



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 01:31 PM GMT

PDB ID : 3BCQ
Title : Crystal structure of oxy-hemoglobin from Brycon cephalus
Authors : Poy, C.D.; Leopoldino, A.M.; Rahal, P.; de Azevedo, W.F.; Rodriguez, G.O.B.;
Murakami, M.T.
Deposited on : 2007-11-13
Resolution : 2.40 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

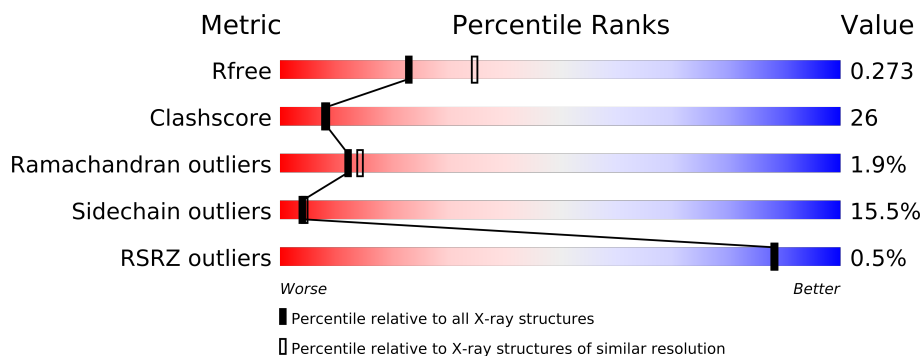
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.40 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}	66092	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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 ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	142	
1	C	142	
2	B	146	
2	D	146	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4772 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Alpha-chain hemoglobin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	142	Total	C	N	O	S	0	1	0
			1124	725	187	209	3			
1	C	142	Total	C	N	O	S	0	2	0
			1128	729	186	210	3			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	SER	ALA	SEE REMARK 999	UNP A1YZP4
A	32	MET	THR	SEE REMARK 999	UNP A1YZP4
A	65	GLY	SER	SEE REMARK 999	UNP A1YZP4
A	66	SER	GLY	SEE REMARK 999	UNP A1YZP4
A	112	GLN	ILE	SEE REMARK 999	UNP A1YZP4
A	132	LEU	ALA	SEE REMARK 999	UNP A1YZP4
A	134	SER	ALA	SEE REMARK 999	UNP A1YZP4
A	135	TRP	LEU	SEE REMARK 999	UNP A1YZP4
A	136	SER	ALA	SEE REMARK 999	UNP A1YZP4
C	9	SER	ALA	SEE REMARK 999	UNP A1YZP4
C	32	MET	THR	SEE REMARK 999	UNP A1YZP4
C	65	GLY	SER	SEE REMARK 999	UNP A1YZP4
C	66	SER	GLY	SEE REMARK 999	UNP A1YZP4
C	112	GLN	ILE	SEE REMARK 999	UNP A1YZP4
C	132	LEU	ALA	SEE REMARK 999	UNP A1YZP4
C	134	SER	ALA	SEE REMARK 999	UNP A1YZP4
C	135	TRP	LEU	SEE REMARK 999	UNP A1YZP4
C	136	SER	ALA	SEE REMARK 999	UNP A1YZP4

- Molecule 2 is a protein called Beta-chain hemoglobin.

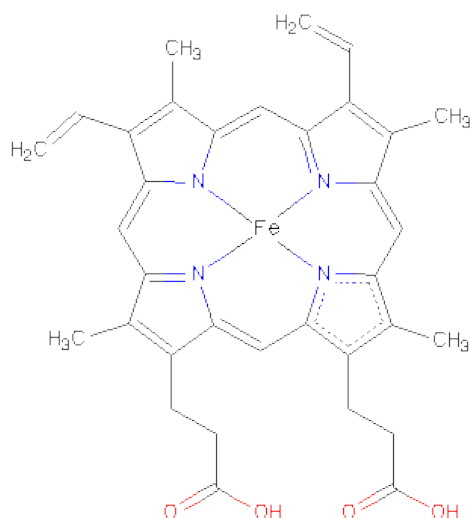
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1125	721	198	200	6			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	146	Total	C	N	O	S	0	0	0
			1125	721	198	200	6			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 4 is OXYGEN MOLECULE (three-letter code: OXY) (formula: O_2).

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

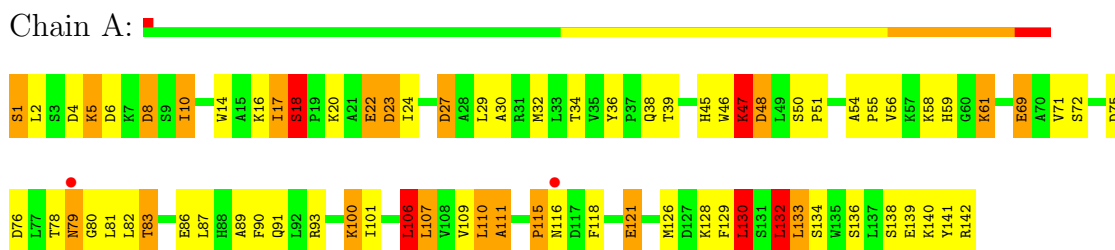
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	31	Total O 31 31	0	0
5	B	13	Total O 13 13	0	0
5	C	25	Total O 25 25	0	0
5	D	21	Total O 21 21	0	0

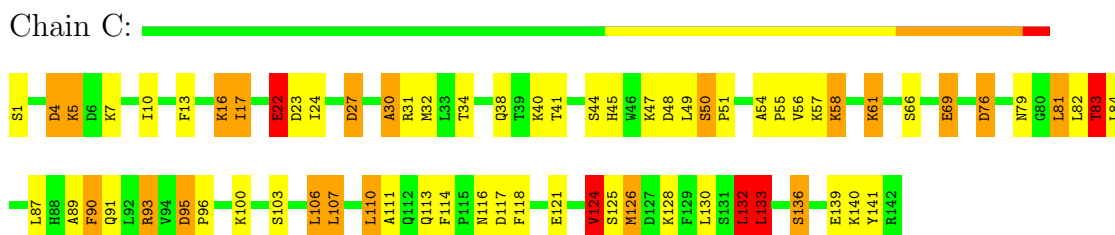
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

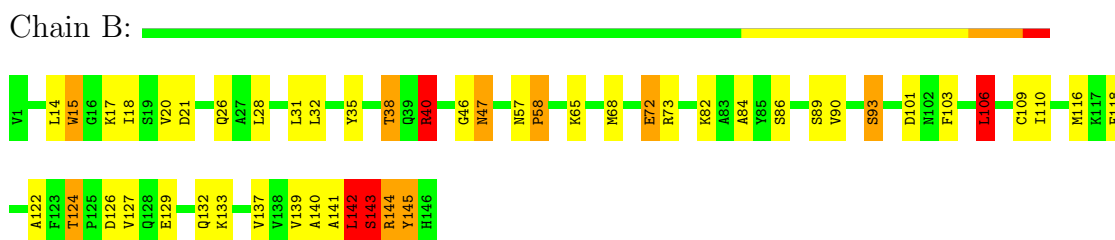
- Molecule 1: Alpha-chain hemoglobin



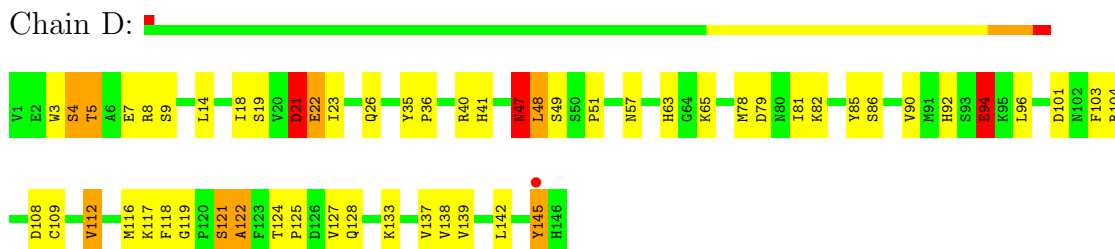
- Molecule 1: Alpha-chain hemoglobin



- Molecule 2: Beta-chain hemoglobin



- Molecule 2: Beta-chain hemoglobin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	119.37Å 54.04Å 101.91Å 90.00° 113.89° 90.00°	Depositor
Resolution (Å)	19.89 – 2.40 19.89 – 2.40	Depositor EDS
% Data completeness (in resolution range)	86.9 (19.89-2.40) 86.9 (19.89-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.61 (at 2.41Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.166 , 0.279 0.164 , 0.273	Depositor DCC
R_{free} test set	1042 reflections (5.39%)	DCC
Wilson B-factor (Å ²)	40.4	Xtriage
Anisotropy	0.093	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 33.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 20358 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4772	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.69% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	1.42	5/1154 (0.4%)	1.45	16/1561 (1.0%)
1	C	1.44	9/1161 (0.8%)	1.32	17/1572 (1.1%)
2	B	1.22	2/1154 (0.2%)	1.15	6/1567 (0.4%)
2	D	1.22	4/1154 (0.3%)	1.10	6/1567 (0.4%)
All	All	1.33	20/4623 (0.4%)	1.26	45/6267 (0.7%)

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	1	3
2	B	0	2
2	D	0	1
All	All	1	6

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	30	ALA	CA-CB	-8.68	1.34	1.52
1	C	22	GLU	CG-CD	-6.91	1.41	1.51
1	C	27	ASP	CB-CG	6.89	1.66	1.51
1	C	4	ASP	CB-CG	6.83	1.66	1.51
1	A	27	ASP	CB-CG	6.68	1.65	1.51
1	A	22	GLU	CG-CD	-6.58	1.42	1.51
1	C	136	SER	CB-OG	6.51	1.50	1.42
2	B	109	CYS	CB-SG	-6.24	1.71	1.82
1	C	136	SER	CA-CB	6.13	1.62	1.52
1	A	69	GLU	CB-CG	5.91	1.63	1.52
1	A	93	ARG	CG-CD	5.78	1.66	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	94	GLU	CB-CG	5.75	1.63	1.52
1	C	83	THR	CA-CB	5.61	1.68	1.53
1	C	121	GLU	CB-CG	5.57	1.62	1.52
2	D	127	VAL	CB-CG2	5.22	1.63	1.52
2	B	101	ASP	CB-CG	5.13	1.62	1.51
2	D	112	VAL	CB-CG2	5.12	1.63	1.52
2	D	109	CYS	CB-SG	-5.11	1.73	1.81
1	C	124	VAL	CB-CG1	-5.07	1.42	1.52
1	A	36	TYR	CD2-CE2	5.04	1.47	1.39

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	106	LEU	CB-CG-CD2	-18.14	80.15	111.00
1	C	106	LEU	CB-CG-CD2	-10.85	92.56	111.00
1	A	29	LEU	CB-CG-CD2	-9.98	94.04	111.00
1	C	27	ASP	CB-CG-OD1	9.62	126.96	118.30
1	A	93	ARG	NE-CZ-NH1	9.26	124.93	120.30
1	A	93	ARG	NE-CZ-NH2	-8.08	116.26	120.30
1	A	107	LEU	CB-CG-CD1	7.99	124.59	111.00
1	A	110	LEU	CA-CB-CG	-7.82	97.31	115.30
1	A	110	LEU	C-N-CA	-7.60	102.70	121.70
1	C	133	LEU	CB-CG-CD2	-7.57	98.13	111.00
1	A	17	ILE	C-N-CA	-7.33	103.36	121.70
2	D	121	SER	C-N-CA	6.99	139.18	121.70
1	C	132	LEU	CA-CB-CG	-6.97	99.28	115.30
1	A	48	ASP	N-CA-C	6.66	128.97	111.00
1	A	130	LEU	CA-CB-CG	6.43	130.09	115.30
2	B	106	LEU	CB-CG-CD2	6.31	121.72	111.00
1	C	110	LEU	CB-CG-CD1	6.23	121.60	111.00
1	A	23	ASP	CB-CG-OD1	-6.13	112.78	118.30
2	D	121	SER	O-C-N	-6.00	113.11	122.70
2	D	121	SER	N-CA-C	5.97	127.11	111.00
2	B	40	ARG	NE-CZ-NH1	-5.75	117.43	120.30
2	D	121	SER	CA-C-N	5.66	129.64	117.20
2	B	145	TYR	CA-CB-CG	-5.65	102.67	113.40
1	C	76	ASP	CB-CG-OD1	-5.58	113.28	118.30
1	C	133	LEU	CA-CB-CG	5.54	128.05	115.30
1	A	130	LEU	CB-CG-CD1	5.52	120.38	111.00
1	A	132	LEU	CA-CB-CG	-5.51	102.62	115.30
1	A	27	ASP	CB-CG-OD2	5.44	123.19	118.30
1	C	132	LEU	CB-CG-CD1	5.43	120.23	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	142	LEU	N-CA-C	5.38	125.52	111.00
2	D	47	ASN	C-N-CA	5.36	135.11	121.70
1	C	95	ASP	CB-CG-OD1	5.32	123.09	118.30
1	C	107[A]	LEU	CB-CG-CD1	5.32	120.04	111.00
1	C	107[B]	LEU	CB-CG-CD1	5.32	120.04	111.00
1	A	47	LYS	C-N-CA	5.22	134.76	121.70
1	C	107[A]	LEU	CB-CG-CD2	-5.21	102.14	111.00
1	C	107[B]	LEU	CB-CG-CD2	-5.21	102.14	111.00
2	B	47	ASN	N-CA-CB	5.20	119.95	110.60
2	B	116	MET	CG-SD-CE	5.15	108.44	100.20
1	A	27	ASP	CB-CA-C	5.14	120.68	110.40
1	C	106	LEU	CA-CB-CG	-5.10	103.57	115.30
2	D	21	ASP	N-CA-C	-5.09	97.26	111.00
1	C	133	LEU	N-CA-CB	-5.07	100.26	110.40
1	C	126	MET	CB-CG-SD	-5.05	97.26	112.40
1	C	69	GLU	N-CA-CB	5.04	119.67	110.60

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	48	ASP	CA

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	115	PRO	Peptide
1	A	18	SER	Peptide
1	A	47	LYS	Peptide
2	B	142	LEU	Peptide
2	B	46	GLY	Peptide
2	D	47	ASN	Peptide

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1124	0	1144	80	0
1	C	1128	0	1145	81	0
2	B	1125	0	1126	36	0
2	D	1125	0	1126	45	0
3	A	43	0	30	5	0
3	B	43	0	30	0	0
3	C	43	0	30	10	0
3	D	43	0	30	1	0
4	A	2	0	0	3	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	1	0
5	A	31	0	0	10	0
5	B	13	0	0	2	0
5	C	25	0	0	11	0
5	D	21	0	0	3	0
All	All	4772	0	4661	246	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 26.

All (246) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:110:LEU:O	1:A:111:ALA:CB	1.80	1.15
3:C:143:HEM:HBA2	3:C:143:HEM:HHA	1.23	1.12
1:A:4:ASP:O	1:A:8:ASP:OD2	1.67	1.10
1:A:110:LEU:H	1:A:110:LEU:HD12	1.10	1.09
2:B:21:ASP:OD1	2:B:65:LYS:HE2	1.56	1.05
1:A:106:LEU:O	1:A:110:LEU:CD1	2.05	1.04
1:A:79:ASN:HD22	1:A:80:GLY:N	1.61	0.98
1:C:30:ALA:HB2	1:C:56:VAL:HG11	1.45	0.96
1:C:22:GLU:OE1	1:C:61:LYS:HG3	1.66	0.95
2:B:144:ARG:HH11	2:B:144:ARG:HG3	1.29	0.94
1:A:110:LEU:HD12	1:A:110:LEU:N	1.73	0.92
2:B:21:ASP:OD1	2:B:65:LYS:CE	2.19	0.90
1:A:110:LEU:O	1:A:111:ALA:HB3	1.08	0.89
3:C:143:HEM:HBA2	3:C:143:HEM:CHA	1.97	0.88
2:D:82:LYS:HE3	2:D:145:TYR:HD2	1.38	0.88
1:C:107[A]:LEU:HD22	1:C:126:MET:CE	2.04	0.88
1:A:121:GLU:HG3	5:A:157:HOH:O	1.75	0.86
1:A:86:GLU:OE2	1:A:90:PHE:CD1	2.30	0.85
2:B:15:TRP:CH2	2:B:72:GLU:HB2	2.11	0.85
1:A:106:LEU:O	1:A:110:LEU:HD12	1.76	0.84

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:106:LEU:O	1:A:110:LEU:HD13	1.78	0.84
2:D:82:LYS:HE3	2:D:145:TYR:CD2	2.13	0.83
2:B:139:VAL:O	2:B:143:SER:HB2	1.79	0.82
1:C:136:SER:O	1:C:139[A]:GLU:HG2	1.80	0.81
1:A:100:LYS:HE2	1:A:100:LYS:HA	1.62	0.81
1:C:111:ALA:HB2	1:C:118:PHE:CD1	2.15	0.81
1:A:87:LEU:HD12	1:A:91:GLN:HB3	1.63	0.81
1:C:107[A]:LEU:HD22	1:C:126:MET:HE1	1.60	0.81
1:A:6:ASP:O	1:A:10:ILE:HG23	1.81	0.81
1:C:107[B]:LEU:HD12	1:C:118:PHE:CE1	2.16	0.81
3:C:143:HEM:HBB2	3:C:143:HEM:CMB	2.09	0.80
2:B:144:ARG:NH1	2:B:144:ARG:HG3	1.95	0.80
2:D:21:ASP:O	2:D:22:GLU:HB3	1.81	0.80
2:B:35:TYR:O	2:B:38:THR:HB	1.82	0.80
1:A:86:GLU:OE2	1:A:90:PHE:HD1	1.66	0.78
1:C:47:LYS:N	1:C:47:LYS:HD3	1.97	0.78
1:C:22:GLU:OE1	1:C:61:LYS:CG	2.31	0.78
1:C:17:ILE:HD11	1:C:114:PHE:CG	2.19	0.78
2:B:89:SER:HA	2:B:142:LEU:O	1.85	0.77
3:A:143:HEM:NB	4:A:144:OXY:O1	2.17	0.76
2:D:47:ASN:O	2:D:57:ASN:ND2	2.18	0.76
2:D:117:LYS:HD3	2:D:118:PHE:CZ	2.21	0.76
2:D:92:HIS:HA	2:D:96:LEU:HD12	1.66	0.76
1:A:32:MET:HE2	1:A:39:THR:HG21	1.66	0.76
2:B:141:ALA:HA	2:B:145:TYR:HE1	1.50	0.76
1:A:79:ASN:HD22	1:A:79:ASN:C	1.90	0.75
2:D:22:GLU:HA	5:D:162:HOH:O	1.87	0.74
1:A:10:ILE:HD11	1:A:71:VAL:HG22	1.70	0.74
1:C:107[B]:LEU:CD1	1:C:118:PHE:CE1	2.72	0.72
1:A:30:ALA:HB1	1:A:51:PRO:HA	1.71	0.72
1:C:124:VAL:O	1:C:128:LYS:HG3	1.88	0.72
2:B:140:ALA:O	2:B:143:SER:HB3	1.90	0.72
1:A:61:LYS:HD3	5:A:153:HOH:O	1.88	0.72
1:C:107[B]:LEU:CD1	1:C:118:PHE:HE1	2.03	0.71
1:A:61:LYS:CD	5:A:153:HOH:O	2.41	0.69
1:A:106:LEU:C	1:A:106:LEU:HD23	2.13	0.69
1:C:111:ALA:HB1	2:D:116:MET:CG	2.23	0.69
2:B:140:ALA:O	2:B:143:SER:CB	2.41	0.69
1:C:128:LYS:O	1:C:132:LEU:HD12	1.93	0.68
1:A:100:LYS:HE2	1:A:100:LYS:CA	2.24	0.67
2:B:141:ALA:HA	2:B:145:TYR:CE1	2.29	0.66
1:A:45:HIS:CE1	3:A:143:HEM:O1D	2.49	0.66

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:136:SER:O	1:A:139:GLU:HG2	1.94	0.66
1:C:31:ARG:HD3	2:D:128:GLN:OE1	1.95	0.66
1:A:110:LEU:CD1	1:A:110:LEU:H	1.98	0.65
2:B:124:THR:HG21	5:B:161:HOH:O	1.97	0.65
2:B:18:ILE:HD13	2:B:118:PHE:HZ	1.62	0.64
2:D:133:LYS:O	2:D:137:VAL:HG23	1.98	0.64
1:A:30:ALA:HB2	1:A:56:VAL:HG11	1.80	0.63
1:A:32:MET:CE	1:A:39:THR:HG21	2.28	0.63
2:D:21:ASP:O	2:D:22:GLU:CB	2.40	0.63
2:B:73:ARG:NH2	2:B:84:ALA:O	2.30	0.63
1:C:118:PHE:CZ	1:C:126:MET:CE	2.82	0.62
1:A:45:HIS:HE1	3:A:143:HEM:O1D	1.81	0.62
1:C:111:ALA:HB1	2:D:116:MET:HG3	1.82	0.61
3:C:143:HEM:HBB2	3:C:143:HEM:HMB1	1.81	0.61
3:C:143:HEM:HHA	3:C:143:HEM:CBA	2.11	0.60
1:A:32:MET:HG3	1:A:101:ILE:HG22	1.82	0.60
2:D:63:HIS:NE2	4:D:148:OXY:O1	2.31	0.60
1:C:118:PHE:CZ	1:C:126:MET:HE1	2.37	0.60
2:D:19:SER:H	2:D:117:LYS:HE3	1.65	0.60
1:A:128:LYS:O	1:A:132:LEU:HD12	2.03	0.59
1:A:107:LEU:CD2	1:A:126:MET:HG3	2.32	0.59
1:C:81:LEU:HD23	1:C:136:SER:HB2	1.84	0.59
2:D:51:PRO:HD3	5:D:166:HOH:O	2.03	0.59
1:C:87:LEU:CD2	3:C:143:HEM:HBA1	2.32	0.58
1:A:79:ASN:ND2	1:A:80:GLY:N	2.43	0.58
2:D:8:ARG:NH2	2:D:79:ASP:OD1	2.32	0.58
3:A:143:HEM:HAA1	5:A:160:HOH:O	2.03	0.58
1:C:58:LYS:CG	5:C:166:HOH:O	2.51	0.58
2:B:129:GLU:O	2:B:129:GLU:HG2	2.01	0.57
1:C:93:ARG:HH11	1:C:93:ARG:HG2	1.69	0.57
2:B:118:PHE:HB3	2:B:122:ALA:HB3	1.86	0.57
2:D:21:ASP:HA	2:D:65:LYS:HE2	1.87	0.57
2:D:47:ASN:HD21	2:D:49:SER:CB	2.18	0.57
1:C:58:LYS:HG2	5:C:166:HOH:O	2.04	0.57
3:C:143:HEM:CBA	3:C:143:HEM:CHA	2.76	0.56
2:B:124:THR:HG22	2:B:127:VAL:H	1.70	0.56
1:C:40:LYS:HG2	1:C:49:LEU:HD23	1.87	0.56
1:C:17:ILE:HD11	1:C:114:PHE:CD2	2.40	0.56
1:C:13:PHE:HE2	1:C:126:MET:SD	2.28	0.56
1:A:111:ALA:HB2	1:A:118:PHE:CD1	2.40	0.56
1:C:107[A]:LEU:H	1:C:107[A]:LEU:HD23	1.69	0.55
1:C:17:ILE:HG12	1:C:110:LEU:HD22	1.87	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:14:LEU:C	2:B:14:LEU:HD23	2.26	0.55
1:C:83:THR:HG22	5:C:161:HOH:O	2.05	0.55
1:A:86:GLU:OE2	1:A:90:PHE:CE1	2.60	0.55
1:C:107[A]:LEU:HD22	1:C:126:MET:HE3	1.88	0.55
1:C:93:ARG:CG	1:C:93:ARG:HH11	2.20	0.55
1:C:118:PHE:CE2	1:C:126:MET:HE2	2.42	0.55
2:D:19:SER:O	2:D:21:ASP:O	2.25	0.55
1:A:81:LEU:HD11	1:A:133:LEU:HD22	1.89	0.54
2:B:21:ASP:OD1	2:B:65:LYS:HE3	2.04	0.54
1:A:89:ALA:HB2	1:A:141:TYR:CD1	2.42	0.54
1:A:110:LEU:HD22	1:A:126:MET:HE2	1.90	0.54
1:C:118:PHE:CZ	1:C:126:MET:HE2	2.43	0.54
1:C:118:PHE:HZ	1:C:126:MET:CE	2.21	0.53
2:D:121:SER:N	2:D:122:ALA:HB3	2.23	0.53
1:A:115:PRO:HG3	5:A:175:HOH:O	2.08	0.53
1:A:83:THR:HG23	5:A:146:HOH:O	2.08	0.53
2:D:101:ASP:OD2	2:D:104:ARG:NH2	2.38	0.52
1:C:118:PHE:CE2	1:C:126:MET:CE	2.93	0.52
1:A:87:LEU:CD1	1:A:91:GLN:HB3	2.37	0.52
2:B:103:PHE:HE2	2:B:143:SER:H	1.54	0.52
1:C:89:ALA:HB1	1:C:141:TYR:CD1	2.45	0.52
2:D:23:ILE:HD11	2:D:117:LYS:HD2	1.91	0.52
1:A:139:GLU:HG3	5:A:145:HOH:O	2.10	0.51
2:D:4:SER:OG	2:D:7:GLU:HG3	2.10	0.51
1:A:86:GLU:HG2	1:A:140:LYS:NZ	2.26	0.51
1:A:20:LYS:O	1:A:23:ASP:HB2	2.10	0.51
1:C:16:LYS:NZ	1:C:117:ASP:OD2	2.44	0.51
1:A:87:LEU:CD1	1:A:91:GLN:HE21	2.24	0.51
1:A:142:ARG:HB3	2:D:36:PRO:HG3	1.93	0.51
1:A:111:ALA:HB2	1:A:118:PHE:CG	2.46	0.51
1:C:13:PHE:O	1:C:17:ILE:HB	2.11	0.51
1:A:100:LYS:HE2	1:A:100:LYS:O	2.11	0.51
1:A:1:SER:HB3	1:A:132:LEU:HD11	1.93	0.50
1:C:1:SER:HA	1:C:132:LEU:HD21	1.92	0.50
1:A:2:LEU:HG	1:A:132:LEU:HD13	1.93	0.50
1:C:34:THR:HG23	5:C:150:HOH:O	2.11	0.50
2:D:22:GLU:O	2:D:26:GLN:HG3	2.11	0.50
1:A:107:LEU:HD21	1:A:130:LEU:CD2	2.42	0.49
1:A:8:ASP:OD2	1:A:8:ASP:N	2.33	0.49
1:A:59:HIS:NE2	4:A:144:OXY:O2	2.32	0.49
2:B:140:ALA:O	2:B:143:SER:HB2	2.12	0.49
1:A:107:LEU:HD22	1:A:126:MET:HG3	1.95	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:50:SER:HB2	1:A:51:PRO:HD2	1.93	0.49
1:A:106:LEU:O	1:A:106:LEU:HD23	2.12	0.49
2:D:103:PHE:HB3	2:D:139:VAL:HG22	1.95	0.48
1:C:45:HIS:CD2	1:C:45:HIS:H	2.32	0.48
2:D:85:TYR:CD2	2:D:142:LEU:HG	2.48	0.48
1:C:100:LYS:HB2	5:C:149:HOH:O	2.12	0.48
1:C:17:ILE:CD1	1:C:114:PHE:CG	2.95	0.48
1:C:54:ALA:HA	1:C:57:LYS:HE3	1.94	0.48
1:A:86:GLU:OE1	1:A:140:LYS:HD2	2.14	0.48
2:B:133:LYS:O	2:B:137:VAL:HG23	2.13	0.48
3:C:143:HEM:HHB	3:C:143:HEM:HMB2	1.35	0.47
1:A:30:ALA:CB	1:A:51:PRO:HA	2.42	0.47
1:C:30:ALA:HB2	1:C:56:VAL:CG1	2.31	0.47
1:A:1:SER:CB	1:A:132:LEU:HD11	2.44	0.47
1:C:51:PRO:HD3	5:C:150:HOH:O	2.14	0.47
2:D:47:ASN:ND2	2:D:49:SER:H	2.13	0.47
1:C:13:PHE:CE2	1:C:126:MET:SD	3.06	0.47
1:C:107[A]:LEU:HD21	1:C:130:LEU:HD12	1.97	0.47
1:A:46:TRP:HA	1:A:47:LYS:HD3	1.96	0.47
1:A:79:ASN:ND2	1:A:79:ASN:C	2.64	0.47
1:C:47:LYS:N	1:C:47:LYS:CD	2.70	0.47
2:D:85:TYR:HD2	2:D:142:LEU:HG	1.80	0.47
1:A:24:ILE:HA	1:A:109:VAL:HG11	1.97	0.47
2:D:124:THR:O	2:D:125:PRO:C	2.50	0.47
2:B:32:LEU:HD23	2:B:38:THR:CG2	2.45	0.46
1:C:132:LEU:HD22	5:C:163:HOH:O	2.14	0.46
2:B:18:ILE:HD13	2:B:118:PHE:CZ	2.45	0.46
1:A:86:GLU:HG2	1:A:140:LYS:HZ3	1.80	0.46
1:C:118:PHE:HZ	1:C:126:MET:HE1	1.76	0.46
1:C:24:ILE:HD13	1:C:113:GLN:OE1	2.16	0.46
2:B:32:LEU:HD23	2:B:38:THR:HG23	1.98	0.46
2:D:117:LYS:HD3	2:D:118:PHE:CE2	2.51	0.45
1:C:87:LEU:CD2	3:C:143:HEM:CBA	2.95	0.45
2:D:5:THR:HG23	5:D:150:HOH:O	2.16	0.45
1:A:86:GLU:CG	1:A:140:LYS:HZ3	2.29	0.45
1:C:31:ARG:O	1:C:32:MET:C	2.53	0.45
2:D:119:GLY:O	2:D:122:ALA:HB3	2.17	0.45
2:D:3:TRP:CE2	2:D:78:MET:HE2	2.52	0.45
1:C:95:ASP:HA	1:C:96:PRO:HD2	1.82	0.45
1:A:5:LYS:HA	1:A:8:ASP:OD2	2.17	0.45
1:A:87:LEU:HD13	1:A:91:GLN:HE21	1.82	0.45
2:B:31:LEU:HG	2:B:38:THR:HG21	1.99	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:18:SER:N	1:A:20:LYS:H	2.15	0.45
1:A:2:LEU:HD21	1:A:129:PHE:HA	1.99	0.44
1:A:14:TRP:O	1:A:17:ILE:O	2.35	0.44
3:A:143:HEM:NA	4:A:144:OXY:O1	2.50	0.44
2:B:40:ARG:NH1	1:C:41:THR:HG23	2.33	0.44
2:D:90:VAL:O	2:D:94:GLU:HB2	2.18	0.44
1:C:66:SER:HB2	3:C:143:HEM:HMA2	2.00	0.44
2:B:28:LEU:O	2:B:32:LEU:HG	2.17	0.44
1:C:81:LEU:HD21	1:C:133:LEU:HD22	1.99	0.44
2:B:15:TRP:CZ3	2:B:72:GLU:OE1	2.71	0.44
1:C:10:ILE:HD11	1:C:125:SER:C	2.38	0.44
1:C:47:LYS:HE2	5:C:162:HOH:O	2.17	0.44
2:B:86:SER:O	2:B:90:VAL:HG23	2.18	0.43
2:D:18:ILE:N	2:D:18:ILE:HD13	2.33	0.43
1:C:130:LEU:O	1:C:133:LEU:HB3	2.18	0.43
2:B:106:LEU:O	2:B:110:ILE:HG13	2.19	0.43
1:C:90:PHE:CD1	1:C:140:LYS:HB3	2.53	0.43
1:A:54:ALA:HB3	1:A:55:PRO:HD3	2.01	0.43
1:C:107[B]:LEU:HD11	2:D:112:VAL:HG11	2.00	0.43
1:A:47:LYS:N	1:A:47:LYS:HD3	2.34	0.43
1:A:46:TRP:CH2	1:A:59:HIS:CB	3.02	0.43
2:B:20:VAL:HG22	2:B:68:MET:HB3	2.00	0.43
1:C:54:ALA:HB3	1:C:55:PRO:HD3	2.01	0.42
1:A:4:ASP:O	1:A:8:ASP:CG	2.52	0.42
1:C:5:LYS:O	1:C:5:LYS:HE2	2.19	0.42
1:C:7:LYS:HZ2	1:C:7:LYS:HG3	1.63	0.42
1:C:10:ILE:HD11	1:C:126:MET:HA	2.00	0.42
1:A:128:LYS:O	1:A:132:LEU:CD1	2.67	0.42
2:D:47:ASN:ND2	2:D:49:SER:CB	2.83	0.42
2:D:137:VAL:O	2:D:138:VAL:C	2.58	0.42
1:A:30:ALA:CB	5:A:149:HOH:O	2.67	0.42
1:C:34:THR:CG2	5:C:150:HOH:O	2.67	0.42
1:A:61:LYS:HD2	5:A:153:HOH:O	2.10	0.42
1:C:10:ILE:CD1	1:C:126:MET:HA	2.49	0.42
1:C:58:LYS:HG3	5:C:166:HOH:O	2.19	0.42
1:C:93:ARG:CG	1:C:93:ARG:NH1	2.82	0.42
1:C:103:SER:HB2	1:C:130:LEU:HD13	2.02	0.42
2:B:118:PHE:N	2:B:118:PHE:CD1	2.89	0.41
1:C:48:ASP:OD1	1:C:50:SER:OG	2.33	0.41
1:A:121:GLU:HG2	1:A:121:GLU:H	1.64	0.41
1:C:30:ALA:HB3	5:C:145:HOH:O	2.20	0.41
1:C:22:GLU:OE1	1:C:61:LYS:CB	2.68	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:D:35:TYR:HA	2:D:36:PRO:HD2	1.93	0.41
2:D:78:MET:HA	2:D:81:ILE:HG13	2.03	0.41
2:B:132:GLN:NE2	5:B:151:HOH:O	2.54	0.41
2:B:57:ASN:HA	2:B:58:PRO:HD3	1.95	0.41
1:C:57:LYS:HB2	1:C:58:LYS:HE3	2.01	0.41
2:D:108:ASP:O	2:D:112:VAL:HG13	2.21	0.41
1:C:66:SER:OG	1:C:84:LEU:HD11	2.20	0.41
1:A:61:LYS:HG2	5:A:156:HOH:O	2.20	0.41
2:D:41:HIS:CD2	3:D:147:HEM:HBC1	2.56	0.41
1:A:111:ALA:CB	1:A:118:PHE:CD1	3.04	0.40
1:A:50:SER:HA	1:A:51:PRO:HD3	1.91	0.40
2:D:19:SER:H	2:D:117:LYS:CE	2.31	0.40
2:D:85:TYR:CE2	2:D:138:VAL:HG13	2.57	0.40
1:C:82:LEU:HD21	1:C:140:LYS:CE	2.51	0.40
1:C:76:ASP:O	1:C:79:ASN:O	2.40	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	141/142 (99%)	128 (91%)	9 (6%)	4 (3%)	8	6
1	C	142/142 (100%)	134 (94%)	7 (5%)	1 (1%)	30	43
2	B	144/146 (99%)	136 (94%)	6 (4%)	2 (1%)	16	22
2	D	144/146 (99%)	135 (94%)	5 (4%)	4 (3%)	8	6
All	All	571/576 (99%)	533 (93%)	27 (5%)	11 (2%)	12	14

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	18	SER
1	A	48	ASP
1	A	116	ASN

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Mol	Chain	Res	Type
2	B	143	SER
2	D	122	ALA
2	D	22	GLU
2	D	48	LEU
1	A	111	ALA
2	B	93	SER
1	C	116	ASN
2	D	4	SER

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	126/125 (101%)	96 (76%)	30 (24%)	1	1
1	C	126/125 (101%)	104 (82%)	22 (18%)	3	3
2	B	118/118 (100%)	103 (87%)	15 (13%)	6	7
2	D	118/118 (100%)	109 (92%)	9 (8%)	19	28
All	All	488/486 (100%)	412 (84%)	76 (16%)	4	4

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	5	LYS
1	A	8	ASP
1	A	10	ILE
1	A	16	LYS
1	A	18	SER
1	A	22	GLU
1	A	27	ASP
1	A	34	THR
1	A	38	GLN
1	A	47	LYS
1	A	58[A]	LYS
1	A	58[B]	LYS
1	A	61	LYS

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Mol	Chain	Res	Type
1	A	69	GLU
1	A	72	SER
1	A	75	ASP
1	A	76	ASP
1	A	78	THR
1	A	79	ASN
1	A	82	LEU
1	A	83	THR
1	A	100	LYS
1	A	106	LEU
1	A	121	GLU
1	A	130	LEU
1	A	132	LEU
1	A	133	LEU
1	A	134	SER
1	A	138	SER
2	B	15	TRP
2	B	17	LYS
2	B	26	GLN
2	B	38	THR
2	B	40	ARG
2	B	47	ASN
2	B	58	PRO
2	B	72	GLU
2	B	82	LYS
2	B	93	SER
2	B	106	LEU
2	B	124	THR
2	B	126	ASP
2	B	143	SER
2	B	144	ARG
1	C	4	ASP
1	C	5	LYS
1	C	16	LYS
1	C	17	ILE
1	C	22	GLU
1	C	23	ASP
1	C	27	ASP
1	C	38	GLN
1	C	44	SER
1	C	50	SER
1	C	58	LYS

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Mol	Chain	Res	Type
1	C	61	LYS
1	C	69	GLU
1	C	81	LEU
1	C	83	THR
1	C	90	PHE
1	C	91	GLN
1	C	93	ARG
1	C	106	LEU
1	C	124	VAL
1	C	132	LEU
1	C	133	LEU
2	D	5	THR
2	D	9	SER
2	D	14	LEU
2	D	21	ASP
2	D	40	ARG
2	D	48	LEU
2	D	86	SER
2	D	94	GLU
2	D	145	TYR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	45	HIS
1	A	79	ASN
1	A	91	GLN
1	A	105	ASN
2	B	132	GLN
1	C	38	GLN
1	C	45	HIS
2	D	47	ASN
2	D	132	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 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	143	1,4	49,50,50	2.45	22 (44%)	46,82,82	2.83	17 (36%)
4	OXY	A	144	3	1,1,1	0.39	0	0,0,0	0.00	-
3	HEM	B	147	2,4	49,50,50	2.53	17 (34%)	46,82,82	2.67	18 (39%)
4	OXY	B	148	3	1,1,1	0.50	0	0,0,0	0.00	-
3	HEM	C	143	1,4	49,50,50	6.16	26 (53%)	46,82,82	4.35	22 (47%)
4	OXY	C	144	3	1,1,1	0.35	0	0,0,0	0.00	-
3	HEM	D	147	2,4	49,50,50	2.61	14 (28%)	46,82,82	2.89	15 (32%)
4	OXY	D	148	3	1,1,1	0.40	0	0,0,0	0.00	-

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	143	1,4	-	0/14/114/114	0/0/8/8
4	OXY	A	144	3	-	0/0/0/0	0/0/0/0
3	HEM	B	147	2,4	-	0/14/114/114	0/0/8/8
4	OXY	B	148	3	-	0/0/0/0	0/0/0/0
3	HEM	C	143	1,4	-	0/14/114/114	0/0/8/8
4	OXY	C	144	3	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	D	147	2,4	-	0/14/114/114	0/0/8/8
4	OXY	D	148	3	-	0/0/0/0	0/0/0/0

All (79) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	143	HEM	C2B-C1B	-31.04	1.36	1.44
3	C	143	HEM	C3D-C4D	19.52	1.49	1.44
3	D	147	HEM	C3D-C4D	-8.93	1.42	1.44
3	C	143	HEM	C4A-NA	6.97	1.50	1.36
3	D	147	HEM	C3D-C2D	6.56	1.55	1.43
3	C	143	HEM	C4A-C3A	6.31	1.48	1.40
3	B	147	HEM	FE-NA	6.31	2.19	1.92
3	C	143	HEM	CMC-C2C	6.16	1.66	1.47
3	B	147	HEM	C3D-C4D	6.02	1.46	1.44
3	C	143	HEM	FE-NB	5.66	2.18	1.97
3	B	147	HEM	C3B-CAB	5.58	1.58	1.40
3	C	143	HEM	CHA-C4D	5.36	1.43	1.35
3	A	143	HEM	C3C-C2C	-5.30	1.34	1.43
3	C	143	HEM	C3D-C2D	5.14	1.52	1.43
3	C	143	HEM	C3B-CAB	5.12	1.56	1.40
3	A	143	HEM	C4C-NC	-4.91	1.31	1.38
3	B	147	HEM	C3C-CAC	4.88	1.55	1.40
3	B	147	HEM	C2D-C1D	-4.86	1.43	1.44
3	C	143	HEM	CMB-C2B	4.79	1.62	1.47
3	C	143	HEM	C4B-NB	4.77	1.49	1.37
3	C	143	HEM	C2D-C1D	-4.73	1.43	1.44
3	A	143	HEM	C3B-CAB	4.72	1.55	1.40
3	A	143	HEM	C3C-CAC	4.70	1.55	1.40
3	C	143	HEM	FE-NA	4.69	2.12	1.92
3	D	147	HEM	C3B-C2B	-4.68	1.35	1.43
3	D	147	HEM	C3B-CAB	4.68	1.55	1.40
3	C	143	HEM	FE-NC	4.65	2.15	1.97
3	D	147	HEM	FE-NB	4.48	2.14	1.97
3	D	147	HEM	C3C-CAC	4.47	1.54	1.40
3	D	147	HEM	FE-ND	4.32	2.13	1.97
3	B	147	HEM	CHA-C4D	4.26	1.41	1.35
3	A	143	HEM	C3D-C2D	4.17	1.51	1.43
3	A	143	HEM	FE-NC	4.12	2.13	1.97
3	D	147	HEM	C3C-C2C	-4.11	1.36	1.43
3	C	143	HEM	CHD-C4C	4.03	1.43	1.36
3	C	143	HEM	C3C-CAC	3.97	1.52	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	143	HEM	CMD-C2D	3.92	1.59	1.47
3	B	147	HEM	C3D-C2D	3.91	1.50	1.43
3	B	147	HEM	C3C-C2C	-3.91	1.36	1.43
3	C	143	HEM	CMA-C3A	3.85	1.59	1.51
3	D	147	HEM	C4A-C3A	3.85	1.45	1.40
3	A	143	HEM	FE-NA	3.77	2.08	1.92
3	B	147	HEM	C3B-C2B	-3.68	1.37	1.43
3	A	143	HEM	CHA-C4D	3.67	1.41	1.35
3	A	143	HEM	CMC-C2C	3.62	1.58	1.47
3	B	147	HEM	C4A-C3A	3.59	1.44	1.40
3	C	143	HEM	CBB-CAB	3.57	1.49	1.28
3	D	147	HEM	C2D-C1D	-3.51	1.43	1.44
3	A	143	HEM	C3B-C2B	-3.49	1.37	1.43
3	C	143	HEM	C4C-NC	3.41	1.42	1.38
3	A	143	HEM	C3D-C4D	3.35	1.45	1.44
3	C	143	HEM	CHB-C1B	3.34	1.40	1.35
3	A	143	HEM	CMD-C2D	3.22	1.57	1.47
3	D	147	HEM	CMB-C2B	3.16	1.57	1.47
3	A	143	HEM	CHC-C1C	3.10	1.42	1.36
3	B	147	HEM	FE-NB	3.07	2.09	1.97
3	C	143	HEM	CHC-C4B	2.99	1.46	1.39
3	B	147	HEM	CMC-C2C	2.93	1.56	1.47
3	B	147	HEM	FE-NC	2.93	2.08	1.97
3	A	143	HEM	C1C-NC	2.84	1.42	1.38
3	C	143	HEM	CHC-C1C	2.83	1.41	1.36
3	D	147	HEM	CMC-C2C	2.69	1.55	1.47
3	A	143	HEM	C1A-C2A	2.66	1.48	1.43
3	B	147	HEM	CAA-C2A	2.62	1.56	1.52
3	A	143	HEM	CMB-C2B	2.60	1.55	1.47
3	A	143	HEM	C4A-C3A	2.57	1.43	1.40
3	C	143	HEM	O1A-CGA	2.56	1.31	1.22
3	A	143	HEM	C1A-NA	2.46	1.41	1.36
3	B	147	HEM	CMD-C2D	2.36	1.54	1.47
3	D	147	HEM	CMA-C3A	2.33	1.56	1.51
3	D	147	HEM	CMD-C2D	2.29	1.54	1.47
3	C	143	HEM	C3B-C2B	-2.27	1.39	1.43
3	A	143	HEM	CMA-C3A	2.23	1.56	1.51
3	B	147	HEM	FE-ND	2.21	2.05	1.97
3	B	147	HEM	CMB-C2B	2.17	1.54	1.47
3	A	143	HEM	O1A-CGA	2.17	1.30	1.22
3	C	143	HEM	C3C-C2C	-2.14	1.40	1.43
3	A	143	HEM	C1B-NB	2.13	1.44	1.39

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	143	HEM	CAA-C2A	2.00	1.55	1.52

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	143	HEM	C3B-C4B-NB	-12.26	105.23	114.00
3	C	143	HEM	C3A-C4A-NA	-10.75	101.29	109.41
3	D	147	HEM	C3B-C4B-NB	-10.28	106.64	114.00
3	A	143	HEM	C3B-C4B-NB	-9.04	107.53	114.00
3	C	143	HEM	CMB-C2B-C3B	8.51	146.20	126.16
3	C	143	HEM	C4A-CHB-C1B	-7.81	117.19	127.47
3	C	143	HEM	C1A-C2A-C3A	7.61	114.80	106.92
3	B	147	HEM	C4D-ND-C1D	7.48	112.82	105.16
3	D	147	HEM	CBA-CAA-C2A	-6.79	100.72	112.69
3	D	147	HEM	C4D-ND-C1D	6.79	112.11	105.16
3	C	143	HEM	CBD-CAD-C3D	-6.34	100.53	114.37
3	A	143	HEM	CMA-C3A-C4A	-6.30	118.92	128.62
3	B	147	HEM	C3A-C4A-NA	-6.06	104.83	109.41
3	C	143	HEM	CAD-C3D-C4D	6.03	135.37	124.53
3	B	147	HEM	CHC-C4B-NB	5.99	129.56	124.58
3	B	147	HEM	C3B-C4B-NB	-5.99	109.71	114.00
3	A	143	HEM	CHC-C4B-NB	5.82	129.42	124.58
3	A	143	HEM	C4A-CHB-C1B	-5.81	119.82	127.47
3	C	143	HEM	CAD-C3D-C2D	-5.79	114.35	127.25
3	A	143	HEM	CHA-C4D-ND	5.59	131.99	124.31
3	A	143	HEM	C4D-ND-C1D	5.56	110.85	105.16
3	C	143	HEM	C4D-ND-C1D	5.54	110.83	105.16
3	C	143	HEM	CHD-C1D-ND	5.33	129.02	124.58
3	D	147	HEM	C4C-NC-C1C	5.25	110.99	105.53
3	C	143	HEM	C3A-C4A-CHB	5.06	135.60	126.00
3	C	143	HEM	CMA-C3A-C4A	5.00	136.31	128.62
3	D	147	HEM	CBD-CAD-C3D	-4.98	103.51	114.37
3	C	143	HEM	CMA-C3A-C2A	-4.97	115.57	124.94
3	A	143	HEM	CMA-C3A-C2A	4.75	133.90	124.94
3	C	143	HEM	CHB-C1B-NB	4.71	130.78	124.31
3	C	143	HEM	C2A-C1A-NA	-4.58	103.37	109.73
3	C	143	HEM	CHA-C1A-NA	4.50	132.09	124.58
3	C	143	HEM	CMC-C2C-C3C	4.43	136.60	126.16
3	C	143	HEM	C4A-NA-C1A	4.35	112.50	106.76
3	B	147	HEM	CBA-CAA-C2A	-4.08	105.50	112.69
3	D	147	HEM	CHD-C4C-NC	4.08	128.27	124.73
3	B	147	HEM	C2D-C1D-ND	-3.96	108.25	112.93

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	147	HEM	C4A-C3A-C2A	3.83	109.66	107.00
3	C	143	HEM	CHC-C4B-NB	3.82	127.76	124.58
3	D	147	HEM	C1B-NB-C4B	3.73	108.97	105.16
3	B	147	HEM	CAD-C3D-C4D	3.69	131.16	124.53
3	C	143	HEM	CMB-C2B-C1B	-3.64	104.14	126.30
3	B	147	HEM	C4C-NC-C1C	3.61	109.29	105.53
3	A	143	HEM	CAD-CBD-CGD	-3.60	102.24	113.48
3	B	147	HEM	C4A-NA-C1A	3.34	111.16	106.76
3	B	147	HEM	CMA-C3A-C4A	-3.22	123.66	128.62
3	D	147	HEM	C2D-C1D-ND	-3.13	109.24	112.93
3	A	143	HEM	CBD-CAD-C3D	-3.08	107.65	114.37
3	D	147	HEM	CMB-C2B-C3B	3.08	133.41	126.16
3	D	147	HEM	CMA-C3A-C4A	-3.00	124.00	128.62
3	B	147	HEM	CHD-C1D-ND	3.00	127.08	124.58
3	D	147	HEM	CHA-C1A-NA	3.00	129.58	124.58
3	B	147	HEM	C1B-NB-C4B	2.87	108.09	105.16
3	D	147	HEM	C3A-C4A-NA	-2.84	107.27	109.41
3	D	147	HEM	CMD-C2D-C3D	2.79	131.91	125.60
3	A	143	HEM	CBA-CAA-C2A	-2.76	107.83	112.69
3	D	147	HEM	CHC-C1C-NC	2.54	126.94	124.73
3	A	143	HEM	C1B-NB-C4B	2.54	107.76	105.16
3	C	143	HEM	CBA-CAA-C2A	-2.39	108.48	112.69
3	A	143	HEM	C3A-C4A-CHB	-2.37	121.50	126.00
3	A	143	HEM	C2D-C1D-ND	-2.31	110.20	112.93
3	A	143	HEM	C2A-C1A-CHA	2.25	130.27	126.00
3	A	143	HEM	C2A-C1A-NA	-2.22	106.65	109.73
3	B	147	HEM	O2D-CGD-CBD	2.22	122.07	114.22
3	A	143	HEM	C3A-C4A-NA	2.21	111.08	109.41
3	C	143	HEM	C1D-CHD-C4C	-2.18	120.83	126.57
3	B	147	HEM	CAD-CBD-CGD	-2.13	106.83	113.48
3	B	147	HEM	CHB-C4A-NA	2.09	128.08	124.58
3	B	147	HEM	O1A-CGA-CBA	-2.08	115.86	123.03
3	D	147	HEM	CHB-C4A-NA	2.08	128.05	124.58
3	A	143	HEM	CMB-C2B-C3B	2.08	131.06	126.16
3	B	147	HEM	CAD-C3D-C2D	-2.04	122.70	127.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	142/142 (100%)	-0.54	2 (1%) 72 71	25, 37, 52, 61	1 (0%)
1	C	142/142 (100%)	-0.53	0 100 100	22, 36, 51, 58	4 (2%)
2	B	146/146 (100%)	-0.49	0 100 100	24, 40, 59, 71	1 (0%)
2	D	146/146 (100%)	-0.57	1 (0%) 84 84	24, 40, 57, 69	1 (0%)
All	All	576/576 (100%)	-0.53	3 (0%) 88 88	22, 39, 56, 71	7 (1%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	D	145	TYR	5.5
1	A	116	ASN	2.3
1	A	79	ASN	2.1

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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, 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	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	HEM	C	143	43/43	0.13	0.68	20,27,48,57	0
3	HEM	D	147	43/43	0.12	0.54	26,36,55,63	0
3	HEM	B	147	43/43	0.13	0.29	31,37,58,61	0
3	HEM	A	143	43/43	0.11	0.12	23,29,48,60	0
4	OXY	D	148	2/2	0.09	-0.81	26,26,26,30	2
4	OXY	C	144	2/2	0.07	-1.06	36,36,36,40	0
4	OXY	B	148	2/2	0.07	-1.77	48,48,48,52	0
4	OXY	A	144	2/2	0.06	-2.55	22,22,22,26	0

6.5 Other polymers

There are no such residues in this entry.