

wwPDB X-ray Structure Validation Summary Report (i)

Nov 14, 2023 – 05:51 PM JST

PDB ID	:	5ZUR
Title	:	Achromobacter Dh1f Bacterioferritin
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Deposited on	:	2018-05-08
Resolution	:	1.95 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 (i) 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 (1)) were used in the production of this report:

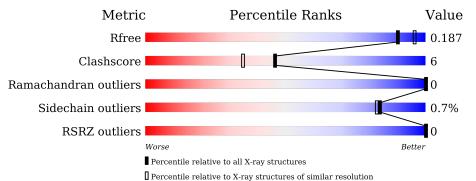
MolProbity Mogul Xtriage (Phenix) EDS	:	4.02b-467 1.8.5 (274361), CSD as541be (2020) 1.13 2.36
buster-report Percentile statistics Refmac	: : :	1.1.7 (2018) 20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		Parkinson et al. (1996) 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: $X\text{-}RAY \, DIFFRACTION$

The reported resolution of this entry is 1.95 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R _{free}	130704	2580 (1.96-1.96)
Clashscore	141614	2705 (1.96-1.96)
Ramachandran outliers	138981	2678 (1.96-1.96)
Sidechain outliers	138945	2678 (1.96-1.96)
RSRZ outliers	127900	2539 (1.96-1.96)

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 <=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	А	158	91%	6%	•••
1	В	158	89%	8%	•••
1	С	158	89%	8%	•



2 Entry composition (i)

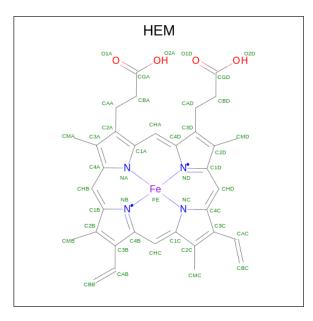
There are 7 unique types of molecules in this entry. The entry contains 4593 atoms, of which 0 are hydrogens and 0 are deuteriums.

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.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Δ	154	Total	С	Ν	0	S	0	1	0
1	Л	104	1250	792	211	240	7	0		
1	р	155	Total	С	Ν	0	S	0	1	0
	D	155	1260	798	216	239	7	0		
1	С	154	Total	С	Ν	0	S	0	0	0
		154	1250	791	211	241	7	0	0	0

• Molecule 1 is a protein called Bacterioferritin.

• Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



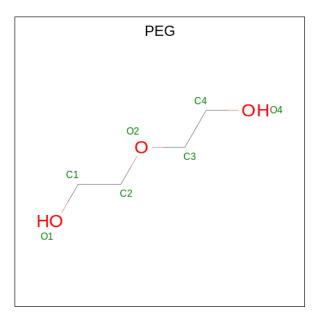
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	Δ	1	Total	С	Fe	Ν	Ο	0	0
2	Π		43	34	1	4	4	0	
2	В	1	Total	С	Fe	Ν	Ο	0	0
	D	1	43	34	1	4	4	0	
2	С	1	Total	С	Fe	Ν	Ο	0	0
			43	34	1	4	4	0	0



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	1	Total Fe 1 1	0	0
3	В	1	Total Fe 1 1	0	0
3	С	1	Total Fe 1 1	0	0

• Molecule 3 is FE (II) ION (three-letter code: FE2) (formula: Fe).

• Molecule 4 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 7 4 3 \end{array}$	0	0
4	А	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 7 & 4 & 3 \end{array}$	0	0
4	В	1	$\begin{array}{ccc} \text{Total} & \text{C} & \text{O} \\ 7 & 4 & 3 \end{array}$	0	0
4	С	1	$\begin{array}{ccc} \text{Total} \text{C} \text{O} \\ 7 4 3 \end{array}$	0	0

• Molecule 5 is BARIUM ION (three-letter code: BA) (formula: Ba).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	А	1	Total Ba 1 1	0	0
5	В	1	Total Ba 1 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	С	1	Total Ba 1 1	0	0

• Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	1	Total Cl 1 1	0	0
6	В	1	Total Cl 1 1	0	0
6	С	1	Total Cl 1 1	0	0

• Molecule 7 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	А	230	Total O 230 230	0	0
7	В	232	Total O 232 232	0	0
7	С	205	Total O 205 205	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.

Chain A:	91%	6%	••
M1 K2 K5 K5 K3 K30 B51 M52 M52 M52 M52 M52 M52 M52 M52 M52 M52			
• Molecule 1: Bacterioferritin			
Chain B:	89%	8%	•••
M1 K2 K3 K39 K39 E44 E51 E65 E85 E85 E85 E85 E85 E85 E85 E85 K92 K92 K92 K92 K92 K92 K92 K92 K92 K92	8155 VAL PRO GLU GLU		
• Molecule 1: Bacterioferritin			
Chain C:	89%	8%	·
M1 K6 K6 K39 E44 E51 M62 E51 M64 M64 M64 M64 M64 E94 M154 S17 S17 S17 S17 S17 S17 S17 S17 S17 S17			

• Molecule 1: Bacterioferritin



4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 4 2 2	Depositor
Cell constants	158.66Å 158.66Å 118.06Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.70 - 1.95	Depositor
Resolution (A)	40.66 - 1.95	EDS
% Data completeness	99.9 (40.70-1.95)	Depositor
(in resolution range)	99.9 (40.66 - 1.95)	EDS
R _{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.93 (at 1.95 Å)	Xtriage
Refinement program	REFMAC 5.8.0222	Depositor
D D.	0.142 , 0.187	Depositor
R, R_{free}	0.142 , 0.187	DCC
R_{free} test set	2714 reflections $(4.97%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	15.7	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35 , 57.5	EDS
L-test for twinning ²	$ \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	0.96	EDS
Total number of atoms	4593	wwPDB-VP
Average B, all atoms $(Å^2)$	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 4.80% of the height of the origin peak. No significant pseudotranslation is detected.

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.



¹Intensities estimated from amplitudes.

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: PEG, CL, FE2, HEM, BA

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).

T	Mol	Chain	Bond lengths		Bond angles	
	VIOI	Unam	RMSZ	# Z > 5	RMSZ	# Z > 5
	1	А	0.80	0/1274	0.77	0/1720
	1	В	0.79	0/1284	0.82	0/1732
	1	С	0.80	0/1271	0.85	3/1716~(0.2%)
	All	All	0.80	0/3829	0.81	3/5168~(0.1%)

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	В	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
1	С	102	ARG	NE-CZ-NH1	9.59	125.09	120.30
1	С	102	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	С	102	ARG	CD-NE-CZ	5.56	131.38	123.60

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	В	102	ARG	Sidechain



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	А	1250	0	1221	10	0
1	В	1260	0	1236	16	0
1	С	1250	0	1218	11	0
2	А	43	0	30	4	0
2	В	43	0	30	6	0
2	С	43	0	30	1	0
3	А	1	0	0	0	0
3	В	1	0	0	0	0
3	С	1	0	0	0	0
4	А	14	0	20	1	0
4	В	7	0	10	3	0
4	С	7	0	10	0	0
5	А	1	0	0	0	0
5	В	1	0	0	0	0
5	С	1	0	0	0	0
6	А	1	0	0	0	0
6	В	1	0	0	0	0
6	С	1	0	0	0	0
7	А	230	0	0	3	1
7	В	232	0	0	8	3
7	С	205	0	0	6	4
All	All	4593	0	3805	49	4

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 6.

The worst 5 of 49 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)
2:B:202:HEM:HBC2	2:B:202:HEM:HHD	1.31	1.12
1:B:39:LYS:NZ	1:B:155:SER:HA	1.68	1.08
2:B:202:HEM:HHC	2:B:202:HEM:CBB	1.86	1.04
2:B:202:HEM:HBB2	2:B:202:HEM:CHC	1.78	1.01
2:B:202:HEM:HHC	2:B:202:HEM:HBB2	0.94	0.93



Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:B:418:HOH:O	7:C:302:HOH:O[4_565]	2.01	0.19
7:B:484:HOH:O	7:C:469:HOH:O[4_565]	2.07	0.13
7:A:507:HOH:O	7:C:455:HOH:O[7_555]	2.17	0.03
7:B:418:HOH:O	7:C:454:HOH:O[4_565]	2.17	0.03

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

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	А	153/158~(97%)	151 (99%)	2(1%)	0	100 100
1	В	154/158~(98%)	152 (99%)	2(1%)	0	100 100
1	С	152/158~(96%)	151 (99%)	1 (1%)	0	100 100
All	All	459/474~(97%)	454 (99%)	5 (1%)	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 side chain 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	А	134/141~(95%)	133~(99%)	1 (1%)	84 82
1	В	135/141~(96%)	134 (99%)	1 (1%)	84 82
1	\mathbf{C}	134/141~(95%)	133~(99%)	1 (1%)	84 82

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	403/423~(95%)	400 (99%)	3(1%)	84 82

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	2	LYS
1	В	2	LYS
1	С	64	MET

Sometimes side chains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such side chains are listed below:

Mol	Chain	Res	Type
1	В	107	HIS

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 16 ligands modelled in this entry, 9 are monoatomic - leaving 7 for Mogul analysis.

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	ol Type Chain H		Res	Link	Bo	Bond lengths			Bond angles		
10101	туре	Unam	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
2	HEM	В	202	1,2	$41,\!50,\!50$	1.25	5 (12%)	45,82,82	2.24	15 (33%)	
4	PEG	А	203	-	$6,\!6,\!6$	0.75	0	$5,\!5,\!5$	0.54	0	
2	HEM	С	202	1,2	$41,\!50,\!50$	1.28	6 (14%)	45,82,82	2.04	12 (26%)	
4	PEG	А	204	-	$6,\!6,\!6$	0.56	0	$5,\!5,\!5$	0.16	0	
4	PEG	С	201	-	$6,\!6,\!6$	1.01	0	$5,\!5,\!5$	0.88	0	
2	HEM	А	201	1	41,50,50	1.35	7 (17%)	45,82,82	2.03	11 (24%)	
4	PEG	В	203	-	$6,\!6,\!6$	0.38	0	$5,\!5,\!5$	0.77	0	

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
2	HEM	В	202	1,2	-	5/12/54/54	-
4	PEG	А	203	-	-	1/4/4/4	-
2	HEM	С	202	1,2	-	4/12/54/54	-
4	PEG	А	204	-	-	3/4/4/4	-
4	PEG	С	201	-	-	1/4/4/4	-
2	HEM	А	201	1	-	7/12/54/54	-
4	PEG	В	203	-	-	3/4/4/4	-

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	201	HEM	CHA-C4D	3.01	1.42	1.35
2	А	201	HEM	FE-NB	2.75	2.10	1.96
2	С	202	HEM	FE-NB	2.74	2.10	1.96
2	В	202	HEM	CHA-C4D	2.71	1.41	1.35
2	В	202	HEM	FE-NB	2.70	2.10	1.96

The worst 5 of 38 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	$Observed(^{o})$	$Ideal(^{o})$
2	В	202	HEM	CHC-C4B-NB	6.56	131.56	124.43
2	С	202	HEM	CHC-C4B-NB	5.99	130.94	124.43
2	А	201	HEM	CHC-C4B-NB	4.82	129.67	124.43
2	С	202	HEM	CHD-C1D-ND	4.58	129.41	124.43
2	В	202	HEM	C1B-NB-C4B	4.53	109.75	105.07



There are no chirality outliers.

Mol	Chain	Res	Type	Atoms
4	С	201	PEG	O1-C1-C2-O2
2	А	201	HEM	C4D-C3D-CAD-CBD
4	В	203	PEG	O2-C3-C4-O4
4	А	203	PEG	O2-C3-C4-O4
2	А	201	HEM	C2D-C3D-CAD-CBD

5 of 24 torsion outliers are listed below:

There are no ring outliers.

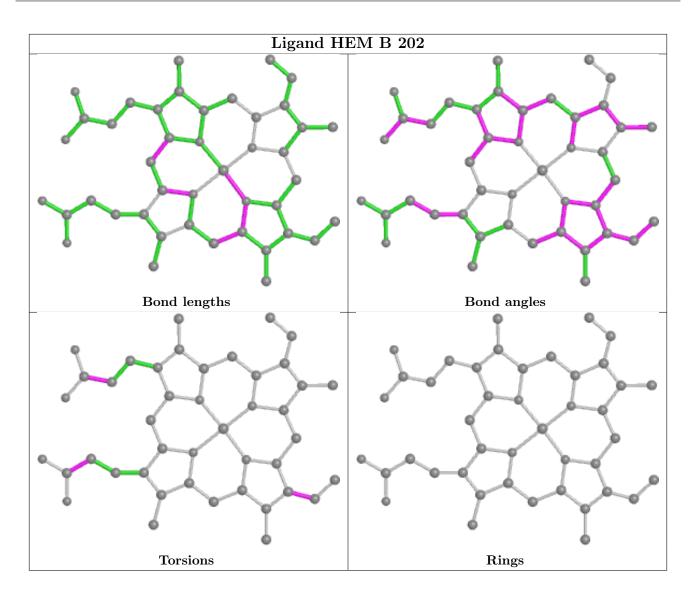
5 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	В	202	HEM	6	0
2	С	202	HEM	1	0
4	А	204	PEG	1	0
2	А	201	HEM	4	0
4	В	203	PEG	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and sufficient the outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

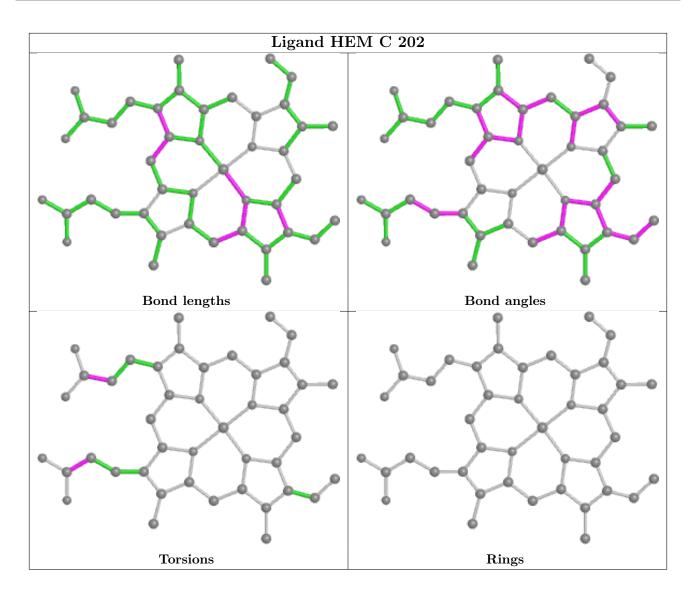




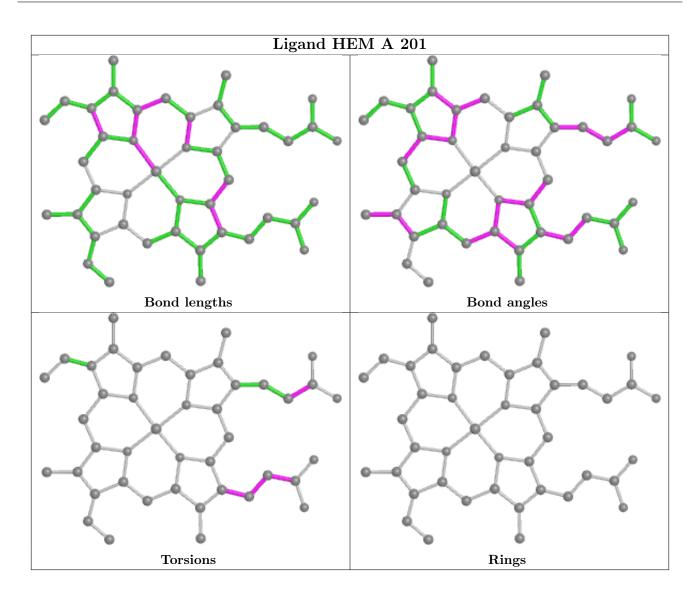












5.7 Other polymers (i)

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^{th} 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	$\langle RSRZ \rangle$	#RSRZ>2	$OWAB(Å^2)$	Q < 0.9
1	А	154/158~(97%)	-0.78	0 100 100	8, 13, 25, 32	1 (0%)
1	В	155/158~(98%)	-0.91	0 100 100	9, 14, 28, 46	2 (1%)
1	С	154/158~(97%)	-0.93	0 100 100	9, 14, 26, 33	2 (1%)
All	All	463/474 (97%)	-0.87	0 100 100	8, 14, 26, 46	5 (1%)

There are no RSRZ outliers to report.

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^{th} 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
4	PEG	А	203	7/7	0.90	0.12	$29,\!31,\!35,\!36$	0
6	CL	А	206	1/1	0.91	0.20	66,66,66,66	1
6	CL	С	205	1/1	0.91	0.20	66,66,66,66	1
4	PEG	С	201	7/7	0.92	0.14	28,30,32,35	0
4	PEG	А	204	7/7	0.94	0.16	27,30,31,32	7

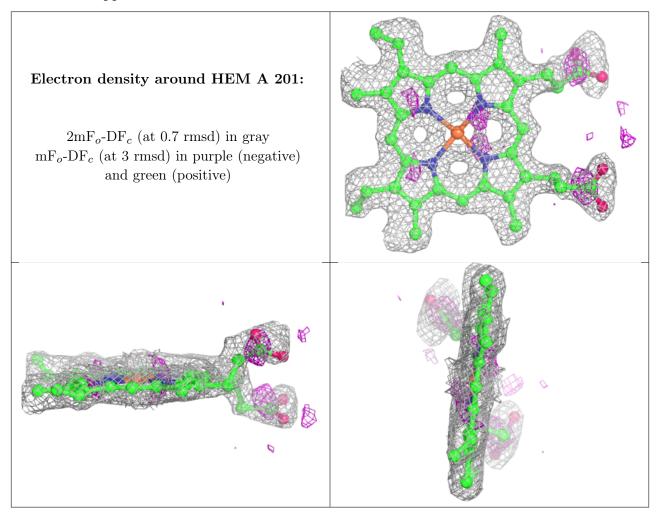
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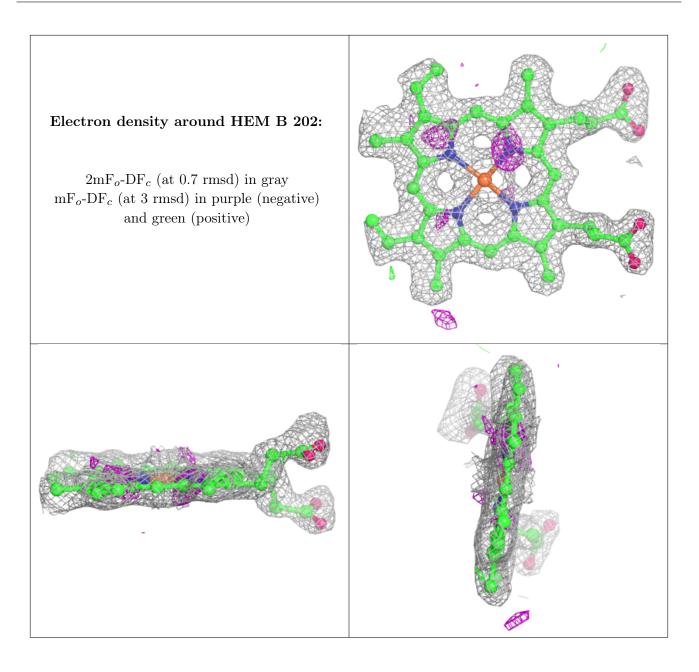
Mol	Type	Chain	Res	Atoms	RSCC	RSR	$B-factors(Å^2)$	Q < 0.9
6	CL	В	205	1/1	0.95	0.08	$66,\!66,\!66,\!66$	1
2	HEM	А	201	43/43	0.96	0.16	14,17,29,41	43
2	HEM	В	202	43/43	0.96	0.16	$14,\!17,\!24,\!27$	43
2	HEM	С	202	43/43	0.96	0.16	17,21,34,46	43
4	PEG	В	203	7/7	0.97	0.09	31,31,35,36	0
5	BA	А	205	1/1	1.00	0.07	12,12,12,12	1
5	BA	В	204	1/1	1.00	0.06	12,12,12,12	1
5	BA	С	204	1/1	1.00	0.07	7, 7, 7, 7	1
3	FE2	В	201	1/1	1.00	0.04	16,16,16,16	0
3	FE2	С	203	1/1	1.00	0.04	16,16,16,16	0
3	FE2	А	202	1/1	1.00	0.06	14,14,14,14	0

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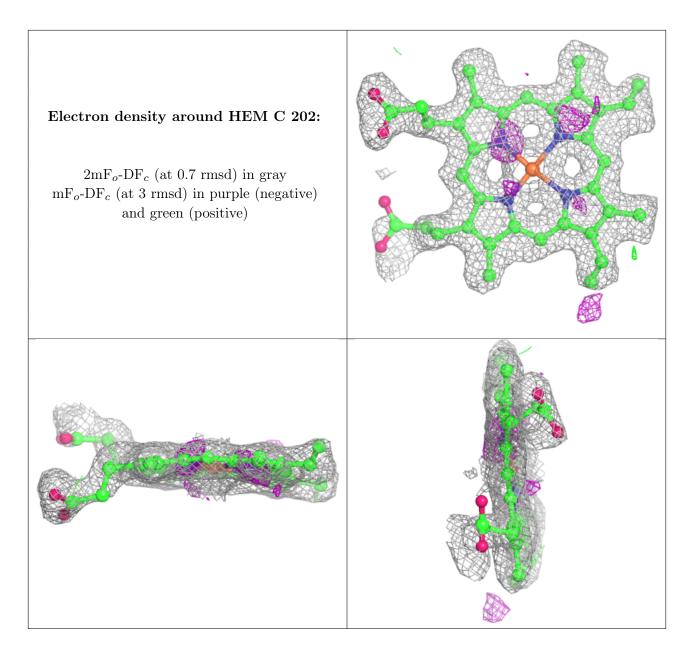
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.











6.5 Other polymers (i)

There are no such residues in this entry.

