



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 07:40 AM GMT

PDB ID : 3BKN
Title : The structure of Mycobacterial bacterioferritin
Authors : Janowski, R.; Auerbach-Nevo, T.; Weiss, M.S.
Deposited on : 2007-12-07
Resolution : 2.72 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
A user guide is available at
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

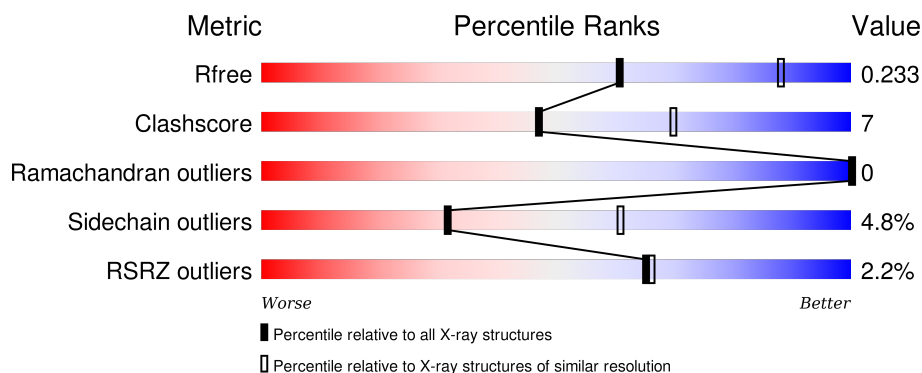
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.72 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}	91344	2439 (2.74-2.70)
Clashscore	102246	2771 (2.74-2.70)
Ramachandran outliers	100387	2726 (2.74-2.70)
Sidechain outliers	100360	2727 (2.74-2.70)
RSRZ outliers	91569	2443 (2.74-2.70)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	161	<div> <div>2%</div> <div>81%19%</div> </div>
1	B	161	<div> <div>2%</div> <div>84%12%••</div> </div>
1	C	161	<div> <div>2%</div> <div>86%13%•</div> </div>
1	D	161	<div> <div>2%</div> <div>89%10%•</div> </div>
1	E	161	<div> <div>3%</div> <div>84%14%•</div> </div>

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Mol	Chain	Length	Quality of chain
1	F	161	
1	G	161	
1	H	161	
1	I	161	
1	J	161	
1	K	161	
1	L	161	

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	MG	A	401	-	-	-	X
5	EPE	J	501	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 16205 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 Bacterioferritin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	161	Total	C	N	O	S	0	1	0
			1303	817	221	259	6			
1	B	161	Total	C	N	O	S	0	1	0
			1303	817	221	259	6			
1	C	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	D	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	E	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	F	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	G	161	Total	C	N	O	S	0	1	0
			1303	817	221	259	6			
1	H	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	I	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	J	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	K	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			
1	L	161	Total	C	N	O	S	0	0	0
			1297	813	221	257	6			

- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	G	2	Total	Zn	0	0
			2	2		
2	J	2	Total	Zn	0	0
			2	2		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	D	2	Total 2	Zn 2	0	0
2	K	2	Total 2	Zn 2	0	0
2	E	2	Total 2	Zn 2	0	0
2	H	2	Total 2	Zn 2	0	0
2	B	2	Total 2	Zn 2	0	0
2	I	2	Total 2	Zn 2	0	0
2	C	2	Total 2	Zn 2	0	0
2	A	2	Total 2	Zn 2	0	0
2	L	2	Total 2	Zn 2	0	0
2	F	2	Total 2	Zn 2	0	0

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

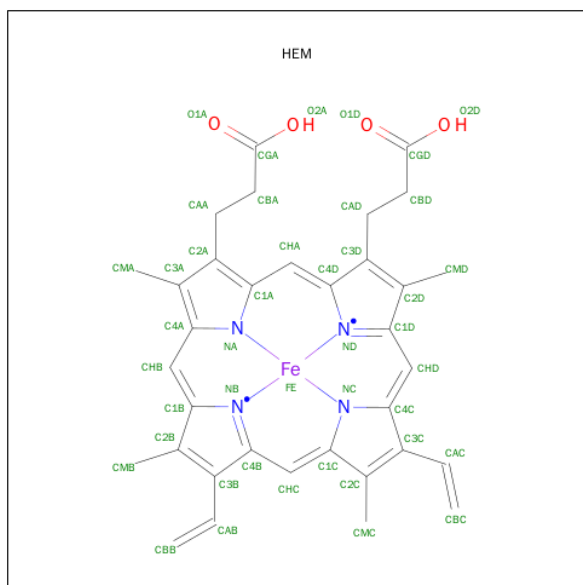
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	J	1	Total 1	Mg 1	0	0
3	K	1	Total 1	Mg 1	0	0
3	E	1	Total 1	Mg 1	0	0
3	H	1	Total 1	Mg 1	0	0
3	B	2	Total 2	Mg 2	0	0
3	I	1	Total 1	Mg 1	0	0
3	C	2	Total 2	Mg 2	0	0
3	A	1	Total 1	Mg 1	0	0
3	L	1	Total 1	Mg 1	0	0

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	F	1	Total	Mg	0	0
			1	1		

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	B	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
4	C	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
4	E	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
4	H	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
4	I	1	Total	C	Fe	N	O	0	1
			86	68	2	8	8		
4	K	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
4	L	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	J	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
5	D	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	5	Total	O	0	0
			5	5		
6	B	3	Total	O	0	0
			3	3		
6	C	1	Total	O	0	0
			1	1		
6	D	2	Total	O	0	0
			2	2		
6	F	3	Total	O	0	0
			3	3		
6	G	9	Total	O	0	0
			9	9		
6	H	3	Total	O	0	0
			3	3		
6	I	6	Total	O	0	0
			6	6		
6	J	3	Total	O	0	0
			3	3		
6	K	3	Total	O	0	0
			3	3		

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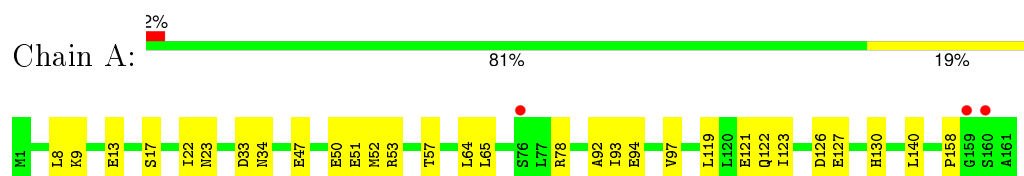
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	L	3	Total	O	0	0
			3	3		

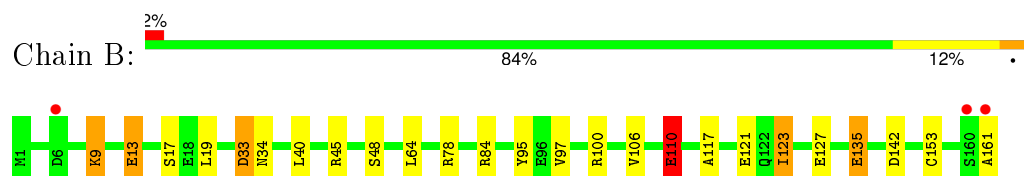
3 Residue-property plots

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 ($\text{RSRZ} > 2$). 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.

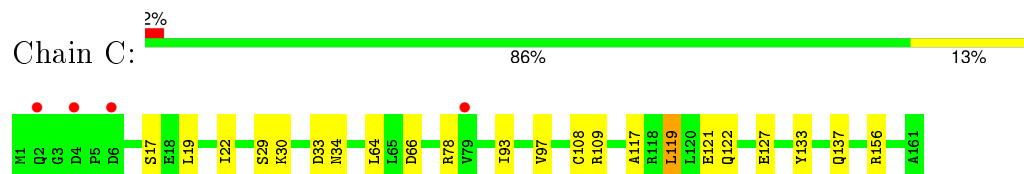
- Molecule 1: Bacterioferritin



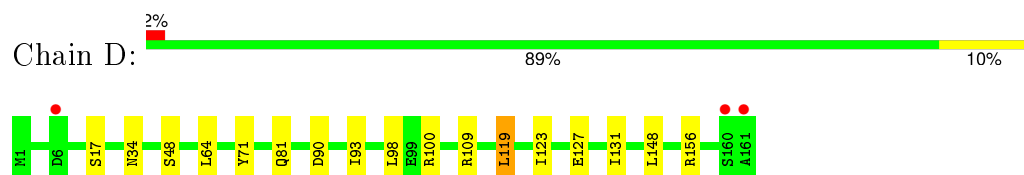
- Molecule 1: Bacterioferritin



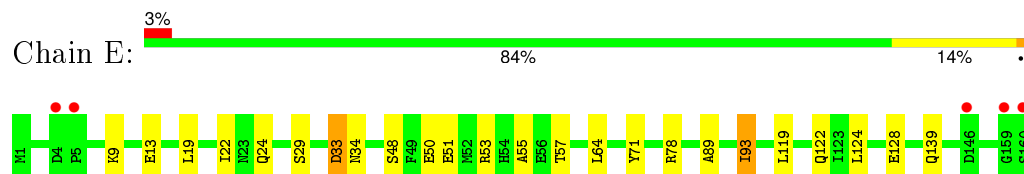
- Molecule 1: Bacterioferritin



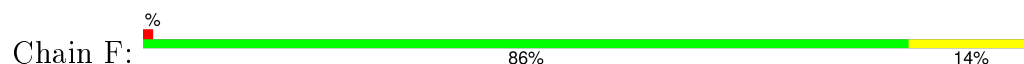
- Molecule 1: Bacterioferritin



- Molecule 1: Bacterioferritin

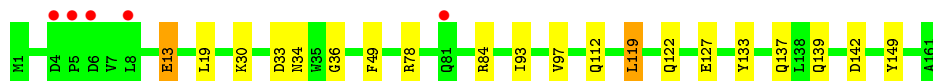
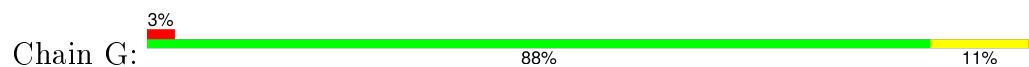


- Molecule 1: Bacterioferritin

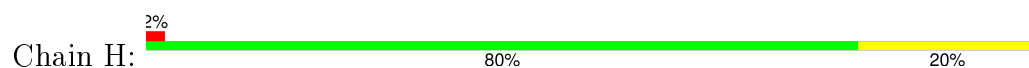




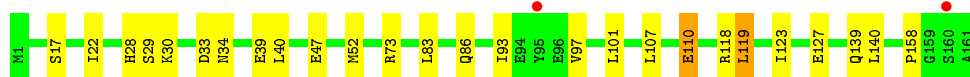
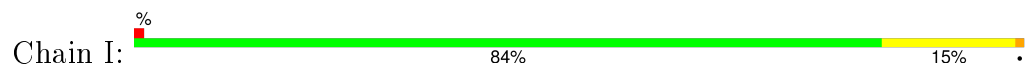
- Molecule 1: Bacterioferritin



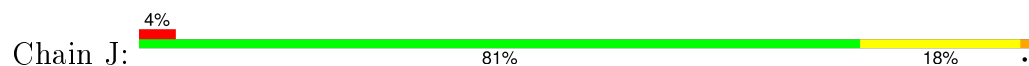
- Molecule 1: Bacterioferritin



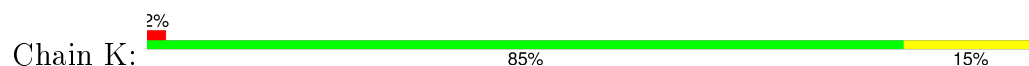
- Molecule 1: Bacterioferritin



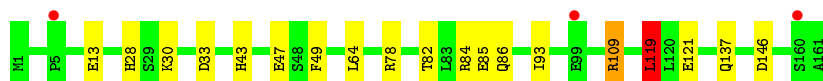
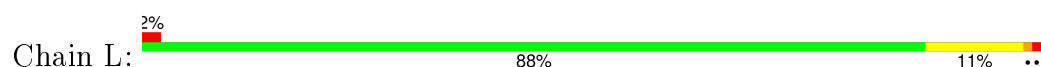
- Molecule 1: Bacterioferritin



- Molecule 1: Bacterioferritin



- Molecule 1: Bacterioferritin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	181.52Å 151.52Å 116.72Å 90.00° 128.08° 90.00°	Depositor
Resolution (Å)	29.64 – 2.72 29.48 – 2.72	Depositor EDS
% Data completeness (in resolution range)	98.8 (29.64-2.72) 98.8 (29.48-2.72)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 2.72Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.179 , 0.228 0.186 , 0.233	Depositor DCC
R_{free} test set	2649 reflections (4.20%)	DCC
Wilson B-factor (Å ²)	33.9	Xtriage
Anisotropy	0.257	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.8	EDS
Estimated twinning fraction	0.020 for -h-2*k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	3 of 65778 reflections (0.005%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16205	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.84	0/1326	0.85	1/1793 (0.1%)
1	B	0.82	1/1326 (0.1%)	0.79	0/1793
1	C	0.81	1/1317 (0.1%)	0.84	1/1781 (0.1%)
1	D	0.79	0/1317	0.84	0/1781
1	E	0.82	0/1317	0.85	0/1781
1	F	0.74	0/1317	0.74	0/1781
1	G	0.86	1/1326 (0.1%)	0.83	0/1793
1	H	0.78	0/1317	0.81	0/1781
1	I	0.86	2/1317 (0.2%)	0.85	1/1781 (0.1%)
1	J	0.88	2/1317 (0.2%)	0.89	1/1781 (0.1%)
1	K	0.80	1/1317 (0.1%)	0.85	0/1781
1	L	0.74	0/1317	0.80	1/1781 (0.1%)
All	All	0.81	8/15831 (0.1%)	0.83	5/21408 (0.0%)

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	J	110	GLU	CG-CD	5.92	1.60	1.51
1	I	110	GLU	CG-CD	5.83	1.60	1.51
1	J	75	PHE	CE2-CZ	5.55	1.47	1.37
1	K	42	GLU	CG-CD	5.44	1.60	1.51
1	I	47	GLU	CG-CD	5.12	1.59	1.51
1	G	13	GLU	CG-CD	5.11	1.59	1.51
1	C	108	CYS	CB-SG	-5.10	1.73	1.81
1	B	110	GLU	CG-CD	5.08	1.59	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	126	ASP	CB-CG-OD1	5.86	123.57	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	66	ASP	CB-CG-OD1	5.79	123.51	118.30
1	I	119	LEU	CA-CB-CG	5.68	128.37	115.30
1	L	119	LEU	CA-CB-CG	5.10	127.03	115.30
1	J	66	ASP	CB-CG-OD1	5.02	122.81	118.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1303	0	1283	23	0
1	B	1303	0	1283	16	0
1	C	1297	0	1277	12	0
1	D	1297	0	1277	9	0
1	E	1297	0	1277	19	0
1	F	1297	0	1277	19	0
1	G	1303	0	1283	12	0
1	H	1297	0	1277	24	0
1	I	1297	0	1277	12	0
1	J	1297	0	1277	23	0
1	K	1297	0	1277	16	0
1	L	1297	0	1277	12	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0
2	K	2	0	0	0	0
2	L	2	0	0	0	0
3	A	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
3	H	1	0	0	0	0
3	I	1	0	0	0	0
3	J	1	0	0	0	0
3	K	1	0	0	0	0
3	L	1	0	0	0	0
4	B	86	0	60	20	0
4	C	86	0	60	13	0
4	E	86	0	60	12	0
4	H	86	0	60	9	0
4	I	86	0	60	4	0
4	K	43	0	30	15	0
4	L	43	0	30	3	0
5	D	15	0	17	0	0
5	J	15	0	18	0	0
6	A	5	0	0	0	0
6	B	3	0	0	0	0
6	C	1	0	0	0	0
6	D	2	0	0	0	0
6	F	3	0	0	0	0
6	G	9	0	0	0	0
6	H	3	0	0	0	0
6	I	6	0	0	0	0
6	J	3	0	0	0	0
6	K	3	0	0	1	0
6	L	3	0	0	0	0
All	All	16205	0	15737	239	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (239) 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:C:119:LEU:HD12	1:C:119:LEU:O	1.38	1.20
1:F:119:LEU:O	1:F:119:LEU:HD12	1.42	1.17
1:D:119:LEU:HD12	1:D:119:LEU:O	1.44	1.13
1:F:119:LEU:C	1:F:119:LEU:HD12	1.75	1.04

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:118:ARG:NH2	1:J:122:GLN:HE21	1.68	0.92
1:B:33:ASP:HB3	1:B:161:ALA:HB2	1.60	0.84
4:B:301[A]:HEM:HHD	4:B:301[A]:HEM:HBC2	1.60	0.83
1:A:23:ASN:OD1	4:B:301[A]:HEM:HBB1	1.77	0.83
1:G:119:LEU:O	1:G:119:LEU:HD12	1.79	0.83
4:C:301[B]:HEM:CMB	4:C:301[B]:HEM:HBB2	2.08	0.82
1:A:23:ASN:OD1	4:B:301[A]:HEM:CBB	2.29	0.81
1:H:78:ARG:HH21	1:H:93:ILE:CG2	1.94	0.80
4:B:301[B]:HEM:CBD	4:B:301[B]:HEM:HHA	2.14	0.78
1:E:24:GLN:NE2	1:E:93:ILE:HD11	1.99	0.78
1:E:55:ALA:HB3	4:E:301[B]:HEM:HBC2	1.67	0.77
1:F:119:LEU:CD1	1:F:119:LEU:C	2.51	0.75
1:A:64:LEU:O	1:A:64:LEU:HD23	1.87	0.74
4:K:301:HEM:CMB	4:K:301:HEM:HBB2	2.17	0.74
1:L:28:HIS:CD2	1:L:86:GLN:HG2	2.24	0.73
1:F:119:LEU:O	1:F:119:LEU:CD1	2.31	0.72
1:C:119:LEU:O	1:C:119:LEU:CD1	2.30	0.71
1:H:109:ARG:HH11	1:H:109:ARG:HG2	1.57	0.70
4:C:301[A]:HEM:CBD	4:C:301[A]:HEM:HHA	2.22	0.70
1:D:17:SER:OG	1:D:100:ARG:NH2	2.24	0.70
4:K:301:HEM:HBA1	4:K:301:HEM:HHA	1.75	0.69
4:C:301[B]:HEM:HMB1	4:C:301[B]:HEM:HBB2	1.75	0.69
1:G:84:ARG:NH2	1:G:142:ASP:OD1	2.27	0.68
4:K:301:HEM:HMB1	4:K:301:HEM:HBB2	1.75	0.68
1:F:118:ARG:HH21	1:J:122:GLN:HE21	1.42	0.67
1:H:78:ARG:HH21	1:H:93:ILE:HG22	1.59	0.67
1:D:123:ILE:O	1:D:127:GLU:HG2	1.95	0.67
1:J:45:ARG:O	1:J:48:SER:HB3	1.95	0.67
1:H:97:VAL:HG11	1:H:127:GLU:HG3	1.77	0.65
1:D:119:LEU:HD12	1:D:119:LEU:C	2.17	0.65
1:B:97:VAL:HG11	1:B:127:GLU:HG3	1.77	0.65
4:C:301[A]:HEM:HBA2	4:C:301[A]:HEM:CMA	2.27	0.65
1:H:78:ARG:NH2	1:H:93:ILE:CG2	2.60	0.64
1:A:9:LYS:O	1:A:13:GLU:HG2	1.96	0.64
1:K:19:LEU:HD22	4:K:301:HEM:HBB1	1.79	0.64
1:A:78:ARG:NH1	1:A:92:ALA:HB1	2.13	0.64
4:B:301[B]:HEM:HBD1	4:B:301[B]:HEM:HHA	1.80	0.64
4:E:301[A]:HEM:HBB2	4:E:301[A]:HEM:CMB	2.28	0.63
4:E:301[B]:HEM:CMB	4:E:301[B]:HEM:HBB2	2.28	0.63
1:A:121:GLU:OE2	1:D:109:ARG:NH2	2.31	0.63
1:L:84:ARG:NH1	1:L:146:ASP:OD1	2.32	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:K:301:HEM:CHA	4:K:301:HEM:HBA1	2.28	0.63
4:C:301[B]:HEM:CMC	4:C:301[B]:HEM:HBC2	2.30	0.62
4:C:301[B]:HEM:HBC2	4:C:301[B]:HEM:HMC1	1.82	0.62
4:B:301[B]:HEM:HBD1	4:B:301[B]:HEM:CHA	2.29	0.62
1:E:55:ALA:HB3	4:E:301[B]:HEM:CBC	2.29	0.62
1:J:57:THR:CG2	1:J:119:LEU:CD1	2.78	0.61
1:E:57:THR:HG22	1:E:119:LEU:HD11	1.81	0.61
1:H:78:ARG:NH2	1:H:93:ILE:HG22	2.14	0.61
4:I:301[A]:HEM:HMC1	4:I:301[A]:HEM:HBC2	1.83	0.61
1:A:64:LEU:HD23	1:A:64:LEU:C	2.21	0.60
1:C:119:LEU:HD12	1:C:119:LEU:C	2.14	0.60
4:K:301:HEM:CBD	4:K:301:HEM:HHA	2.32	0.59
1:J:97:VAL:HG11	1:J:127:GLU:HG3	1.84	0.59
1:D:119:LEU:O	1:D:119:LEU:CD1	2.36	0.59
1:J:57:THR:CG2	1:J:119:LEU:HD11	2.33	0.59
1:J:57:THR:HG21	1:J:119:LEU:CD1	2.32	0.59
1:G:119:LEU:HD12	1:G:119:LEU:C	2.22	0.58
1:C:117:ALA:O	1:C:121:GLU:HG3	2.04	0.58
1:E:50:GLU:HG3	1:E:53:ARG:NH2	2.19	0.58
1:G:97:VAL:HG11	1:G:127:GLU:HG3	1.86	0.57
1:A:57:THR:HG21	1:A:119:LEU:HD11	1.87	0.57
1:B:33:ASP:CB	1:B:161:ALA:HB2	2.33	0.57
1:A:57:THR:CG2	1:A:119:LEU:HD11	2.34	0.57
1:C:64:LEU:HD23	1:C:64:LEU:O	2.05	0.56
1:J:57:THR:HG22	1:J:119:LEU:HD11	1.88	0.56
4:I:301[A]:HEM:CMC	4:I:301[A]:HEM:HBC2	2.36	0.56
1:I:97:VAL:HG11	1:I:127:GLU:HG3	1.88	0.56
1:H:27:LEU:O	1:H:31:MET:HG3	2.05	0.56
4:H:301[A]:HEM:HHD	4:H:301[A]:HEM:HBC2	1.87	0.55
1:B:19:LEU:HD22	4:B:301[A]:HEM:HAB	1.87	0.55
1:H:109:ARG:NH1	1:H:109:ARG:HG2	2.22	0.55
1:H:119:LEU:HD12	1:H:119:LEU:O	2.07	0.55
4:B:301[A]:HEM:HHA	4:B:301[A]:HEM:CBA	2.36	0.55
4:B:301[B]:HEM:HBD2	4:B:301[B]:HEM:HHA	1.89	0.55
4:B:301[B]:HEM:CBD	4:B:301[B]:HEM:CHA	2.80	0.55
4:K:301:HEM:HBC2	4:K:301:HEM:CMC	2.37	0.55
1:H:84:ARG:NH2	1:H:142:ASP:OD1	2.39	0.55
1:G:49:PHE:CE1	4:H:301[B]:HEM:HAD2	2.41	0.54
4:L:301:HEM:HHC	4:L:301:HEM:CBB	2.38	0.54
1:B:64:LEU:HD23	1:B:64:LEU:O	2.07	0.54
1:F:97:VAL:HG11	1:F:127:GLU:HG3	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:C:301[B]:HEM:CBB	4:C:301[B]:HEM:HMB1	2.37	0.54
4:K:301:HEM:CBD	4:K:301:HEM:CHA	2.86	0.54
1:K:64:LEU:HD23	1:K:64:LEU:O	2.08	0.54
1:I:39:GLU:OE1	1:I:39:GLU:HA	2.08	0.54
1:K:45:ARG:O	1:K:48:SER:HB3	2.09	0.53
1:B:19:LEU:CD2	4:B:301[A]:HEM:HAB	2.37	0.53
1:F:119:LEU:CD1	1:F:123:ILE:HD12	2.38	0.53
1:J:57:THR:CG2	1:J:119:LEU:HD13	2.39	0.53
1:B:40:LEU:HB2	1:B:153:CYS:HB3	1.90	0.53
1:F:118:ARG:NH2	1:J:122:GLN:NE2	2.50	0.53
1:E:55:ALA:CB	4:E:301[B]:HEM:HBC2	2.37	0.53
1:H:57:THR:HG22	1:H:119:LEU:HD11	1.89	0.53
1:D:90:ASP:O	1:D:93:ILE:HG22	2.08	0.53
1:K:123:ILE:O	1:K:127:GLU:HG2	2.07	0.53
4:K:301:HEM:CHA	4:K:301:HEM:HBD1	2.39	0.53
1:E:57:THR:HG22	1:E:119:LEU:CD1	2.38	0.52
1:F:119:LEU:HD11	1:F:123:ILE:HD12	1.91	0.52
4:H:301[B]:HEM:HHD	4:H:301[B]:HEM:HBC2	1.91	0.52
4:B:301[A]:HEM:CHD	4:B:301[A]:HEM:HBC2	2.35	0.52
1:E:51:GLU:OE1	1:E:51:GLU:HA	2.09	0.52
1:L:28:HIS:NE2	1:L:86:GLN:HG2	2.25	0.52
4:H:301[A]:HEM:HHD	4:H:301[A]:HEM:CBC	2.39	0.52
1:K:102:LYS:NZ	6:K:402:HOH:O	2.42	0.52
1:K:18:GLU:O	1:K:22:ILE:HG13	2.10	0.51
4:K:301:HEM:HBC2	4:K:301:HEM:HMC1	1.91	0.51
1:J:19:LEU:HA	1:J:22:ILE:HD12	1.93	0.51
1:J:57:THR:HG21	1:J:119:LEU:HD13	1.93	0.51
1:A:51:GLU:OE1	1:A:51:GLU:HA	2.11	0.51
1:G:36:GLY:HA3	1:I:140:LEU:HD11	1.93	0.51
4:C:301[A]:HEM:HHA	4:C:301[A]:HEM:HBD1	1.93	0.50
1:B:45:ARG:O	1:B:48:SER:HB3	2.11	0.50
1:D:98:LEU:HD13	1:D:131:ILE:CD1	2.41	0.50
1:H:28:HIS:CD2	1:H:86:GLN:HG2	2.46	0.50
1:F:118:ARG:HH21	1:J:122:GLN:NE2	2.08	0.50
1:E:55:ALA:CB	4:E:301[B]:HEM:CBC	2.90	0.49
4:H:301[B]:HEM:HBB2	4:H:301[B]:HEM:CMB	2.43	0.49
1:L:49:PHE:CZ	4:L:301:HEM:HAD2	2.47	0.49
4:C:301[A]:HEM:CHA	4:C:301[A]:HEM:HBD1	2.41	0.49
1:L:43:HIS:CE1	1:L:47:GLU:OE2	2.65	0.49
4:B:301[A]:HEM:CHA	4:B:301[A]:HEM:HBA1	2.41	0.49
4:B:301[B]:HEM:HBB2	4:B:301[B]:HEM:HHC	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:119:LEU:HD12	1:L:119:LEU:C	2.32	0.48
1:B:9:LYS:NZ	1:B:13[A]:GLU:OE1	2.46	0.48
1:L:82:THR:OG1	1:L:85:GLU:HB2	2.13	0.48
1:H:84:ARG:NH1	1:H:146:ASP:OD1	2.47	0.48
1:L:64:LEU:HD23	1:L:64:LEU:O	2.14	0.48
4:C:301[B]:HEM:HBD2	4:C:301[B]:HEM:CMD	2.42	0.48
1:C:19:LEU:HA	1:C:22:ILE:HD12	1.94	0.48
4:B:301[B]:HEM:CBB	4:B:301[B]:HEM:HHC	2.44	0.48
4:I:301[B]:HEM:HBC2	4:I:301[B]:HEM:CMC	2.44	0.48
1:J:57:THR:HG22	1:J:119:LEU:CD1	2.43	0.48
1:H:57:THR:CG2	1:H:119:LEU:HD11	2.44	0.48
1:E:78:ARG:HB2	1:E:89:ALA:HB1	1.96	0.48
1:L:109:ARG:HH11	1:L:109:ARG:HG2	1.78	0.48
4:C:301[A]:HEM:HBB2	4:C:301[A]:HEM:CMB	2.43	0.48
1:I:30:LYS:HE2	1:J:56:GLU:HG3	1.96	0.48
4:C:301[A]:HEM:HHA	4:C:301[A]:HEM:HBD2	1.94	0.47
1:H:78:ARG:HH22	1:H:96:GLU:CD	2.18	0.47
1:B:106:VAL:O	1:B:110:GLU:HG3	2.14	0.47
1:E:9:LYS:O	1:E:13:GLU:HG2	2.14	0.47
1:H:4:ASP:CG	1:H:5:PRO:HD2	2.35	0.47
1:F:149:TYR:C	1:F:149:TYR:CD2	2.87	0.47
1:K:40:LEU:HB2	1:K:153:CYS:HB3	1.96	0.47
1:I:22:ILE:HD11	1:I:52:MET:HA	1.95	0.47
1:B:123:ILE:O	1:B:127:GLU:HG2	2.15	0.47
1:K:47:GLU:OE1	1:K:130:HIS:CD2	2.67	0.47
1:G:49:PHE:CE1	4:H:301[A]:HEM:HAA1	2.50	0.47
1:J:50:GLU:CG	1:J:53:ARG:HH21	2.28	0.47
4:K:301:HEM:CBB	4:K:301:HEM:HMB1	2.44	0.46
1:K:19:LEU:CD2	4:K:301:HEM:HBB1	2.43	0.46
1:A:22:ILE:HG22	4:B:301[A]:HEM:HBB2	1.98	0.46
1:K:19:LEU:CD2	4:K:301:HEM:CBB	2.93	0.46
1:A:78:ARG:NH1	1:A:92:ALA:CB	2.79	0.46
1:H:18:GLU:O	1:H:22:ILE:HG13	2.16	0.46
1:E:50:GLU:HG3	1:E:53:ARG:HH21	1.80	0.46
4:E:301[A]:HEM:CBB	4:E:301[A]:HEM:CMB	2.94	0.46
1:E:64:LEU:HD23	1:E:64:LEU:O	2.16	0.46
1:C:97:VAL:HG11	1:C:127:GLU:HG3	1.97	0.46
1:B:84:ARG:NH2	1:B:142:ASP:OD1	2.48	0.45
1:F:45:ARG:O	1:F:48:SER:HB3	2.16	0.45
1:A:94:GLU:OE1	1:A:127:GLU:HB3	2.17	0.45
1:F:14:GLN:NE2	1:F:18:GLU:OE2	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:VAL:HG11	1:A:127:GLU:HG3	1.99	0.45
4:K:301:HEM:HHA	4:K:301:HEM:CBA	2.45	0.45
1:G:19:LEU:HD22	4:H:301[B]:HEM:HBB1	1.98	0.45
4:C:301[A]:HEM:CBB	1:D:71:TYR:HE2	2.31	0.44
4:B:301[A]:HEM:HHA	4:B:301[A]:HEM:HBA1	2.00	0.44
4:B:301[B]:HEM:HBB2	4:B:301[B]:HEM:CHC	2.47	0.44
1:H:57:THR:CG2	1:H:119:LEU:CD1	2.95	0.44
1:K:58:ILE:HD11	1:K:123:ILE:HD13	1.97	0.44
1:C:133:TYR:O	1:C:137:GLN:HG2	2.17	0.44
1:A:8:LEU:HD21	1:A:65:LEU:HB2	1.98	0.44
1:A:52:MET:HB3	4:B:301[B]:HEM:CHB	2.47	0.44
1:F:78:ARG:HH21	1:F:93:ILE:HG22	1.83	0.44
1:J:158:PRO:HB2	1:L:137:GLN:OE1	2.18	0.44
1:E:71:TYR:HE2	4:E:301[A]:HEM:CBB	2.31	0.43
1:B:17:SER:HB2	1:B:100:ARG:NH2	2.33	0.43
1:C:93:ILE:HD13	1:C:93:ILE:HG21	1.74	0.43
1:K:19:LEU:HD22	4:K:301:HEM:CBB	2.46	0.43
1:H:109:ARG:NH1	1:H:109:ARG:CG	2.81	0.43
1:E:57:THR:CG2	1:E:119:LEU:CD1	2.96	0.43
1:G:30:LYS:HD2	1:G:30:LYS:HA	1.65	0.43
1:C:30:LYS:HD2	1:C:30:LYS:HA	1.81	0.43
1:H:124:LEU:O	1:H:128:GLU:HG3	2.18	0.43
1:E:124:LEU:O	1:E:128:GLU:HG3	2.17	0.43
4:E:301[B]:HEM:CMB	4:E:301[B]:HEM:CBB	2.96	0.43
1:F:19:LEU:HA	1:F:22:ILE:HD12	2.01	0.43
4:H:301[A]:HEM:CHD	4:H:301[A]:HEM:HBC2	2.48	0.43
1:J:133:TYR:O	1:J:137:GLN:HG2	2.18	0.43
1:H:45:ARG:O	1:H:48:SER:HB3	2.18	0.43
1:H:78:ARG:NH2	1:H:93:ILE:HG23	2.32	0.43
4:I:301[B]:HEM:HMC1	4:I:301[B]:HEM:HBC2	2.00	0.43
1:L:30:LYS:HA	1:L:30:LYS:HD2	1.81	0.43
1:J:36:GLY:O	1:J:154:VAL:HG12	2.19	0.43
1:A:47[A]:GLU:OE1	1:A:130:HIS:NE2	2.51	0.43
1:C:109:ARG:HH21	1:C:121:GLU:CD	2.23	0.43
1:H:30:LYS:HD2	1:H:30:LYS:HA	1.73	0.43
1:E:33:ASP:OD1	1:E:161:ALA:HA	2.18	0.43
1:J:24:GLN:OE1	1:J:93:ILE:HD11	2.19	0.42
1:K:37:PHE:CE1	1:K:150:ALA:HB1	2.55	0.42
1:I:101:LEU:HD23	1:I:101:LEU:HA	1.73	0.42
1:K:97:VAL:HG11	1:K:127:GLU:HG3	2.02	0.42
1:A:140:LEU:HD23	1:A:140:LEU:HA	1.89	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:71:TYR:HE2	4:E:301[A]:HEM:HBB2	1.85	0.42
4:E:301[B]:HEM:HMB1	4:E:301[B]:HEM:HBB2	1.98	0.42
1:J:123:ILE:O	1:J:127:GLU:HG2	2.19	0.42
1:G:19:LEU:CD2	4:H:301[B]:HEM:CBB	2.98	0.42
1:J:64:LEU:O	1:J:64:LEU:HD23	2.19	0.42
1:J:148:LEU:HA	1:J:148:LEU:HD12	1.72	0.42
1:B:95:TYR:OH	1:B:135:GLU:OE2	2.20	0.42
1:E:19:LEU:HA	1:E:22:ILE:HD12	2.02	0.42
1:A:50:GLU:HG3	1:A:53:ARG:NH2	2.35	0.42
1:B:64:LEU:C	1:B:64:LEU:HD23	2.40	0.41
4:B:301[A]:HEM:CHA	4:B:301[A]:HEM:CBA	2.96	0.41
1:A:158:PRO:HG2	1:G:149:TYR:CE1	2.56	0.41
4:E:301[A]:HEM:HBC2	1:F:55:ALA:HB3	2.02	0.41
1:A:123:ILE:O	1:A:127:GLU:HG2	2.21	0.41
1:I:93:ILE:HG21	1:I:93:ILE:HD13	1.74	0.41
1:F:94:GLU:HA	1:F:94:GLU:OE1	2.20	0.41
1:A:57:THR:HG21	1:A:119:LEU:CD1	2.50	0.41
1:K:93:ILE:HG21	1:K:93:ILE:HD13	1.76	0.41
1:I:28:HIS:CD2	1:I:86:GLN:HG2	2.55	0.41
1:H:140:LEU:HD23	1:H:140:LEU:HA	1.95	0.41
1:F:39:GLU:OE1	1:F:39:GLU:HA	2.21	0.41
1:B:117:ALA:O	1:B:121:GLU:HG3	2.20	0.41
4:L:301:HEM:CHC	4:L:301:HEM:CBB	2.98	0.41
1:C:137:GLN:OE1	1:I:158:PRO:HB2	2.21	0.41
1:I:40:LEU:CD2	1:I:83:LEU:HD22	2.51	0.41
1:A:93:ILE:HG21	1:A:93:ILE:HD13	1.84	0.41
1:I:123:ILE:O	1:I:127:GLU:HG2	2.21	0.40
1:I:118:ARG:HD2	1:L:121:GLU:OE1	2.22	0.40
1:G:133:TYR:O	1:G:137:GLN:HG2	2.20	0.40
1:K:30:LYS:HD2	1:K:30:LYS:HA	1.84	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	160/161 (99%)	159 (99%)	1 (1%)	0	100	100
1	B	160/161 (99%)	158 (99%)	2 (1%)	0	100	100
1	C	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
1	D	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
1	E	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
1	F	159/161 (99%)	157 (99%)	2 (1%)	0	100	100
1	G	160/161 (99%)	158 (99%)	2 (1%)	0	100	100
1	H	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
1	I	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
1	J	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
1	K	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
1	L	159/161 (99%)	158 (99%)	1 (1%)	0	100	100
All	All	1911/1932 (99%)	1896 (99%)	15 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	142/141 (101%)	138 (97%)	4 (3%)	51	80
1	B	142/141 (101%)	133 (94%)	9 (6%)	22	47
1	C	141/141 (100%)	133 (94%)	8 (6%)	25	52
1	D	141/141 (100%)	134 (95%)	7 (5%)	30	59
1	E	141/141 (100%)	134 (95%)	7 (5%)	30	59
1	F	141/141 (100%)	134 (95%)	7 (5%)	30	59
1	G	142/141 (101%)	133 (94%)	9 (6%)	22	47

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	141/141 (100%)	134 (95%)	7 (5%)	30	59
1	I	141/141 (100%)	132 (94%)	9 (6%)	22	46
1	J	141/141 (100%)	135 (96%)	6 (4%)	35	65
1	K	141/141 (100%)	137 (97%)	4 (3%)	51	80
1	L	141/141 (100%)	135 (96%)	6 (4%)	35	65
All	All	1695/1692 (100%)	1612 (95%)	83 (5%)	31	60

All (83) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	17	SER
1	A	33	ASP
1	A	34	ASN
1	A	122	GLN
1	B	9	LYS
1	B	13[A]	GLU
1	B	13[B]	GLU
1	B	33	ASP
1	B	34	ASN
1	B	78	ARG
1	B	110	GLU
1	B	123	ILE
1	B	135	GLU
1	C	17	SER
1	C	29	SER
1	C	33	ASP
1	C	34	ASN
1	C	78	ARG
1	C	119	LEU
1	C	122	GLN
1	C	156	ARG
1	D	34	ASN
1	D	48	SER
1	D	64	LEU
1	D	81	GLN
1	D	119	LEU
1	D	148	LEU
1	D	156	ARG
1	E	29	SER
1	E	33	ASP

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Mol	Chain	Res	Type
1	E	34	ASN
1	E	48	SER
1	E	93	ILE
1	E	122	GLN
1	E	139	GLN
1	F	9	LYS
1	F	13	GLU
1	F	17	SER
1	F	33	ASP
1	F	107	LEU
1	F	119	LEU
1	F	122	GLN
1	G	13	GLU
1	G	33	ASP
1	G	34	ASN
1	G	78	ARG
1	G	93	ILE
1	G	112	GLN
1	G	119	LEU
1	G	122	GLN
1	G	139	GLN
1	H	2	GLN
1	H	29	SER
1	H	33	ASP
1	H	34	ASN
1	H	42	GLU
1	H	107	LEU
1	H	122	GLN
1	I	17	SER
1	I	29	SER
1	I	33	ASP
1	I	34	ASN
1	I	73	ARG
1	I	107	LEU
1	I	110	GLU
1	I	119	LEU
1	I	139	GLN
1	J	9	LYS
1	J	17	SER
1	J	29	SER
1	J	76	SER
1	J	122	GLN

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Mol	Chain	Res	Type
1	J	139	GLN
1	K	13	GLU
1	K	33	ASP
1	K	107	LEU
1	K	122	GLN
1	L	13	GLU
1	L	33	ASP
1	L	78	ARG
1	L	93	ILE
1	L	109	ARG
1	L	119	LEU

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

Mol	Chain	Res	Type
1	A	43	HIS
1	J	34	ASN
1	J	122	GLN
1	L	43	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 50 ligands modelled in this entry, 36 are monoatomic - leaving 14 for Mogul analysis.

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
4	HEM	B	301[A]	1	30,50,50	4.11	14 (46%)	24,82,82	2.66	12 (50%)
4	HEM	B	301[B]	1	30,50,50	4.07	15 (50%)	24,82,82	2.87	9 (37%)
4	HEM	C	301[A]	1	30,50,50	4.38	16 (53%)	24,82,82	2.68	10 (41%)
4	HEM	C	301[B]	1	30,50,50	4.48	17 (56%)	24,82,82	2.88	8 (33%)
5	EPE	D	502	-	14,15,15	0.46	0	18,20,20	2.09	4 (22%)
4	HEM	E	301[A]	1	30,50,50	4.39	14 (46%)	24,82,82	3.06	13 (54%)
4	HEM	E	301[B]	1	30,50,50	4.34	16 (53%)	24,82,82	2.91	12 (50%)
4	HEM	H	301[A]	1	30,50,50	4.27	15 (50%)	24,82,82	2.77	11 (45%)
4	HEM	H	301[B]	1	30,50,50	4.24	14 (46%)	24,82,82	2.92	10 (41%)
4	HEM	I	301[A]	1	30,50,50	4.42	14 (46%)	24,82,82	2.84	9 (37%)
4	HEM	I	301[B]	1	30,50,50	4.35	14 (46%)	24,82,82	2.75	8 (33%)
5	EPE	J	501	-	14,15,15	0.58	0	18,20,20	2.02	4 (22%)
4	HEM	K	301	1	30,50,50	4.49	17 (56%)	24,82,82	2.87	10 (41%)
4	HEM	L	301	1	30,50,50	4.40	14 (46%)	24,82,82	2.79	10 (41%)

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
4	HEM	B	301[A]	1	-	0/10/54/54	0/0/8/8
4	HEM	B	301[B]	1	-	0/10/54/54	0/0/8/8
4	HEM	C	301[A]	1	-	0/10/54/54	0/0/8/8
4	HEM	C	301[B]	1	-	0/10/54/54	0/0/8/8
5	EPE	D	502	-	-	0/9/19/19	0/1/1/1
4	HEM	E	301[A]	1	-	0/10/54/54	0/0/8/8
4	HEM	E	301[B]	1	-	0/10/54/54	0/0/8/8
4	HEM	H	301[A]	1	-	0/10/54/54	0/0/8/8
4	HEM	H	301[B]	1	-	0/10/54/54	0/0/8/8
4	HEM	I	301[A]	1	-	0/10/54/54	0/0/8/8
4	HEM	I	301[B]	1	-	0/10/54/54	0/0/8/8
5	EPE	J	501	-	-	0/9/19/19	0/1/1/1
4	HEM	K	301	1	-	0/10/54/54	0/0/8/8
4	HEM	L	301	1	-	0/10/54/54	0/0/8/8

All (180) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	I	301[A]	HEM	C3B-C4B	-18.17	1.35	1.51
4	K	301	HEM	C3B-C4B	-18.09	1.35	1.51
4	C	301[B]	HEM	C3B-C4B	-17.76	1.36	1.51
4	I	301[B]	HEM	C3B-C4B	-17.68	1.36	1.51
4	E	301[A]	HEM	C3B-C4B	-17.67	1.36	1.51
4	C	301[A]	HEM	C3B-C4B	-17.19	1.36	1.51
4	E	301[B]	HEM	C3B-C4B	-17.12	1.36	1.51
4	L	301	HEM	C3B-C4B	-17.10	1.36	1.51
4	H	301[A]	HEM	C3B-C4B	-16.90	1.36	1.51
4	H	301[B]	HEM	C3B-C4B	-16.77	1.36	1.51
4	B	301[A]	HEM	C3B-C4B	-15.89	1.37	1.51
4	B	301[B]	HEM	C3B-C4B	-15.39	1.38	1.51
4	C	301[B]	HEM	C3D-C4D	-12.44	1.35	1.51
4	L	301	HEM	C3D-C4D	-12.29	1.35	1.51
4	C	301[A]	HEM	C3D-C4D	-11.96	1.36	1.51
4	H	301[A]	HEM	C3D-C4D	-11.62	1.36	1.51
4	K	301	HEM	C3D-C4D	-11.57	1.36	1.51
4	E	301[B]	HEM	C3D-C4D	-11.39	1.36	1.51
4	E	301[A]	HEM	C3D-C4D	-11.39	1.36	1.51
4	B	301[B]	HEM	C3D-C4D	-11.33	1.37	1.51
4	I	301[A]	HEM	C3D-C4D	-11.30	1.37	1.51
4	B	301[A]	HEM	C3D-C4D	-11.15	1.37	1.51
4	H	301[B]	HEM	C3D-C4D	-11.12	1.37	1.51
4	I	301[B]	HEM	C3D-C4D	-11.10	1.37	1.51
4	L	301	HEM	C2C-C1C	-3.87	1.45	1.52
4	E	301[B]	HEM	C2D-C3D	-3.40	1.44	1.54
4	C	301[A]	HEM	C2D-C3D	-3.14	1.45	1.54
4	C	301[A]	HEM	C2C-C1C	-3.09	1.46	1.52
4	I	301[B]	HEM	C2C-C1C	-3.09	1.46	1.52
4	H	301[A]	HEM	C2D-C3D	-3.08	1.45	1.54
4	K	301	HEM	C2D-C3D	-3.08	1.45	1.54
4	H	301[B]	HEM	C2C-C1C	-3.01	1.46	1.52
4	E	301[A]	HEM	C2D-C3D	-2.99	1.45	1.54
4	B	301[B]	HEM	C2D-C3D	-2.96	1.45	1.54
4	C	301[B]	HEM	C2D-C3D	-2.94	1.45	1.54
4	I	301[A]	HEM	C2D-C3D	-2.92	1.45	1.54
4	B	301[A]	HEM	C2D-C3D	-2.87	1.46	1.54
4	C	301[B]	HEM	C3B-CAB	-2.86	1.46	1.51
4	E	301[A]	HEM	C3B-CAB	-2.82	1.46	1.51
4	I	301[B]	HEM	C2D-C3D	-2.81	1.46	1.54
4	E	301[B]	HEM	C2C-C1C	-2.79	1.47	1.52
4	H	301[B]	HEM	C2D-C3D	-2.78	1.46	1.54

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	301[A]	HEM	C2C-C1C	-2.76	1.47	1.52
4	E	301[B]	HEM	C4C-NC	-2.75	1.32	1.36
4	H	301[A]	HEM	C2C-C1C	-2.70	1.47	1.52
4	I	301[A]	HEM	C2C-C1C	-2.70	1.47	1.52
4	H	301[B]	HEM	C3B-CAB	-2.69	1.46	1.51
4	L	301	HEM	C2D-C3D	-2.67	1.46	1.54
4	E	301[B]	HEM	C3B-CAB	-2.66	1.46	1.51
4	K	301	HEM	C2C-C1C	-2.64	1.47	1.52
4	B	301[B]	HEM	C2C-C1C	-2.63	1.47	1.52
4	I	301[B]	HEM	C3B-CAB	-2.63	1.46	1.51
4	C	301[B]	HEM	C3C-CAC	-2.61	1.46	1.51
4	K	301	HEM	C3B-CAB	-2.59	1.46	1.51
4	E	301[B]	HEM	C3C-CAC	-2.52	1.46	1.51
4	E	301[B]	HEM	CAD-C3D	-2.49	1.49	1.54
4	K	301	HEM	C3C-CAC	-2.44	1.46	1.51
4	I	301[A]	HEM	C3B-CAB	-2.40	1.46	1.51
4	C	301[B]	HEM	C2C-C1C	-2.39	1.48	1.52
4	C	301[A]	HEM	C4C-NC	-2.37	1.33	1.36
4	K	301	HEM	CAD-C3D	-2.33	1.49	1.54
4	H	301[A]	HEM	C3B-CAB	-2.32	1.47	1.51
4	E	301[A]	HEM	C2C-C1C	-2.32	1.48	1.52
4	H	301[B]	HEM	C2B-C1B	-2.31	1.44	1.51
4	L	301	HEM	C3B-CAB	-2.31	1.47	1.51
4	K	301	HEM	C4C-NC	-2.26	1.33	1.36
4	C	301[B]	HEM	C4C-NC	-2.25	1.33	1.36
4	I	301[A]	HEM	C3C-CAC	-2.24	1.47	1.51
4	B	301[B]	HEM	C3B-CAB	-2.17	1.47	1.51
4	C	301[A]	HEM	C3C-CAC	-2.16	1.47	1.51
4	C	301[B]	HEM	C2D-C1D	-2.14	1.44	1.51
4	B	301[B]	HEM	C3C-CAC	-2.13	1.47	1.51
4	H	301[A]	HEM	CAD-C3D	-2.10	1.49	1.54
4	I	301[B]	HEM	C3C-CAC	-2.09	1.47	1.51
4	E	301[A]	HEM	C3C-CAC	-2.07	1.47	1.51
4	B	301[A]	HEM	CAD-C3D	-2.05	1.50	1.54
4	C	301[B]	HEM	CAD-C3D	-2.03	1.50	1.54
4	C	301[A]	HEM	C3B-CAB	-2.03	1.47	1.51
4	H	301[B]	HEM	FE-NB	2.00	2.08	1.97
4	E	301[A]	HEM	C4A-CHB	2.04	1.45	1.39
4	B	301[A]	HEM	FE-NB	2.05	2.08	1.97
4	B	301[B]	HEM	C4A-CHB	2.07	1.45	1.39
4	C	301[A]	HEM	C4A-CHB	2.08	1.45	1.39
4	K	301	HEM	FE-NB	2.09	2.08	1.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	301[A]	HEM	C4A-CHB	2.11	1.45	1.39
4	H	301[A]	HEM	FE-NB	2.11	2.08	1.97
4	C	301[B]	HEM	CHD-C1D	2.14	1.44	1.38
4	E	301[A]	HEM	C2A-C3A	2.19	1.44	1.37
4	C	301[A]	HEM	FE-NB	2.20	2.09	1.97
4	E	301[B]	HEM	C4A-CHB	2.20	1.45	1.39
4	L	301	HEM	FE-NB	2.21	2.09	1.97
4	B	301[B]	HEM	FE-NB	2.27	2.09	1.97
4	C	301[B]	HEM	C1A-CHA	2.27	1.46	1.39
4	C	301[A]	HEM	CHC-C4B	2.33	1.45	1.38
4	L	301	HEM	C4A-CHB	2.34	1.46	1.39
4	C	301[B]	HEM	C4A-CHB	2.34	1.46	1.39
4	B	301[A]	HEM	C4A-CHB	2.35	1.46	1.39
4	I	301[B]	HEM	C4A-CHB	2.37	1.46	1.39
4	H	301[B]	HEM	CHC-C4B	2.40	1.45	1.38
4	K	301	HEM	C1A-CHA	2.41	1.46	1.39
4	H	301[B]	HEM	C2A-C3A	2.42	1.44	1.37
4	L	301	HEM	C1A-CHA	2.42	1.46	1.39
4	E	301[B]	HEM	CHD-C1D	2.43	1.45	1.38
4	I	301[A]	HEM	C4A-CHB	2.44	1.46	1.39
4	C	301[A]	HEM	C1A-CHA	2.45	1.46	1.39
4	H	301[A]	HEM	CHD-C1D	2.45	1.45	1.38
4	I	301[A]	HEM	CHD-C1D	2.46	1.45	1.38
4	I	301[B]	HEM	C1A-CHA	2.46	1.46	1.39
4	B	301[A]	HEM	CHC-C4B	2.47	1.45	1.38
4	L	301	HEM	CHC-C4B	2.48	1.45	1.38
4	I	301[A]	HEM	C1A-CHA	2.52	1.46	1.39
4	E	301[A]	HEM	C1A-CHA	2.53	1.46	1.39
4	H	301[B]	HEM	C1A-CHA	2.55	1.46	1.39
4	E	301[B]	HEM	C1A-CHA	2.56	1.46	1.39
4	B	301[A]	HEM	C1A-CHA	2.56	1.46	1.39
4	C	301[B]	HEM	C2A-C3A	2.58	1.45	1.37
4	B	301[B]	HEM	CHC-C4B	2.59	1.46	1.38
4	I	301[A]	HEM	CHC-C4B	2.60	1.46	1.38
4	I	301[B]	HEM	CHD-C1D	2.61	1.46	1.38
4	E	301[B]	HEM	C2A-C3A	2.63	1.45	1.37
4	B	301[A]	HEM	CHD-C1D	2.64	1.46	1.38
4	C	301[A]	HEM	CHD-C1D	2.66	1.46	1.38
4	I	301[A]	HEM	C2A-C3A	2.67	1.45	1.37
4	B	301[A]	HEM	C2A-C3A	2.67	1.45	1.37
4	K	301	HEM	CHC-C4B	2.68	1.46	1.38
4	H	301[A]	HEM	CHC-C4B	2.68	1.46	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	301[B]	HEM	CHC-C4B	2.68	1.46	1.38
4	K	301	HEM	C2A-C3A	2.71	1.45	1.37
4	C	301[A]	HEM	C2A-C3A	2.71	1.45	1.37
4	H	301[A]	HEM	C2A-C3A	2.75	1.45	1.37
4	I	301[B]	HEM	CHC-C4B	2.75	1.46	1.38
4	K	301	HEM	C4A-CHB	2.76	1.47	1.39
4	E	301[A]	HEM	CHD-C1D	2.77	1.46	1.38
4	E	301[A]	HEM	CHC-C4B	2.77	1.46	1.38
4	B	301[B]	HEM	C2A-C3A	2.77	1.45	1.37
4	I	301[B]	HEM	C2A-C3A	2.78	1.45	1.37
4	H	301[A]	HEM	C1A-CHA	2.78	1.47	1.39
4	E	301[B]	HEM	CHC-C4B	2.80	1.46	1.38
4	K	301	HEM	CHD-C1D	2.80	1.46	1.38
4	H	301[B]	HEM	CHD-C1D	2.82	1.46	1.38
4	B	301[B]	HEM	CHD-C1D	2.84	1.46	1.38
4	L	301	HEM	C2A-C3A	2.95	1.46	1.37
4	I	301[A]	HEM	FE-NC	2.99	2.07	1.95
4	B	301[B]	HEM	C1A-CHA	3.03	1.48	1.39
4	C	301[A]	HEM	FE-NC	3.04	2.07	1.95
4	L	301	HEM	CHD-C1D	3.04	1.47	1.38
4	L	301	HEM	FE-NC	3.05	2.07	1.95
4	B	301[B]	HEM	FE-NC	3.08	2.07	1.95
4	I	301[B]	HEM	FE-NC	3.09	2.08	1.95
4	E	301[A]	HEM	FE-NC	3.19	2.08	1.95
4	H	301[A]	HEM	FE-NC	3.20	2.08	1.95
4	E	301[B]	HEM	FE-NC	3.20	2.08	1.95
4	C	301[B]	HEM	FE-NC	3.22	2.08	1.95
4	H	301[B]	HEM	FE-NC	3.30	2.08	1.95
4	K	301	HEM	FE-NC	3.37	2.09	1.95
4	B	301[A]	HEM	FE-NC	3.38	2.09	1.95
4	E	301[B]	HEM	CHD-C4C	3.62	1.44	1.36
4	H	301[A]	HEM	CHD-C4C	3.65	1.45	1.36
4	C	301[B]	HEM	CHD-C4C	3.85	1.45	1.36
4	B	301[A]	HEM	CHC-C1C	3.93	1.45	1.36
4	I	301[B]	HEM	CHC-C1C	4.06	1.45	1.36
4	C	301[A]	HEM	CHC-C1C	4.08	1.46	1.36
4	H	301[B]	HEM	CHC-C1C	4.09	1.46	1.36
4	E	301[A]	HEM	CHD-C4C	4.14	1.46	1.36
4	B	301[B]	HEM	CHD-C4C	4.16	1.46	1.36
4	I	301[A]	HEM	CHD-C4C	4.16	1.46	1.36
4	K	301	HEM	CHC-C1C	4.19	1.46	1.36
4	C	301[A]	HEM	CHD-C4C	4.20	1.46	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	K	301	HEM	CHD-C4C	4.20	1.46	1.36
4	I	301[B]	HEM	CHD-C4C	4.21	1.46	1.36
4	H	301[A]	HEM	CHC-C1C	4.22	1.46	1.36
4	L	301	HEM	CHC-C1C	4.24	1.46	1.36
4	I	301[A]	HEM	CHC-C1C	4.30	1.46	1.36
4	C	301[B]	HEM	CHC-C1C	4.31	1.46	1.36
4	E	301[B]	HEM	CHC-C1C	4.34	1.46	1.36
4	B	301[A]	HEM	CHD-C4C	4.40	1.46	1.36
4	B	301[B]	HEM	CHC-C1C	4.42	1.46	1.36
4	E	301[A]	HEM	CHC-C1C	4.43	1.46	1.36
4	L	301	HEM	CHD-C4C	4.54	1.47	1.36
4	H	301[B]	HEM	CHD-C4C	4.60	1.47	1.36

All (130) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	301[B]	HEM	C3B-CAB-CBB	-6.97	113.77	124.46
4	E	301[A]	HEM	C3B-CAB-CBB	-6.06	115.16	124.46
4	E	301[A]	HEM	CAA-CBA-CGA	-5.94	101.86	112.75
4	I	301[B]	HEM	C3B-CAB-CBB	-5.69	115.73	124.46
4	K	301	HEM	C3B-CAB-CBB	-5.64	115.80	124.46
4	E	301[B]	HEM	C3B-CAB-CBB	-5.63	115.82	124.46
4	I	301[A]	HEM	CBA-CAA-C2A	-5.37	102.90	112.53
4	C	301[A]	HEM	C3B-CAB-CBB	-5.13	116.58	124.46
4	H	301[B]	HEM	C3B-CAB-CBB	-5.07	116.67	124.46
4	H	301[A]	HEM	C3B-CAB-CBB	-4.83	117.04	124.46
4	L	301	HEM	C3B-CAB-CBB	-4.66	117.31	124.46
4	I	301[A]	HEM	C3B-CAB-CBB	-4.45	117.63	124.46
4	E	301[B]	HEM	CBA-CAA-C2A	-4.35	104.73	112.53
4	C	301[B]	HEM	C3C-CAC-CBC	-4.11	118.15	124.46
4	H	301[B]	HEM	CBA-CAA-C2A	-4.04	105.29	112.53
4	B	301[B]	HEM	C3B-CAB-CBB	-3.82	118.59	124.46
4	K	301	HEM	C3C-CAC-CBC	-3.55	119.01	124.46
4	B	301[A]	HEM	C3C-CAC-CBC	-3.52	119.05	124.46
4	L	301	HEM	CBA-CAA-C2A	-3.40	106.43	112.53
4	K	301	HEM	CBD-CAD-C3D	-3.36	103.76	113.55
4	E	301[B]	HEM	C3C-CAC-CBC	-3.31	119.37	124.46
4	C	301[A]	HEM	C3C-CAC-CBC	-3.26	119.45	124.46
4	I	301[A]	HEM	C3C-CAC-CBC	-3.15	119.63	124.46
4	I	301[B]	HEM	C3C-CAC-CBC	-3.03	119.81	124.46
4	E	301[A]	HEM	CBA-CAA-C2A	-2.94	107.26	112.53
4	E	301[A]	HEM	CBD-CAD-C3D	-2.90	105.13	113.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	301[A]	HEM	C3B-CAB-CBB	-2.88	120.05	124.46
4	L	301	HEM	C2C-C1C-CHC	-2.81	119.40	123.68
4	E	301[B]	HEM	CAD-CBD-CGD	-2.76	101.77	113.02
4	H	301[B]	HEM	C3C-CAC-CBC	-2.66	120.37	124.46
4	B	301[A]	HEM	CBA-CAA-C2A	-2.66	107.76	112.53
4	E	301[A]	HEM	C3C-CAC-CBC	-2.52	120.59	124.46
4	H	301[A]	HEM	CBD-CAD-C3D	-2.47	106.37	113.55
4	C	301[A]	HEM	CBA-CAA-C2A	-2.43	108.17	112.53
4	B	301[A]	HEM	CBD-CAD-C3D	-2.00	107.72	113.55
4	E	301[A]	HEM	CMA-C3A-C2A	2.00	129.43	125.24
4	H	301[A]	HEM	CMA-C3A-C2A	2.04	129.50	125.24
4	H	301[B]	HEM	C4B-CHC-C1C	2.05	129.25	125.82
4	E	301[B]	HEM	CMA-C3A-C2A	2.12	129.67	125.24
4	B	301[A]	HEM	C2D-C3D-C4D	2.13	105.11	101.50
4	K	301	HEM	CAA-C2A-C1A	2.17	129.37	127.01
4	H	301[A]	HEM	C2D-C3D-C4D	2.20	105.22	101.50
5	J	501	EPE	C7-N4-C3	2.22	116.97	111.27
4	L	301	HEM	CMA-C3A-C2A	2.25	129.94	125.24
4	E	301[B]	HEM	C2D-C3D-C4D	2.30	105.40	101.50
4	B	301[A]	HEM	C4B-CHC-C1C	2.45	129.91	125.82
5	D	502	EPE	C7-N4-C3	2.58	117.88	111.27
4	C	301[A]	HEM	CMA-C3A-C2A	2.59	130.65	125.24
4	E	301[A]	HEM	CAA-C2A-C1A	2.66	129.90	127.01
4	B	301[B]	HEM	CMA-C3A-C2A	2.71	130.90	125.24
4	H	301[A]	HEM	C2C-C1C-NC	2.72	114.79	110.21
5	J	501	EPE	C7-N4-C5	2.82	118.50	111.27
4	C	301[A]	HEM	C2C-C1C-NC	2.88	115.06	110.21
4	C	301[A]	HEM	CMD-C2D-C3D	2.95	127.42	114.35
4	B	301[B]	HEM	C2C-C1C-NC	2.98	115.23	110.21
4	I	301[A]	HEM	C2C-C1C-NC	2.99	115.25	110.21
4	E	301[A]	HEM	C2C-C1C-NC	3.02	115.30	110.21
4	C	301[B]	HEM	C2C-C1C-NC	3.14	115.51	110.21
4	B	301[B]	HEM	CMD-C2D-C3D	3.19	128.46	114.35
4	L	301	HEM	CMD-C2D-C3D	3.20	128.53	114.35
4	B	301[A]	HEM	C2C-C1C-NC	3.23	115.66	110.21
4	E	301[B]	HEM	C2C-C1C-NC	3.26	115.70	110.21
4	H	301[B]	HEM	C2C-C1C-NC	3.27	115.72	110.21
5	D	502	EPE	C7-N4-C5	3.29	119.69	111.27
4	K	301	HEM	C2C-C1C-NC	3.35	115.86	110.21
4	H	301[A]	HEM	CMD-C2D-C3D	3.36	129.22	114.35
4	E	301[A]	HEM	CMD-C2D-C3D	3.37	129.24	114.35
4	E	301[B]	HEM	CMD-C2D-C3D	3.37	129.24	114.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	D	502	EPE	O2S-S-C10	3.38	109.79	106.91
4	I	301[B]	HEM	C2C-C1C-NC	3.41	115.96	110.21
4	I	301[A]	HEM	CMD-C2D-C3D	3.48	129.72	114.35
4	H	301[B]	HEM	CMD-C2D-C3D	3.48	129.74	114.35
4	K	301	HEM	CMD-C2D-C3D	3.51	129.86	114.35
4	I	301[B]	HEM	CMD-C2D-C3D	3.61	130.31	114.35
4	L	301	HEM	C2C-C1C-NC	3.65	116.36	110.21
4	C	301[B]	HEM	CMD-C2D-C3D	3.70	130.73	114.35
4	B	301[A]	HEM	CMD-C2D-C3D	3.71	130.78	114.35
4	B	301[A]	HEM	CMC-C2C-C3C	3.80	126.03	116.53
4	H	301[B]	HEM	CAD-C3D-C4D	3.82	125.94	112.47
4	K	301	HEM	CAD-C3D-C4D	3.85	126.06	112.47
4	L	301	HEM	CAD-C3D-C4D	3.96	126.42	112.47
4	E	301[A]	HEM	CMC-C2C-C3C	3.97	126.44	116.53
4	C	301[B]	HEM	CAD-C3D-C4D	3.98	126.51	112.47
4	I	301[A]	HEM	CAD-C3D-C4D	4.05	126.74	112.47
4	I	301[B]	HEM	CAD-C3D-C4D	4.06	126.78	112.47
4	H	301[A]	HEM	CAD-C3D-C4D	4.10	126.94	112.47
4	B	301[A]	HEM	CAD-C3D-C4D	4.18	127.20	112.47
4	E	301[B]	HEM	CAD-C3D-C2D	4.22	125.34	113.22
4	B	301[B]	HEM	CAD-C3D-C4D	4.30	127.64	112.47
4	C	301[A]	HEM	CMC-C2C-C3C	4.36	127.42	116.53
4	E	301[A]	HEM	CAD-C3D-C4D	4.38	127.91	112.47
4	H	301[A]	HEM	CMC-C2C-C3C	4.40	127.51	116.53
5	J	501	EPE	C5-N4-C3	4.42	118.48	108.90
4	E	301[B]	HEM	CMC-C2C-C3C	4.47	127.69	116.53
4	C	301[A]	HEM	CAD-C3D-C4D	4.53	128.46	112.47
4	B	301[B]	HEM	CMC-C2C-C3C	4.54	127.87	116.53
4	C	301[B]	HEM	CMC-C2C-C3C	4.57	127.94	116.53
4	H	301[B]	HEM	CMC-C2C-C3C	4.59	127.98	116.53
4	H	301[A]	HEM	CAA-C2A-C1A	4.68	132.09	127.01
4	I	301[A]	HEM	CMC-C2C-C3C	4.72	128.32	116.53
4	E	301[B]	HEM	CAD-C3D-C4D	4.76	129.26	112.47
4	C	301[A]	HEM	CMB-C2B-C3B	4.84	128.62	116.53
4	C	301[A]	HEM	CAD-C3D-C2D	4.93	127.39	113.22
5	J	501	EPE	O2S-S-C10	5.02	111.19	106.91
4	B	301[A]	HEM	CAD-C3D-C2D	5.03	127.67	113.22
4	L	301	HEM	CMC-C2C-C3C	5.05	129.13	116.53
4	K	301	HEM	CMC-C2C-C3C	5.06	129.16	116.53
4	H	301[A]	HEM	CAD-C3D-C2D	5.08	127.82	113.22
4	E	301[A]	HEM	CAD-C3D-C2D	5.10	127.87	113.22
4	L	301	HEM	CMB-C2B-C3B	5.13	129.34	116.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	301[B]	HEM	CMB-C2B-C3B	5.15	129.38	116.53
4	I	301[B]	HEM	CMC-C2C-C3C	5.22	129.55	116.53
4	K	301	HEM	CMB-C2B-C3B	5.26	129.67	116.53
4	C	301[B]	HEM	CMB-C2B-C3B	5.29	129.73	116.53
4	B	301[B]	HEM	CAD-C3D-C2D	5.30	128.45	113.22
4	I	301[A]	HEM	CMB-C2B-C3B	5.30	129.76	116.53
4	E	301[A]	HEM	CMB-C2B-C3B	5.39	129.98	116.53
4	B	301[B]	HEM	CAA-C2A-C1A	5.45	132.93	127.01
4	K	301	HEM	CAD-C3D-C2D	5.50	129.04	113.22
4	I	301[A]	HEM	CAD-C3D-C2D	5.50	129.04	113.22
4	I	301[B]	HEM	CAD-C3D-C2D	5.51	129.07	113.22
4	C	301[B]	HEM	CAD-C3D-C2D	5.61	129.35	113.22
4	L	301	HEM	CAD-C3D-C2D	5.66	129.48	113.22
4	E	301[B]	HEM	CMB-C2B-C3B	5.71	130.79	116.53
4	H	301[B]	HEM	CAD-C3D-C2D	5.75	129.74	113.22
4	H	301[A]	HEM	CMB-C2B-C3B	5.79	130.99	116.53
4	B	301[A]	HEM	CMB-C2B-C3B	5.92	131.31	116.53
4	B	301[B]	HEM	CMB-C2B-C3B	5.99	131.49	116.53
5	D	502	EPE	C5-N4-C3	6.09	122.08	108.90
4	H	301[B]	HEM	CMB-C2B-C3B	6.79	133.47	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

12 monomers are involved in 76 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	301[A]	HEM	11	0
4	B	301[B]	HEM	9	0
4	C	301[A]	HEM	7	0
4	C	301[B]	HEM	6	0
4	E	301[A]	HEM	5	0
4	E	301[B]	HEM	7	0
4	H	301[A]	HEM	4	0
4	H	301[B]	HEM	5	0
4	I	301[A]	HEM	2	0
4	I	301[B]	HEM	2	0
4	K	301	HEM	15	0
4	L	301	HEM	3	0

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	161/161 (100%)	0.01	3 (1%) 70 71	38, 46, 53, 63	10 (6%)
1	B	161/161 (100%)	-0.08	3 (1%) 70 71	38, 46, 54, 62	8 (4%)
1	C	161/161 (100%)	0.00	4 (2%) 61 61	38, 46, 53, 63	9 (5%)
1	D	161/161 (100%)	-0.05	3 (1%) 70 71	38, 46, 53, 62	11 (6%)
1	E	161/161 (100%)	0.06	5 (3%) 52 53	38, 46, 53, 63	12 (7%)
1	F	161/161 (100%)	0.06	2 (1%) 81 81	37, 46, 53, 62	12 (7%)
1	G	161/161 (100%)	-0.02	5 (3%) 52 53	38, 46, 53, 65	9 (5%)
1	H	161/161 (100%)	-0.04	4 (2%) 61 61	38, 46, 53, 63	15 (9%)
1	I	161/161 (100%)	0.02	2 (1%) 81 81	38, 45, 53, 63	9 (5%)
1	J	161/161 (100%)	0.03	6 (3%) 45 46	38, 45, 53, 65	10 (6%)
1	K	161/161 (100%)	0.00	3 (1%) 70 71	38, 46, 54, 63	10 (6%)
1	L	161/161 (100%)	0.01	3 (1%) 70 71	37, 46, 53, 63	11 (6%)
All	All	1932/1932 (100%)	-0.00	43 (2%) 65 66	37, 46, 54, 65	126 (6%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	161	ALA	5.1
1	J	161	ALA	5.1
1	I	160	SER	3.9
1	H	161	ALA	3.3
1	B	6	ASP	3.3
1	C	4	ASP	3.2
1	H	46	ALA	3.2
1	H	6	ASP	3.2
1	A	76	SER	3.1
1	D	160	SER	3.1
1	C	6	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	I	95	TYR	3.0
1	K	6	ASP	3.0
1	E	146	ASP	2.9
1	J	3	GLY	2.8
1	C	2	GLN	2.8
1	J	2	GLN	2.8
1	G	4	ASP	2.8
1	H	5	PRO	2.7
1	G	6	ASP	2.7
1	L	5	PRO	2.7
1	E	159	GLY	2.6
1	J	6	ASP	2.6
1	G	8	LEU	2.5
1	E	160	SER	2.5
1	B	160	SER	2.5
1	F	146	ASP	2.5
1	E	5	PRO	2.4
1	J	160	SER	2.4
1	G	5	PRO	2.4
1	L	160	SER	2.4
1	F	142	ASP	2.4
1	J	4	ASP	2.4
1	A	159	GLY	2.3
1	L	99	GLU	2.3
1	K	5	PRO	2.2
1	E	4	ASP	2.2
1	K	161	ALA	2.2
1	B	161	ALA	2.2
1	D	6	ASP	2.2
1	C	79	VAL	2.2
1	A	160	SER	2.1
1	G	81	GLN	2.1

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	EPE	J	501	15/15	0.78	0.28	2.49	57,61,71,73	15
3	MG	A	401	1/1	0.93	0.21	2.12	51,51,51,51	0
3	MG	B	401	1/1	0.67	0.23	1.80	58,58,58,58	0
4	HEM	H	301[B]	43/43	0.88	0.26	1.52	50,51,59,63	43
4	HEM	H	301[A]	43/43	0.88	0.26	1.52	30,32,44,46	43
5	EPE	D	502	15/15	0.83	0.26	1.51	49,55,66,67	15
4	HEM	B	301[A]	43/43	0.88	0.28	1.40	40,43,57,61	43
4	HEM	B	301[B]	43/43	0.88	0.28	1.40	48,49,59,62	43
4	HEM	C	301[B]	43/43	0.90	0.22	0.81	40,42,55,58	43
4	HEM	C	301[A]	43/43	0.90	0.22	0.78	41,45,56,59	43
4	HEM	K	301	43/43	0.86	0.21	0.54	46,46,55,56	43
4	HEM	E	301[B]	43/43	0.87	0.23	0.50	45,48,55,57	43
4	HEM	E	301[A]	43/43	0.87	0.23	0.50	47,49,58,60	43
4	HEM	I	301[A]	43/43	0.88	0.22	0.50	40,44,53,59	43
4	HEM	I	301[B]	43/43	0.88	0.22	0.45	56,57,64,70	43
4	HEM	L	301	43/43	0.88	0.21	0.42	52,54,57,60	43
3	MG	C	401	1/1	0.91	0.19	0.42	57,57,57,57	0
3	MG	F	401	1/1	0.94	0.17	0.20	46,46,46,46	0
3	MG	L	401	1/1	0.90	0.16	-0.22	50,50,50,50	0
3	MG	H	401	1/1	0.87	0.15	-0.30	47,47,47,47	0
3	MG	B	402	1/1	0.83	0.14	-0.52	50,50,50,50	0
3	MG	I	401	1/1	0.94	0.13	-0.55	51,51,51,51	0
3	MG	E	401	1/1	0.78	0.12	-1.15	60,60,60,60	0
2	ZN	D	202	1/1	0.99	0.10	-1.63	48,48,48,48	0
2	ZN	H	202	1/1	0.99	0.12	-1.65	52,52,52,52	0
2	ZN	C	201	1/1	0.99	0.07	-1.88	44,44,44,44	0
2	ZN	A	201	1/1	0.99	0.12	-1.94	39,39,39,39	0
3	MG	K	401	1/1	0.85	0.10	-1.94	40,40,40,40	0
2	ZN	J	202	1/1	0.99	0.10	-2.08	48,48,48,48	0
2	ZN	K	201	1/1	1.00	0.08	-2.12	43,43,43,43	0
2	ZN	I	202	1/1	0.99	0.11	-2.12	53,53,53,53	0
3	MG	J	401	1/1	0.92	0.10	-2.20	57,57,57,57	0
2	ZN	A	202	1/1	0.98	0.10	-2.53	53,53,53,53	0
2	ZN	I	201	1/1	0.99	0.08	-2.64	44,44,44,44	0
2	ZN	F	202	1/1	0.99	0.09	-2.75	60,60,60,60	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	ZN	L	201	1/1	0.99	0.07	-2.95	43,43,43,43	0
3	MG	C	402	1/1	0.91	0.09	-3.28	55,55,55,55	0
2	ZN	G	201	1/1	0.99	0.05	-3.86	37,37,37,37	0
2	ZN	B	201	1/1	0.99	0.07	-4.04	41,41,41,41	0
2	ZN	E	202	1/1	0.99	0.07	-4.32	55,55,55,55	0
2	ZN	J	201	1/1	0.99	0.06	-4.46	41,41,41,41	0
2	ZN	K	202	1/1	0.99	0.06	-4.96	55,55,55,55	0
2	ZN	G	202	1/1	0.99	0.06	-5.57	49,49,49,49	0
2	ZN	D	201	1/1	0.99	0.09	-5.88	39,39,39,39	0
2	ZN	F	201	1/1	0.97	0.07	-6.00	48,48,48,48	0
2	ZN	E	201	1/1	0.99	0.08	-9.30	44,44,44,44	0
2	ZN	L	202	1/1	0.97	0.08	-	65,65,65,65	0
2	ZN	C	202	1/1	0.99	0.09	-	49,49,49,49	0
2	ZN	B	202	1/1	0.99	0.11	-	53,53,53,53	0
2	ZN	H	201	1/1	0.99	0.09	-	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.