



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 11:44 PM GMT

PDB ID : 1YG5  
Title : T-To-T(High) quaternary transitions in human hemoglobin: betaW37H OXY  
(2MM IHP, 20% PEG) (10 test sets)  
Authors : Kavanaugh, J.S.; Rogers, P.H.; Arnone, A.  
Deposited on : 2005-01-04  
Resolution : 2.70 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](mailto: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.

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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

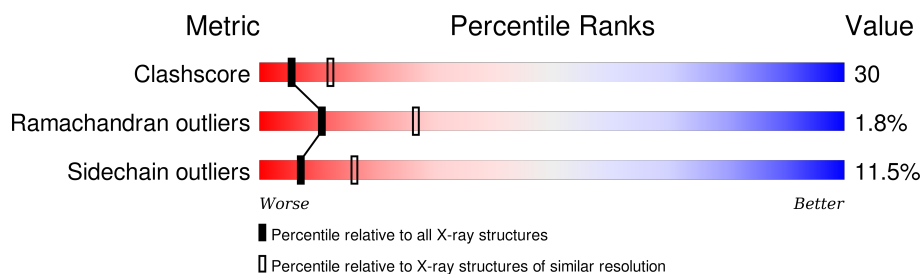
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.70 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	2422 (2.70-2.70)
Ramachandran outliers	100387	2382 (2.70-2.70)
Sidechain outliers	100360	2382 (2.70-2.70)

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  $\geq 3$ , 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 4672 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 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 beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1120	719	196	201	4			
2	D	146	Total	C	N	O	S	0	0	0
			1120	719	196	201	4			

There are 4 discrepancies between the modelled and reference sequences:

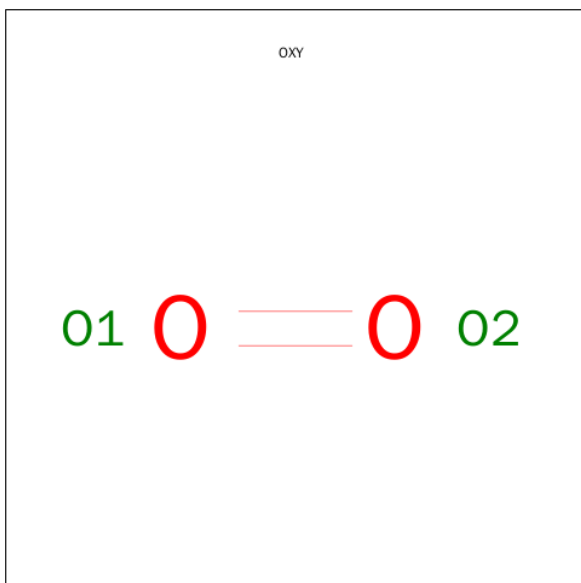
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	VAL	ENGINEERED	UNP P68871
B	37	HIS	TRP	ENGINEERED	UNP P68871
D	1	MET	VAL	ENGINEERED	UNP P68871
D	37	HIS	TRP	ENGINEERED	UNP P68871

- 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 2 2	0	0
4	B	1	Total O 2 2	0	0
4	C	1	Total O 2 2	0	0
4	D	1	Total O 2 2	0	0

- Molecule 5 is water.

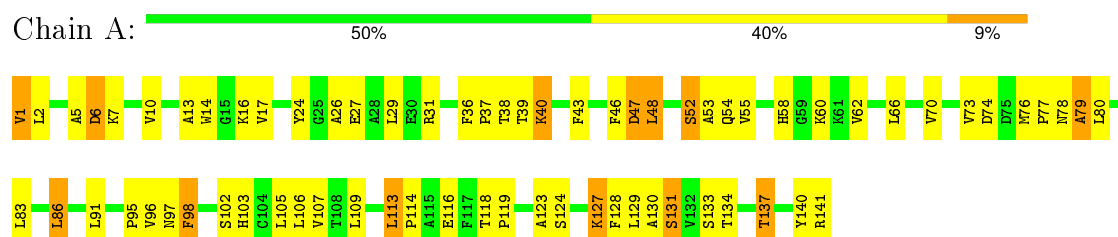
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	33	Total O 33 33	0	0
5	B	31	Total O 31 31	0	0
5	C	27	Total O 27 27	0	0
5	D	23	Total O 23 23	0	0

### 3 Residue-property plots

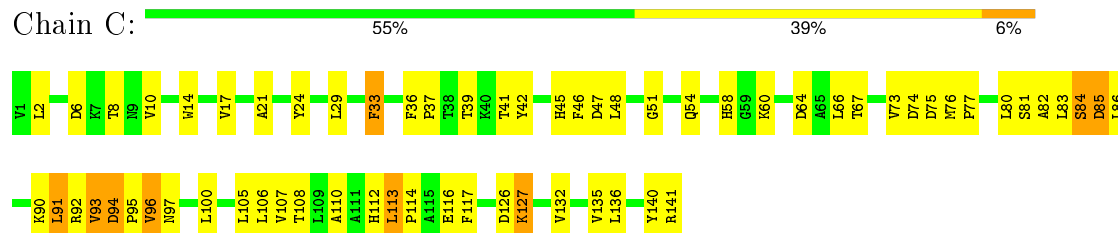
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according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

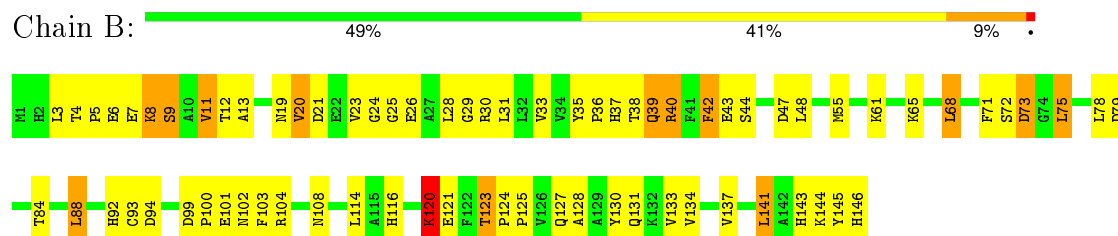
#### • Molecule 1: Hemoglobin alpha chain



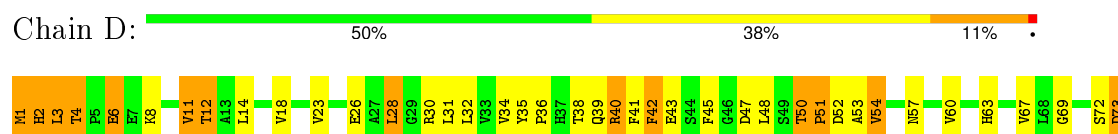
#### • Molecule 1: Hemoglobin alpha chain



#### • Molecule 2: Hemoglobin beta chain



#### • Molecule 2: Hemoglobin 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 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	95.57Å 98.59Å 66.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.70	Depositor
% Data completeness (in resolution range)	87.3 (10.00-2.70)	Depositor
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	REFMAC 5	Depositor
R, $R_{free}$	0.179 , 0.273	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4672	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	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	1/1097 (0.1%)	1.22	4/1491 (0.3%)
1	C	1.06	1/1097 (0.1%)	1.20	10/1491 (0.7%)
2	B	1.19	2/1149 (0.2%)	1.22	11/1558 (0.7%)
2	D	1.06	1/1149 (0.1%)	1.19	7/1558 (0.4%)
All	All	1.10	5/4492 (0.1%)	1.20	32/6098 (0.5%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	73	ASP	CB-CG	7.65	1.67	1.51
2	D	111	VAL	CB-CG2	-5.83	1.40	1.52
1	A	98	PHE	CB-CG	-5.63	1.41	1.51
1	C	33	PHE	CE1-CZ	-5.54	1.26	1.37
2	B	130	TYR	CD1-CE1	-5.07	1.31	1.39

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	94	ASP	CB-CG-OD2	8.83	126.24	118.30
2	D	28	LEU	CA-CB-CG	-8.34	96.11	115.30
2	D	114	LEU	CB-CG-CD2	-8.09	97.25	111.00
2	B	73	ASP	CB-CG-OD1	7.22	124.80	118.30
2	B	79	ASP	CB-CG-OD2	7.12	124.71	118.30
2	D	112	CYS	CA-CB-SG	-7.04	101.33	114.00
2	B	100	PRO	N-CD-CG	-6.99	92.71	103.20
1	A	6	ASP	CB-CG-OD2	6.94	124.55	118.30
2	D	94	ASP	CB-CG-OD2	6.53	124.17	118.30
1	C	64	ASP	CB-CG-OD2	6.31	123.98	118.30
1	A	47	ASP	CB-CG-OD2	6.13	123.81	118.30
2	D	52	ASP	CB-CG-OD2	6.05	123.74	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	113	LEU	CA-CB-CG	-5.88	101.77	115.30
1	C	85	ASP	CB-CG-OD2	5.88	123.59	118.30
2	D	14	LEU	CB-CG-CD2	-5.63	101.42	111.00
2	B	75	LEU	CA-CB-CG	5.61	128.20	115.30
2	B	28	LEU	CA-CB-CG	-5.59	102.44	115.30
2	B	114	LEU	CA-CB-CG	5.58	128.14	115.30
2	B	73	ASP	CB-CA-C	5.55	121.51	110.40
2	B	47	ASP	CB-CG-OD2	5.49	123.24	118.30
1	A	74	ASP	CB-CG-OD2	5.46	123.22	118.30
1	C	91	LEU	CB-CG-CD1	-5.45	101.74	111.00
1	C	96	VAL	CB-CA-C	-5.43	101.09	111.40
1	A	134	THR	OG1-CB-CG2	-5.39	97.60	110.00
2	D	73	ASP	CB-CG-OD2	5.39	123.15	118.30
2	B	120	LYS	CD-CE-NZ	5.39	124.09	111.70
1	C	93	VAL	CB-CA-C	-5.35	101.24	111.40
1	C	126	ASP	CB-CG-OD1	5.27	123.04	118.30
2	B	73	ASP	OD1-CG-OD2	-5.26	113.30	123.30
1	C	106	LEU	CB-CG-CD1	-5.25	102.07	111.00
2	B	99	ASP	CB-CG-OD1	5.13	122.92	118.30
1	C	47	ASP	CB-CG-OD2	5.07	122.87	118.30

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	70	0
1	C	1069	0	1073	54	0
2	B	1120	0	1115	67	0
2	D	1120	0	1115	68	0
3	A	43	0	30	8	0
3	B	43	0	30	10	0
3	C	43	0	30	6	0
3	D	43	0	30	2	0
4	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	2	0	0	0	0
4	C	2	0	0	0	0
4	D	2	0	0	0	0
5	A	33	0	0	4	0
5	B	31	0	0	15	0
5	C	27	0	0	2	0
5	D	23	0	0	3	0
All	All	4672	0	4496	269	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 30.

All (269) 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:38:THR:HG23	5:A:287:HOH:O	1.21	1.29
3:B:147:HEM:HBC2	3:B:147:HEM:HMC2	1.13	1.06
2:B:73:ASP:HB3	5:B:269:HOH:O	1.55	1.05
3:A:142:HEM:HMB1	3:A:142:HEM:HBB2	1.40	1.01
2:B:144:LYS:CE	5:B:298:HOH:O	2.07	1.00
2:B:108:ASN:CG	5:B:283:HOH:O	2.00	0.99
1:C:39:THR:HG22	1:C:97:ASN:HD22	1.29	0.96
1:A:123:ALA:O	1:A:127:LYS:HG3	1.65	0.94
1:C:46:PHE:HA	1:C:54:GLN:NE2	1.82	0.94
1:C:14:TRP:HE1	1:C:67:THR:HG1	1.16	0.94
2:B:144:LYS:HE2	5:B:298:HOH:O	1.67	0.93
3:B:147:HEM:HBC2	3:B:147:HEM:CMC	1.91	0.93
1:A:7:LYS:HD3	1:A:73:VAL:HG11	1.49	0.93
2:D:124:PRO:HB2	2:D:125:PRO:HD3	1.50	0.92
2:B:73:ASP:CB	5:B:269:HOH:O	2.14	0.90
2:B:31:LEU:HD21	2:B:38:THR:HG21	1.56	0.87
3:B:147:HEM:CBC	3:B:147:HEM:HMC2	2.02	0.86
1:C:114:PRO:O	2:D:116:HIS:HE1	1.58	0.85
1:C:39:THR:HG22	1:C:97:ASN:ND2	1.91	0.84
3:A:142:HEM:CMB	3:A:142:HEM:HBB2	2.07	0.84
1:A:70:VAL:O	1:A:73:VAL:HG23	1.76	0.83
2:B:104:ARG:NH1	5:B:248:HOH:O	1.99	0.81
1:C:94:ASP:OD1	1:C:96:VAL:HG23	1.81	0.81
2:D:32:LEU:HD21	2:D:42:PHE:CE2	2.17	0.79
1:A:46:PHE:HB3	1:A:48:LEU:CD2	2.14	0.77
2:B:144:LYS:HE3	5:B:298:HOH:O	1.79	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:142:HEM:O1D	5:C:437:HOH:O	2.04	0.75
1:A:83:LEU:HD21	3:A:142:HEM:HMA3	1.68	0.75
2:B:40:ARG:HD2	1:C:92:ARG:NE	2.00	0.75
2:D:1:MET:O	2:D:3:LEU:N	2.21	0.74
2:B:21:ASP:OD1	2:B:65:LYS:HD2	1.88	0.72
1:C:29:LEU:HD11	1:C:58:HIS:HD2	1.56	0.71
1:A:46:PHE:CB	1:A:48:LEU:CD2	2.68	0.70
2:D:18:VAL:HG13	2:D:23:VAL:HG21	1.73	0.70
2:D:47:ASP:O	2:D:48:LEU:HD23	1.91	0.70
2:D:32:LEU:HD21	2:D:42:PHE:HE2	1.56	0.68
1:A:46:PHE:CB	1:A:48:LEU:HD21	2.23	0.68
2:B:88:LEU:HB2	5:B:409:HOH:O	1.93	0.68
1:A:16:LYS:O	1:A:17:VAL:C	2.31	0.68
2:D:124:PRO:HB2	2:D:125:PRO:CD	2.24	0.68
2:B:37:HIS:O	2:B:40:ARG:HG2	1.94	0.67
2:D:31:LEU:HG	2:D:31:LEU:O	1.93	0.67
2:B:38:THR:HG22	2:B:102:ASN:ND2	2.10	0.67
1:C:114:PRO:O	2:D:116:HIS:CE1	2.46	0.67
1:A:46:PHE:HB2	1:A:48:LEU:HD21	1.76	0.66
2:B:104:ARG:HG2	2:B:104:ARG:HH11	1.61	0.65
1:A:128:PHE:O	1:A:131:SER:HB2	1.96	0.65
2:D:4:THR:CG2	2:D:6:GLU:HG2	2.27	0.65
3:B:147:HEM:CMB	3:B:147:HEM:HBB2	2.26	0.65
1:A:37:PRO:O	1:A:40:LYS:HG2	1.95	0.65
3:A:142:HEM:HMB1	3:A:142:HEM:CBB	2.24	0.65
2:D:32:LEU:CD2	2:D:42:PHE:HE2	2.09	0.65
3:C:142:HEM:HBB2	3:C:142:HEM:CMB	2.27	0.64
2:B:143:HIS:CG	2:B:143:HIS:O	2.50	0.64
1:C:33:PHE:CE2	1:C:48:LEU:HD22	2.33	0.64
2:D:109:VAL:O	2:D:112:CYS:HB2	1.98	0.64
1:A:95:PRO:O	1:A:96:VAL:C	2.35	0.64
1:C:80:LEU:HD22	1:C:83:LEU:HD12	1.79	0.64
1:A:119:PRO:HD2	5:A:311:HOH:O	1.97	0.64
2:D:69:GLY:O	2:D:72:SER:HB2	1.98	0.63
2:D:36:PRO:O	2:D:39:GLN:HB2	1.98	0.63
2:D:92:HIS:HE1	3:D:147:HEM:NA	1.97	0.63
2:D:40:ARG:HG2	2:D:41:PHE:N	2.13	0.63
2:B:120:LYS:HD3	2:B:120:LYS:H	1.63	0.63
1:A:39:THR:HG22	1:A:97:ASN:HD22	1.63	0.63
1:C:46:PHE:HA	1:C:54:GLN:HE22	1.63	0.62
2:D:50:THR:O	2:D:54:VAL:HG23	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:92:HIS:HE1	3:B:147:HEM:C4A	2.16	0.62
2:B:124:PRO:HB2	2:B:125:PRO:HD3	1.81	0.62
1:A:140:TYR:O	1:A:141:ARG:HB3	2.00	0.62
1:A:113:LEU:HG	1:A:116:GLU:HG3	1.82	0.61
2:D:40:ARG:NH1	2:D:41:PHE:CZ	2.68	0.61
2:D:35:TYR:CZ	2:D:105:LEU:HD22	2.36	0.61
2:D:109:VAL:O	2:D:112:CYS:CB	2.49	0.60
1:C:42:TYR:CE1	1:C:93:VAL:HA	2.37	0.60
1:C:97:ASN:HA	1:C:100:LEU:HD12	1.84	0.60
2:B:3:LEU:HD23	2:B:78:LEU:HD13	1.82	0.59
3:A:142:HEM:HBA1	3:A:142:HEM:HMA1	1.83	0.59
3:B:147:HEM:CMC	3:B:147:HEM:CBC	2.62	0.58
2:D:35:TYR:CE2	2:D:105:LEU:HD22	2.39	0.58
2:B:143:HIS:CD2	2:B:143:HIS:O	2.57	0.58
1:A:123:ALA:O	1:A:127:LYS:CG	2.48	0.58
1:C:36:PHE:C	5:C:268:HOH:O	2.43	0.57
1:C:116:GLU:OE1	1:C:116:GLU:N	2.37	0.57
2:D:32:LEU:CD2	2:D:42:PHE:CE2	2.87	0.57
2:D:109:VAL:O	2:D:112:CYS:N	2.38	0.57
2:D:11:VAL:HG22	2:D:130:TYR:CZ	2.39	0.57
2:B:75:LEU:CD2	2:B:133:VAL:HG11	2.34	0.57
1:A:6:ASP:HA	1:A:124:SER:OG	2.03	0.57
1:A:83:LEU:HD11	3:A:142:HEM:HMA3	1.87	0.56
2:D:40:ARG:O	2:D:43:GLU:HG2	2.05	0.56
2:D:63:HIS:O	2:D:67:VAL:HG23	2.04	0.56
1:A:103:HIS:O	1:A:107:VAL:HG23	2.06	0.56
1:A:1:VAL:HG12	1:A:1:VAL:O	2.05	0.56
2:B:39:GLN:HG2	2:B:48:LEU:HD13	1.87	0.55
2:B:108:ASN:OD1	5:B:283:HOH:O	2.14	0.55
2:B:92:HIS:HE1	3:B:147:HEM:NA	1.96	0.55
2:B:3:LEU:CD2	2:B:78:LEU:HD13	2.36	0.55
2:B:11:VAL:O	2:B:12:THR:C	2.45	0.55
1:A:113:LEU:HB3	1:A:116:GLU:HG2	1.88	0.55
1:C:112:HIS:C	1:C:114:PRO:HD3	2.27	0.54
2:D:40:ARG:HD3	2:D:41:PHE:CE2	2.41	0.54
1:C:45:HIS:O	1:C:54:GLN:NE2	2.40	0.54
3:B:147:HEM:O1A	5:B:262:HOH:O	2.18	0.53
1:A:78:ASN:O	1:A:80:LEU:N	2.41	0.53
1:A:127:LYS:HD3	5:A:226:HOH:O	2.09	0.53
1:A:78:ASN:O	1:A:79:ALA:C	2.47	0.53
2:D:99:ASP:CB	5:D:207:HOH:O	2.55	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:40:ARG:HD2	1:C:92:ARG:HE	1.71	0.52
1:C:80:LEU:CD2	1:C:83:LEU:HD12	2.39	0.52
1:A:127:LYS:HD2	1:C:141:ARG:NH1	2.25	0.52
1:A:76:MET:N	1:A:77:PRO:CD	2.71	0.52
2:B:123:THR:OG1	2:B:125:PRO:HD2	2.08	0.52
2:B:103:PHE:CE2	2:B:141:LEU:HB3	2.44	0.52
2:B:7:GLU:O	2:B:8:LYS:C	2.46	0.52
2:D:93:CYS:HB2	2:D:145:TYR:CZ	2.45	0.52
2:D:48:LEU:HD22	2:D:54:VAL:HG13	1.90	0.52
1:C:110:ALA:HB2	1:C:117:PHE:CZ	2.45	0.52
1:C:73:VAL:O	1:C:75:ASP:N	2.43	0.52
2:B:101:GLU:OE2	2:B:104:ARG:HD2	2.10	0.51
1:A:66:LEU:HD11	1:A:105:LEU:HD11	1.91	0.51
2:D:50:THR:O	2:D:53:ALA:HB3	2.10	0.51
2:D:93:CYS:HB2	2:D:145:TYR:CE1	2.45	0.51
2:B:25:GLY:O	2:B:26:GLU:C	2.47	0.51
2:B:5:PRO:O	2:B:9:SER:HB2	2.10	0.51
2:D:99:ASP:HB2	5:D:207:HOH:O	2.10	0.51
2:B:55:MET:O	2:B:61:LYS:HE2	2.11	0.51
2:D:30:ARG:HG2	2:D:30:ARG:NH1	2.26	0.51
2:D:77:HIS:HD1	2:D:77:HIS:N	2.08	0.50
1:A:95:PRO:O	1:A:97:ASN:N	2.45	0.50
1:C:90:LYS:O	1:C:92:ARG:NE	2.45	0.50
1:C:132:VAL:HG12	1:C:136:LEU:HD12	1.93	0.50
3:B:147:HEM:CMB	3:B:147:HEM:CBB	2.89	0.50
1:A:114:PRO:O	2:B:116:HIS:NE2	2.39	0.50
2:D:124:PRO:CB	2:D:125:PRO:HD3	2.34	0.50
2:B:93:CYS:HB2	2:B:145:TYR:CE1	2.47	0.50
1:A:46:PHE:CB	1:A:48:LEU:HD23	2.41	0.50
1:C:73:VAL:C	1:C:75:ASP:H	2.15	0.50
2:B:121:GLU:CD	5:B:289:HOH:O	2.50	0.49
1:C:33:PHE:CD2	1:C:48:LEU:HD22	2.48	0.49
2:B:42:PHE:N	2:B:42:PHE:CD2	2.80	0.49
2:D:103:PHE:CE2	2:D:141:LEU:HB3	2.47	0.49
1:C:86:LEU:O	1:C:91:LEU:HB2	2.12	0.49
1:C:91:LEU:HD11	3:C:142:HEM:C3D	2.47	0.48
1:A:105:LEU:O	1:A:109:LEU:HG	2.13	0.48
1:A:95:PRO:O	1:A:98:PHE:N	2.44	0.48
1:A:83:LEU:HD21	3:A:142:HEM:CMA	2.39	0.48
2:B:4:THR:HB	2:B:5:PRO:HD2	1.95	0.48
1:C:82:ALA:O	1:C:85:ASP:N	2.47	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:142:HEM:CBB	3:C:142:HEM:CMB	2.92	0.48
2:B:93:CYS:HB2	2:B:145:TYR:CZ	2.48	0.48
2:D:73:ASP:O	2:D:77:HIS:ND1	2.47	0.48
1:A:26:ALA:O	1:A:27:GLU:C	2.50	0.48
1:C:82:ALA:O	1:C:84:SER:N	2.47	0.48
1:C:6:ASP:O	1:C:10:VAL:HG23	2.13	0.48
3:C:142:HEM:HMB1	3:C:142:HEM:CBB	2.44	0.47
3:C:142:HEM:HMB1	3:C:142:HEM:HBB2	1.94	0.47
1:A:46:PHE:HB3	1:A:48:LEU:HD21	1.91	0.47
2:D:42:PHE:O	2:D:45:PHE:HB2	2.14	0.47
2:B:11:VAL:HG12	2:B:12:THR:N	2.29	0.47
1:A:86:LEU:HA	1:A:86:LEU:HD22	1.63	0.47
1:A:102:SER:HB3	1:A:129:LEU:HD13	1.96	0.47
2:D:141:LEU:HA	2:D:141:LEU:HD23	1.59	0.47
2:D:127:GLN:O	2:D:131:GLN:HG2	2.14	0.47
2:B:11:VAL:O	2:B:13:ALA:N	2.47	0.47
5:B:242:HOH:O	1:C:41:THR:HG22	2.14	0.47
1:C:83:LEU:HB3	1:C:136:LEU:HD21	1.97	0.47
1:A:113:LEU:HG	1:A:116:GLU:CG	2.45	0.47
1:A:29:LEU:HD13	1:A:58:HIS:HD2	1.80	0.47
2:B:93:CYS:HB2	2:B:145:TYR:CD1	2.50	0.47
2:B:125:PRO:O	2:B:128:ALA:HB3	2.14	0.47
1:A:5:ALA:O	1:A:6:ASP:C	2.53	0.47
2:B:3:LEU:HA	2:B:3:LEU:HD12	1.70	0.46
1:A:114:PRO:CD	5:A:352:HOH:O	2.63	0.46
2:B:44:SER:HB3	5:B:251:HOH:O	2.14	0.46
2:B:37:HIS:HE1	2:B:102:ASN:HD21	1.63	0.46
2:D:93:CYS:SG	2:D:93:CYS:O	2.73	0.46
2:B:131:GLN:HA	2:B:134:VAL:HG22	1.98	0.46
1:A:58:HIS:O	1:A:62:VAL:HG23	2.16	0.46
1:A:7:LYS:HD3	1:A:73:VAL:CG1	2.34	0.45
1:C:136:LEU:HD23	1:C:136:LEU:HA	1.72	0.45
2:D:11:VAL:HG12	2:D:12:THR:N	2.31	0.45
2:D:28:LEU:HA	2:D:28:LEU:HD12	1.74	0.45
2:B:23:VAL:HG12	2:B:68:LEU:HD11	1.98	0.45
2:D:30:ARG:HG2	2:D:30:ARG:HH11	1.82	0.45
1:C:76:MET:N	1:C:77:PRO:CD	2.79	0.45
2:D:11:VAL:HG21	2:D:133:VAL:HG21	1.99	0.45
2:D:57:ASN:CG	2:D:60:VAL:HG23	2.37	0.45
2:B:29:GLY:O	2:B:33:VAL:HG23	2.16	0.45
1:A:10:VAL:HG21	1:A:128:PHE:HB2	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:ALA:O	1:A:17:VAL:HG23	2.16	0.45
1:A:37:PRO:O	1:A:38:THR:C	2.54	0.45
1:A:76:MET:HB2	1:A:77:PRO:HD3	1.98	0.45
2:D:77:HIS:N	2:D:77:HIS:ND1	2.64	0.45
1:A:133:SER:O	1:A:137:THR:HB	2.17	0.45
1:C:91:LEU:HA	1:C:91:LEU:HD23	1.78	0.45
1:C:81:SER:O	1:C:84:SER:HB2	2.17	0.45
1:C:112:HIS:C	1:C:113:LEU:HG	2.38	0.44
2:D:41:PHE:CE1	2:D:98:VAL:HG22	2.52	0.44
2:B:24:GLY:HA2	2:B:68:LEU:HG	1.98	0.44
2:D:50:THR:O	2:D:51:PRO:C	2.56	0.44
1:A:118:THR:HB	1:A:119:PRO:CD	2.47	0.44
1:A:46:PHE:HB2	1:A:48:LEU:CD2	2.38	0.44
2:D:106:LEU:HD23	3:D:147:HEM:HAB	1.98	0.44
2:B:71:PHE:CE2	2:B:137:VAL:HG21	2.52	0.44
1:C:90:LYS:O	1:C:92:ARG:CZ	2.66	0.44
1:C:24:TYR:OH	1:C:112:HIS:HB3	2.18	0.44
2:D:115:ALA:HB2	2:D:122:PHE:CD2	2.53	0.44
1:C:94:ASP:O	1:C:95:PRO:C	2.54	0.44
1:C:82:ALA:C	1:C:84:SER:N	2.70	0.43
1:A:6:ASP:OD1	1:A:124:SER:OG	2.34	0.43
2:D:106:LEU:HD12	2:D:106:LEU:HA	1.72	0.43
1:A:106:LEU:HA	1:A:106:LEU:HD23	1.76	0.43
2:D:124:PRO:CB	2:D:125:PRO:CD	2.91	0.43
2:D:8:LYS:NZ	2:D:79:ASP:OD2	2.38	0.43
2:B:30:ARG:O	2:B:31:LEU:C	2.57	0.43
2:B:108:ASN:ND2	5:B:283:HOH:O	2.37	0.43
2:B:38:THR:CG2	2:B:102:ASN:ND2	2.79	0.43
2:D:40:ARG:O	2:D:43:GLU:CG	2.67	0.43
2:B:35:TYR:N	2:B:36:PRO:CD	2.82	0.43
1:A:43:PHE:HD2	1:A:46:PHE:CD2	2.36	0.43
2:D:1:MET:O	2:D:2:HIS:C	2.57	0.42
1:A:38:THR:O	1:A:39:THR:C	2.57	0.42
2:D:35:TYR:O	2:D:38:THR:HG23	2.19	0.42
2:B:123:THR:O	2:B:124:PRO:C	2.57	0.42
2:B:94:ASP:OD1	2:B:146:HIS:NE2	2.46	0.42
2:D:57:ASN:C	2:D:57:ASN:OD1	2.58	0.42
2:D:109:VAL:O	2:D:112:CYS:CA	2.67	0.42
1:C:36:PHE:O	1:C:37:PRO:C	2.58	0.42
1:A:86:LEU:HD13	1:A:91:LEU:CD1	2.50	0.42
1:A:47:ASP:N	1:A:54:GLN:OE1	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:83:LEU:HD21	3:A:142:HEM:C3A	2.55	0.42
1:C:17:VAL:HG13	1:C:24:TYR:CD2	2.55	0.42
1:C:83:LEU:HA	1:C:83:LEU:HD23	1.69	0.42
1:A:66:LEU:HD11	1:A:105:LEU:CD1	2.49	0.42
2:B:104:ARG:HG2	2:B:104:ARG:NH1	2.32	0.42
1:C:73:VAL:C	1:C:75:ASP:N	2.73	0.42
1:A:36:PHE:N	1:A:37:PRO:HD3	2.35	0.41
2:B:143:HIS:HD2	5:B:298:HOH:O	2.03	0.41
2:B:84:THR:O	2:B:84:THR:HG22	2.20	0.41
2:D:99:ASP:HB3	5:D:207:HOH:O	2.19	0.41
2:B:4:THR:HB	2:B:5:PRO:CD	2.50	0.41
2:B:5:PRO:O	2:B:6:GLU:C	2.58	0.41
1:C:82:ALA:O	1:C:83:LEU:C	2.56	0.41
1:A:17:VAL:HG13	1:A:24:TYR:CD2	2.56	0.41
1:C:6:ASP:OD2	1:C:127:LYS:HE2	2.21	0.41
1:C:76:MET:HE3	1:C:76:MET:HB3	1.74	0.41
1:A:14:TRP:HA	1:A:17:VAL:HG23	2.03	0.41
1:C:113:LEU:HA	1:C:113:LEU:HD23	1.89	0.41
3:B:147:HEM:HAA1	3:B:147:HEM:HHA	1.74	0.41
1:A:127:LYS:HE2	1:A:127:LYS:HB3	0.95	0.41
1:C:91:LEU:O	1:C:92:ARG:HB2	2.21	0.41
2:D:93:CYS:HB2	2:D:145:TYR:CD1	2.56	0.41
2:D:4:THR:HG21	2:D:6:GLU:HG2	2.03	0.41
1:A:118:THR:HB	1:A:119:PRO:HD2	2.02	0.41
2:D:11:VAL:HG22	2:D:130:TYR:CE1	2.56	0.41
1:A:52:SER:OG	1:A:53:ALA:N	2.53	0.41
1:C:66:LEU:HD23	1:C:66:LEU:HA	1.84	0.41
2:D:93:CYS:HB2	2:D:145:TYR:CE2	2.55	0.40
1:A:137:THR:O	1:A:137:THR:CG2	2.69	0.40
1:A:31:ARG:HG2	2:B:127:GLN:OE1	2.20	0.40
2:B:19:ASN:O	2:B:20:VAL:C	2.58	0.40
2:B:37:HIS:CE1	2:B:102:ASN:HD21	2.39	0.40
2:D:34:VAL:HG12	2:D:34:VAL:O	2.20	0.40
2:B:120:LYS:H	2:B:120:LYS:CD	2.32	0.40
1:A:102:SER:HA	1:A:129:LEU:HD13	2.03	0.40
1:A:130:ALA:O	1:A:133:SER:HB2	2.22	0.40
2:D:105:LEU:HD23	2:D:105:LEU:HA	1.95	0.40
1:C:107:VAL:O	1:C:108:THR:C	2.57	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	139/141 (99%)	119 (86%)	18 (13%)	2 (1%)	14	35
1	C	139/141 (99%)	121 (87%)	15 (11%)	3 (2%)	8	22
2	B	144/146 (99%)	130 (90%)	11 (8%)	3 (2%)	9	23
2	D	144/146 (99%)	130 (90%)	12 (8%)	2 (1%)	14	35
All	All	566/574 (99%)	500 (88%)	56 (10%)	10 (2%)	11	27

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	2	HIS
1	A	79	ALA
2	B	39	GLN
1	C	21	ALA
1	C	74	ASP
2	B	120	LYS
2	D	51	PRO
2	B	11	VAL
1	C	51	GLY
1	A	55	VAL

### 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%)	102 (90%)	11 (10%)	10	23

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	113/113 (100%)	105 (93%)	8 (7%)	18	41
2	B	118/118 (100%)	106 (90%)	12 (10%)	9	21
2	D	118/118 (100%)	96 (81%)	22 (19%)	2	5
All	All	462/462 (100%)	409 (88%)	53 (12%)	7	16

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	VAL
1	A	2	LEU
1	A	40	LYS
1	A	48	LEU
1	A	52	SER
1	A	60	LYS
1	A	86	LEU
1	A	113	LEU
1	A	127	LYS
1	A	131	SER
1	A	137	THR
2	B	8	LYS
2	B	9	SER
2	B	20	VAL
2	B	40	ARG
2	B	42	PHE
2	B	43	GLU
2	B	68	LEU
2	B	72	SER
2	B	88	LEU
2	B	120	LYS
2	B	123	THR
2	B	141	LEU
1	C	2	LEU
1	C	8	THR
1	C	60	LYS
1	C	84	SER
1	C	105	LEU
1	C	127	LYS
1	C	135	VAL
1	C	140	TYR
2	D	1	MET
2	D	3	LEU

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Mol	Chain	Res	Type
2	D	4	THR
2	D	6	GLU
2	D	11	VAL
2	D	12	THR
2	D	26	GLU
2	D	40	ARG
2	D	42	PHE
2	D	50	THR
2	D	54	VAL
2	D	77	HIS
2	D	82	LYS
2	D	87	THR
2	D	88	LEU
2	D	90	GLU
2	D	95	LYS
2	D	110	LEU
2	D	112	CYS
2	D	117	HIS
2	D	120	LYS
2	D	124	PRO

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	58	HIS
1	A	97	ASN
2	B	37	HIS
2	B	97	HIS
2	B	102	ASN
2	B	143	HIS
1	C	58	HIS
1	C	89	HIS
1	C	97	ASN
2	D	37	HIS
2	D	102	ASN
2	D	116	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	142	1,4	30,50,50	2.76	6 (20%)	24,82,82	2.48	9 (37%)
4	OXY	A	150	3	1,1,1	0.13	0	0,0,0	0.00	-
3	HEM	B	147	2,4	30,50,50	2.64	10 (33%)	24,82,82	2.96	9 (37%)
4	OXY	B	150	3	1,1,1	0.23	0	0,0,0	0.00	-
3	HEM	C	142	1,4	30,50,50	2.40	7 (23%)	24,82,82	2.77	12 (50%)
4	OXY	C	150	3	1,1,1	0.21	0	0,0,0	0.00	-
3	HEM	D	147	2,4	30,50,50	2.28	7 (23%)	24,82,82	3.75	14 (58%)
4	OXY	D	150	3	1,1,1	0.12	0	0,0,0	0.00	-

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	-	0/10/54/54	0/0/8/8
4	OXY	A	150	3	-	0/0/0/0	0/0/0/0
3	HEM	B	147	2,4	-	0/10/54/54	0/0/8/8
4	OXY	B	150	3	-	0/0/0/0	0/0/0/0
3	HEM	C	142	1,4	-	0/10/54/54	0/0/8/8
4	OXY	C	150	3	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	D	147	2,4	-	0/10/54/54	0/0/8/8
4	OXY	D	150	3	-	0/0/0/0	0/0/0/0

All (30) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	142	HEM	C3B-C4B	-11.34	1.41	1.51
3	B	147	HEM	C3B-C4B	-8.73	1.44	1.51
3	C	142	HEM	C3B-C4B	-8.22	1.44	1.51
3	D	147	HEM	C3B-C4B	-6.04	1.46	1.51
3	D	147	HEM	C3D-C4D	-5.91	1.44	1.51
3	B	147	HEM	C3D-C4D	-5.75	1.44	1.51
3	A	142	HEM	C3D-C4D	-5.68	1.44	1.51
3	D	147	HEM	C2C-C1C	-5.57	1.42	1.52
3	C	142	HEM	C2C-C1C	-5.06	1.43	1.52
3	C	142	HEM	C3D-C4D	-4.98	1.45	1.51
3	B	147	HEM	C2C-C1C	-4.70	1.43	1.52
3	A	142	HEM	C2C-C1C	-4.63	1.43	1.52
3	B	147	HEM	FE-NC	-3.49	1.82	1.95
3	B	147	HEM	C1A-CHA	-2.74	1.32	1.39
3	D	147	HEM	C2A-C3A	-2.63	1.29	1.37
3	B	147	HEM	C2B-C1B	-2.63	1.43	1.51
3	D	147	HEM	C1A-CHA	-2.48	1.33	1.39
3	C	142	HEM	FE-NC	-2.38	1.86	1.95
3	C	142	HEM	CHD-C1D	-2.36	1.32	1.38
3	D	147	HEM	C2D-C1D	-2.29	1.44	1.51
3	C	142	HEM	C2D-C1D	-2.27	1.44	1.51
3	B	147	HEM	C2A-C3A	-2.24	1.31	1.37
3	B	147	HEM	C2D-C1D	-2.09	1.45	1.51
3	A	142	HEM	C2D-C1D	-2.07	1.45	1.51
3	A	142	HEM	C2B-C1B	-2.04	1.45	1.51
3	C	142	HEM	C2B-C1B	-2.01	1.45	1.51
3	A	142	HEM	CAA-C2A	2.75	1.56	1.52
3	B	147	HEM	CAA-C2A	2.77	1.56	1.52
3	D	147	HEM	CAA-C2A	4.02	1.58	1.52
3	B	147	HEM	FE-ND	4.91	2.23	1.97

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	147	HEM	C3B-CAB-CBB	-7.28	113.29	124.46
3	C	142	HEM	C3C-CAC-CBC	-6.70	114.18	124.46

*Continued on next page...*

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	D	147	HEM	C3C-CAC-CBC	-6.47	114.53	124.46
3	B	147	HEM	C3C-CAC-CBC	-6.13	115.06	124.46
3	D	147	HEM	C3B-CAB-CBB	-5.22	116.45	124.46
3	A	142	HEM	C3C-CAC-CBC	-4.43	117.66	124.46
3	A	142	HEM	C3B-CAB-CBB	-4.08	118.20	124.46
3	C	142	HEM	C3B-CAB-CBB	-4.02	118.29	124.46
3	D	147	HEM	CAA-C2A-C3A	-3.64	118.62	129.00
3	B	147	HEM	CAA-CBA-CGA	-2.90	107.43	112.75
3	C	142	HEM	CAA-CBA-CGA	-2.71	107.78	112.75
3	D	147	HEM	C1D-CHD-C4C	-2.64	121.41	125.82
3	C	142	HEM	CBD-CAD-C3D	-2.64	105.88	113.55
3	A	142	HEM	CAA-C2A-C1A	-2.53	124.26	127.01
3	A	142	HEM	C4B-CHC-C1C	-2.43	121.77	125.82
3	D	147	HEM	C3B-C4B-NB	-2.42	107.00	111.63
3	B	147	HEM	CBA-CAA-C2A	-2.40	108.22	112.53
3	C	142	HEM	CMA-C3A-C4A	-2.36	124.46	128.36
3	D	147	HEM	CMD-C2D-C3D	2.16	123.90	114.35
3	C	142	HEM	C2D-C3D-C4D	2.31	105.42	101.50
3	A	142	HEM	CMD-C2D-C3D	2.36	124.78	114.35
3	D	147	HEM	CMB-C2B-C3B	2.59	122.99	116.53
3	B	147	HEM	CMC-C2C-C3C	2.62	123.07	116.53
3	D	147	HEM	CHD-C1D-ND	2.77	131.19	124.52
3	C	142	HEM	CBA-CAA-C2A	2.78	117.51	112.53
3	C	142	HEM	CMD-C2D-C3D	2.87	127.02	114.35
3	A	142	HEM	CMB-C2B-C3B	3.13	124.34	116.53
3	C	142	HEM	CMB-C2B-C3B	3.22	124.57	116.53
3	B	147	HEM	CAD-C3D-C2D	3.41	123.03	113.22
3	C	142	HEM	CMC-C2C-C3C	3.52	125.32	116.53
3	D	147	HEM	C3B-C4B-CHC	3.52	128.12	123.16
3	B	147	HEM	C3B-C4B-CHC	3.67	128.33	123.16
3	D	147	HEM	CMC-C2C-C3C	3.80	126.02	116.53
3	C	142	HEM	CAD-C3D-C4D	3.95	126.41	112.47
3	A	142	HEM	CAD-C3D-C2D	4.12	125.05	113.22
3	D	147	HEM	CAD-C3D-C4D	4.74	129.17	112.47
3	A	142	HEM	CAD-C3D-C4D	5.00	130.11	112.47
3	B	147	HEM	CMB-C2B-C3B	5.18	129.46	116.53
3	C	142	HEM	CAD-C3D-C2D	5.20	128.16	113.22
3	A	142	HEM	CBA-CAA-C2A	5.21	121.87	112.53
3	D	147	HEM	CAD-C3D-C2D	5.32	128.51	113.22
3	B	147	HEM	CAD-C3D-C4D	5.69	132.55	112.47
3	D	147	HEM	CBA-CAA-C2A	6.33	123.87	112.53
3	D	147	HEM	CAA-C2A-C1A	9.63	137.46	127.01

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	142	HEM	8	0
3	B	147	HEM	10	0
3	C	142	HEM	6	0
3	D	147	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 [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.