



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 10:00 am GMT

PDB ID : 3B98  
Title : Crystal structure of zebrafish prostacyclin synthase (cytochrome P450 8A1)  
Authors : Li, Y.-C.; Chiang, C.-W.; Yeh, H.-C.; Hsu, P.-Y.; Whitby, F.G.; Wang, L.-H.;  
Chan, N.-L.  
Deposited on : 2007-11-03  
Resolution : 2.08 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : recalc28949

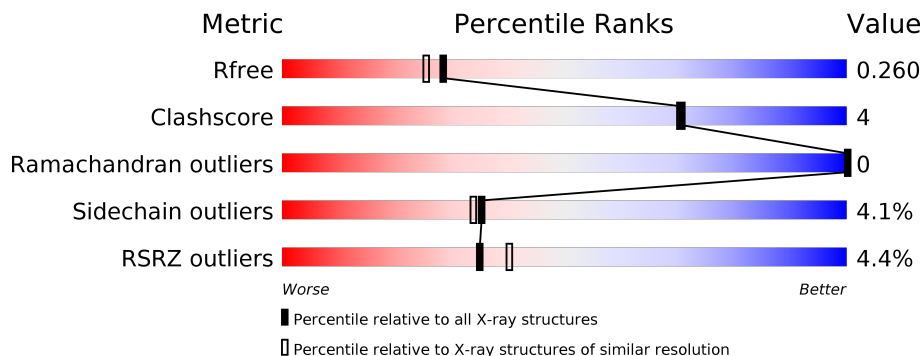
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.08 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))
$R_{free}$	100719	4955 (2.10-2.06)
Clashscore	112137	5537 (2.10-2.06)
Ramachandran outliers	110173	5483 (2.10-2.06)
Sidechain outliers	110143	5484 (2.10-2.06)
RSRZ outliers	101464	4991 (2.10-2.06)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	475	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%;"></div> <div style="position: absolute; bottom: 0; left: 0; width: 100%; height: 10px;"> <div style="width: 1%; height: 10px; background-color: red;"></div> <div style="width: 83%, height: 10px; background-color: green;"></div> <div style="width: 10%, height: 10px; background-color: yellow;"></div> <div style="width: 6%, height: 10px; background-color: orange;"></div> <div style="width: 2%, height: 10px; background-color: grey;"></div> </div> </div> </div>
1	B	475	<div> <div style="width: 100%; height: 10px; position: relative;"> <div style="position: absolute; top: -10px; left: 0; width: 100%;"></div> <div style="position: absolute; bottom: 0; left: 0; width: 100%; height: 10px;"> <div style="width: 7%, height: 10px; background-color: red;"></div> <div style="width: 80%, height: 10px; background-color: green;"></div> <div style="width: 10%, height: 10px; background-color: yellow;"></div> <div style="width: 3%, height: 10px; background-color: orange;"></div> <div style="width: 0%, height: 10px; background-color: grey;"></div> </div> </div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7443 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.

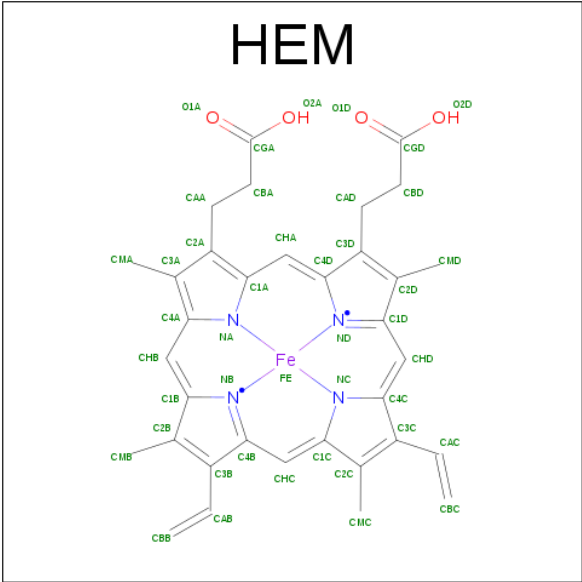
- Molecule 1 is a protein called Prostaglandin I2 synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	0	0	0
			3613	2300	641	653	19			
1	B	427	Total	C	N	O	S	0	0	0
			3462	2214	611	618	19			

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	10	MET	-	INITIATING METHIONINE	UNP A9LLA5
A	11	ALA	-	EXPRESSION TAG	UNP A9LLA5
A	12	LYS	-	EXPRESSION TAG	UNP A9LLA5
A	13	LYS	-	EXPRESSION TAG	UNP A9LLA5
A	14	THR	-	EXPRESSION TAG	UNP A9LLA5
A	15	SER	-	EXPRESSION TAG	UNP A9LLA5
A	16	SER	-	EXPRESSION TAG	UNP A9LLA5
A	481	HIS	-	EXPRESSION TAG	UNP A9LLA5
A	482	HIS	-	EXPRESSION TAG	UNP A9LLA5
A	483	HIS	-	EXPRESSION TAG	UNP A9LLA5
A	484	HIS	-	EXPRESSION TAG	UNP A9LLA5
B	10	MET	-	INITIATING METHIONINE	UNP A9LLA5
B	11	ALA	-	EXPRESSION TAG	UNP A9LLA5
B	12	LYS	-	EXPRESSION TAG	UNP A9LLA5
B	13	LYS	-	EXPRESSION TAG	UNP A9LLA5
B	14	THR	-	EXPRESSION TAG	UNP A9LLA5
B	15	SER	-	EXPRESSION TAG	UNP A9LLA5
B	16	SER	-	EXPRESSION TAG	UNP A9LLA5
B	481	HIS	-	EXPRESSION TAG	UNP A9LLA5
B	482	HIS	-	EXPRESSION TAG	UNP A9LLA5
B	483	HIS	-	EXPRESSION TAG	UNP A9LLA5
B	484	HIS	-	EXPRESSION TAG	UNP A9LLA5

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula:  $C_{34}H_{32}FeN_4O_4$ ).



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

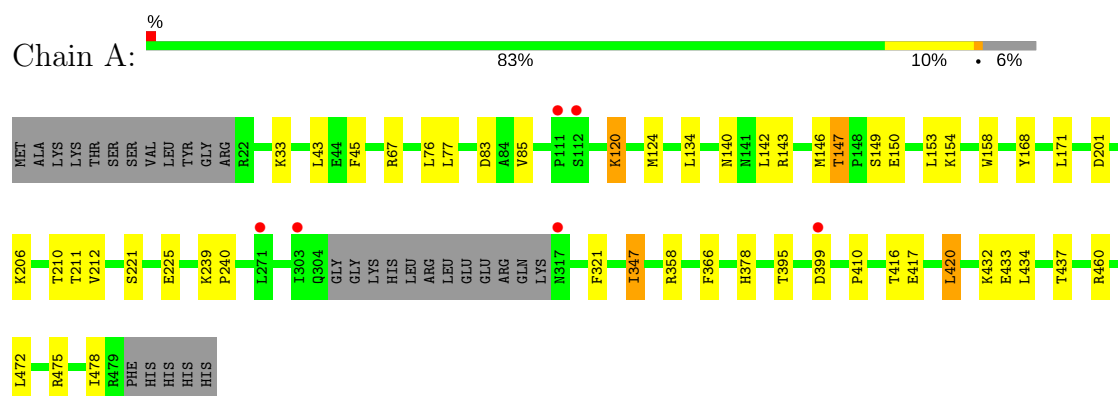
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	198	Total	O	0	0
			198	198		
3	B	84	Total	O	0	0
			84	84		

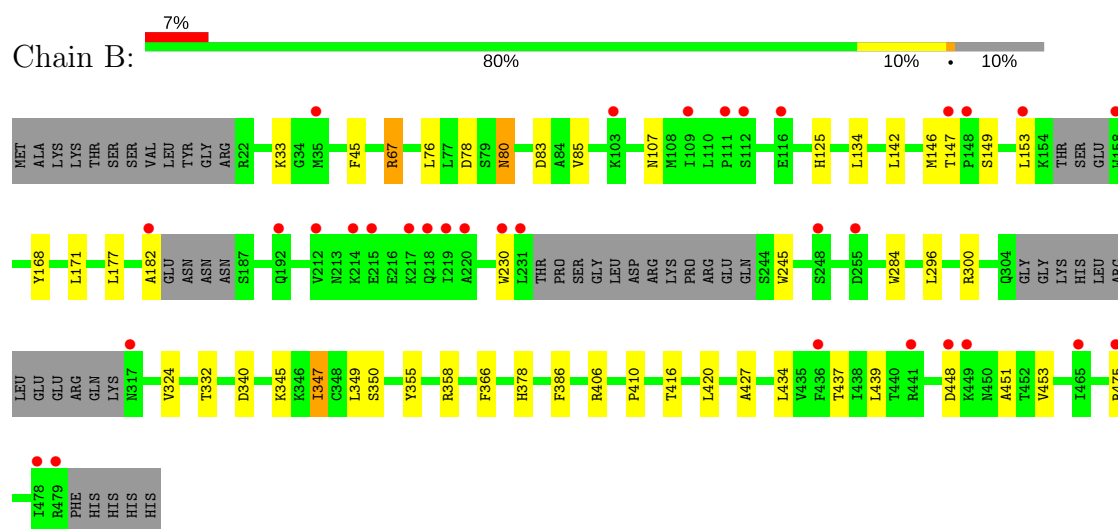
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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: Prostaglandin I2 synthase



#### • Molecule 1: Prostaglandin I2 synthase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	58.67Å 87.90Å 190.85Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.08 28.99 – 2.08	Depositor EDS
% Data completeness (in resolution range)	99.8 (30.00-2.08) 99.9 (28.99-2.08)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	0.05	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.58 (at 2.08Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, $R_{free}$	0.227 , 0.264 0.229 , 0.260	Depositor DCC
$R_{free}$ test set	3031 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	27.0	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 40.9	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.93	EDS
Total number of atoms	7443	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.58% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.42	0/3695	0.57	0/4993
1	B	0.38	0/3539	0.53	0/4777
All	All	0.40	0/7234	0.55	0/9770

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	3613	0	3614	30	0
1	B	3462	0	3472	25	0
2	A	43	0	30	2	0
2	B	43	0	30	2	0
3	A	198	0	0	4	0
3	B	84	0	0	0	0
All	All	7443	0	7146	58	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 4.

All (58) 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:146:MET:CE	1:A:437:THR:HG21	1.98	0.94
1:A:146:MET:HE3	1:A:437:THR:HG21	1.63	0.79
1:A:146:MET:HE1	1:A:437:THR:HG21	1.68	0.72
1:B:142:LEU:HG	1:B:146:MET:HE2	1.72	0.72
1:B:85:VAL:HG22	1:B:347:ILE:HG12	1.74	0.69
1:A:83:ASP:OD1	1:A:416:THR:HG23	1.94	0.67
2:B:600:HEM:HHC	2:B:600:HEM:HBB2	1.78	0.66
1:B:146:MET:HE3	1:B:437:THR:HG21	1.78	0.65
1:A:201:ASP:O	1:A:460:ARG:NH2	2.24	0.65
1:B:78:ASP:OD1	1:B:80:ASN:ND2	2.32	0.63
1:A:147:THR:HG22	1:A:150:GLU:H	1.64	0.62
1:B:378:HIS:HE1	1:B:410:PRO:O	1.83	0.62
1:A:146:MET:CE	3:A:711:HOH:O	2.49	0.60
1:A:146:MET:HE1	3:A:711:HOH:O	2.01	0.60
1:A:210:THR:O	1:A:210:THR:HG22	2.01	0.60
1:A:149:SER:HA	1:A:154:LYS:HD2	1.84	0.60
1:A:211:THR:HG22	1:A:211:THR:O	2.01	0.59
1:B:300:ARG:NH2	1:B:439:LEU:O	2.40	0.55
1:B:33:LYS:O	1:B:67:ARG:HD2	2.08	0.53
1:A:33:LYS:O	1:A:67:ARG:HD2	2.09	0.53
1:B:142:LEU:HG	1:B:146:MET:CE	2.39	0.52
1:A:378:HIS:HE1	1:A:410:PRO:O	1.93	0.52
1:B:146:MET:CE	1:B:437:THR:HG21	2.40	0.52
2:A:600:HEM:HMB2	2:A:600:HEM:HBB2	1.91	0.51
1:B:448:ASP:HB2	1:B:451:ALA:HB2	1.92	0.51
2:A:600:HEM:CMB	2:A:600:HEM:HBB2	2.41	0.51
1:A:410:PRO:HG2	3:A:705:HOH:O	2.11	0.50
1:A:85:VAL:HG22	1:A:347:ILE:HG12	1.94	0.50
1:A:321:PHE:HD2	1:A:432:LYS:HZ3	1.58	0.49
1:B:284:TRP:CG	1:B:332:THR:HG21	2.49	0.48
1:B:125:HIS:CG	1:B:245:TRP:HB2	2.48	0.48
1:A:221:SER:O	1:A:225:GLU:HG3	2.14	0.47
1:A:206:LYS:HB3	1:A:212:VAL:HG13	1.95	0.47
1:A:211:THR:O	1:A:211:THR:CG2	2.61	0.47
1:A:158:TRP:CE3	1:A:475:ARG:HG3	2.49	0.47
1:B:324:VAL:HG22	1:B:386:PHE:HB2	1.97	0.47
1:B:296:LEU:O	1:B:300:ARG:HG3	2.14	0.47
1:A:171:LEU:HG	1:A:434:LEU:HD22	1.96	0.46
1:B:171:LEU:HG	1:B:434:LEU:HD22	1.97	0.46
1:B:147:THR:HG22	1:B:149:SER:H	1.80	0.46
1:B:80:ASN:HD21	1:B:350:SER:HB3	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:177:LEU:HD23	1:B:182:ALA:HB2	1.98	0.45
1:B:80:ASN:ND2	1:B:350:SER:HB3	2.31	0.45
1:A:153:LEU:HD13	1:A:478:ILE:HG13	1.98	0.45
1:A:321:PHE:CE2	1:A:432:LYS:HG2	2.51	0.44
1:A:416:THR:HG22	3:A:794:HOH:O	2.16	0.43
1:A:120:LYS:O	1:A:124:MET:HG2	2.18	0.43
1:B:85:VAL:HG13	1:B:347:ILE:HD11	1.99	0.43
1:A:140:ASN:OD1	1:A:143:ARG:NH2	2.45	0.43
1:A:239:LYS:HD3	1:A:240:PRO:HD2	2.01	0.43
1:A:417:GLU:O	1:A:420:LEU:HD22	2.19	0.42
1:B:349:LEU:HD22	1:B:355:TYR:CE1	2.53	0.42
1:B:125:HIS:CD2	1:B:134:LEU:HD11	2.55	0.42
1:B:427:ALA:HB2	2:B:600:HEM:HMC2	2.02	0.42
1:A:85:VAL:HG13	1:A:347:ILE:HD11	2.02	0.41
1:A:142:LEU:HD23	1:A:433:GLU:HG3	2.03	0.41
1:B:345:LYS:HE3	1:B:347:ILE:HD11	2.03	0.40
1:B:83:ASP:OD1	1:B:416:THR:HG23	2.22	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	442/475 (93%)	434 (98%)	8 (2%)	0	100	100
1	B	417/475 (88%)	414 (99%)	3 (1%)	0	100	100
All	All	859/950 (90%)	848 (99%)	11 (1%)	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	389/414 (94%)	374 (96%)	15 (4%)	37	36
1	B	371/414 (90%)	355 (96%)	16 (4%)	33	31
All	All	760/828 (92%)	729 (96%)	31 (4%)	35	34

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	LEU
1	A	45	PHE
1	A	76	LEU
1	A	77	LEU
1	A	120	LYS
1	A	134	LEU
1	A	147	THR
1	A	168	TYR
1	A	347	ILE
1	A	358	ARG
1	A	366	PHE
1	A	395	THR
1	A	399	ASP
1	A	420	LEU
1	A	472	LEU
1	B	45	PHE
1	B	67	ARG
1	B	76	LEU
1	B	80	ASN
1	B	107	ASN
1	B	153	LEU
1	B	168	TYR
1	B	230	TRP
1	B	340	ASP
1	B	347	ILE
1	B	358	ARG
1	B	366	PHE

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Mol	Chain	Res	Type
1	B	406	ARG
1	B	420	LEU
1	B	453	VAL
1	B	475	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	139	GLN
1	A	184	ASN
1	A	185	ASN
1	A	378	HIS
1	A	385	GLN
1	A	467	GLN
1	B	80	ASN
1	B	376	GLN
1	B	378	HIS
1	B	380	GLN
1	B	385	GLN
1	B	467	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 [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
2	HEM	A	600	1,3	28,50,50	2.13	8 (28%)	17,82,82	1.53	5 (29%)
2	HEM	B	600	1	28,50,50	2.27	6 (21%)	17,82,82	1.43	3 (17%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	600	1,3	-	0/6/54/54	0/0/8/8
2	HEM	B	600	1	-	0/6/54/54	0/0/8/8

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	600	HEM	C3B-C2B	-4.72	1.34	1.40
2	B	600	HEM	C3C-C2C	-3.98	1.35	1.40
2	A	600	HEM	C3B-C2B	-3.90	1.35	1.40
2	A	600	HEM	C3C-C2C	-3.63	1.35	1.40
2	A	600	HEM	CMB-C2B	2.03	1.55	1.51
2	A	600	HEM	CAA-C2A	2.58	1.56	1.52
2	A	600	HEM	C4D-ND	2.69	1.39	1.36
2	B	600	HEM	C4D-ND	2.73	1.40	1.36
2	A	600	HEM	C3B-CAB	3.45	1.54	1.47
2	B	600	HEM	C3B-CAB	3.77	1.55	1.47
2	A	600	HEM	C3C-CAC	3.94	1.55	1.47
2	B	600	HEM	C3C-CAC	4.20	1.56	1.47
2	A	600	HEM	C3D-C2D	5.30	1.53	1.37
2	B	600	HEM	C3D-C2D	5.53	1.54	1.37

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	HEM	C1D-C2D-C3D	-2.53	105.23	107.00
2	B	600	HEM	CMA-C3A-C4A	-2.31	124.92	128.46
2	A	600	HEM	C1D-C2D-C3D	-2.30	105.39	107.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	600	HEM	CBD-CAD-C3D	-2.25	108.18	112.47
2	A	600	HEM	CMA-C3A-C4A	-2.24	125.02	128.46
2	A	600	HEM	CAD-CBD-CGD	-2.13	109.02	112.66
2	A	600	HEM	C4A-C3A-C2A	2.21	108.53	107.00
2	A	600	HEM	CMC-C2C-C3C	2.43	129.40	124.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	600	HEM	2	0
2	B	600	HEM	2	0

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	446/475 (93%)	-0.05	6 (1%) 77 80	10, 21, 41, 55	0
1	B	427/475 (89%)	0.33	32 (7%) 15 18	15, 32, 57, 73	0
All	All	873/950 (91%)	0.13	38 (4%) 35 40	10, 26, 53, 73	0

All (38) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	215	GLU	4.7
1	B	218	GLN	3.8
1	B	147	THR	3.7
1	B	248	SER	3.7
1	B	449	LYS	3.7
1	B	182	ALA	3.6
1	B	112	SER	3.2
1	B	153	LEU	3.1
1	B	478	ILE	2.9
1	A	112	SER	2.8
1	B	148	PRO	2.7
1	B	441	ARG	2.7
1	B	219	ILE	2.7
1	B	192	GLN	2.7
1	B	212	VAL	2.6
1	B	317	ASN	2.6
1	B	214	LYS	2.5
1	B	111	PRO	2.5
1	A	111	PRO	2.4
1	A	399	ASP	2.4
1	B	479	ARG	2.4
1	B	109	ILE	2.3
1	B	158	TRP	2.3
1	B	465	ILE	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	448	ASP	2.2
1	B	230	TRP	2.1
1	B	103	LYS	2.1
1	A	317	ASN	2.1
1	B	231	LEU	2.1
1	A	303	ILE	2.1
1	B	116	GLU	2.1
1	B	475	ARG	2.1
1	A	271	LEU	2.1
1	B	217	LYS	2.1
1	B	255	ASP	2.0
1	B	436	PHE	2.0
1	B	220	ALA	2.0
1	B	35	MET	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 carbohydrates 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	HEM	B	600	43/43	0.93	0.20	1.55	20,22,37,39	0
2	HEM	A	600	43/43	0.93	0.20	1.53	10,15,33,36	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.