



Full wwPDB X-ray Structure Validation Report ⓘ

May 22, 2020 – 04:25 am BST

PDB ID : 5U1B
Title : Ferritin with Gc MtrE loop2 inserted at the N-terminus
Authors : Wang, S.
Deposited on : 2016-11-28
Resolution : 2.81 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The 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 : 2.11
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) : 2.11

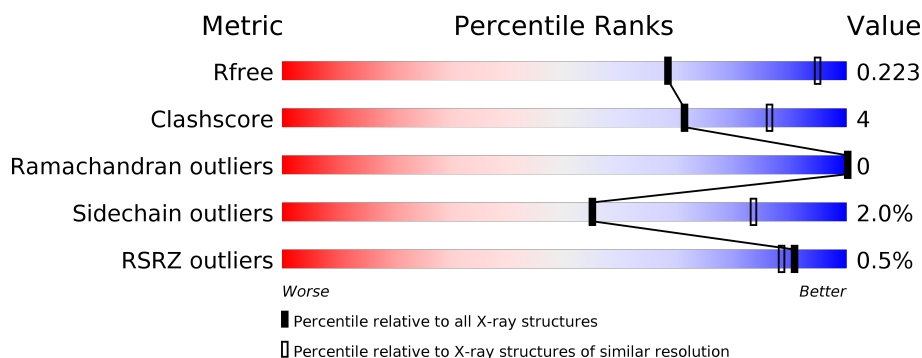
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.81 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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	189	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>11%</div> <div>11%</div> </div> </div>
1	B	189	<div> <div>%</div> <div> <div></div> <div>81%</div> <div>5%</div> <div>12%</div> </div> </div>
1	C	189	<div> <div></div> <div> <div>76%</div> <div>12%</div> <div>12%</div> </div> </div>
1	D	189	<div> <div></div> <div> <div>82%</div> <div>6%</div> <div>12%</div> </div> </div>
1	E	189	<div> <div>%</div> <div> <div></div> <div>79%</div> <div>8%</div> <div>12%</div> </div> </div>
1	F	189	<div> <div></div> <div> <div>81%</div> <div>6%</div> <div>12%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	G	189	<div><div>%</div><div><div></div></div><div>83%5% • 12%</div></div>
1	H	189	<div><div>%</div><div><div></div></div><div>80%8% • 11%</div></div>

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10895 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 MtrE protein, Ferritin chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	168	Total	C	N	O	S	0	0	0
			1374	875	229	265	5			
1	B	166	Total	C	N	O	S	0	0	0
			1356	865	225	261	5			
1	C	166	Total	C	N	O	S	0	0	0
			1356	865	225	261	5			
1	D	166	Total	C	N	O	S	0	0	0
			1356	865	225	261	5			
1	E	166	Total	C	N	O	S	0	0	0
			1356	865	225	261	5			
1	F	166	Total	C	N	O	S	0	0	0
			1356	865	225	261	5			
1	G	166	Total	C	N	O	S	0	0	0
			1356	865	225	261	5			
1	H	168	Total	C	N	O	S	0	0	0
			1374	875	229	265	5			

There are 64 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	initiating methionine	UNP Q51006
A	-20	ALA	-	expression tag	UNP Q51006
A	-19	ARG	-	expression tag	UNP Q51006
A	-18	CYS	-	expression tag	UNP Q51006
A	-2	CYS	-	linker	UNP Q51006
A	-1	ASP	-	linker	UNP Q51006
A	0	HIS	-	linker	UNP Q51006
A	125	SER	ALA	conflict	UNP O69434
B	-21	MET	-	initiating methionine	UNP Q51006
B	-20	ALA	-	expression tag	UNP Q51006
B	-19	ARG	-	expression tag	UNP Q51006
B	-18	CYS	-	expression tag	UNP Q51006
B	-2	CYS	-	linker	UNP Q51006

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-1	ASP	-	linker	UNP Q51006
B	0	HIS	-	linker	UNP Q51006
B	125	SER	ALA	conflict	UNP O69434
C	-21	MET	-	initiating methionine	UNP Q51006
C	-20	ALA	-	expression tag	UNP Q51006
C	-19	ARG	-	expression tag	UNP Q51006
C	-18	CYS	-	expression tag	UNP Q51006
C	-2	CYS	-	linker	UNP Q51006
C	-1	ASP	-	linker	UNP Q51006
C	0	HIS	-	linker	UNP Q51006
C	125	SER	ALA	conflict	UNP O69434
D	-21	MET	-	initiating methionine	UNP Q51006
D	-20	ALA	-	expression tag	UNP Q51006
D	-19	ARG	-	expression tag	UNP Q51006
D	-18	CYS	-	expression tag	UNP Q51006
D	-2	CYS	-	linker	UNP Q51006
D	-1	ASP	-	linker	UNP Q51006
D	0	HIS	-	linker	UNP Q51006
D	125	SER	ALA	conflict	UNP O69434
E	-21	MET	-	initiating methionine	UNP Q51006
E	-20	ALA	-	expression tag	UNP Q51006
E	-19	ARG	-	expression tag	UNP Q51006
E	-18	CYS	-	expression tag	UNP Q51006
E	-2	CYS	-	linker	UNP Q51006
E	-1	ASP	-	linker	UNP Q51006
E	0	HIS	-	linker	UNP Q51006
E	125	SER	ALA	conflict	UNP O69434
F	-21	MET	-	initiating methionine	UNP Q51006
F	-20	ALA	-	expression tag	UNP Q51006
F	-19	ARG	-	expression tag	UNP Q51006
F	-18	CYS	-	expression tag	UNP Q51006
F	-2	CYS	-	linker	UNP Q51006
F	-1	ASP	-	linker	UNP Q51006
F	0	HIS	-	linker	UNP Q51006
F	125	SER	ALA	conflict	UNP O69434
G	-21	MET	-	initiating methionine	UNP Q51006
G	-20	ALA	-	expression tag	UNP Q51006
G	-19	ARG	-	expression tag	UNP Q51006
G	-18	CYS	-	expression tag	UNP Q51006
G	-2	CYS	-	linker	UNP Q51006
G	-1	ASP	-	linker	UNP Q51006
G	0	HIS	-	linker	UNP Q51006

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Chain	Residue	Modelled	Actual	Comment	Reference
G	125	SER	ALA	conflict	UNP O69434
H	-21	MET	-	initiating methionine	UNP Q51006
H	-20	ALA	-	expression tag	UNP Q51006
H	-19	ARG	-	expression tag	UNP Q51006
H	-18	CYS	-	expression tag	UNP Q51006
H	-2	CYS	-	linker	UNP Q51006
H	-1	ASP	-	linker	UNP Q51006
H	0	HIS	-	linker	UNP Q51006
H	125	SER	ALA	conflict	UNP O69434

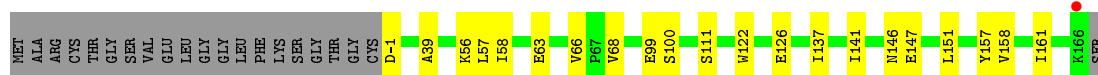
- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	5	Total O 5 5	0	0
2	B	1	Total O 1 1	0	0
2	C	1	Total O 1 1	0	0
2	G	3	Total O 3 3	0	0
2	H	1	Total O 1 1	0	0

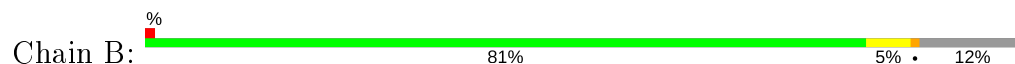
3 Residue-property plots [i](#)

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.

- Molecule 1: MtrE protein,Ferritin chimera



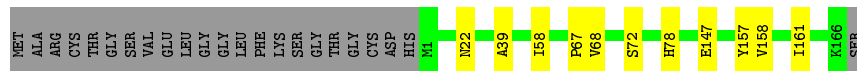
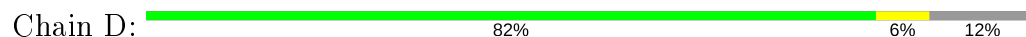
- Molecule 1: MtrE protein,Ferritin chimera



- Molecule 1: MtrE protein,Ferritin chimera



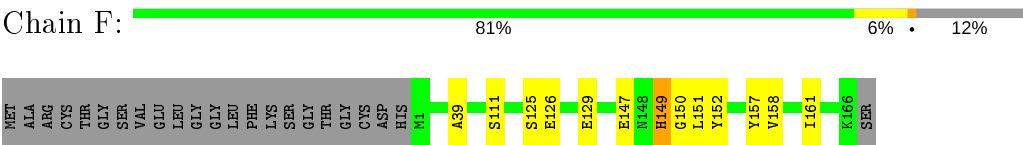
- Molecule 1: MtrE protein,Ferritin chimera



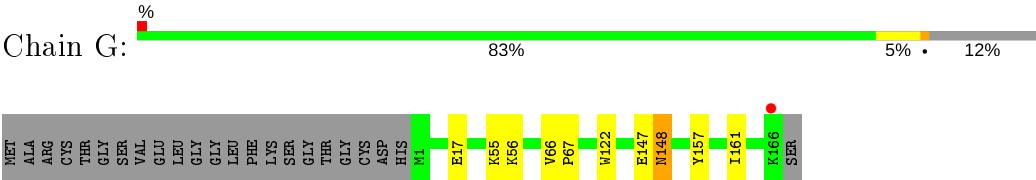
- Molecule 1: MtrE protein,Ferritin chimera



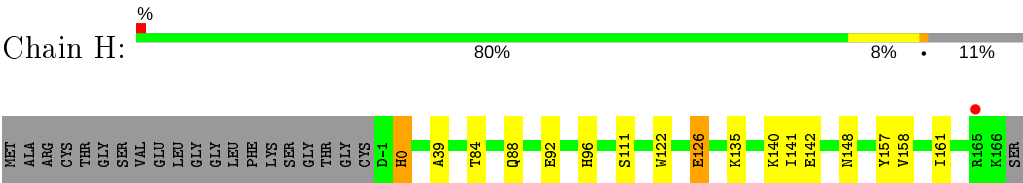
● Molecule 1: MtrE protein,Ferritin chimera



● Molecule 1: MtrE protein,Ferritin chimera



● Molecule 1: MtrE protein,Ferritin chimera



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	124.51Å 124.51Å 314.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.89 – 2.81 19.89 – 2.81	Depositor EDS
% Data completeness (in resolution range)	92.4 (19.89-2.81) 92.7 (19.89-2.81)	Depositor EDS
R_{merge}	0.17	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 2.79Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.177 , 0.225 0.178 , 0.223	Depositor DCC
R_{free} test set	2273 reflections (5.52%)	wwPDB-VP
Wilson B-factor (Å ²)	53.1	Xtriage
Anisotropy	0.408	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 9.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.056 for $-1/3^*h+1/3^*k+1/3^*l, -k, 8/3^*h+4/3^*k+1/3^*l$ 0.079 for $-2/3^*h-1/3^*k-1/3^*l, -1/3^*h-2/3^*k+1/3^*l, -4/3^*h+4/3^*k+1/3^*l$ 0.057 for $-h, 1/3^*h-1/3^*k-1/3^*l, -4/3^*h-8/3^*k+1/3^*l$ 0.056 for $-h, 2/3^*h+1/3^*k+1/3^*l, 4/3^*h+8/3^*k-1/3^*l$ 0.056 for $1/3^*h+2/3^*k-1/3^*l, -k, -8/3^*h-4/3^*k-1/3^*l$ 0.080 for $-1/3^*h-2/3^*k+1/3^*l, -2/3^*h-1/3^*k-1/3^*l, 4/3^*h-4/3^*k-1/3^*l$ 0.437 for $-h-k, k, -l$	Xtriage
Reported twinning fraction	0.545 for H, K, L 0.455 for K, H, -L	Depositor
Outliers	0 of 41193 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	10895	wwPDB-VP
Average B, all atoms (Å ²)	62.0	wwPDB-VP

¹Intensities estimated from amplitudes.

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

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.*

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 >5	RMSZ	# Z >5
1	A	1.06	1/1405 (0.1%)	0.84	1/1895 (0.1%)
1	B	0.98	0/1386	0.80	0/1869
1	C	0.99	0/1386	0.80	0/1869
1	D	0.92	0/1386	0.83	0/1869
1	E	0.90	0/1386	0.80	0/1869
1	F	0.90	0/1386	0.79	1/1869 (0.1%)
1	G	1.05	0/1386	0.81	0/1869
1	H	1.04	0/1405	0.78	0/1895
All	All	0.98	1/11126 (0.0%)	0.81	2/15004 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	63	GLU	CG-CD	5.17	1.59	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	147	GLU	N-CA-C	5.13	124.86	111.00
1	F	149	HIS	C-N-CA	-5.03	111.74	122.30

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	1374	0	1326	8	0
1	B	1356	0	1315	11	0
1	C	1356	0	1315	13	0
1	D	1356	0	1315	5	0
1	E	1356	0	1315	10	0
1	F	1356	0	1315	7	0
1	G	1356	0	1315	8	0
1	H	1374	0	1326	14	0
2	A	5	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	G	3	0	0	0	0
2	H	1	0	0	0	0
All	All	10895	0	10542	75	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (75) 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:E:137:ILE:O	1:E:141:ILE:HG13	1.75	0.86
1:C:18:MET:HE1	1:D:22:ASN:OD1	1.81	0.80
1:F:149:HIS:O	1:F:150:GLY:C	2.17	0.80
1:F:149:HIS:O	1:F:152:TYR:N	2.20	0.74
1:H:96:HIS:O	1:H:96:HIS:ND1	2.25	0.69
1:H:122:TRP:HE3	1:H:126:GLU:OE1	1.77	0.67
1:C:77:GLU:OE2	1:C:78:HIS:N	2.26	0.65
1:G:148:ASN:OD1	1:G:148:ASN:N	2.30	0.63
1:B:148:ASN:HD22	1:B:148:ASN:N	1.95	0.63
1:H:122:TRP:CE3	1:H:126:GLU:OE1	2.53	0.61
1:B:135:LYS:HD3	1:B:135:LYS:C	2.21	0.61
1:B:122:TRP:CE3	1:B:126:GLU:OE1	2.56	0.58
1:A:-1:ASP:OD2	1:A:66:VAL:HG22	2.03	0.57
1:C:142:GLU:OE1	1:C:142:GLU:HA	2.04	0.57
1:C:94:GLU:OE2	1:C:127:GLN:NE2	2.34	0.57
1:E:111:SER:O	1:E:112:LYS:HB2	2.05	0.57
1:A:58:ILE:HG23	1:A:68:VAL:HG21	1.87	0.56
1:H:142:GLU:HA	1:H:142:GLU:OE1	2.05	0.56
1:A:137:ILE:O	1:A:141:ILE:HG13	2.05	0.56
1:H:88:GLN:O	1:H:92:GLU:HG3	2.05	0.56
1:E:142:GLU:OE1	1:E:142:GLU:HA	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:147:GLU:O	1:D:147:GLU:HG3	2.07	0.54
1:H:96:HIS:C	1:H:96:HIS:HD1	2.12	0.52
1:G:147:GLU:H	1:G:147:GLU:CD	2.12	0.52
1:C:148:ASN:HB3	1:C:149:HIS:HD2	1.76	0.51
1:F:125:SER:O	1:F:129:GLU:HG3	2.10	0.51
1:H:140:LYS:HG3	1:H:157:TYR:CZ	2.46	0.51
1:F:149:HIS:O	1:F:151:LEU:N	2.44	0.50
1:B:135:LYS:O	1:B:135:LYS:HD3	2.12	0.49
1:F:147:GLU:HA	1:F:147:GLU:OE2	2.13	0.49
1:C:148:ASN:HB3	1:C:149:HIS:CD2	2.48	0.49
1:G:56:LYS:HD2	1:G:122:TRP:CZ2	2.48	0.48
1:B:148:ASN:HD22	1:B:148:ASN:H	1.59	0.48
1:F:157:TYR:CZ	1:F:161:ILE:HD11	2.49	0.48
1:G:157:TYR:CZ	1:G:161:ILE:HD11	2.49	0.47
1:B:157:TYR:CZ	1:B:161:ILE:HD11	2.50	0.47
1:C:152:TYR:O	1:C:156:GLN:HG2	2.14	0.47
1:C:157:TYR:CZ	1:C:161:ILE:HD11	2.50	0.46
1:A:157:TYR:CZ	1:A:161:ILE:HD11	2.50	0.46
1:D:58:ILE:HG23	1:D:68:VAL:HG21	1.96	0.46
1:H:157:TYR:CZ	1:H:161:ILE:HD11	2.51	0.46
1:C:1:MET:HG2	1:C:2:LEU:HG	1.97	0.46
1:E:157:TYR:CZ	1:E:161:ILE:HD11	2.51	0.46
1:E:149:HIS:O	1:E:150:GLY:C	2.52	0.45
1:H:148:ASN:ND2	1:H:148:ASN:N	2.62	0.45
1:D:157:TYR:CZ	1:D:161:ILE:HD11	2.52	0.45
1:H:39:ALA:HA	1:H:158:VAL:HG11	1.99	0.45
1:A:57:LEU:HD23	1:A:57:LEU:HA	1.63	0.45
1:E:140:LYS:HG2	1:E:157:TYR:CZ	2.52	0.45
1:B:39:ALA:HA	1:B:158:VAL:HG11	2.00	0.44
1:C:58:ILE:HG23	1:C:68:VAL:HG21	1.99	0.44
1:H:0:HIS:CD2	1:H:0:HIS:H	2.34	0.44
1:B:94:GLU:OE2	1:B:127:GLN:NE2	2.38	0.43
1:E:58:ILE:HG23	1:E:68:VAL:HG21	2.01	0.43
1:A:56:LYS:HD3	1:A:122:TRP:CZ2	2.54	0.43
1:C:39:ALA:HA	1:C:158:VAL:HG11	2.01	0.43
1:G:147:GLU:N	1:G:147:GLU:CD	2.72	0.43
1:E:39:ALA:HA	1:E:158:VAL:HG11	2.01	0.42
1:A:39:ALA:HA	1:A:158:VAL:HG11	2.01	0.42
1:F:39:ALA:HA	1:F:158:VAL:HG11	2.02	0.42
1:E:148:ASN:N	1:E:148:ASN:ND2	2.65	0.42
1:G:56:LYS:HD2	1:G:122:TRP:CH2	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:149:HIS:N	1:C:149:HIS:CD2	2.88	0.42
1:D:39:ALA:HA	1:D:158:VAL:HG11	2.02	0.42
1:B:148:ASN:ND2	1:B:148:ASN:N	2.60	0.41
1:H:84:THR:HG22	1:H:141:ILE:CG2	2.50	0.41
1:H:84:THR:HG22	1:H:141:ILE:HG21	2.02	0.41
1:B:58:ILE:HG23	1:B:68:VAL:HG21	2.02	0.41
1:H:96:HIS:C	1:H:96:HIS:ND1	2.71	0.40
1:B:122:TRP:HE3	1:B:126:GLU:OE1	2.02	0.40
1:A:151:LEU:HA	1:A:151:LEU:HD23	1.82	0.40
1:E:151:LEU:HA	1:E:151:LEU:HD23	1.94	0.40
1:G:55:LYS:HB3	1:G:55:LYS:HE2	1.83	0.40
1:C:1:MET:CG	1:C:2:LEU:N	2.83	0.40
1:G:66:VAL:HA	1:G:67:PRO:HD3	1.92	0.40

There are no symmetry-related clashes.

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	166/189 (88%)	162 (98%)	4 (2%)	0	100	100
1	B	164/189 (87%)	160 (98%)	4 (2%)	0	100	100
1	C	164/189 (87%)	160 (98%)	4 (2%)	0	100	100
1	D	164/189 (87%)	159 (97%)	5 (3%)	0	100	100
1	E	164/189 (87%)	160 (98%)	4 (2%)	0	100	100
1	F	164/189 (87%)	159 (97%)	5 (3%)	0	100	100
1	G	164/189 (87%)	160 (98%)	4 (2%)	0	100	100
1	H	166/189 (88%)	162 (98%)	4 (2%)	0	100	100
All	All	1316/1512 (87%)	1282 (97%)	34 (3%)	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	152/167 (91%)	147 (97%)	5 (3%)	38	70
1	B	150/167 (90%)	148 (99%)	2 (1%)	69	90
1	C	150/167 (90%)	146 (97%)	4 (3%)	44	77
1	D	150/167 (90%)	147 (98%)	3 (2%)	55	83
1	E	150/167 (90%)	148 (99%)	2 (1%)	69	90
1	F	150/167 (90%)	148 (99%)	2 (1%)	69	90
1	G	150/167 (90%)	148 (99%)	2 (1%)	69	90
1	H	152/167 (91%)	148 (97%)	4 (3%)	46	78
All	All	1204/1336 (90%)	1180 (98%)	24 (2%)	55	83

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	99	GLU
1	A	100	SER
1	A	111	SER
1	A	126	GLU
1	A	146	ASN
1	B	126	GLU
1	B	135	LYS
1	C	37	ASP
1	C	67	PRO
1	C	99	GLU
1	C	111	SER
1	D	67	PRO
1	D	72	SER
1	D	78	HIS
1	E	67	PRO
1	E	111	SER

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Mol	Chain	Res	Type
1	F	111	SER
1	F	126	GLU
1	G	17	GLU
1	G	148	ASN
1	H	0	HIS
1	H	111	SER
1	H	126	GLU
1	H	135	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	148	ASN
1	C	15	ASN
1	C	128	HIS
1	C	148	ASN
1	C	149	HIS
1	E	148	ASN
1	F	11	ASN
1	F	65	ASN
1	H	0	HIS
1	H	148	ASN

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.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	168/189 (88%)	-0.55	1 (0%) 89 86	36, 56, 92, 137	2 (1%)
1	B	166/189 (87%)	-0.51	1 (0%) 89 86	31, 57, 87, 109	2 (1%)
1	C	166/189 (87%)	-0.53	0 100 100	37, 60, 89, 122	2 (1%)
1	D	166/189 (87%)	-0.54	0 100 100	37, 59, 82, 123	0
1	E	166/189 (87%)	-0.52	2 (1%) 79 73	44, 64, 93, 126	0
1	F	166/189 (87%)	-0.46	0 100 100	44, 64, 93, 120	4 (2%)
1	G	166/189 (87%)	-0.53	1 (0%) 89 86	35, 55, 92, 136	2 (1%)
1	H	168/189 (88%)	-0.54	1 (0%) 89 86	39, 55, 98, 124	1 (0%)
All	All	1332/1512 (88%)	-0.52	6 (0%) 91 88	31, 59, 93, 137	13 (0%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	G	166	LYS	4.1
1	E	166	LYS	2.7
1	H	165	ARG	2.2
1	B	1	MET	2.2
1	E	1	MET	2.1
1	A	166	LYS	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.