



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 6, 2024 – 03:33 AM EST

PDB ID : 1YMB  
Title : HIGH RESOLUTION STUDY OF THE THREE-DIMENSIONAL STRUCTURE OF HORSE HEART METMYOGLOBIN  
Authors : Evans, S.V.; Brayer, G.D.  
Deposited on : 1993-09-27  
Resolution : 1.90 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtrriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

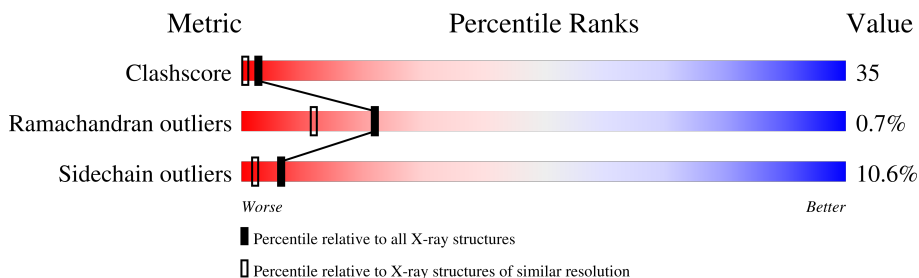
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
Clashscore	141614	6847 (1.90-1.90)
Ramachandran outliers	138981	6760 (1.90-1.90)
Sidechain outliers	138945	6760 (1.90-1.90)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$

Mol	Chain	Length	Quality of chain
1	A	153	 41% 44% 13% .

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1401 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 METMYOGLOBIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	153	1199	769	210	218	2	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



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

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C<sub>34</sub>H<sub>32</sub>FeN<sub>4</sub>O<sub>4</sub>).



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

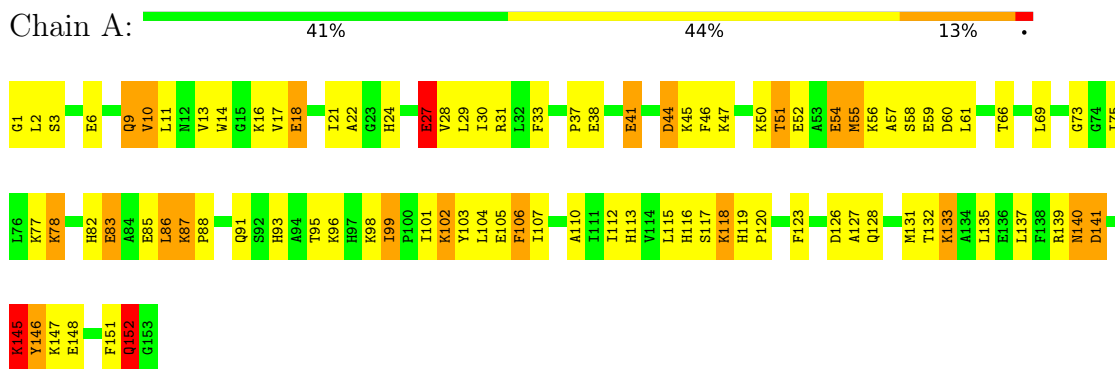
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	154	Total	O	0	0
			154	154		

### 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. 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. 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: METMYOGLOBIN



## 4 Data and refinement statistics i

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	64.27Å 28.92Å 35.93Å 90.00° 107.10° 90.00°	Depositor
Resolution (Å)	(Not available) – 1.90 21.06 – 1.90	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-1.90) 85.4 (21.06-1.90)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.14 (at 1.90Å)	Xtrriage
Refinement program	PROLSQ	Depositor
R, $R_{free}$	0.155 , (Not available) 0.222 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	12.9	Xtrriage
Anisotropy	0.211	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 91.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.46$ , $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	0.045 for -h-l,-k,l	Xtrriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	1401	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	16.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.35% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality i

### 5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	1.39	2/1226 (0.2%)	2.06	41/1645 (2.5%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	85	GLU	CD-OE1	-6.40	1.18	1.25
1	A	18	GLU	CD-OE1	5.25	1.31	1.25

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	31	ARG	NE-CZ-NH1	14.78	127.69	120.30
1	A	31	ARG	NE-CZ-NH2	-12.37	114.12	120.30
1	A	103	TYR	CB-CG-CD1	10.14	127.09	121.00
1	A	102	LYS	CB-CG-CD	9.15	135.39	111.60
1	A	86	LEU	CB-CA-C	8.19	125.76	110.20
1	A	140	ASN	CA-CB-CG	7.61	130.14	113.40
1	A	103	TYR	CB-CG-CD2	-7.46	116.53	121.00
1	A	106	PHE	CB-CG-CD1	7.25	125.88	120.80
1	A	148	GLU	CA-CB-CG	7.06	128.94	113.40
1	A	41	GLU	CA-CB-CG	6.80	128.37	113.40
1	A	27	GLU	CG-CD-OE2	6.79	131.88	118.30
1	A	29	LEU	CB-CA-C	6.43	122.41	110.20
1	A	147	LYS	O-C-N	6.37	132.89	122.70
1	A	132	THR	CA-CB-CG2	6.27	121.18	112.40
1	A	44	ASP	C-N-CA	6.20	137.20	121.70
1	A	38	GLU	OE1-CD-OE2	-6.17	115.89	123.30
1	A	10	VAL	CA-CB-CG2	6.17	120.16	110.90
1	A	58	SER	N-CA-CB	-6.12	101.33	110.50
1	A	55	MET	CG-SD-CE	6.06	109.90	100.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	99	ILE	CB-CG1-CD1	6.04	130.82	113.90
1	A	106	PHE	CB-CG-CD2	-6.03	116.58	120.80
1	A	22	ALA	CB-CA-C	5.98	119.06	110.10
1	A	37	PRO	N-CD-CG	-5.98	94.23	103.20
1	A	146	TYR	CB-CG-CD1	5.78	124.47	121.00
1	A	69	LEU	CB-CG-CD1	-5.62	101.44	111.00
1	A	145	LYS	CA-CB-CG	5.62	125.76	113.40
1	A	93	HIS	CA-CB-CG	-5.60	104.08	113.60
1	A	10	VAL	CG1-CB-CG2	-5.59	101.96	110.90
1	A	59	GLU	OE1-CD-OE2	5.52	129.93	123.30
1	A	51	THR	CA-C-O	5.32	131.27	120.10
1	A	141	ASP	CB-CG-OD1	5.31	123.08	118.30
1	A	44	ASP	CB-CG-OD1	5.30	123.07	118.30
1	A	139	ARG	NE-CZ-NH2	-5.30	117.65	120.30
1	A	145	LYS	CB-CG-CD	5.27	125.30	111.60
1	A	46	PHE	CB-CG-CD1	5.22	124.45	120.80
1	A	27	GLU	CG-CD-OE1	-5.17	107.96	118.30
1	A	11	LEU	CB-CA-C	5.15	119.99	110.20
1	A	113	HIS	CA-C-O	5.10	130.80	120.10
1	A	41	GLU	CG-CD-OE1	5.03	128.37	118.30
1	A	105	GLU	CA-C-O	5.03	130.67	120.10
1	A	91	GLN	CG-CD-OE1	-5.02	111.56	121.60

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	1199	0	1211	87	0
2	A	5	0	0	1	0
3	A	43	0	30	3	0
4	A	154	0	0	36	1
All	All	1401	0	1241	88	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including



hydrogen atoms). The all-atom clashscore for this structure is 35.

All (88) 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:1:GLY:HA2	4:A:276:HOH:O	1.61	1.00
1:A:98:LYS:HG3	4:A:225:HOH:O	1.76	0.85
1:A:145:LYS:HE3	4:A:218:HOH:O	1.76	0.85
1:A:21:ILE:HD12	1:A:66:THR:HG23	1.59	0.84
1:A:45:LYS:HE3	4:A:173:HOH:O	1.87	0.74
1:A:78:LYS:CG	4:A:261:HOH:O	2.36	0.73
1:A:13:VAL:HG13	1:A:16:LYS:HZ2	1.53	0.73
1:A:78:LYS:HG2	4:A:261:HOH:O	1.86	0.73
1:A:9:GLN:HG3	4:A:165:HOH:O	1.89	0.73
1:A:44:ASP:HA	1:A:47:LYS:HD2	1.71	0.70
1:A:98:LYS:HB2	4:A:237:HOH:O	1.93	0.69
1:A:98:LYS:HD3	1:A:151:PHE:HE1	1.58	0.69
1:A:21:ILE:CD1	1:A:66:THR:HG23	2.23	0.68
1:A:98:LYS:HD3	1:A:151:PHE:CE1	2.29	0.68
1:A:3:SER:HB2	4:A:284:HOH:O	1.95	0.66
1:A:98:LYS:CB	4:A:237:HOH:O	2.44	0.65
1:A:98:LYS:HE3	4:A:292:HOH:O	1.95	0.65
1:A:96:LYS:HE3	4:A:248:HOH:O	1.97	0.65
1:A:98:LYS:CE	4:A:292:HOH:O	2.46	0.63
1:A:75:ILE:O	1:A:78:LYS:HB2	1.99	0.63
1:A:119:HIS:N	1:A:120:PRO:HD3	2.14	0.63
1:A:123:PHE:HA	4:A:234:HOH:O	1.98	0.62
1:A:101:ILE:HG12	1:A:146:TYR:CD2	2.35	0.62
1:A:96:LYS:CE	4:A:248:HOH:O	2.48	0.62
1:A:118:LYS:C	1:A:120:PRO:HD3	2.21	0.61
1:A:3:SER:OG	1:A:6:GLU:HG3	2.02	0.59
1:A:126:ASP:HB2	4:A:227:HOH:O	2.02	0.59
1:A:30:ILE:HD13	1:A:56:LYS:HG3	1.83	0.59
1:A:131:MET:HA	4:A:264:HOH:O	2.03	0.58
1:A:78:LYS:HE2	4:A:158:HOH:O	2.02	0.58
1:A:99:ILE:HD12	3:A:154:HEM:HAC	1.86	0.57
1:A:101:ILE:H	1:A:152:GLN:HE22	1.52	0.57
1:A:98:LYS:NZ	4:A:292:HOH:O	2.36	0.57
1:A:83:GLU:H	1:A:83:GLU:CD	2.09	0.56
1:A:88:PRO:HD3	4:A:199:HOH:O	2.05	0.55
1:A:78:LYS:CE	4:A:261:HOH:O	2.53	0.55
1:A:82:HIS:HE1	1:A:137:LEU:HD23	1.70	0.55
1:A:51:THR:OG1	1:A:54:GLU:HB2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:86:LEU:HD21	1:A:141:ASP:HB3	1.88	0.54
1:A:13:VAL:HG21	1:A:127:ALA:HB1	1.89	0.53
3:A:154:HEM:CGA	4:A:283:HOH:O	2.55	0.53
1:A:95:THR:HA	1:A:98:LYS:HE2	1.90	0.53
1:A:9:GLN:HG2	1:A:127:ALA:HA	1.91	0.52
1:A:2:LEU:HD12	1:A:133:LYS:HZ2	1.75	0.51
1:A:118:LYS:HG3	4:A:279:HOH:O	2.09	0.51
1:A:10:VAL:HG22	4:A:264:HOH:O	2.09	0.51
1:A:44:ASP:CA	1:A:47:LYS:HD2	2.42	0.49
1:A:3:SER:HB3	4:A:270:HOH:O	2.12	0.49
1:A:73:GLY:O	1:A:77:LYS:HG3	2.12	0.49
1:A:82:HIS:HE1	1:A:137:LEU:CD2	2.25	0.49
1:A:14:TRP:O	1:A:18:GLU:HG3	2.12	0.49
1:A:16:LYS:HE3	4:A:229:HOH:O	2.13	0.49
1:A:60:ASP:HB2	2:A:155:SO4:O4	2.13	0.49
1:A:101:ILE:H	1:A:152:GLN:NE2	2.10	0.49
1:A:131:MET:HE2	4:A:264:HOH:O	2.13	0.49
1:A:146:TYR:HD2	1:A:152:GLN:HE21	1.59	0.49
1:A:33:PHE:HE2	1:A:61:LEU:HD11	1.78	0.48
1:A:146:TYR:HD2	1:A:152:GLN:NE2	2.12	0.47
1:A:99:ILE:CD1	3:A:154:HEM:HAC	2.44	0.46
1:A:2:LEU:HA	1:A:6:GLU:OE1	2.14	0.46
1:A:116:HIS:NE2	1:A:128:GLN:NE2	2.64	0.46
1:A:52:GLU:HA	1:A:55:MET:HE3	1.97	0.46
1:A:101:ILE:HD13	1:A:104:LEU:HD12	1.98	0.46
1:A:95:THR:HG21	4:A:190:HOH:O	2.16	0.45
1:A:51:THR:O	1:A:55:MET:HE2	2.15	0.45
1:A:52:GLU:O	1:A:55:MET:HB2	2.17	0.45
1:A:107:ILE:O	1:A:110:ALA:HB3	2.17	0.45
1:A:112:ILE:HG12	4:A:198:HOH:O	2.17	0.45
1:A:14:TRP:CH2	1:A:17:VAL:HG11	2.52	0.44
1:A:123:PHE:HD1	4:A:234:HOH:O	1.99	0.44
1:A:133:LYS:HD2	4:A:161:HOH:O	2.17	0.44
1:A:117:SER:O	1:A:120:PRO:HG3	2.17	0.44
1:A:52:GLU:HA	1:A:55:MET:CE	2.47	0.44
1:A:151:PHE:C	1:A:152:GLN:HG3	2.39	0.43
1:A:10:VAL:HA	4:A:264:HOH:O	2.19	0.43
1:A:10:VAL:HG13	4:A:264:HOH:O	2.19	0.42
1:A:115:LEU:HD12	1:A:123:PHE:CZ	2.54	0.42
1:A:112:ILE:HD11	1:A:135:LEU:HD12	2.00	0.42
1:A:57:ALA:HB1	4:A:170:HOH:O	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:LYS:HD2	4:A:222:HOH:O	2.20	0.41
1:A:24:HIS:O	1:A:28:VAL:HG23	2.20	0.41
1:A:33:PHE:CE2	1:A:61:LEU:HD11	2.54	0.41
1:A:27:GLU:CD	1:A:118:LYS:HE3	2.40	0.41
1:A:146:TYR:CD2	1:A:152:GLN:NE2	2.88	0.41
1:A:2:LEU:HD11	1:A:133:LYS:HG2	2.02	0.41
1:A:87:LYS:NZ	4:A:282:HOH:O	2.53	0.40
1:A:6:GLU:OE1	1:A:133:LYS:HE3	2.20	0.40
1:A:102:LYS:HE3	1:A:106:PHE:CE1	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:168:HOH:O	4:A:309:HOH:O[2_656]	1.98	0.22

## 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	151/153 (99%)	137 (91%)	13 (9%)	1 (1%)	22 12

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	152	GLN

### 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	123/123 (100%)	110 (89%)	13 (11%)	<b>6</b> <b>2</b>

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

Mol	Chain	Res	Type
1	A	9	GLN
1	A	27	GLU
1	A	41	GLU
1	A	50	LYS
1	A	54	GLU
1	A	78	LYS
1	A	83	GLU
1	A	87	LYS
1	A	118	LYS
1	A	133	LYS
1	A	140	ASN
1	A	145	LYS
1	A	152	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	26	GLN
1	A	82	HIS
1	A	128	GLN
1	A	140	ASN

### 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](#)

2 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	A	154	4,1	41,50,50	1.62	8 (19%)	45,82,82	1.79	11 (24%)
2	SO4	A	155	-	4,4,4	0.88	0	6,6,6	0.76	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
3	HEM	A	154	4,1	-	5/12/54/54	-

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	154	HEM	C3C-C2C	-5.14	1.33	1.40
3	A	154	HEM	CAB-C3B	2.77	1.55	1.47
3	A	154	HEM	C3C-CAC	2.53	1.53	1.47
3	A	154	HEM	C2C-C1C	2.52	1.48	1.42
3	A	154	HEM	C3B-C2B	-2.46	1.32	1.37
3	A	154	HEM	CMD-C2D	2.21	1.55	1.50
3	A	154	HEM	CMB-C2B	2.16	1.55	1.50
3	A	154	HEM	CHA-C4D	2.08	1.40	1.35

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	154	HEM	O1A-CGA-CBA	-4.31	109.25	123.08
3	A	154	HEM	CHC-C4B-NB	4.27	129.07	124.43
3	A	154	HEM	CMA-C3A-C4A	-4.16	122.07	128.46
3	A	154	HEM	CMC-C2C-C3C	3.48	131.19	124.68
3	A	154	HEM	CMA-C3A-C2A	3.13	130.84	124.94
3	A	154	HEM	C4C-CHD-C1D	2.49	125.85	122.56
3	A	154	HEM	O2A-CGA-CBA	2.39	121.72	114.03
3	A	154	HEM	C3B-C2B-C1B	2.29	108.19	106.49
3	A	154	HEM	C1B-NB-C4B	2.23	107.38	105.07
3	A	154	HEM	O2A-CGA-O1A	2.19	128.76	123.30
3	A	154	HEM	CHD-C1D-C2D	2.04	128.16	124.98

There are no chirality outliers.

All (5) torsion outliers are listed below:

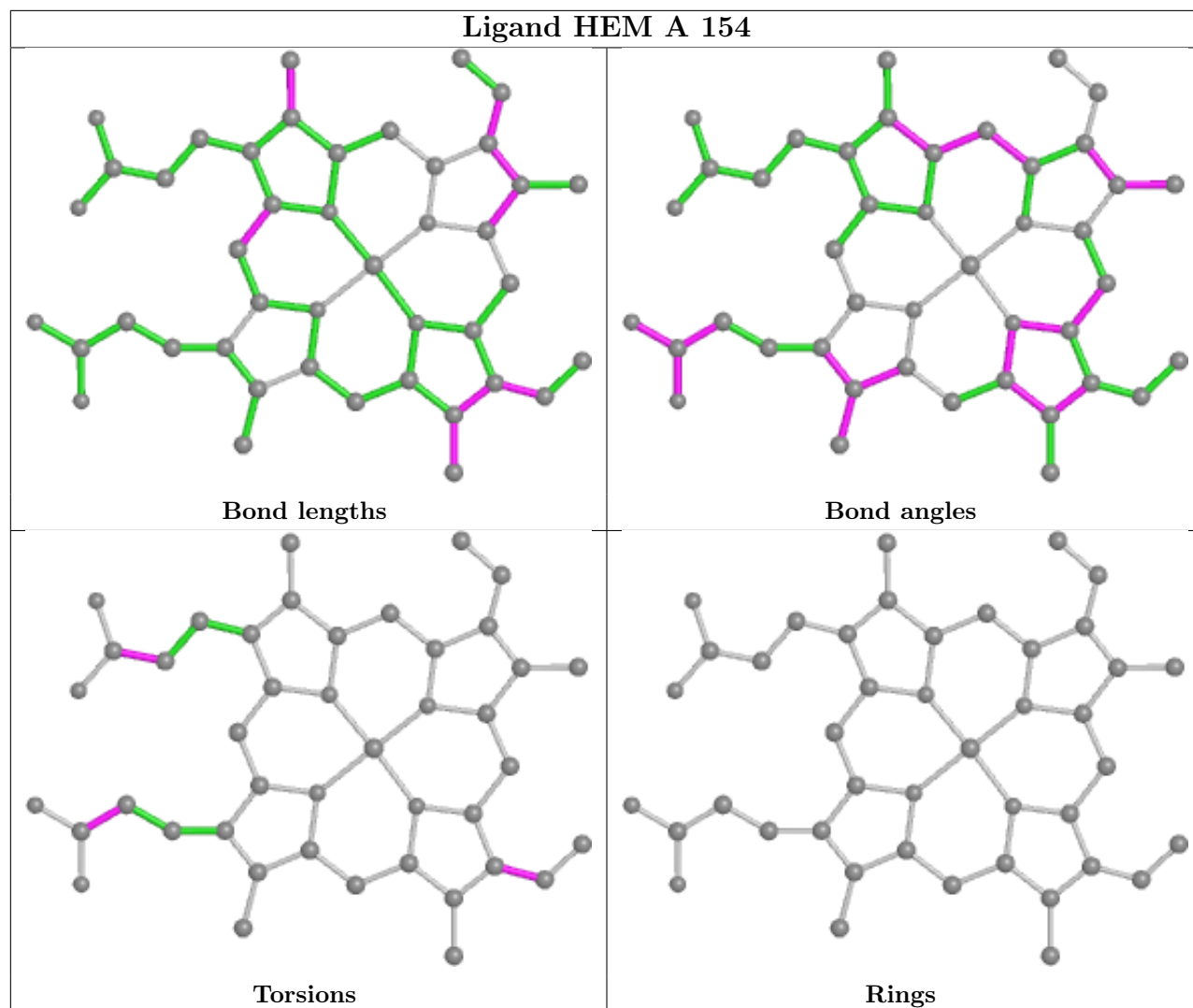
Mol	Chain	Res	Type	Atoms
3	A	154	HEM	C2B-C3B-CAB-CBB
3	A	154	HEM	C4B-C3B-CAB-CBB
3	A	154	HEM	CAA-CBA-CGA-O2A
3	A	154	HEM	CAD-CBD-CGD-O2D
3	A	154	HEM	CAA-CBA-CGA-O1A

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	154	HEM	3	0
2	A	155	SO4	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 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

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.2 Non-standard residues in protein, DNA, RNA chains

Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.3 Carbohydrates

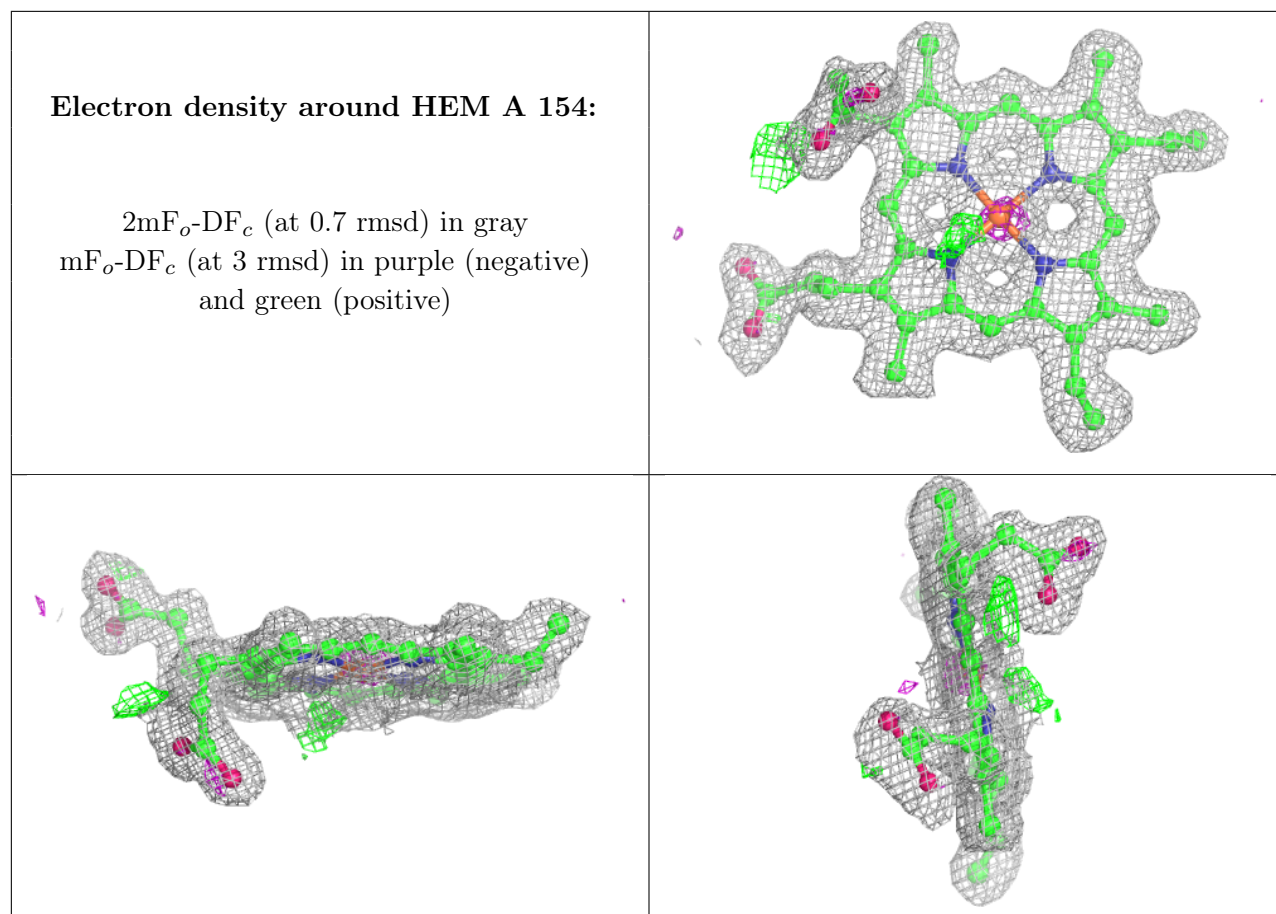
Unable to reproduce the depositors R factor - this section is therefore empty.

### 6.4 Ligands

Unable to reproduce the depositors R factor - this section is therefore empty.

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](#)

Unable to reproduce the depositor's R factor - this section is therefore empty.