



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 06:45 PM GMT

PDB ID : 1CBL
Title : THE 1.9 ANGSTROM STRUCTURE OF DEOXY-BETA4 HEMOGLOBIN:
ANALYSIS OF THE PARTITIONING OF QUATERNARY-ASSOCIATED
AND LIGAND-INDUCED CHANGES IN TERTIARY STRUCTURE
Authors : Borgstahl, G.E.O.; Arnone, A.
Deposited on : 1993-02-18
Resolution : 1.80 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

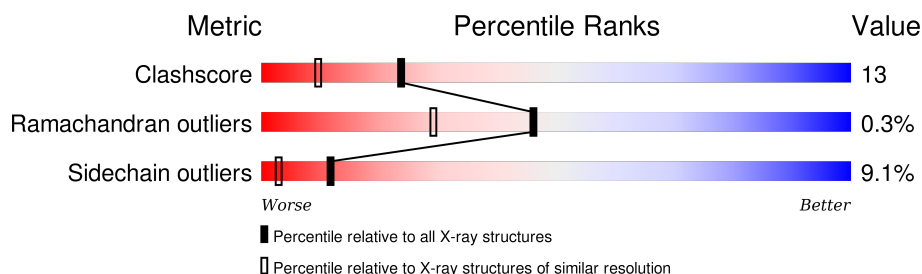
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)

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	 63% 31% 6%
1	B	146	 66% 29% • •
1	C	146	 74% 21% 5%
1	D	146	 62% 30% 7% •

2 Entry composition [i](#)

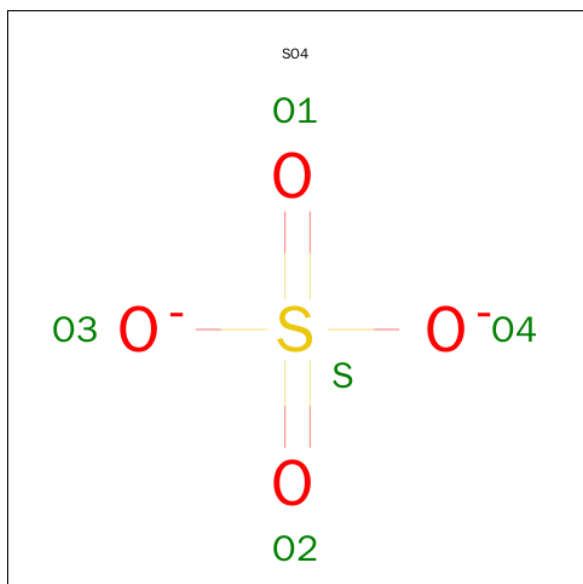
There are 4 unique types of molecules in this entry. The entry contains 4893 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 (DEOXY).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	146	Total	C	N	O	S	0	8	0
			1154	742	201	207	4			
1	B	146	Total	C	N	O	S	0	9	0
			1157	744	200	209	4			
1	C	146	Total	C	N	O	S	0	10	0
			1163	748	202	209	4			
1	D	146	Total	C	N	O	S	0	9	0
			1156	747	202	203	4			

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



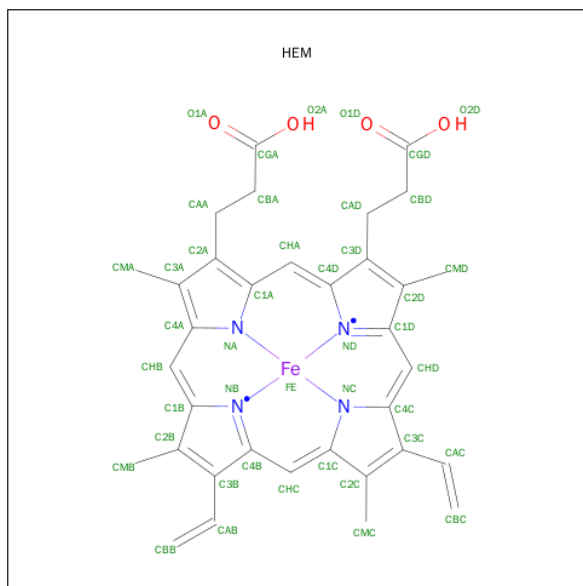
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

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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	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 water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	20	Total	O	0	0
			20	20		
4	B	12	Total	O	0	0
			12	12		
4	C	24	Total	O	0	0
			24	24		

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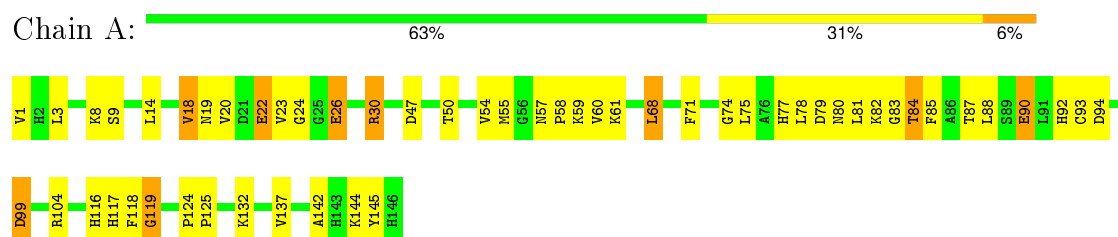
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	15	Total	O	0	0
			15	15		

3 Residue-property plots

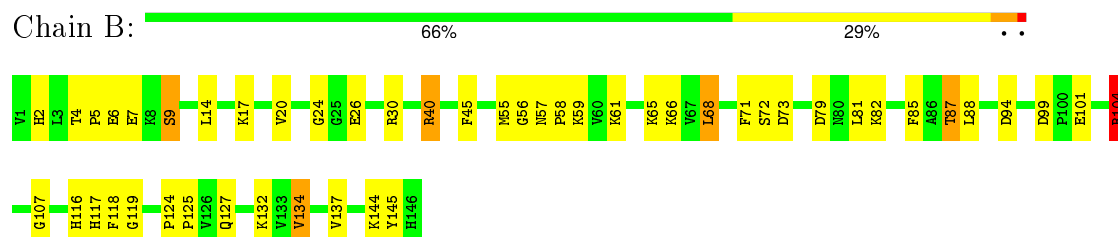
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

Note EDS was not executed.

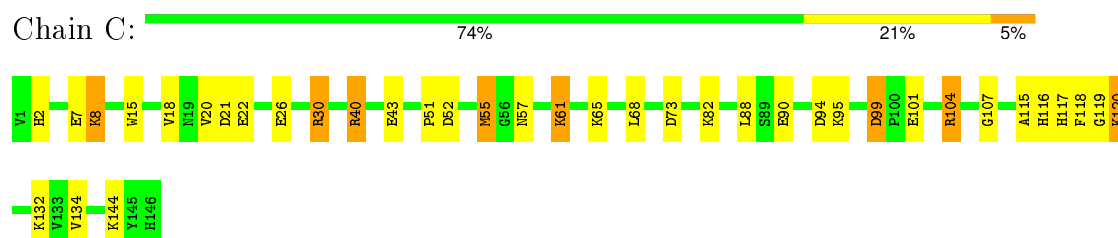
• Molecule 1: HEMOGLOBIN BETA 4 (DEOXY)



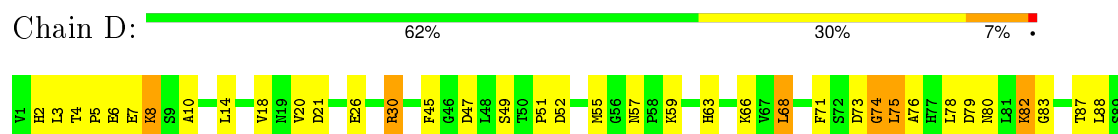
• Molecule 1: HEMOGLOBIN BETA 4 (DEOXY)



• Molecule 1: HEMOGLOBIN BETA 4 (DEOXY)



• Molecule 1: HEMOGLOBIN BETA 4 (DEOXY)





4 Data and refinement statistics

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

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.00 Å 81.80 Å 54.40 Å 90.00° 89.70° 90.00°	Depositor
Resolution (Å)	10.00 – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, TNT	Depositor
R, R_{free}	0.190 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4893	wwPDB-VP
Average B, all atoms (Å ²)	32.0	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: 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	0.82	0/1227	1.54	12/1665 (0.7%)
1	B	0.84	0/1234	1.55	20/1673 (1.2%)
1	C	0.87	0/1246	1.47	17/1689 (1.0%)
1	D	0.81	0/1235	1.50	8/1675 (0.5%)
All	All	0.84	0/4942	1.52	57/6702 (0.9%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

There are no bond length outliers.

All (57) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	30	ARG	NE-CZ-NH1	17.81	129.21	120.30
1	A	99[A]	ASP	CB-CG-OD1	12.21	129.29	118.30
1	A	99[B]	ASP	CB-CG-OD1	12.21	129.29	118.30
1	D	99[A]	ASP	CB-CG-OD1	9.84	127.16	118.30
1	D	99[B]	ASP	CB-CG-OD1	9.84	127.16	118.30
1	B	99[A]	ASP	CB-CG-OD1	9.00	126.40	118.30
1	B	99[B]	ASP	CB-CG-OD1	9.00	126.40	118.30
1	B	30	ARG	NE-CZ-NH1	7.83	124.21	120.30
1	B	104	ARG	NE-CZ-NH1	7.67	124.13	120.30
1	D	30	ARG	NE-CZ-NH1	7.60	124.10	120.30
1	C	22	GLU	CA-CB-CG	7.58	130.07	113.40
1	B	99[A]	ASP	OD1-CG-OD2	-7.47	109.11	123.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	99[B]	ASP	OD1-CG-OD2	-7.47	109.11	123.30
1	A	30	ARG	CD-NE-CZ	7.45	134.03	123.60
1	D	116	HIS	CA-CB-CG	6.94	125.40	113.60
1	B	99[A]	ASP	CB-CG-OD2	6.83	124.44	118.30
1	B	99[B]	ASP	CB-CG-OD2	6.83	124.44	118.30
1	A	30	ARG	NE-CZ-NH2	-6.82	116.89	120.30
1	C	104	ARG	NE-CZ-NH1	6.70	123.65	120.30
1	C	21	ASP	CB-CG-OD1	6.67	124.30	118.30
1	C	30	ARG	NE-CZ-NH1	6.57	123.59	120.30
1	B	30	ARG	NE-CZ-NH2	-6.36	117.12	120.30
1	C	104	ARG	NE-CZ-NH2	-6.25	117.17	120.30
1	C	55	MET	CG-SD-CE	6.12	109.99	100.20
1	C	99[A]	ASP	CB-CG-OD1	5.93	123.64	118.30
1	C	99[B]	ASP	CB-CG-OD1	5.93	123.64	118.30
1	B	55	MET	CG-SD-CE	5.83	109.52	100.20
1	B	73[A]	ASP	CB-CA-C	5.82	122.03	110.40
1	B	73[B]	ASP	CB-CA-C	5.82	122.03	110.40
1	C	73	ASP	CB-CA-C	5.79	121.98	110.40
1	D	47	ASP	CB-CG-OD2	5.79	123.51	118.30
1	C	73	ASP	CB-CG-OD2	-5.76	113.12	118.30
1	D	134	VAL	CB-CA-C	5.75	122.31	111.40
1	A	99[A]	ASP	OD1-CG-OD2	-5.73	112.42	123.30
1	A	99[B]	ASP	OD1-CG-OD2	-5.73	112.42	123.30
1	C	94	ASP	CB-CG-OD2	5.71	123.44	118.30
1	C	52	ASP	O-C-N	5.56	131.59	122.70
1	B	26	GLU	CA-CB-CG	5.50	125.50	113.40
1	C	99[A]	ASP	CB-CG-OD2	-5.50	113.35	118.30
1	C	99[B]	ASP	CB-CG-OD2	-5.50	113.35	118.30
1	A	104	ARG	NE-CZ-NH2	5.47	123.04	120.30
1	D	30	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	B	134	VAL	CB-CA-C	5.41	121.68	111.40
1	A	99[A]	ASP	CA-CB-CG	5.37	125.21	113.40
1	A	99[B]	ASP	CA-CB-CG	5.37	125.21	113.40
1	C	2	HIS	N-CA-CB	5.35	120.23	110.60
1	C	40[A]	ARG	NE-CZ-NH2	-5.31	117.65	120.30
1	C	40[B]	ARG	NE-CZ-NH2	-5.31	117.65	120.30
1	B	94	ASP	CB-CG-OD1	5.21	122.98	118.30
1	B	2	HIS	N-CA-CB	5.18	119.93	110.60
1	B	87	THR	CA-CB-CG2	5.18	119.65	112.40
1	B	145	TYR	CB-CG-CD1	5.11	124.06	121.00
1	B	40[A]	ARG	NE-CZ-NH1	5.08	122.84	120.30
1	B	40[B]	ARG	NE-CZ-NH1	5.08	122.84	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	134	VAL	CA-CB-CG1	5.04	118.47	110.90
1	A	84	THR	CA-CB-CG2	5.03	119.44	112.40
1	A	30	ARG	NH1-CZ-NH2	-5.01	113.89	119.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	104	ARG	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1154	0	1142	45	0
1	B	1157	0	1149	28	0
1	C	1163	0	1154	24	0
1	D	1156	0	1158	37	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	0	0
3	C	43	0	30	0	0
3	D	43	0	30	2	0
4	A	20	0	0	0	0
4	B	12	0	0	1	0
4	C	24	0	0	0	0
4	D	15	0	0	0	0
All	All	4893	0	4723	127	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 (127) 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:D:7:GLU:HG2	1:D:129:ALA:HB2	1.44	0.98
1:D:3:LEU:HD12	1:D:8:LYS:HG2	1.60	0.83
1:D:82:LYS:H	1:D:82:LYS:HD2	1.47	0.79
1:A:90:GLU:HA	1:A:144:LYS:HG3	1.71	0.71
1:B:58:PRO:HA	1:B:61:LYS:HD2	1.74	0.70
1:A:19:ASN:ND2	1:A:22:GLU:HB3	2.11	0.66
1:D:8:LYS:HE3	1:D:78:LEU:HD12	1.77	0.66
1:A:14:LEU:HD11	1:A:118:PHE:CG	2.30	0.66
1:D:143:HIS:HA	1:D:146:HIS:CE1	2.32	0.65
1:B:116:HIS:HD2	1:B:117:HIS:HD2	1.44	0.65
1:C:20:VAL:HB	1:C:68:LEU:HD23	1.79	0.65
1:A:20:VAL:HG13	1:A:68:LEU:HB3	1.79	0.65
1:A:71[A]:PHE:CE1	1:A:137:VAL:HG21	2.32	0.64
1:B:45:PHE:HA	1:B:59:LYS:HD3	1.80	0.64
1:A:90:GLU:HG3	1:A:144:LYS:HD3	1.81	0.63
1:B:59:LYS:HD2	4:B:239:HOH:O	2.01	0.61
1:D:4:THR:HB	1:D:5:PRO:HD2	1.83	0.61
1:A:24:GLY:HA2	1:A:68:LEU:HG	1.84	0.59
1:A:81:LEU:HD22	1:A:85:PHE:HE2	1.66	0.59
1:A:71[A]:PHE:HE1	1:A:137:VAL:HG21	1.68	0.58
1:B:57:ASN:O	1:B:61:LYS:HG3	2.02	0.58
1:A:116:HIS:ND1	1:A:117[A]:HIS:HD2	2.00	0.58
1:A:116:HIS:HB2	1:D:115:ALA:HB1	1.86	0.57
1:D:26:GLU:HG3	1:D:30:ARG:HE	1.69	0.57
1:A:19:ASN:HD22	1:A:22:GLU:HB3	1.69	0.56
1:A:90:GLU:O	1:A:94:ASP:HB2	2.05	0.56
1:A:145:TYR:O	1:B:132[B]:LYS:HD2	2.05	0.56
1:A:1:VAL:HG12	1:A:3:LEU:HG	1.87	0.56
1:D:51:PRO:O	1:D:55:MET:HG2	2.07	0.55
1:D:133:VAL:O	1:D:137:VAL:HG23	2.08	0.54
1:A:88:LEU:HD22	1:A:92:HIS:HE1	1.73	0.54
1:A:142:ALA:O	1:A:145:TYR:HB2	2.07	0.53
1:D:71[B]:PHE:CE1	1:D:137:VAL:HG21	2.44	0.53
1:A:18:VAL:HG23	1:A:118:PHE:CZ	2.44	0.52
1:B:116:HIS:HD2	1:B:117:HIS:CD2	2.26	0.52
1:B:58:PRO:HA	1:B:61:LYS:CD	2.39	0.52
1:B:14:LEU:HD21	1:B:118:PHE:CD2	2.44	0.52
1:D:45:PHE:HA	1:D:59:LYS:HD3	1.91	0.51
1:D:118:PHE:O	1:D:119:GLY:C	2.50	0.51
1:A:18:VAL:HG23	1:A:118:PHE:HZ	1.76	0.50
1:C:15:TRP:O	1:C:18:VAL:HB	2.12	0.50
1:D:3:LEU:HB3	1:D:7:GLU:HB3	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:CYS:SG	1:A:144:LYS:HB2	2.52	0.50
1:B:20:VAL:HA	1:B:68:LEU:HD12	1.94	0.50
1:A:116:HIS:ND1	1:A:117[A]:HIS:CD2	2.80	0.49
1:A:77:HIS:C	1:A:79:ASP:H	2.15	0.49
1:D:90:GLU:HA	1:D:144:LYS:HG2	1.94	0.49
1:C:101:GLU:OE1	1:C:104:ARG:NH2	2.44	0.49
1:A:50:THR:O	1:A:54:VAL:HG23	2.14	0.48
1:A:75:LEU:HD23	1:A:78:LEU:HD21	1.95	0.48
1:A:3:LEU:HB2	1:A:8:LYS:HG2	1.95	0.48
1:D:73:ASP:O	1:D:76:ALA:N	2.45	0.48
1:B:71[B]:PHE:CE1	1:B:137:VAL:HG21	2.49	0.48
1:C:107:GLY:HA3	1:C:134:VAL:HG13	1.96	0.48
1:A:26:GLU:O	1:A:30:ARG:HG3	2.14	0.48
1:C:120:LYS:NZ	1:C:120:LYS:HB2	2.29	0.47
1:A:74:GLY:HA2	1:A:84:THR:HG21	1.96	0.47
1:C:90:GLU:HA	1:C:144:LYS:HG3	1.95	0.47
1:B:4:THR:OG1	1:B:7:GLU:HG3	2.14	0.47
1:A:80:ASN:OD1	1:A:83:GLY:HA3	2.15	0.46
1:D:10:ALA:O	1:D:14[B]:LEU:HG	2.15	0.46
1:A:3:LEU:HD12	1:A:78:LEU:HD13	1.98	0.46
1:D:88:LEU:HD22	1:D:92:HIS:HE1	1.81	0.46
1:A:57:ASN:O	1:A:60:VAL:HB	2.15	0.46
1:C:26[A]:GLU:OE1	1:C:117[A]:HIS:NE2	2.49	0.45
1:D:80:ASN:OD1	1:D:83:GLY:HA3	2.16	0.45
1:A:55:MET:HA	1:A:55:MET:HE3	1.99	0.45
1:B:116:HIS:HB2	1:C:115:ALA:HB1	1.99	0.44
1:D:116:HIS:ND1	1:D:117[A]:HIS:HD2	2.15	0.44
1:D:116:HIS:ND1	1:D:117[A]:HIS:CD2	2.86	0.44
1:B:40[A]:ARG:NH2	1:D:99[A]:ASP:OD1	2.47	0.44
1:B:127:GLN:OE1	1:C:30:ARG:HD3	2.17	0.44
1:D:71[A]:PHE:CE1	1:D:75[A]:LEU:HD21	2.53	0.43
1:A:23:VAL:HG12	1:A:68:LEU:HD11	2.00	0.43
1:A:116:HIS:CD2	1:D:119:GLY:O	2.71	0.43
1:D:68:LEU:HA	1:D:68:LEU:HD22	1.84	0.43
1:B:144:LYS:HB3	1:B:144:LYS:HE2	1.80	0.43
1:A:118:PHE:O	1:A:119:GLY:C	2.57	0.43
1:B:124:PRO:N	1:B:125:PRO:CD	2.81	0.43
1:D:4:THR:O	1:D:5:PRO:C	2.56	0.43
1:D:5:PRO:HG2	1:D:6:GLU:H	1.83	0.43
1:D:73:ASP:O	1:D:74:GLY:C	2.57	0.43
1:B:101:GLU:OE1	1:B:104:ARG:NH1	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:5:PRO:O	1:B:9:SER:HB3	2.19	0.43
1:D:82:LYS:CD	1:D:82:LYS:H	2.16	0.42
1:C:20:VAL:CB	1:C:68:LEU:HD23	2.48	0.42
1:D:71[B]:PHE:CE2	3:D:147:HEM:HBB1	2.55	0.42
1:C:118:PHE:O	1:C:119:GLY:C	2.58	0.42
1:B:68:LEU:HA	1:B:71[B]:PHE:HB2	2.01	0.42
1:C:57:ASN:O	1:C:61:LYS:HG3	2.20	0.42
1:B:119:GLY:O	1:C:116:HIS:NE2	2.52	0.42
1:C:51:PRO:O	1:C:55:MET:HG2	2.20	0.42
1:B:24:GLY:HA2	1:B:68:LEU:HG	2.02	0.41
1:A:57:ASN:O	1:A:61:LYS:HG3	2.20	0.41
1:C:88:LEU:HD23	1:C:88:LEU:HA	1.85	0.41
1:A:88:LEU:HD22	1:A:92:HIS:CE1	2.55	0.41
1:A:3:LEU:CD1	1:A:78:LEU:HD13	2.51	0.41
1:A:77:HIS:C	1:A:79:ASP:N	2.74	0.41
1:A:124:PRO:HB2	1:A:125:PRO:HD3	2.02	0.41
1:B:81:LEU:HD22	1:B:85:PHE:HE2	1.85	0.41
1:C:8:LYS:HB2	1:C:8:LYS:HE3	1.84	0.41
1:B:107:GLY:HA3	1:B:134:VAL:HG13	2.03	0.41
1:D:63:HIS:HE1	3:D:147:HEM:CHA	2.34	0.41
1:D:57:ASN:OD1	1:D:59:LYS:HB2	2.21	0.41
1:B:56:GLY:O	1:B:61:LYS:HE3	2.21	0.40
1:D:88:LEU:HD22	1:D:92:HIS:CE1	2.55	0.40
1:B:88:LEU:HD23	1:B:88:LEU:HA	1.81	0.40
1:B:61:LYS:H	1:B:61:LYS:HG3	1.65	0.40
1:A:80:ASN:ND2	1:A:83:GLY:H	2.20	0.40
1:C:7:GLU:OE2	1:C:132[A]:LYS:NZ	2.49	0.40
1:A:57:ASN:HA	1:A:58:PRO:HD3	1.82	0.40
1:D:123:THR:OG1	1:D:126:VAL:HG23	2.22	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	152/146 (104%)	146 (96%)	5 (3%)	1 (1%)	26	11
1	B	153/146 (105%)	148 (97%)	5 (3%)	0	100	100
1	C	154/146 (106%)	152 (99%)	2 (1%)	0	100	100
1	D	153/146 (105%)	147 (96%)	5 (3%)	1 (1%)	26	11
All	All	612/584 (105%)	593 (97%)	17 (3%)	2 (0%)	46	29

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	119	GLY
1	D	74	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	126/118 (107%)	116 (92%)	10 (8%)	15	4
1	B	127/118 (108%)	115 (91%)	12 (9%)	11	3
1	C	128/118 (108%)	119 (93%)	9 (7%)	19	6
1	D	127/118 (108%)	109 (86%)	18 (14%)	4	1
All	All	508/472 (108%)	459 (90%)	49 (10%)	12	2

All (49) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	A	18	VAL
1	A	22	GLU
1	A	26	GLU
1	A	47	ASP
1	A	59	LYS
1	A	68	LEU
1	A	82	LYS
1	A	87	THR

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Mol	Chain	Res	Type
1	A	90	GLU
1	B	6[A]	GLU
1	B	6[B]	GLU
1	B	9	SER
1	B	17	LYS
1	B	65	LYS
1	B	66	LYS
1	B	68	LEU
1	B	72	SER
1	B	79	ASP
1	B	82[A]	LYS
1	B	82[B]	LYS
1	B	87	THR
1	C	8	LYS
1	C	40[A]	ARG
1	C	40[B]	ARG
1	C	61	LYS
1	C	65	LYS
1	C	95	LYS
1	C	99[A]	ASP
1	C	99[B]	ASP
1	C	120	LYS
1	D	2	HIS
1	D	8	LYS
1	D	20	VAL
1	D	21	ASP
1	D	49	SER
1	D	52	ASP
1	D	66	LYS
1	D	68	LEU
1	D	75[A]	LEU
1	D	75[B]	LEU
1	D	79	ASP
1	D	82	LYS
1	D	87	THR
1	D	99[A]	ASP
1	D	99[B]	ASP
1	D	134	VAL
1	D	144	LYS
1	D	146	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	ASN
1	B	80	ASN
1	B	116	HIS
1	B	117	HIS
1	B	131	GLN
1	C	80	ASN
1	C	131	GLN
1	D	146	HIS

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	147	1	30,50,50	2.75	8 (26%)	24,82,82	2.74	12 (50%)
2	SO4	A	150	-	4,4,4	0.85	0	6,6,6	0.15	0
3	HEM	B	147	1	30,50,50	2.69	9 (30%)	24,82,82	3.16	14 (58%)
2	SO4	B	151	-	4,4,4	0.85	0	6,6,6	0.18	0
3	HEM	C	147	1	30,50,50	2.63	10 (33%)	24,82,82	3.13	12 (50%)
2	SO4	C	153	-	4,4,4	0.80	0	6,6,6	0.16	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HEM	D	147	1	30,50,50	2.87	9 (30%)	24,82,82	3.03	12 (50%)
2	SO4	D	152	-	4,4,4	0.80	0	6,6,6	0.12	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	-	0/10/54/54	0/0/8/8
2	SO4	A	150	-	-	0/0/0/0	0/0/0/0
3	HEM	B	147	1	-	0/10/54/54	0/0/8/8
2	SO4	B	151	-	-	0/0/0/0	0/0/0/0
3	HEM	C	147	1	-	0/10/54/54	0/0/8/8
2	SO4	C	153	-	-	0/0/0/0	0/0/0/0
3	HEM	D	147	1	-	0/10/54/54	0/0/8/8
2	SO4	D	152	-	-	0/0/0/0	0/0/0/0

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	147	HEM	C3B-C4B	-9.67	1.43	1.51
3	B	147	HEM	C3B-C4B	-8.87	1.43	1.51
3	A	147	HEM	C3B-C4B	-8.67	1.44	1.51
3	C	147	HEM	C2D-C3D	-7.09	1.33	1.54
3	A	147	HEM	C2D-C3D	-7.06	1.33	1.54
3	D	147	HEM	C2D-C3D	-6.89	1.33	1.54
3	C	147	HEM	C3B-C4B	-6.86	1.45	1.51
3	B	147	HEM	C2D-C3D	-6.83	1.34	1.54
3	D	147	HEM	C3D-C4D	-6.54	1.43	1.51
3	C	147	HEM	C3D-C4D	-6.41	1.43	1.51
3	A	147	HEM	C3D-C4D	-6.15	1.43	1.51
3	B	147	HEM	C3D-C4D	-5.74	1.44	1.51
3	C	147	HEM	C2C-C1C	-4.06	1.44	1.52
3	A	147	HEM	C2C-C1C	-3.90	1.45	1.52
3	D	147	HEM	C2C-C1C	-3.67	1.45	1.52
3	B	147	HEM	C2C-C1C	-3.53	1.45	1.52
3	A	147	HEM	C2B-C1B	-2.52	1.43	1.51
3	C	147	HEM	C2D-C1D	-2.49	1.43	1.51
3	D	147	HEM	C2D-C1D	-2.44	1.43	1.51
3	D	147	HEM	C2B-C1B	-2.44	1.43	1.51
3	B	147	HEM	C2B-C1B	-2.39	1.44	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	147	HEM	C2D-C1D	-2.22	1.44	1.51
3	C	147	HEM	C2B-C1B	-2.10	1.45	1.51
3	C	147	HEM	C2A-C3A	-2.08	1.31	1.37
3	B	147	HEM	C2D-C1D	-2.06	1.45	1.51
3	C	147	HEM	C4C-NC	2.01	1.38	1.36
3	B	147	HEM	CAA-C2A	2.04	1.55	1.52
3	C	147	HEM	C1C-NC	2.29	1.38	1.36
3	A	147	HEM	CAA-C2A	2.32	1.56	1.52
3	B	147	HEM	C1C-NC	2.33	1.38	1.36
3	D	147	HEM	C3B-CAB	2.50	1.56	1.51
3	D	147	HEM	C1C-NC	2.58	1.39	1.36
3	B	147	HEM	FE-NC	2.76	2.06	1.95
3	D	147	HEM	FE-NC	2.91	2.07	1.95
3	A	147	HEM	FE-NC	3.14	2.08	1.95
3	C	147	HEM	FE-NC	3.33	2.08	1.95

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	147	HEM	C3B-CAB-CBB	-6.63	114.29	124.46
3	C	147	HEM	CAA-C2A-C1A	-5.68	120.84	127.01
3	B	147	HEM	CAA-C2A-C1A	-5.39	121.16	127.01
3	B	147	HEM	CMA-C3A-C4A	-4.41	121.07	128.36
3	C	147	HEM	CAA-CBA-CGA	-4.20	105.05	112.75
3	C	147	HEM	C3B-CAB-CBB	-3.76	118.69	124.46
3	D	147	HEM	CMA-C3A-C4A	-3.62	122.37	128.36
3	A	147	HEM	CMA-C3A-C4A	-3.32	122.86	128.36
3	C	147	HEM	CMA-C3A-C4A	-2.96	123.47	128.36
3	A	147	HEM	CAA-C2A-C1A	-2.56	124.22	127.01
3	D	147	HEM	CAA-C2A-C1A	-2.56	124.23	127.01
3	C	147	HEM	CMA-C3A-C2A	2.11	129.65	125.24
3	C	147	HEM	C3B-C4B-CHC	2.12	126.15	123.16
3	A	147	HEM	CMA-C3A-C2A	2.21	129.86	125.24
3	A	147	HEM	C2C-C1C-CHC	2.30	127.17	123.68
3	B	147	HEM	C4B-CHC-C1C	2.30	129.67	125.82
3	C	147	HEM	CAD-C3D-C4D	2.40	120.92	112.47
3	A	147	HEM	CAA-CBA-CGA	2.44	117.22	112.75
3	B	147	HEM	CAD-C3D-C4D	2.44	121.08	112.47
3	D	147	HEM	C2C-C1C-CHC	2.47	127.43	123.68
3	D	147	HEM	CMA-C3A-C2A	2.50	130.46	125.24
3	D	147	HEM	CAD-C3D-C4D	2.91	122.73	112.47
3	A	147	HEM	C3B-C4B-CHC	2.98	127.36	123.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	147	HEM	CAD-C3D-C4D	3.02	123.13	112.47
3	D	147	HEM	C3B-C4B-CHC	3.20	127.66	123.16
3	B	147	HEM	C3C-CAC-CBC	3.24	129.42	124.46
3	B	147	HEM	CMA-C3A-C2A	3.28	132.09	125.24
3	B	147	HEM	CAA-CBA-CGA	3.28	118.75	112.75
3	B	147	HEM	C2C-C1C-CHC	3.29	128.69	123.68
3	D	147	HEM	CMD-C2D-C3D	3.46	129.65	114.35
3	B	147	HEM	CMD-C2D-C3D	3.50	129.81	114.35
3	A	147	HEM	CMD-C2D-C3D	3.56	130.10	114.35
3	C	147	HEM	CMD-C2D-C3D	3.86	131.42	114.35
3	B	147	HEM	C3B-C4B-CHC	4.09	128.91	123.16
3	A	147	HEM	C2D-C3D-C4D	4.42	108.99	101.50
3	D	147	HEM	C2D-C3D-C4D	4.42	109.00	101.50
3	B	147	HEM	CMB-C2B-C3B	4.56	127.91	116.53
3	B	147	HEM	C2D-C3D-C4D	4.58	109.26	101.50
3	B	147	HEM	CMC-C2C-C3C	4.77	128.43	116.53
3	D	147	HEM	CMB-C2B-C3B	4.86	128.67	116.53
3	A	147	HEM	CAD-C3D-C2D	5.01	127.63	113.22
3	D	147	HEM	CMC-C2C-C3C	5.12	129.31	116.53
3	C	147	HEM	C2D-C3D-C4D	5.18	110.28	101.50
3	D	147	HEM	CAD-C3D-C2D	5.20	128.17	113.22
3	A	147	HEM	CMB-C2B-C3B	5.29	129.74	116.53
3	A	147	HEM	CMC-C2C-C3C	5.37	129.92	116.53
3	C	147	HEM	CAD-C3D-C2D	5.37	128.65	113.22
3	C	147	HEM	CMC-C2C-C3C	5.55	130.39	116.53
3	B	147	HEM	CAD-C3D-C2D	5.72	129.66	113.22
3	C	147	HEM	CMB-C2B-C3B	6.09	131.72	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	147	HEM	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

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

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

6.3 Carbohydrates [i](#)

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

6.4 Ligands [i](#)

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

6.5 Other polymers [i](#)

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