



# wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 20, 2025 – 11:58 AM EDT

PDB ID : 4IWK  
Title : Crystal structure of iron soaked (overnight) ferritin from Pseudo-nitzschia multiseria  
Authors : Pfaffen, S.; Murphy, M.E.P.  
Deposited on : 2013-01-23  
Resolution : 1.65 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.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

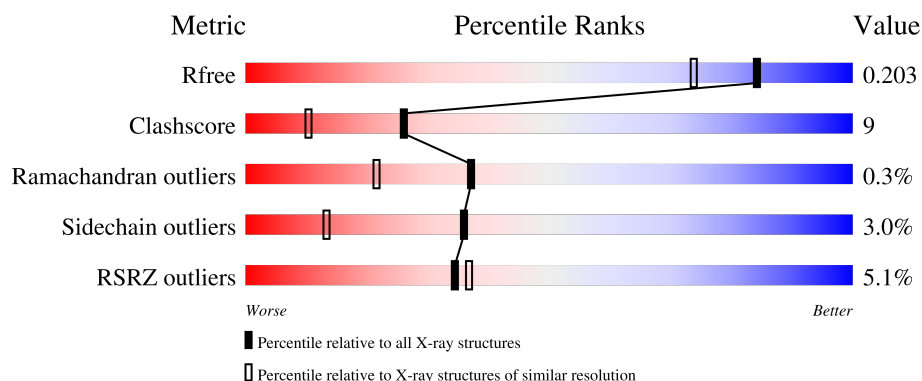
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.65 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))
$R_{free}$	164625	2328 (1.66-1.66)
Clashscore	180529	2515 (1.66-1.66)
Ramachandran outliers	177936	2475 (1.66-1.66)
Sidechain outliers	177891	2475 (1.66-1.66)
RSRZ outliers	164620	2328 (1.66-1.66)

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	168	<div> <div>4%</div> <div> <div></div> <div>76%</div> <div>15%</div> <div>• • 5%</div> </div> </div>
1	B	168	<div> <div>4%</div> <div> <div></div> <div>77%</div> <div>13%</div> <div>• 7%</div> </div> </div>
1	C	168	<div> <div>4%</div> <div> <div></div> <div>80%</div> <div>13%</div> <div>• • 6%</div> </div> </div>
1	D	168	<div> <div>5%</div> <div> <div></div> <div>80%</div> <div>11%</div> <div>• 5%</div> </div> </div>
1	E	168	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>16%</div> <div>• • 6%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
1	F	168	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
1	YCM	A	77[A]	-	-	X	-
1	YCM	B	77[C]	-	-	X	-
1	YCM	F	77[A]	-	X	-	-

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8599 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	A	159	Total	C	N	O	S	0	8	0
			1303	826	213	258	6			
1	B	157	Total	C	N	O	S	0	6	0
			1272	803	209	254	6			
1	C	158	Total	C	N	O	S	0	4	0
			1271	800	212	254	5			
1	D	159	Total	C	N	O	S	0	7	0
			1299	818	215	260	6			
1	E	158	Total	C	N	O	S	0	5	0
			1276	804	209	258	5			
1	F	159	Total	C	N	O	S	0	5	0
			1282	809	210	257	6			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP B6DMH6
B	0	GLY	-	expression tag	UNP B6DMH6
C	0	GLY	-	expression tag	UNP B6DMH6
D	0	GLY	-	expression tag	UNP B6DMH6
E	0	GLY	-	expression tag	UNP B6DMH6
F	0	GLY	-	expression tag	UNP B6DMH6

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	4	Total	Fe	0	0
			4	4		
2	B	3	Total	Fe	0	0
			3	3		
2	C	4	Total	Fe	0	0
			4	4		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	5	Total 5	Fe 5	0	0
2	E	4	Total 4	Fe 4	0	0
2	F	4	Total 4	Fe 4	0	0

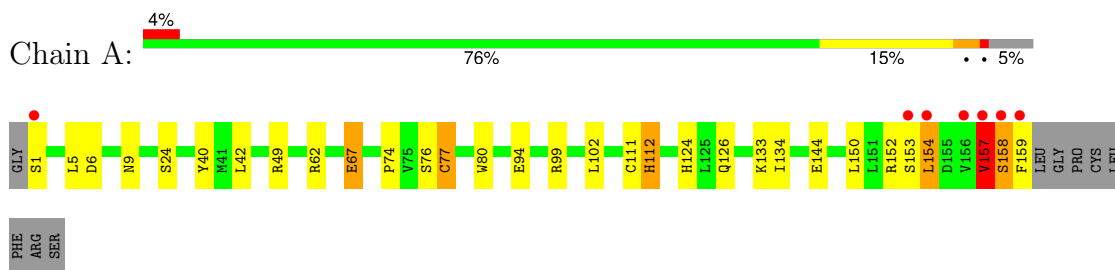
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	133	Total 133	O 133	0	0
3	B	150	Total 150	O 150	0	0
3	C	154	Total 154	O 154	0	0
3	D	150	Total 150	O 150	0	0
3	E	155	Total 155	O 155	0	0
3	F	130	Total 130	O 130	0	0

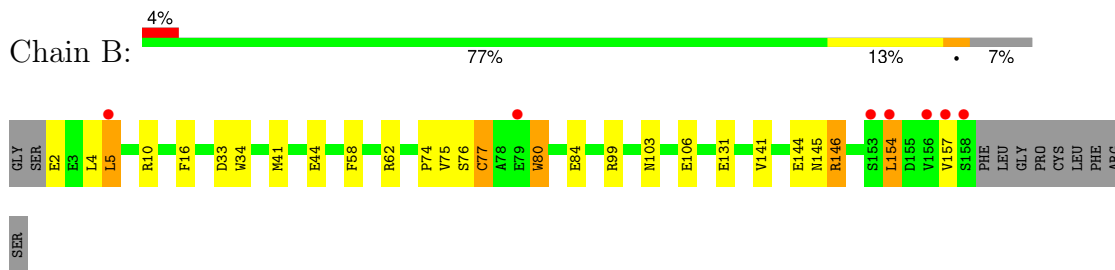
### 3 Residue-property plots [i](#)

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.

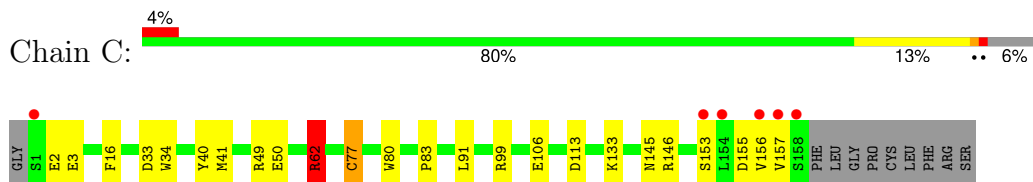
- Molecule 1: Ferritin



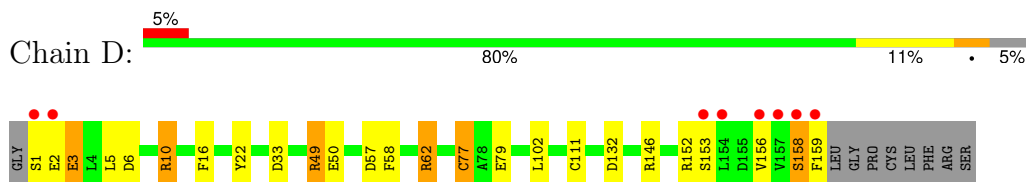
- Molecule 1: Ferritin



- Molecule 1: Ferritin



- Molecule 1: Ferritin

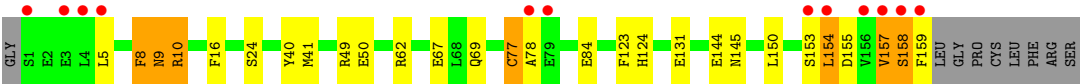
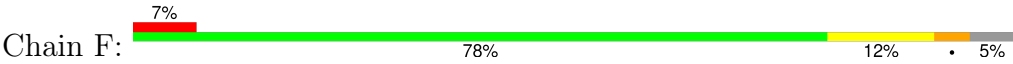


- Molecule 1: Ferritin





● Molecule 1: Ferritin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	126.25Å 126.25Å 170.30Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.95 – 1.65 47.95 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.2 (47.95-1.65) 99.2 (47.95-1.65)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.08 (at 1.65Å)	Xtriage
Refinement program	REFMAC	Depositor
R, $R_{free}$	0.176 , 0.205 0.174 , 0.203	Depositor DCC
$R_{free}$ test set	8205 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	22.4	Xtriage
Anisotropy	0.029	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 38.4	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	8599	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.72% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: YCM, 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	1.18	0/1330	1.17	6/1809 (0.3%)
1	B	1.32	3/1295 (0.2%)	1.32	11/1761 (0.6%)
1	C	1.32	4/1298 (0.3%)	1.50	11/1764 (0.6%)
1	D	1.35	3/1323 (0.2%)	1.43	14/1798 (0.8%)
1	E	1.42	3/1303 (0.2%)	1.42	15/1772 (0.8%)
1	F	1.32	4/1307 (0.3%)	1.26	9/1777 (0.5%)
All	All	1.32	17/7856 (0.2%)	1.35	66/10681 (0.6%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	1

The worst 5 of 17 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	62	ARG	CZ-NH1	7.66	1.43	1.33
1	F	9	ASN	N-CA	-6.08	1.34	1.46
1	C	16	PHE	CG-CD2	5.90	1.47	1.38
1	F	69	GLN	CG-CD	5.58	1.63	1.51
1	B	80	TRP	CZ3-CH2	5.57	1.49	1.40

The worst 5 of 66 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	62	ARG	NE-CZ-NH2	-27.64	106.48	120.30
1	C	62	ARG	NE-CZ-NH1	21.06	130.83	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	62	ARG	NE-CZ-NH2	-16.51	112.05	120.30
1	D	49	ARG	NE-CZ-NH1	-11.40	114.60	120.30
1	A	62	ARG	NE-CZ-NH1	-10.20	115.20	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	1	SER	Peptide

## 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	1303	0	1260	38	0
1	B	1272	0	1225	31	0
1	C	1271	0	1222	17	0
1	D	1299	0	1243	10	0
1	E	1276	0	1221	22	0
1	F	1282	0	1226	22	1
2	A	4	0	0	0	0
2	B	3	0	0	1	0
2	C	4	0	0	1	0
2	D	5	0	0	0	0
2	E	4	0	0	1	0
2	F	4	0	0	0	0
3	A	133	0	0	7	0
3	B	150	0	0	11	0
3	C	154	0	0	7	0
3	D	150	0	0	3	0
3	E	155	0	0	10	0
3	F	130	0	0	8	0
All	All	8599	0	7397	141	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 141 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:74:PRO:HB2	1:A:77[A]:YCM:HD2	1.29	1.15
1:E:2:GLU:HA	1:E:2:GLU:OE1	1.47	1.15
1:E:77:YCM:HB3	3:E:447:HOH:O	1.51	1.10
1:E:157:VAL:HG23	1:E:158:SER:N	1.79	0.92
1:A:74:PRO:CB	1:A:77[A]:YCM:HD2	1.99	0.92

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)
1:F:10:ARG:NH1	1:F:10:ARG:NH1[7_554]	1.89	0.31

## 5.3 Torsion angles ⓘ

### 5.3.1 Protein backbone ⓘ

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	163/168 (97%)	161 (99%)	1 (1%)	1 (1%)	22	8
1	B	159/168 (95%)	157 (99%)	2 (1%)	0	100	100
1	C	159/168 (95%)	158 (99%)	1 (1%)	0	100	100
1	D	162/168 (96%)	161 (99%)	1 (1%)	0	100	100
1	E	160/168 (95%)	155 (97%)	4 (2%)	1 (1%)	22	8
1	F	160/168 (95%)	155 (97%)	4 (2%)	1 (1%)	22	8
All	All	963/1008 (96%)	947 (98%)	13 (1%)	3 (0%)	37	21

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	157	VAL
1	F	78	ALA

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Mol	Chain	Res	Type
1	E	157	VAL

### 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	143/143 (100%)	138 (96%)	5 (4%)	31	9
1	B	139/143 (97%)	137 (99%)	2 (1%)	62	43
1	C	139/143 (97%)	135 (97%)	4 (3%)	37	14
1	D	142/143 (99%)	139 (98%)	3 (2%)	48	26
1	E	140/143 (98%)	134 (96%)	6 (4%)	25	6
1	F	140/143 (98%)	136 (97%)	4 (3%)	37	14
All	All	843/858 (98%)	819 (97%)	24 (3%)	36	15

5 of 24 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	2	GLU
1	E	79	GLU
1	E	62	ARG
1	E	153	SER
1	B	154	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19 such sidechains are listed below:

Mol	Chain	Res	Type
1	E	145	ASN
1	F	124	HIS
1	F	145	ASN
1	F	121	ASN
1	C	145	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

10 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
1	YCM	E	77	1	7,9,10	1.52	1 (14%)	5,10,12	0.53	0
1	YCM	B	77[C]	-	7,9,10	1.37	1 (14%)	5,10,12	1.69	1 (20%)
1	YCM	D	77[B]	1	7,9,10	2.21	3 (42%)	5,10,12	0.88	0
1	YCM	A	77[B]	1	7,9,10	2.07	2 (28%)	5,10,12	0.66	0
1	YCM	B	77[B]	-	7,9,10	2.57	4 (57%)	5,10,12	1.20	0
1	YCM	A	77[A]	1	7,9,10	2.40	2 (28%)	5,10,12	3.60	1 (20%)
1	YCM	D	77[A]	1	7,9,10	1.86	2 (28%)	5,10,12	2.77	1 (20%)
1	YCM	C	77	1	7,9,10	2.48	3 (42%)	5,10,12	0.66	0
1	YCM	F	77[B]	-	7,9,10	3.99	4 (57%)	5,10,12	2.22	2 (40%)
1	YCM	F	77[A]	-	7,5,10	35.12	2 (28%)	5,5,12	5.02	2 (40%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	YCM	E	77	1	-	3/6/8/10	-
1	YCM	B	77[C]	-	-	3/6/8/10	-
1	YCM	D	77[B]	1	-	1/6/8/10	-
1	YCM	A	77[B]	1	-	2/6/8/10	-
1	YCM	B	77[B]	-	-	1/6/8/10	-
1	YCM	A	77[A]	1	-	4/6/8/10	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	YCM	D	77[A]	1	-	1/6/8/10	-
1	YCM	C	77	1	-	0/6/8/10	-
1	YCM	F	77[B]	-	-	0/6/8/10	-
1	YCM	F	77[A]	-	-	3/6/4/10	-

The worst 5 of 24 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	F	77[A]	YCM	CD-SG	92.77	4.13	1.81
1	F	77[B]	YCM	CD-SG	7.05	1.99	1.81
1	A	77[A]	YCM	CD-SG	5.30	1.94	1.81
1	F	77[B]	YCM	CB-SG	5.13	2.00	1.81
1	C	77	YCM	CD-SG	5.08	1.94	1.81

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	77[A]	YCM	CE-CD-SG	-9.50	83.82	113.81
1	A	77[A]	YCM	CE-CD-SG	7.92	138.83	113.81
1	D	77[A]	YCM	CE-CD-SG	-5.63	96.05	113.81
1	F	77[A]	YCM	CB-SG-CD	-5.62	54.09	104.24
1	F	77[B]	YCM	CA-CB-SG	4.22	129.09	113.51

There are no chirality outliers.

5 of 18 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	77[A]	YCM	C-CA-CB-SG
1	A	77[A]	YCM	CE-CD-SG-CB
1	B	77[C]	YCM	N-CA-CB-SG
1	B	77[C]	YCM	C-CA-CB-SG
1	B	77[C]	YCM	CE-CD-SG-CB

There are no ring outliers.

9 monomers are involved in 37 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	E	77	YCM	2	0
1	B	77[C]	YCM	10	0
1	A	77[B]	YCM	4	0
1	B	77[B]	YCM	4	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	77[A]	YCM	8	0
1	D	77[A]	YCM	1	0
1	C	77	YCM	3	0
1	F	77[B]	YCM	4	0
1	F	77[A]	YCM	1	0

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 24 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	158/168 (94%)	-0.15	7 (4%) 39 43	12, 22, 46, 66	7 (4%)
1	B	156/168 (92%)	-0.12	7 (4%) 39 42	11, 22, 42, 69	5 (3%)
1	C	157/168 (93%)	-0.20	6 (3%) 44 47	13, 21, 39, 64	4 (2%)
1	D	158/168 (94%)	-0.15	8 (5%) 34 37	9, 21, 45, 70	6 (3%)
1	E	157/168 (93%)	-0.17	8 (5%) 34 37	10, 20, 43, 66	5 (3%)
1	F	158/168 (94%)	-0.03	12 (7%) 21 23	12, 22, 47, 70	4 (2%)
All	All	944/1008 (93%)	-0.14	48 (5%) 34 37	9, 22, 45, 70	31 (3%)

The worst 5 of 48 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	157	VAL	7.0
1	C	156	VAL	6.4
1	A	156	VAL	5.5
1	C	157	VAL	5.5
1	D	157	VAL	5.0

### 6.2 Non-standard residues in protein, DNA, RNA chains [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
1	YCM	F	77[A]	6/11	0.72	0.19	33,35,43,49	3
1	YCM	F	77[B]	10/11	0.72	0.19	20,33,43,49	7
1	YCM	A	77[A]	10/11	0.86	0.17	22,30,33,35	10
1	YCM	A	77[B]	10/11	0.86	0.17	18,22,29,30	10

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
1	YCM	D	77[A]	10/11	0.87	0.16	19,25,31,34	10
1	YCM	D	77[B]	10/11	0.87	0.16	15,17,24,24	10
1	YCM	C	77	10/11	0.91	0.12	24,27,31,32	0
1	YCM	B	77[C]	10/11	0.92	0.11	20,26,32,39	7
1	YCM	B	77[B]	10/11	0.92	0.11	21,22,31,32	7
1	YCM	E	77	10/11	0.94	0.17	22,27,52,56	0

## 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( $\text{\AA}^2$ )	Q<0.9
2	FE	E	204	1/1	0.82	0.12	38,38,38,38	1
2	FE	D	203	1/1	0.85	0.14	40,40,40,40	1
2	FE	A	202	1/1	0.85	0.13	40,40,40,40	1
2	FE	F	204	1/1	0.89	0.12	35,35,35,35	1
2	FE	C	204	1/1	0.93	0.09	36,36,36,36	1
2	FE	C	203	1/1	0.93	0.09	38,38,38,38	1
2	FE	B	201	1/1	0.94	0.09	38,38,38,38	1
2	FE	D	202	1/1	0.95	0.12	35,35,35,35	1
2	FE	D	205	1/1	0.96	0.09	34,34,34,34	1
2	FE	E	203	1/1	0.97	0.09	35,35,35,35	1
2	FE	B	202	1/1	0.98	0.04	22,22,22,22	1
2	FE	A	204	1/1	0.98	0.07	39,39,39,39	1
2	FE	F	203	1/1	0.98	0.12	32,32,32,32	1
2	FE	E	201	1/1	0.98	0.04	22,22,22,22	1
2	FE	D	204	1/1	0.99	0.08	21,21,21,21	1
2	FE	C	202	1/1	0.99	0.06	21,21,21,21	1
2	FE	A	203	1/1	0.99	0.07	23,23,23,23	1
2	FE	E	202	1/1	0.99	0.06	20,20,20,20	1
2	FE	A	201	1/1	0.99	0.05	24,24,24,24	0
2	FE	D	201	1/1	0.99	0.04	21,21,21,21	0
2	FE	F	201	1/1	0.99	0.05	23,23,23,23	0
2	FE	F	202	1/1	0.99	0.06	22,22,22,22	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FE	B	203	1/1	0.99	0.05	21,21,21,21	1
2	FE	C	201	1/1	0.99	0.03	22,22,22,22	1

## 6.5 Other polymers [i](#)

There are no such residues in this entry.