



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 7, 2024 – 10:44 am GMT

PDB ID : 5N5E
Title : Crystal structure of encapsulated ferritin domain from *Pyrococcus furiosus* PFC_05175
Authors : Marles-Wright, J.; He, D.
Deposited on : 2017-02-13
Resolution : 2.03 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

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

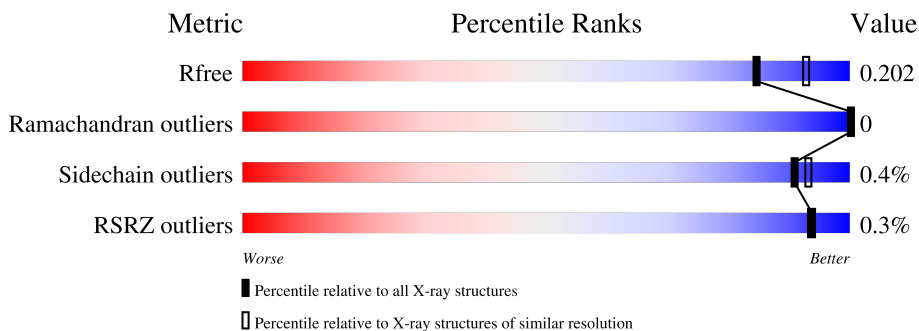
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.03 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.



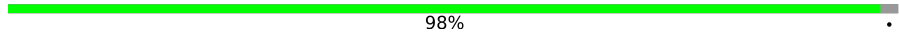
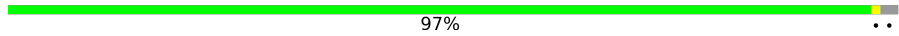
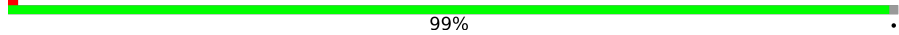
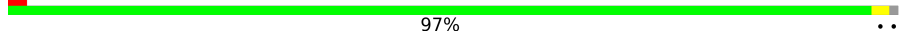
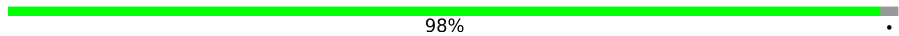
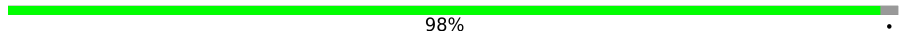
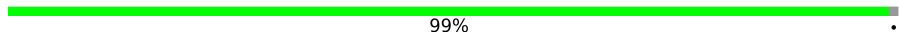
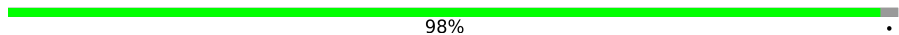
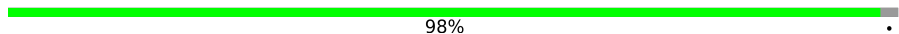
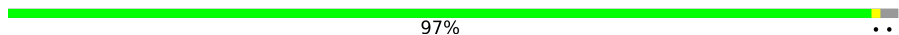
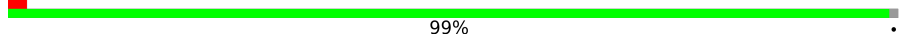
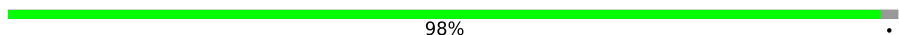
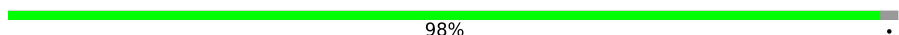
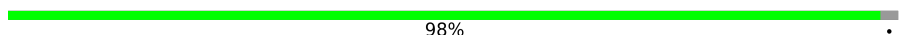
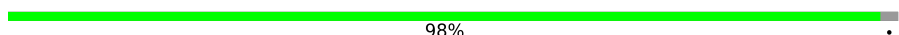
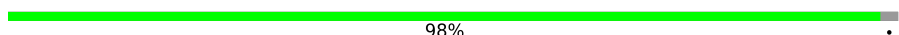
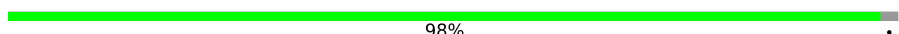
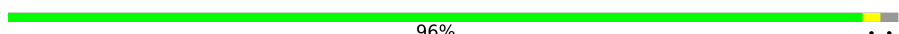
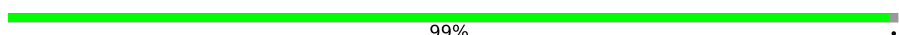
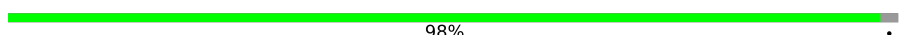
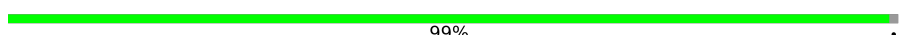
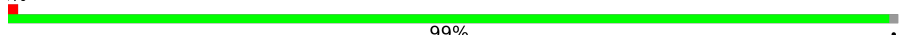
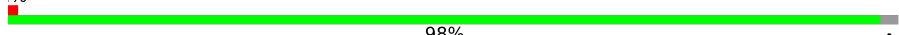
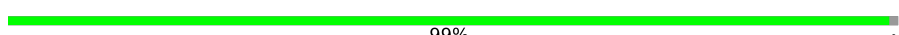
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	10434 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 ≥ 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	99	
1	B	99	
1	C	99	
1	D	99	
1	E	99	
1	F	99	

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Mol	Chain	Length	Quality of chain
1	G	99	 98%
1	H	99	 97%
1	I	99	 99%
1	J	99	 97%
1	K	99	 98%
1	L	99	 98%
1	M	99	 99%
1	N	99	 98%
1	O	99	 98%
1	P	99	 97%
1	Q	99	 99%
1	R	99	 98%
1	T	99	 98%
1	U	99	 98%
1	V	99	 98%
1	W	99	 98%
1	X	99	 98%
1	Y	99	 96%
1	Z	99	 99%
1	a	99	 98%
1	b	99	 99%
1	c	99	 99%
1	d	99	 98%
1	e	99	 99%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 49851 atoms, of which 24116 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 PFC_05175.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	e	98	1591	500	803	128	156	4	0	1	0
1	A	98	1588	500	799	128	158	3	0	1	0
1	B	98	1592	500	804	128	156	4	0	1	0
1	C	98	1598	501	807	131	156	3	0	1	0
1	D	97	1596	502	804	128	158	4	0	2	0
1	E	98	1606	505	809	131	158	3	0	2	0
1	F	97	1579	497	795	127	157	3	0	1	0
1	G	97	1585	498	799	128	156	4	0	1	0
1	H	97	1579	497	795	127	157	3	0	1	0
1	I	98	1589	500	800	128	158	3	0	1	0
1	J	98	1619	506	820	134	156	3	0	2	0
1	K	97	1612	506	815	130	157	4	0	3	0
1	L	97	1625	510	823	130	157	5	0	4	0
1	M	98	1602	504	808	128	158	4	0	2	0
1	N	97	1592	501	803	127	157	4	0	2	0
1	O	97	1603	503	808	131	158	3	0	2	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
			Total	C	H	N	O				S
1	P	97	Total 1587	C 499	H 801	N 127	O 156	S 4	0	2	0
1	Q	98	Total 1591	C 500	H 803	N 128	O 156	S 4	0	1	0
1	R	97	Total 1581	C 497	H 798	N 127	O 155	S 4	0	1	0
1	T	97	Total 1609	C 506	H 812	N 128	O 158	S 5	0	3	0
1	U	97	Total 1581	C 497	H 798	N 127	O 155	S 4	0	1	0
1	V	97	Total 1587	C 499	H 800	N 127	O 158	S 3	0	2	0
1	W	97	Total 1592	C 501	H 803	N 127	O 157	S 4	0	2	0
1	X	97	Total 1579	C 497	H 795	N 127	O 157	S 3	0	1	0
1	Y	97	Total 1599	C 502	H 807	N 130	O 157	S 3	0	2	0
1	Z	98	Total 1589	C 500	H 800	N 128	O 158	S 3	0	1	0
1	a	97	Total 1579	C 497	H 795	N 127	O 157	S 3	0	1	0
1	b	98	Total 1578	C 496	H 795	N 128	O 156	S 3	0	0	0
1	c	98	Total 1631	C 510	H 827	N 134	O 156	S 4	0	3	0
1	d	97	Total 1568	C 493	H 790	N 127	O 155	S 3	0	0	0

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
e	1	MET	-	initiating methionine	UNP I6U7J4
e	2	GLY	-	expression tag	UNP I6U7J4
A	1	MET	-	initiating methionine	UNP I6U7J4
A	2	GLY	-	expression tag	UNP I6U7J4
B	1	MET	-	initiating methionine	UNP I6U7J4
B	2	GLY	-	expression tag	UNP I6U7J4
C	1	MET	-	initiating methionine	UNP I6U7J4
C	2	GLY	-	expression tag	UNP I6U7J4
D	1	MET	-	initiating methionine	UNP I6U7J4
D	2	GLY	-	expression tag	UNP I6U7J4
E	1	MET	-	initiating methionine	UNP I6U7J4

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Chain	Residue	Modelled	Actual	Comment	Reference
E	2	GLY	-	expression tag	UNP I6U7J4
F	1	MET	-	initiating methionine	UNP I6U7J4
F	2	GLY	-	expression tag	UNP I6U7J4
G	1	MET	-	initiating methionine	UNP I6U7J4
G	2	GLY	-	expression tag	UNP I6U7J4
H	1	MET	-	initiating methionine	UNP I6U7J4
H	2	GLY	-	expression tag	UNP I6U7J4
I	1	MET	-	initiating methionine	UNP I6U7J4
I	2	GLY	-	expression tag	UNP I6U7J4
J	1	MET	-	initiating methionine	UNP I6U7J4
J	2	GLY	-	expression tag	UNP I6U7J4
K	1	MET	-	initiating methionine	UNP I6U7J4
K	2	GLY	-	expression tag	UNP I6U7J4
L	1	MET	-	initiating methionine	UNP I6U7J4
L	2	GLY	-	expression tag	UNP I6U7J4
M	1	MET	-	initiating methionine	UNP I6U7J4
M	2	GLY	-	expression tag	UNP I6U7J4
N	1	MET	-	initiating methionine	UNP I6U7J4
N	2	GLY	-	expression tag	UNP I6U7J4
O	1	MET	-	initiating methionine	UNP I6U7J4
O	2	GLY	-	expression tag	UNP I6U7J4
P	1	MET	-	initiating methionine	UNP I6U7J4
P	2	GLY	-	expression tag	UNP I6U7J4
Q	1	MET	-	initiating methionine	UNP I6U7J4
Q	2	GLY	-	expression tag	UNP I6U7J4
R	1	MET	-	initiating methionine	UNP I6U7J4
R	2	GLY	-	expression tag	UNP I6U7J4
T	1	MET	-	initiating methionine	UNP I6U7J4
T	2	GLY	-	expression tag	UNP I6U7J4
U	1	MET	-	initiating methionine	UNP I6U7J4
U	2	GLY	-	expression tag	UNP I6U7J4
V	1	MET	-	initiating methionine	UNP I6U7J4
V	2	GLY	-	expression tag	UNP I6U7J4
W	1	MET	-	initiating methionine	UNP I6U7J4
W	2	GLY	-	expression tag	UNP I6U7J4
X	1	MET	-	initiating methionine	UNP I6U7J4
X	2	GLY	-	expression tag	UNP I6U7J4
Y	1	MET	-	initiating methionine	UNP I6U7J4
Y	2	GLY	-	expression tag	UNP I6U7J4
Z	1	MET	-	initiating methionine	UNP I6U7J4
Z	2	GLY	-	expression tag	UNP I6U7J4
a	1	MET	-	initiating methionine	UNP I6U7J4

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Chain	Residue	Modelled	Actual	Comment	Reference
a	2	GLY	-	expression tag	UNP I6U7J4
b	1	MET	-	initiating methionine	UNP I6U7J4
b	2	GLY	-	expression tag	UNP I6U7J4
c	1	MET	-	initiating methionine	UNP I6U7J4
c	2	GLY	-	expression tag	UNP I6U7J4
d	1	MET	-	initiating methionine	UNP I6U7J4
d	2	GLY	-	expression tag	UNP I6U7J4

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	e	2	Total Fe 2 2	0	0
2	A	2	Total Fe 2 2	0	0
2	B	2	Total Fe 2 2	0	0
2	C	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	F	2	Total Fe 2 2	0	0
2	G	2	Total Fe 2 2	0	0
2	L	2	Total Fe 2 2	0	0
2	M	2	Total Fe 2 2	0	0
2	P	2	Total Fe 2 2	0	0
2	Q	2	Total Fe 2 2	0	0
2	T	2	Total Fe 2 2	0	0
2	X	2	Total Fe 2 2	0	0
2	a	2	Total Fe 2 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	e	62	Total O 62 62	0	0
3	A	76	Total O 76 76	0	0
3	B	68	Total O 68 68	0	0
3	C	57	Total O 57 57	0	0
3	D	59	Total O 59 59	0	0
3	E	58	Total O 58 58	0	0
3	F	70	Total O 70 70	0	0
3	G	72	Total O 72 72	0	0
3	H	59	Total O 59 59	0	0
3	I	73	Total O 73 73	0	0
3	J	61	Total O 61 61	0	0
3	K	52	Total O 52 52	0	0
3	L	72	Total O 72 72	0	0
3	M	91	Total O 91 91	0	0
3	N	74	Total O 74 74	0	0
3	O	62	Total O 62 62	0	0
3	P	56	Total O 56 56	0	0
3	Q	59	Total O 59 59	0	0
3	R	53	Total O 53 53	0	0
3	T	92	Total O 92 92	0	0
3	U	69	Total O 69 69	0	0
3	V	80	Total O 80 80	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	W	66	Total O 66 66	0	0
3	X	70	Total O 70 70	0	0
3	Y	71	Total O 71 71	0	0
3	Z	74	Total O 74 74	0	0
3	a	71	Total O 71 71	0	0
3	b	69	Total O 69 69	0	0
3	c	68	Total O 68 68	0	0
3	d	50	Total O 50 50	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: PFC_05175

Chain e:  99%



- Molecule 1: PFC_05175

Chain A:  % 99%



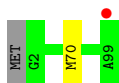
- Molecule 1: PFC_05175

Chain B:  99%



- Molecule 1: PFC_05175

Chain C:  % 98%



- Molecule 1: PFC_05175

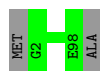
Chain D:  97%



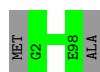
- Molecule 1: PFC_05175



- Molecule 1: PFC_05175



- Molecule 1: PFC_05175



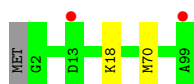
- Molecule 1: PFC_05175



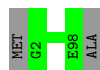
- Molecule 1: PFC_05175



- Molecule 1: PFC_05175

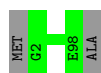


- Molecule 1: PFC_05175



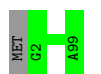
- Molecule 1: PFC_05175

Chain L:  98%



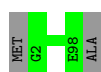
• Molecule 1: PFC_05175

Chain M:  99%



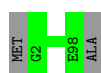
• Molecule 1: PFC_05175

Chain N:  98%



• Molecule 1: PFC_05175

Chain O:  98%



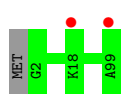
• Molecule 1: PFC_05175

Chain P:  97%



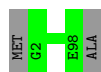
• Molecule 1: PFC_05175

Chain Q:  99% 2%



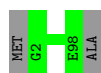
• Molecule 1: PFC_05175

Chain R:  98%



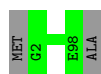
• Molecule 1: PFC_05175

Chain T:  98% .



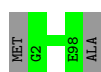
• Molecule 1: PFC_05175

Chain U:  98% .



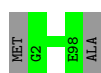
• Molecule 1: PFC_05175

Chain V:  98% .



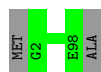
• Molecule 1: PFC_05175

Chain W:  98% .



• Molecule 1: PFC_05175

Chain X:  98% .



• Molecule 1: PFC_05175

Chain Y:  96% ..



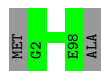
• Molecule 1: PFC_05175

Chain Z:  99% .



• Molecule 1: PFC_05175

Chain a:  98%



• Molecule 1: PFC_05175

Chain b:  99%



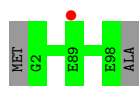
• Molecule 1: PFC_05175

Chain c:  99%



• Molecule 1: PFC_05175

Chain d:  98%



4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	99.85Å 110.06Å 136.27Å 90.00° 91.22° 90.00°	Depositor
Resolution (Å)	47.19 – 2.03 49.71 – 2.03	Depositor EDS
% Data completeness (in resolution range)	97.0 (47.19-2.03) 99.7 (49.71-2.03)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.52 (at 2.03Å)	Xtrriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.174 , 0.202 0.174 , 0.202	Depositor DCC
R_{free} test set	9396 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	22.6	Xtrriage
Anisotropy	0.375	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.000 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	49851	wwPDB-VP
Average B, all atoms (Å ²)	31.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 65.55 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.9620e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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.

5 Model quality

5.1 Standard geometry

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	0.25	0/800	0.39	0/1078
1	B	0.25	0/799	0.39	0/1076
1	C	0.24	0/802	0.39	0/1080
1	D	0.25	0/803	0.39	0/1081
1	E	0.25	0/811	0.41	0/1092
1	F	0.25	0/795	0.40	0/1071
1	G	0.25	0/794	0.40	0/1069
1	H	0.24	0/795	0.39	0/1071
1	I	0.24	0/800	0.39	0/1078
1	J	0.24	0/813	0.39	0/1094
1	K	0.25	0/814	0.40	0/1095
1	L	0.24	0/822	0.39	0/1105
1	M	0.25	0/808	0.39	0/1088
1	N	0.24	0/803	0.39	0/1081
1	O	0.24	0/806	0.38	0/1085
1	P	0.24	0/800	0.38	0/1077
1	Q	0.24	0/799	0.38	0/1076
1	R	0.24	0/794	0.38	0/1069
1	T	0.24	0/811	0.38	0/1091
1	U	0.24	0/794	0.39	0/1069
1	V	0.25	0/801	0.39	0/1079
1	W	0.25	0/803	0.39	0/1081
1	X	0.24	0/795	0.39	0/1071
1	Y	0.24	0/806	0.40	0/1085
1	Z	0.24	0/800	0.38	0/1078
1	a	0.25	0/795	0.40	0/1071
1	b	0.24	0/791	0.38	0/1066
1	c	0.24	0/821	0.39	0/1104
1	d	0.24	0/786	0.37	0/1059
1	e	0.24	0/799	0.39	0/1076
All	All	0.24	0/24060	0.39	0/32396

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

Due to software issues we are unable to calculate clashes - this section is therefore empty.

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	97/99 (98%)	97 (100%)	0	0	100	100
1	B	97/99 (98%)	97 (100%)	0	0	100	100
1	C	97/99 (98%)	97 (100%)	0	0	100	100
1	D	97/99 (98%)	97 (100%)	0	0	100	100
1	E	98/99 (99%)	98 (100%)	0	0	100	100
1	F	96/99 (97%)	96 (100%)	0	0	100	100
1	G	96/99 (97%)	96 (100%)	0	0	100	100
1	H	96/99 (97%)	96 (100%)	0	0	100	100
1	I	97/99 (98%)	97 (100%)	0	0	100	100
1	J	98/99 (99%)	98 (100%)	0	0	100	100
1	K	98/99 (99%)	98 (100%)	0	0	100	100
1	L	99/99 (100%)	99 (100%)	0	0	100	100
1	M	98/99 (99%)	98 (100%)	0	0	100	100
1	N	97/99 (98%)	97 (100%)	0	0	100	100
1	O	97/99 (98%)	97 (100%)	0	0	100	100
1	P	97/99 (98%)	96 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	Q	97/99 (98%)	97 (100%)	0	0	100	100
1	R	96/99 (97%)	96 (100%)	0	0	100	100
1	T	98/99 (99%)	98 (100%)	0	0	100	100
1	U	96/99 (97%)	96 (100%)	0	0	100	100
1	V	97/99 (98%)	97 (100%)	0	0	100	100
1	W	97/99 (98%)	97 (100%)	0	0	100	100
1	X	96/99 (97%)	95 (99%)	1 (1%)	0	100	100
1	Y	97/99 (98%)	97 (100%)	0	0	100	100
1	Z	97/99 (98%)	97 (100%)	0	0	100	100
1	a	96/99 (97%)	96 (100%)	0	0	100	100
1	b	96/99 (97%)	96 (100%)	0	0	100	100
1	c	99/99 (100%)	98 (99%)	1 (1%)	0	100	100
1	d	95/99 (96%)	95 (100%)	0	0	100	100
1	e	97/99 (98%)	97 (100%)	0	0	100	100
All	All	2909/2970 (98%)	2906 (100%)	3 (0%)	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	86/86 (100%)	86 (100%)	0	100	100
1	B	86/86 (100%)	86 (100%)	0	100	100
1	C	86/86 (100%)	85 (99%)	1 (1%)	71	75
1	D	87/86 (101%)	86 (99%)	1 (1%)	73	77
1	E	87/86 (101%)	86 (99%)	1 (1%)	73	77
1	F	86/86 (100%)	86 (100%)	0	100	100
1	G	86/86 (100%)	86 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	86/86 (100%)	85 (99%)	1 (1%)	71	75
1	I	86/86 (100%)	86 (100%)	0	100	100
1	J	87/86 (101%)	85 (98%)	2 (2%)	50	51
1	K	88/86 (102%)	88 (100%)	0	100	100
1	L	89/86 (104%)	89 (100%)	0	100	100
1	M	87/86 (101%)	87 (100%)	0	100	100
1	N	87/86 (101%)	87 (100%)	0	100	100
1	O	87/86 (101%)	87 (100%)	0	100	100
1	P	87/86 (101%)	86 (99%)	1 (1%)	73	77
1	Q	86/86 (100%)	86 (100%)	0	100	100
1	R	86/86 (100%)	86 (100%)	0	100	100
1	T	88/86 (102%)	88 (100%)	0	100	100
1	U	86/86 (100%)	86 (100%)	0	100	100
1	V	87/86 (101%)	87 (100%)	0	100	100
1	W	87/86 (101%)	87 (100%)	0	100	100
1	X	86/86 (100%)	86 (100%)	0	100	100
1	Y	87/86 (101%)	85 (98%)	2 (2%)	50	51
1	Z	86/86 (100%)	86 (100%)	0	100	100
1	a	86/86 (100%)	86 (100%)	0	100	100
1	b	85/86 (99%)	85 (100%)	0	100	100
1	c	88/86 (102%)	88 (100%)	0	100	100
1	d	85/86 (99%)	85 (100%)	0	100	100
1	e	86/86 (100%)	86 (100%)	0	100	100
All	All	2597/2580 (101%)	2588 (100%)	9 (0%)	91	94

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	Y	47	GLU
1	Y	70	MET
1	H	70	MET
1	J	18	LYS
1	J	70	MET

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

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 30 ligands modelled in this entry, 30 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, 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	98/99 (98%)	-0.40	1 (1%) 82 82	18, 27, 46, 62	0
1	B	98/99 (98%)	-0.47	0 100 100	15, 25, 44, 62	0
1	C	98/99 (98%)	-0.42	1 (1%) 82 82	16, 30, 49, 63	0
1	D	97/99 (97%)	-0.29	0 100 100	19, 27, 48, 58	0
1	E	98/99 (98%)	-0.50	1 (1%) 82 82	14, 26, 45, 59	0
1	F	97/99 (97%)	-0.39	0 100 100	16, 25, 44, 56	0
1	G	97/99 (97%)	-0.36	0 100 100	14, 23, 41, 53	0
1	H	97/99 (97%)	-0.47	0 100 100	18, 26, 46, 63	0
1	I	98/99 (98%)	-0.29	1 (1%) 82 82	15, 25, 45, 62	0
1	J	98/99 (98%)	-0.40	2 (2%) 65 64	17, 27, 46, 63	0
1	K	97/99 (97%)	-0.39	0 100 100	17, 26, 48, 59	0
1	L	97/99 (97%)	-0.40	0 100 100	14, 23, 44, 52	0
1	M	98/99 (98%)	-0.46	0 100 100	15, 25, 47, 58	0
1	N	97/99 (97%)	-0.44	0 100 100	16, 26, 44, 61	0
1	O	97/99 (97%)	-0.41	0 100 100	18, 25, 44, 59	0
1	P	97/99 (97%)	-0.36	0 100 100	19, 29, 51, 76	0
1	Q	98/99 (98%)	-0.33	2 (2%) 65 64	19, 30, 52, 70	0
1	R	97/99 (97%)	-0.42	0 100 100	19, 30, 47, 58	0
1	T	97/99 (97%)	-0.42	0 100 100	14, 23, 42, 56	0
1	U	97/99 (97%)	-0.41	0 100 100	16, 25, 44, 60	0
1	V	97/99 (97%)	-0.42	0 100 100	14, 23, 40, 57	0
1	W	97/99 (97%)	-0.45	0 100 100	16, 24, 44, 62	0
1	X	97/99 (97%)	-0.47	0 100 100	14, 23, 43, 52	0
1	Y	97/99 (97%)	-0.49	0 100 100	15, 24, 43, 58	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	Z	98/99 (98%)	-0.46	0 100 100	17, 27, 48, 62	0
1	a	97/99 (97%)	-0.45	0 100 100	17, 26, 43, 61	0
1	b	98/99 (98%)	-0.47	0 100 100	16, 26, 49, 66	0
1	c	98/99 (98%)	-0.31	1 (1%) 82 82	15, 24, 49, 67	0
1	d	97/99 (97%)	-0.33	1 (1%) 82 82	18, 29, 48, 63	0
1	e	98/99 (98%)	-0.36	0 100 100	17, 28, 48, 62	0
All	All	2922/2970 (98%)	-0.41	10 (0%) 94 93	14, 26, 47, 76	0

The worst 5 of 10 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	99	ALA	6.7
1	c	99	ALA	6.5
1	Q	99	ALA	3.0
1	J	99	ALA	2.9
1	A	99	ALA	2.8

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, 95th 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 ²)	Q<0.9
2	FE	D	101	1/1	0.91	0.32	19,19,19,19	0
2	FE	P	101	1/1	0.92	0.36	20,20,20,20	1
2	FE	D	102	1/1	0.93	0.29	20,20,20,20	0
2	FE	G	102	1/1	0.93	0.29	15,15,15,15	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	FE	A	101	1/1	0.93	0.38	17,17,17,17	0
2	FE	e	101	1/1	0.94	0.30	19,19,19,19	1
2	FE	P	102	1/1	0.94	0.33	19,19,19,19	0
2	FE	a	102	1/1	0.94	0.32	16,16,16,16	0
2	FE	B	102	1/1	0.95	0.31	20,20,20,20	0
2	FE	L	102	1/1	0.95	0.26	21,21,21,21	0
2	FE	C	101	1/1	0.95	0.32	19,19,19,19	0
2	FE	e	102	1/1	0.95	0.34	16,16,16,16	0
2	FE	Q	102	1/1	0.95	0.30	20,20,20,20	0
2	FE	X	102	1/1	0.95	0.29	18,18,18,18	0
2	FE	a	101	1/1	0.95	0.32	18,18,18,18	0
2	FE	B	101	1/1	0.95	0.30	21,21,21,21	0
2	FE	G	101	1/1	0.96	0.31	14,14,14,14	0
2	FE	Q	101	1/1	0.96	0.28	27,27,27,27	0
2	FE	A	102	1/1	0.96	0.32	18,18,18,18	0
2	FE	E	101	1/1	0.96	0.30	23,23,23,23	0
2	FE	M	101	1/1	0.96	0.30	20,20,20,20	0
2	FE	F	102	1/1	0.96	0.30	16,16,16,16	0
2	FE	C	102	1/1	0.97	0.28	23,23,23,23	0
2	FE	E	102	1/1	0.97	0.29	20,20,20,20	0
2	FE	T	101	1/1	0.97	0.27	21,21,21,21	0
2	FE	T	102	1/1	0.97	0.25	19,19,19,19	0
2	FE	X	101	1/1	0.97	0.29	24,24,24,24	0
2	FE	M	102	1/1	0.97	0.27	22,22,22,22	0
2	FE	F	101	1/1	0.97	0.35	14,14,14,14	0
2	FE	L	101	1/1	0.97	0.24	20,20,20,20	0

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

There are no such residues in this entry.