



# Full wwPDB X-ray Structure Validation Report ⓘ

May 19, 2020 – 04:01 am BST

PDB ID : 1LH2  
Title : X-RAY STRUCTURAL INVESTIGATION OF LEGHEMOGLOBIN. VI. STRUCTURE OF ACETATE-FERRILEGHEMOGLOBIN AT A RESOLUTION OF 2.0 ANGSTROMS (RUSSIAN)  
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Deposited on : 1982-04-23  
Resolution : 2.00 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.

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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)  
Xtriage (Phenix) : 1.13  
EDS : 2.11  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.11

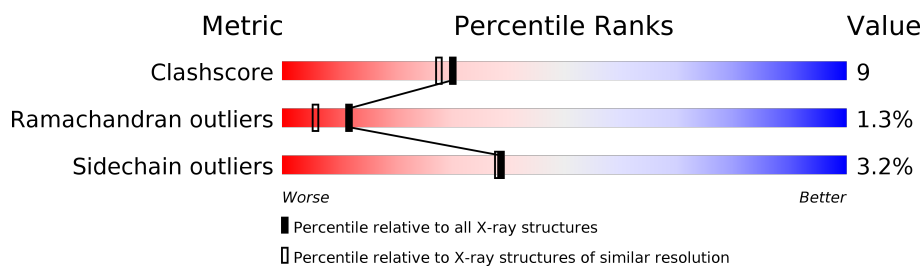
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)

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 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	

## 2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1291 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.

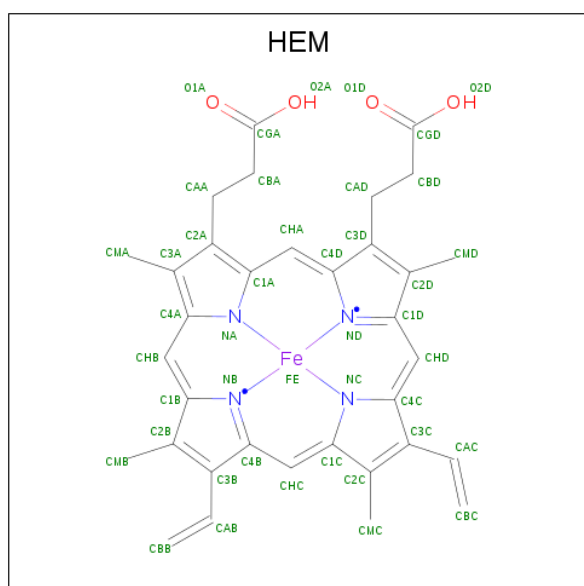
- Molecule 1 is a protein called LEGHEMOGLOBIN (AQUO MET).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	153	1180	761	193	225	1	36	1	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	79	GLU	GLN	CONFLICT	UNP P02240
A	150	ASP	ASN	CONFLICT	UNP P02240

- 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
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	3	0

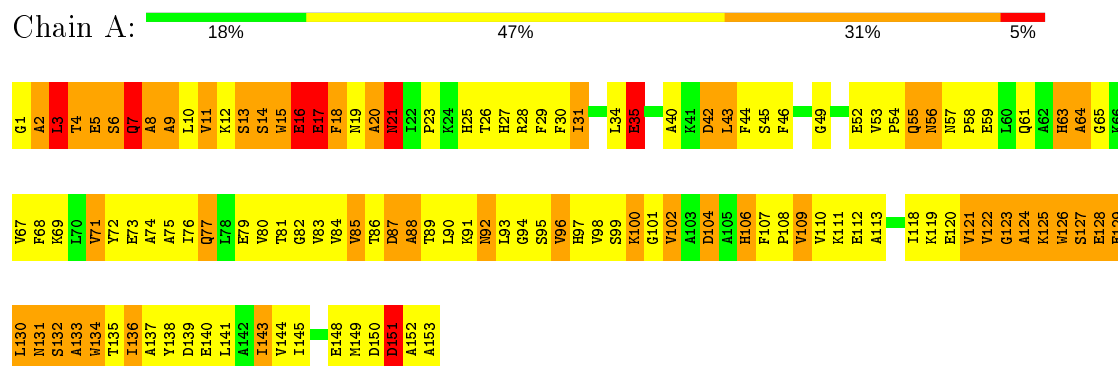
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	68	Total	O	0	0
			68	68		

### 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. 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: LEGHEMOGLOBIN (AQUO MET)



## 4 Data and refinement statistics

Property	Value	Source
Space group	B 1 1 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	93.22Å 38.26Å 52.01Å 90.00° 90.00° 98.80°	Depositor
Resolution (Å)	(Not available) – 2.00 9.93 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00) 96.4 (9.93-2.00)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	unknown	Depositor
R, $R_{free}$	(Not available) , (Not available) 0.500 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	19.3	Xtriage
Anisotropy	0.156	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	1.25 , 282.4	EDS
L-test for twinning <sup>1</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.51	EDS
Total number of atoms	1291	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	20.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.87% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</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	3.36	165/1214 (13.6%)	2.34	52/1648 (3.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	13

All (165) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	15	TRP	CD2-CE2	10.80	1.54	1.41
1	A	138	TYR	CB-CG	10.65	1.67	1.51
1	A	95	SER	CB-OG	10.15	1.55	1.42
1	A	112	GLU	CG-CD	9.39	1.66	1.51
1	A	35	GLU	CD-OE2	9.09	1.35	1.25
1	A	13	SER	CB-OG	-9.04	1.30	1.42
1	A	106	HIS	CB-CG	8.98	1.66	1.50
1	A	44	PHE	CB-CG	8.90	1.66	1.51
1	A	121	VAL	CB-CG2	8.85	1.71	1.52
1	A	132[A]	SER	CA-CB	8.85	1.66	1.52
1	A	132[B]	SER	CA-CB	8.85	1.66	1.52
1	A	132[C]	SER	CA-CB	8.85	1.66	1.52
1	A	120	GLU	CG-CD	8.77	1.65	1.51
1	A	13	SER	CA-CB	8.77	1.66	1.52
1	A	123	GLY	CA-C	8.67	1.65	1.51
1	A	15	TRP	CB-CG	8.40	1.65	1.50
1	A	45	SER	CB-OG	8.35	1.53	1.42
1	A	5	GLU	CD-OE2	8.34	1.34	1.25

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	11	VAL	CB-CG1	8.27	1.70	1.52
1	A	15	TRP	CZ3-CH2	8.17	1.53	1.40
1	A	138	TYR	CZ-OH	8.15	1.51	1.37
1	A	72	TYR	CE1-CZ	8.04	1.49	1.38
1	A	94	GLY	CA-C	8.02	1.64	1.51
1	A	15	TRP	CG-CD1	7.96	1.47	1.36
1	A	72	TYR	CG-CD2	7.89	1.49	1.39
1	A	140	GLU	CB-CG	7.83	1.67	1.52
1	A	138	TYR	CD2-CE2	7.79	1.51	1.39
1	A	35	GLU	CD-OE1	-7.61	1.17	1.25
1	A	65	GLY	C-O	7.59	1.35	1.23
1	A	46	PHE	CB-CG	7.51	1.64	1.51
1	A	18	PHE	CB-CG	7.50	1.64	1.51
1	A	79	GLU	CB-CG	7.42	1.66	1.52
1	A	27	HIS	CE1-NE2	7.41	1.49	1.32
1	A	128	GLU	CD-OE2	-7.30	1.17	1.25
1	A	111	LYS	N-CA	7.26	1.60	1.46
1	A	101	GLY	CA-C	7.25	1.63	1.51
1	A	73	GLU	CG-CD	7.16	1.62	1.51
1	A	129	GLU	CD-OE1	7.14	1.33	1.25
1	A	72	TYR	C-O	7.11	1.36	1.23
1	A	16	GLU	CG-CD	7.04	1.62	1.51
1	A	109	VAL	CA-CB	7.04	1.69	1.54
1	A	113	ALA	CA-CB	7.03	1.67	1.52
1	A	102	VAL	CB-CG1	6.95	1.67	1.52
1	A	88	ALA	N-CA	6.91	1.60	1.46
1	A	99	SER	N-CA	6.88	1.60	1.46
1	A	68	PHE	CG-CD2	6.87	1.49	1.38
1	A	148	GLU	CD-OE2	6.83	1.33	1.25
1	A	85	VAL	CB-CG2	6.82	1.67	1.52
1	A	97	HIS	CG-CD2	-6.78	1.24	1.35
1	A	124	ALA	N-CA	6.78	1.59	1.46
1	A	97	HIS	CA-CB	6.75	1.68	1.53
1	A	67	VAL	CB-CG2	6.73	1.67	1.52
1	A	152	ALA	C-O	6.69	1.36	1.23
1	A	134	TRP	N-CA	6.68	1.59	1.46
1	A	45	SER	N-CA	6.60	1.59	1.46
1	A	95	SER	N-CA	6.54	1.59	1.46
1	A	68	PHE	C-O	6.54	1.35	1.23
1	A	124	ALA	C-O	6.52	1.35	1.23
1	A	53	VAL	CA-CB	6.47	1.68	1.54
1	A	144	VAL	CB-CG2	6.46	1.66	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	79	GLU	CD-OE2	6.40	1.32	1.25
1	A	15	TRP	C-O	6.39	1.35	1.23
1	A	15	TRP	CD1-NE1	6.39	1.48	1.38
1	A	40	ALA	CA-CB	6.38	1.65	1.52
1	A	149	MET	C-O	6.38	1.35	1.23
1	A	14	SER	CB-OG	-6.38	1.33	1.42
1	A	6	SER	CA-CB	6.37	1.62	1.52
1	A	71	VAL	CB-CG1	6.33	1.66	1.52
1	A	61	GLN	C-O	6.30	1.35	1.23
1	A	92	ASN	CB-CG	6.29	1.65	1.51
1	A	27	HIS	CG-ND1	6.29	1.52	1.38
1	A	148	GLU	CB-CG	6.28	1.64	1.52
1	A	11	VAL	N-CA	6.27	1.58	1.46
1	A	59	GLU	CD-OE2	6.27	1.32	1.25
1	A	92	ASN	C-O	6.26	1.35	1.23
1	A	86	THR	N-CA	6.26	1.58	1.46
1	A	102	VAL	N-CA	6.25	1.58	1.46
1	A	68	PHE	CA-CB	-6.21	1.40	1.53
1	A	90	LEU	CA-CB	6.21	1.68	1.53
1	A	138	TYR	CD1-CE1	6.20	1.48	1.39
1	A	143	ILE	CA-CB	6.19	1.69	1.54
1	A	79	GLU	CD-OE1	6.18	1.32	1.25
1	A	145	ILE	N-CA	6.18	1.58	1.46
1	A	16	GLU	CD-OE1	6.17	1.32	1.25
1	A	110	VAL	CB-CG2	6.11	1.65	1.52
1	A	106	HIS	ND1-CE1	6.09	1.50	1.34
1	A	68	PHE	CE1-CZ	6.09	1.49	1.37
1	A	30	PHE	CA-CB	6.08	1.67	1.53
1	A	75	ALA	C-O	6.04	1.34	1.23
1	A	2	ALA	CA-CB	6.01	1.65	1.52
1	A	18	PHE	N-CA	6.01	1.58	1.46
1	A	122	VAL	N-CA	6.01	1.58	1.46
1	A	76	ILE	C-O	5.96	1.34	1.23
1	A	130	LEU	N-CA	5.96	1.58	1.46
1	A	127	SER	C-O	5.96	1.34	1.23
1	A	152	ALA	N-CA	5.92	1.58	1.46
1	A	140	GLU	CD-OE2	5.90	1.32	1.25
1	A	63	HIS	CA-C	5.90	1.68	1.52
1	A	108	PRO	N-CD	-5.89	1.39	1.47
1	A	120	GLU	CA-CB	5.88	1.66	1.53
1	A	4	THR	C-O	5.87	1.34	1.23
1	A	71	VAL	N-CA	5.87	1.58	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	29	PHE	CD1-CE1	-5.76	1.27	1.39
1	A	55	GLN	C-O	5.76	1.34	1.23
1	A	81	THR	N-CA	-5.74	1.34	1.46
1	A	79	GLU	CA-CB	-5.74	1.41	1.53
1	A	84	VAL	N-CA	5.72	1.57	1.46
1	A	133	ALA	CA-C	5.70	1.67	1.52
1	A	77	GLN	CG-CD	5.69	1.64	1.51
1	A	99	SER	C-O	5.65	1.34	1.23
1	A	46	PHE	N-CA	-5.64	1.35	1.46
1	A	110	VAL	CA-C	5.62	1.67	1.52
1	A	68	PHE	N-CA	5.61	1.57	1.46
1	A	28	ARG	CZ-NH1	5.59	1.40	1.33
1	A	111	LYS	CB-CG	5.58	1.67	1.52
1	A	148	GLU	CD-OE1	5.57	1.31	1.25
1	A	148	GLU	N-CA	5.56	1.57	1.46
1	A	52	GLU	CG-CD	-5.54	1.43	1.51
1	A	20	ALA	N-CA	-5.54	1.35	1.46
1	A	44	PHE	CD2-CE2	5.51	1.50	1.39
1	A	46	PHE	CD2-CE2	5.51	1.50	1.39
1	A	9	ALA	CA-CB	5.51	1.64	1.52
1	A	134	TRP	CD2-CE2	-5.50	1.34	1.41
1	A	58	PRO	CA-C	-5.48	1.41	1.52
1	A	141	LEU	CB-CG	5.46	1.68	1.52
1	A	29	PHE	CG-CD2	5.45	1.47	1.38
1	A	141	LEU	N-CA	5.41	1.57	1.46
1	A	87	ASP	CA-C	5.40	1.67	1.52
1	A	81	THR	CA-CB	5.40	1.67	1.53
1	A	104	ASP	N-CA	5.38	1.57	1.46
1	A	75	ALA	N-CA	5.36	1.57	1.46
1	A	1	GLY	C-O	5.33	1.32	1.23
1	A	43	LEU	CA-CB	5.32	1.66	1.53
1	A	125	LYS	CD-CE	5.32	1.64	1.51
1	A	136	ILE	CA-CB	5.32	1.67	1.54
1	A	137	ALA	CA-C	5.29	1.66	1.52
1	A	2	ALA	C-N	-5.27	1.22	1.34
1	A	108	PRO	C-N	5.26	1.46	1.34
1	A	97	HIS	C-N	-5.25	1.22	1.34
1	A	100	LYS	C-N	-5.25	1.23	1.33
1	A	82	GLY	CA-C	5.22	1.60	1.51
1	A	144	VAL	CA-C	5.22	1.66	1.52
1	A	49	GLY	CA-C	-5.22	1.43	1.51
1	A	65	GLY	N-CA	5.21	1.53	1.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	52	GLU	N-CA	5.21	1.56	1.46
1	A	79	GLU	C-O	5.18	1.33	1.23
1	A	64	ALA	C-O	5.15	1.33	1.23
1	A	54	PRO	C-N	-5.14	1.22	1.34
1	A	23	PRO	N-CA	5.14	1.55	1.47
1	A	111	LYS	CD-CE	5.14	1.64	1.51
1	A	8	ALA	N-CA	5.13	1.56	1.46
1	A	56	ASN	C-N	-5.12	1.22	1.34
1	A	7	GLN	N-CA	5.12	1.56	1.46
1	A	64	ALA	N-CA	5.11	1.56	1.46
1	A	29	PHE	CE1-CZ	5.10	1.47	1.37
1	A	18	PHE	CD2-CE2	5.09	1.49	1.39
1	A	153	ALA	C-OXT	5.08	1.33	1.23
1	A	98	VAL	CB-CG2	5.08	1.63	1.52
1	A	29	PHE	N-CA	-5.08	1.36	1.46
1	A	67	VAL	CA-C	5.07	1.66	1.52
1	A	8	ALA	C-O	5.06	1.32	1.23
1	A	134	TRP	C-O	5.05	1.32	1.23
1	A	84	VAL	CB-CG1	5.03	1.63	1.52
1	A	107	PHE	CE2-CZ	5.02	1.46	1.37
1	A	131	ASN	C-N	5.00	1.45	1.34

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	72	TYR	CB-CG-CD1	-11.33	114.20	121.00
1	A	16	GLU	OE1-CD-OE2	-10.41	110.80	123.30
1	A	15	TRP	CG-CD2-CE3	-9.13	125.68	133.90
1	A	87	ASP	CB-CG-OD2	-8.99	110.21	118.30
1	A	112	GLU	OE1-CD-OE2	-8.91	112.61	123.30
1	A	73	GLU	OE1-CD-OE2	-8.65	112.92	123.30
1	A	17	GLU	OE1-CD-OE2	-8.06	113.63	123.30
1	A	29	PHE	CD1-CG-CD2	7.33	127.82	118.30
1	A	29	PHE	CB-CG-CD1	-7.31	115.68	120.80
1	A	122	VAL	CA-CB-CG1	-7.30	99.94	110.90
1	A	68	PHE	CB-CG-CD1	-7.29	115.69	120.80
1	A	15	TRP	CD1-NE1-CE2	-7.26	102.47	109.00
1	A	138	TYR	CB-CG-CD2	7.18	125.31	121.00
1	A	15	TRP	CB-CG-CD1	6.96	136.04	127.00
1	A	72	TYR	CG-CD1-CE1	-6.90	115.78	121.30
1	A	128	GLU	OE1-CD-OE2	-6.87	115.06	123.30
1	A	46	PHE	CB-CG-CD2	6.86	125.60	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	80	VAL	CA-CB-CG2	-6.75	100.78	110.90
1	A	126	TRP	NE1-CE2-CD2	6.63	113.93	107.30
1	A	30	PHE	CB-CG-CD2	-6.59	116.19	120.80
1	A	15	TRP	CE2-CD2-CE3	6.52	126.53	118.70
1	A	126	TRP	CE2-CD2-CG	-6.51	102.09	107.30
1	A	134	TRP	CG-CD1-NE1	-6.47	103.63	110.10
1	A	138	TYR	CD1-CE1-CZ	6.35	125.52	119.80
1	A	124	ALA	CB-CA-C	-6.31	100.63	110.10
1	A	23	PRO	N-CA-CB	6.21	110.75	103.30
1	A	29	PHE	CB-CG-CD2	-6.19	116.46	120.80
1	A	153	ALA	N-CA-CB	-6.14	101.50	110.10
1	A	120	GLU	OE1-CD-OE2	-6.05	116.03	123.30
1	A	15	TRP	CH2-CZ2-CE2	-5.93	111.47	117.40
1	A	15	TRP	CD2-CE3-CZ3	-5.85	111.19	118.80
1	A	31	ILE	CB-CA-C	-5.85	99.91	111.60
1	A	69	LYS	O-C-N	5.82	132.01	122.70
1	A	72	TYR	N-CA-CB	-5.80	100.16	110.60
1	A	15	TRP	NE1-CE2-CD2	5.78	113.08	107.30
1	A	21	ASN	CB-CG-OD1	-5.66	110.29	121.60
1	A	151	ASP	CB-CG-OD2	-5.60	113.26	118.30
1	A	127	SER	O-C-N	5.49	131.49	122.70
1	A	121	VAL	CA-C-N	5.45	129.19	117.20
1	A	144	VAL	O-C-N	-5.40	114.05	122.70
1	A	3	LEU	N-CA-CB	5.40	121.19	110.40
1	A	72	TYR	CZ-CE2-CD2	-5.37	114.97	119.80
1	A	77	GLN	CB-CA-C	5.34	121.07	110.40
1	A	42	ASP	CB-CG-OD1	-5.32	113.51	118.30
1	A	35	GLU	CB-CA-C	-5.31	99.79	110.40
1	A	89	THR	N-CA-CB	-5.23	100.37	110.30
1	A	54	PRO	N-CA-CB	5.18	109.52	103.30
1	A	15	TRP	NE1-CE2-CZ2	-5.12	124.76	130.40
1	A	76	ILE	O-C-N	5.12	130.88	122.70
1	A	126	TRP	CE2-CD2-CE3	5.05	124.77	118.70
1	A	104	ASP	CB-CA-C	-5.04	100.32	110.40
1	A	110	VAL	O-C-N	-5.04	114.64	122.70

There are no chirality outliers.

All (13) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	104	ASP	Sidechain
1	A	150	ASP	Sidechain

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Mol	Chain	Res	Type	Group
1	A	151	ASP	Sidechain
1	A	16	GLU	Sidechain
1	A	17	GLU	Sidechain
1	A	19	ASN	Sidechain
1	A	21	ASN	Sidechain
1	A	35	GLU	Sidechain
1	A	42	ASP	Sidechain
1	A	57	ASN	Sidechain
1	A	7	GLN	Sidechain
1	A	83	VAL	Mainchain
1	A	87	ASP	Sidechain

## 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	1180	0	1200	22	206
2	A	43	0	30	3	0
3	A	68	0	0	0	30
All	All	1291	0	1230	22	207

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 9.

All (22) 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:92:ASN:O	1:A:96:VAL:HG12	1.94	0.68
1:A:31:ILE:O	1:A:35:GLU:HG3	2.02	0.58
1:A:77:GLN:NE2	1:A:85:VAL:H	2.03	0.56
1:A:139:ASP:O	1:A:143:ILE:HG13	2.06	0.56
1:A:21:ASN:C	1:A:21:ASN:HD22	2.13	0.52
1:A:21:ASN:C	1:A:21:ASN:ND2	2.64	0.51
1:A:106:HIS:O	1:A:109:VAL:HB	2.12	0.50
1:A:17:GLU:OE2	1:A:122:VAL:HG12	2.12	0.50
1:A:18:PHE:CE1	1:A:25:HIS:HB3	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:LEU:HA	1:A:93:LEU:HD23	1.50	0.48
1:A:63:HIS:HE1	2:A:154:HEM:C4D	2.31	0.48
1:A:126:TRP:CZ2	1:A:131:ASN:HB2	2.48	0.48
1:A:100:LYS:HG3	2:A:154:HEM:HAD2	1.97	0.47
1:A:71:VAL:O	1:A:74:ALA:HB3	2.17	0.45
1:A:88:ALA:O	1:A:91:LYS:HB2	2.17	0.45
1:A:26:THR:HB	1:A:64:ALA:HB3	1.99	0.43
1:A:102:VAL:HG13	2:A:154:HEM:HAC	2.00	0.43
1:A:43:LEU:HD23	1:A:43:LEU:HA	1.87	0.43
1:A:18:PHE:O	1:A:25:HIS:HD2	2.03	0.41
1:A:143:ILE:HG21	1:A:143:ILE:HD13	1.88	0.41
1:A:12:LYS:HE3	1:A:16:GLU:OE2	2.21	0.40

All (207) 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)
1:A:4:THR:OG1	1:A:17:GLU:CB[2_675]	0.09	2.11
1:A:126:TRP:CD1	1:A:128:GLU:C[2_675]	0.37	1.83
1:A:130:LEU:CB	1:A:130:LEU:CB[2_675]	0.44	1.76
1:A:9:ALA:CB	1:A:12:LYS:C[2_675]	0.48	1.72
1:A:6:SER:O	1:A:14:SER:N[2_675]	0.56	1.64
1:A:8:ALA:CA	1:A:13:SER:OG[2_675]	0.58	1.62
1:A:127:SER:N	1:A:131:ASN:CA[2_675]	0.61	1.59
1:A:126:TRP:NE1	1:A:128:GLU:CA[2_675]	0.66	1.54
1:A:126:TRP:C	1:A:131:ASN:CA[2_675]	0.69	1.51
1:A:9:ALA:N	1:A:13:SER:CA[2_675]	0.71	1.49
1:A:119:LYS:O	3:A:182:HOH:O[2_675]	0.74	1.46
1:A:6:SER:OG	1:A:14:SER:O[2_675]	0.77	1.43
1:A:119:LYS:CE	1:A:128:GLU:CG[2_675]	0.78	1.42
1:A:119:LYS:CD	1:A:128:GLU:OE2[2_675]	0.81	1.39
1:A:10:LEU:CA	1:A:10:LEU:C[2_675]	0.82	1.38
1:A:126:TRP:C	1:A:131:ASN:C[2_675]	0.83	1.37
1:A:8:ALA:C	1:A:13:SER:CB[2_675]	0.89	1.31
1:A:9:ALA:CA	1:A:13:SER:N[2_675]	0.89	1.31
1:A:10:LEU:CA	1:A:10:LEU:O[2_675]	0.93	1.27
1:A:126:TRP:CG	1:A:128:GLU:O[2_675]	0.96	1.24
1:A:10:LEU:O	1:A:10:LEU:CB[2_675]	1.00	1.20
1:A:126:TRP:NE1	1:A:128:GLU:N[2_675]	1.02	1.18
1:A:125:LYS:CB	1:A:133:ALA:CA[2_675]	1.06	1.14

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:GLN:CA	3:A:203:HOH:O[2_675]	1.11	1.09
1:A:127:SER:N	1:A:131:ASN:N[2_675]	1.11	1.09
1:A:130:LEU:N	1:A:130:LEU:CD2[2_675]	1.11	1.09
1:A:7:GLN:CG	1:A:17:GLU:OE2[2_675]	1.13	1.07
1:A:126:TRP:CD1	1:A:128:GLU:O[2_675]	1.16	1.04
1:A:7:GLN:OE1	1:A:17:GLU:OE1[2_675]	1.19	1.01
1:A:7:GLN:NE2	1:A:122:VAL:CA[2_675]	1.19	1.01
1:A:9:ALA:CB	1:A:13:SER:N[2_675]	1.19	1.01
1:A:119:LYS:CD	1:A:128:GLU:CD[2_675]	1.19	1.01
1:A:126:TRP:O	1:A:131:ASN:O[2_675]	1.19	1.01
1:A:7:GLN:CB	3:A:203:HOH:O[2_675]	1.20	1.00
1:A:130:LEU:O	3:A:192:HOH:O[2_675]	1.23	0.97
1:A:134:TRP:N	3:A:219:HOH:O[2_675]	1.24	0.96
1:A:6:SER:N	1:A:17:GLU:N[2_675]	1.25	0.95
1:A:9:ALA:N	1:A:13:SER:CB[2_675]	1.25	0.95
1:A:6:SER:CB	1:A:14:SER:O[2_675]	1.25	0.95
1:A:6:SER:O	1:A:14:SER:CA[2_675]	1.27	0.93
1:A:8:ALA:C	1:A:13:SER:OG[2_675]	1.28	0.92
1:A:7:GLN:C	3:A:203:HOH:O[2_675]	1.28	0.92
1:A:126:TRP:O	1:A:131:ASN:C[2_675]	1.29	0.91
1:A:123:GLY:O	1:A:132[B]:SER:OG[2_675]	1.29	0.91
1:A:122:VAL:N	1:A:129:GLU:OE2[2_675]	1.32	0.88
1:A:125:LYS:C	1:A:133:ALA:N[2_675]	1.32	0.88
1:A:9:ALA:O	1:A:9:ALA:O[2_675]	1.34	0.86
1:A:126:TRP:CD1	1:A:128:GLU:CA[2_675]	1.35	0.85
1:A:130:LEU:CB	1:A:130:LEU:CG[2_675]	1.36	0.84
1:A:130:LEU:CA	1:A:130:LEU:CD2[2_675]	1.37	0.83
1:A:4:THR:CB	1:A:17:GLU:CB[2_675]	1.39	0.81
1:A:126:TRP:CE2	1:A:128:GLU:CA[2_675]	1.39	0.81
1:A:8:ALA:N	1:A:13:SER:OG[2_675]	1.42	0.78
1:A:7:GLN:CB	1:A:17:GLU:OE2[2_675]	1.43	0.77
1:A:126:TRP:CH2	1:A:126:TRP:CH2[2_675]	1.43	0.77
1:A:6:SER:C	1:A:14:SER:CA[2_675]	1.43	0.77
1:A:126:TRP:CA	1:A:131:ASN:C[2_675]	1.44	0.76
1:A:9:ALA:CB	1:A:12:LYS:CA[2_675]	1.44	0.76
1:A:6:SER:O	1:A:13:SER:C[2_675]	1.45	0.75
1:A:9:ALA:N	1:A:13:SER:N[2_675]	1.48	0.72
1:A:123:GLY:C	1:A:132[C]:SER:OG[2_675]	1.49	0.71
1:A:17:GLU:O	3:A:222:HOH:O[2_675]	1.49	0.71
1:A:126:TRP:CB	1:A:128:GLU:O[2_675]	1.49	0.71
1:A:6:SER:C	1:A:13:SER:O[2_675]	1.50	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:4:THR:OG1	1:A:17:GLU:CA[2_675]	1.50	0.70
1:A:7:GLN:CD	1:A:17:GLU:OE2[2_675]	1.51	0.69
1:A:4:THR:N	1:A:17:GLU:CG[2_675]	1.52	0.68
1:A:15:TRP:CB	3:A:221:HOH:O[2_675]	1.52	0.68
1:A:126:TRP:NE1	1:A:128:GLU:C[2_675]	1.53	0.67
1:A:10:LEU:CG	1:A:14:SER:OG[2_675]	1.53	0.67
1:A:10:LEU:CD1	1:A:14:SER:OG[2_675]	1.57	0.63
1:A:7:GLN:OE1	1:A:17:GLU:CD[2_675]	1.58	0.62
1:A:4:THR:CA	1:A:17:GLU:CG[2_675]	1.59	0.61
1:A:125:LYS:O	1:A:133:ALA:N[2_675]	1.59	0.61
1:A:7:GLN:OE1	1:A:17:GLU:OE2[2_675]	1.60	0.60
1:A:9:ALA:CB	1:A:12:LYS:O[2_675]	1.60	0.60
1:A:4:THR:CB	1:A:17:GLU:CA[2_675]	1.60	0.60
1:A:6:SER:C	1:A:14:SER:N[2_675]	1.61	0.59
1:A:7:GLN:N	1:A:13:SER:O[2_675]	1.61	0.59
1:A:125:LYS:CA	1:A:133:ALA:N[2_675]	1.62	0.58
1:A:126:TRP:O	1:A:131:ASN:CA[2_675]	1.63	0.57
1:A:56:ASN:OD1	3:A:171:HOH:O[1_545]	1.63	0.57
1:A:125:LYS:CB	1:A:133:ALA:N[2_675]	1.63	0.57
1:A:118:ILE:CG2	1:A:129:GLU:CB[2_675]	1.64	0.56
1:A:124:ALA:O	1:A:135:THR:CB[2_675]	1.64	0.56
1:A:122:VAL:CG2	1:A:132[A]:SER:OG[2_675]	1.65	0.55
1:A:126:TRP:C	1:A:131:ASN:CB[2_675]	1.65	0.55
1:A:10:LEU:CD2	1:A:11:VAL:CA[2_675]	1.66	0.54
1:A:10:LEU:N	1:A:10:LEU:O[2_675]	1.66	0.54
1:A:126:TRP:CG	1:A:128:GLU:C[2_675]	1.66	0.54
1:A:126:TRP:NE1	1:A:127:SER:C[2_675]	1.67	0.53
1:A:9:ALA:CA	1:A:13:SER:CA[2_675]	1.67	0.53
1:A:56:ASN:N	3:A:170:HOH:O[1_545]	1.67	0.53
1:A:4:THR:OG1	1:A:17:GLU:CG[2_675]	1.68	0.52
1:A:6:SER:CB	1:A:14:SER:C[2_675]	1.68	0.52
1:A:118:ILE:O	1:A:129:GLU:OE2[2_675]	1.69	0.51
1:A:10:LEU:CA	1:A:10:LEU:CA[2_675]	1.69	0.51
1:A:4:THR:O	3:A:206:HOH:O[2_675]	1.70	0.50
1:A:6:SER:C	1:A:13:SER:C[2_675]	1.72	0.48
1:A:126:TRP:CD1	1:A:129:GLU:N[2_675]	1.72	0.48
1:A:127:SER:N	1:A:131:ASN:CB[2_675]	1.72	0.48
1:A:126:TRP:N	1:A:131:ASN:C[2_675]	1.72	0.48
1:A:9:ALA:C	1:A:13:SER:N[2_675]	1.73	0.47
1:A:119:LYS:CG	1:A:128:GLU:CB[2_675]	1.73	0.47
1:A:6:SER:C	1:A:14:SER:C[2_675]	1.73	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:15:TRP:N	3:A:221:HOH:O[2_675]	1.74	0.46
1:A:8:ALA:CA	1:A:13:SER:CB[2_675]	1.75	0.45
1:A:6:SER:CA	1:A:16:GLU:N[2_675]	1.75	0.45
1:A:134:TRP:CA	3:A:219:HOH:O[2_675]	1.76	0.44
1:A:5:GLU:CB	1:A:16:GLU:CB[2_675]	1.76	0.44
1:A:6:SER:CB	1:A:15:TRP:CA[2_675]	1.76	0.44
1:A:119:LYS:CD	1:A:128:GLU:CG[2_675]	1.77	0.43
1:A:9:ALA:C	1:A:9:ALA:O[2_675]	1.77	0.43
1:A:127:SER:CA	1:A:131:ASN:CA[2_675]	1.78	0.42
1:A:5:GLU:CG	1:A:16:GLU:CB[2_675]	1.78	0.42
1:A:128:GLU:N	3:A:212:HOH:O[2_675]	1.78	0.42
1:A:129:GLU:C	1:A:130:LEU:CD2[2_675]	1.78	0.42
1:A:7:GLN:NE2	1:A:122:VAL:C[2_675]	1.80	0.40
1:A:119:LYS:C	3:A:182:HOH:O[2_675]	1.80	0.40
1:A:130:LEU:CA	1:A:130:LEU:CG[2_675]	1.80	0.40
1:A:6:SER:CB	1:A:15:TRP:C[2_675]	1.81	0.39
1:A:8:ALA:O	1:A:13:SER:CB[2_675]	1.81	0.39
1:A:125:LYS:O	1:A:130:LEU:O[2_675]	1.81	0.39
1:A:10:LEU:O	1:A:10:LEU:CG[2_675]	1.81	0.39
1:A:17:GLU:C	3:A:222:HOH:O[2_675]	1.82	0.38
1:A:126:TRP:C	1:A:131:ASN:N[2_675]	1.83	0.37
3:A:159:HOH:O	3:A:177:HOH:O[2_675]	1.83	0.37
1:A:10:LEU:C	1:A:10:LEU:CB[2_675]	1.84	0.36
1:A:123:GLY:O	1:A:132[C]:SER:OG[2_675]	1.84	0.36
1:A:122:VAL:O	1:A:132[C]:SER:OG[2_675]	1.84	0.36
1:A:4:THR:N	1:A:17:GLU:CD[2_675]	1.85	0.35
1:A:130:LEU:CA	1:A:130:LEU:CB[2_675]	1.85	0.35
1:A:4:THR:C	1:A:17:GLU:CG[2_675]	1.86	0.34
1:A:6:SER:CB	1:A:15:TRP:N[2_675]	1.86	0.34
1:A:56:ASN:CB	3:A:189:HOH:O[1_545]	1.86	0.34
1:A:126:TRP:C	1:A:131:ASN:O[2_675]	1.86	0.34
1:A:118:ILE:O	1:A:129:GLU:CD[2_675]	1.88	0.32
1:A:10:LEU:CD1	1:A:14:SER:CB[2_675]	1.88	0.32
1:A:135:THR:OG1	3:A:159:HOH:O[2_675]	1.89	0.31
1:A:7:GLN:N	1:A:14:SER:CA[2_675]	1.90	0.30
1:A:126:TRP:CE2	1:A:128:GLU:N[2_675]	1.90	0.30
1:A:5:GLU:C	1:A:16:GLU:CB[2_675]	1.90	0.30
1:A:15:TRP:CA	3:A:221:HOH:O[2_675]	1.93	0.27
1:A:122:VAL:CG2	1:A:129:GLU:O[2_675]	1.93	0.27
1:A:6:SER:N	1:A:16:GLU:C[2_675]	1.93	0.27
1:A:7:GLN:NE2	1:A:122:VAL:CB[2_675]	1.93	0.27

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:6:SER:CA	1:A:14:SER:C[2_675]	1.94	0.26
1:A:7:GLN:CG	1:A:122:VAL:CG1[2_675]	1.94	0.26
1:A:126:TRP:CZ3	1:A:126:TRP:CH2[2_675]	1.95	0.25
1:A:119:LYS:CG	1:A:128:GLU:OE2[2_675]	1.95	0.25
1:A:127:SER:N	1:A:131:ASN:C[2_675]	1.96	0.24
1:A:10:LEU:C	1:A:10:LEU:C[2_675]	1.97	0.23
1:A:9:ALA:CA	1:A:12:LYS:C[2_675]	1.97	0.23
1:A:8:ALA:C	1:A:13:SER:CA[2_675]	1.97	0.23
1:A:6:SER:OG	1:A:14:SER:C[2_675]	1.97	0.23
1:A:6:SER:CA	1:A:15:TRP:N[2_675]	1.97	0.23
1:A:6:SER:N	1:A:16:GLU:N[2_675]	1.97	0.23
1:A:122:VAL:CG2	1:A:129:GLU:CA[2_675]	1.98	0.22
1:A:126:TRP:CB	1:A:132[B]:SER:OG[2_675]	1.98	0.22
1:A:56:ASN:OD1	3:A:170:HOH:O[1_545]	1.98	0.22
1:A:56:ASN:CB	3:A:170:HOH:O[1_545]	2.01	0.19
1:A:5:GLU:CA	1:A:16:GLU:CB[2_675]	2.03	0.17
1:A:119:LYS:CB	1:A:128:GLU:CB[2_675]	2.03	0.17
1:A:125:LYS:CB	1:A:133:ALA:C[2_675]	2.04	0.16
1:A:10:LEU:N	1:A:10:LEU:C[2_675]	2.04	0.16
1:A:3:LEU:CG	1:A:125:LYS:CE[2_675]	2.04	0.16
1:A:7:GLN:O	3:A:203:HOH:O[2_675]	2.04	0.16
1:A:11:VAL:O	3:A:221:HOH:O[2_675]	2.05	0.15
1:A:5:GLU:O	1:A:13:SER:O[2_675]	2.05	0.15
1:A:126:TRP:O	1:A:131:ASN:CG[2_675]	2.06	0.14
1:A:4:THR:CB	1:A:17:GLU:CG[2_675]	2.06	0.14
1:A:122:VAL:CG2	1:A:129:GLU:C[2_675]	2.07	0.13
1:A:4:THR:N	1:A:17:GLU:OE1[2_675]	2.08	0.12
1:A:7:GLN:C	1:A:13:SER:OG[2_675]	2.08	0.12
1:A:4:THR:O	1:A:17:GLU:CG[2_675]	2.08	0.12
1:A:130:LEU:N	1:A:130:LEU:CG[2_675]	2.09	0.11
1:A:55:GLN:OE1	3:A:172:HOH:O[1_545]	2.09	0.11
1:A:126:TRP:CG	1:A:128:GLU:CA[2_675]	2.09	0.11
1:A:8:ALA:CB	1:A:13:SER:OG[2_675]	2.09	0.11
1:A:7:GLN:NE2	1:A:122:VAL:O[2_675]	2.10	0.10
1:A:9:ALA:N	1:A:13:SER:OG[2_675]	2.11	0.09
1:A:12:LYS:CB	3:A:205:HOH:O[2_675]	2.11	0.09
1:A:122:VAL:CG1	1:A:129:GLU:CG[2_675]	2.11	0.09
1:A:6:SER:O	1:A:13:SER:O[2_675]	2.11	0.09
1:A:125:LYS:CA	1:A:133:ALA:CA[2_675]	2.12	0.08
1:A:126:TRP:CB	3:A:191:HOH:O[2_675]	2.12	0.08
1:A:126:TRP:CA	1:A:131:ASN:CA[2_675]	2.12	0.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:ALA:N	1:A:132[C]:SER:OG[2_675]	2.13	0.07
1:A:6:SER:O	1:A:14:SER:C[2_675]	2.13	0.07
1:A:6:SER:N	1:A:16:GLU:CA[2_675]	2.13	0.07
1:A:121:VAL:CB	1:A:129:GLU:OE1[2_675]	2.13	0.07
1:A:125:LYS:C	1:A:131:ASN:C[2_675]	2.14	0.06
1:A:8:ALA:N	3:A:203:HOH:O[2_675]	2.14	0.06
1:A:130:LEU:C	3:A:192:HOH:O[2_675]	2.15	0.05
1:A:20:ALA:N	3:A:222:HOH:O[2_675]	2.15	0.05
1:A:5:GLU:N	1:A:17:GLU:N[2_675]	2.15	0.05
1:A:130:LEU:CB	1:A:130:LEU:CD2[2_675]	2.15	0.05
1:A:5:GLU:O	1:A:16:GLU:CB[2_675]	2.16	0.04
1:A:127:SER:N	1:A:130:LEU:C[2_675]	2.16	0.04
1:A:5:GLU:C	1:A:13:SER:O[2_675]	2.16	0.04
1:A:126:TRP:N	1:A:133:ALA:N[2_675]	2.17	0.03
1:A:9:ALA:CB	1:A:12:LYS:CB[2_675]	2.17	0.03
1:A:119:LYS:NZ	1:A:128:GLU:CG[2_675]	2.18	0.02
1:A:126:TRP:CD2	1:A:128:GLU:CA[2_675]	2.18	0.02
1:A:6:SER:CA	1:A:13:SER:O[2_675]	2.19	0.01
1:A:126:TRP:CE3	1:A:131:ASN:CB[2_675]	2.19	0.01
1:A:119:LYS:CE	1:A:128:GLU:CB[2_675]	2.19	0.01

## 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	153/153 (100%)	148 (97%)	3 (2%)	2 (1%)	12 6

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	2	ALA
1	A	3	LEU

### 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	127/125 (102%)	123 (97%)	4 (3%)	40	40

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	34	LEU
1	A	96	VAL
1	A	151	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	ASN
1	A	25	HIS
1	A	27	HIS
1	A	61	GLN
1	A	63	HIS
1	A	77	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

1 ligand is 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	154	1,3	27,50,50	4.07	20 (74%)	17,82,82	2.97	9 (52%)

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	154	1,3	-	0/6/54/54	-

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	154	HEM	C3B-CAB	8.61	1.65	1.47
2	A	154	HEM	CAD-C3D	6.68	1.63	1.52
2	A	154	HEM	C4D-C3D	6.06	1.56	1.42
2	A	154	HEM	C1B-C2B	5.81	1.55	1.42
2	A	154	HEM	CAA-C2A	5.54	1.60	1.52
2	A	154	HEM	C1A-NA	5.34	1.47	1.36
2	A	154	HEM	C3C-C2C	5.23	1.47	1.40
2	A	154	HEM	C4A-NA	4.84	1.46	1.36
2	A	154	HEM	C3C-CAC	4.83	1.57	1.47
2	A	154	HEM	CMD-C2D	4.83	1.61	1.51
2	A	154	HEM	C4A-CHB	4.50	1.53	1.41
2	A	154	HEM	CMC-C2C	3.48	1.59	1.51
2	A	154	HEM	C2A-C3A	3.24	1.47	1.37
2	A	154	HEM	CMB-C2B	3.18	1.59	1.51
2	A	154	HEM	C1A-CHA	2.99	1.49	1.41
2	A	154	HEM	C3B-C2B	-2.93	1.36	1.40
2	A	154	HEM	CMA-C3A	2.58	1.57	1.51
2	A	154	HEM	C4B-NB	2.49	1.41	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	154	HEM	C1D-ND	2.46	1.41	1.36
2	A	154	HEM	C1C-C2C	2.19	1.47	1.42

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	154	HEM	CMA-C3A-C4A	-6.03	119.20	128.46
2	A	154	HEM	CBA-CAA-C2A	-4.81	103.61	112.49
2	A	154	HEM	C4A-C3A-C2A	4.58	110.18	107.00
2	A	154	HEM	C4C-C3C-C2C	4.48	110.03	106.90
2	A	154	HEM	C3C-C4C-NC	-3.63	104.09	110.94
2	A	154	HEM	CMA-C3A-C2A	2.99	130.57	124.94
2	A	154	HEM	C3B-C4B-NB	2.69	112.69	109.21
2	A	154	HEM	CAA-CBA-CGA	-2.61	108.29	112.67
2	A	154	HEM	CBD-CAD-C3D	-2.15	108.51	112.48

There are no chirality outliers.

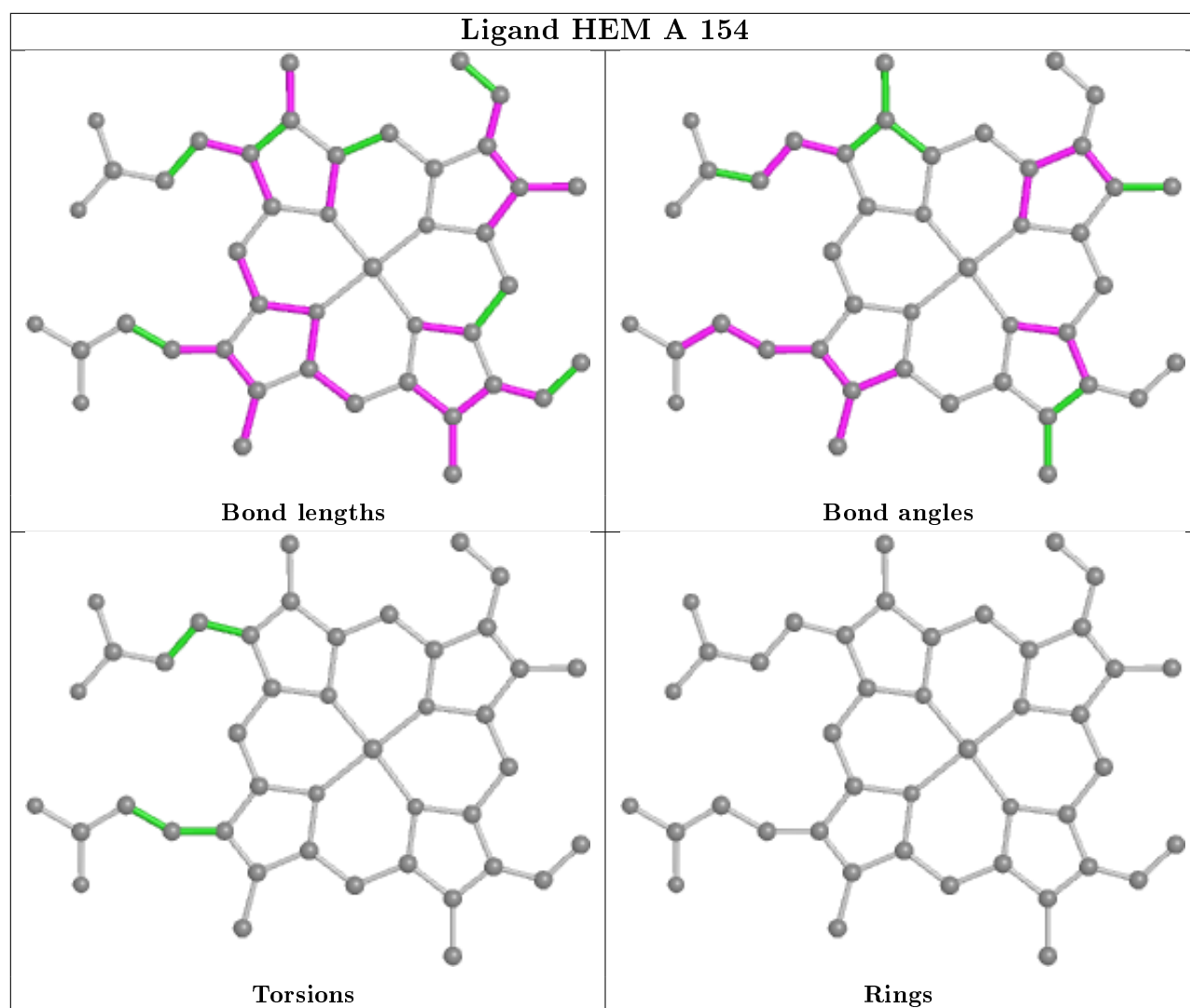
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 3 short contacts:

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

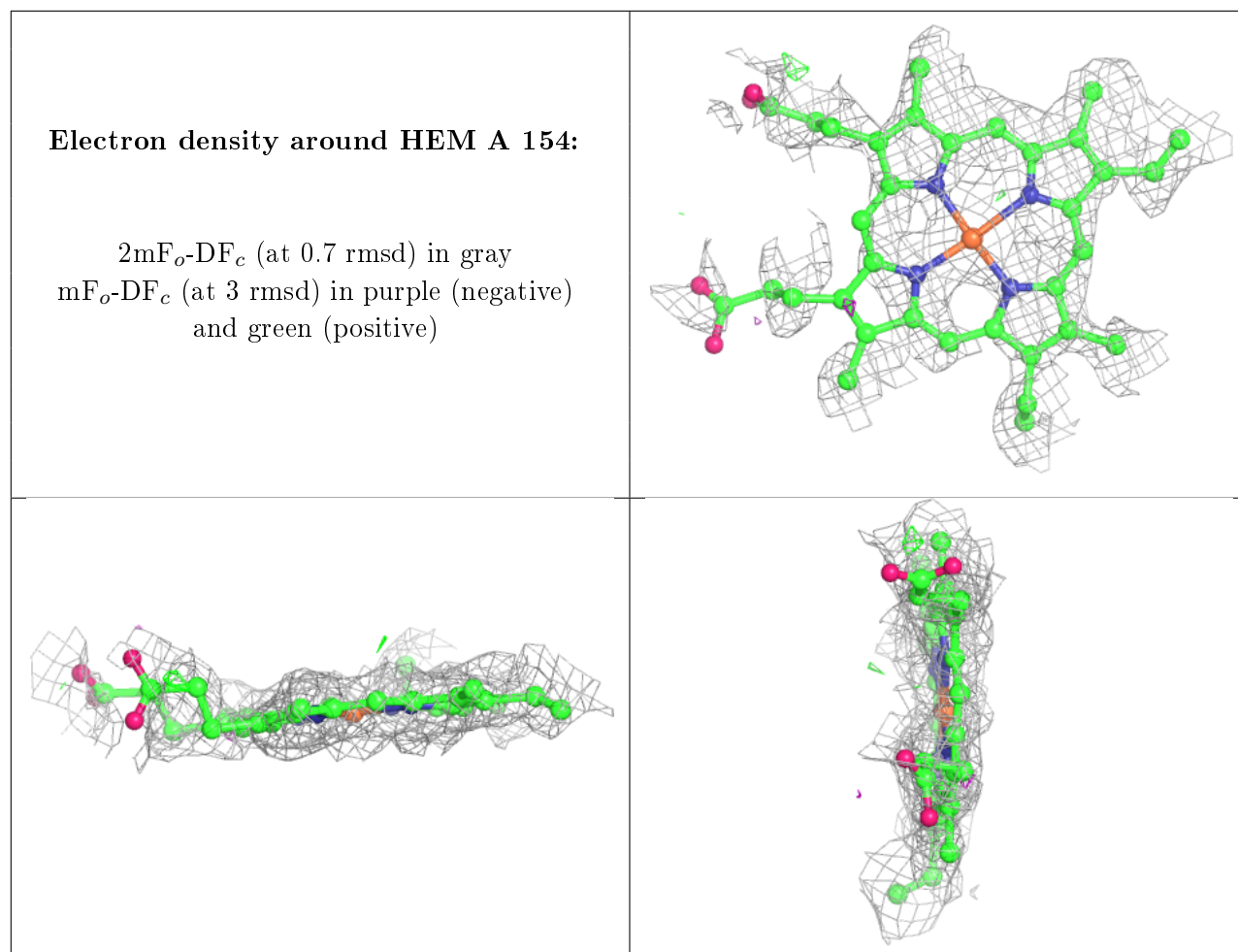
Unable to reproduce the depositors R factor - this section is therefore empty.

### 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 depositors R factor - this section is therefore empty.