



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 04:07 pm GMT

PDB ID : 1GBV
Title : (ALPHA-OXY, BETA-(C112G)DEOXY) T-STATE HUMAN
HEMOGLOBIN
Authors : Vasquez, G.B.; Ji, X.; Pechik, I.; Fronticelli, C.; Gilliland, G.L.
Deposited on : 1995-12-20
Resolution : 2.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7.2 (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

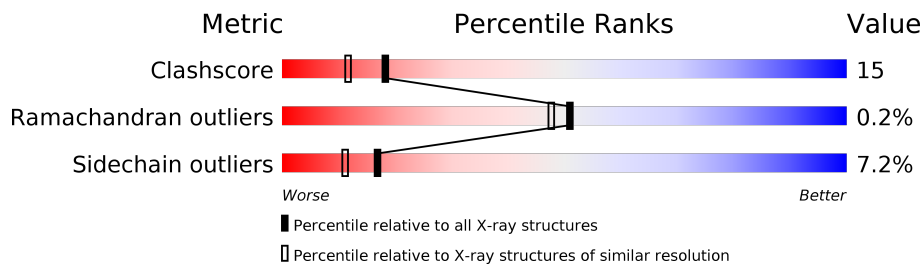
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	112137	7775 (2.00-2.00)
Ramachandran outliers	110173	7679 (2.00-2.00)
Sidechain outliers	110143	7678 (2.00-2.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	141	
1	C	141	
2	B	146	
2	D	146	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4923 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.

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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1121	723	195	201	2			
2	D	146	Total	C	N	O	S	0	0	0
			1121	723	195	201	2			

There are 2 discrepancies between the modelled and reference sequences:

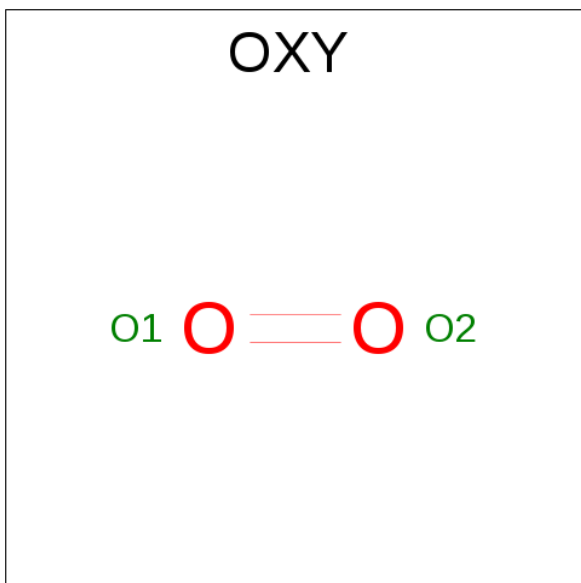
Chain	Residue	Modelled	Actual	Comment	Reference
B	112	GLY	CYS	ENGINEERED	UNP P68871
D	112	GLY	CYS	ENGINEERED	UNP P68871

- 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 OXYGEN MOLECULE (three-letter code: OXY) (formula: O_2).



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

- Molecule 5 is water.

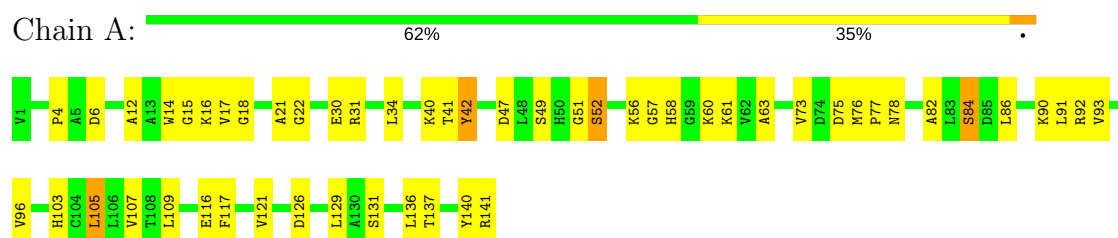
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	98	Total	O	0	0
			98	98		
5	B	94	Total	O	0	0
			94	94		
5	C	95	Total	O	0	0
			95	95		
5	D	80	Total	O	0	0
			80	80		

3 Residue-property plots [i](#)

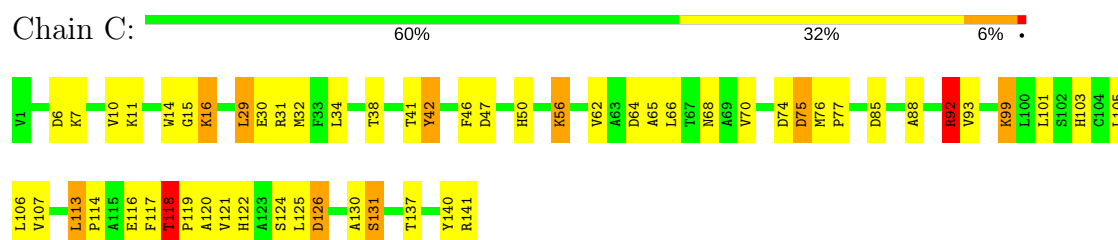
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are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

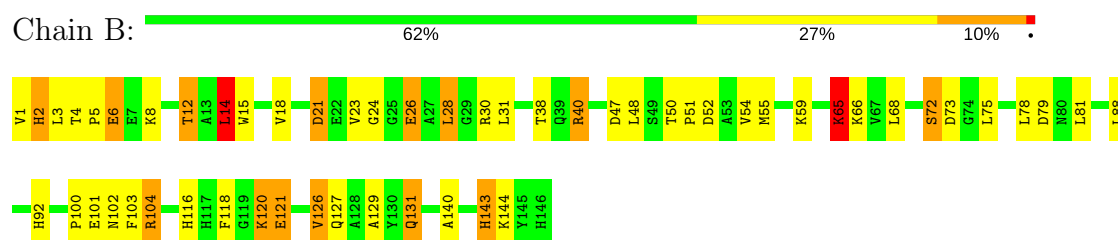
• Molecule 1: HEMOGLOBIN



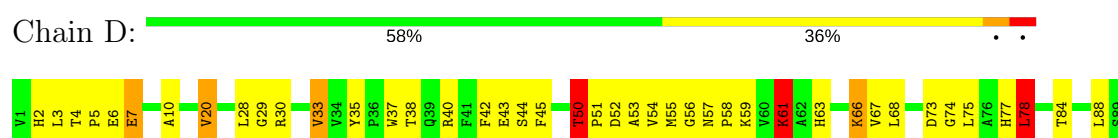
• Molecule 1: HEMOGLOBIN



• Molecule 2: HEMOGLOBIN



• Molecule 2: HEMOGLOBIN



E90	L91	H92	C93	D94	K95	Y98	E101	N102	F103	R104	L105	H116	P124	P125	V126	A129	Y130	Q131	N139	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	63.18Å 83.56Å 53.73Å 90.00° 99.60° 90.00°	Depositor
Resolution (Å)	6.00 – 2.00	Depositor
% Data completeness (in resolution range)	92.8 (6.00-2.00)	Depositor
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	GPRLSA, PROLSQ	Depositor
R, R_{free}	0.184 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4923	wwPDB-VP
Average B, all atoms (Å ²)	19.0	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: HEM, OXY

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	1.08	0/1097	1.78	22/1491 (1.5%)
1	C	1.12	0/1097	1.92	30/1491 (2.0%)
2	B	1.06	0/1151	1.80	22/1563 (1.4%)
2	D	1.09	0/1151	1.75	18/1563 (1.2%)
All	All	1.09	0/4496	1.81	92/6108 (1.5%)

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	C	0	3
2	B	0	3
2	D	0	1
All	All	0	7

There are no bond length outliers.

All (92) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	30	ARG	NE-CZ-NH1	15.31	127.96	120.30
1	C	6	ASP	CB-CG-OD2	-14.37	105.37	118.30
1	C	141	ARG	NE-CZ-NH1	13.21	126.91	120.30
1	C	105	LEU	CA-CB-CG	13.20	145.66	115.30
2	B	121	GLU	OE1-CD-OE2	12.15	137.88	123.30
1	A	47	ASP	CB-CG-OD1	10.74	127.96	118.30
1	A	31	ARG	NE-CZ-NH2	9.68	125.14	120.30
2	B	26	GLU	OE1-CD-OE2	9.38	134.55	123.30
1	C	42	TYR	CB-CG-CD1	-9.30	115.42	121.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	40	ARG	NE-CZ-NH2	-9.08	115.76	120.30
1	C	141	ARG	NE-CZ-NH2	-8.77	115.91	120.30
1	A	42	TYR	CB-CG-CD1	-8.61	115.84	121.00
1	A	126	ASP	CB-CG-OD2	-8.34	110.80	118.30
1	C	118	THR	CA-CB-CG2	8.32	124.05	112.40
2	B	30	ARG	CD-NE-CZ	8.31	135.24	123.60
2	D	104	ARG	NE-CZ-NH2	-8.27	116.16	120.30
1	C	92	ARG	NE-CZ-NH1	8.09	124.35	120.30
1	A	116	GLU	CG-CD-OE1	8.05	134.40	118.30
1	A	6	ASP	CB-CG-OD2	-8.03	111.07	118.30
1	C	126	ASP	CB-CA-C	7.98	126.36	110.40
1	C	92	ARG	NE-CZ-NH2	-7.91	116.34	120.30
2	D	73	ASP	CB-CG-OD1	7.86	125.37	118.30
1	C	64	ASP	CB-CG-OD2	-7.72	111.35	118.30
1	A	96	VAL	CA-CB-CG2	7.62	122.33	110.90
2	B	73	ASP	CB-CG-OD1	7.55	125.10	118.30
1	C	75	ASP	CB-CG-OD1	7.54	125.09	118.30
1	C	64	ASP	CB-CG-OD1	7.45	125.01	118.30
1	A	92	ARG	NE-CZ-NH2	-6.99	116.80	120.30
1	A	6	ASP	O-C-N	6.77	133.53	122.70
1	C	140	TYR	CB-CG-CD2	-6.73	116.96	121.00
2	B	14	LEU	CA-CB-CG	6.71	130.72	115.30
2	D	40	ARG	NE-CZ-NH2	-6.67	116.97	120.30
2	D	78	LEU	CA-CB-CG	6.48	130.20	115.30
1	C	75	ASP	CB-CG-OD2	-6.47	112.48	118.30
1	A	6	ASP	N-CA-CB	6.43	122.18	110.60
1	C	113	LEU	CA-CB-CG	6.32	129.84	115.30
2	B	30	ARG	NH1-CZ-NH2	-6.22	112.56	119.40
1	C	6	ASP	OD1-CG-OD2	6.19	135.06	123.30
1	C	130	ALA	CB-CA-C	6.13	119.30	110.10
1	C	42	TYR	CB-CG-CD2	6.12	124.67	121.00
1	C	131	SER	N-CA-CB	6.02	119.53	110.50
1	A	121	VAL	CB-CA-C	5.98	122.77	111.40
2	B	65	LYS	CA-CB-CG	5.96	126.50	113.40
1	C	126	ASP	CB-CG-OD2	5.96	123.66	118.30
2	B	72	SER	CB-CA-C	5.90	121.31	110.10
2	D	84	THR	CA-CB-CG2	5.88	120.63	112.40
1	C	88	ALA	O-C-N	5.78	131.94	122.70
2	B	28	LEU	CB-CA-C	5.77	121.17	110.20
2	D	101	GLU	OE1-CD-OE2	5.74	130.19	123.30
2	B	126	VAL	CB-CA-C	5.73	122.29	111.40
2	B	104	ARG	NE-CZ-NH1	5.71	123.16	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	35	TYR	CB-CG-CD2	-5.70	117.58	121.00
2	B	26	GLU	CG-CD-OE1	-5.70	106.90	118.30
2	B	52	ASP	CB-CG-OD1	5.69	123.42	118.30
2	D	50	THR	CA-CB-CG2	5.69	120.36	112.40
1	C	31	ARG	NE-CZ-NH1	5.68	123.14	120.30
2	D	20	VAL	CB-CA-C	5.64	122.12	111.40
1	C	105	LEU	CB-CG-CD1	5.63	120.58	111.00
2	D	30	ARG	NE-CZ-NH2	-5.56	117.52	120.30
2	B	143	HIS	N-CA-CB	5.56	120.60	110.60
2	D	30	ARG	CD-NE-CZ	5.51	131.31	123.60
1	C	41	THR	CA-CB-CG2	5.46	120.04	112.40
1	A	141	ARG	NE-CZ-NH2	-5.42	117.59	120.30
1	A	30	GLU	CA-CB-CG	5.41	125.31	113.40
2	B	79	ASP	CB-CG-OD1	-5.40	113.44	118.30
1	A	116	GLU	CG-CD-OE2	-5.39	107.52	118.30
1	A	105	LEU	CA-CB-CG	5.38	127.66	115.30
2	D	130	TYR	CB-CG-CD1	-5.35	117.79	121.00
1	A	126	ASP	CB-CG-OD1	5.34	123.11	118.30
2	B	21	ASP	CB-CG-OD1	5.32	123.09	118.30
1	C	137	THR	CA-CB-CG2	5.32	119.84	112.40
2	B	129	ALA	O-C-N	5.25	131.11	122.70
2	D	61	LYS	N-CA-CB	5.25	120.05	110.60
1	A	12	ALA	N-CA-CB	5.23	117.42	110.10
1	A	141	ARG	NE-CZ-NH1	5.23	122.91	120.30
1	C	66	LEU	CB-CA-C	5.21	120.10	110.20
1	C	47	ASP	N-CA-C	-5.18	97.01	111.00
2	D	33	VAL	CA-CB-CG1	5.15	118.63	110.90
1	C	34	LEU	O-C-N	5.15	130.93	122.70
1	A	47	ASP	N-CA-C	-5.14	97.13	111.00
1	A	47	ASP	O-C-N	5.14	130.92	122.70
1	C	46	PHE	O-C-N	5.13	130.91	122.70
1	A	6	ASP	OD1-CG-OD2	5.12	133.03	123.30
1	C	29	LEU	CA-C-O	5.12	130.86	120.10
2	B	131	GLN	N-CA-CB	5.12	119.81	110.60
2	B	104	ARG	NH1-CZ-NH2	-5.10	113.79	119.40
1	A	107	VAL	CA-CB-CG1	5.09	118.54	110.90
2	B	103	PHE	O-C-N	5.09	130.85	122.70
2	D	130	TYR	CB-CG-CD2	5.09	124.06	121.00
2	D	93	CYS	CA-CB-SG	-5.07	104.88	114.00
2	D	7	GLU	OE1-CD-OE2	5.06	129.37	123.30
2	D	105	LEU	N-CA-CB	5.04	120.48	110.40

There are no chirality outliers.

All (7) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	104	ARG	Sidechain,Mainchain
2	B	126	VAL	Mainchain
1	C	101	LEU	Mainchain
1	C	120	ALA	Mainchain
1	C	92	ARG	Sidechain
2	D	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	1069	0	1073	29	0
1	C	1069	0	1073	29	0
2	B	1121	0	1116	35	0
2	D	1121	0	1116	39	0
3	A	43	0	30	4	0
3	B	43	0	30	4	0
3	C	43	0	30	2	0
3	D	43	0	30	3	0
4	A	2	0	0	0	0
4	C	2	0	0	0	0
5	A	98	0	0	4	0
5	B	94	0	0	2	0
5	C	95	0	0	8	0
5	D	80	0	0	2	0
All	All	4923	0	4498	136	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 15.

All (136) 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:118:THR:HG22	1:C:121:VAL:H	1.35	0.91
3:D:147:HEM:HH C	3:D:147:HEM:HBB2	1.55	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:B:147:HEM:HMB1	3:B:147:HEM:HBB2	1.62	0.82
2:D:104:ARG:HH21	2:D:139:ASN:HD21	1.26	0.81
2:D:75:LEU:O	2:D:78:LEU:HD22	1.87	0.73
2:B:24:GLY:HA2	2:B:68:LEU:HG	1.72	0.71
1:A:21:ALA:HB1	1:A:63:ALA:HB1	1.73	0.69
2:D:50:THR:HG22	2:D:53:ALA:H	1.59	0.68
1:A:76:MET:N	1:A:77:PRO:HD2	2.09	0.67
3:B:147:HEM:CMB	3:B:147:HEM:HBB2	2.23	0.67
2:D:104:ARG:HH21	2:D:139:ASN:ND2	1.92	0.67
1:C:7:LYS:HE3	1:C:74:ASP:OD1	1.95	0.66
2:B:50:THR:O	2:B:54:VAL:HG23	1.96	0.66
1:C:16:LYS:N	1:C:16:LYS:HE3	2.11	0.65
1:C:113:LEU:HB3	1:C:116:GLU:HG2	1.80	0.64
2:B:143:HIS:CD2	2:B:144:LYS:HD3	2.33	0.63
1:C:118:THR:HG22	1:C:121:VAL:N	2.12	0.62
2:B:140:ALA:O	2:B:143:HIS:ND1	2.31	0.60
1:A:73:VAL:HG23	1:A:76:MET:HE2	1.83	0.60
3:C:142:HEM:HBB2	3:C:142:HEM:HMB2	1.83	0.60
2:B:1:VAL:HB	2:B:2:HIS:CE1	2.38	0.58
3:C:142:HEM:CMB	3:C:142:HEM:HBB2	2.34	0.58
3:A:142:HEM:HMB2	3:A:142:HEM:HBB2	1.86	0.57
2:B:88:LEU:HD22	2:B:92:HIS:HE1	1.69	0.57
2:D:4:THR:O	2:D:7:GLU:N	2.37	0.57
1:A:51:GLY:O	1:A:56:LYS:HE2	2.05	0.56
2:B:51:PRO:O	2:B:55:MET:HG2	2.04	0.56
2:D:63:HIS:ND1	2:D:66:LYS:HD2	2.21	0.55
2:D:10:ALA:HB1	2:D:126:VAL:HG22	1.88	0.55
1:A:86:LEU:HD12	1:A:90:LYS:HB2	1.89	0.55
2:D:50:THR:HG22	2:D:53:ALA:HB2	1.88	0.55
1:A:84:SER:HB3	1:A:136:LEU:HA	1.90	0.54
1:C:106:LEU:HD21	1:C:126:ASP:HB2	1.90	0.54
1:A:57:GLY:O	1:A:61:LYS:HG3	2.08	0.54
2:D:63:HIS:O	2:D:67:VAL:HG23	2.08	0.54
2:D:57:ASN:O	2:D:61:LYS:HD3	2.08	0.53
2:D:42:PHE:O	2:D:45:PHE:HB2	2.09	0.53
2:D:63:HIS:O	2:D:66:LYS:HG3	2.08	0.53
2:D:50:THR:O	2:D:54:VAL:HG23	2.09	0.52
1:A:73:VAL:HG23	1:A:76:MET:CE	2.39	0.52
2:B:21:ASP:CG	2:B:65:LYS:HD3	2.31	0.51
1:C:68:ASN:HB2	5:C:203:HOH:O	2.11	0.51
1:A:61:LYS:HE3	5:A:192:HOH:O	2.09	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76:MET:N	1:C:77:PRO:CD	2.74	0.51
2:D:38:THR:HG22	2:D:102:ASN:OD1	2.11	0.51
1:C:15:GLY:C	1:C:16:LYS:HE3	2.30	0.51
1:A:16:LYS:NZ	5:A:229:HOH:O	2.43	0.51
1:C:99:LYS:HD2	5:C:153:HOH:O	2.09	0.50
2:D:104:ARG:NH2	2:D:139:ASN:OD1	2.44	0.50
1:A:77:PRO:HG2	1:A:78:ASN:H	1.75	0.50
2:D:57:ASN:OD1	2:D:59:LYS:HB2	2.12	0.50
2:D:3:LEU:HD13	2:D:7:GLU:HB3	1.93	0.50
2:B:8:LYS:O	2:B:12:THR:HG23	2.11	0.50
2:B:4:THR:OG1	2:B:6:GLU:HG2	2.12	0.49
2:D:51:PRO:O	2:D:55:MET:HG2	2.11	0.49
1:C:92:ARG:HD2	5:C:196:HOH:O	2.13	0.49
2:B:1:VAL:HG13	2:B:81:LEU:HD12	1.94	0.49
1:A:117:PHE:HB3	2:B:116:HIS:CD2	2.47	0.49
1:A:14:TRP:O	1:A:15:GLY:C	2.51	0.49
2:D:50:THR:HG22	2:D:53:ALA:CB	2.42	0.49
2:D:88:LEU:HG	5:D:150:HOH:O	2.12	0.49
1:C:85:ASP:HB3	5:C:171:HOH:O	2.12	0.48
1:A:14:TRP:O	1:A:17:VAL:HB	2.13	0.48
2:B:38:THR:HG22	2:B:102:ASN:OD1	2.14	0.47
1:C:117:PHE:HB3	2:D:116:HIS:CD2	2.48	0.47
2:D:52:ASP:O	2:D:56:GLY:N	2.43	0.47
1:C:103:HIS:O	1:C:107:VAL:HG23	2.15	0.47
2:D:104:ARG:NH2	2:D:139:ASN:ND2	2.61	0.46
2:B:102:ASN:N	5:C:177:HOH:O	2.48	0.46
1:A:61:LYS:NZ	3:A:142:HEM:O1A	2.47	0.46
1:C:30:GLU:OE2	1:C:50:HIS:ND1	2.49	0.46
3:A:142:HEM:CMB	3:A:142:HEM:HBB2	2.46	0.46
1:A:34:LEU:HD13	5:B:196:HOH:O	2.16	0.46
2:B:127:GLN:O	2:B:131:GLN:HG2	2.16	0.46
1:C:122:HIS:O	1:C:122:HIS:HD2	1.99	0.45
2:B:75:LEU:O	2:B:78:LEU:HG	2.17	0.45
2:B:1:VAL:HG13	2:B:81:LEU:CD1	2.46	0.45
1:C:119:PRO:HD3	5:D:195:HOH:O	2.15	0.45
1:A:42:TYR:CE2	1:A:93:VAL:HA	2.52	0.45
1:C:62:VAL:O	1:C:65:ALA:HB3	2.17	0.45
1:A:40:LYS:O	1:A:42:TYR:N	2.50	0.44
1:A:82:ALA:HB3	5:A:173:HOH:O	2.16	0.44
2:B:18:VAL:HG13	2:B:23:VAL:HG21	2.00	0.44
1:A:49:SER:O	1:A:52:SER:HB3	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:100:PRO:HD2	1:C:38:THR:HB	1.99	0.44
2:B:15:TRP:CZ2	2:B:72:SER:HB2	2.53	0.44
2:B:31:LEU:HD11	2:B:38:THR:HG21	1.98	0.44
2:D:124:PRO:N	2:D:125:PRO:HD2	2.32	0.44
1:C:76:MET:N	1:C:77:PRO:HD3	2.33	0.43
2:D:20:VAL:HG13	2:D:68:LEU:HB3	2.00	0.43
2:B:24:GLY:CA	2:B:68:LEU:HG	2.45	0.43
2:D:28:LEU:CD2	2:D:63:HIS:HB3	2.47	0.43
1:A:75:ASP:C	1:A:77:PRO:HD2	2.39	0.43
2:B:102:ASN:HB3	3:B:147:HEM:HMC1	2.01	0.43
2:D:29:GLY:O	2:D:33:VAL:HG23	2.18	0.43
2:D:91:LEU:CD1	2:D:95:LYS:HD2	2.49	0.43
3:B:147:HEM:HBB1	5:B:188:HOH:O	2.19	0.43
2:B:47:ASP:O	2:B:48:LEU:HD23	2.19	0.43
1:C:114:PRO:HB3	5:C:151:HOH:O	2.19	0.43
1:A:140:TYR:HD2	2:D:37:TRP:CE3	2.37	0.43
1:A:22:GLY:HA2	1:A:60:LYS:HA	2.00	0.42
2:D:7:GLU:HG2	2:D:129:ALA:HB2	2.00	0.42
2:D:4:THR:OG1	2:D:6:GLU:HB2	2.19	0.42
2:B:101:GLU:HB3	5:C:177:HOH:O	2.20	0.42
1:A:52:SER:O	1:A:56:LYS:HG3	2.19	0.42
1:C:56:LYS:HG2	5:C:202:HOH:O	2.20	0.42
1:C:75:ASP:OD1	1:C:77:PRO:HD2	2.18	0.42
1:C:7:LYS:O	1:C:11:LYS:HG3	2.19	0.42
1:C:42:TYR:CE2	1:C:93:VAL:HA	2.54	0.42
2:D:2:HIS:O	2:D:3:LEU:HD23	2.19	0.42
1:A:58:HIS:HA	1:A:61:LYS:HD3	2.00	0.42
2:B:21:ASP:HA	2:B:65:LYS:HB3	2.02	0.41
2:B:6:GLU:CD	2:B:6:GLU:H	2.23	0.41
1:A:91:LEU:HD13	3:A:142:HEM:C4D	2.56	0.41
1:A:129:LEU:HA	1:A:129:LEU:HD23	1.81	0.41
1:C:29:LEU:HD23	1:C:32:MET:CE	2.51	0.41
2:B:26:GLU:O	2:B:26:GLU:HG3	2.21	0.41
3:D:147:HEM:CHC	3:D:147:HEM:HBB2	2.36	0.41
2:B:100:PRO:HG2	1:C:38:THR:HG21	2.01	0.41
2:D:43:GLU:OE1	2:D:43:GLU:HA	2.21	0.41
2:D:50:THR:HG22	2:D:53:ALA:N	2.32	0.41
2:B:1:VAL:HG23	2:B:3:LEU:HD12	2.03	0.41
2:D:98:VAL:O	2:D:145:TYR:OH	2.32	0.41
1:C:10:VAL:HG22	1:C:125:LEU:HD23	2.03	0.41
2:D:50:THR:O	2:D:53:ALA:HB3	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:103:HIS:NE2	2:B:131:GLN:OE1	2.40	0.40
1:C:14:TRP:CD1	1:C:70:VAL:HG21	2.55	0.40
3:D:147:HEM:HHA	3:D:147:HEM:HAA1	1.90	0.40
2:D:57:ASN:HA	2:D:58:PRO:HD3	1.62	0.40
2:D:74:GLY:O	2:D:77:HIS:N	2.42	0.40
1:A:15:GLY:O	1:A:18:GLY:N	2.54	0.40
5:A:190:HOH:O	2:B:120:LYS:HD2	2.21	0.40
2:B:5:PRO:HG2	2:B:6:GLU:OE1	2.21	0.40
2:B:131:GLN:HB3	2:B:131:GLN:HE21	1.65	0.40
2:B:14:LEU:HD21	2:B:118:PHE:CG	2.56	0.40
2:D:75:LEU:HD23	2:D:75:LEU:HA	1.88	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%)	127 (91%)	11 (8%)	1 (1%)	25	18
1	C	139/141 (99%)	132 (95%)	7 (5%)	0	100	100
2	B	144/146 (99%)	137 (95%)	7 (5%)	0	100	100
2	D	144/146 (99%)	136 (94%)	8 (6%)	0	100	100
All	All	566/574 (99%)	532 (94%)	33 (6%)	1 (0%)	51	48

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	41	THR

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%)	106 (94%)	7 (6%)	21	16
1	C	113/113 (100%)	107 (95%)	6 (5%)	26	21
2	B	117/117 (100%)	106 (91%)	11 (9%)	10	6
2	D	117/117 (100%)	108 (92%)	9 (8%)	15	9
All	All	460/460 (100%)	427 (93%)	33 (7%)	17	11

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	PRO
1	A	52	SER
1	A	84	SER
1	A	105	LEU
1	A	109	LEU
1	A	131	SER
1	A	137	THR
2	B	2	HIS
2	B	6	GLU
2	B	12	THR
2	B	14	LEU
2	B	28	LEU
2	B	40	ARG
2	B	59	LYS
2	B	65	LYS
2	B	66	LYS
2	B	120	LYS
2	B	121	GLU
1	C	16	LYS
1	C	56	LYS
1	C	99	LYS
1	C	118	THR
1	C	124	SER
1	C	131	SER

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Mol	Chain	Res	Type
2	D	5	PRO
2	D	44	SER
2	D	50	THR
2	D	61	LYS
2	D	66	LYS
2	D	78	LEU
2	D	90	GLU
2	D	131	GLN
2	D	146	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	77	HIS
2	D	117	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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
3	HEM	A	142	1,4	28,50,50	2.30	9 (32%)	17,82,82	2.13	8 (47%)
4	OXY	A	143	3	1,1,1	0.55	0	0,0,0	0.00	-
3	HEM	B	147	2	28,50,50	2.35	7 (25%)	17,82,82	2.30	6 (35%)
3	HEM	C	142	1,4	28,50,50	2.42	12 (42%)	17,82,82	2.01	6 (35%)
4	OXY	C	143	3	1,1,1	0.05	0	0,0,0	0.00	-
3	HEM	D	147	2	28,50,50	1.98	7 (25%)	17,82,82	2.54	4 (23%)

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	142	1,4	-	0/6/54/54	0/0/8/8
4	OXY	A	143	3	-	0/0/0/0	0/0/0/0
3	HEM	B	147	2	-	0/6/54/54	0/0/8/8
3	HEM	C	142	1,4	-	0/6/54/54	0/0/8/8
4	OXY	C	143	3	-	0/0/0/0	0/0/0/0
3	HEM	D	147	2	-	0/6/54/54	0/0/8/8

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	142	HEM	C3B-C2B	-6.21	1.32	1.40
3	D	147	HEM	C3B-C2B	-5.14	1.33	1.40
3	B	147	HEM	C3B-C2B	-4.81	1.34	1.40
3	C	142	HEM	C3B-C2B	-4.79	1.34	1.40
3	A	142	HEM	C3C-C2C	-4.52	1.34	1.40
3	B	147	HEM	C3C-C2C	-4.14	1.34	1.40
3	C	142	HEM	C3C-C2C	-3.75	1.35	1.40
3	D	147	HEM	C3C-C2C	-3.36	1.35	1.40
3	C	142	HEM	C2A-C3A	-2.10	1.31	1.37
3	C	142	HEM	C1B-NB	2.00	1.39	1.36
3	D	147	HEM	C4C-NC	2.04	1.39	1.36
3	B	147	HEM	C1A-NA	2.10	1.40	1.36
3	D	147	HEM	C4A-NA	2.11	1.40	1.36
3	C	142	HEM	CMC-C2C	2.28	1.56	1.51
3	B	147	HEM	C3C-CAC	2.37	1.52	1.47
3	C	142	HEM	C4C-NC	2.43	1.39	1.36
3	C	142	HEM	CAD-C3D	2.43	1.56	1.52
3	A	142	HEM	CAD-C3D	2.51	1.57	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	142	HEM	CMC-C2C	2.66	1.57	1.51
3	A	142	HEM	C1C-NC	2.66	1.39	1.36
3	C	142	HEM	C3C-CAC	2.74	1.53	1.47
3	A	142	HEM	C4C-NC	2.95	1.40	1.36
3	C	142	HEM	C4B-NB	2.98	1.42	1.36
3	A	142	HEM	C4D-ND	3.00	1.40	1.36
3	A	142	HEM	C3C-CAC	3.26	1.54	1.47
3	D	147	HEM	C3C-CAC	3.30	1.54	1.47
3	B	147	HEM	CAA-C2A	3.78	1.58	1.52
3	D	147	HEM	C3B-CAB	3.83	1.55	1.47
3	D	147	HEM	C1C-NC	3.96	1.41	1.36
3	C	142	HEM	CAA-C2A	4.18	1.59	1.52
3	A	142	HEM	C3B-CAB	4.54	1.56	1.47
3	C	142	HEM	C3B-CAB	4.87	1.57	1.47
3	C	142	HEM	C1C-NC	5.07	1.42	1.36
3	B	147	HEM	C3B-CAB	5.26	1.58	1.47
3	B	147	HEM	C1C-NC	6.13	1.44	1.36

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	147	HEM	CMA-C3A-C4A	-3.95	122.39	128.46
3	B	147	HEM	CMA-C3A-C4A	-3.72	122.74	128.46
3	C	142	HEM	CMD-C2D-C1D	-3.71	122.77	128.46
3	B	147	HEM	CMD-C2D-C1D	-3.48	123.11	128.46
3	A	142	HEM	CMA-C3A-C4A	-3.29	123.41	128.46
3	A	142	HEM	CAD-CBD-CGD	-3.16	107.26	112.66
3	A	142	HEM	CAA-CBA-CGA	-3.04	107.46	112.66
3	C	142	HEM	CMA-C3A-C4A	-2.94	123.94	128.46
3	A	142	HEM	CMD-C2D-C1D	-2.63	124.43	128.46
3	C	142	HEM	CBA-CAA-C2A	-2.02	108.61	112.48
3	A	142	HEM	CMD-C2D-C3D	2.56	129.77	124.94
3	D	147	HEM	CMA-C3A-C2A	2.61	129.86	124.94
3	A	142	HEM	CMA-C3A-C2A	2.79	130.21	124.94
3	C	142	HEM	CMA-C3A-C2A	2.83	130.28	124.94
3	A	142	HEM	CBA-CAA-C2A	2.87	117.98	112.48
3	B	147	HEM	CMD-C2D-C3D	2.95	130.50	124.94
3	D	147	HEM	CMB-C2B-C3B	3.03	130.51	124.89
3	C	142	HEM	CMD-C2D-C3D	3.17	130.92	124.94
3	A	142	HEM	CMB-C2B-C3B	3.26	130.94	124.89
3	C	142	HEM	CMB-C2B-C3B	3.27	130.96	124.89
3	B	147	HEM	CMA-C3A-C2A	3.36	131.27	124.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	147	HEM	CAA-CBA-CGA	3.75	119.06	112.66
3	B	147	HEM	CMB-C2B-C3B	3.84	132.02	124.89
3	D	147	HEM	CBA-CAA-C2A	7.83	127.45	112.48

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	142	HEM	4	0
3	B	147	HEM	4	0
3	C	142	HEM	2	0
3	D	147	HEM	3	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 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.