



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 05:02 pm GMT

PDB ID : 1CBM
Title : THE 1.8 ANGSTROM STRUCTURE OF CARBONMONOXY-BETA4
HEMOGLOBIN: ANALYSIS OF A HOMOTETRAMER WITH THE R QUATERNARY
STRUCTURE OF LIGANDED ALPHA2BETA2 HEMOGLOBIN
Authors : Borgstahl, G.E.O.; Arnone, A.
Deposited on : 1993-02-18
Resolution : 1.74 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 : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

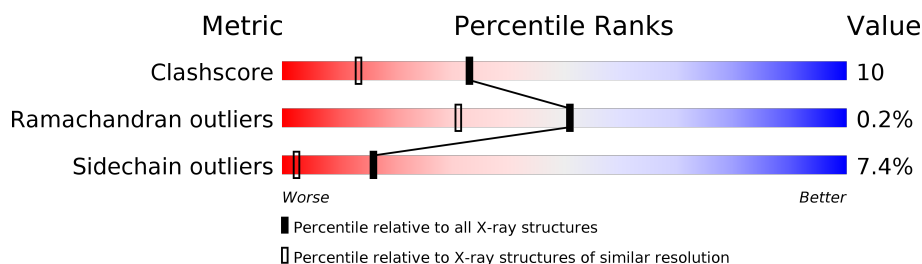
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.74 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	112137	2854 (1.76-1.72)
Ramachandran outliers	110173	2824 (1.76-1.72)
Sidechain outliers	110143	2824 (1.76-1.72)

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	146	
1	B	146	
1	C	146	
1	D	146	

2 Entry composition [i](#)

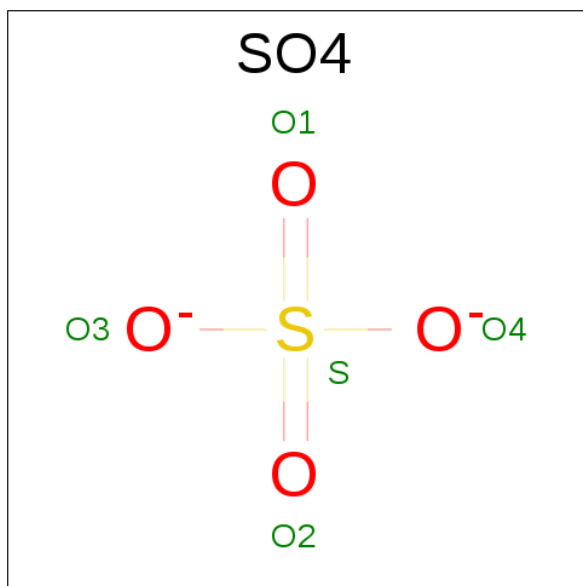
There are 5 unique types of molecules in this entry. The entry contains 4953 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BETA 4 (CARBONMONOXY).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	6	0
			1144	739	198	203	4			
1	B	146	Total	C	N	O	S	0	11	0
			1162	750	199	209	4			
1	C	146	Total	C	N	O	S	0	8	0
			1153	744	200	205	4			
1	D	146	Total	C	N	O	S	0	10	0
			1161	748	200	209	4			

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



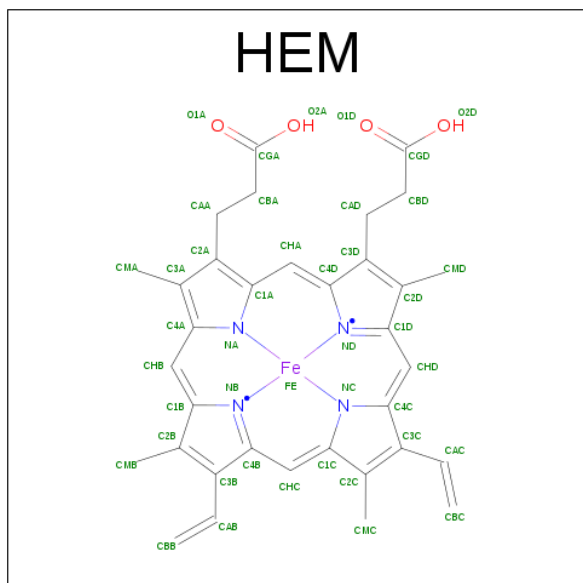
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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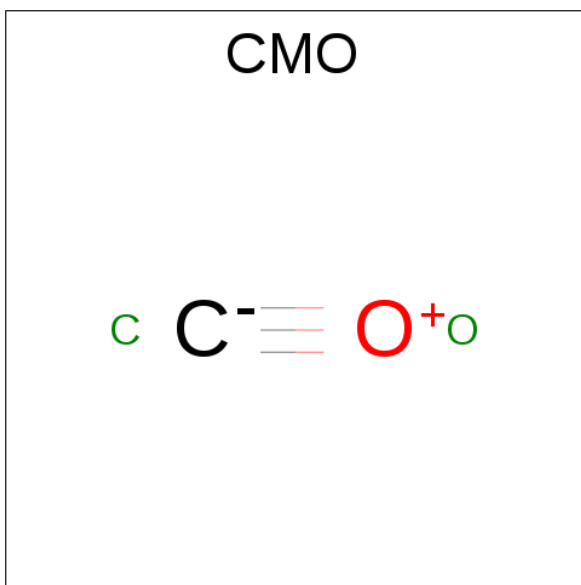
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	D	1	Total	O	S	0	0
			5	4	1		
2	C	1	Total	O	S	0	0
			5	4	1		

- 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 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- 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.

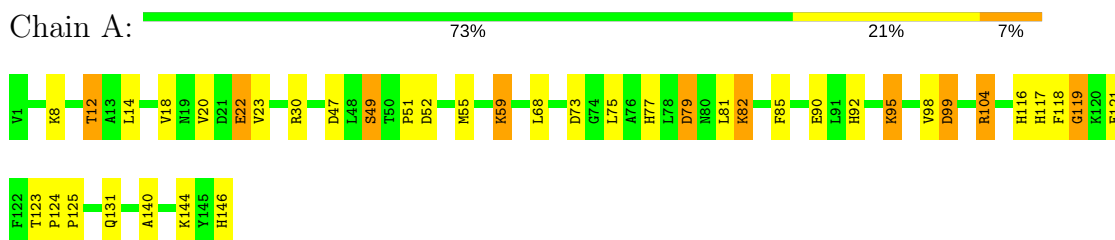
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	31	Total	O	0	0
			31	31		
5	B	34	Total	O	0	0
			34	34		
5	C	46	Total	O	0	0
			46	46		
5	D	22	Total	O	0	0
			22	22		

3 Residue-property plots [i](#)

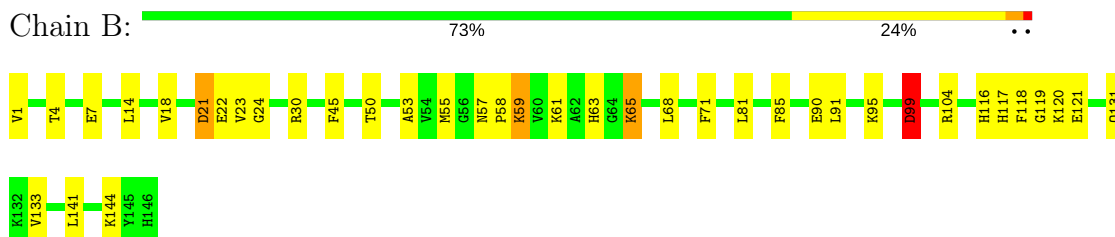
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

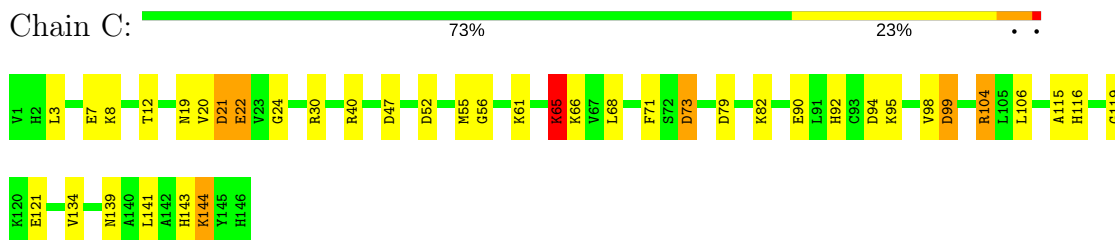
• Molecule 1: HEMOGLOBIN BETA 4 (CARBONMONOXY)



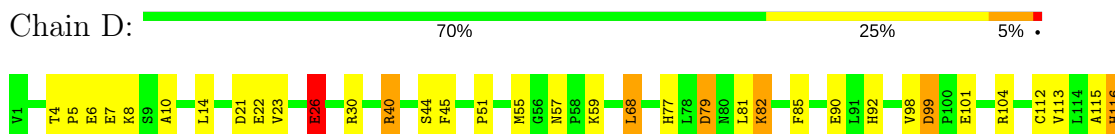
• Molecule 1: HEMOGLOBIN BETA 4 (CARBONMONOXY)



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H117	F118	G119	K120	E121	Y130	A140	L141	A142	H143	K144	Y145	H146
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4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.30 Å 82.40 Å 53.70 Å 90.00° 90.10° 90.00°	Depositor
Resolution (Å)	10.00 – 1.74	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.74)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, TNT	Depositor
R, R_{free}	0.177 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4953	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CMO, HEM, SO4

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	0.88	0/1205	1.47	13/1637 (0.8%)
1	B	0.89	0/1249	1.42	10/1695 (0.6%)
1	C	0.93	0/1225	1.64	20/1664 (1.2%)
1	D	0.92	2/1243 (0.2%)	1.52	15/1688 (0.9%)
All	All	0.91	2/4922 (0.0%)	1.52	58/6684 (0.9%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	112[A]	CYS	CB-SG	-5.22	1.73	1.81
1	D	112[B]	CYS	CB-SG	-5.22	1.73	1.81

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	30	ARG	NE-CZ-NH1	18.99	129.80	120.30
1	C	30	ARG	NE-CZ-NH2	-12.19	114.21	120.30
1	A	30	ARG	NE-CZ-NH1	11.29	125.95	120.30
1	D	104	ARG	NE-CZ-NH1	11.11	125.85	120.30
1	C	104	ARG	NE-CZ-NH1	10.84	125.72	120.30
1	A	99[A]	ASP	CB-CG-OD1	9.90	127.21	118.30
1	A	99[B]	ASP	CB-CG-OD1	9.90	127.21	118.30
1	B	104	ARG	NE-CZ-NH1	9.84	125.22	120.30
1	B	104	ARG	CD-NE-CZ	9.71	137.20	123.60
1	D	26	GLU	CA-CB-CG	9.09	133.39	113.40
1	C	22	GLU	CA-CB-CG	8.93	133.05	113.40
1	C	99[A]	ASP	CB-CG-OD2	-7.97	111.12	118.30
1	C	99[B]	ASP	CB-CG-OD2	-7.97	111.12	118.30
1	D	99[A]	ASP	CB-CG-OD1	7.91	125.42	118.30
1	D	99[B]	ASP	CB-CG-OD1	7.91	125.42	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99[A]	ASP	CB-CG-OD1	7.90	125.41	118.30
1	B	99[B]	ASP	CB-CG-OD1	7.90	125.41	118.30
1	C	121	GLU	CA-CB-CG	7.86	130.68	113.40
1	C	21	ASP	CB-CG-OD1	7.64	125.17	118.30
1	A	30	ARG	NE-CZ-NH2	-7.58	116.51	120.30
1	D	30	ARG	NE-CZ-NH1	7.29	123.95	120.30
1	A	104	ARG	NE-CZ-NH1	7.08	123.84	120.30
1	D	118	PHE	CA-CB-CG	7.07	130.87	113.90
1	C	99[A]	ASP	CB-CG-OD1	7.01	124.61	118.30
1	C	99[B]	ASP	CB-CG-OD1	7.01	124.61	118.30
1	D	99[A]	ASP	CB-CG-OD2	-6.97	112.03	118.30
1	D	99[B]	ASP	CB-CG-OD2	-6.97	112.03	118.30
1	C	116	HIS	CA-CB-CG	6.81	125.17	113.60
1	B	99[A]	ASP	CB-CG-OD2	-6.76	112.21	118.30
1	B	99[B]	ASP	CB-CG-OD2	-6.76	112.21	118.30
1	B	21	ASP	CB-CG-OD1	6.69	124.32	118.30
1	D	40[A]	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	D	40[B]	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	C	73	ASP	CB-CG-OD2	-6.17	112.74	118.30
1	D	130	TYR	CB-CG-CD2	-6.12	117.33	121.00
1	C	65	LYS	CB-CA-C	6.05	122.50	110.40
1	C	52	ASP	O-C-N	5.91	132.16	122.70
1	B	104	ARG	NH1-CZ-NH2	-5.80	113.02	119.40
1	B	30	ARG	NE-CZ-NH1	5.77	123.19	120.30
1	D	118	PHE	N-CA-CB	5.72	120.90	110.60
1	C	40[A]	ARG	NE-CZ-NH2	-5.65	117.47	120.30
1	C	40[B]	ARG	NE-CZ-NH2	-5.65	117.47	120.30
1	C	52	ASP	N-CA-CB	5.63	120.74	110.60
1	D	116	HIS	CA-CB-CG	5.57	123.06	113.60
1	D	101	GLU	CG-CD-OE2	-5.46	107.37	118.30
1	A	59	LYS	CA-CB-CG	5.45	125.39	113.40
1	A	121	GLU	CA-CB-CG	5.43	125.35	113.40
1	A	59	LYS	N-CA-CB	5.43	120.37	110.60
1	A	75[A]	LEU	CA-CB-CG	5.40	127.73	115.30
1	A	75[B]	LEU	CA-CB-CG	5.40	127.73	115.30
1	B	21	ASP	CB-CG-OD2	-5.33	113.51	118.30
1	C	65	LYS	CA-CB-CG	5.31	125.08	113.40
1	A	52	ASP	CB-CG-OD2	-5.24	113.59	118.30
1	A	79	ASP	CB-CG-OD2	5.17	122.95	118.30
1	D	5	PRO	O-C-N	5.10	130.86	122.70
1	A	18	VAL	CA-CB-CG1	5.09	118.53	110.90
1	C	47	ASP	CB-CG-OD2	-5.05	113.76	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	52	ASP	CA-C-O	-5.04	109.53	120.10

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	1144	0	1145	19	0
1	B	1162	0	1160	27	0
1	C	1153	0	1150	22	0
1	D	1161	0	1151	28	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
2	C	5	0	0	0	0
2	D	5	0	0	0	0
3	A	43	0	30	0	0
3	B	43	0	30	4	0
3	C	43	0	30	1	0
3	D	43	0	30	0	0
4	A	2	0	0	0	0
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
5	A	31	0	0	1	0
5	B	34	0	0	2	0
5	C	46	0	0	0	0
5	D	22	0	0	0	0
All	All	4953	0	4726	94	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 10.

All (94) 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:B:90:GLU:HG3	1:B:144:LYS:HE2	1.63	0.81
1:A:82:LYS:HD2	1:A:140:ALA:HA	1.71	0.72
1:D:23:VAL:HG12	1:D:68:LEU:HD11	1.72	0.72
1:B:55:MET:HE3	5:B:214:HOH:O	1.92	0.69
1:D:14:LEU:HD21	1:D:121[B]:GLU:HG3	1.75	0.68
1:B:58:PRO:HA	1:B:61:LYS:HD2	1.76	0.66
1:A:8:LYS:O	1:A:12:THR:HG23	1.95	0.66
1:B:24:GLY:CA	1:B:68[B]:LEU:HG	2.25	0.65
1:B:91:LEU:CD1	1:B:95:LYS:HE2	2.26	0.65
1:C:92:HIS:HB3	1:C:98[B]:VAL:CG1	2.27	0.64
1:B:24:GLY:HA2	1:B:68[B]:LEU:HG	1.83	0.60
1:C:71[B]:PHE:CZ	1:C:134:VAL:HG12	2.37	0.60
1:B:45:PHE:HA	1:B:59:LYS:HE2	1.83	0.59
1:C:20:VAL:HA	1:C:68[A]:LEU:CD2	2.32	0.59
3:B:147:HEM:HBB2	3:B:147:HEM:HHC	1.85	0.59
1:D:82:LYS:HE2	1:D:143:HIS:HB2	1.86	0.58
1:C:104:ARG:HH22	1:C:139:ASN:HD22	1.51	0.58
1:A:92:HIS:O	1:A:98[B]:VAL:HG12	2.04	0.58
1:C:82:LYS:HD2	1:C:143:HIS:ND1	2.20	0.57
1:D:116:HIS:ND1	1:D:117[A]:HIS:CD2	2.73	0.56
1:B:23:VAL:HG22	1:B:117:HIS:CE1	2.41	0.55
1:D:82:LYS:HD2	1:D:140:ALA:HA	1.87	0.55
1:A:47:ASP:OD1	1:A:49:SER:OG	2.25	0.54
1:D:57:ASN:OD1	1:D:59:LYS:HB2	2.06	0.54
1:B:116:HIS:HB2	1:C:115:ALA:HB1	1.89	0.54
1:C:8:LYS:O	1:C:12:THR:HG23	2.07	0.54
1:D:4:THR:OG1	1:D:6[A]:GLU:HG2	2.06	0.54
1:C:92:HIS:O	1:C:98[B]:VAL:HG12	2.09	0.53
1:B:4:THR:OG1	1:B:7:GLU:HG3	2.10	0.52
1:D:82:LYS:HE3	1:D:140:ALA:HA	1.92	0.52
1:D:92:HIS:O	1:D:98[B]:VAL:HG12	2.10	0.51
1:B:57:ASN:OD1	1:B:59:LYS:HB2	2.10	0.51
1:A:92:HIS:HB3	1:A:98[B]:VAL:CG1	2.40	0.51
1:C:24:GLY:CA	1:C:68[B]:LEU:HG	2.41	0.51
1:B:18:VAL:HG13	1:B:23:VAL:HG21	1.93	0.50
1:D:7:GLU:O	1:D:10:ALA:HB3	2.11	0.50
1:C:20:VAL:HA	1:C:68[A]:LEU:HD21	1.93	0.50
1:D:81:LEU:HD22	1:D:85:PHE:HE2	1.75	0.50
1:C:56:GLY:HA2	1:C:61:LYS:CE	2.42	0.49
1:A:81:LEU:HD22	1:A:85:PHE:HE2	1.77	0.49
1:B:116:HIS:CD2	1:C:119:GLY:O	2.66	0.49
1:B:50:THR:H	1:B:53:ALA:HB3	1.77	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:LEU:HD12	3:B:147:HEM:CBB	2.43	0.48
1:C:144:LYS:HA	1:C:144:LYS:HD2	1.69	0.48
1:B:118:PHE:O	1:B:119:GLY:C	2.53	0.47
1:C:20:VAL:HA	1:C:68[A]:LEU:HD23	1.96	0.47
1:C:19:ASN:ND2	1:C:22:GLU:OE1	2.48	0.47
1:D:113:VAL:O	1:D:116:HIS:HB3	2.15	0.47
1:A:116:HIS:HB2	1:D:115:ALA:HB1	1.97	0.46
1:D:82:LYS:HE2	1:D:143:HIS:CG	2.51	0.46
1:B:14[B]:LEU:HD21	1:B:121[B]:GLU:CG	2.46	0.46
1:D:14:LEU:HD21	1:D:121[B]:GLU:CG	2.42	0.46
1:C:71[B]:PHE:HZ	1:C:134:VAL:HG12	1.81	0.46
1:D:22:GLU:O	1:D:26:GLU:HB2	2.16	0.46
1:A:123:THR:OG1	1:A:125:PRO:HD2	2.17	0.45
1:A:20:VAL:HA	1:A:68:LEU:CD2	2.47	0.45
1:D:23:VAL:CG1	1:D:68:LEU:HD11	2.44	0.45
1:D:82:LYS:CE	1:D:140:ALA:HA	2.47	0.45
1:A:51:PRO:O	1:A:55:MET:HG2	2.17	0.45
1:D:90:GLU:HG2	1:D:144:LYS:HE3	1.98	0.45
1:A:95:LYS:HE2	1:A:95:LYS:HB3	1.79	0.45
1:D:23:VAL:HG22	1:D:117[A]:HIS:ND1	2.32	0.45
1:A:22:GLU:HG2	1:A:23:VAL:N	2.33	0.44
1:D:144:LYS:HA	1:D:144:LYS:HD2	1.63	0.43
1:B:57:ASN:O	1:B:61:LYS:HD2	2.18	0.43
1:B:91:LEU:HD11	1:B:95:LYS:HE2	1.99	0.43
1:B:21:ASP:OD1	1:B:65:LYS:HE3	2.18	0.43
1:A:124:PRO:N	1:A:125:PRO:CD	2.82	0.43
1:D:92:HIS:HB3	1:D:98[B]:VAL:CG1	2.49	0.43
1:D:45:PHE:HA	1:D:59:LYS:HD3	2.01	0.42
1:C:92:HIS:HB3	1:C:98[B]:VAL:HG11	2.01	0.42
1:B:116:HIS:ND1	1:B:117:HIS:HD2	2.16	0.42
1:C:106:LEU:HD23	3:C:147:HEM:HBB2	2.02	0.42
3:B:147:HEM:CBB	3:B:147:HEM:HHC	2.49	0.42
1:A:73:ASP:O	1:A:77:HIS:ND1	2.52	0.42
1:C:3:LEU:HA	1:C:7:GLU:OE1	2.20	0.41
1:D:82:LYS:HE2	1:D:143:HIS:CB	2.50	0.41
1:B:81:LEU:HD22	1:B:85:PHE:HE2	1.85	0.41
1:A:116:HIS:ND1	1:A:117:HIS:HD2	2.18	0.41
1:D:77:HIS:C	1:D:79:ASP:H	2.23	0.41
1:A:118:PHE:O	1:A:119:GLY:C	2.58	0.41
1:A:131:GLN:HG3	5:A:276:HOH:O	2.20	0.41
1:B:63:HIS:CE1	3:B:147:HEM:HBD2	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:146:HIS:OXT	1:B:1:VAL:HA	2.20	0.41
1:B:71[B]:PHE:CZ	1:B:133:VAL:HG12	2.56	0.41
1:C:55:MET:O	1:C:61:LYS:HE2	2.21	0.41
1:C:90:GLU:O	1:C:94:ASP:HB2	2.20	0.41
1:A:90:GLU:HA	1:A:144:LYS:HG3	2.03	0.40
1:C:21:ASP:HA	1:C:65:LYS:HD3	2.04	0.40
1:D:82:LYS:HE2	1:D:143:HIS:ND1	2.36	0.40
1:D:51:PRO:O	1:D:55:MET:HG2	2.21	0.40
1:D:81:LEU:HA	1:D:81:LEU:HD23	1.97	0.40
1:B:131:GLN:NE2	5:B:255:HOH:O	2.54	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	150/146 (103%)	146 (97%)	3 (2%)	1 (1%)	25	9
1	B	155/146 (106%)	151 (97%)	4 (3%)	0	100	100
1	C	152/146 (104%)	148 (97%)	4 (3%)	0	100	100
1	D	154/146 (106%)	150 (97%)	4 (3%)	0	100	100
All	All	611/584 (105%)	595 (97%)	15 (2%)	1 (0%)	51	31

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	GLY

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	124/118 (105%)	113 (91%)	11 (9%)	11	1
1	B	129/118 (109%)	123 (95%)	6 (5%)	30	8
1	C	126/118 (107%)	117 (93%)	9 (7%)	17	3
1	D	128/118 (108%)	114 (89%)	14 (11%)	7	1
All	All	507/472 (107%)	467 (92%)	40 (8%)	16	2

All (40) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	THR
1	A	14	LEU
1	A	22	GLU
1	A	49	SER
1	A	59	LYS
1	A	79	ASP
1	A	82	LYS
1	A	95	LYS
1	A	99[A]	ASP
1	A	99[B]	ASP
1	A	104	ARG
1	B	22	GLU
1	B	59	LYS
1	B	65	LYS
1	B	99[A]	ASP
1	B	99[B]	ASP
1	B	120	LYS
1	C	65	LYS
1	C	66	LYS
1	C	73	ASP
1	C	79	ASP
1	C	95	LYS
1	C	99[A]	ASP
1	C	99[B]	ASP

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Mol	Chain	Res	Type
1	C	141	LEU
1	C	144	LYS
1	D	8	LYS
1	D	21	ASP
1	D	26	GLU
1	D	40[A]	ARG
1	D	40[B]	ARG
1	D	44	SER
1	D	68	LEU
1	D	79	ASP
1	D	82	LYS
1	D	99[A]	ASP
1	D	99[B]	ASP
1	D	120	LYS
1	D	141	LEU
1	D	144	LYS

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	117	HIS
1	B	131	GLN
1	C	131	GLN
1	C	139	ASN

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

12 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	147	1,4	28,50,50	1.94	9 (32%)	17,82,82	1.78	4 (23%)
4	CMO	A	148	3	0,1,1	0.00	-	0,0,0	0.00	-
2	SO4	A	150	-	4,4,4	0.74	0	6,6,6	0.23	0
3	HEM	B	147	1,4	28,50,50	2.10	7 (25%)	17,82,82	2.10	7 (41%)
4	CMO	B	148	3	0,1,1	0.00	-	0,0,0	0.00	-
2	SO4	B	151	-	4,4,4	0.76	0	6,6,6	0.32	0
3	HEM	C	147	1,4	28,50,50	2.00	8 (28%)	17,82,82	2.15	8 (47%)
4	CMO	C	148	3	0,1,1	0.00	-	0,0,0	0.00	-
2	SO4	C	153	-	4,4,4	0.73	0	6,6,6	0.10	0
3	HEM	D	147	1,4	28,50,50	2.14	9 (32%)	17,82,82	2.17	5 (29%)
4	CMO	D	148	3	0,1,1	0.00	-	0,0,0	0.00	-
2	SO4	D	152	-	4,4,4	0.73	0	6,6,6	0.08	0

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	147	1,4	-	0/6/54/54	0/0/8/8
4	CMO	A	148	3	-	0/0/0/0	0/0/0/0
2	SO4	A	150	-	-	0/0/0/0	0/0/0/0
3	HEM	B	147	1,4	-	0/6/54/54	0/0/8/8
4	CMO	B	148	3	-	0/0/0/0	0/0/0/0
2	SO4	B	151	-	-	0/0/0/0	0/0/0/0
3	HEM	C	147	1,4	-	0/6/54/54	0/0/8/8
4	CMO	C	148	3	-	0/0/0/0	0/0/0/0
2	SO4	C	153	-	-	0/0/0/0	0/0/0/0
3	HEM	D	147	1,4	-	0/6/54/54	0/0/8/8
4	CMO	D	148	3	-	0/0/0/0	0/0/0/0
2	SO4	D	152	-	-	0/0/0/0	0/0/0/0

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	147	HEM	C3C-C2C	-5.45	1.33	1.40
3	C	147	HEM	C3C-C2C	-5.07	1.33	1.40
3	B	147	HEM	C3B-C2B	-4.80	1.34	1.40
3	B	147	HEM	C3C-C2C	-4.35	1.34	1.40
3	A	147	HEM	C3C-C2C	-4.23	1.34	1.40
3	D	147	HEM	C3B-C2B	-3.99	1.35	1.40
3	C	147	HEM	C3B-C2B	-3.77	1.35	1.40
3	A	147	HEM	C3B-C2B	-3.26	1.36	1.40
3	B	147	HEM	C4B-NB	2.01	1.40	1.36
3	B	147	HEM	CAD-C3D	2.06	1.56	1.52
3	A	147	HEM	C4C-NC	2.11	1.39	1.36
3	C	147	HEM	C1C-NC	2.16	1.39	1.36
3	A	147	HEM	CAD-C3D	2.20	1.56	1.52
3	D	147	HEM	CAA-C2A	2.21	1.55	1.52
3	C	147	HEM	C1B-NB	2.25	1.39	1.36
3	A	147	HEM	CMB-C2B	2.28	1.56	1.51
3	D	147	HEM	C4B-NB	2.29	1.41	1.36
3	D	147	HEM	CAD-C3D	2.33	1.56	1.52
3	D	147	HEM	C1B-NB	2.41	1.39	1.36
3	C	147	HEM	CAA-C2A	2.44	1.56	1.52
3	C	147	HEM	CMB-C2B	2.87	1.57	1.51
3	C	147	HEM	C4C-NC	2.89	1.40	1.36
3	A	147	HEM	C3C-CAC	2.90	1.53	1.47
3	A	147	HEM	CAA-C2A	3.09	1.57	1.52
3	C	147	HEM	C3C-CAC	3.10	1.53	1.47
3	A	147	HEM	C4D-ND	3.22	1.40	1.36
3	D	147	HEM	C4D-ND	3.35	1.40	1.36
3	A	147	HEM	C3B-CAB	3.48	1.54	1.47
3	B	147	HEM	C3B-CAB	3.52	1.54	1.47
3	D	147	HEM	C3C-CAC	3.52	1.54	1.47
3	B	147	HEM	C4D-ND	4.02	1.41	1.36
3	B	147	HEM	C3C-CAC	4.03	1.55	1.47
3	D	147	HEM	C3B-CAB	4.21	1.56	1.47

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	147	HEM	CMA-C3A-C4A	-5.36	120.22	128.46
3	C	147	HEM	CMD-C2D-C1D	-4.41	121.68	128.46
3	B	147	HEM	CMA-C3A-C4A	-4.15	122.09	128.46
3	A	147	HEM	CMA-C3A-C4A	-3.55	123.01	128.46
3	D	147	HEM	CMD-C2D-C1D	-3.30	123.39	128.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	147	HEM	CMD-C2D-C1D	-3.29	123.41	128.46
3	B	147	HEM	CMD-C2D-C1D	-3.10	123.70	128.46
3	C	147	HEM	CAA-CBA-CGA	-2.86	107.77	112.66
3	C	147	HEM	CMA-C3A-C4A	-2.85	124.09	128.46
3	B	147	HEM	CAA-CBA-CGA	-2.57	108.27	112.66
3	B	147	HEM	CBA-CAA-C2A	-2.48	107.74	112.48
3	C	147	HEM	CBA-CAA-C2A	-2.21	108.27	112.48
3	C	147	HEM	CMB-C2B-C3B	2.09	128.76	124.89
3	A	147	HEM	CMD-C2D-C3D	2.15	129.00	124.94
3	D	147	HEM	CMD-C2D-C3D	2.20	129.09	124.94
3	B	147	HEM	CMC-C2C-C3C	2.40	129.35	124.89
3	C	147	HEM	CMA-C3A-C2A	2.46	129.59	124.94
3	B	147	HEM	CMD-C2D-C3D	2.47	129.59	124.94
3	D	147	HEM	CMC-C2C-C3C	2.49	129.50	124.89
3	C	147	HEM	CMC-C2C-C3C	2.61	129.73	124.89
3	A	147	HEM	CMA-C3A-C2A	3.04	130.67	124.94
3	C	147	HEM	CMD-C2D-C3D	3.67	131.87	124.94
3	B	147	HEM	CMA-C3A-C2A	3.74	131.99	124.94
3	D	147	HEM	CMA-C3A-C2A	4.06	132.59	124.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	147	HEM	4	0
3	C	147	HEM	1	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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.