



# wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 11, 2021 – 02:15 PM JST

PDB ID : 6LS3  
Title : cubic-shaped crystal of TmFtn mutant-T4FY stimulated by Mg  
Authors : Zhao, G.; Zhang, X.  
Deposited on : 2020-01-16  
Resolution : 2.81 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

<https://www.wwpdb.org/validation/2017/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:

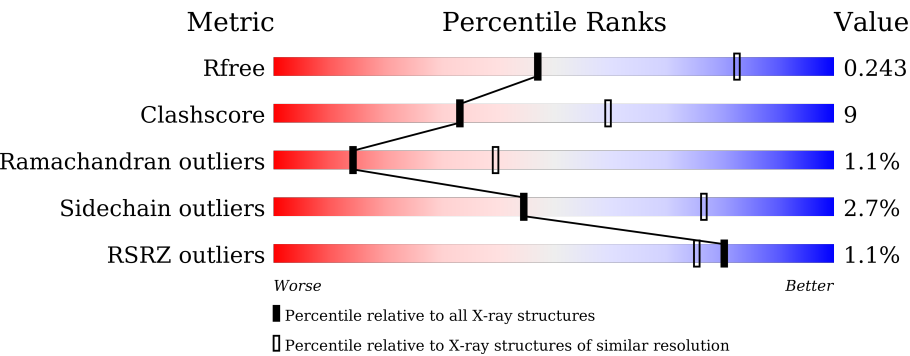
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.16
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.16

# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:  
*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.81 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	3617 (2.84-2.80)
Clashscore	141614	4060 (2.84-2.80)
Ramachandran outliers	138981	3978 (2.84-2.80)
Sidechain outliers	138945	3980 (2.84-2.80)
RSRZ outliers	127900	3552 (2.84-2.80)

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 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  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	164	<div><div>%</div><div><div></div><div>76%</div><div>23%</div><div>.</div></div></div>
1	B	164	<div><div>%</div><div><div></div><div>76%</div><div>23%</div><div>.</div></div></div>
1	C	164	<div><div>2%</div><div><div></div><div>81%</div><div>17%</div><div>.</div></div></div>
1	D	164	<div><div>2%</div><div><div></div><div>79%</div><div>19%</div><div>.</div></div></div>
1	E	164	<div><div>%</div><div><div></div><div>81%</div><div>18%</div><div>.</div></div></div>
1	F	164	<div><div>%</div><div><div></div><div>76%</div><div>22%</div><div>...</div></div></div>

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Mol	Chain	Length	Quality of chain
1	G	164	<div><div>%</div><div><div></div><div>76%</div><div>23%</div><div>.</div></div></div>
1	H	164	<div><div></div><div><div>80%</div><div>16%</div><div>..</div></div></div>

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 10977 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called Ferritin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	164	Total	C	N	O	S	0	0	0
			1368	878	221	263	6			
1	E	164	Total	C	N	O	S	0	0	0
			1368	878	221	263	6			
1	A	164	Total	C	N	O	S	0	0	0
			1368	878	221	263	6			
1	C	164	Total	C	N	O	S	0	0	0
			1368	878	221	263	6			
1	D	164	Total	C	N	O	S	0	0	0
			1368	878	221	263	6			
1	F	163	Total	C	N	O	S	0	0	0
			1360	873	220	262	5			
1	G	164	Total	C	N	O	S	0	0	0
			1368	878	221	263	6			
1	H	163	Total	C	N	O	S	0	0	0
			1360	873	220	262	5			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	83	ALA	ASN	conflict	UNP Q9X0L2
B	147	TYR	ASN	conflict	UNP Q9X0L2
E	83	ALA	ASN	conflict	UNP Q9X0L2
E	147	TYR	ASN	conflict	UNP Q9X0L2
A	83	ALA	ASN	conflict	UNP Q9X0L2
A	147	TYR	ASN	conflict	UNP Q9X0L2
C	83	ALA	ASN	conflict	UNP Q9X0L2
C	147	TYR	ASN	conflict	UNP Q9X0L2
D	83	ALA	ASN	conflict	UNP Q9X0L2
D	147	TYR	ASN	conflict	UNP Q9X0L2
F	83	ALA	ASN	conflict	UNP Q9X0L2
F	147	TYR	ASN	conflict	UNP Q9X0L2
G	83	ALA	ASN	conflict	UNP Q9X0L2

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Chain	Residue	Modelled	Actual	Comment	Reference
G	147	TYR	ASN	conflict	UNP Q9X0L2
H	83	ALA	ASN	conflict	UNP Q9X0L2
H	147	TYR	ASN	conflict	UNP Q9X0L2

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	G	2	Total Fe 2 2	0	0
2	D	2	Total Fe 2 2	0	0
2	E	2	Total Fe 2 2	0	0
2	H	1	Total Fe 1 1	0	0
2	B	2	Total Fe 2 2	0	0
2	C	2	Total Fe 2 2	0	0
2	A	2	Total Fe 2 2	0	0
2	F	2	Total Fe 2 2	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	H	1	Total Mg 1 1	0	0
3	B	1	Total Mg 1 1	0	0
3	A	1	Total Mg 1 1	0	0
3	D	1	Total Mg 1 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	5	Total O 5 5	0	0

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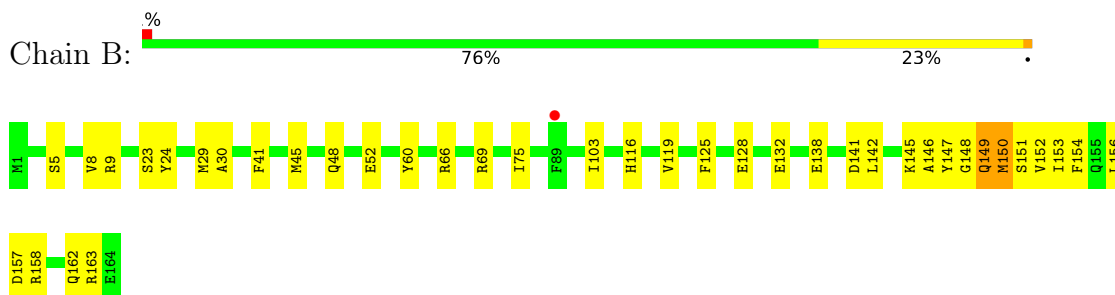
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	E	5	Total 5	O 5	0	0
4	A	5	Total 5	O 5	0	0
4	C	3	Total 3	O 3	0	0
4	D	2	Total 2	O 2	0	0
4	F	6	Total 6	O 6	0	0
4	G	3	Total 3	O 3	0	0
4	H	1	Total 1	O 1	0	0

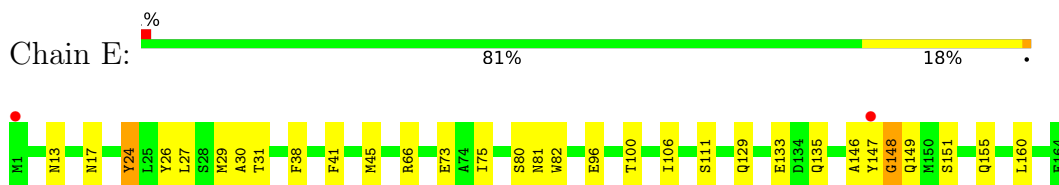
### 3 Residue-property plots [i](#)

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 and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

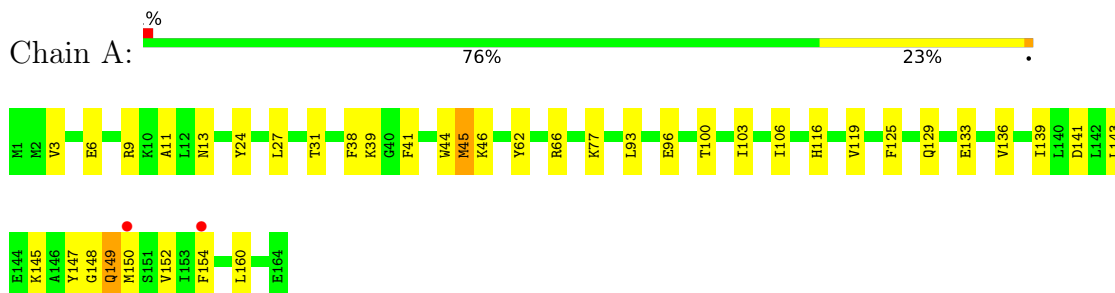
- Molecule 1: Ferritin



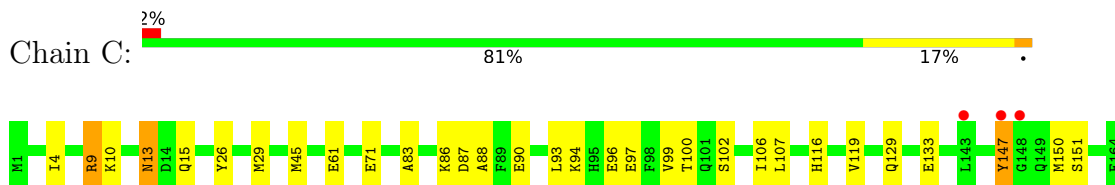
- Molecule 1: Ferritin



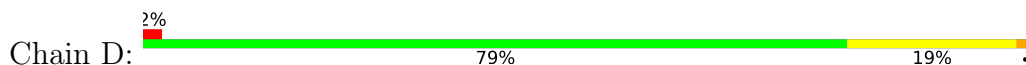
- Molecule 1: Ferritin

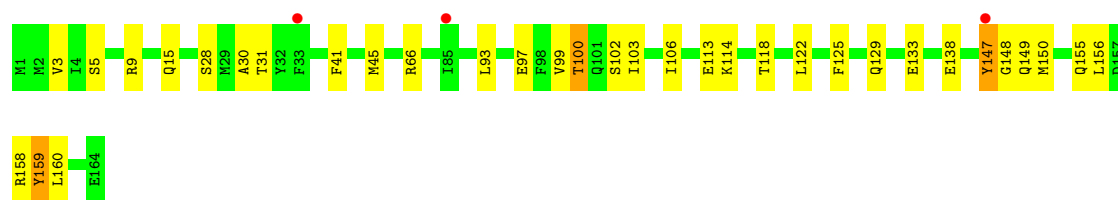


- Molecule 1: Ferritin

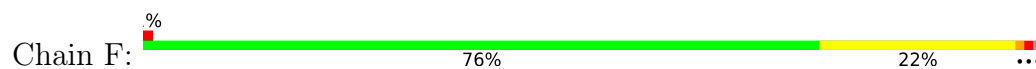


- Molecule 1: Ferritin

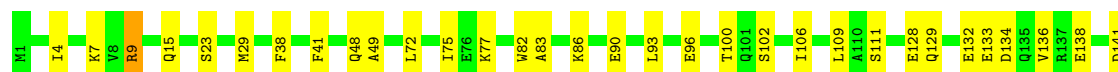
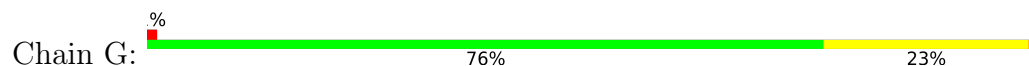




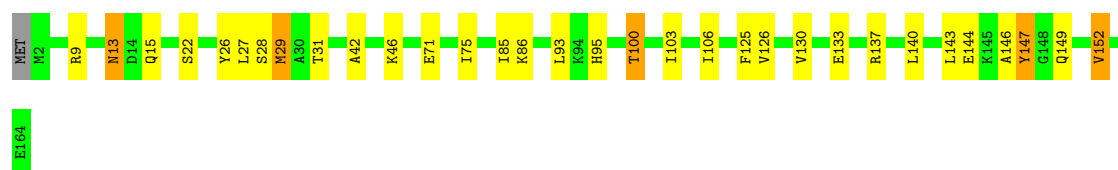
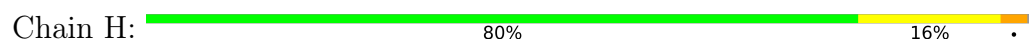
• Molecule 1: Ferritin



• Molecule 1: Ferritin



• Molecule 1: Ferritin





## 4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	171.74Å 171.74Å 384.29Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	138.70 – 2.81 138.70 – 2.82	Depositor EDS
% Data completeness (in resolution range)	94.8 (138.70-2.81) 89.4 (138.70-2.82)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.53 (at 2.82Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, $R_{free}$	0.187 , 0.243 0.189 , 0.243	Depositor DCC
$R_{free}$ test set	2431 reflections (4.82%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	70.8	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 35.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	10977	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.86% 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: MG, 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	0.47	0/1398	0.57	0/1879
1	B	0.47	0/1398	0.59	0/1879
1	C	0.45	0/1398	0.56	0/1879
1	D	0.45	0/1398	0.55	0/1879
1	E	0.49	0/1398	0.59	0/1879
1	F	0.43	0/1390	0.55	0/1869
1	G	0.50	0/1398	0.57	0/1879
1	H	0.56	1/1390 (0.1%)	0.63	0/1869
All	All	0.48	1/11168 (0.0%)	0.58	0/15012

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	H	152	VAL	CB-CG1	-5.51	1.41	1.52

There are no bond angle outliers.

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	1368	0	1321	26	0
1	B	1368	0	1321	30	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	1368	0	1321	23	1
1	D	1368	0	1321	21	0
1	E	1368	0	1321	23	0
1	F	1360	0	1309	26	0
1	G	1368	0	1321	24	0
1	H	1360	0	1309	21	0
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	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	D	1	0	0	0	0
3	H	1	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
4	C	3	0	0	0	0
4	D	2	0	0	0	0
4	E	5	0	0	1	0
4	F	6	0	0	1	0
4	G	3	0	0	0	0
4	H	1	0	0	0	0
All	All	10977	0	10544	185	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 9.

The worst 5 of 185 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:6:GLU:OE2	1:A:9:ARG:NH1	1.95	0.99
1:C:9:ARG:NH1	1:C:10:LYS:HA	1.95	0.80
1:C:4:ILE:HD12	1:C:9:ARG:HB2	1.67	0.75
1:C:26:TYR:CD1	1:C:45:MET:HE1	2.22	0.74
1:E:146:ALA:O	1:E:148:GLY:N	2.19	0.73

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:MET:N	1:C:147:TYR:OH[6_455]	2.05	0.15

## 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	162/164 (99%)	156 (96%)	4 (2%)	2 (1%)	13	37
1	B	162/164 (99%)	158 (98%)	2 (1%)	2 (1%)	13	37
1	C	162/164 (99%)	156 (96%)	4 (2%)	2 (1%)	13	37
1	D	162/164 (99%)	158 (98%)	2 (1%)	2 (1%)	13	37
1	E	162/164 (99%)	159 (98%)	1 (1%)	2 (1%)	13	37
1	F	161/164 (98%)	152 (94%)	7 (4%)	2 (1%)	13	37
1	G	162/164 (99%)	157 (97%)	4 (2%)	1 (1%)	25	54
1	H	161/164 (98%)	157 (98%)	3 (2%)	1 (1%)	25	54
All	All	1294/1312 (99%)	1253 (97%)	27 (2%)	14 (1%)	14	39

5 of 14 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	149	GLN
1	E	147	TYR
1	A	149	GLN
1	D	149	GLN
1	F	151	SER

### 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	143/143 (100%)	140 (98%)	3 (2%)	53	82
1	B	143/143 (100%)	139 (97%)	4 (3%)	43	76
1	C	143/143 (100%)	140 (98%)	3 (2%)	53	82
1	D	143/143 (100%)	139 (97%)	4 (3%)	43	76
1	E	143/143 (100%)	138 (96%)	5 (4%)	36	68
1	F	142/143 (99%)	138 (97%)	4 (3%)	43	76
1	G	143/143 (100%)	139 (97%)	4 (3%)	43	76
1	H	142/143 (99%)	138 (97%)	4 (3%)	43	76
All	All	1142/1144 (100%)	1111 (97%)	31 (3%)	44	77

5 of 31 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	150	MET
1	D	147	TYR
1	H	13	ASN
1	D	9	ARG
1	D	159	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	E	17	ASN
1	E	105	ASN
1	F	135	GLN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 19 ligands modelled in this entry, 19 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](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ > 2	OWAB(Å <sup>2</sup> )	Q < 0.9
1	A	164/164 (100%)	0.25	2 (1%) 79 73	49, 65, 93, 114	0
1	B	164/164 (100%)	0.29	1 (0%) 89 86	50, 62, 92, 116	0
1	C	164/164 (100%)	0.32	3 (1%) 68 61	50, 63, 98, 124	0
1	D	164/164 (100%)	0.28	3 (1%) 68 61	56, 67, 93, 125	0
1	E	164/164 (100%)	0.28	2 (1%) 79 73	51, 64, 93, 125	0
1	F	163/164 (99%)	0.21	2 (1%) 79 73	53, 68, 100, 130	0
1	G	164/164 (100%)	0.27	2 (1%) 79 73	50, 63, 94, 116	0
1	H	163/164 (99%)	0.25	0 100 100	51, 63, 83, 109	0
All	All	1310/1312 (99%)	0.27	15 (1%) 80 75	49, 64, 95, 130	0

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	147	TYR	4.4
1	E	147	TYR	3.5
1	A	150	MET	3.2
1	C	148	GLY	3.1
1	C	147	TYR	2.9

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	MG	D	203	1/1	0.82	0.11	83,83,83,83	0
2	FE	D	202	1/1	0.85	0.14	98,98,98,98	0
3	MG	A	203	1/1	0.89	0.13	81,81,81,81	0
3	MG	B	203	1/1	0.93	0.12	81,81,81,81	0
3	MG	H	202	1/1	0.94	0.15	79,79,79,79	0
2	FE	F	202	1/1	0.97	0.07	106,106,106,106	0
2	FE	C	201	1/1	0.97	0.16	78,78,78,78	0
2	FE	C	202	1/1	0.97	0.11	94,94,94,94	0
2	FE	E	201	1/1	0.98	0.21	75,75,75,75	0
2	FE	E	202	1/1	0.98	0.10	98,98,98,98	0
2	FE	B	201	1/1	0.98	0.16	76,76,76,76	0
2	FE	A	202	1/1	0.98	0.09	98,98,98,98	0
2	FE	A	201	1/1	0.98	0.14	77,77,77,77	0
2	FE	G	202	1/1	0.98	0.09	94,94,94,94	0
2	FE	H	201	1/1	0.99	0.15	84,84,84,84	0
2	FE	F	201	1/1	0.99	0.14	81,81,81,81	0
2	FE	B	202	1/1	0.99	0.10	89,89,89,89	0
2	FE	G	201	1/1	0.99	0.17	74,74,74,74	0
2	FE	D	201	1/1	0.99	0.17	79,79,79,79	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.