



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:46 PM GMT

PDB ID : 1V4U
Title : Crystal structure of bluefin tuna carbonmonoxy-hemoglobin
Authors : Yokoyama, T.; Chong, K.T.; Miyazaki, Y.; Nakatsukasa, T.; Unzai, S.;
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Deposited on : 2003-11-19
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

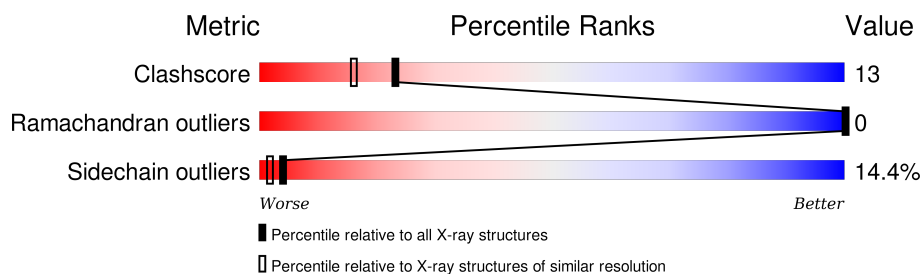
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	144	 65% 27% 7% •
1	C	144	 62% 28% 10%
2	B	146	 71% 21% 5% • •
2	D	146	 70% 21% 5% • •

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
4	CMO	A	145	-	-	X	-
4	CMO	B	148	-	-	X	-
4	CMO	D	148	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 4694 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 hemoglobin alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	144	Total	C	N	O	S	0	0	0
			1088	699	184	199	6			
1	C	144	Total	C	N	O	S	0	0	0
			1088	699	184	199	6			

- Molecule 2 is a protein called hemoglobin beta chain.

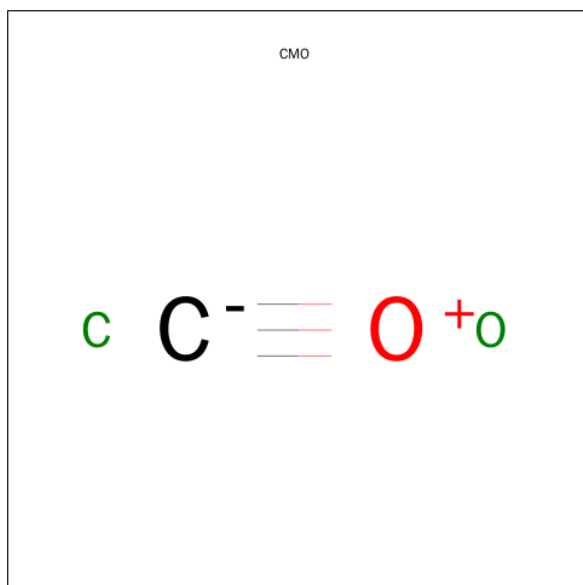
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	142	Total	C	N	O	S	0	0	0
			1100	706	187	203	4			
2	D	142	Total	C	N	O	S	0	0	0
			1100	706	187	203	4			

- 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 CARBON MONOXIDE (three-letter code: CMO) (formula: CO).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	42	Total	O	0	0
			42	42		
5	B	42	Total	O	0	0
			42	42		

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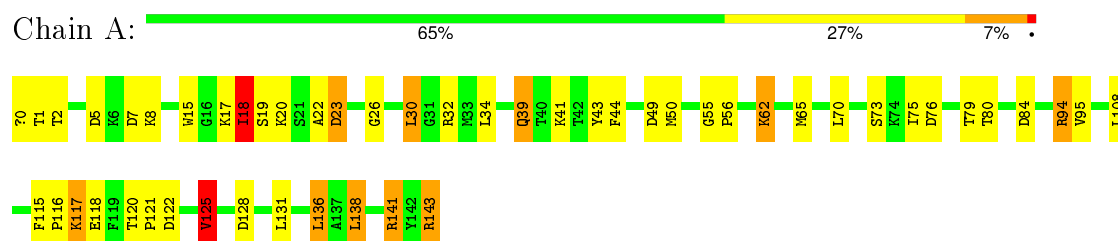
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	C	35	Total	O	0	0
			35	35		
5	D	19	Total	O	0	0
			19	19		

3 Residue-property plots [i](#)

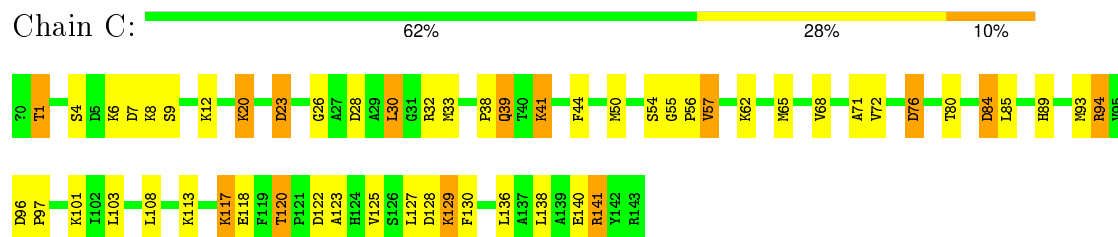
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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.

Note EDS was not executed.

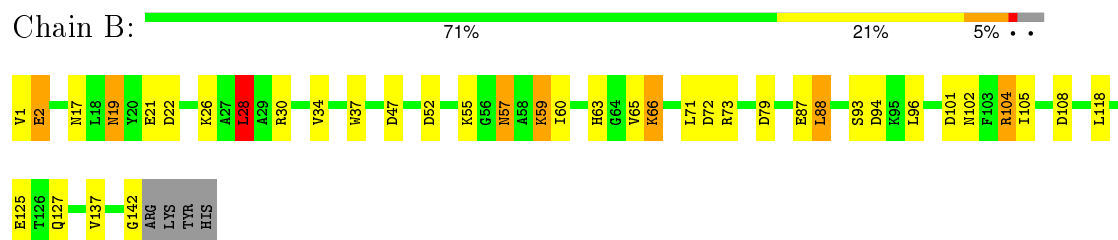
- Molecule 1: hemoglobin alpha chain



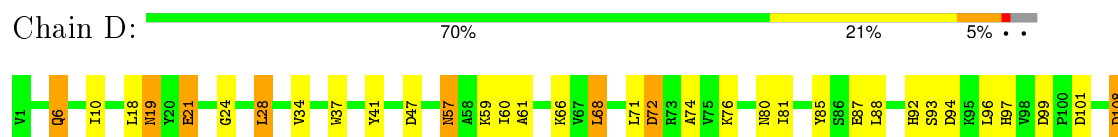
- Molecule 1: hemoglobin alpha chain



- Molecule 2: hemoglobin beta chain



- Molecule 2: hemoglobin beta chain



G109	L110	F111	V112	G119	E126	F126	Q127	V137	G142	ARG	LYS	TYR	HIS
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4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section will therefore be incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.02Å 102.54Å 108.45Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.00	Depositor
% Data completeness (in resolution range)	100.0 (20.00-2.00)	Depositor
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.1.22	Depositor
R, R_{free}	0.204 , 0.264	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4694	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

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: CMO, HEM, ACE

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	1.28	2/1110 (0.2%)	1.32	18/1499 (1.2%)
1	C	1.25	0/1110	1.29	10/1499 (0.7%)
2	B	1.25	1/1123 (0.1%)	1.17	9/1527 (0.6%)
2	D	1.18	1/1123 (0.1%)	1.12	7/1527 (0.5%)
All	All	1.24	4/4466 (0.1%)	1.23	44/6052 (0.7%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	125	GLU	CD-OE2	5.27	1.31	1.25
1	A	22	ALA	CA-CB	5.16	1.63	1.52
2	D	61	ALA	CA-CB	5.14	1.63	1.52
1	A	44	PHE	CE1-CZ	5.06	1.47	1.37

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	141	ARG	NE-CZ-NH2	10.38	125.49	120.30
1	A	128	ASP	CB-CG-OD2	8.95	126.36	118.30
1	C	76	ASP	CB-CG-OD2	8.84	126.26	118.30
1	C	28	ASP	CB-CG-OD1	8.82	126.24	118.30
1	C	23	ASP	CB-CG-OD2	8.28	125.75	118.30
2	D	72	ASP	CB-CG-OD2	8.18	125.67	118.30
2	B	52	ASP	CB-CG-OD2	7.79	125.31	118.30
1	A	125	VAL	CB-CA-C	-7.49	97.17	111.40
2	D	99	ASP	CB-CG-OD2	7.47	125.03	118.30
2	B	72	ASP	CB-CG-OD2	7.25	124.83	118.30
1	A	30	LEU	CB-CG-CD2	7.13	123.12	111.00
1	A	143	ARG	NE-CZ-NH1	-6.86	116.87	120.30
2	D	112	VAL	CG1-CB-CG2	6.72	121.65	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	141	ARG	NE-CZ-NH1	-6.67	116.97	120.30
1	A	1	THR	CA-CB-CG2	6.58	121.62	112.40
2	B	30	ARG	NE-CZ-NH2	-6.42	117.09	120.30
1	A	141	ARG	NE-CZ-NH1	-6.38	117.11	120.30
1	C	128	ASP	CB-CG-OD2	6.34	124.00	118.30
2	B	108	ASP	CB-CG-OD1	6.33	124.00	118.30
1	A	1	THR	N-CA-CB	6.22	122.11	110.30
1	A	49	ASP	CB-CG-OD2	6.18	123.87	118.30
1	C	76	ASP	CB-CG-OD1	-6.14	112.78	118.30
1	A	23	ASP	CB-CG-OD2	6.12	123.81	118.30
2	B	47	ASP	CB-CG-OD2	5.94	123.65	118.30
1	C	129	LYS	CD-CE-NZ	-5.90	98.13	111.70
1	C	7	ASP	CB-CG-OD2	5.90	123.61	118.30
1	A	18	ILE	CG1-CB-CG2	5.81	124.18	111.40
2	D	47	ASP	CB-CG-OD2	5.80	123.52	118.30
2	B	94	ASP	CB-CG-OD2	5.63	123.36	118.30
2	D	108	ASP	CB-CG-OD1	5.60	123.34	118.30
1	A	84	ASP	CB-CG-OD2	5.57	123.31	118.30
1	A	138	LEU	CB-CG-CD1	5.56	120.46	111.00
1	C	85	LEU	CB-CG-CD1	-5.44	101.75	111.00
2	B	79	ASP	CB-CG-OD2	5.38	123.14	118.30
2	D	110	LEU	CB-CG-CD2	-5.32	101.96	111.00
1	A	136	LEU	CB-CG-CD2	-5.31	101.97	111.00
1	A	5	ASP	CB-CG-OD1	5.18	122.96	118.30
1	C	84	ASP	CB-CG-OD2	5.16	122.94	118.30
2	D	94	ASP	CB-CG-OD2	5.14	122.93	118.30
1	A	7	ASP	CB-CG-OD2	5.11	122.90	118.30
2	B	28	LEU	CB-CG-CD2	5.11	119.69	111.00
2	B	104	ARG	NE-CZ-NH2	5.10	122.85	120.30
1	A	76	ASP	CB-CG-OD2	5.09	122.89	118.30
1	A	0	ACE	C-N-CA	-5.06	109.05	121.70

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	1088	0	1120	39	0
1	C	1088	0	1120	39	0
2	B	1100	0	1093	22	0
2	D	1100	0	1093	22	0
3	A	43	0	30	4	0
3	B	43	0	30	5	0
3	C	43	0	30	2	0
3	D	43	0	30	6	0
4	A	2	0	0	4	0
4	B	2	0	0	4	0
4	C	2	0	0	0	0
4	D	2	0	0	4	0
5	A	42	0	0	0	0
5	B	42	0	0	0	0
5	C	35	0	0	0	0
5	D	19	0	0	0	0
All	All	4694	0	4546	119	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 13.

All (119) 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:117:LYS:HD2	1:C:117:LYS:H	1.31	0.93
3:A:144:HEM:ND	4:A:145:CMO:C	2.32	0.92
1:C:32:ARG:HH11	2:D:127:GLN:HE21	1.11	0.92
1:A:117:LYS:HD2	1:A:117:LYS:H	1.37	0.90
3:B:147:HEM:ND	4:B:148:CMO:C	2.34	0.90
1:A:41:LYS:HD3	1:A:50:MET:HE2	1.51	0.90
1:A:41:LYS:CG	1:A:50:MET:HE1	2.05	0.87
2:B:57:ASN:HD21	2:B:59:LYS:HD2	1.41	0.86
1:A:41:LYS:HD3	1:A:50:MET:CE	2.05	0.85
1:C:32:ARG:HH11	2:D:127:GLN:NE2	1.76	0.84
1:A:117:LYS:CD	1:A:117:LYS:H	1.87	0.83
3:A:144:HEM:NC	4:A:145:CMO:C	2.43	0.81
3:B:147:HEM:NC	4:B:148:CMO:C	2.43	0.81
1:A:32:ARG:HH11	2:B:127:GLN:HE21	1.28	0.80
1:A:39:GLN:HE21	1:A:39:GLN:H	1.29	0.80
1:A:117:LYS:HD3	1:A:118:GLU:OE1	1.84	0.77
1:C:117:LYS:HD3	1:C:118:GLU:OE1	1.85	0.76
3:D:147:HEM:ND	4:D:148:CMO:C	2.49	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:120:THR:CG2	1:C:123:ALA:H	1.99	0.76
1:A:32:ARG:HH11	2:B:127:GLN:NE2	1.83	0.76
1:A:136:LEU:HD11	1:C:1:THR:HB	1.69	0.74
1:C:120:THR:HG22	1:C:123:ALA:H	1.54	0.73
1:C:39:GLN:H	1:C:39:GLN:HE21	1.36	0.73
1:C:117:LYS:CD	1:C:117:LYS:H	2.04	0.71
1:A:41:LYS:HG2	1:A:50:MET:HE1	1.74	0.68
3:A:144:HEM:NB	4:A:145:CMO:C	2.57	0.68
2:B:57:ASN:ND2	2:B:59:LYS:HD2	2.07	0.68
1:A:41:LYS:CD	1:A:50:MET:CE	2.72	0.68
3:B:147:HEM:NA	4:B:148:CMO:C	2.57	0.67
1:A:41:LYS:CD	1:A:50:MET:HE1	2.25	0.67
2:B:1:VAL:CG1	2:B:2:GLU:H	2.07	0.67
3:D:147:HEM:NA	4:D:148:CMO:C	2.57	0.66
3:A:144:HEM:NA	4:A:145:CMO:C	2.60	0.65
1:A:18:ILE:HG13	1:A:115:PHE:CE2	2.34	0.63
1:C:32:ARG:NH1	2:D:127:GLN:HE21	1.92	0.62
1:A:8:LYS:HD3	1:A:75:ILE:HG23	1.81	0.62
2:B:1:VAL:HG12	2:B:2:GLU:H	1.64	0.62
1:C:108:LEU:HD23	1:C:127:LEU:HD23	1.81	0.61
2:D:57:ASN:HD22	2:D:60:ILE:H	1.48	0.61
1:A:41:LYS:CB	1:A:50:MET:HE1	2.31	0.60
1:C:44:PHE:HB2	1:C:50:MET:CE	2.30	0.60
1:C:125:VAL:HG22	2:D:34:VAL:HA	1.82	0.59
3:D:147:HEM:NC	4:D:148:CMO:C	2.66	0.59
3:D:147:HEM:NB	4:D:148:CMO:C	2.66	0.58
2:D:108:ASP:O	2:D:112:VAL:HG13	2.04	0.57
2:B:37:TRP:HA	1:C:94:ARG:HG3	1.86	0.57
2:B:19:ASN:HD22	2:B:19:ASN:C	2.07	0.57
1:C:23:ASP:OD1	1:C:62:LYS:HE2	2.05	0.56
1:A:23:ASP:OD1	1:A:62:LYS:HE3	2.05	0.56
2:B:28:LEU:HD13	2:B:60:ILE:HG23	1.86	0.56
1:C:120:THR:HG23	1:C:122:ASP:N	2.21	0.55
1:A:15:TRP:O	1:A:19:SER:HB3	2.06	0.54
1:A:32:ARG:HD3	2:B:127:GLN:HE22	1.71	0.54
1:A:116:PRO:HD2	1:A:117:LYS:HE3	1.89	0.54
1:A:39:GLN:NE2	1:A:39:GLN:H	2.02	0.54
1:C:1:THR:HG21	1:C:136:LEU:HD22	1.89	0.54
3:C:144:HEM:CMB	3:C:144:HEM:HBB2	2.37	0.53
1:A:32:ARG:NH1	2:B:127:GLN:HE21	2.03	0.53
3:B:147:HEM:NB	4:B:148:CMO:C	2.71	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:92:HIS:HE1	3:D:147:HEM:NA	2.05	0.53
1:A:117:LYS:HD2	1:A:117:LYS:N	2.17	0.52
1:C:120:THR:HG23	1:C:122:ASP:H	1.75	0.52
1:C:108:LEU:CD2	1:C:127:LEU:HD23	2.40	0.52
2:D:74:ALA:HB2	2:D:85:TYR:CE2	2.45	0.51
2:B:17:ASN:ND2	2:B:118:LEU:HD21	2.26	0.50
2:D:68:LEU:HD13	2:D:71:LEU:HD12	1.94	0.50
1:A:108:LEU:HD21	1:A:131:LEU:HD12	1.94	0.50
2:B:22:ASP:OD1	2:B:26:LYS:HE2	2.12	0.49
1:A:116:PRO:N	1:A:117:LYS:HD2	2.27	0.49
2:D:72:ASP:OD2	2:D:76:LYS:NZ	2.43	0.49
2:D:6:GLN:O	2:D:10:ILE:HG13	2.12	0.49
1:A:120:THR:HB	1:A:121:PRO:HD2	1.95	0.48
1:A:2:THR:HB	1:C:140:GLU:OE1	2.14	0.48
2:D:19:ASN:HD21	2:D:21:GLU:HG3	1.77	0.48
1:A:17:LYS:HD3	1:A:118:GLU:HG2	1.95	0.47
2:B:1:VAL:CG1	2:B:2:GLU:N	2.74	0.47
3:C:144:HEM:HMB1	3:C:144:HEM:HBB2	1.96	0.47
1:A:55:GLY:N	1:A:56:PRO:CD	2.78	0.47
1:C:55:GLY:N	1:C:56:PRO:CD	2.78	0.47
1:C:26:GLY:CA	1:C:65:MET:HG2	2.44	0.47
2:D:24:GLY:CA	2:D:68:LEU:HD23	2.45	0.46
1:C:113:LYS:HE2	2:D:119:GLY:O	2.16	0.46
2:B:88:LEU:HD11	3:B:147:HEM:HBA1	1.98	0.46
1:A:26:GLY:CA	1:A:65:MET:HG2	2.45	0.46
1:A:117:LYS:CD	1:A:117:LYS:N	2.66	0.46
1:C:54:SER:OG	1:C:57:VAL:HG13	2.14	0.46
1:A:34:LEU:O	1:A:41:LYS:HE2	2.16	0.46
1:C:4:SER:O	1:C:8:LYS:HG3	2.16	0.46
2:B:102:ASN:ND2	1:C:96:ASP:OD2	2.26	0.46
1:A:70:LEU:O	1:A:73:SER:HB2	2.17	0.45
2:D:81:ILE:O	2:D:85:TYR:HB2	2.16	0.45
2:B:93:SER:OG	2:B:142:GLY:HA2	2.17	0.45
1:C:89:HIS:HA	1:C:93:MET:HE2	1.99	0.44
1:C:118:GLU:N	1:C:118:GLU:OE1	2.50	0.44
1:C:32:ARG:HD3	2:D:127:GLN:HE22	1.83	0.44
1:C:26:GLY:HA3	1:C:65:MET:HG2	1.99	0.44
1:C:33:MET:SD	1:C:103:LEU:HB2	2.58	0.44
1:C:38:PRO:O	1:C:41:LYS:HG3	2.17	0.44
1:A:23:ASP:OD1	1:A:62:LYS:CE	2.65	0.44
1:A:94:ARG:HG3	2:D:37:TRP:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:VAL:HG13	2:B:34:VAL:HA	2.00	0.43
1:C:20:LYS:HG2	1:C:20:LYS:O	2.18	0.43
2:B:71:LEU:HD23	2:B:71:LEU:HA	1.90	0.43
2:D:6:GLN:HE21	2:D:6:GLN:HA	1.85	0.42
1:A:143:ARG:HG3	1:C:129:LYS:HD2	2.02	0.42
1:C:68:VAL:O	1:C:72:VAL:HG23	2.20	0.42
2:D:28:LEU:HD13	2:D:60:ILE:HG23	2.01	0.42
2:D:68:LEU:HD13	2:D:68:LEU:HA	1.85	0.42
2:D:57:ASN:HB3	2:D:60:ILE:HB	2.01	0.41
1:C:96:ASP:HA	1:C:97:PRO:HD3	1.85	0.41
2:B:63:HIS:HA	2:B:66:LYS:HG3	2.03	0.41
2:B:57:ASN:C	2:B:57:ASN:HD22	2.24	0.41
1:C:120:THR:HG22	1:C:123:ALA:CB	2.50	0.41
1:A:41:LYS:CG	1:A:50:MET:CE	2.89	0.41
2:D:41:TYR:CD1	3:D:147:HEM:HBC1	2.56	0.40
1:A:43:TYR:CE2	1:A:95:VAL:HA	2.56	0.40
2:B:60:ILE:HD13	2:B:60:ILE:HA	1.95	0.40
1:C:26:GLY:O	1:C:30:LEU:HB2	2.21	0.40
1:C:71:ALA:HB3	1:C:130:PHE:HZ	1.86	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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	142/144 (99%)	141 (99%)	1 (1%)	0	100	100
1	C	142/144 (99%)	135 (95%)	7 (5%)	0	100	100
2	B	140/146 (96%)	136 (97%)	4 (3%)	0	100	100
2	D	140/146 (96%)	136 (97%)	4 (3%)	0	100	100
All	All	564/580 (97%)	548 (97%)	16 (3%)	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	117/117 (100%)	104 (89%)	13 (11%)	8	4
1	C	117/117 (100%)	99 (85%)	18 (15%)	3	1
2	B	115/119 (97%)	98 (85%)	17 (15%)	4	1
2	D	115/119 (97%)	96 (84%)	19 (16%)	3	1
All	All	464/472 (98%)	397 (86%)	67 (14%)	4	2

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	ILE
1	A	20	LYS
1	A	30	LEU
1	A	39	GLN
1	A	62	LYS
1	A	79	THR
1	A	80	THR
1	A	94	ARG
1	A	117	LYS
1	A	122	ASP
1	A	125	VAL
1	A	138	LEU
1	A	141	ARG
2	B	2	GLU
2	B	19	ASN
2	B	21	GLU
2	B	28	LEU
2	B	55	LYS
2	B	57	ASN
2	B	59	LYS
2	B	65	VAL
2	B	66	LYS

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Mol	Chain	Res	Type
2	B	73	ARG
2	B	87	GLU
2	B	88	LEU
2	B	96	LEU
2	B	101	ASP
2	B	104	ARG
2	B	105	ILE
2	B	137	VAL
1	C	1	THR
1	C	6	LYS
1	C	9	SER
1	C	12	LYS
1	C	20	LYS
1	C	30	LEU
1	C	39	GLN
1	C	41	LYS
1	C	57	VAL
1	C	76	ASP
1	C	80	THR
1	C	84	ASP
1	C	94	ARG
1	C	101	LYS
1	C	117	LYS
1	C	120	THR
1	C	138	LEU
1	C	141	ARG
2	D	6	GLN
2	D	18	LEU
2	D	19	ASN
2	D	21	GLU
2	D	28	LEU
2	D	57	ASN
2	D	59	LYS
2	D	66	LYS
2	D	68	LEU
2	D	80	ASN
2	D	87	GLU
2	D	88	LEU
2	D	93	SER
2	D	96	LEU
2	D	97	HIS
2	D	101	ASP

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Mol	Chain	Res	Type
2	D	112	VAL
2	D	125	GLU
2	D	137	VAL

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	39	GLN
2	B	17	ASN
2	B	19	ASN
2	B	39	GLN
2	B	57	ASN
2	B	127	GLN
2	B	131	GLN
1	C	39	GLN
2	D	6	GLN
2	D	17	ASN
2	D	19	ASN
2	D	39	GLN
2	D	57	ASN
2	D	69	HIS
2	D	77	ASN
2	D	82	ASN
2	D	127	GLN
2	D	131	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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	144	1,4	30,50,50	2.18	9 (30%)	24,82,82	3.00	12 (50%)
4	CMO	A	145	3	0,1,1	0.00	-	0,0,0	0.00	-
3	HEM	B	147	2,4	30,50,50	2.00	8 (26%)	24,82,82	2.59	10 (41%)
4	CMO	B	148	3	0,1,1	0.00	-	0,0,0	0.00	-
3	HEM	C	144	1,4	30,50,50	2.40	10 (33%)	24,82,82	3.09	13 (54%)
4	CMO	C	145	3	0,1,1	0.00	-	0,0,0	0.00	-
3	HEM	D	147	2,4	30,50,50	2.46	11 (36%)	24,82,82	2.65	9 (37%)
4	CMO	D	148	3	0,1,1	0.00	-	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	144	1,4	-	0/10/54/54	0/0/8/8
4	CMO	A	145	3	-	0/0/0/0	0/0/0/0
3	HEM	B	147	2,4	-	0/10/54/54	0/0/8/8
4	CMO	B	148	3	-	0/0/0/0	0/0/0/0
3	HEM	C	144	1,4	-	0/10/54/54	0/0/8/8
4	CMO	C	145	3	-	0/0/0/0	0/0/0/0
3	HEM	D	147	2,4	-	0/10/54/54	0/0/8/8
4	CMO	D	148	3	-	0/0/0/0	0/0/0/0

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	144	HEM	C3B-C4B	-8.81	1.44	1.51
3	D	147	HEM	C3B-C4B	-8.68	1.44	1.51
3	B	147	HEM	C3B-C4B	-6.17	1.46	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	144	HEM	C3D-C4D	-6.12	1.43	1.51
3	C	144	HEM	C3D-C4D	-5.39	1.44	1.51
3	A	144	HEM	C3B-C4B	-5.22	1.47	1.51
3	D	147	HEM	C3D-C4D	-4.98	1.45	1.51
3	D	147	HEM	C2C-C1C	-3.19	1.46	1.52
3	B	147	HEM	C3D-C4D	-2.95	1.47	1.51
3	A	144	HEM	FE-ND	-2.70	1.83	1.97
3	B	147	HEM	C2C-C1C	-2.65	1.47	1.52
3	A	144	HEM	C2C-C1C	-2.54	1.47	1.52
3	B	147	HEM	C2B-C1B	-2.44	1.43	1.51
3	A	144	HEM	C2D-C1D	-2.42	1.43	1.51
3	C	144	HEM	C2D-C1D	-2.36	1.44	1.51
3	C	144	HEM	C2B-C1B	-2.28	1.44	1.51
3	A	144	HEM	C1A-CHA	-2.22	1.33	1.39
3	C	144	HEM	FE-NB	-2.21	1.85	1.97
3	D	147	HEM	C2B-C1B	-2.13	1.44	1.51
3	A	144	HEM	C2B-C1B	-2.11	1.45	1.51
3	D	147	HEM	C2D-C1D	-2.05	1.45	1.51
3	C	144	HEM	CMA-C3A	2.11	1.56	1.51
3	C	144	HEM	C3C-CAC	2.12	1.55	1.51
3	D	147	HEM	CMA-C3A	2.18	1.56	1.51
3	C	144	HEM	C3B-CAB	2.21	1.55	1.51
3	B	147	HEM	CMA-C3A	2.43	1.56	1.51
3	A	144	HEM	C3B-CAB	2.48	1.56	1.51
3	C	144	HEM	CMB-C2B	2.54	1.59	1.53
3	D	147	HEM	C1C-NC	2.56	1.39	1.36
3	D	147	HEM	FE-NB	2.64	2.11	1.97
3	D	147	HEM	C3B-CAB	2.71	1.56	1.51
3	D	147	HEM	CAA-C2A	2.99	1.57	1.52
3	B	147	HEM	CAA-C2A	3.07	1.57	1.52
3	B	147	HEM	FE-NC	3.10	2.08	1.95
3	C	144	HEM	FE-NC	3.18	2.08	1.95
3	B	147	HEM	C4C-NC	3.24	1.40	1.36
3	D	147	HEM	FE-NC	3.41	2.09	1.95
3	A	144	HEM	CAA-C2A	4.34	1.59	1.52

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	144	HEM	CAA-C2A-C1A	-6.66	119.78	127.01
3	A	144	HEM	C3C-CAC-CBC	-6.33	114.74	124.46
3	A	144	HEM	CAA-C2A-C1A	-4.77	121.83	127.01

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	144	HEM	CMA-C3A-C4A	-4.77	120.48	128.36
3	B	147	HEM	C3C-CAC-CBC	-4.29	117.88	124.46
3	C	144	HEM	C1D-CHD-C4C	-4.01	119.11	125.82
3	D	147	HEM	C3B-CAB-CBB	-3.97	118.36	124.46
3	A	144	HEM	CMA-C3A-C4A	-3.97	121.79	128.36
3	C	144	HEM	C3B-CAB-CBB	-3.38	119.28	124.46
3	A	144	HEM	C3B-CAB-CBB	-2.63	120.42	124.46
3	A	144	HEM	C2C-C1C-NC	-2.50	105.99	110.21
3	D	147	HEM	C4B-CHC-C1C	-2.46	121.71	125.82
3	B	147	HEM	C3B-CAB-CBB	-2.34	120.86	124.46
3	B	147	HEM	CBA-CAA-C2A	-2.05	108.86	112.53
3	A	144	HEM	C4B-CHC-C1C	-2.03	122.43	125.82
3	B	147	HEM	C2C-C1C-NC	-2.02	106.80	110.21
3	B	147	HEM	CBD-CAD-C3D	2.08	119.60	113.55
3	A	144	HEM	CMD-C2D-C3D	2.26	124.35	114.35
3	C	144	HEM	C4B-CHC-C1C	2.46	129.94	125.82
3	D	147	HEM	C2D-C3D-C4D	2.47	105.68	101.50
3	C	144	HEM	C2C-C1C-CHC	2.47	127.44	123.68
3	D	147	HEM	CMD-C2D-C3D	2.81	126.77	114.35
3	B	147	HEM	CMD-C2D-C3D	2.87	127.06	114.35
3	C	144	HEM	CMC-C2C-C3C	3.39	125.00	116.53
3	A	144	HEM	C2C-C1C-CHC	3.42	128.89	123.68
3	C	144	HEM	CMD-C2D-C3D	3.44	129.57	114.35
3	C	144	HEM	CMA-C3A-C2A	3.45	132.44	125.24
3	B	147	HEM	CAD-C3D-C4D	3.55	125.00	112.47
3	B	147	HEM	CMC-C2C-C3C	3.60	125.51	116.53
3	C	144	HEM	C2D-C3D-C4D	3.95	108.20	101.50
3	D	147	HEM	CAA-CBA-CGA	3.99	120.05	112.75
3	C	144	HEM	CAD-C3D-C2D	3.99	124.68	113.22
3	D	147	HEM	CAD-C3D-C4D	4.13	127.04	112.47
3	C	144	HEM	CAD-C3D-C4D	4.16	127.12	112.47
3	A	144	HEM	CMC-C2C-C3C	4.23	127.09	116.53
3	A	144	HEM	CAD-C3D-C2D	4.50	126.15	113.22
3	D	147	HEM	CMC-C2C-C3C	4.62	128.05	116.53
3	D	147	HEM	CAD-C3D-C2D	4.82	127.07	113.22
3	C	144	HEM	CMB-C2B-C3B	4.83	128.60	116.53
3	A	144	HEM	CAD-C3D-C4D	5.03	130.21	112.47
3	A	144	HEM	CMB-C2B-C3B	5.37	129.93	116.53
3	B	147	HEM	CMB-C2B-C3B	5.80	131.01	116.53
3	D	147	HEM	CMB-C2B-C3B	5.91	131.29	116.53
3	B	147	HEM	CAD-C3D-C2D	6.55	132.04	113.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	144	HEM	4	0
4	A	145	CMO	4	0
3	B	147	HEM	5	0
4	B	148	CMO	4	0
3	C	144	HEM	2	0
3	D	147	HEM	6	0
4	D	148	CMO	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.