



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 10:09 PM GMT

PDB ID : 1SDK
Title : CROSS-LINKED, CARBONMONOXY HEMOGLOBIN A
Authors : Schumacher, M.A.; Dixon, M.M.; Kluger, R.; Jones, R.T.; Brennan, R.G.
Deposited on : 1996-02-26
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

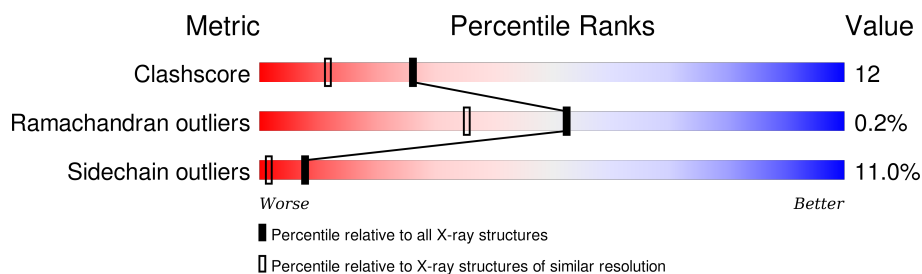
1 Overall quality at a glance

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	141	 68% 26% . .
1	C	141	 75% 21% . .
2	B	146	 75% 20% 5% .
2	D	146	 73% 18% 9%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 5015 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.

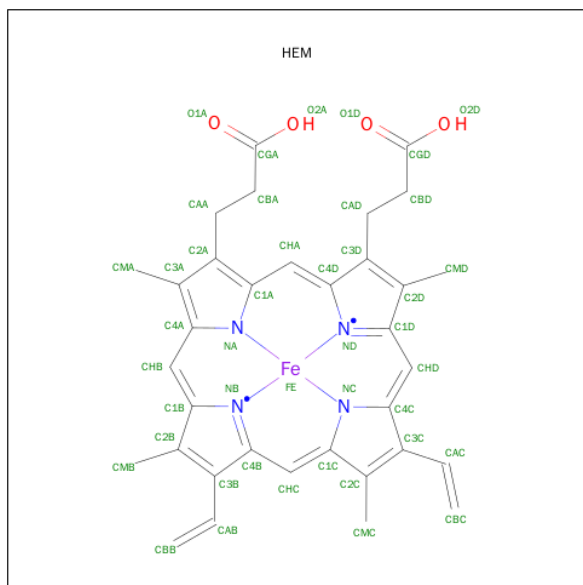
- Molecule 1 is a protein called HEMOGLOBIN A.

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 A.

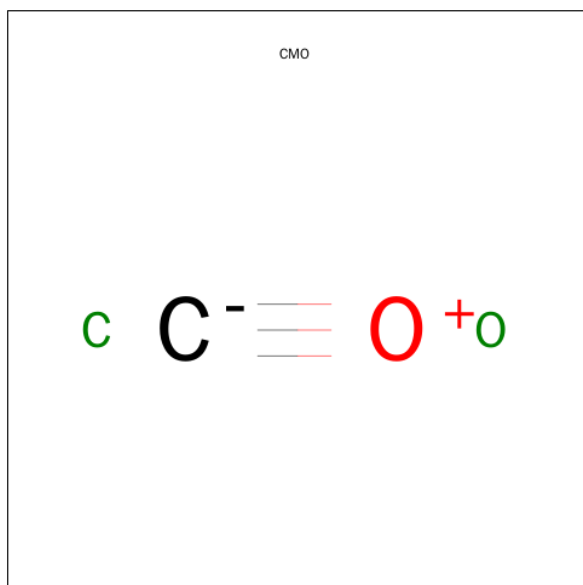
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1123	724	195	201	3			
2	D	146	Total	C	N	O	S	0	0	0
			1123	724	195	201	3			

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



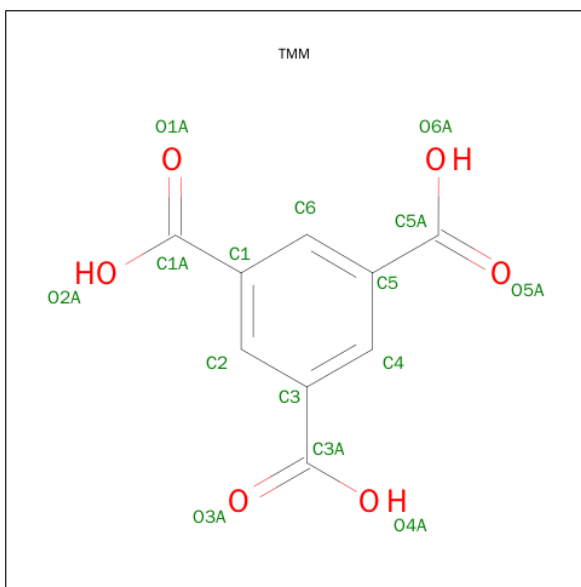
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
3	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
3	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
3	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- 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 1,3,5-BENZENETRICARBOXYLIC ACID (three-letter code: TMM) (formula: C₉H₆O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	O	0	0
			12	9	3		

- Molecule 6 is water.

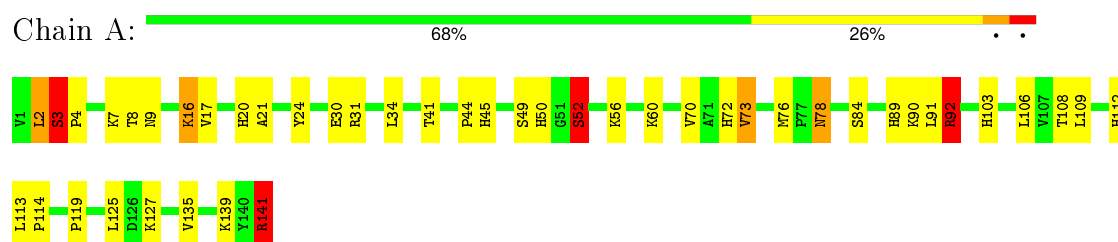
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	102	Total	O	0	0
			102	102		
6	B	118	Total	O	0	0
			118	118		
6	C	99	Total	O	0	0
			99	99		
6	D	120	Total	O	0	0
			120	120		

3 Residue-property plots

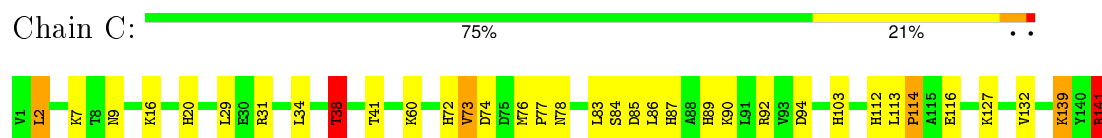
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

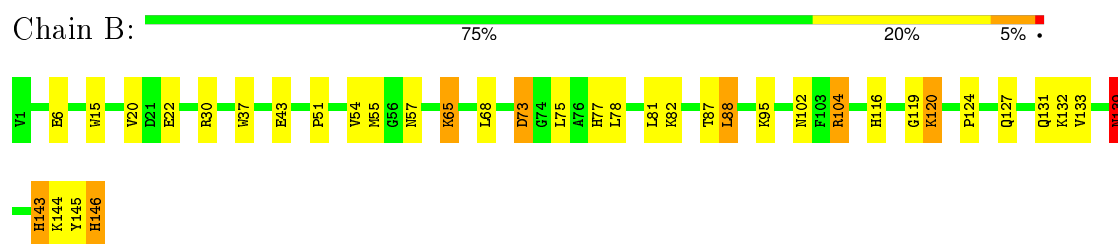
• Molecule 1: HEMOGLOBIN A



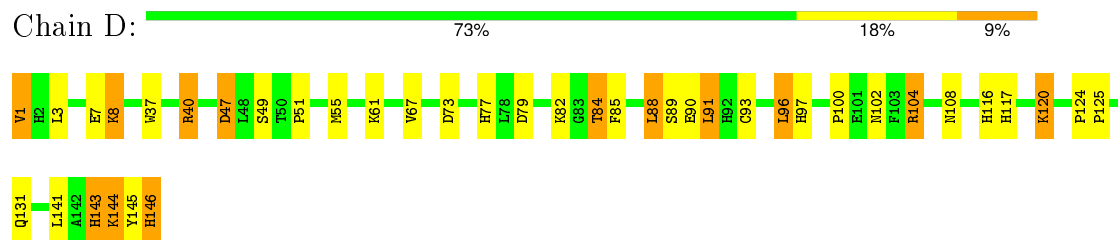
• Molecule 1: HEMOGLOBIN A



• Molecule 2: HEMOGLOBIN A



• Molecule 2: HEMOGLOBIN A



4 Data and refinement statistics

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	104.43Å 72.16Å 88.03Å 90.00° 108.25° 90.00°	Depositor
Resolution (Å)	8.00 – 1.80	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-1.80)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	TNT	Depositor
R, R_{free}	0.139 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5015	wwPDB-VP
Average B, all atoms (Å ²)	24.0	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, TMM

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	1.13	2/1097 (0.2%)	1.34	12/1491 (0.8%)
1	C	1.15	4/1097 (0.4%)	1.22	6/1491 (0.4%)
2	B	1.11	2/1153 (0.2%)	1.27	5/1566 (0.3%)
2	D	1.16	2/1153 (0.2%)	1.23	7/1566 (0.4%)
All	All	1.13	10/4500 (0.2%)	1.27	30/6114 (0.5%)

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	A	0	1
1	C	1	0
2	B	0	1
2	D	2	0
All	All	3	2

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	22	GLU	CG-CD	6.04	1.61	1.51
2	D	8	LYS	CB-CG	-5.77	1.36	1.52
2	B	139	ASN	CB-CG	-5.65	1.38	1.51
1	C	132	VAL	CB-CG1	5.60	1.64	1.52
1	A	20	HIS	CA-CB	-5.58	1.41	1.53
2	D	67	VAL	CB-CG1	5.50	1.64	1.52
1	C	132	VAL	CB-CG2	5.49	1.64	1.52
1	C	87	HIS	CA-CB	5.10	1.65	1.53
1	A	78	ASN	CB-CG	5.02	1.62	1.51
1	C	20	HIS	CA-CB	-5.02	1.43	1.53

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	2	LEU	CA-C-N	-9.15	97.06	117.20
2	B	30	ARG	NE-CZ-NH2	-8.77	115.92	120.30
2	B	30	ARG	NE-CZ-NH1	7.71	124.16	120.30
2	D	1	VAL	CB-CA-C	6.86	124.43	111.40
1	A	21	ALA	CB-CA-C	6.73	120.19	110.10
2	B	82	LYS	CD-CE-NZ	-6.59	96.54	111.70
1	A	141	ARG	NE-CZ-NH1	6.54	123.57	120.30
2	D	40	ARG	NE-CZ-NH1	6.40	123.50	120.30
1	C	94	ASP	CB-CG-OD1	6.29	123.96	118.30
1	A	2	LEU	CB-CA-C	6.25	122.07	110.20
1	C	29	LEU	CB-CG-CD1	-6.05	100.72	111.00
1	A	91	LEU	CB-CG-CD2	-5.75	101.22	111.00
1	A	92	ARG	NE-CZ-NH1	5.72	123.16	120.30
1	C	31	ARG	NE-CZ-NH1	5.68	123.14	120.30
1	C	38	THR	OG1-CB-CG2	5.56	122.78	110.00
2	D	96	LEU	CB-CG-CD2	-5.48	101.69	111.00
1	A	3	SER	N-CA-CB	-5.48	102.29	110.50
2	D	40	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	A	106	LEU	CB-CG-CD2	5.39	120.17	111.00
1	C	34	LEU	CB-CG-CD2	-5.35	101.90	111.00
2	D	47	ASP	CB-CA-C	5.29	120.97	110.40
2	D	96	LEU	CA-CB-CG	-5.28	103.16	115.30
2	D	146	HIS	N-CA-C	5.26	125.20	111.00
1	A	2	LEU	N-CA-CB	5.23	120.87	110.40
2	B	20	VAL	CB-CA-C	-5.22	101.49	111.40
1	C	141	ARG	CB-CA-C	5.20	120.80	110.40
1	A	72	HIS	CB-CA-C	-5.12	100.16	110.40
1	A	141	ARG	NE-CZ-NH2	-5.08	117.76	120.30
1	A	7	LYS	CB-CA-C	-5.07	100.27	110.40
2	B	54	VAL	CB-CA-C	-5.05	101.80	111.40

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	C	38	THR	CB
2	D	1	VAL	CA
2	D	146	HIS	CA

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	2	LEU	Mainchain
2	B	139	ASN	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	1069	0	1073	33	2
1	C	1069	0	1073	26	0
2	B	1123	0	1116	30	1
2	D	1123	0	1113	38	0
3	A	43	0	30	0	0
3	B	43	0	30	0	0
3	C	43	0	30	1	0
3	D	43	0	30	2	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	D	12	0	3	0	0
6	A	102	0	0	10	2
6	B	118	0	0	7	0
6	C	99	0	0	5	1
6	D	120	0	0	10	1
All	All	5015	0	4498	112	4

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 12.

All (112) 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:A:92:ARG:HH11	1:A:92:ARG:HG3	1.07	1.07
1:A:127:LYS:HE3	1:C:141:ARG:HD2	1.53	0.90
1:A:92:ARG:NH1	1:A:92:ARG:HG3	1.80	0.89
2:D:79:ASP:HB3	6:D:352:HOH:O	1.73	0.88
2:B:119:GLY:HA3	2:B:120:LYS:HE2	1.56	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:127:LYS:HE3	1:C:141:ARG:CD	2.07	0.85
6:A:272:HOH:O	2:B:124:PRO:HA	1.75	0.84
2:B:132:LYS:HA	2:D:146:HIS:HE1	1.46	0.81
2:B:120:LYS:H	2:B:120:LYS:CD	1.95	0.78
2:D:120:LYS:HE3	6:D:386:HOH:O	1.86	0.75
1:A:31:ARG:HG2	6:A:272:HOH:O	1.87	0.74
2:D:97:HIS:ND1	6:D:350:HOH:O	2.22	0.73
2:D:117:HIS:NE2	6:D:356:HOH:O	2.22	0.73
1:C:103:HIS:HE1	2:D:131:GLN:OE1	1.71	0.72
2:B:51:PRO:O	2:B:55:MET:HG2	1.90	0.71
2:D:61:LYS:HB3	6:D:331:HOH:O	1.91	0.71
1:A:49:SER:O	1:A:52:SER:HB3	1.91	0.71
2:B:132:LYS:HA	2:D:146:HIS:CE1	2.27	0.70
2:B:120:LYS:HE2	2:B:120:LYS:H	1.56	0.70
1:A:31:ARG:NH1	6:A:272:HOH:O	2.26	0.69
1:A:141:ARG:C	1:C:127:LYS:HD3	2.13	0.68
2:B:120:LYS:CE	2:B:120:LYS:H	2.07	0.66
1:C:112:HIS:O	1:C:113:LEU:HD23	1.96	0.65
2:D:55:MET:HE3	6:D:405:HOH:O	1.95	0.65
2:B:143:HIS:CD2	6:B:212:HOH:O	2.50	0.64
2:B:57:ASN:ND2	6:B:285:HOH:O	2.28	0.64
1:A:103:HIS:HE1	2:B:131:GLN:OE1	1.81	0.64
2:B:143:HIS:HB3	6:B:239:HOH:O	1.98	0.62
2:B:146:HIS:OXT	2:D:1:VAL:HG22	2.00	0.62
1:C:73:VAL:HG12	1:C:74:ASP:OD1	2.00	0.62
2:D:93:CYS:HB2	2:D:145:TYR:CE1	2.34	0.61
1:A:112:HIS:O	1:A:113:LEU:HD23	2.01	0.61
2:D:84:THR:HG23	6:D:353:HOH:O	2.01	0.60
2:B:75:LEU:HD21	2:B:133:VAL:HG11	1.85	0.59
2:D:144:LYS:N	2:D:144:LYS:HZ3	2.00	0.59
1:A:108:THR:O	1:A:112:HIS:HD2	1.87	0.58
2:D:141:LEU:CD1	3:D:200:HEM:CBB	2.83	0.57
2:B:104:ARG:NH1	6:B:282:HOH:O	2.38	0.57
2:D:144:LYS:HA	2:D:144:LYS:NZ	2.19	0.56
2:D:120:LYS:N	2:D:120:LYS:HD3	2.20	0.56
1:A:103:HIS:HD2	6:A:227:HOH:O	1.89	0.56
1:C:7:LYS:HG2	1:C:73:VAL:HG11	1.86	0.56
2:D:91:LEU:HD22	2:D:91:LEU:O	2.05	0.56
2:D:73:ASP:O	2:D:77:HIS:HD2	1.89	0.56
2:D:108:ASN:ND2	6:D:344:HOH:O	2.38	0.55
1:C:86:LEU:CD2	3:C:200:HEM:HBA2	2.38	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:143:HIS:C	2:D:144:LYS:HZ3	2.12	0.53
2:B:43:GLU:OE2	1:C:92:ARG:NH2	2.34	0.53
2:D:3:LEU:HA	2:D:7:GLU:OE1	2.10	0.52
2:B:144:LYS:HA	2:B:144:LYS:HE2	1.90	0.52
6:B:202:HOH:O	1:C:38:THR:HG23	2.09	0.52
2:D:141:LEU:HD12	3:D:200:HEM:CBB	2.40	0.51
2:D:84:THR:HG22	2:D:85:PHE:CD2	2.46	0.51
1:A:34:LEU:HD13	6:A:296:HOH:O	2.11	0.51
1:C:2:LEU:HD23	1:C:2:LEU:N	2.27	0.50
6:A:272:HOH:O	2:B:127:GLN:HB3	2.12	0.50
6:B:202:HOH:O	1:C:38:THR:CG2	2.58	0.50
2:D:89:SER:OG	2:D:144:LYS:HG2	2.12	0.50
2:B:120:LYS:HE2	2:B:120:LYS:N	2.26	0.50
1:A:70:VAL:O	1:A:73:VAL:HG22	2.12	0.50
2:D:144:LYS:HZ2	2:D:144:LYS:HA	1.76	0.49
1:A:89:HIS:NE2	1:A:139:LYS:HG3	2.28	0.49
2:B:119:GLY:CA	2:B:120:LYS:HE2	2.36	0.48
1:C:73:VAL:O	1:C:73:VAL:HG13	2.13	0.48
1:A:17:VAL:HG13	1:A:24:TYR:CD2	2.48	0.47
1:C:38:THR:HG22	6:C:266:HOH:O	2.14	0.47
2:D:100:PRO:HB2	2:D:104:ARG:HH12	1.81	0.46
2:D:91:LEU:HD22	2:D:91:LEU:C	2.36	0.46
2:B:73:ASP:O	2:B:77:HIS:HD2	1.99	0.46
2:B:145:TYR:O	2:B:146:HIS:HB3	2.16	0.46
2:B:65:LYS:HE3	2:B:65:LYS:HB2	1.61	0.46
1:A:56:LYS:HG3	6:A:263:HOH:O	2.15	0.45
1:C:103:HIS:CE1	2:D:131:GLN:OE1	2.61	0.45
1:A:41:THR:HG21	2:D:40:ARG:HH12	1.82	0.44
1:A:41:THR:CG2	2:D:40:ARG:HH12	2.30	0.44
1:A:89:HIS:NE2	1:A:139:LYS:CG	2.80	0.44
2:D:37:TRP:HE1	2:D:102:ASN:HD21	1.64	0.44
1:A:139:LYS:HA	1:A:139:LYS:HD2	1.67	0.43
1:A:16:LYS:H	1:A:16:LYS:HG2	1.02	0.43
1:C:89:HIS:CE1	1:C:139:LYS:HD2	2.53	0.43
2:D:96:LEU:O	2:D:97:HIS:HB2	2.19	0.43
1:A:135:VAL:HG23	6:A:250:HOH:O	2.18	0.43
1:A:44:PRO:HG2	1:A:45:HIS:N	2.34	0.43
1:C:114:PRO:HA	2:D:116:HIS:CD2	2.53	0.43
1:A:44:PRO:HG2	1:A:45:HIS:H	1.83	0.43
1:C:89:HIS:CE1	1:C:139:LYS:CD	3.01	0.43
1:A:76:MET:HE2	6:A:292:HOH:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:37:TRP:HE1	2:B:102:ASN:HD21	1.65	0.43
1:A:114:PRO:HA	2:B:116:HIS:CD2	2.54	0.42
2:D:88:LEU:HD12	2:D:88:LEU:HA	1.85	0.42
2:B:73:ASP:O	2:B:77:HIS:CD2	2.73	0.42
1:A:30:GLU:OE2	1:A:50:HIS:CD2	2.72	0.42
2:D:84:THR:CA	6:D:353:HOH:O	2.67	0.42
2:D:144:LYS:CA	2:D:144:LYS:NZ	2.83	0.42
2:D:146:HIS:HB2	6:D:355:HOH:O	2.19	0.42
1:C:127:LYS:HE2	6:C:290:HOH:O	2.19	0.41
1:A:45:HIS:N	1:A:45:HIS:ND1	2.66	0.41
1:C:9:ASN:HD22	1:C:9:ASN:HA	1.60	0.41
2:B:75:LEU:HA	2:B:75:LEU:HD23	1.86	0.41
1:C:116:GLU:HB2	6:C:217:HOH:O	2.20	0.41
2:D:124:PRO:HB2	2:D:125:PRO:HD3	2.02	0.41
1:A:3:SER:HA	1:A:4:PRO:HD3	1.92	0.41
1:C:89:HIS:HB3	6:C:208:HOH:O	2.20	0.41
2:B:15:TRP:HD1	6:B:314:HOH:O	2.03	0.41
1:C:2:LEU:HA	1:C:2:LEU:HD22	1.84	0.41
1:A:109:LEU:HD12	1:A:125:LEU:HD13	2.03	0.41
2:B:78:LEU:HA	2:B:78:LEU:HD23	1.85	0.41
1:A:127:LYS:NZ	6:A:253:HOH:O	2.29	0.40
1:C:72:HIS:HE1	6:C:262:HOH:O	2.04	0.40
1:C:76:MET:HB2	1:C:77:PRO:HD3	2.04	0.40
1:A:127:LYS:HE3	1:C:141:ARG:HD3	1.99	0.40
2:B:88:LEU:HA	2:B:88:LEU:HD12	1.92	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:8:THR:CG2	6:A:267:HOH:O[4_555]	0.71	1.49
6:D:408:HOH:O	6:D:408:HOH:O[2_655]	1.80	0.40
1:A:8:THR:CB	6:A:267:HOH:O[4_555]	1.95	0.25
2:B:6:GLU:OE1	6:C:228:HOH:O[3_445]	2.13	0.07

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	139/141 (99%)	129 (93%)	9 (6%)	1 (1%)	26	11
1	C	139/141 (99%)	138 (99%)	1 (1%)	0	100	100
2	B	144/146 (99%)	142 (99%)	2 (1%)	0	100	100
2	D	144/146 (99%)	142 (99%)	2 (1%)	0	100	100
All	All	566/574 (99%)	551 (97%)	14 (2%)	1 (0%)	52	35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	52	SER

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%)	101 (89%)	12 (11%)	8	2
1	C	113/113 (100%)	99 (88%)	14 (12%)	6	1
2	B	118/118 (100%)	106 (90%)	12 (10%)	9	2
2	D	118/118 (100%)	105 (89%)	13 (11%)	8	1
All	All	462/462 (100%)	411 (89%)	51 (11%)	8	1

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	9	ASN
1	A	16	LYS
1	A	52	SER
1	A	60	LYS
1	A	73	VAL
1	A	78	ASN
1	A	84	SER
1	A	90	LYS
1	A	92	ARG
1	A	119	PRO
1	A	141	ARG
2	B	65	LYS
2	B	68	LEU
2	B	73	ASP
2	B	81	LEU
2	B	87	THR
2	B	88	LEU
2	B	95	LYS
2	B	104	ARG
2	B	120	LYS
2	B	139	ASN
2	B	143	HIS
2	B	146	HIS
1	C	2	LEU
1	C	16	LYS
1	C	38	THR
1	C	41	THR
1	C	60	LYS
1	C	73	VAL
1	C	78	ASN
1	C	83	LEU
1	C	84	SER
1	C	85	ASP
1	C	90	LYS
1	C	114	PRO
1	C	139	LYS
1	C	141	ARG
2	D	8	LYS
2	D	47	ASP
2	D	49	SER
2	D	51	PRO
2	D	82	LYS

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Mol	Chain	Res	Type
2	D	84	THR
2	D	88	LEU
2	D	90	GLU
2	D	91	LEU
2	D	104	ARG
2	D	120	LYS
2	D	143	HIS
2	D	144	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	50	HIS
1	A	97	ASN
1	A	103	HIS
2	B	77	HIS
2	B	102	ASN
2	B	108	ASN
1	C	9	ASN
1	C	68	ASN
1	C	97	ASN
1	C	103	HIS
2	D	77	HIS
2	D	80	ASN
2	D	97	HIS
2	D	102	ASN
2	D	108	ASN
2	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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

9 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	200	1,4	30,50,50	2.54	6 (20%)	24,82,82	2.73	11 (45%)
4	CMO	A	201	3	0,1,1	0.00	-	0,0,0	0.00	-
3	HEM	B	200	2,4	30,50,50	3.13	8 (26%)	24,82,82	3.05	12 (50%)
4	CMO	B	201	3	0,1,1	0.00	-	0,0,0	0.00	-
3	HEM	C	200	1,4	30,50,50	2.75	10 (33%)	24,82,82	3.27	12 (50%)
4	CMO	C	201	3	0,1,1	0.00	-	0,0,0	0.00	-
3	HEM	D	200	2,4	30,50,50	2.27	8 (26%)	24,82,82	3.15	14 (58%)
4	CMO	D	201	3	0,1,1	0.00	-	0,0,0	0.00	-
5	TMM	D	300	2	12,12,15	1.14	0	15,15,21	1.10	2 (13%)

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	200	1,4	-	0/10/54/54	0/0/8/8
4	CMO	A	201	3	-	0/0/0/0	0/0/0/0
3	HEM	B	200	2,4	-	0/10/54/54	0/0/8/8
4	CMO	B	201	3	-	0/0/0/0	0/0/0/0
3	HEM	C	200	1,4	-	0/10/54/54	0/0/8/8
4	CMO	C	201	3	-	0/0/0/0	0/0/0/0
3	HEM	D	200	2,4	-	0/10/54/54	0/0/8/8
4	CMO	D	201	3	-	0/0/0/0	0/0/0/0
5	TMM	D	300	2	-	0/6/6/12	0/1/1/1

All (32) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	200	HEM	C3B-C4B	-11.53	1.41	1.51
3	A	200	HEM	C2D-C3D	-7.72	1.31	1.54
3	B	200	HEM	C3D-C4D	-7.56	1.41	1.51
3	A	200	HEM	C3B-C4B	-7.55	1.45	1.51
3	D	200	HEM	C2D-C3D	-7.44	1.32	1.54
3	C	200	HEM	C2D-C3D	-7.32	1.32	1.54
3	C	200	HEM	C3B-C4B	-6.88	1.45	1.51
3	B	200	HEM	C2D-C3D	-6.46	1.35	1.54
3	C	200	HEM	C3D-C4D	-6.42	1.43	1.51
3	D	200	HEM	C3D-C4D	-6.01	1.43	1.51
3	A	200	HEM	C3D-C4D	-5.16	1.45	1.51
3	A	200	HEM	C2C-C1C	-4.09	1.44	1.52
3	B	200	HEM	C2C-C1C	-4.00	1.45	1.52
3	D	200	HEM	C2C-C1C	-3.52	1.45	1.52
3	D	200	HEM	C3B-C4B	-3.31	1.48	1.51
3	C	200	HEM	C2C-C1C	-3.11	1.46	1.52
3	B	200	HEM	C2B-C1B	-2.99	1.42	1.51
3	C	200	HEM	C2D-C1D	-2.49	1.43	1.51
3	A	200	HEM	C2B-C1B	-2.48	1.43	1.51
3	B	200	HEM	C3C-CAC	2.02	1.55	1.51
3	C	200	HEM	C3C-CAC	2.03	1.55	1.51
3	D	200	HEM	C1C-NC	2.05	1.38	1.36
3	C	200	HEM	FE-NC	2.26	2.04	1.95
3	B	200	HEM	FE-NC	2.31	2.04	1.95
3	D	200	HEM	CMA-C3A	2.53	1.56	1.51
3	D	200	HEM	C3B-CAB	2.58	1.56	1.51
3	B	200	HEM	C3B-CAB	2.62	1.56	1.51
3	D	200	HEM	FE-NC	2.66	2.06	1.95
3	C	200	HEM	C4C-NC	2.98	1.39	1.36
3	C	200	HEM	CMA-C3A	3.09	1.58	1.51
3	A	200	HEM	CAA-C2A	3.16	1.57	1.52
3	C	200	HEM	C1C-NC	4.10	1.41	1.36

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	200	HEM	C3B-CAB-CBB	-8.21	111.87	124.46
3	C	200	HEM	CAA-C2A-C1A	-5.41	121.14	127.01
3	A	200	HEM	C3B-CAB-CBB	-4.71	117.23	124.46
3	B	200	HEM	CAA-C2A-C1A	-4.24	122.40	127.01
3	D	200	HEM	CBA-CAA-C2A	-4.20	105.01	112.53
3	D	200	HEM	C3C-CAC-CBC	-4.02	118.29	124.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	200	HEM	CMA-C3A-C4A	-3.64	122.34	128.36
3	B	200	HEM	CBA-CAA-C2A	-3.64	106.00	112.53
3	A	200	HEM	CMA-C3A-C4A	-3.40	122.74	128.36
3	D	200	HEM	CAA-C2A-C1A	-3.37	123.35	127.01
3	B	200	HEM	CBD-CAD-C3D	-2.98	104.87	113.55
3	B	200	HEM	C3B-CAB-CBB	-2.59	120.49	124.46
3	D	200	HEM	CBD-CAD-C3D	-2.53	106.19	113.55
5	D	300	TMM	C6-C5-C5A	-2.31	115.35	119.91
3	A	200	HEM	C3B-C4B-NB	-2.27	107.29	111.63
3	A	200	HEM	CBD-CAD-C3D	-2.20	107.16	113.55
3	C	200	HEM	C1D-CHD-C4C	-2.13	122.27	125.82
3	C	200	HEM	CBD-CAD-C3D	-2.08	107.49	113.55
3	C	200	HEM	C3B-C4B-NB	-2.04	107.73	111.63
3	B	200	HEM	C3C-CAC-CBC	-2.00	121.39	124.46
3	D	200	HEM	CMA-C3A-C2A	2.02	129.45	125.24
3	D	200	HEM	C2C-C1C-CHC	2.27	127.14	123.68
3	A	200	HEM	C3B-C4B-CHC	2.27	126.37	123.16
3	C	200	HEM	C3B-C4B-CHC	2.28	126.37	123.16
5	D	300	TMM	C4-C5-C5A	2.28	124.40	119.91
3	D	200	HEM	CAD-C3D-C4D	2.50	121.29	112.47
3	C	200	HEM	CAD-C3D-C4D	2.57	121.55	112.47
3	D	200	HEM	CMD-C2D-C3D	2.67	126.16	114.35
3	A	200	HEM	C2D-C3D-C4D	2.94	106.49	101.50
3	A	200	HEM	CMD-C2D-C3D	3.04	127.78	114.35
3	B	200	HEM	CMA-C3A-C2A	3.05	131.61	125.24
3	B	200	HEM	CMD-C2D-C3D	3.13	128.21	114.35
3	D	200	HEM	C4B-CHC-C1C	3.70	132.01	125.82
3	D	200	HEM	C3B-C4B-CHC	3.76	128.46	123.16
3	C	200	HEM	CMD-C2D-C3D	3.81	131.21	114.35
3	A	200	HEM	CAD-C3D-C2D	3.88	124.38	113.22
3	A	200	HEM	CMC-C2C-C3C	4.69	128.24	116.53
3	A	200	HEM	CAD-C3D-C4D	4.69	129.01	112.47
3	C	200	HEM	CMC-C2C-C3C	4.71	128.28	116.53
3	D	200	HEM	C2D-C3D-C4D	4.73	109.52	101.50
3	C	200	HEM	CAD-C3D-C2D	4.75	126.86	113.22
3	B	200	HEM	C2D-C3D-C4D	4.90	109.81	101.50
3	D	200	HEM	CAD-C3D-C2D	5.31	128.50	113.22
3	C	200	HEM	C2D-C3D-C4D	5.45	110.73	101.50
3	D	200	HEM	CMB-C2B-C3B	5.53	130.35	116.53
3	B	200	HEM	CMC-C2C-C3C	5.63	130.59	116.53
3	C	200	HEM	CMB-C2B-C3B	5.99	131.48	116.53
3	B	200	HEM	CAD-C3D-C2D	6.10	130.75	113.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	200	HEM	CMB-C2B-C3B	6.18	131.96	116.53
3	A	200	HEM	CMB-C2B-C3B	6.41	132.54	116.53
3	D	200	HEM	CMC-C2C-C3C	6.78	133.45	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	200	HEM	1	0
3	D	200	HEM	2	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.