



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 16, 2020 – 12:03 am BST

PDB ID : 5WMX
Title : Structural Insights into Substrate and Inhibitor Binding Sites in Human Indoleamine 2,3-Dioxygenase 1
Authors : Lewis-Ballester, A.; Yeh, S.R.; Pham, K.N.; Batabyal, D.; Karkashon, S.; Bonanno, J.B.; Poulos, T.M.
Deposited on : 2017-07-31
Resolution : 2.69 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
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
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

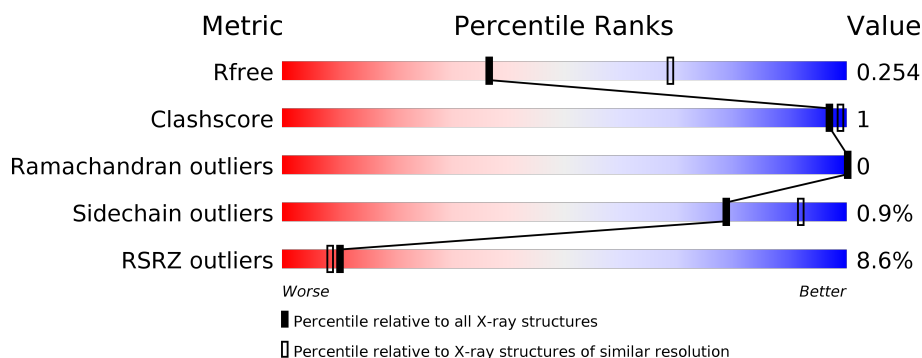
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.69 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	425	<div> <div>9%</div> <div>87%</div> <div>11%</div> </div>
1	B	425	<div> <div>6%</div> <div>87%</div> <div>11%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6277 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 Indoleamine 2,3-dioxygenase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	379	Total	C	N	O	S	0	1	0
			3002	1925	515	545	17			
1	B	378	Total	C	N	O	S	0	1	0
			3000	1926	514	543	17			

There are 68 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	MET	-	initiating methionine	UNP P14902
A	270	GLY	PHE	engineered mutation	UNP P14902
A	404	LYS	-	expression tag	UNP P14902
A	405	GLY	-	expression tag	UNP P14902
A	406	GLU	-	expression tag	UNP P14902
A	407	LEU	-	expression tag	UNP P14902
A	408	ASN	-	expression tag	UNP P14902
A	409	SER	-	expression tag	UNP P14902
A	410	LYS	-	expression tag	UNP P14902
A	411	LEU	-	expression tag	UNP P14902
A	412	GLU	-	expression tag	UNP P14902
A	413	GLY	-	expression tag	UNP P14902
A	414	LYS	-	expression tag	UNP P14902
A	415	PRO	-	expression tag	UNP P14902
A	416	ILE	-	expression tag	UNP P14902
A	417	PRO	-	expression tag	UNP P14902
A	418	ASN	-	expression tag	UNP P14902
A	419	PRO	-	expression tag	UNP P14902
A	420	LEU	-	expression tag	UNP P14902
A	421	LEU	-	expression tag	UNP P14902
A	422	GLY	-	expression tag	UNP P14902
A	423	LEU	-	expression tag	UNP P14902
A	424	ASP	-	expression tag	UNP P14902
A	425	SER	-	expression tag	UNP P14902
A	426	THR	-	expression tag	UNP P14902

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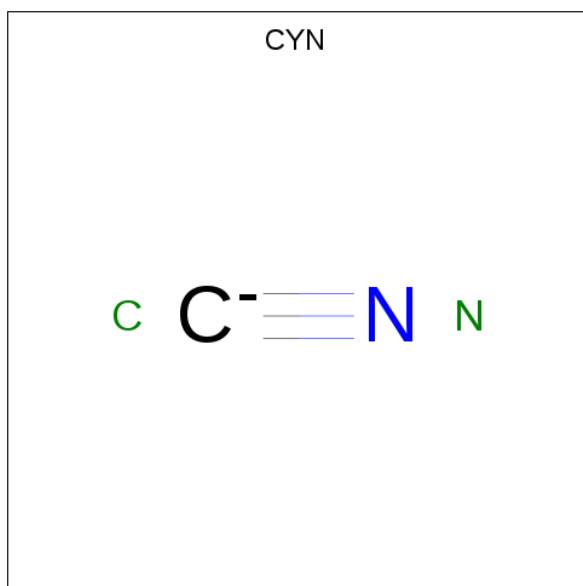
Chain	Residue	Modelled	Actual	Comment	Reference
A	427	ARG	-	expression tag	UNP P14902
A	428	THR	-	expression tag	UNP P14902
A	429	GLY	-	expression tag	UNP P14902
A	430	HIS	-	expression tag	UNP P14902
A	431	HIS	-	expression tag	UNP P14902
A	432	HIS	-	expression tag	UNP P14902
A	433	HIS	-	expression tag	UNP P14902
A	434	HIS	-	expression tag	UNP P14902
A	435	HIS	-	expression tag	UNP P14902
B	11	MET	-	initiating methionine	UNP P14902
B	270	GLY	PHE	engineered mutation	UNP P14902
B	404	LYS	-	expression tag	UNP P14902
B	405	GLY	-	expression tag	UNP P14902
B	406	GLU	-	expression tag	UNP P14902
B	407	LEU	-	expression tag	UNP P14902
B	408	ASN	-	expression tag	UNP P14902
B	409	SER	-	expression tag	UNP P14902
B	410	LYS	-	expression tag	UNP P14902
B	411	LEU	-	expression tag	UNP P14902
B	412	GLU	-	expression tag	UNP P14902
B	413	GLY	-	expression tag	UNP P14902
B	414	LYS	-	expression tag	UNP P14902
B	415	PRO	-	expression tag	UNP P14902
B	416	ILE	-	expression tag	UNP P14902
B	417	PRO	-	expression tag	UNP P14902
B	418	ASN	-	expression tag	UNP P14902
B	419	PRO	-	expression tag	UNP P14902
B	420	LEU	-	expression tag	UNP P14902
B	421	LEU	-	expression tag	UNP P14902
B	422	GLY	-	expression tag	UNP P14902
B	423	LEU	-	expression tag	UNP P14902
B	424	ASP	-	expression tag	UNP P14902
B	425	SER	-	expression tag	UNP P14902
B	426	THR	-	expression tag	UNP P14902
B	427	ARG	-	expression tag	UNP P14902
B	428	THR	-	expression tag	UNP P14902
B	429	GLY	-	expression tag	UNP P14902
B	430	HIS	-	expression tag	UNP P14902
B	431	HIS	-	expression tag	UNP P14902
B	432	HIS	-	expression tag	UNP P14902
B	433	HIS	-	expression tag	UNP P14902
B	434	HIS	-	expression tag	UNP P14902

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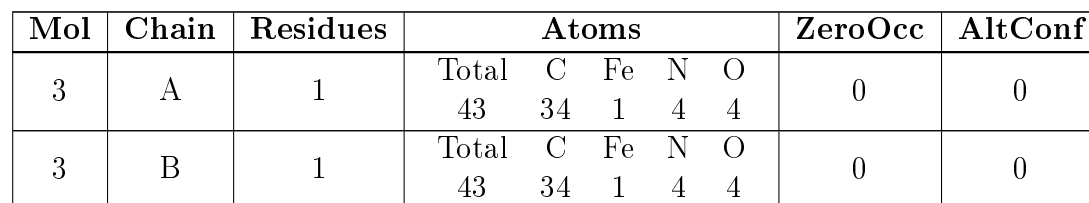
Chain	Residue	Modelled	Actual	Comment	Reference
B	435	HIS	-	expression tag	UNP P14902

- Molecule 2 is CYANIDE ION (three-letter code: CYN) (formula: CN).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			2	1	1		
2	B	1	Total	C	N	0	0
			2	1	1		

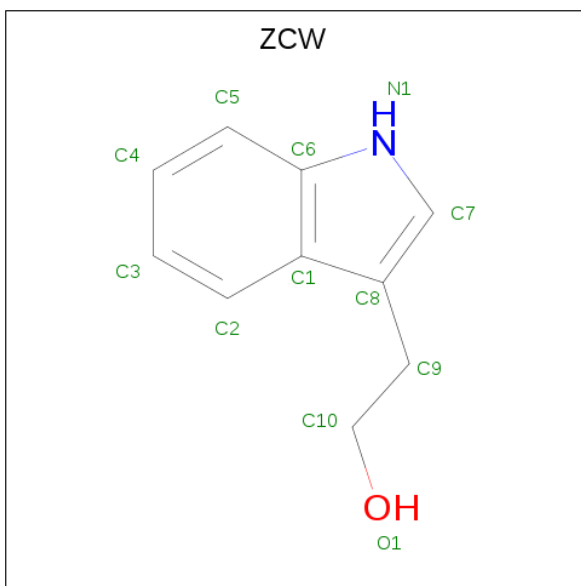
- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



- TRP
-
- The diagram shows the chemical structure of Tryptophan (TRP). The indole ring system is labeled with green text: CZ2 and CE2 for the benzene ring carbons, CE3 for the indole ring carbon, CD1 and CD2 for the indole ring carbons, and CG for the carbon attached to the side chain. The nitrogen atom in the indole ring is labeled NE1 in blue. The side chain consists of a CH2 group (labeled CH2 in green) and a CH group (labeled CB in green). The CH group is bonded to a nitrogen atom (labeled N in blue) which is part of an amino group (NH2). The CH group is also bonded to a carbon atom (labeled C in green) which is part of a carboxylate group (COO-). The carboxylate group is shown with a double bond to an oxygen atom (labeled O in red) and a single bond to a hydroxyl group (OH, labeled OXT in red). The labels are placed around the structure to identify specific atoms.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total 15	C 11	N 2	O 2	0	0
4	B	1	Total 15	C 11	N 2	O 2	0	0

- Molecule 5 is 2-(1H-indol-3-yl)ethanol (three-letter code: ZCW) (formula: $C_{10}H_{11}NO$).

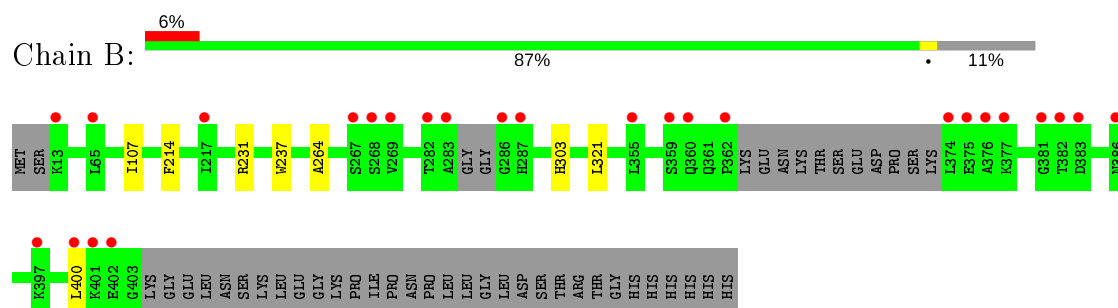


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			12	10	1	1		
5	B	1	Total	C	N	O	0	0
			12	10	1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	69	Total	O	0	0
			69	69		
6	B	62	Total	O	0	0
			62	62		

- Molecule 1: Indoleamine 2,3-dioxygenase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	87.06 Å 97.57 Å 130.42 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.10 – 2.69 39.06 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.1 (39.10-2.69) 99.2 (39.06-2.69)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.41 (at 2.69 Å)	Xtriage
Refinement program	REFMAC 5.8.0158	Depositor
R, R_{free}	0.214 , 0.254 0.214 , 0.254	Depositor DCC
R_{free} test set	1988 reflections (6.34%)	wwPDB-VP
Wilson B-factor (Å ²)	88.0	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 51.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6277	wwPDB-VP
Average B, all atoms (Å ²)	101.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.94% of the height of the origin peak. No significant pseudotranslation is detected.*

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: HEM, ZCW, CYN

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	0.36	0/3072	0.51	0/4151
1	B	0.35	0/3070	0.50	0/4149
All	All	0.36	0/6142	0.50	0/8300

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

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	3002	0	3019	4	0
1	B	3000	0	3022	3	0
2	A	2	0	0	1	0
2	B	2	0	0	1	0
3	A	43	0	30	2	0
3	B	43	0	30	2	0
4	A	15	0	9	1	0
4	B	15	0	9	0	0
5	A	12	0	0	0	0
5	B	12	0	0	0	0
6	A	69	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	B	62	0	0	0	0
All	All	6277	0	6119	11	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 1.

All (11) 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)
3:B:501:HEM:HBB2	3:B:501:HEM:HMB2	1.86	0.56
3:A:501:HEM:HMB1	3:A:501:HEM:HBB2	1.89	0.54
1:A:321:LEU:HD21	1:A:400:LEU:HD22	1.89	0.54
1:A:379:THR:HG1	4:A:502:TRP:N	2.06	0.52
1:B:231[B]:ARG:CZ	1:B:231[B]:ARG:HB3	2.43	0.49
3:B:501:HEM:HBC2	3:B:501:HEM:HHH	1.96	0.48
1:B:264:ALA:HB3	2:B:500:CYN:C	2.46	0.45
1:B:321:LEU:HD21	1:B:400:LEU:HD22	1.99	0.45
3:A:501:HEM:HMC1	3:A:501:HEM:HBC2	1.99	0.45
1:A:264:ALA:HB3	2:A:500:CYN:C	2.48	0.43
1:A:50:ILE:HD11	1:A:89:ALA:HB1	2.03	0.41

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	374/425 (88%)	357 (96%)	17 (4%)	0	100	100
1	B	373/425 (88%)	359 (96%)	14 (4%)	0	100	100
All	All	747/850 (88%)	716 (96%)	31 (4%)	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	327/367 (89%)	325 (99%)	2 (1%)	86	95
1	B	327/367 (89%)	323 (99%)	4 (1%)	71	88
All	All	654/734 (89%)	648 (99%)	6 (1%)	78	92

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	ASP
1	A	214	PHE
1	B	107	ILE
1	B	214	PHE
1	B	237	TRP
1	B	303	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	29	GLN
1	A	290	GLN

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

8 ligands 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
3	HEM	B	501	1,2	27,50,50	0.82	1 (3%)	17,82,82	1.04	0
2	CYN	B	500	3	0,1,1	0.00	-	-		
5	ZCW	B	503	-	12,13,13	1.16	0	12,17,17	1.41	1 (8%)
2	CYN	A	500	3	0,1,1	0.00	-	-		
5	ZCW	A	503	-	12,13,13	1.14	0	12,17,17	1.17	1 (8%)
3	HEM	A	501	1,2	27,50,50	0.80	1 (3%)	17,82,82	0.96	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 Torsions and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	B	501	1,2	-	1/6/54/54	-
5	ZCW	B	503	-	-	2/3/3/3	0/2/2/2
5	ZCW	A	503	-	-	2/3/3/3	0/2/2/2
3	HEM	A	501	1,2	-	0/6/54/54	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	HEM	C4D-C3D	2.42	1.48	1.42
3	B	501	HEM	C4D-C3D	2.33	1.47	1.42

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	503	ZCW	C2-C1-C6	3.01	122.16	118.17
5	A	503	ZCW	C2-C1-C6	2.48	121.45	118.17

There are no chirality outliers.

All (5) torsion outliers are listed below:

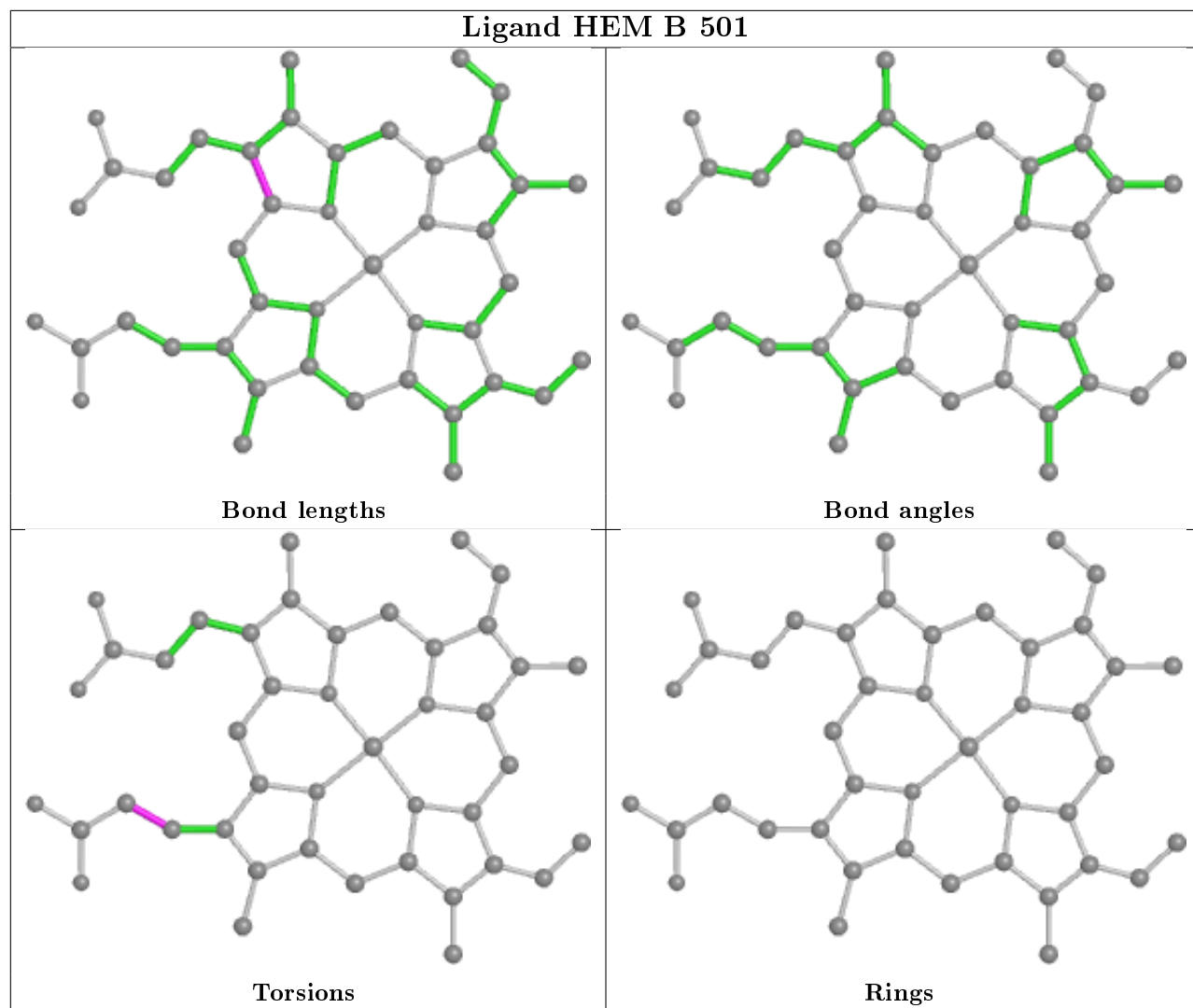
Mol	Chain	Res	Type	Atoms
5	B	503	ZCW	C1-C8-C9-C10
5	A	503	ZCW	O1-C10-C9-C8
5	A	503	ZCW	C1-C8-C9-C10
5	B	503	ZCW	C7-C8-C9-C10
3	B	501	HEM	C2A-CAA-CBA-CGA

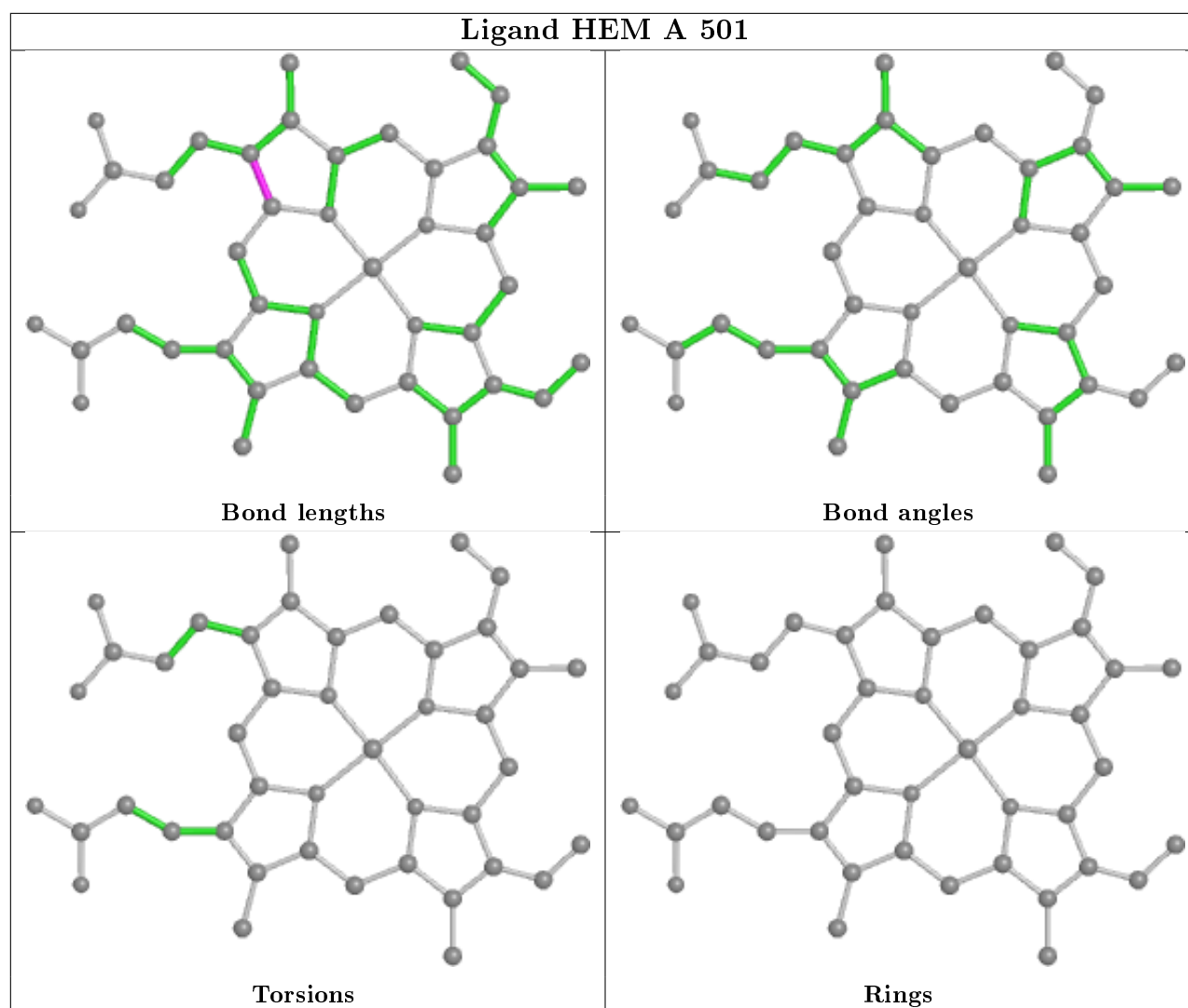
There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	501	HEM	2	0
2	B	500	CYN	1	0
2	A	500	CYN	1	0
3	A	501	HEM	2	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 than 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 any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. 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 ⓘ

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	379/425 (89%)	0.58	39 (10%) 6 5	69, 101, 143, 156	0
1	B	378/425 (88%)	0.42	26 (6%) 16 15	67, 93, 145, 186	0
All	All	757/850 (89%)	0.50	65 (8%) 10 8	67, 97, 143, 186	0

All (65) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	374	LEU	5.8
1	A	67	ILE	5.8
1	A	361	GLN	5.4
1	A	41	PHE	5.4
1	A	362	PRO	4.9
1	A	69	HIS	4.9
1	B	362	PRO	4.5
1	B	376	ALA	4.3
1	A	241	PRO	4.3
1	A	282	THR	4.2
1	B	402	GLU	4.2
1	A	251	GLY	3.9
1	B	375	GLU	3.8
1	A	70	LEU	3.8
1	B	13	LYS	3.8
1	B	355	LEU	3.5
1	B	377	LYS	3.5
1	A	402	GLU	3.4
1	B	381	GLY	3.4
1	B	386	ASN	3.3
1	A	39	TRP	3.3
1	A	355	LEU	3.2
1	A	254	GLU	3.2
1	A	142	PRO	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	32	LEU	3.0
1	A	113	GLN	2.9
1	A	13	LYS	2.9
1	B	65	LEU	2.8
1	A	224	LYS	2.7
1	A	14	GLU	2.7
1	A	383	ASP	2.7
1	A	66	SER	2.7
1	B	397	LYS	2.7
1	A	255	ASP	2.7
1	B	283	ALA	2.7
1	A	359	SER	2.7
1	B	401	LYS	2.7
1	B	268	SER	2.6
1	A	252	PHE	2.6
1	B	383	ASP	2.6
1	B	267	SER	2.6
1	A	141	LYS	2.6
1	B	400	LEU	2.5
1	A	382	THR	2.5
1	B	217	ILE	2.5
1	B	287	HIS	2.5
1	A	403	GLY	2.4
1	A	45	HIS	2.3
1	A	100	ARG	2.3
1	B	360	GLN	2.3
1	B	282	THR	2.3
1	A	38	ASP	2.3
1	A	376	ALA	2.3
1	B	382	THR	2.2
1	B	359	SER	2.2
1	A	65	LEU	2.2
1	A	378	GLY	2.2
1	A	377	LYS	2.2
1	A	285	GLY	2.2
1	A	124	LEU	2.2
1	B	286	GLY	2.1
1	A	85	CYS	2.1
1	A	68	ASP	2.1
1	A	146	GLU	2.0
1	B	269	VAL	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 ⓘ

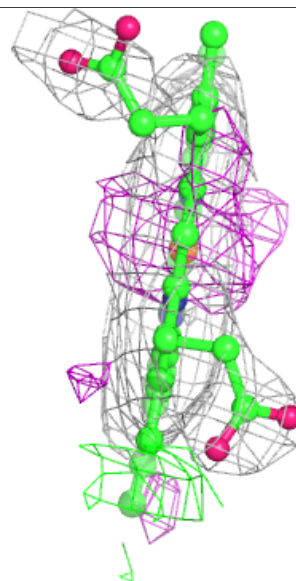
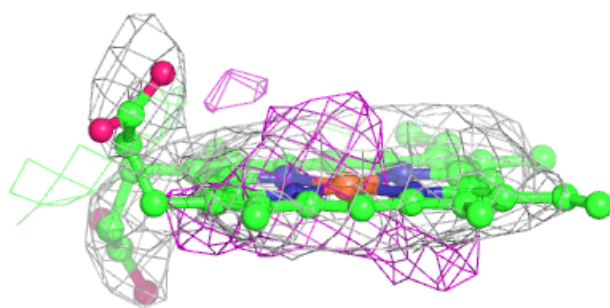
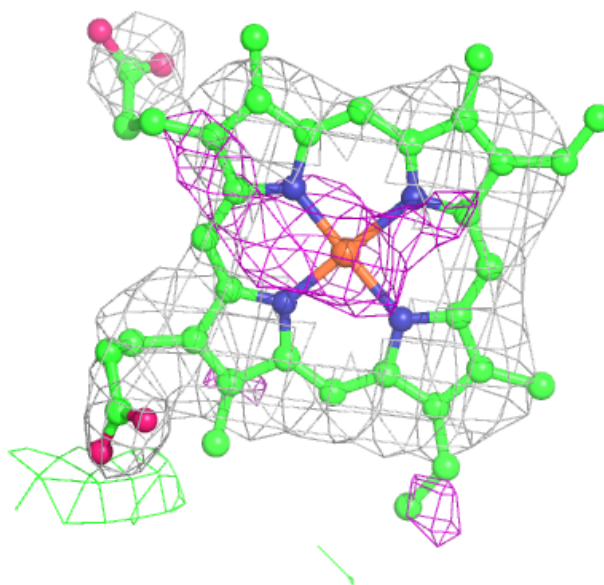
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	ZCW	B	503	12/12	0.89	0.37	89,90,94,95	0
5	ZCW	A	503	12/12	0.93	0.27	82,83,86,87	0
3	HEM	B	501	43/43	0.94	0.20	98,104,116,122	0
4	TRP	A	502	15/15	0.95	0.25	118,120,129,131	0
4	TRP	B	502	15/15	0.95	0.22	92,96,110,112	0
3	HEM	A	501	43/43	0.97	0.15	82,90,97,101	0
2	CYN	B	500	2/2	0.98	0.22	106,106,106,109	0
2	CYN	A	500	2/2	0.99	0.14	93,93,93,97	0

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.

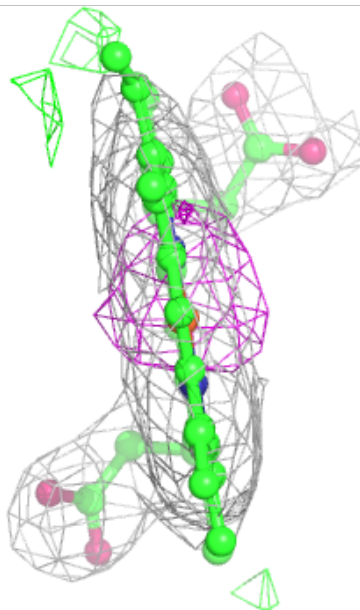
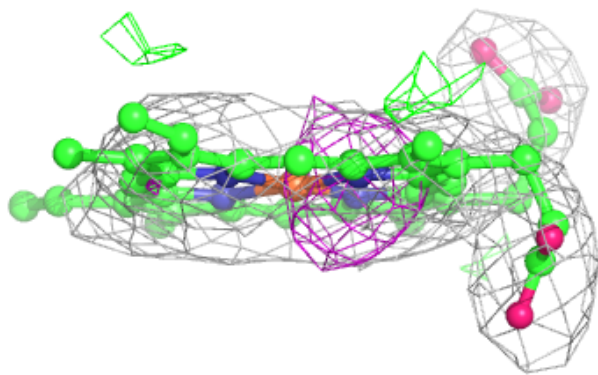
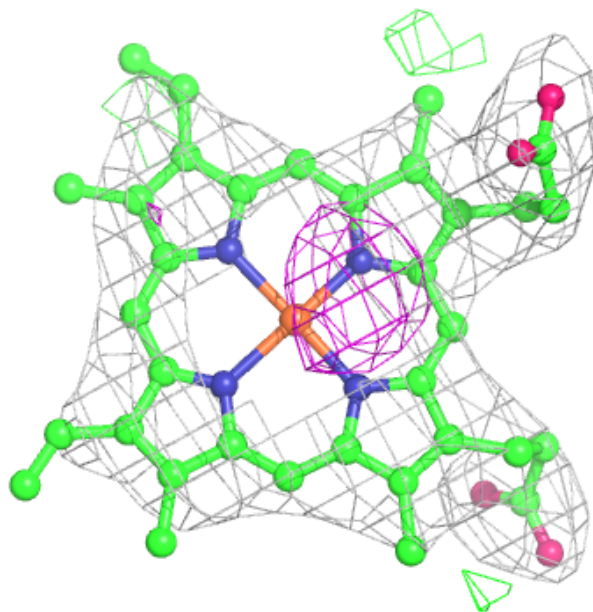
Electron density around HEM B 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 501:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.