	0.53
1:D:170:ASN:ND2	3:D:306:HOH:O	2.40	0.53
1:B:105:LYS:NZ	3:B:303:HOH:O	2.40	0.53
1:F:9:TYR:OH	1:F:17:ILE:HD12	2.11	0.51
1:C:50:LYS:NZ	1:C:54:GLU:OE2	2.44	0.51
1:C:35:ALA:HB1	1:C:53:LYS:HG3	1.94	0.50
1:A:139:VAL:HG12	1:F:72:LYS:HG2	1.95	0.49
1:D:35:ALA:HB1	1:D:53:LYS:HG3	1.93	0.49
1:E:81:GLN:NE2	3:E:309:HOH:O	2.45	0.49
1:F:167:LYS:NZ	3:F:306:HOH:O	2.38	0.49
1:A:72:LYS:HG2	1:D:139:VAL:HG12	1.96	0.47
1:A:80:GLN:HG3	3:A:440:HOH:O	2.13	0.47
1:C:57:ASP:HA	1:C:60:ARG:HH22	1.80	0.47
1:F:170:ASN:HB3	3:F:391:HOH:O	2.14	0.47
1:E:35:ALA:HB1	1:E:53:LYS:HG3	1.98	0.46
1:F:20:GLN:NE2	1:F:23:MET:CE	2.79	0.46
1:D:50:LYS:NZ	1:D:170:ASN:OXT	2.38	0.45
1:B:72:LYS:HD3	1:C:140:ASP:OD1	2.16	0.45
1:D:9:TYR:OH	1:D:17:ILE:HD12	2.16	0.45
1:F:3:SER:OG	1:F:5:VAL:HG22	2.16	0.45
1:A:20:GLN:O	1:A:24:GLU:HG2	2.16	0.45
1:D:72:LYS:HD3	1:F:140:ASP:OD1	2.17	0.45
1:D:2:ALA:N	3:D:310:HOH:O	2.50	0.44
1:D:153:LYS:NZ	3:D:304:HOH:O	2.27	0.44
1:C:76:ARG:HD3	1:C:76:ARG:HA	1.57	0.43
1:B:63:ALA:O	1:B:67:MET:HG3	2.18	0.43
1:E:80:GLN:NE2	3:E:307:HOH:O	2.40	0.43
1:A:139:VAL:CG1	1:F:72:LYS:HG2	2.50	0.42
1:A:80:GLN:NE2	3:A:313:HOH:O	2.52	0.42
1:A:72:LYS:HD3	1:D:140:ASP:OD1	2.19	0.42
1:B:137:GLU:OE2	1:B:137:GLU:N	2.44	0.41
1:B:139:VAL:HG12	1:E:72:LYS:HG2	2.03	0.41
1:F:20:GLN:HE22	1:F:23:MET:CE	2.34	0.41
1:E:76:ARG:HD3	1:E:76:ARG:HA	1.91	0.40
1:F:23:MET:HE1	1:F:110:SER:HB3	2.03	0.40

All (1) 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)
3:A:307:HOH:O	3:A:409:HOH:O[2_585]	2.19	0.01

## 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	168/170 (99%)	166 (99%)	2 (1%)	0	100	100
1	B	167/170 (98%)	163 (98%)	4 (2%)	0	100	100
1	C	168/170 (99%)	166 (99%)	2 (1%)	0	100	100
1	D	167/170 (98%)	165 (99%)	2 (1%)	0	100	100
1	E	167/170 (98%)	164 (98%)	3 (2%)	0	100	100
1	F	168/170 (99%)	164 (98%)	4 (2%)	0	100	100
All	All	1005/1020 (98%)	988 (98%)	17 (2%)	0	100	100

There are no Ramachandran outliers to report.

### 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	146/146 (100%)	144 (99%)	2 (1%)	67	59
1	B	145/146 (99%)	144 (99%)	1 (1%)	84	81
1	C	146/146 (100%)	143 (98%)	3 (2%)	53	42
1	D	145/146 (99%)	144 (99%)	1 (1%)	84	81
1	E	145/146 (99%)	144 (99%)	1 (1%)	84	81
1	F	146/146 (100%)	145 (99%)	1 (1%)	84	81
All	All	873/876 (100%)	864 (99%)	9 (1%)	76	71

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	TYR
1	A	128	LYS
1	B	29	TYR
1	C	29	TYR
1	C	119	SER
1	C	168	GLU
1	D	29	TYR
1	E	29	TYR
1	F	76	ARG

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

Mol	Chain	Res	Type
1	D	80	GLN
1	F	20	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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 6 ligands modelled in this entry, 6 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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.5 Other polymers [i](#)

Unable to reproduce the depositors R factor - this section is therefore empty.