



Full wwPDB X-ray Structure Validation Report ⓘ

Nov 25, 2024 – 04:47 PM EST

PDB ID : 2LH7
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.

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A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

The following versions of software and data (see [references](#) ①) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

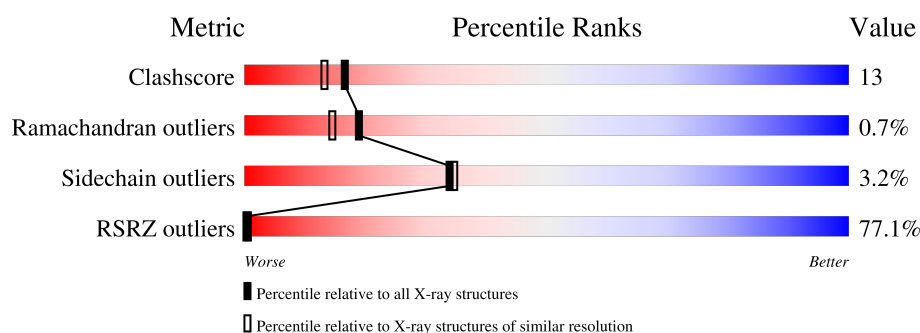
1 Overall quality at a glance

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	180529	10737 (2.00-2.00)
Ramachandran outliers	177936	10628 (2.00-2.00)
Sidechain outliers	177891	10627 (2.00-2.00)
RSRZ outliers	164620	9409 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	153	<div> <div>77%</div> <div>27%</div> <div>54%</div> <div>15%</div> <div>.</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NBE	A	155	-	-	X	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 1295 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.

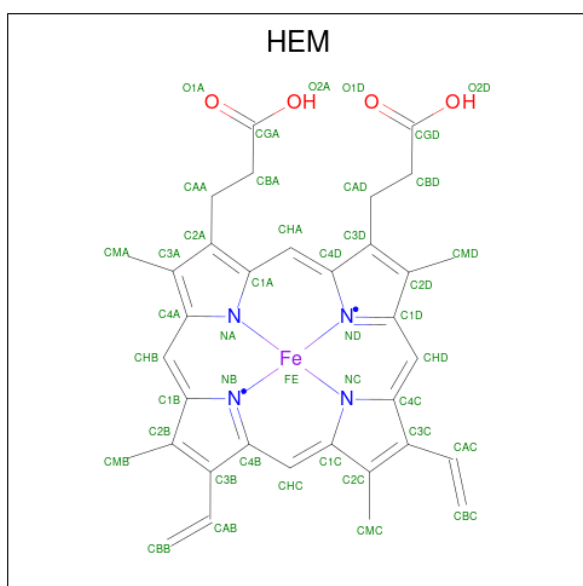
- Molecule 1 is a protein called LEGHEMOGLOBIN (NITROSOBENZENE).

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

There are 2 discrepancies between the modelled and reference sequences:

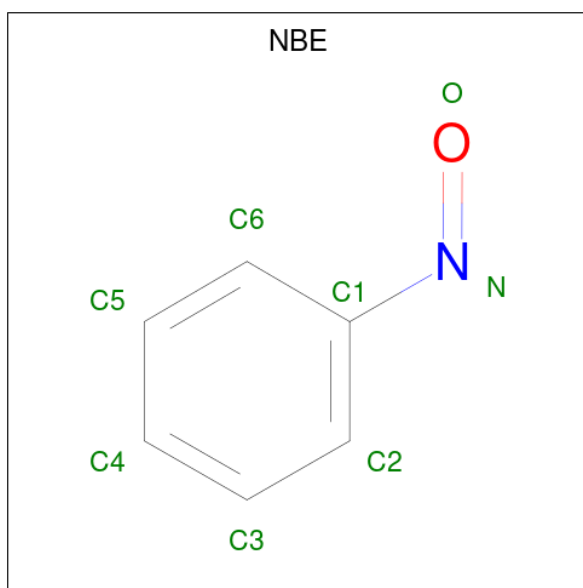
Chain	Residue	Modelled	Actual	Comment	Reference
A	79	GLU	GLN	conflict	UNP P02240
A	150	ASP	ASN	conflict	UNP P02240

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

- Molecule 3 is NITROSOBENZENE (three-letter code: NBE) (formula: C_6H_5NO).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			8	6	1	1		

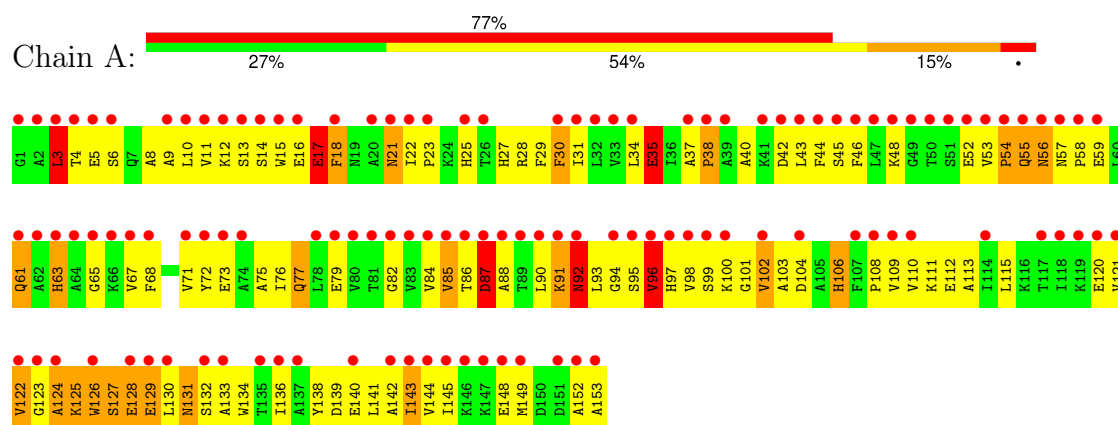
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	64	Total	O	0	0
			64	64		

3 Residue-property plots

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: LEGHEMOGLOBIN (NITROSOBENZENE)



4 Data and refinement statistics

Property	Value	Source
Space group	B 1 1 2	Depositor
Cell constants a, b, c, α , β , γ	93.23Å 38.25Å 51.88Å 90.00° 90.00° 98.70°	Depositor
Resolution (Å)	(Not available) – 2.00 46.08 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00) 92.9 (46.08-2.00)	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) 0.423 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	20.9	Xtriage
Anisotropy	0.256	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.45 , 114.9	EDS
L-test for twinning ¹	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.67	EDS
Total number of atoms	1295	wwPDB-VP
Average B, all atoms (Å ²)	21.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.07% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Theoretical 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, NBE

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	3.17	128/1214 (10.5%)	2.08	37/1648 (2.2%)

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	8

All (128) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	95	SER	CB-OG	10.30	1.55	1.42
1	A	138	TYR	CB-CG	9.74	1.66	1.51
1	A	138	TYR	CZ-OH	9.43	1.53	1.37
1	A	15	TRP	CD2-CE2	9.38	1.52	1.41
1	A	112	GLU	CG-CD	9.14	1.65	1.51
1	A	72	TYR	CE1-CZ	8.56	1.49	1.38
1	A	132[A]	SER	CA-CB	8.55	1.65	1.52
1	A	132[B]	SER	CA-CB	8.55	1.65	1.52
1	A	132[C]	SER	CA-CB	8.55	1.65	1.52
1	A	120	GLU	CG-CD	8.54	1.64	1.51
1	A	35	GLU	CD-OE2	8.32	1.34	1.25
1	A	106	HIS	CB-CG	8.19	1.64	1.50
1	A	123	GLY	CA-C	8.18	1.65	1.51
1	A	15	TRP	CB-CG	8.17	1.65	1.50
1	A	65	GLY	C-O	8.00	1.36	1.23
1	A	72	TYR	CG-CD2	7.97	1.49	1.39
1	A	13	SER	CA-CB	7.97	1.65	1.52
1	A	18	PHE	CB-CG	7.89	1.64	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	94	GLY	CA-C	7.76	1.64	1.51
1	A	129	GLU	CD-OE1	7.31	1.33	1.25
1	A	5	GLU	CD-OE2	7.13	1.33	1.25
1	A	138	TYR	CD2-CE2	7.04	1.50	1.39
1	A	15	TRP	CZ3-CH2	7.03	1.51	1.40
1	A	144	VAL	CB-CG2	7.00	1.67	1.52
1	A	11	VAL	CB-CG1	6.99	1.67	1.52
1	A	140	GLU	CB-CG	6.96	1.65	1.52
1	A	44	PHE	CB-CG	6.93	1.63	1.51
1	A	110	VAL	CB-CG2	6.90	1.67	1.52
1	A	102	VAL	CB-CG1	6.87	1.67	1.52
1	A	46	PHE	CB-CG	6.85	1.62	1.51
1	A	128	GLU	CD-OE2	-6.82	1.18	1.25
1	A	27	HIS	CE1-NE2	6.78	1.48	1.32
1	A	85	VAL	CB-CG2	6.75	1.67	1.52
1	A	152	ALA	C-O	6.75	1.36	1.23
1	A	61	GLN	C-O	6.73	1.36	1.23
1	A	111	LYS	N-CA	6.72	1.59	1.46
1	A	104	ASP	N-CA	6.71	1.59	1.46
1	A	72	TYR	C-O	6.70	1.36	1.23
1	A	15	TRP	NE1-CE2	-6.63	1.28	1.37
1	A	28	ARG	CZ-NH1	6.62	1.41	1.33
1	A	134	TRP	N-CA	6.60	1.59	1.46
1	A	14	SER	CB-OG	-6.58	1.33	1.42
1	A	121	VAL	CB-CG2	6.57	1.66	1.52
1	A	101	GLY	CA-C	6.56	1.62	1.51
1	A	113	ALA	CA-CB	6.56	1.66	1.52
1	A	79	GLU	CB-CG	6.55	1.64	1.52
1	A	88	ALA	N-CA	6.54	1.59	1.46
1	A	68	PHE	C-O	6.53	1.35	1.23
1	A	138	TYR	CD1-CE1	6.51	1.49	1.39
1	A	35	GLU	CD-OE1	-6.49	1.18	1.25
1	A	13	SER	CB-OG	-6.40	1.33	1.42
1	A	73	GLU	CG-CD	6.40	1.61	1.51
1	A	124	ALA	N-CA	6.31	1.58	1.46
1	A	27	HIS	CG-ND1	6.30	1.52	1.38
1	A	11	VAL	N-CA	6.29	1.58	1.46
1	A	149	MET	C-O	6.26	1.35	1.23
1	A	86	THR	N-CA	6.24	1.58	1.46
1	A	145	ILE	N-CA	6.21	1.58	1.46
1	A	109	VAL	CA-CB	6.15	1.67	1.54
1	A	111	LYS	CD-CE	6.14	1.66	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	95	SER	N-CA	6.10	1.58	1.46
1	A	68	PHE	CG-CD2	6.05	1.47	1.38
1	A	6	SER	CA-CB	6.01	1.61	1.52
1	A	16	GLU	CG-CD	6.01	1.60	1.51
1	A	15	TRP	C-O	5.98	1.34	1.23
1	A	124	ALA	C-O	5.98	1.34	1.23
1	A	97	HIS	CA-CB	5.95	1.67	1.53
1	A	148	GLU	CB-CG	5.95	1.63	1.52
1	A	40	ALA	CA-CB	5.93	1.65	1.52
1	A	63	HIS	CB-CG	5.92	1.60	1.50
1	A	55	GLN	C-O	5.89	1.34	1.23
1	A	68	PHE	CE1-CZ	5.89	1.48	1.37
1	A	99	SER	N-CA	5.88	1.58	1.46
1	A	67	VAL	CB-CG2	5.88	1.65	1.52
1	A	122	VAL	N-CA	5.88	1.58	1.46
1	A	52	GLU	CG-CD	-5.83	1.43	1.51
1	A	97	HIS	CG-CD2	-5.83	1.25	1.35
1	A	75	ALA	N-CA	5.83	1.58	1.46
1	A	45	SER	CB-OG	5.81	1.49	1.42
1	A	102	VAL	N-CA	5.77	1.57	1.46
1	A	15	TRP	CD1-NE1	5.74	1.47	1.38
1	A	152	ALA	N-CA	5.72	1.57	1.46
1	A	84	VAL	CB-CG1	5.68	1.64	1.52
1	A	8	ALA	C-O	5.66	1.34	1.23
1	A	82	GLY	CA-C	5.64	1.60	1.51
1	A	90	LEU	CA-CB	5.64	1.66	1.53
1	A	77	GLN	CG-CD	5.64	1.64	1.51
1	A	120	GLU	CA-CB	5.57	1.66	1.53
1	A	111	LYS	CB-CG	5.55	1.67	1.52
1	A	18	PHE	N-CA	5.55	1.57	1.46
1	A	58	PRO	N-CD	5.53	1.55	1.47
1	A	53	VAL	CA-CB	5.50	1.66	1.54
1	A	115	LEU	N-CA	5.49	1.57	1.46
1	A	9	ALA	CA-CB	5.45	1.64	1.52
1	A	16	GLU	CD-OE1	5.45	1.31	1.25
1	A	59	GLU	CD-OE2	5.42	1.31	1.25
1	A	71	VAL	CB-CG1	5.41	1.64	1.52
1	A	125	LYS	CD-CE	5.40	1.64	1.51
1	A	99	SER	C-O	5.38	1.33	1.23
1	A	92	ASN	CB-CG	5.38	1.63	1.51
1	A	76	ILE	C-O	5.37	1.33	1.23
1	A	153	ALA	C-OXT	5.34	1.33	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	140	GLU	CD-OE2	5.34	1.31	1.25
1	A	96	VAL	CB-CG1	5.33	1.64	1.52
1	A	75	ALA	C-O	5.31	1.33	1.23
1	A	98	VAL	CB-CG2	5.28	1.64	1.52
1	A	92	ASN	C-O	5.26	1.33	1.23
1	A	4	THR	C-O	5.26	1.33	1.23
1	A	106	HIS	ND1-CE1	5.26	1.47	1.34
1	A	45	SER	N-CA	5.24	1.56	1.46
1	A	79	GLU	C-O	5.23	1.33	1.23
1	A	63	HIS	CE1-NE2	5.21	1.44	1.32
1	A	3	LEU	CA-CB	5.17	1.65	1.53
1	A	15	TRP	N-CA	5.17	1.56	1.46
1	A	91	LYS	CD-CE	5.14	1.64	1.51
1	A	68	PHE	N-CA	5.13	1.56	1.46
1	A	43	LEU	CA-CB	5.11	1.65	1.53
1	A	148	GLU	CG-CD	-5.10	1.44	1.51
1	A	133	ALA	CA-C	5.09	1.66	1.52
1	A	141	LEU	N-CA	5.08	1.56	1.46
1	A	138	TYR	N-CA	5.07	1.56	1.46
1	A	127	SER	C-O	5.05	1.32	1.23
1	A	143	ILE	CA-CB	5.05	1.66	1.54
1	A	130	LEU	N-CA	5.04	1.56	1.46
1	A	38	PRO	N-CA	5.03	1.55	1.47
1	A	85	VAL	CA-C	5.02	1.66	1.52
1	A	84	VAL	N-CA	5.01	1.56	1.46
1	A	52	GLU	N-CA	5.00	1.56	1.46

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	72	TYR	CB-CG-CD1	-10.21	114.87	121.00
1	A	28	ARG	NE-CZ-NH2	-8.96	115.82	120.30
1	A	112	GLU	OE1-CD-OE2	-8.76	112.79	123.30
1	A	15	TRP	CG-CD2-CE3	-7.64	127.03	133.90
1	A	16	GLU	OE1-CD-OE2	-7.38	114.45	123.30
1	A	73	GLU	OE1-CD-OE2	-7.24	114.61	123.30
1	A	138	TYR	CB-CG-CD2	7.16	125.30	121.00
1	A	128	GLU	OE1-CD-OE2	-6.96	114.94	123.30
1	A	108	PRO	N-CA-CB	6.92	111.61	103.30
1	A	17	GLU	OE1-CD-OE2	-6.75	115.20	123.30
1	A	29	PHE	CB-CG-CD1	-6.75	116.08	120.80
1	A	46	PHE	CB-CG-CD2	6.59	125.42	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	54	PRO	N-CA-CB	6.55	111.16	103.30
1	A	23	PRO	N-CA-CB	6.53	111.14	103.30
1	A	30	PHE	CB-CG-CD2	-6.52	116.23	120.80
1	A	120	GLU	OE1-CD-OE2	-6.41	115.61	123.30
1	A	72	TYR	CG-CD1-CE1	-6.12	116.40	121.30
1	A	29	PHE	CD1-CG-CD2	6.11	126.25	118.30
1	A	15	TRP	CH2-CZ2-CE2	-6.04	111.36	117.40
1	A	126	TRP	CE2-CD2-CG	-5.95	102.54	107.30
1	A	134	TRP	CG-CD1-NE1	-5.84	104.26	110.10
1	A	15	TRP	CD2-CE3-CZ3	-5.76	111.31	118.80
1	A	68	PHE	CB-CG-CD1	-5.73	116.79	120.80
1	A	48	LYS	CB-CA-C	-5.67	99.07	110.40
1	A	5	GLU	OE1-CD-OE2	-5.64	116.53	123.30
1	A	142	ALA	O-C-N	5.62	131.70	122.70
1	A	138	TYR	CD1-CE1-CZ	5.62	124.85	119.80
1	A	12	LYS	O-C-N	5.51	131.52	122.70
1	A	126	TRP	NE1-CE2-CD2	5.47	112.77	107.30
1	A	87	ASP	CB-CG-OD2	-5.43	113.41	118.30
1	A	149	MET	O-C-N	5.37	131.29	122.70
1	A	131	ASN	O-C-N	5.35	131.26	122.70
1	A	15	TRP	CD1-NE1-CE2	-5.33	104.20	109.00
1	A	15	TRP	NE1-CE2-CZ2	-5.31	124.56	130.40
1	A	8	ALA	O-C-N	5.26	131.11	122.70
1	A	124	ALA	CB-CA-C	-5.14	102.39	110.10
1	A	15	TRP	CD1-CG-CD2	-5.02	102.28	106.30

There are no chirality outliers.

All (8) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	17	GLU	Sidechain
1	A	35	GLU	Sidechain
1	A	42	ASP	Sidechain
1	A	55	GLN	Sidechain
1	A	56	ASN	Sidechain
1	A	61	GLN	Sidechain
1	A	87	ASP	Sidechain
1	A	92	ASN	Sidechain

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	1180	0	1200	25	16
2	A	43	0	30	7	0
3	A	8	0	5	12	0
4	A	64	0	0	0	4
All	All	1295	0	1235	30	18

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 13.

All (30) 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:30:PHE:HZ	3:A:155:NBE:H5	1.25	1.00
1:A:30:PHE:CZ	3:A:155:NBE:H5	2.17	0.73
1:A:63:HIS:HB3	3:A:155:NBE:H3	1.72	0.69
1:A:63:HIS:HB3	3:A:155:NBE:C3	2.24	0.68
2:A:154:HEM:NA	3:A:155:NBE:H2	2.10	0.66
1:A:92:ASN:O	1:A:96:VAL:HG12	2.01	0.60
1:A:30:PHE:HZ	3:A:155:NBE:C5	2.08	0.60
1:A:63:HIS:HB3	3:A:155:NBE:C4	2.32	0.59
1:A:18:PHE:CE1	1:A:25:HIS:HB3	2.40	0.56
1:A:77:GLN:NE2	1:A:85:VAL:H	2.03	0.56
1:A:63:HIS:HB3	3:A:155:NBE:H4	1.89	0.54
1:A:21:ASN:C	1:A:21:ASN:HD22	2.11	0.54
2:A:154:HEM:C4A	3:A:155:NBE:H2	2.43	0.53
1:A:139:ASP:O	1:A:143:ILE:HG13	2.13	0.48
1:A:54:PRO:HB2	1:A:57:ASN:HB2	1.95	0.48
1:A:102:VAL:HG13	2:A:154:HEM:HAC	1.95	0.47
1:A:103:ALA:O	1:A:106:HIS:HB2	2.14	0.47
2:A:154:HEM:C1A	3:A:155:NBE:H2	2.50	0.47
1:A:87:ASP:O	1:A:91:LYS:HG3	2.16	0.46
1:A:31:ILE:O	1:A:35:GLU:HG3	2.16	0.46
2:A:154:HEM:NA	3:A:155:NBE:C2	2.79	0.46
2:A:154:HEM:C1A	3:A:155:NBE:C2	2.99	0.45
1:A:100:LYS:HG3	2:A:154:HEM:HAD2	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:21:ASN:C	1:A:21:ASN:ND2	2.72	0.43
1:A:22:ILE:HD13	1:A:22:ILE:HA	1.81	0.43
1:A:93:LEU:HD23	1:A:93:LEU:HA	1.59	0.42
1:A:37:ALA:HA	1:A:38:PRO:HD2	1.97	0.42
1:A:17:GLU:OE2	1:A:122:VAL:HG12	2.20	0.42
1:A:126:TRP:CZ2	1:A:131:ASN:HB2	2.56	0.41

All (18) 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:126:TRP:O	1:A:126:TRP:O[2_555]	1.08	1.12
1:A:124:ALA:O	1:A:127:SER:CB[2_555]	1.23	0.97
1:A:124:ALA:O	1:A:127:SER:CA[2_555]	1.37	0.83
1:A:124:ALA:CB	1:A:129:GLU:N[2_555]	1.42	0.78
1:A:124:ALA:O	1:A:127:SER:C[2_555]	1.52	0.68
1:A:124:ALA:C	1:A:127:SER:CB[2_555]	1.61	0.59
1:A:124:ALA:O	1:A:128:GLU:N[2_555]	1.77	0.43
4:A:215:HOH:O	4:A:215:HOH:O[2_555]	1.77	0.43
1:A:125:LYS:CA	1:A:127:SER:CB[2_555]	1.80	0.40
4:A:190:HOH:O	4:A:215:HOH:O[2_555]	1.81	0.39
1:A:124:ALA:O	1:A:127:SER:OG[2_555]	1.83	0.37
1:A:125:LYS:CA	1:A:127:SER:OG[2_555]	1.83	0.37
1:A:10:LEU:CD1	1:A:125:LYS:CD[2_555]	1.84	0.36
1:A:125:LYS:N	1:A:127:SER:CB[2_555]	1.92	0.28
1:A:56:ASN:OD1	4:A:171:HOH:O[1_565]	1.95	0.25
1:A:125:LYS:O	4:A:215:HOH:O[2_555]	2.09	0.11
1:A:125:LYS:C	1:A:127:SER:CB[2_555]	2.10	0.10
1:A:124:ALA:C	1:A:127:SER:OG[2_555]	2.14	0.06

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%)	149 (97%)	3 (2%)	1 (1%)	19	14

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	3	LEU

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	127/125 (102%)	123 (97%)	4 (3%)	35	36

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

Mol	Chain	Res	Type
1	A	3	LEU
1	A	21	ASN
1	A	34	LEU
1	A	96	VAL

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	21	ASN
1	A	25	HIS
1	A	61	GLN
1	A	77	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	NBE	A	155	2	8,8,8	0.65	0	9,9,9	1.00	0
2	HEM	A	154	3,1	42,50,50	3.80	30 (71%)	46,82,82	2.35	20 (43%)

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	NBE	A	155	2	-	2/2/2/2	0/1/1/1
2	HEM	A	154	3,1	-	1/12/54/54	-

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	154	HEM	C1D-C2D	7.13	1.59	1.44
2	A	154	HEM	C4D-C3D	6.81	1.56	1.45
2	A	154	HEM	C3B-C4B	6.54	1.57	1.44
2	A	154	HEM	CBD-CGD	6.38	1.65	1.50
2	A	154	HEM	CAB-C3B	6.35	1.64	1.47
2	A	154	HEM	C1B-C2B	6.14	1.57	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	154	HEM	FE-NB	6.05	2.31	1.98
2	A	154	HEM	C1A-NA	5.58	1.47	1.36
2	A	154	HEM	CAD-C3D	5.05	1.64	1.51
2	A	154	HEM	CMD-C2D	4.76	1.60	1.50
2	A	154	HEM	C3C-C4C	4.69	1.48	1.41
2	A	154	HEM	C3C-CAC	4.41	1.57	1.47
2	A	154	HEM	CMB-C2B	4.38	1.59	1.50
2	A	154	HEM	C3C-C2C	3.96	1.45	1.40
2	A	154	HEM	C4A-NA	3.88	1.44	1.36
2	A	154	HEM	CAA-C2A	3.75	1.61	1.52
2	A	154	HEM	CBA-CGA	3.72	1.59	1.50
2	A	154	HEM	FE-ND	3.47	2.17	1.98
2	A	154	HEM	C4A-CHB	3.40	1.50	1.41
2	A	154	HEM	C2A-C3A	3.34	1.47	1.37
2	A	154	HEM	CMC-C2C	3.25	1.59	1.51
2	A	154	HEM	CMA-C3A	3.04	1.57	1.51
2	A	154	HEM	C1A-CHA	2.82	1.48	1.41
2	A	154	HEM	O2A-CGA	-2.53	1.22	1.30
2	A	154	HEM	O1D-CGD	2.49	1.30	1.22
2	A	154	HEM	CHB-C1B	2.47	1.40	1.34
2	A	154	HEM	O2D-CGD	-2.47	1.22	1.30
2	A	154	HEM	CHA-C4D	2.10	1.39	1.34
2	A	154	HEM	C4D-ND	2.07	1.43	1.40
2	A	154	HEM	CBD-CAD	-2.00	1.45	1.51

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	154	HEM	CMA-C3A-C4A	-5.64	120.19	128.46
2	A	154	HEM	C3B-C4B-NB	4.94	113.01	109.47
2	A	154	HEM	C4C-CHD-C1D	4.68	128.73	122.56
2	A	154	HEM	C4A-C3A-C2A	3.86	109.68	107.00
2	A	154	HEM	O2D-CGD-O1D	-3.64	113.97	123.33
2	A	154	HEM	C4B-C3B-C2B	-3.40	104.16	107.28
2	A	154	HEM	CHC-C4B-NB	-3.22	120.97	124.44
2	A	154	HEM	CHD-C1D-ND	-3.17	121.03	124.44
2	A	154	HEM	C2C-C3C-C4C	3.13	109.08	106.90
2	A	154	HEM	C3C-C4C-NC	-3.10	105.10	110.94
2	A	154	HEM	O2A-CGA-O1A	-3.04	115.51	123.33
2	A	154	HEM	C4B-CHC-C1C	2.80	126.25	122.56
2	A	154	HEM	CMA-C3A-C2A	2.75	130.12	124.94
2	A	154	HEM	CAD-C3D-C4D	2.70	129.41	124.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	154	HEM	C3D-C4D-ND	2.49	112.90	110.17
2	A	154	HEM	CAA-CBA-CGA	-2.47	107.18	113.83
2	A	154	HEM	CHA-C4D-ND	-2.44	121.35	124.37
2	A	154	HEM	CMC-C2C-C3C	2.35	129.38	124.68
2	A	154	HEM	C2B-C1B-NB	2.26	112.44	109.84
2	A	154	HEM	CHB-C1B-NB	-2.15	121.71	124.37

There are no chirality outliers.

All (3) torsion outliers are listed below:

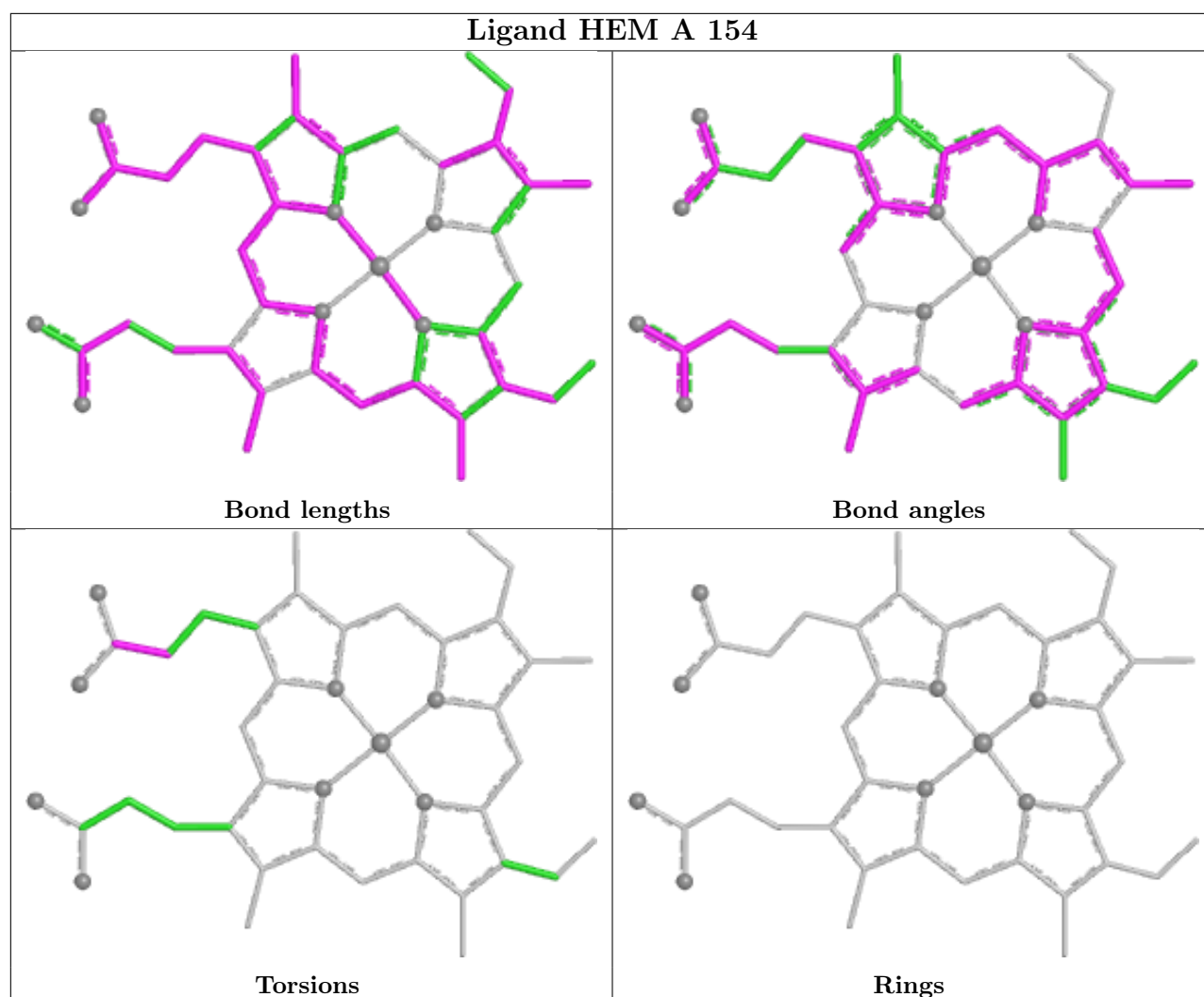
Mol	Chain	Res	Type	Atoms
3	A	155	NBE	C2-C1-N-O
3	A	155	NBE	C6-C1-N-O
2	A	154	HEM	CAD-CBD-CGD-O1D

There are no ring outliers.

2 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	155	NBE	12	0
2	A	154	HEM	7	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 ⓘ

Warning: The R factor obtained from EDS is 0.4272, which does not match the depositor's R factor of 0.0. Please interpret the results in this section carefully.

In the following table, the column labelled '#RSRZ > 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q < 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	153/153 (100%)	3.30	118 (77%) 0 0	6, 18, 41, 54	20 (13%)

All (118) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	152	ALA	10.5
1	A	123	GLY	8.6
1	A	45	SER	8.3
1	A	83	VAL	7.1
1	A	149	MET	7.1
1	A	9	ALA	7.1
1	A	2	ALA	6.9
1	A	126	TRP	6.9
1	A	50	THR	6.8
1	A	48	LYS	6.5
1	A	153	ALA	6.1
1	A	49	GLY	6.1
1	A	88	ALA	6.1
1	A	58	PRO	5.9
1	A	86	THR	5.9
1	A	1	GLY	5.7
1	A	128	GLU	5.5
1	A	143	ILE	5.5
1	A	53	VAL	5.5
1	A	92	ASN	5.4
1	A	99	SER	5.2
1	A	59	GLU	5.2
1	A	144	VAL	5.1
1	A	4	THR	5.1

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Mol	Chain	Res	Type	RSRZ
1	A	14	SER	5.0
1	A	42	ASP	4.9
1	A	79	GLU	4.9
1	A	147	LYS	4.9
1	A	23	PRO	4.8
1	A	20	ALA	4.8
1	A	15	TRP	4.7
1	A	136	ILE	4.7
1	A	151	ASP	4.6
1	A	91	LYS	4.4
1	A	5	GLU	4.4
1	A	89	THR	4.3
1	A	3	LEU	4.3
1	A	108	PRO	4.2
1	A	85	VAL	4.2
1	A	61	GLN	4.1
1	A	44	PHE	4.1
1	A	81	THR	3.9
1	A	11	VAL	3.9
1	A	6	SER	3.8
1	A	64	ALA	3.8
1	A	98	VAL	3.8
1	A	107	PHE	3.7
1	A	146	LYS	3.7
1	A	130	LEU	3.7
1	A	87	ASP	3.7
1	A	90	LEU	3.7
1	A	104	ASP	3.7
1	A	18	PHE	3.4
1	A	97	HIS	3.4
1	A	117	THR	3.4
1	A	96	VAL	3.4
1	A	54	PRO	3.4
1	A	129	GLU	3.3
1	A	124	ALA	3.3
1	A	68	PHE	3.2
1	A	62	ALA	3.2
1	A	63	HIS	3.2
1	A	56	ASN	3.2
1	A	12	LYS	3.1
1	A	32	LEU	3.1
1	A	65	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	145	ILE	3.1
1	A	72	TYR	3.1
1	A	13	SER	3.1
1	A	21	ASN	3.0
1	A	78	LEU	3.0
1	A	51	SER	3.0
1	A	25	HIS	3.0
1	A	10	LEU	3.0
1	A	109	VAL	2.9
1	A	142	ALA	2.9
1	A	39	ALA	2.9
1	A	133	ALA	2.8
1	A	57	ASN	2.8
1	A	121	VAL	2.8
1	A	132[A]	SER	2.8
1	A	41	LYS	2.8
1	A	38	PRO	2.8
1	A	47	LEU	2.8
1	A	95	SER	2.8
1	A	119	LYS	2.8
1	A	31	ILE	2.8
1	A	82	GLY	2.8
1	A	43	LEU	2.7
1	A	114	ILE	2.7
1	A	118	ILE	2.7
1	A	16	GLU	2.7
1	A	66	LYS	2.7
1	A	22	ILE	2.7
1	A	55	GLN	2.6
1	A	33	VAL	2.6
1	A	74	ALA	2.5
1	A	71	VAL	2.5
1	A	34	LEU	2.5
1	A	80	VAL	2.5
1	A	46	PHE	2.5
1	A	94	GLY	2.4
1	A	73	GLU	2.3
1	A	120	GLU	2.3
1	A	148	GLU	2.3
1	A	30	PHE	2.3
1	A	137	ALA	2.3
1	A	67	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	100	LYS	2.2
1	A	26	THR	2.2
1	A	135	THR	2.2
1	A	37	ALA	2.2
1	A	102	VAL	2.2
1	A	84	VAL	2.1
1	A	110	VAL	2.1
1	A	122	VAL	2.1
1	A	52	GLU	2.0
1	A	140	GLU	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](#)

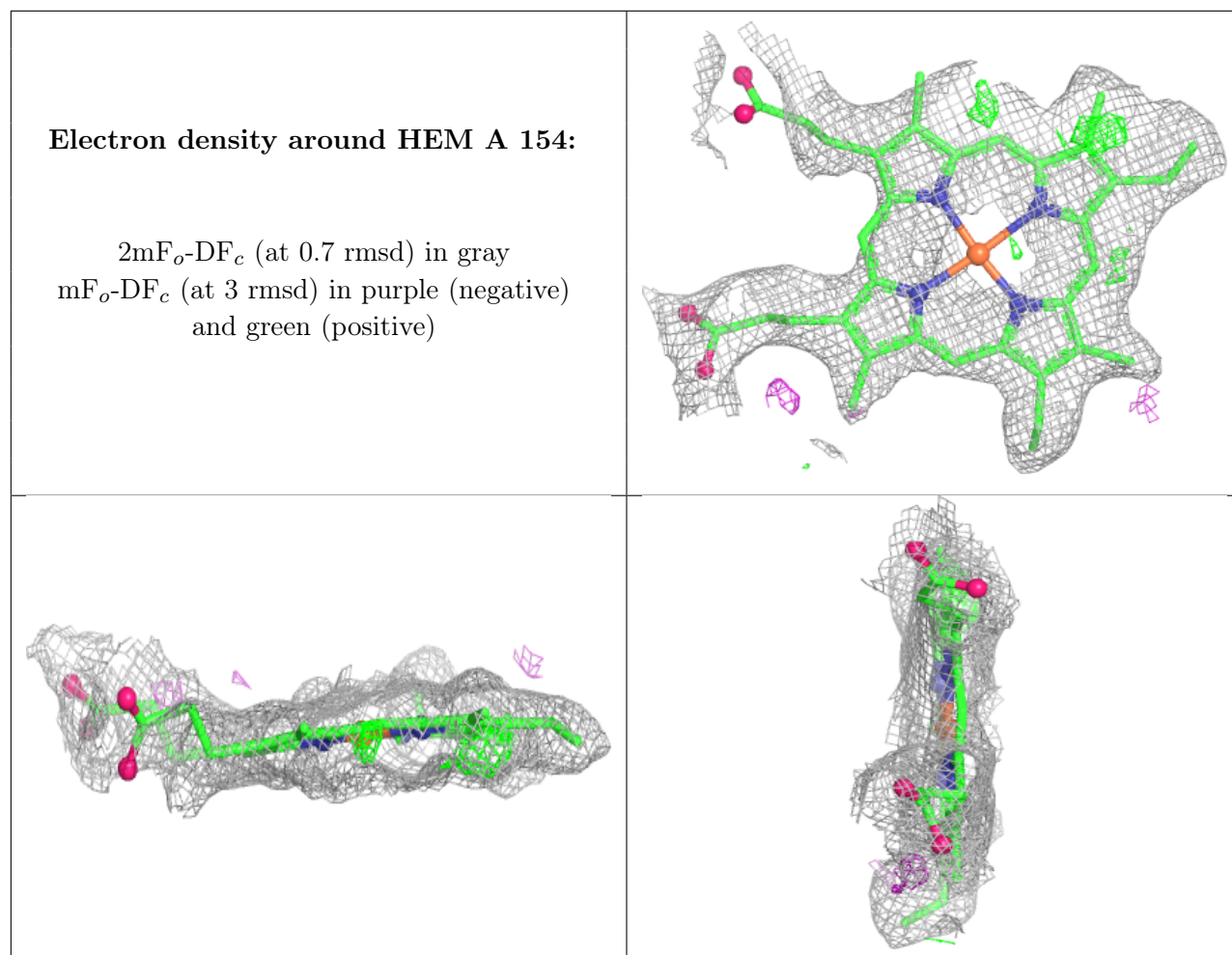
There are no monosaccharides in this entry.

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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NBE	A	155	8/8	0.53	0.26	19,33,53,58	0
2	HEM	A	154	43/43	0.83	0.16	0,18,43,56	3

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 ⓘ

There are no such residues in this entry.