Full wwPDB X-ray Structure Validation Report (i)

Aug 5, 2026 – 03:20 PM EDT

PDB ID	:	2IG3 / pdb_00002ig3
Title	:	Crystal structure of group III truncated hemoglobin from Campylobacter jejuni
Authors	:	Nardini, M.; Pesce, A.; Labarre, M.; Ascenzi, P.; Guertin, M.; Bolognesi, M.
Deposited on	:	2006-09-22
Resolution	:	2.15 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 (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 (i)) were used in the production of this report:

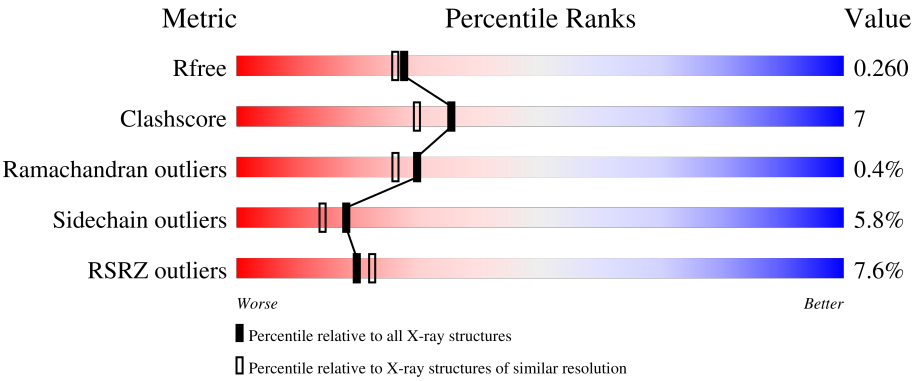
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	A	127	10% 83% 13% ·
1	B	127	5% 88% 8% · ·

2 Entry composition (i)

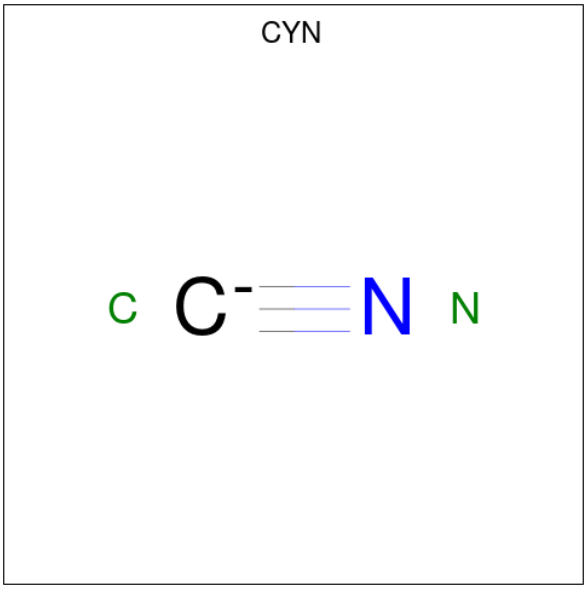
There are 6 unique types of molecules in this entry. The entry contains 2315 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Group III truncated haemoglobin.

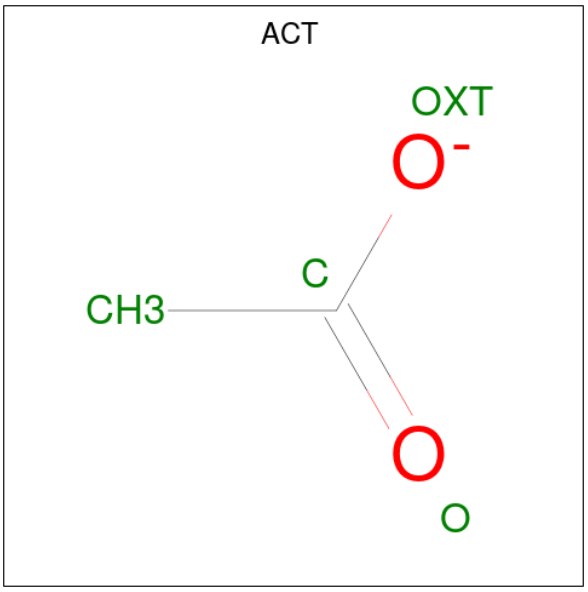
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	127	Total	C	N	O	S	0	0	0
			1056	687	172	191	6			
1	B	124	Total	C	N	O	S	0	1	0
			1044	681	169	188	6			

- Molecule 2 is CYANIDE ION (CCD ID: CYN) (formula: CN).



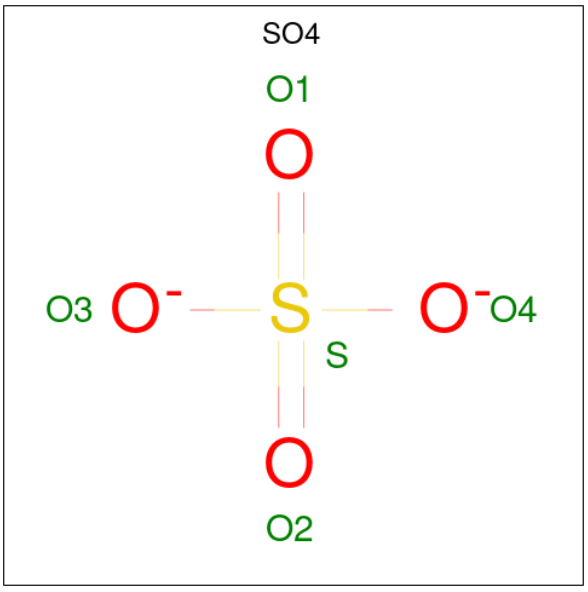
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	N	0	0
			2	1	1		
2	B	1	Total	C	N	0	0
			2	1	1		

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O_4S).



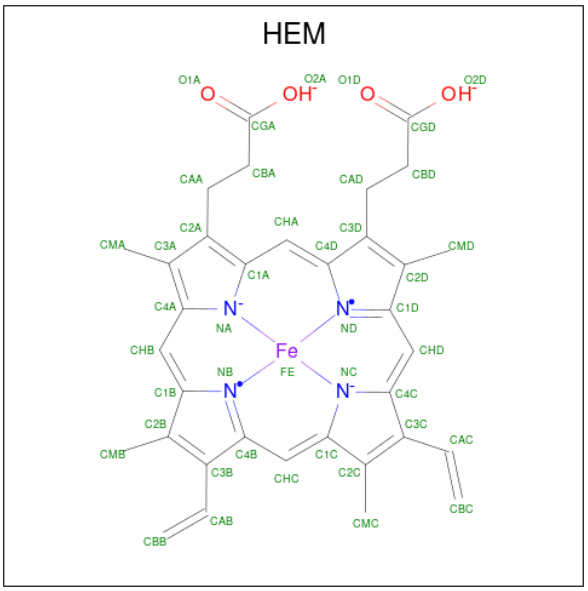
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	O	S	0	0
			5	4	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

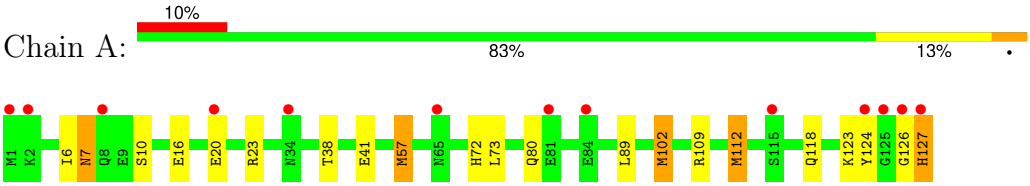
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	50	Total O 50 50	0	0
6	B	42	Total O 43 43	0	1

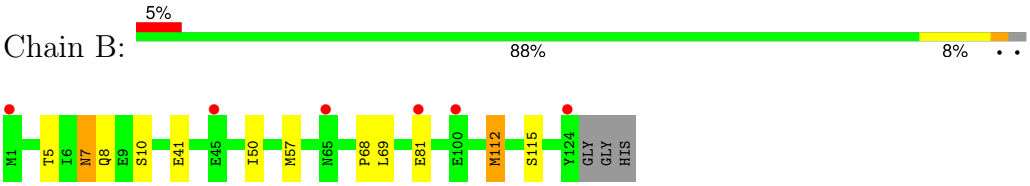
3 Residue-property plots (i)

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: Group III truncated haemoglobin



• Molecule 1: Group III truncated haemoglobin



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α, β, γ	95.59Å 95.59Å 65.42Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	31.30 – 2.15 31.30 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (31.30-2.15) 99.7 (31.30-2.15)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	7.18 (at 2.16Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, R_{free}	0.206 , 0.262 0.208 , 0.260	Depositor DCC
R_{free} test set	1935 reflections (10.15%)	wwPDB-VP
Wilson B-factor (Å^2)	35.4	Xtriage
Anisotropy	0.358	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 37.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.040 for -h,-k,l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	2315	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.11% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: ACT, SO4, CYN, HEM

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	0.62	0/1084	0.89	0/1456
1	B	0.60	0/1076	0.91	2/1446 (0.1%)
All	All	0.61	0/2160	0.90	2/2902 (0.1%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	8	GLN	OE1-CD-NE2	-10.26	112.34	122.60
1	B	8	GLN	CG-CD-NE2	6.58	126.27	116.40

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	1056	0	1036	23	0
1	B	1044	0	1026	16	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	4	0	3	0	0
3	B	8	0	6	0	0
4	A	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	15	0	0	0	0
5	A	43	0	30	1	0
5	B	43	0	30	2	0
6	A	50	0	0	1	0
6	B	43	0	0	0	0
All	All	2315	0	2131	32	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 7.

All (32) 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:16:GLU:O	1:A:20:GLU:HG2	1.59	1.01
1:A:127:HIS:HE1	1:B:68:PRO:HG2	1.31	0.94
1:A:127:HIS:O	1:B:69:LEU:HB3	1.72	0.88
1:A:127:HIS:CE1	1:B:68:PRO:HG2	2.11	0.85
1:A:6:ILE:HG13	1:A:102:MET:HE3	1.63	0.78
1:A:127:HIS:HB3	1:B:69:LEU:HB2	1.66	0.77
1:B:112:MET:HE2	1:B:112:MET:O	1.93	0.68
1:A:6:ILE:CG1	1:A:102:MET:HE3	2.29	0.63
1:B:81:GLU:HA	1:B:81:GLU:OE1	2.04	0.58
1:A:127:HIS:CE1	1:B:68:PRO:CG	2.86	0.57
1:A:23:ARG:NH2	1:A:38:THR:O	2.38	0.56
1:B:7:ASN:ND2	1:B:10:SER:H	2.03	0.56
1:A:6:ILE:CD1	1:A:102:MET:HE3	2.36	0.56
1:A:80:GLN:HE21	1:A:118:GLN:HE21	1.54	0.55
1:A:6:ILE:HG13	1:A:102:MET:CE	2.36	0.54
1:A:127:HIS:C	1:B:69:LEU:HB3	2.35	0.52
1:A:7:ASN:ND2	1:A:10:SER:H	2.08	0.51
5:B:700:HEM:HBB2	5:B:700:HEM:HMB2	1.91	0.51
1:B:112:MET:HE2	1:B:112:MET:C	2.36	0.50
1:B:7:ASN:C	1:B:7:ASN:HD22	2.20	0.49
1:B:81:GLU:OE1	1:B:81:GLU:CA	2.58	0.49
1:A:7:ASN:HD22	1:A:10:SER:H	1.61	0.48
5:B:700:HEM:HBB2	5:B:700:HEM:CMB	2.48	0.43
1:A:123:LYS:HB2	1:B:112:MET:HE3	2.00	0.43
1:A:109:ARG:HD2	6:A:822:HOH:O	2.17	0.43
1:A:127:HIS:HB3	1:B:69:LEU:CB	2.45	0.42
1:A:112:MET:HA	1:A:112:MET:CE	2.50	0.42
1:A:57:MET:HG3	1:A:109:ARG:NH2	2.35	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:73:LEU:HD13	1:B:115:SER:HB2	2.02	0.41
1:A:123:LYS:NZ	1:A:127:HIS:CD2	2.89	0.40
1:A:72:HIS:CE1	5:A:700:HEM:ND	2.90	0.40

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	125/127 (98%)	123 (98%)	1 (1%)	1 (1%)	16 10
1	B	123/127 (97%)	122 (99%)	1 (1%)	0	100 100
All	All	248/254 (98%)	245 (99%)	2 (1%)	1 (0%)	30 26

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	126	GLY

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	112/112 (100%)	104 (93%)	8 (7%)	13 8
1	B	112/112 (100%)	107 (96%)	5 (4%)	24 22

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	224/224 (100%)	211 (94%)	13 (6%)	18 14

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	41	GLU
1	A	57	MET
1	A	89	LEU
1	A	102	MET
1	A	112	MET
1	A	124	TYR
1	A	127	HIS
1	B	5	THR
1	B	7	ASN
1	B	41	GLU
1	B	57	MET
1	B	112	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	7	ASN
1	A	34	ASN
1	A	67	GLN
1	A	80	GLN
1	A	127	HIS
1	B	7	ASN
1	B	8	GLN
1	B	33	ASN
1	B	34	ASN
1	B	108	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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)

11 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	SO4	B	811	-	4,4,4	0.22	0	6,6,6	0.13	0
3	ACT	B	822	-	3,3,3	0.81	0	3,3,3	1.57	0
2	CYN	B	800	-	1,1,1	0.28	0	-		
2	CYN	A	800	-	1,1,1	0.27	0	-		
4	SO4	B	810	-	4,4,4	0.29	0	6,6,6	0.22	0
5	HEM	B	700	1	50,50,50	1.91	11 (22%)	67,82,82	1.49	7 (10%)
3	ACT	B	823	-	3,3,3	0.82	0	3,3,3	1.51	0
4	SO4	A	812	-	4,4,4	0.21	0	6,6,6	0.21	0
4	SO4	B	813	-	4,4,4	0.22	0	6,6,6	0.12	0
3	ACT	A	820	-	3,3,3	0.82	0	3,3,3	1.36	0
5	HEM	A	700	1	50,50,50	1.83	9 (18%)	67,82,82	1.30	5 (7%)

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
5	HEM	B	700	1	-	2/14/54/54	-
5	HEM	A	700	1	-	3/14/54/54	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	700	HEM	C3D-C2D	7.76	1.53	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	700	HEM	C3D-C2D	7.44	1.52	1.36
5	A	700	HEM	FE-NC	4.75	2.10	1.95
5	B	700	HEM	FE-ND	4.55	2.08	1.94
5	B	700	HEM	FE-NC	4.14	2.08	1.95
5	A	700	HEM	FE-ND	3.53	2.05	1.94
5	B	700	HEM	FE-NA	3.36	2.06	1.95
5	B	700	HEM	FE-NB	3.31	2.05	1.94
5	B	700	HEM	CAC-C3C	3.02	1.55	1.47
5	A	700	HEM	FE-NA	3.02	2.05	1.95
5	B	700	HEM	CAB-C3B	2.85	1.55	1.47
5	A	700	HEM	CAB-C3B	2.69	1.54	1.47
5	A	700	HEM	FE-NB	2.68	2.03	1.94
5	A	700	HEM	CAC-C3C	2.67	1.54	1.47
5	A	700	HEM	CMB-C2B	2.43	1.55	1.50
5	B	700	HEM	CMA-C3A	2.30	1.55	1.50
5	A	700	HEM	CMC-C2C	2.25	1.55	1.50
5	B	700	HEM	CAA-C2A	2.16	1.56	1.51
5	B	700	HEM	CMB-C2B	2.08	1.55	1.50
5	B	700	HEM	CMC-C2C	2.08	1.55	1.50

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
5	B	700	HEM	C4D-ND-C1D	5.85	112.14	105.21
5	A	700	HEM	C4D-ND-C1D	5.23	111.41	105.21
5	B	700	HEM	C1B-NB-C4B	3.57	109.44	105.21
5	B	700	HEM	CAD-CBD-CGD	-3.34	104.80	113.67
5	B	700	HEM	C3B-C4B-NB	-3.29	107.10	109.47
5	B	700	HEM	CHA-C4D-ND	2.81	127.84	124.37
5	A	700	HEM	CAD-CBD-CGD	-2.71	106.49	113.67
5	A	700	HEM	CHA-C4D-ND	2.35	127.27	124.37
5	A	700	HEM	CMD-C2D-C1D	2.33	128.67	125.03
5	B	700	HEM	CHD-C4C-NC	2.10	126.74	124.45
5	A	700	HEM	C4C-C3C-C2C	2.09	108.62	106.81
5	B	700	HEM	C2B-C1B-NB	-2.07	107.45	109.84

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	700	HEM	CAD-CBD-CGD-O1D
5	A	700	HEM	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
5	B	700	HEM	CAD-CBD-CGD-O2D
5	B	700	HEM	CAD-CBD-CGD-O1D
5	A	700	HEM	CAA-CBA-CGA-O1A

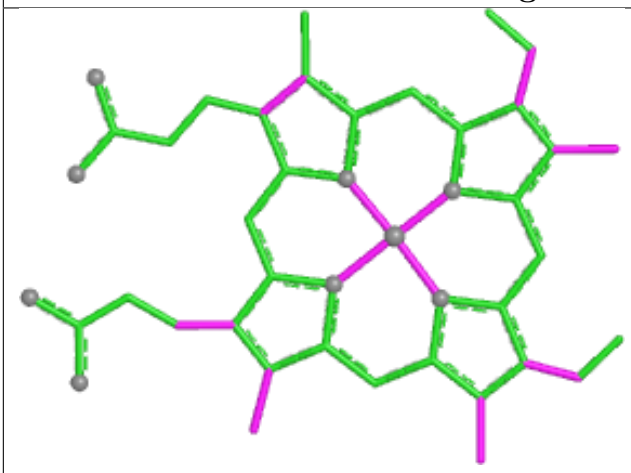
There are no ring outliers.

2 monomers are involved in 3 short contacts:

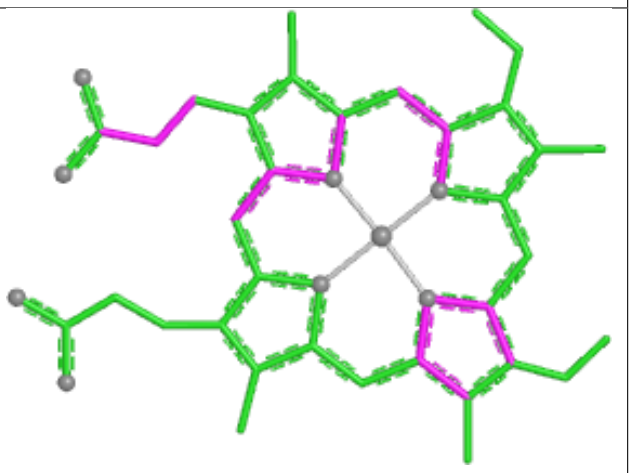
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	700	HEM	2	0
5	A	700	HEM	1	0

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 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.

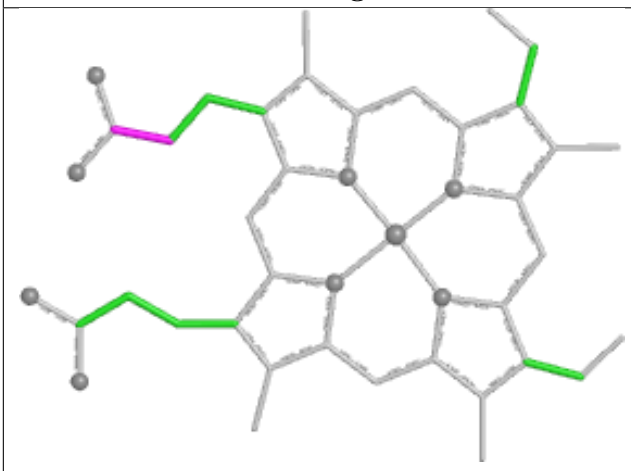
Ligand HEM B 700



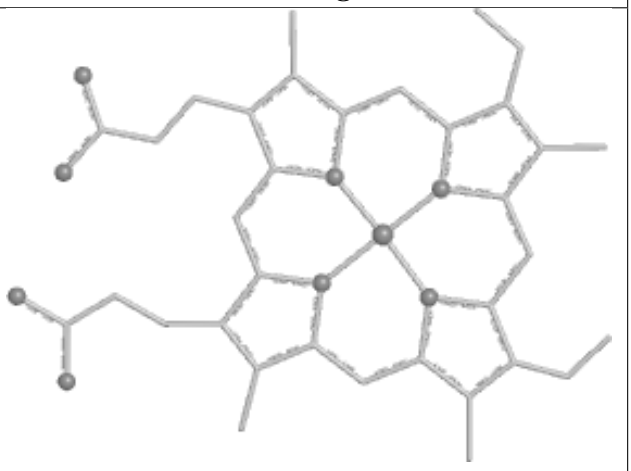
Bond lengths



Bond angles

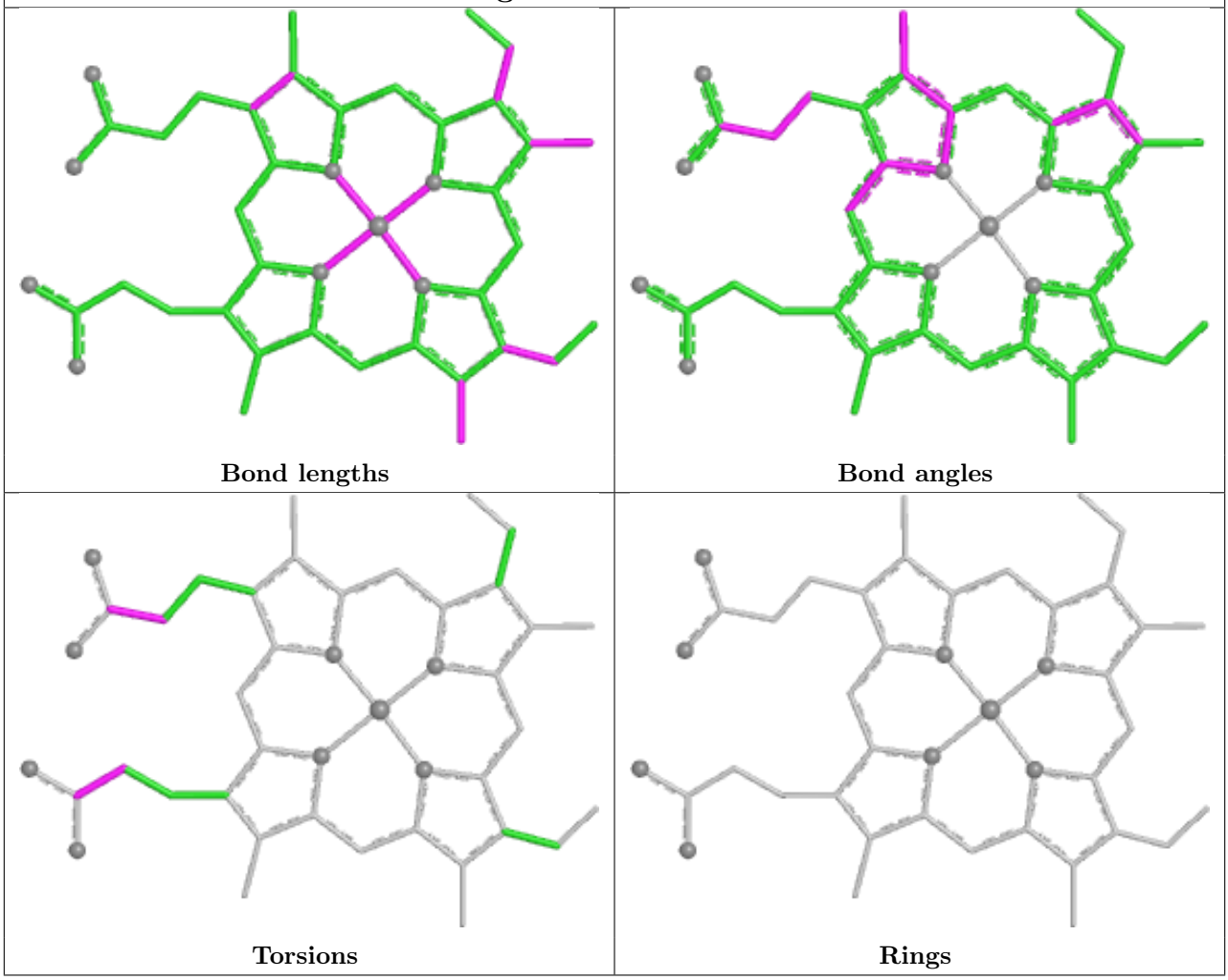


Torsions



Rings

Ligand HEM A 700



5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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, 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	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	127/127 (100%)	0.64	13 (10%) 12 13	24, 33, 52, 63	0
1	B	124/127 (97%)	0.52	6 (4%) 35 40	17, 33, 47, 65	1 (0%)
All	All	251/254 (98%)	0.58	19 (7%) 20 22	17, 33, 48, 65	1 (0%)

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	124	TYR	6.4
1	A	125	GLY	6.0
1	A	127	HIS	5.4
1	B	65	ASN	5.3
1	B	1	MET	5.2
1	A	1	MET	4.0
1	A	126	GLY	3.5
1	B	45	GLU	3.5
1	B	124	TYR	2.9
1	A	65	ASN	2.6
1	B	81	GLU	2.5
1	A	8	GLN	2.5
1	A	115	SER	2.5
1	A	84	GLU	2.3
1	A	2	LYS	2.3
1	B	100	GLU	2.3
1	A	20	GLU	2.2
1	A	34	ASN	2.0
1	A	81	GLU	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 (i)

There are no oligosaccharides in this entry.

6.4 Ligands (i)

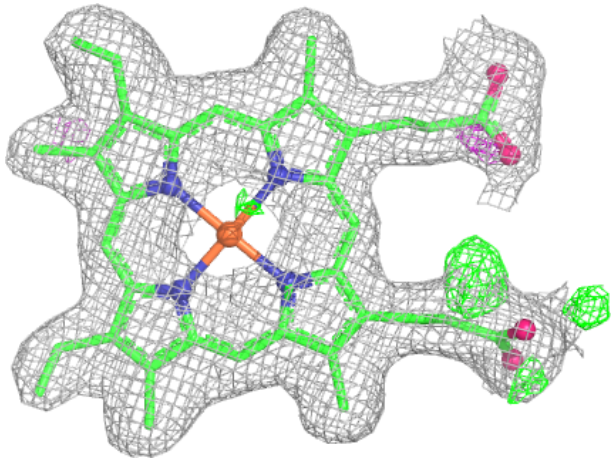
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9'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

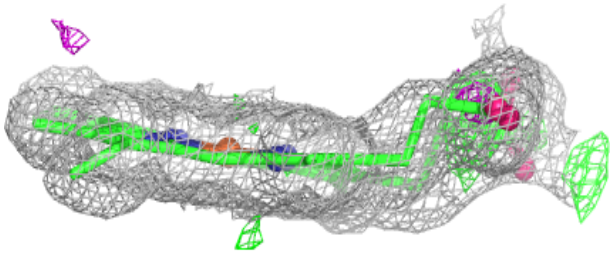
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	ACT	B	822	4/4	0.44	0.27	63,63,63,63	0
3	ACT	A	820	4/4	0.64	0.24	68,68,68,68	0
3	ACT	B	823	4/4	0.81	0.15	59,60,60,60	0
4	SO4	B	813	5/5	0.83	0.13	80,80,81,81	0
2	CYN	A	800	2/2	0.88	0.17	20,20,20,20	0
4	SO4	B	811	5/5	0.89	0.09	81,81,81,81	0
4	SO4	A	812	5/5	0.92	0.14	48,49,50,50	0
2	CYN	B	800	2/2	0.93	0.10	21,21,21,22	0
4	SO4	B	810	5/5	0.96	0.08	35,36,37,38	0
5	HEM	B	700	43/43	0.96	0.09	24,26,33,39	4
5	HEM	A	700	43/43	0.97	0.08	22,24,35,41	0

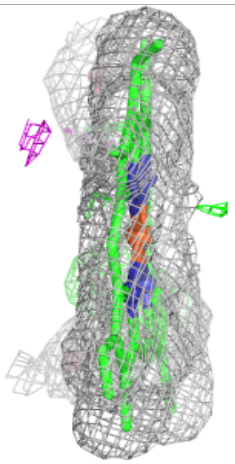
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HEM B 700:

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

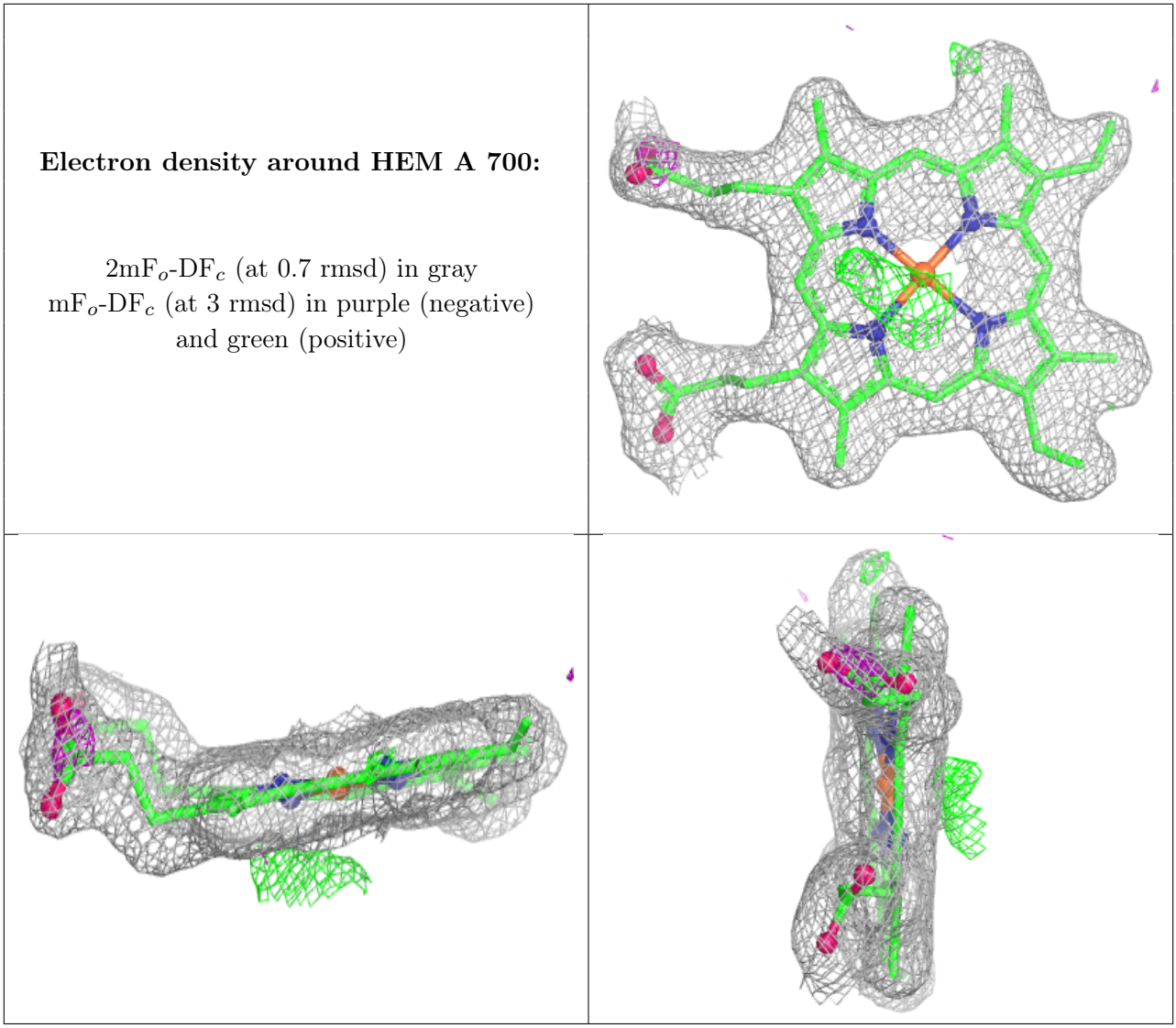






Electron density around HEM A 700:

$2mF_o$-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.