



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 03:37 AM UTC

PDB ID : 6E35 / pdb_00006e35
Title : Crystal structure of human indoleamine 2,3-dioxygenase (IDO1) in complex with L-Trp and cyanide, Northeast Structural Genomics Target HR6160
Authors : Forouhar, F.; Lewis-Ballester, A.; Lew, S.; Karkashon, S.; Seetharaman, J.; Lu, C.; Hussain, M.; Yeh, S.-R.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2018-07-13
Resolution : 2.41 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 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	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

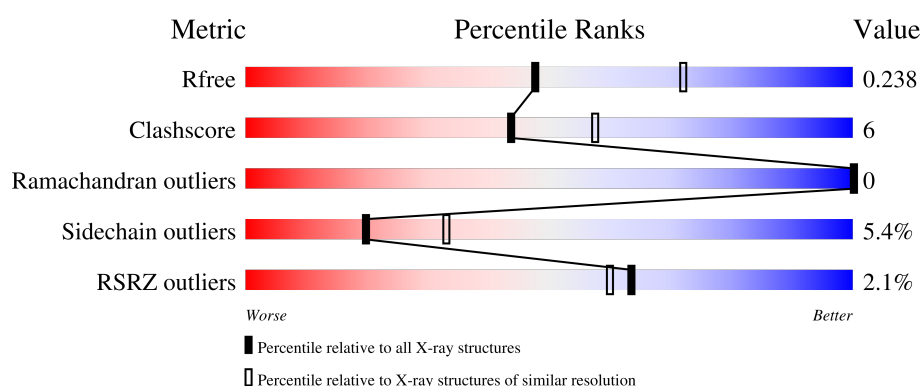
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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}	180053	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	425	2% 72% 13% • 14%
1	B	425	% 75% 10% • 14%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 6146 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	365	Total	C	N	O	S	0	0	0
			2887	1855	496	519	17			
1	B	366	Total	C	N	O	S	0	0	0
			2894	1859	496	522	17			

There are 66 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	11	MET	-	initiating methionine	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
A	427	ARG	-	expression tag	UNP P14902

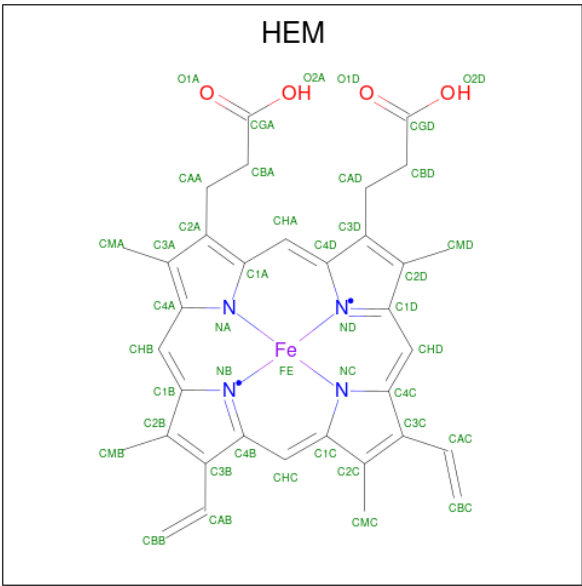
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Chain	Residue	Modelled	Actual	Comment	Reference
A	428	THR	-	expression tag	UNP P14902
A	429	GLY	-	expression tag	UNP P14902
A	430	HIS	-	expression tag	UNP P14902
A	431	HIS	-	expression tag	UNP P14902
A	432	HIS	-	expression tag	UNP P14902
A	433	HIS	-	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	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
B	435	HIS	-	expression tag	UNP P14902

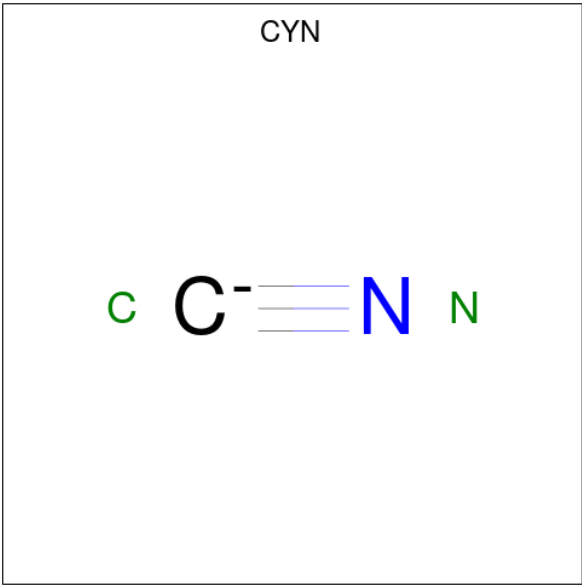
- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula:

C₃₄H₃₂FeN₄O₄).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is CYANIDE ION (CCD ID: CYN) (formula: CN).



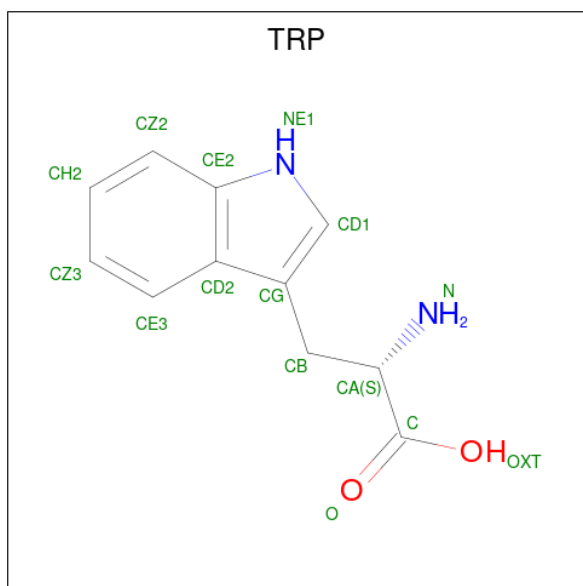
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	N	0	0
			2	1	1		

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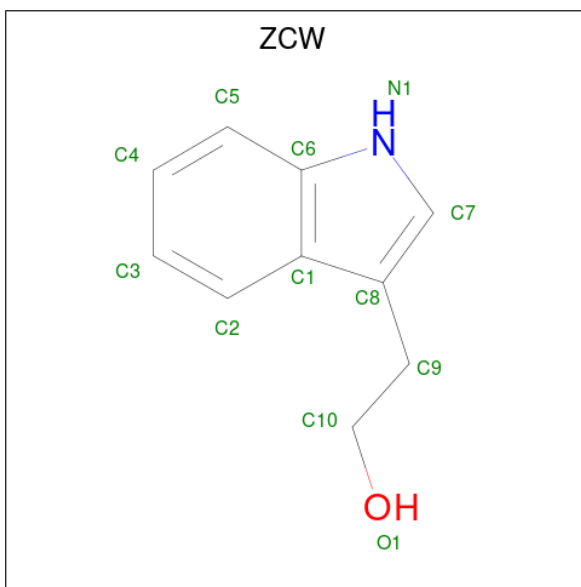
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	N	0	0
			2	1	1		

- Molecule 4 is TRYPTOPHAN (CCD ID: TRP) (formula: $C_{11}H_{12}N_2O_2$).



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

- Molecule 5 is 2-(1H-indol-3-yl)ethanol (CCD ID: 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		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	105	Total	O	0	0
			105	105		
6	B	128	Total	O	0	0
			128	128		

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	85.77Å 90.30Å 136.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.70 – 2.41 40.70 – 2.41	Depositor EDS
% Data completeness (in resolution range)	93.4 (40.70-2.41) 93.4 (40.70-2.41)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.28 (at 2.42Å)	Xtriage
Refinement program	PHENIX (1.12_2829: ???)	Depositor
R, R_{free}	0.182 , 0.236 0.184 , 0.238	Depositor DCC
R_{free} test set	3882 reflections (9.96%)	wwPDB-VP
Wilson B-factor (Å ²)	45.0	Xtriage
Anisotropy	0.391	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 42.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.012 for k,h,-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6146	wwPDB-VP
Average B, all atoms (Å ²)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.09% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CYN, HEM, ZCW

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	0.44	0/2952	0.60	0/3990
1	B	0.48	0/2960	0.66	0/4002
All	All	0.46	0/5912	0.63	0/7992

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](#)

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	2887	0	2920	37	0
1	B	2894	0	2921	33	0
2	A	43	0	30	0	0
2	B	43	0	30	3	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	15	0	9	2	0
4	B	15	0	9	3	0
5	A	12	0	0	1	0
6	A	105	0	0	2	0
6	B	128	0	0	3	0
All	All	6146	0	5919	72	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 6.

All (72) 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:55:LEU:HD21	1:A:89:ALA:HB1	1.65	0.78
1:A:363:LYS:HD3	1:A:363:LYS:H	1.54	0.73
1:B:245:ASP:O	6:B:601:HOH:O	2.10	0.70
1:B:102:VAL:HA	1:B:248:VAL:HG22	1.74	0.69
1:B:155:ARG:HH21	1:B:155:ARG:HG3	1.68	0.57
1:A:94:LYS:HA	1:A:155:ARG:HH11	1.70	0.56
1:B:155:ARG:NH1	6:B:604:HOH:O	2.38	0.55
1:B:126:TYR:OH	1:B:171:GLU:HG3	2.06	0.54
1:A:48:ASP:OD2	1:A:48:ASP:N	2.40	0.54
1:B:379:THR:HG1	4:B:503:TRP:N	2.07	0.52
2:B:501:HEM:HMB1	2:B:501:HEM:HBB2	1.92	0.51
1:A:50:ILE:HD11	1:A:93:GLY:C	2.35	0.51
1:A:379:THR:HG1	4:A:503:TRP:N	2.09	0.51
1:A:237:TRP:HB2	1:A:246:GLY:HA2	1.93	0.50
1:A:50:ILE:HG21	1:A:155:ARG:HD3	1.93	0.50
1:B:163:PHE:HZ	4:B:503:TRP:CD1	2.29	0.50
1:A:138:ASP:HB3	1:A:141:LYS:HB2	1.94	0.49
1:A:343:ARG:HG2	5:A:504:ZCW:C10	2.43	0.49
1:A:126:TYR:OH	1:A:171:GLU:HG3	2.12	0.48
1:B:321:LEU:HD21	1:B:400:LEU:HD22	1.95	0.48
1:B:44:LYS:HE2	1:B:152:PHE:CE1	2.49	0.48
1:A:321:LEU:HD21	1:A:400:LEU:HD22	1.95	0.47
1:B:356:ILE:HB	1:B:357:PRO:HD3	1.95	0.47
1:A:52:SER:HB2	1:A:54:GLN:HG2	1.97	0.47
1:B:198:LYS:HB3	1:B:198:LYS:HE3	1.69	0.47
1:B:55:LEU:O	1:B:55:LEU:HG	1.96	0.47
1:B:155:ARG:HG3	1:B:155:ARG:NH2	2.29	0.47
1:A:102:VAL:HA	1:A:248:VAL:HG13	1.97	0.47
1:A:223:PRO:HG3	1:A:352:LYS:HE2	1.98	0.46
1:A:72:ASP:OD2	1:A:72:ASP:N	2.47	0.46
1:A:94:LYS:O	1:A:155:ARG:NH1	2.49	0.46
1:B:155:ARG:HD2	6:B:647:HOH:O	2.17	0.45
1:A:163:PHE:HZ	4:A:503:TRP:CD1	2.34	0.45
1:A:46:LEU:HD22	1:A:154:PHE:HA	1.99	0.45
1:B:27:ASN:HD21	1:B:136:LYS:HD2	1.82	0.45
1:B:141:LYS:HA	1:B:141:LYS:HE2	1.98	0.45
1:A:28:PRO:HG2	1:A:134:TRP:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:321:LEU:HD23	1:A:321:LEU:HA	1.60	0.44
1:A:34:ASP:O	1:A:37:ASN:HB2	2.18	0.44
1:A:56:ARG:HG3	1:A:90:TYR:CE1	2.53	0.44
1:A:232:ILE:H	1:A:232:ILE:HG13	1.58	0.44
1:A:224:LYS:HD2	1:A:224:LYS:HA	1.75	0.44
1:B:177:ALA:O	1:B:180:VAL:HG12	2.18	0.44
1:B:214:PHE:HD1	1:B:214:PHE:O	2.01	0.43
1:A:319:PHE:O	1:A:323:LYS:HG2	2.18	0.43
2:B:501:HEM:HBB2	2:B:501:HEM:CMB	2.49	0.43
1:A:196:LEU:HD13	1:A:328:LEU:HD11	2.01	0.43
1:A:198:LYS:HG2	6:A:634:HOH:O	2.19	0.43
1:B:115:SER:HB3	1:B:120:LEU:O	2.19	0.42
1:B:310:LEU:HD23	1:B:310:LEU:HA	1.84	0.42
1:A:72:ASP:O	1:A:76:GLN:HG3	2.19	0.42
1:A:245:ASP:O	6:A:601:HOH:O	2.21	0.42
1:B:58:ARG:HE	1:B:58:ARG:HB3	1.46	0.42
1:B:100:ARG:HA	1:B:100:ARG:HD3	1.78	0.42
1:B:97:GLY:O	1:B:99:VAL:HG13	2.20	0.41
1:B:393:SER:O	1:B:397:LYS:HG3	2.20	0.41
1:B:323:LYS:HD3	1:B:323:LYS:HA	1.76	0.41
1:B:189:GLN:HE21	1:B:189:GLN:HB2	1.60	0.41
1:B:225:ALA:O	1:B:229:VAL:HB	2.21	0.41
1:A:120:LEU:HD23	1:A:120:LEU:HA	1.79	0.41
1:A:56:ARG:NH2	1:A:100:ARG:HG2	2.35	0.41
1:B:388:LEU:HD21	2:B:501:HEM:C1A	2.56	0.41
1:A:61:LYS:HD3	1:A:61:LYS:HA	1.94	0.41
1:A:108:ALA:HB1	1:A:249:TYR:HB3	2.02	0.41
1:A:109:VAL:HB	1:A:110:PRO:HD3	2.03	0.41
1:A:310:LEU:HD23	1:A:310:LEU:HA	1.82	0.41
1:B:155:ARG:HH21	1:B:155:ARG:CG	2.32	0.41
1:B:328:LEU:HD12	1:B:328:LEU:HA	1.78	0.40
1:B:120:LEU:HD23	1:B:120:LEU:HA	1.95	0.40
1:A:118:LEU:HD23	1:A:118:LEU:HA	1.87	0.40
1:B:101:LYS:O	1:B:247:LEU:HD12	2.22	0.40
1:B:163:PHE:CZ	4:B:503:TRP:CD1	3.09	0.40

There are no symmetry-related clashes.

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	361/425 (85%)	344 (95%)	17 (5%)	0	100	100
1	B	362/425 (85%)	348 (96%)	14 (4%)	0	100	100
All	All	723/850 (85%)	692 (96%)	31 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	315/368 (86%)	299 (95%)	16 (5%)	21	35
1	B	316/368 (86%)	298 (94%)	18 (6%)	18	31
All	All	631/736 (86%)	597 (95%)	34 (5%)	20	33

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	63	ASN
1	A	78	LEU
1	A	106	ASN
1	A	135	LYS
1	A	141	LYS
1	A	149	ASP
1	A	171	GLU

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Mol	Chain	Res	Type
1	A	180	VAL
1	A	196	LEU
1	A	237	TRP
1	A	243	LEU
1	A	248	VAL
1	A	308	CYS
1	A	333	ASP
1	A	363	LYS
1	B	54	GLN
1	B	55	LEU
1	B	58	ARG
1	B	62	LEU
1	B	78	LEU
1	B	153	SER
1	B	155	ARG
1	B	171	GLU
1	B	176	SER
1	B	189	GLN
1	B	196	LEU
1	B	198	LYS
1	B	202	GLU
1	B	209	LYS
1	B	212	GLN
1	B	248	VAL
1	B	361	GLN
1	B	393	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	96	HIS
1	B	27	ASN
1	B	29	GLN
1	B	191	GLN

5.3.3 RNA ⓘ

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 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$
4	TRP	B	503	-	15,16,16	0.98	0	18,22,22	0.76	1 (5%)
2	HEM	A	501	1,3	50,50,50	1.76	9 (18%)	67,82,82	2.36	28 (41%)
3	CYN	A	502	2	1,1,1	0.49	0	-		
5	ZCW	A	504	-	13,13,13	2.05	4 (30%)	16,17,17	0.88	1 (6%)
2	HEM	B	501	1,3	50,50,50	1.75	8 (16%)	67,82,82	2.12	25 (37%)
3	CYN	B	502	2	1,1,1	0.58	0	-		
4	TRP	A	503	-	15,16,16	1.02	1 (6%)	18,22,22	0.71	1 (5%)

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
4	TRP	B	503	-	-	4/8/8/8	0/2/2/2
2	HEM	A	501	1,3	-	7/14/54/54	-
5	ZCW	A	504	-	-	2/3/3/3	0/2/2/2
2	HEM	B	501	1,3	-	4/14/54/54	-
4	TRP	A	503	-	-	2/8/8/8	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	FE-ND	7.20	2.17	1.94
2	A	501	HEM	FE-ND	6.67	2.15	1.94
2	A	501	HEM	FE-NA	4.45	2.09	1.95
2	B	501	HEM	FE-NA	4.18	2.09	1.95
5	A	504	ZCW	C1-C6	-4.14	1.36	1.41
5	A	504	ZCW	C5-C6	4.14	1.46	1.39
2	B	501	HEM	FE-NC	3.97	2.08	1.95
2	A	501	HEM	FE-NB	3.53	2.05	1.94
2	A	501	HEM	CAB-C3B	3.06	1.55	1.47
5	A	504	ZCW	C3-C2	3.01	1.44	1.38
2	A	501	HEM	FE-NC	2.94	2.04	1.95
2	B	501	HEM	CAB-C3B	2.87	1.55	1.47
2	B	501	HEM	CAC-C3C	2.62	1.54	1.47
5	A	504	ZCW	C7-N1	-2.50	1.33	1.37
2	A	501	HEM	CAC-C3C	2.23	1.53	1.47
2	B	501	HEM	CMB-C2B	2.23	1.55	1.50
4	A	503	TRP	OXT-C	-2.12	1.23	1.30
2	A	501	HEM	CMA-C3A	2.12	1.55	1.50
2	A	501	HEM	CMD-C2D	2.10	1.55	1.50
2	B	501	HEM	CMC-C2C	2.09	1.55	1.50
2	B	501	HEM	CMA-C3A	2.08	1.55	1.50
2	A	501	HEM	O1D-CGD	2.05	1.28	1.22

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	HEM	CHD-C1D-ND	5.53	130.38	124.42
2	A	501	HEM	C4D-ND-C1D	5.36	111.56	105.21
2	A	501	HEM	C3D-C4D-ND	-5.32	104.33	110.17
2	A	501	HEM	CHC-C1C-NC	4.70	129.57	124.45
2	A	501	HEM	CHD-C1D-ND	4.47	129.23	124.42
2	A	501	HEM	C2D-C1D-ND	-4.44	104.77	109.90
2	A	501	HEM	CHA-C4D-ND	4.42	129.83	124.37
2	B	501	HEM	C2A-C1A-NA	-4.11	105.59	110.15
2	B	501	HEM	C2D-C1D-ND	-3.98	105.30	109.90
2	A	501	HEM	C3B-C2B-C1B	3.92	109.35	106.41
2	A	501	HEM	C1B-NB-C4B	3.91	109.84	105.21
2	B	501	HEM	CHC-C4B-NB	3.85	128.57	124.42
2	B	501	HEM	C1A-CHA-C4D	-3.78	117.37	126.25
2	B	501	HEM	CHC-C1C-NC	3.74	128.53	124.45
2	A	501	HEM	C2A-C1A-NA	-3.70	106.05	110.15
2	B	501	HEM	C4D-ND-C1D	3.69	109.57	105.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CBD-CAD-C3D	-3.68	102.34	112.53
2	A	501	HEM	C3B-C4B-NB	-3.53	106.93	109.47
2	A	501	HEM	C4C-C3C-C2C	3.49	109.84	106.81
2	B	501	HEM	CHB-C4A-NA	3.47	130.15	123.86
2	A	501	HEM	CHB-C4A-NA	3.43	130.09	123.86
2	B	501	HEM	CHA-C4D-ND	3.43	128.60	124.37
2	B	501	HEM	C3D-C4D-ND	-3.42	106.42	110.17
2	A	501	HEM	C2B-C1B-NB	-3.26	106.10	109.84
2	A	501	HEM	CHB-C1B-NB	3.13	128.24	124.37
2	B	501	HEM	C1B-NB-C4B	3.10	108.88	105.21
2	B	501	HEM	C3B-C4B-NB	-3.01	107.31	109.47
2	B	501	HEM	C1D-C2D-C3D	2.92	110.05	106.98
2	A	501	HEM	C1A-CHA-C4D	-2.79	119.69	126.25
2	B	501	HEM	CAC-C3C-C4C	2.74	131.36	124.82
2	B	501	HEM	CHD-C4C-C3C	2.74	129.82	125.21
2	B	501	HEM	C3A-C4A-NA	-2.66	105.88	110.14
4	B	503	TRP	OXT-C-O	-2.65	118.06	124.08
2	B	501	HEM	C4A-NA-C1A	2.65	110.14	105.82
2	B	501	HEM	CAC-C3C-C2C	-2.60	119.99	128.43
2	B	501	HEM	C2B-C1B-NB	-2.56	106.90	109.84
2	A	501	HEM	CHC-C4B-NB	2.51	127.12	124.42
2	A	501	HEM	CAC-C3C-C2C	-2.51	120.28	128.43
2	A	501	HEM	C3A-C4A-NA	-2.45	106.22	110.14
2	B	501	HEM	C1C-CHC-C4B	-2.38	120.95	126.02
2	B	501	HEM	CHA-C1A-NA	2.38	128.18	123.86
2	B	501	HEM	C3B-C2B-C1B	2.34	108.17	106.41
2	A	501	HEM	CMB-C2B-C1B	-2.33	121.40	125.03
2	A	501	HEM	O2A-CGA-CBA	2.26	121.14	114.00
2	A	501	HEM	C4D-C3D-C2D	2.26	110.18	106.89
2	B	501	HEM	C4A-C3A-C2A	2.24	109.38	106.82
2	A	501	HEM	C4A-C3A-C2A	2.19	109.33	106.82
2	A	501	HEM	C4A-NA-C1A	2.19	109.39	105.82
2	A	501	HEM	C4C-NC-C1C	2.14	109.32	105.82
2	B	501	HEM	C3C-C2C-C1C	2.12	109.06	107.05
2	A	501	HEM	CAC-C3C-C4C	2.11	129.87	124.82
2	B	501	HEM	C4C-C3C-C2C	2.11	108.64	106.81
2	A	501	HEM	C1D-C2D-C3D	2.07	109.16	106.98
2	A	501	HEM	C3C-C2C-C1C	2.07	109.00	107.05
4	A	503	TRP	OXT-C-O	-2.04	119.46	124.08
5	A	504	ZCW	C5-C6-C1	-2.01	120.24	122.19

There are no chirality outliers.

All (19) torsion outliers are listed below:

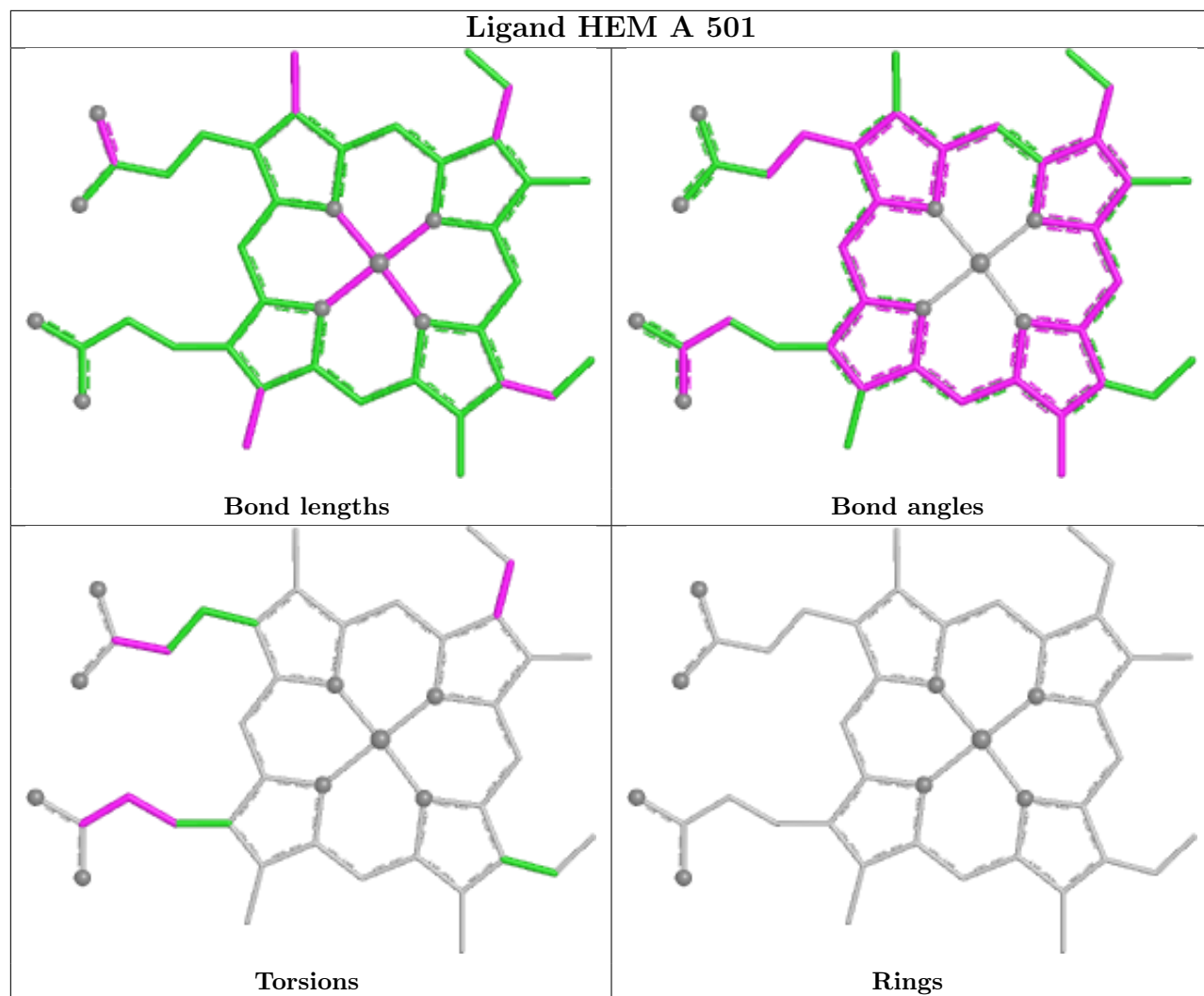
Mol	Chain	Res	Type	Atoms
5	A	504	ZCW	O1-C10-C9-C8
4	A	503	TRP	CA-CB-CG-CD2
2	A	501	HEM	C2A-CAA-CBA-CGA
2	A	501	HEM	C2C-C3C-CAC-CBC
2	A	501	HEM	C4C-C3C-CAC-CBC
4	A	503	TRP	CA-CB-CG-CD1
4	B	503	TRP	CA-CB-CG-CD2
2	B	501	HEM	CAD-CBD-CGD-O1D
2	A	501	HEM	CAA-CBA-CGA-O2A
2	B	501	HEM	CAD-CBD-CGD-O2D
2	A	501	HEM	CAA-CBA-CGA-O1A
5	A	504	ZCW	C1-C8-C9-C10
2	B	501	HEM	CAA-CBA-CGA-O2A
2	B	501	HEM	CAA-CBA-CGA-O1A
4	B	503	TRP	CA-CB-CG-CD1
4	B	503	TRP	OXT-C-CA-CB
2	A	501	HEM	CAD-CBD-CGD-O1D
4	B	503	TRP	O-C-CA-CB
2	A	501	HEM	CAD-CBD-CGD-O2D

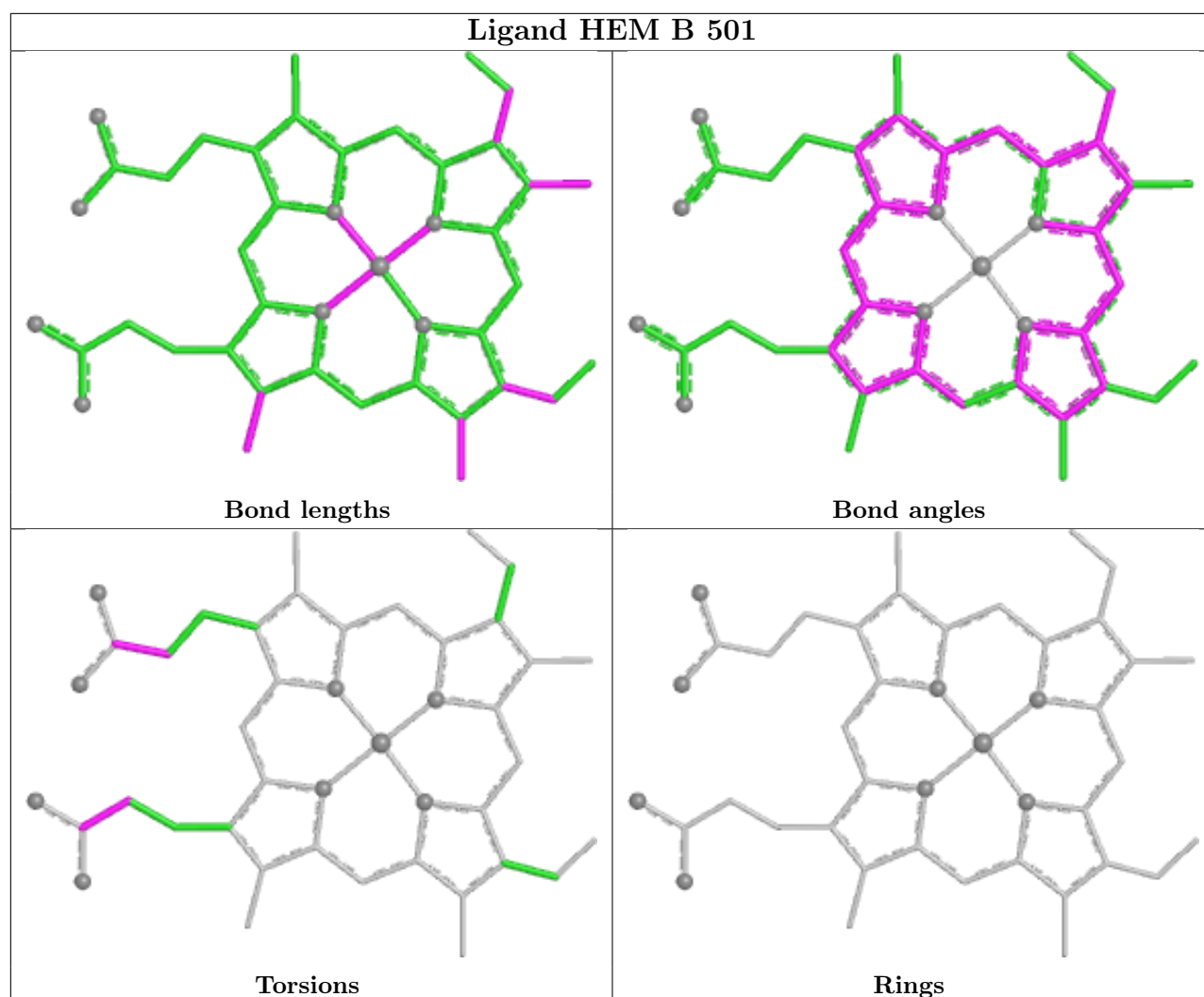
There are no ring outliers.

4 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	503	TRP	3	0
5	A	504	ZCW	1	0
2	B	501	HEM	3	0
4	A	503	TRP	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 [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	365/425 (85%)	-0.02	9 (2%) 58 55	28, 54, 102, 145	0
1	B	366/425 (86%)	-0.27	6 (1%) 70 67	29, 47, 72, 146	0
All	All	731/850 (86%)	-0.14	15 (2%) 63 60	28, 49, 95, 146	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	374	LEU	4.2
1	B	26	PRO	3.5
1	B	401	LYS	3.3
1	B	374	LEU	3.1
1	A	50	ILE	3.1
1	A	401	LYS	2.9
1	A	63	ASN	2.6
1	B	54	GLN	2.6
1	B	402	GLU	2.6
1	B	362	PRO	2.4
1	A	41	PHE	2.3
1	A	362	PRO	2.1
1	A	55	LEU	2.0
1	A	57	GLU	2.0
1	A	46	LEU	2.0

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 oligosaccharides in this entry.

6.4 Ligands

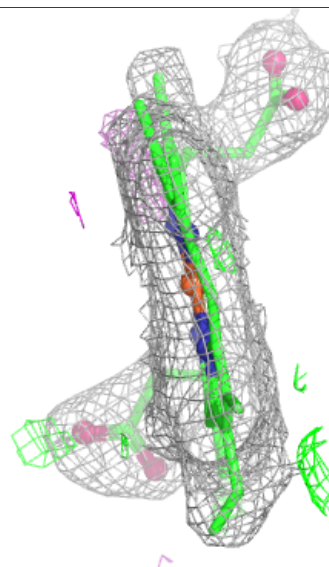
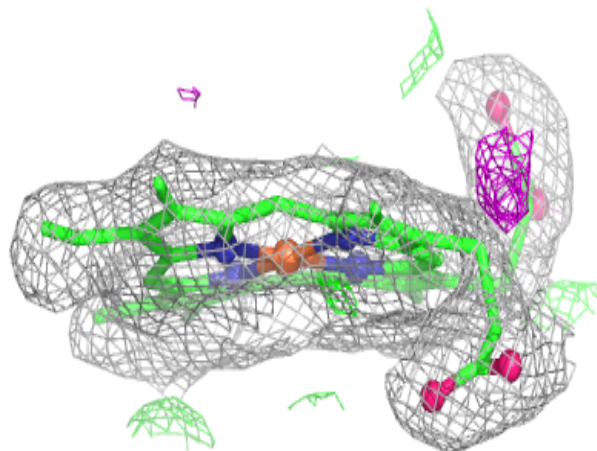
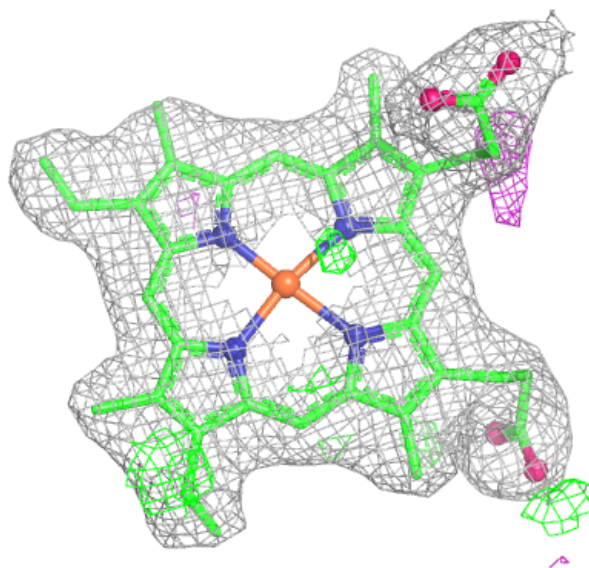
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	TRP	A	503	15/15	0.93	0.10	39,43,48,52	0
5	ZCW	A	504	12/12	0.93	0.14	46,73,84,86	0
3	CYN	B	502	2/2	0.94	0.12	31,31,31,38	0
4	TRP	B	503	15/15	0.95	0.07	33,40,43,43	0
3	CYN	A	502	2/2	0.97	0.12	31,31,31,32	0
2	HEM	B	501	43/43	0.98	0.07	31,37,45,51	0
2	HEM	A	501	43/43	0.98	0.07	27,38,48,58	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.

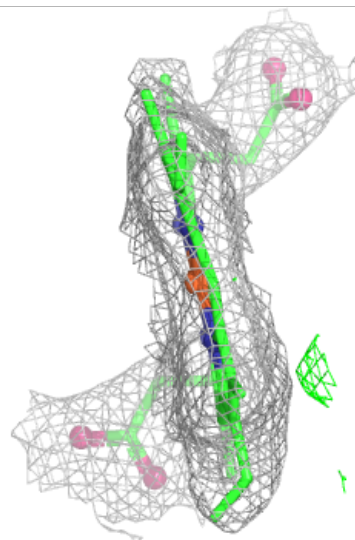
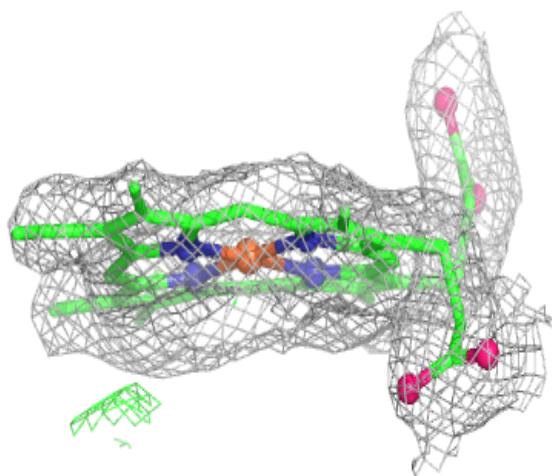
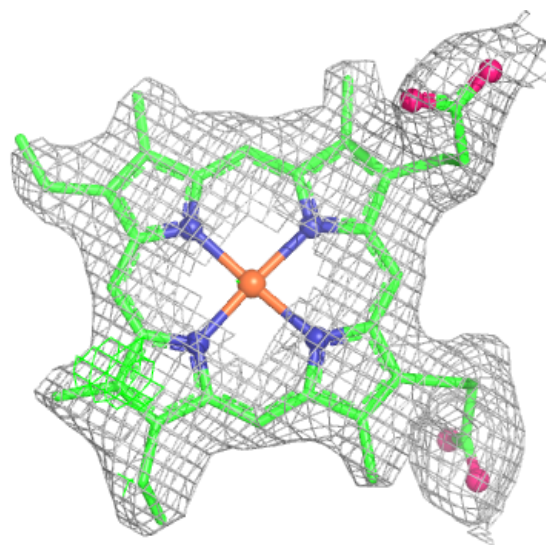
Electron density around HEM B 501:

$2mF_o-DF_c$ (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 [i](#)

There are no such residues in this entry.