



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:31 AM GMT

PDB ID : 2HHD
Title : OXYGEN AFFINITY MODULATION BY THE N-TERMINI OF THE BETA-CHAINS IN HUMAN AND BOVINE HEMOGLOBIN
Authors : Gilliland, G.L.; Pechik, I.; Fronticelli, C.; Ji, X.
Deposited on : 1994-09-29
Resolution : 2.20 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

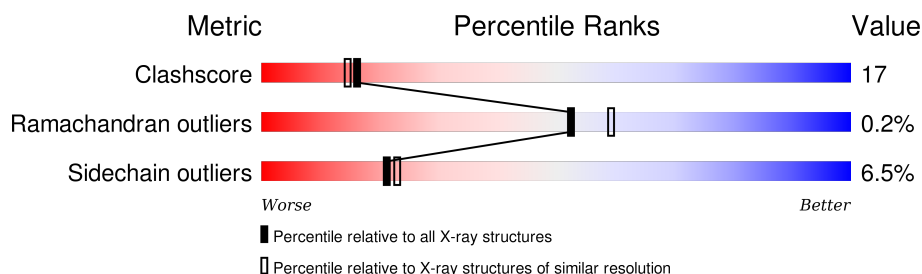
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	4477 (2.20-2.20)
Ramachandran outliers	100387	4404 (2.20-2.20)
Sidechain outliers	100360	4405 (2.20-2.20)

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	141	 65% 29% 5%
1	C	141	 66% 28% 5%
2	B	146	 60% 36% 2%
2	D	146	 54% 38% 8%

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	SO4	B	147	-	X	-	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 5040 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.

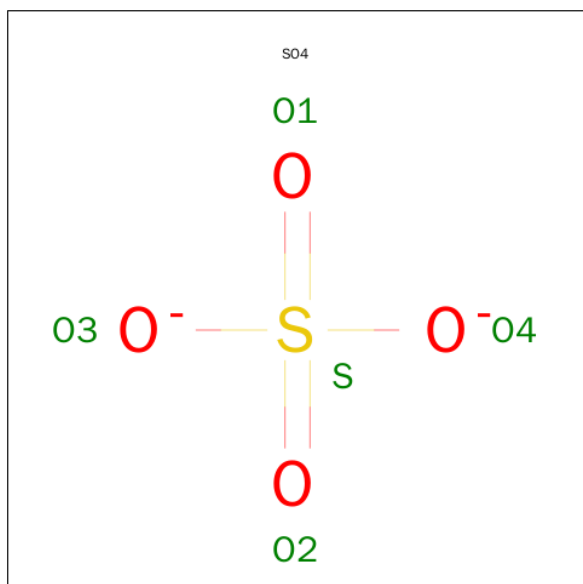
- Molecule 1 is a protein called HEMOGLOBIN (DEOXY) (ALPHA CHAIN).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	141	Total	C	N	O	S	0	0	0
			1069	685	187	194	3			
1	C	141	Total	C	N	O	S	0	0	0
			1069	685	187	194	3			

- Molecule 2 is a protein called HEMOGLOBIN (DEOXY) (BETA CHAIN).

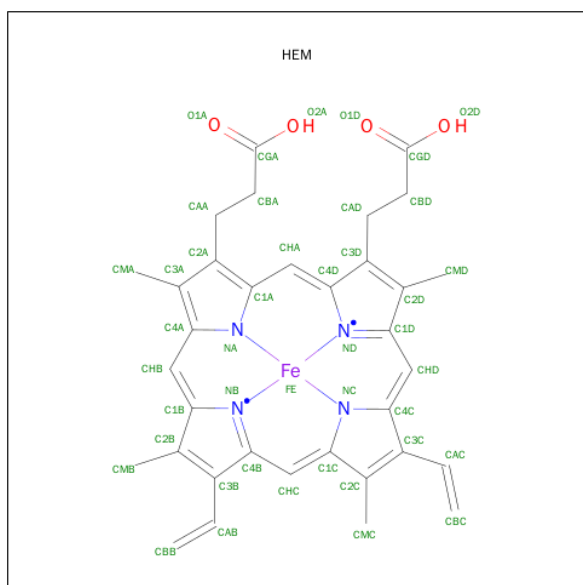
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1123	724	195	201	3			
2	D	146	Total	C	N	O	S	0	0	0
			1123	724	195	201	3			

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	132	Total O 132 132	0	0
5	B	116	Total O 116 116	0	0
5	C	126	Total O 126 126	0	0

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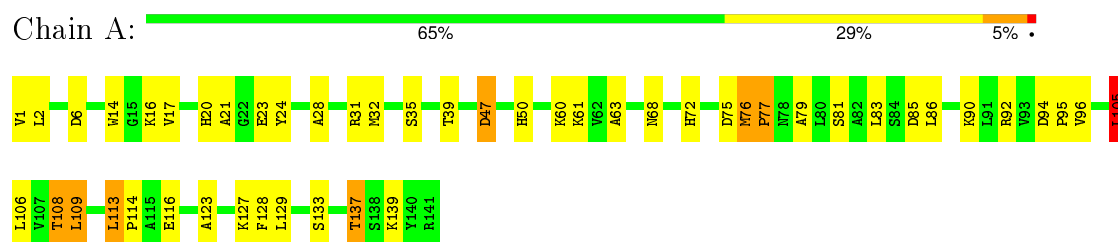
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	100	Total	O	0	0
			100	100		

3 Residue-property plots

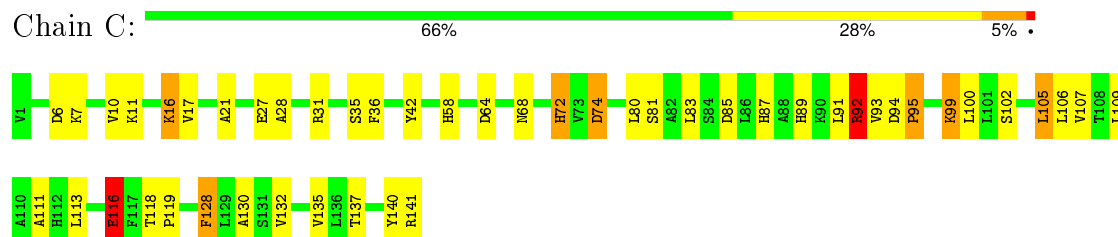
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

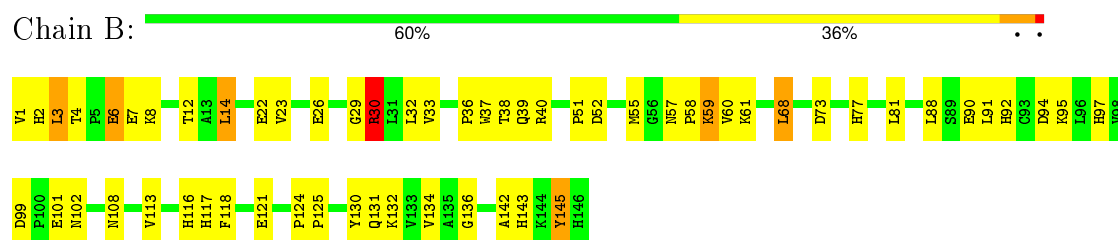
• Molecule 1: HEMOGLOBIN (DEOXY) (ALPHA CHAIN)



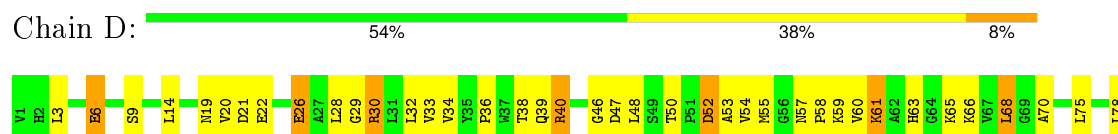
• Molecule 1: HEMOGLOBIN (DEOXY) (ALPHA CHAIN)



• Molecule 2: HEMOGLOBIN (DEOXY) (BETA CHAIN)



• Molecule 2: HEMOGLOBIN (DEOXY) (BETA CHAIN)





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	62.45Å 82.13Å 53.76Å 90.00° 98.87° 90.00°	Depositor
Resolution (Å)	6.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (6.00-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.137 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5040	wwPDB-VP
Average B, all atoms (Å ²)	16.0	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	0.94	0/1097	1.63	12/1491 (0.8%)
1	C	0.97	0/1097	1.76	20/1491 (1.3%)
2	B	0.95	0/1153	1.72	19/1566 (1.2%)
2	D	0.95	0/1153	1.64	17/1566 (1.1%)
All	All	0.95	0/4500	1.69	68/6114 (1.1%)

There are no bond length outliers.

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	92	ARG	NE-CZ-NH2	-13.25	113.67	120.30
1	A	92	ARG	NE-CZ-NH2	-12.88	113.86	120.30
2	B	30	ARG	NE-CZ-NH2	-11.68	114.46	120.30
2	B	40	ARG	NE-CZ-NH1	11.38	125.99	120.30
1	C	92	ARG	NE-CZ-NH1	10.64	125.62	120.30
1	C	6	ASP	CB-CG-OD2	-9.99	109.31	118.30
2	D	30	ARG	NE-CZ-NH2	-9.91	115.35	120.30
1	C	64	ASP	CB-CG-OD1	9.53	126.88	118.30
2	B	94	ASP	CB-CG-OD2	-9.52	109.73	118.30
1	C	92	ARG	CD-NE-CZ	8.86	136.00	123.60
1	C	74	ASP	CB-CG-OD2	8.73	126.16	118.30
1	A	31	ARG	NE-CZ-NH1	8.65	124.62	120.30
2	D	40	ARG	NE-CZ-NH1	8.46	124.53	120.30
1	C	31	ARG	NE-CZ-NH1	8.18	124.39	120.30
1	C	141	ARG	NE-CZ-NH2	-8.18	116.21	120.30
1	A	31	ARG	NE-CZ-NH2	-7.83	116.38	120.30
1	C	31	ARG	NE-CZ-NH2	-7.70	116.45	120.30
2	B	145	TYR	CB-CG-CD1	-7.64	116.42	121.00
2	B	94	ASP	CB-CG-OD1	7.37	124.93	118.30
2	D	99	ASP	CB-CG-OD2	7.16	124.75	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	130	ALA	CB-CA-C	6.89	120.44	110.10
1	A	6	ASP	O-C-N	6.79	133.56	122.70
2	B	30	ARG	NH1-CZ-NH2	6.73	126.80	119.40
1	C	64	ASP	CB-CG-OD2	-6.44	112.50	118.30
2	B	14	LEU	CA-CB-CG	6.43	130.09	115.30
1	C	85	ASP	CB-CG-OD2	-6.41	112.53	118.30
1	A	105	LEU	CA-CB-CG	6.39	130.00	115.30
2	D	70	ALA	CB-CA-C	6.35	119.63	110.10
2	B	68	LEU	CA-CB-CG	6.29	129.75	115.30
1	A	76	MET	CG-SD-CE	6.22	110.15	100.20
2	D	79	ASP	CB-CG-OD2	6.18	123.87	118.30
1	A	92	ARG	NE-CZ-NH1	6.16	123.38	120.30
1	C	89	HIS	CA-CB-CG	-6.06	103.30	113.60
2	B	52	ASP	CB-CG-OD2	-6.03	112.88	118.30
1	C	27	GLU	CG-CD-OE1	-5.78	106.74	118.30
2	D	28	LEU	CB-CA-C	5.76	121.15	110.20
1	A	47	ASP	N-CA-C	-5.72	95.56	111.00
2	B	6	GLU	CA-CB-CG	5.71	125.97	113.40
2	B	101	GLU	OE1-CD-OE2	5.71	130.15	123.30
2	D	6	GLU	CA-CB-CG	5.63	125.79	113.40
2	B	121	GLU	CG-CD-OE2	-5.62	107.05	118.30
2	D	47	ASP	CB-CG-OD2	-5.60	113.26	118.30
2	B	73	ASP	CB-CG-OD1	5.56	123.30	118.30
2	D	26	GLU	CG-CD-OE2	5.55	129.41	118.30
1	A	79	ALA	CB-CA-C	5.55	118.42	110.10
2	B	99	ASP	CB-CG-OD2	-5.52	113.34	118.30
2	D	92	HIS	CA-CB-CG	-5.50	104.25	113.60
1	A	113	LEU	CA-CB-CG	5.49	127.93	115.30
2	D	129	ALA	CB-CA-C	5.45	118.28	110.10
1	C	102	SER	N-CA-CB	5.43	118.65	110.50
2	B	134	VAL	CA-CB-CG2	5.42	119.04	110.90
1	C	27	GLU	CG-CD-OE2	5.30	128.90	118.30
1	C	116	GLU	CA-CB-CG	5.28	125.02	113.40
1	C	72	HIS	CA-CB-CG	-5.26	104.66	113.60
1	C	109	LEU	CA-CB-CG	5.14	127.12	115.30
2	D	40	ARG	CD-NE-CZ	5.14	130.79	123.60
2	D	79	ASP	CB-CG-OD1	-5.11	113.70	118.30
1	C	128	PHE	CB-CG-CD2	-5.10	117.23	120.80
2	B	130	TYR	CB-CG-CD1	-5.10	117.94	121.00
2	B	90	GLU	CG-CD-OE1	5.09	128.49	118.30
2	D	86	ALA	CB-CA-C	-5.09	102.47	110.10
2	B	101	GLU	CG-CD-OE2	-5.09	108.13	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	108	THR	CA-CB-CG2	5.06	119.49	112.40
2	D	47	ASP	CB-CG-OD1	5.04	122.84	118.30
2	D	52	ASP	CB-CG-OD1	5.03	122.83	118.30
2	D	60	VAL	CA-C-O	5.01	130.63	120.10
1	A	50	HIS	CA-CB-CG	5.00	122.11	113.60
2	B	77	HIS	CA-CB-CG	-5.00	105.09	113.60

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	1069	0	1073	31	0
1	C	1069	0	1073	34	0
2	B	1123	0	1118	42	0
2	D	1123	0	1118	46	0
3	B	5	0	0	1	0
3	D	5	0	0	1	0
4	A	43	0	30	4	0
4	B	43	0	30	4	0
4	C	43	0	30	1	0
4	D	43	0	30	1	0
5	A	132	0	0	4	0
5	B	116	0	0	6	0
5	C	126	0	0	4	0
5	D	100	0	0	8	0
All	All	5040	0	4502	152	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 17.

All (152) 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:113:LEU:HB3	1:C:116:GLU:HG2	1.30	1.10
1:C:16:LYS:HG3	1:C:116:GLU:HG3	1.36	1.03
1:A:90:LYS:HE3	5:A:211:HOH:O	1.70	0.90
1:A:16:LYS:HG3	1:A:116:GLU:HG2	1.56	0.86
2:B:143:HIS:HB2	5:B:184:HOH:O	1.79	0.82
2:D:50:THR:HG23	2:D:53:ALA:H	1.51	0.75
2:D:46:GLY:HA3	5:D:243:HOH:O	1.89	0.73
2:D:19:ASN:ND2	2:D:22:GLU:HG2	2.04	0.72
1:C:118:THR:HG21	5:C:159:HOH:O	1.91	0.69
2:D:20:VAL:HG13	2:D:68:LEU:HB3	1.74	0.69
2:D:55:MET:O	2:D:61:LYS:HE2	1.94	0.68
2:B:57:ASN:O	2:B:61:LYS:HG3	1.94	0.67
4:B:148:HEM:CMB	4:B:148:HEM:HBB2	2.25	0.67
2:B:23:VAL:HG13	2:B:113:VAL:HG11	1.79	0.64
4:B:148:HEM:HBC2	4:B:148:HEM:HMC2	1.80	0.64
2:D:19:ASN:HB2	5:D:193:HOH:O	1.97	0.63
2:B:58:PRO:HA	2:B:61:LYS:HZ2	1.62	0.62
2:B:38:THR:HG22	2:B:102:ASN:OD1	1.98	0.62
4:B:148:HEM:HMB1	4:B:148:HEM:HBB2	1.81	0.62
1:C:7:LYS:O	1:C:11:LYS:HG3	2.02	0.59
2:D:3:LEU:CD2	2:D:132:LYS:HG3	2.32	0.59
1:C:81:SER:HB3	5:C:162:HOH:O	2.02	0.59
1:A:83:LEU:HD21	4:A:142:HEM:HMA3	1.85	0.59
2:B:36:PRO:O	2:B:39:GLN:HG3	2.03	0.58
2:D:50:THR:CG2	2:D:53:ALA:HB2	2.34	0.57
2:D:75:LEU:HA	2:D:78:LEU:HD13	1.86	0.57
1:A:133:SER:O	1:A:137:THR:HB	2.05	0.57
2:D:82:LYS:HG2	3:D:147:SO4:O4	2.05	0.57
2:B:58:PRO:HA	2:B:61:LYS:CE	2.35	0.56
2:B:58:PRO:HA	2:B:61:LYS:NZ	2.19	0.56
2:B:14:LEU:HB2	5:B:220:HOH:O	2.06	0.56
1:A:24:TYR:O	1:A:108:THR:HG21	2.05	0.56
1:A:86:LEU:HD21	4:A:142:HEM:HBA2	1.88	0.56
1:C:28:ALA:CB	1:C:105:LEU:HD13	2.36	0.55
2:D:50:THR:HG23	2:D:53:ALA:N	2.21	0.55
1:C:99:LYS:HE2	5:D:225:HOH:O	2.06	0.55
1:C:111:ALA:HA	2:D:115:ALA:O	2.07	0.55
2:B:4:THR:OG1	2:B:7:GLU:HG3	2.06	0.55
1:C:118:THR:HG23	1:C:119:PRO:HD2	1.88	0.55
2:B:26:GLU:O	2:B:30:ARG:HB2	2.08	0.54
1:A:2:LEU:HD11	1:A:128:PHE:HD1	1.72	0.54
1:C:35:SER:HB3	2:D:131:GLN:HG3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:29:GLY:O	2:D:33:VAL:HG23	2.07	0.53
1:A:20:HIS:HB3	1:A:24:TYR:CE1	2.44	0.53
2:D:124:PRO:HB2	2:D:125:PRO:HD3	1.90	0.53
2:D:53:ALA:O	2:D:57:ASN:HB2	2.09	0.52
2:B:1:VAL:N	3:B:147:SO4:O4	2.42	0.52
1:A:16:LYS:O	1:A:17:VAL:C	2.48	0.52
4:B:148:HEM:CMC	4:B:148:HEM:HBC2	2.39	0.52
5:A:153:HOH:O	2:D:40:ARG:HD3	2.09	0.51
2:D:9:SER:HB3	5:D:224:HOH:O	2.10	0.51
1:A:21:ALA:HB1	1:A:63:ALA:HB1	1.93	0.51
2:D:50:THR:HG22	2:D:53:ALA:HB2	1.92	0.51
1:A:114:PRO:O	2:B:116:HIS:NE2	2.38	0.51
1:C:80:LEU:HD22	1:C:83:LEU:HD12	1.93	0.51
1:C:11:LYS:NZ	5:C:164:HOH:O	2.40	0.51
2:D:3:LEU:HD21	2:D:132:LYS:HG3	1.92	0.51
2:B:4:THR:HG22	5:B:261:HOH:O	2.11	0.51
2:D:21:ASP:HA	2:D:65:LYS:HG3	1.92	0.51
1:A:94:ASP:HB2	5:D:187:HOH:O	2.11	0.50
2:D:104:ARG:H	2:D:104:ARG:HD2	1.76	0.50
2:B:8:LYS:HD3	5:B:174:HOH:O	2.11	0.50
1:A:14:TRP:O	1:A:17:VAL:HB	2.10	0.50
2:D:66:LYS:HG3	5:D:203:HOH:O	2.12	0.50
2:D:14:LEU:HD11	2:D:118:PHE:CD2	2.47	0.50
1:A:76:MET:N	1:A:77:PRO:HD2	2.27	0.49
1:C:42:TYR:CE1	1:C:93:VAL:HA	2.48	0.49
1:C:16:LYS:HA	1:C:16:LYS:HE3	1.94	0.49
1:C:137:THR:HA	1:C:140:TYR:CE1	2.48	0.49
2:B:3:LEU:HD23	2:B:7:GLU:HB3	1.94	0.49
2:B:1:VAL:HG13	2:B:81:LEU:HD12	1.95	0.49
2:B:92:HIS:O	2:B:97:HIS:N	2.46	0.49
2:D:143:HIS:HB3	5:D:184:HOH:O	2.13	0.49
2:B:37:TRP:HA	1:C:92:ARG:HD2	1.95	0.48
1:A:35:SER:HB3	2:B:131:GLN:HG3	1.95	0.48
4:D:148:HEM:CMC	4:D:148:HEM:HBC2	2.43	0.48
2:B:91:LEU:HD12	2:B:95:LYS:HB2	1.96	0.48
1:C:80:LEU:CD2	1:C:83:LEU:HD12	2.43	0.47
2:D:117:HIS:HD2	2:D:118:PHE:CZ	2.31	0.47
2:B:51:PRO:O	2:B:55:MET:HG2	2.14	0.47
1:C:68:ASN:O	1:C:72:HIS:HD2	1.97	0.47
2:B:7:GLU:OE2	2:B:132:LYS:NZ	2.47	0.47
2:D:14:LEU:HD11	2:D:118:PHE:CG	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:38:THR:HG22	2:D:102:ASN:OD1	2.14	0.46
2:B:2:HIS:O	2:B:2:HIS:ND1	2.49	0.46
2:B:14:LEU:HD21	2:B:118:PHE:CG	2.51	0.46
1:A:76:MET:N	1:A:77:PRO:CD	2.78	0.46
2:B:57:ASN:HB3	2:B:60:VAL:HG23	1.98	0.46
2:D:48:LEU:HA	2:D:54:VAL:HG22	1.98	0.46
1:C:28:ALA:HB1	1:C:105:LEU:HD13	1.98	0.46
1:C:17:VAL:HG12	1:C:21:ALA:HB2	1.97	0.46
2:B:29:GLY:O	2:B:33:VAL:HG23	2.16	0.46
1:C:87:HIS:CG	1:C:91:LEU:HD12	2.51	0.46
2:D:129:ALA:O	2:D:133:VAL:HG23	2.15	0.46
2:B:58:PRO:HA	2:B:61:LYS:HE3	1.98	0.45
2:B:1:VAL:HG11	2:B:136:GLY:HA3	1.97	0.45
1:A:83:LEU:HD11	4:A:142:HEM:HMA3	1.97	0.45
4:A:142:HEM:HMD2	4:A:142:HEM:HBD1	1.97	0.45
1:A:16:LYS:HG3	1:A:116:GLU:CG	2.39	0.45
2:B:32:LEU:HG	2:B:39:GLN:HG2	1.99	0.45
2:D:111:VAL:HG13	2:D:122:PHE:HZ	1.82	0.45
2:B:108:ASN:ND2	5:B:195:HOH:O	2.50	0.44
1:C:7:LYS:HE3	1:C:74:ASP:OD2	2.18	0.44
2:D:92:HIS:HB3	2:D:145:TYR:OH	2.18	0.44
1:A:94:ASP:HA	1:A:95:PRO:HD3	1.75	0.44
1:C:137:THR:HA	1:C:140:TYR:CD1	2.52	0.44
1:A:96:VAL:HG11	2:D:101:GLU:HG2	1.98	0.44
1:C:128:PHE:O	1:C:132:VAL:HG23	2.17	0.44
2:D:19:ASN:HD22	2:D:22:GLU:HG2	1.82	0.44
2:B:57:ASN:HA	2:B:58:PRO:HD3	1.81	0.44
1:C:58:HIS:HE1	4:C:142:HEM:C1A	2.36	0.43
2:D:96:LEU:HD21	5:D:229:HOH:O	2.17	0.43
2:B:88:LEU:HD23	2:B:91:LEU:HD23	1.99	0.43
2:D:36:PRO:O	2:D:39:GLN:HB2	2.18	0.43
2:B:57:ASN:HB3	2:B:60:VAL:CG2	2.48	0.43
2:D:127:GLN:O	2:D:131:GLN:HG2	2.19	0.43
2:B:142:ALA:O	2:B:145:TYR:HB2	2.19	0.43
2:D:123:THR:OG1	2:D:126:VAL:HG23	2.18	0.43
1:A:123:ALA:O	1:A:127:LYS:HG3	2.18	0.43
1:A:85:ASP:OD1	1:A:139:LYS:HD3	2.19	0.43
1:A:32:MET:CG	1:A:39:THR:HG21	2.48	0.43
2:D:104:ARG:HD2	2:D:104:ARG:N	2.34	0.42
2:D:57:ASN:HA	2:D:58:PRO:HD3	1.64	0.42
2:D:34:VAL:C	2:D:36:PRO:HD3	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:87:HIS:HA	1:C:91:LEU:HD12	2.00	0.42
1:A:90:LYS:HB3	5:A:174:HOH:O	2.18	0.42
1:C:92:ARG:HH11	1:C:92:ARG:HG3	1.84	0.42
1:C:10:VAL:HG21	1:C:128:PHE:HB2	2.02	0.42
1:C:135:VAL:HG23	5:C:235:HOH:O	2.19	0.42
2:B:2:HIS:H	2:B:2:HIS:CD2	2.37	0.42
1:A:106:LEU:HD13	1:A:129:LEU:HD12	2.02	0.42
2:D:131:GLN:O	2:D:134:VAL:HG22	2.20	0.42
2:B:32:LEU:HD12	2:B:38:THR:OG1	2.20	0.41
1:A:68:ASN:OD1	1:A:72:HIS:CD2	2.73	0.41
1:A:83:LEU:HD12	1:A:83:LEU:HA	1.79	0.41
1:A:20:HIS:ND1	1:A:23:GLU:CD	2.74	0.41
1:C:94:ASP:HA	1:C:95:PRO:HD3	1.94	0.41
2:B:57:ASN:OD1	2:B:59:LYS:HB2	2.20	0.41
1:C:107:VAL:HG13	2:D:115:ALA:HB2	2.02	0.41
1:A:28:ALA:CB	1:A:105:LEU:HD13	2.50	0.41
2:D:105:LEU:HD23	2:D:105:LEU:HA	1.94	0.41
1:A:81:SER:HB3	5:A:228:HOH:O	2.21	0.41
2:D:30:ARG:HD2	2:D:113:VAL:HG22	2.02	0.41
2:B:1:VAL:HG23	2:B:3:LEU:HD12	2.03	0.41
1:C:83:LEU:HA	1:C:83:LEU:HD23	1.82	0.41
1:A:105:LEU:O	1:A:109:LEU:HB2	2.21	0.41
2:B:12:THR:HG22	5:B:188:HOH:O	2.19	0.41
1:C:137:THR:HG22	1:C:140:TYR:CE1	2.56	0.40
2:B:22:GLU:CD	2:B:117:HIS:HE2	2.24	0.40
2:B:124:PRO:N	2:B:125:PRO:CD	2.84	0.40
2:D:63:HIS:O	2:D:66:LYS:HB2	2.22	0.40
1:C:36:PHE:CD1	1:C:100:LEU:HD22	2.57	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	139/141 (99%)	134 (96%)	4 (3%)	1 (1%)	26	25
1	C	139/141 (99%)	137 (99%)	2 (1%)	0	100	100
2	B	144/146 (99%)	139 (96%)	5 (4%)	0	100	100
2	D	144/146 (99%)	137 (95%)	7 (5%)	0	100	100
All	All	566/574 (99%)	547 (97%)	18 (3%)	1 (0%)	52	59

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	75	ASP

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	113/113 (100%)	104 (92%)	9 (8%)	15	15
1	C	113/113 (100%)	106 (94%)	7 (6%)	23	25
2	B	118/118 (100%)	113 (96%)	5 (4%)	36	44
2	D	118/118 (100%)	109 (92%)	9 (8%)	16	16
All	All	462/462 (100%)	432 (94%)	30 (6%)	21	23

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	VAL
1	A	47	ASP
1	A	60	LYS
1	A	61	LYS
1	A	77	PRO
1	A	105	LEU
1	A	109	LEU
1	A	113	LEU
1	A	137	THR
2	B	3	LEU

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Mol	Chain	Res	Type
2	B	6	GLU
2	B	30	ARG
2	B	59	LYS
2	B	68	LEU
1	C	16	LYS
1	C	92	ARG
1	C	95	PRO
1	C	99	LYS
1	C	105	LEU
1	C	106	LEU
1	C	116	GLU
2	D	6	GLU
2	D	26	GLU
2	D	32	LEU
2	D	52	ASP
2	D	59	LYS
2	D	61	LYS
2	D	68	LEU
2	D	104	ARG
2	D	132	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	HIS
2	B	108	ASN
1	C	72	HIS
2	D	63	HIS
2	D	117	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

6 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
4	HEM	A	142	1	30,50,50	2.47	7 (23%)	24,82,82	2.61	11 (45%)
3	SO4	B	147	-	4,4,4	2.13	4 (100%)	6,6,6	0.65	0
4	HEM	B	148	2	30,50,50	2.49	5 (16%)	24,82,82	3.21	11 (45%)
4	HEM	C	142	1	30,50,50	2.65	9 (30%)	24,82,82	2.42	8 (33%)
3	SO4	D	147	-	4,4,4	2.15	2 (50%)	6,6,6	0.82	0
4	HEM	D	148	2	30,50,50	2.13	8 (26%)	24,82,82	3.66	15 (62%)

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	A	142	1	-	0/10/54/54	0/0/8/8
3	SO4	B	147	-	-	0/0/0/0	0/0/0/0
4	HEM	B	148	2	-	0/10/54/54	0/0/8/8
4	HEM	C	142	1	-	0/10/54/54	0/0/8/8
3	SO4	D	147	-	-	0/0/0/0	0/0/0/0
4	HEM	D	148	2	-	0/10/54/54	0/0/8/8

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	142	HEM	C3B-C4B	-8.51	1.44	1.51
4	B	148	HEM	C3B-C4B	-8.15	1.44	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	142	HEM	C2D-C3D	-7.22	1.32	1.54
4	C	142	HEM	C2D-C3D	-6.92	1.33	1.54
4	D	148	HEM	C2D-C3D	-6.53	1.34	1.54
4	B	148	HEM	C2D-C3D	-6.51	1.35	1.54
4	A	142	HEM	C3B-C4B	-6.20	1.46	1.51
4	C	142	HEM	C3D-C4D	-5.59	1.44	1.51
4	D	148	HEM	C3B-C4B	-5.35	1.47	1.51
4	B	148	HEM	C3D-C4D	-5.09	1.45	1.51
4	B	148	HEM	C2C-C1C	-4.29	1.44	1.52
4	A	142	HEM	C3D-C4D	-3.87	1.46	1.51
4	D	148	HEM	C3D-C4D	-3.78	1.46	1.51
4	A	142	HEM	C2C-C1C	-3.55	1.45	1.52
4	D	148	HEM	C2C-C1C	-3.40	1.46	1.52
4	C	142	HEM	C2C-C1C	-3.14	1.46	1.52
4	A	142	HEM	C2D-C1D	-2.58	1.43	1.51
4	C	142	HEM	C3C-CAC	2.01	1.55	1.51
3	B	147	SO4	O4-S	2.01	1.54	1.47
4	C	142	HEM	CMA-C3A	2.01	1.55	1.51
4	B	148	HEM	C3B-CAB	2.02	1.55	1.51
3	B	147	SO4	O3-S	2.04	1.54	1.47
4	C	142	HEM	C4C-NC	2.12	1.38	1.36
3	B	147	SO4	O1-S	2.15	1.54	1.47
4	D	148	HEM	C1C-NC	2.22	1.38	1.36
4	C	142	HEM	C3B-CAB	2.29	1.55	1.51
3	B	147	SO4	O2-S	2.31	1.55	1.47
4	D	148	HEM	C3C-CAC	2.33	1.55	1.51
4	A	142	HEM	FE-NC	2.39	2.05	1.95
4	D	148	HEM	C4C-NC	2.49	1.39	1.36
3	D	147	SO4	O3-S	2.55	1.56	1.47
3	D	147	SO4	O1-S	2.60	1.56	1.47
4	D	148	HEM	C3B-CAB	2.66	1.56	1.51
4	C	142	HEM	CAA-C2A	2.93	1.57	1.52
4	A	142	HEM	C4C-NC	4.67	1.41	1.36

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	148	HEM	C3C-CAC-CBC	-5.92	115.38	124.46
4	B	148	HEM	CMA-C3A-C4A	-5.57	119.15	128.36
4	B	148	HEM	CAA-C2A-C1A	-5.50	121.04	127.01
4	B	148	HEM	C3B-CAB-CBB	-5.26	116.39	124.46
4	A	142	HEM	CMA-C3A-C4A	-5.04	120.03	128.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	148	HEM	CMA-C3A-C4A	-4.83	120.36	128.36
4	D	148	HEM	C3B-CAB-CBB	-4.43	117.66	124.46
4	C	142	HEM	CAA-C2A-C1A	-4.10	122.56	127.01
4	B	148	HEM	CBA-CAA-C2A	-3.54	106.18	112.53
4	A	142	HEM	CAA-C2A-C1A	-3.11	123.63	127.01
4	A	142	HEM	C1D-CHD-C4C	-2.83	121.09	125.82
4	D	148	HEM	CAA-CBA-CGA	-2.65	107.89	112.75
4	C	142	HEM	C3B-CAB-CBB	-2.52	120.59	124.46
4	D	148	HEM	C4B-CHC-C1C	-2.48	121.68	125.82
4	A	142	HEM	CAA-CBA-CGA	2.04	116.49	112.75
4	D	148	HEM	CAD-CBD-CGD	2.22	122.08	113.02
4	A	142	HEM	CMD-C2D-C3D	2.57	125.71	114.35
4	D	148	HEM	CBD-CAD-C3D	2.74	121.54	113.55
4	B	148	HEM	CAD-C3D-C4D	2.87	122.61	112.47
4	D	148	HEM	CMD-C2D-C3D	2.90	127.20	114.35
4	D	148	HEM	CMA-C3A-C2A	2.97	131.44	125.24
4	A	142	HEM	CMA-C3A-C2A	3.03	131.56	125.24
4	D	148	HEM	C2D-C3D-C4D	3.05	106.67	101.50
4	A	142	HEM	CAD-C3D-C4D	3.09	123.39	112.47
4	C	142	HEM	CAD-C3D-C4D	3.23	123.85	112.47
4	A	142	HEM	C2D-C3D-C4D	3.28	107.05	101.50
4	B	148	HEM	CMD-C2D-C3D	3.44	129.55	114.35
4	C	142	HEM	CMD-C2D-C3D	3.46	129.64	114.35
4	D	148	HEM	CAD-C3D-C4D	3.55	124.98	112.47
4	C	142	HEM	CMC-C2C-C3C	3.87	126.18	116.53
4	C	142	HEM	C2D-C3D-C4D	3.94	108.17	101.50
4	A	142	HEM	CMB-C2B-C3B	3.95	126.38	116.53
4	B	148	HEM	C2D-C3D-C4D	4.28	108.75	101.50
4	B	148	HEM	CMC-C2C-C3C	4.38	127.45	116.53
4	B	148	HEM	CMA-C3A-C2A	4.55	134.75	125.24
4	D	148	HEM	CMC-C2C-C3C	4.75	128.39	116.53
4	D	148	HEM	CMB-C2B-C3B	4.81	128.54	116.53
4	A	142	HEM	CMC-C2C-C3C	4.98	128.97	116.53
4	C	142	HEM	CAD-C3D-C2D	5.13	127.96	113.22
4	D	148	HEM	CAD-C3D-C2D	5.26	128.33	113.22
4	B	148	HEM	CMB-C2B-C3B	5.27	129.68	116.53
4	B	148	HEM	CAD-C3D-C2D	5.30	128.45	113.22
4	C	142	HEM	CMB-C2B-C3B	5.44	130.11	116.53
4	A	142	HEM	CAD-C3D-C2D	5.65	129.45	113.22
4	D	148	HEM	CBA-CAA-C2A	9.91	130.30	112.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	142	HEM	4	0
3	B	147	SO4	1	0
4	B	148	HEM	4	0
4	C	142	HEM	1	0
3	D	147	SO4	1	0
4	D	148	HEM	1	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 [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.