



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 08:12 PM GMT

PDB ID : 1J7Y
Title : Crystal structure of partially ligated mutant of HbA
Authors : Miele, A.E.; Draghi, F.; Arcovito, A.; Bellelli, A.; Brunori, M.; Travaglini-All
ocatelli, C.; Vallone, B.
Deposited on : 2001-05-19
Resolution : 1.70 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
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

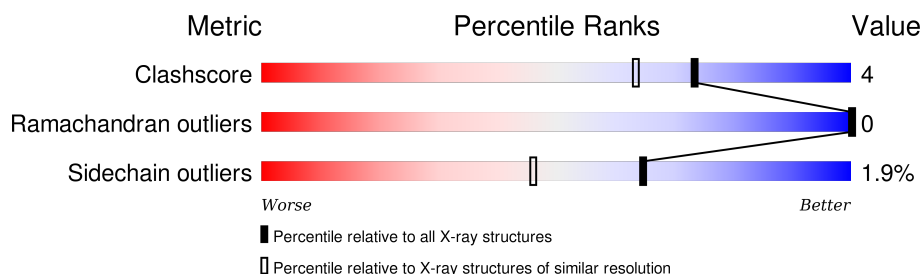
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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	3585 (1.70-1.70)
Ramachandran outliers	100387	3527 (1.70-1.70)
Sidechain outliers	100360	3527 (1.70-1.70)

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 ≥ 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	

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
5	CMO	B	148	-	-	X	-
5	CMO	D	148	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 5041 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	1	0
			1077	690	187	196	4			
1	C	141	Total	C	N	O	S	0	1	0
			1075	689	186	196	4			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	VAL	ENGINEERED	UNP P69905
A	29	TYR	LEU	ENGINEERED	UNP P69905
A	58	GLN	HIS	ENGINEERED	UNP P69905
C	1	MET	VAL	ENGINEERED	UNP P69905
C	29	TYR	LEU	ENGINEERED	UNP P69905
C	58	GLN	HIS	ENGINEERED	UNP P69905

- Molecule 2 is a protein called Hemoglobin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	2	0
			1136	733	194	205	4			
2	D	146	Total	C	N	O	S	0	0	0
			1127	726	194	203	4			

There are 6 discrepancies between the modelled and reference sequences:

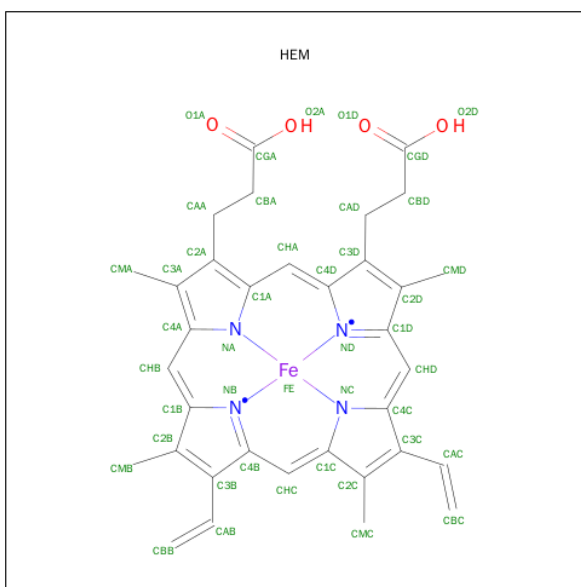
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	VAL	ENGINEERED	UNP P68871
B	28	TYR	LEU	ENGINEERED	UNP P68871
B	63	GLN	HIS	ENGINEERED	UNP P68871
D	1	MET	VAL	ENGINEERED	UNP P68871
D	28	TYR	LEU	ENGINEERED	UNP P68871
D	63	GLN	HIS	ENGINEERED	UNP P68871

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



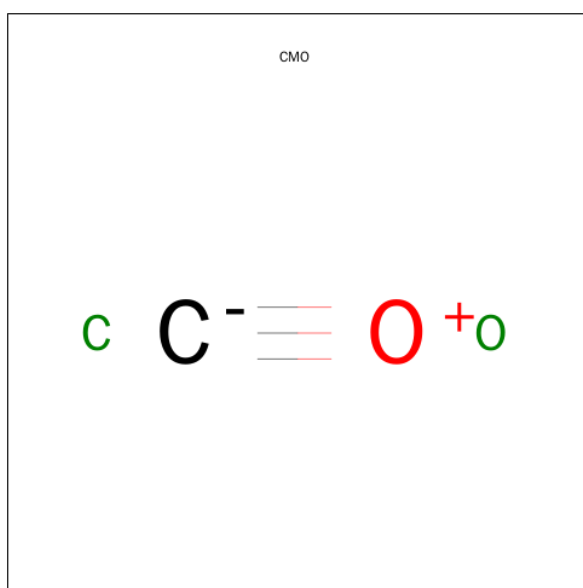
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: C₃₄H₃₂FeN₄O₄).



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

- Molecule 5 is CARBON MONOXIDE (three-letter code: CMO) (formula: CO).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	B	1	Total	C	O	0	0
			2	1	1		
5	D	1	Total	C	O	0	0
			2	1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	149	Total	O	0	0
			149	149		
6	B	106	Total	O	0	0
			106	106		
6	C	91	Total	O	0	0
			91	91		
6	D	94	Total	O	0	0
			94	94		

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 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.

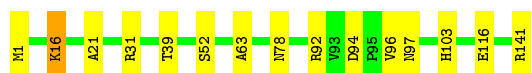
• Molecule 1: Hemoglobin

Chain A:  89% 10%




• Molecule 1: Hemoglobin

Chain C:  89% 10%




• Molecule 2: Hemoglobin

Chain B:  88% 12%



• Molecule 2: Hemoglobin

Chain D:  86% 12%



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	62.57Å 82.47Å 53.69Å 90.00° 99.70° 90.00°	Depositor
Resolution (Å)	14.90 – 1.70	Depositor
% Data completeness (in resolution range)	(Not available) (14.90-1.70)	Depositor
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC	Depositor
R, R_{free}	0.155 , 0.201	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	5041	wwPDB-VP
Average B, all atoms (Å ²)	24.0	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: CMO, 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	0.55	0/1110	1.19	11/1506 (0.7%)
1	C	0.53	0/1108	1.12	4/1505 (0.3%)
2	B	0.51	0/1177	1.13	6/1598 (0.4%)
2	D	0.54	0/1157	1.20	10/1570 (0.6%)
All	All	0.53	0/4552	1.16	31/6179 (0.5%)

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	28	TYR	CB-CG-CD1	-9.43	115.34	121.00
2	D	40	ARG	CD-NE-CZ	9.41	136.77	123.60
1	C	31	ARG	NE-CZ-NH1	8.97	124.78	120.30
1	C	31	ARG	NE-CZ-NH2	-8.96	115.82	120.30
1	C	141	ARG	NE-CZ-NH2	-8.55	116.03	120.30
2	D	73	ASP	CB-CG-OD1	8.49	125.94	118.30
1	A	92	ARG	NE-CZ-NH1	-8.31	116.14	120.30
2	D	28	TYR	CB-CG-CD2	7.82	125.69	121.00
1	A	31	ARG	NE-CZ-NH2	-7.38	116.61	120.30
2	B	30	ARG	NE-CZ-NH1	7.16	123.88	120.30
2	D	30	ARG	NE-CZ-NH1	7.01	123.81	120.30
2	B	52	ASP	CB-CG-OD1	6.97	124.58	118.30
1	A	31	ARG	NE-CZ-NH1	6.72	123.66	120.30
1	A	141	ARG	NE-CZ-NH1	6.67	123.63	120.30
1	A	128	PHE	CB-CG-CD2	-6.52	116.24	120.80
2	D	40	ARG	NE-CZ-NH2	6.43	123.52	120.30
1	A	141	ARG	NE-CZ-NH2	-6.40	117.10	120.30
2	D	40	ARG	NE-CZ-NH1	-6.30	117.15	120.30
2	B	7	GLU	OE1-CD-OE2	-6.15	115.92	123.30
1	A	47	ASP	CB-CG-OD1	6.07	123.76	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	42	TYR	CB-CG-CD2	5.93	124.56	121.00
2	B	21	ASP	CB-CG-OD1	5.86	123.57	118.30
1	C	92	ARG	NE-CZ-NH1	-5.84	117.38	120.30
2	B	40	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	A	42	TYR	CB-CG-CD1	-5.55	117.67	121.00
1	A	92	ARG	NH1-CZ-NH2	5.50	125.44	119.40
2	D	45	PHE	CB-CG-CD1	-5.49	116.96	120.80
1	A	29	TYR	CB-CG-CD2	-5.37	117.78	121.00
2	D	1	MET	C-N-CA	5.30	134.96	121.70
2	D	67	VAL	CB-CA-C	-5.18	101.56	111.40
2	B	131	GLN	CG-CD-OE1	-5.13	111.34	121.60

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	1077	0	1081	5	0
1	C	1075	0	1077	6	0
2	B	1136	0	1125	11	0
2	D	1127	0	1117	15	0
3	B	10	0	0	0	0
4	A	43	0	30	0	0
4	B	43	0	30	1	0
4	C	43	0	30	0	0
4	D	43	0	30	1	0
5	B	2	0	0	2	0
5	D	2	0	0	4	0
6	A	149	0	0	1	0
6	B	106	0	0	2	0
6	C	91	0	0	1	0
6	D	94	0	0	1	0
All	All	5041	0	4520	37	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 4.

All (37) 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)
2:D:67:VAL:HG22	5:D:148:CMO:C	2.17	0.75
2:B:67:VAL:HG22	5:B:148:CMO:C	2.20	0.72
2:D:38:THR:HG22	2:D:102:ASN:ND2	2.09	0.68
2:D:38:THR:HG22	2:D:102:ASN:HD22	1.58	0.67
2:B:67:VAL:CG2	5:B:148:CMO:C	2.82	0.58
1:A:103:HIS:HE1	2:B:131:GLN:OE1	1.87	0.58
4:D:147:HEM:HMC2	4:D:147:HEM:HBC2	1.86	0.57
1:C:103:HIS:HD2	6:C:146:HOH:O	1.87	0.57
1:C:103:HIS:HE1	2:D:131:GLN:OE1	1.89	0.55
1:C:39:THR:HG22	1:C:97:ASN:HD22	1.72	0.55
1:C:16:LYS:HE2	1:C:116:GLU:HG2	1.89	0.54
2:D:28:TYR:CE1	2:D:67:VAL:HG21	2.43	0.54
1:A:103:HIS:HD2	6:A:149:HOH:O	1.90	0.53
2:D:1:MET:HE2	2:D:3:LEU:HG	1.90	0.53
2:B:12[B]:THR:HG23	6:B:1261:HOH:O	2.09	0.53
1:C:21:ALA:HB1	1:C:63:ALA:HB1	1.93	0.51
2:D:43:GLU:HG2	6:D:198:HOH:O	2.12	0.49
2:D:67:VAL:HG11	2:D:106:LEU:HD21	1.94	0.48
2:B:65:LYS:HE3	6:B:1265:HOH:O	2.14	0.47
2:B:28[A]:TYR:CE1	2:B:67:VAL:HG21	2.50	0.47
2:B:86:ALA:HB1	2:B:144:LYS:HE2	1.97	0.47
2:B:51:PRO:O	2:B:55:MET:HG2	2.18	0.44
2:D:42:PHE:O	2:D:45:PHE:HB2	2.18	0.44
1:C:94:ASP:OD2	1:C:96[A]:VAL:HG13	2.18	0.44
2:D:1:MET:HE2	2:D:3:LEU:CG	2.49	0.43
2:B:61:LYS:O	2:B:65:LYS:HG3	2.17	0.43
2:D:67:VAL:CG2	5:D:148:CMO:O	2.67	0.43
2:D:61:LYS:O	2:D:65:LYS:HG3	2.19	0.42
2:D:67:VAL:CG2	5:D:148:CMO:C	2.94	0.42
2:D:63:GLN:NE2	5:D:148:CMO:O	2.53	0.41
1:A:76:MET:N	1:A:77:PRO:CD	2.84	0.41
4:B:147:HEM:HBC2	4:B:147:HEM:HMC2	2.03	0.41
2:B:28[B]:TYR:CD1	2:B:60:VAL:HG13	2.56	0.41
1:A:68:ASN:O	1:A:72:HIS:HD2	2.04	0.41
2:D:50:THR:HB	2:D:51:PRO:HD2	2.03	0.40
1:A:42:TYR:CE1	1:A:93:VAL:HA	2.56	0.40
2:B:63:GLN:O	2:B:67:VAL:HG23	2.21	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	140/141 (99%)	138 (99%)	2 (1%)	0	100	100
1	C	140/141 (99%)	138 (99%)	2 (1%)	0	100	100
2	B	146/146 (100%)	144 (99%)	2 (1%)	0	100	100
2	D	144/146 (99%)	142 (99%)	2 (1%)	0	100	100
All	All	570/574 (99%)	562 (99%)	8 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	114/113 (101%)	112 (98%)	2 (2%)	66	49
1	C	114/113 (101%)	110 (96%)	4 (4%)	43	20
2	B	120/118 (102%)	119 (99%)	1 (1%)	86	79
2	D	118/118 (100%)	116 (98%)	2 (2%)	68	51
All	All	466/462 (101%)	457 (98%)	9 (2%)	65	46

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	LYS
1	A	109	LEU

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Mol	Chain	Res	Type
2	B	104	ARG
1	C	1	MET
1	C	16	LYS
1	C	52	SER
1	C	78	ASN
2	D	66	LYS
2	D	68	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	HIS
1	A	103	HIS
2	B	63	GLN
2	B	102	ASN
2	B	108	ASN
1	C	72	HIS
1	C	97	ASN
1	C	103	HIS
2	D	102	ASN

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](#)

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
4	HEM	A	142	1	30,50,50	2.45	9 (30%)	24,82,82	2.52	10 (41%)
3	SO4	B	1148	-	4,4,4	1.03	0	6,6,6	0.14	0
3	SO4	B	1200	-	4,4,4	1.08	0	6,6,6	0.37	0
4	HEM	B	147	2,5	30,50,50	2.54	9 (30%)	24,82,82	2.64	11 (45%)
5	CMO	B	148	4	0,1,1	0.00	-	0,0,0	0.00	-
4	HEM	C	142	1	30,50,50	2.63	9 (30%)	24,82,82	2.59	11 (45%)
4	HEM	D	147	2,5	30,50,50	2.53	7 (23%)	24,82,82	2.80	10 (41%)
5	CMO	D	148	4	0,1,1	0.00	-	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
4	HEM	A	142	1	-	0/10/54/54	0/0/8/8
3	SO4	B	1148	-	-	0/0/0/0	0/0/0/0
3	SO4	B	1200	-	-	0/0/0/0	0/0/0/0
4	HEM	B	147	2,5	-	0/10/54/54	0/0/8/8
5	CMO	B	148	4	-	0/0/0/0	0/0/0/0
4	HEM	C	142	1	-	0/10/54/54	0/0/8/8
4	HEM	D	147	2,5	-	0/10/54/54	0/0/8/8
5	CMO	D	148	4	-	0/0/0/0	0/0/0/0

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	142	HEM	C2D-C3D	-7.24	1.32	1.54
4	D	147	HEM	C2D-C3D	-7.13	1.33	1.54
4	B	147	HEM	C3B-C4B	-7.07	1.45	1.51
4	C	142	HEM	C3B-C4B	-6.98	1.45	1.51
4	D	147	HEM	C3B-C4B	-6.84	1.45	1.51
4	C	142	HEM	C2D-C3D	-6.50	1.35	1.54
4	B	147	HEM	C2D-C3D	-6.37	1.35	1.54
4	C	142	HEM	C3D-C4D	-6.23	1.43	1.51
4	A	142	HEM	C3B-C4B	-6.19	1.46	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	147	HEM	C3D-C4D	-5.51	1.44	1.51
4	B	147	HEM	C3D-C4D	-5.41	1.44	1.51
4	A	142	HEM	C3D-C4D	-5.35	1.44	1.51
4	B	147	HEM	C2C-C1C	-4.16	1.44	1.52
4	D	147	HEM	C2C-C1C	-4.10	1.44	1.52
4	C	142	HEM	C2C-C1C	-3.94	1.45	1.52
4	A	142	HEM	C2C-C1C	-3.39	1.46	1.52
4	C	142	HEM	C2B-C1B	-2.54	1.43	1.51
4	A	142	HEM	C2B-C1B	-2.47	1.43	1.51
4	B	147	HEM	C2D-C1D	-2.21	1.44	1.51
4	B	147	HEM	C2B-C1B	-2.12	1.44	1.51
4	C	142	HEM	C2D-C1D	-2.02	1.45	1.51
4	A	142	HEM	C4C-NC	2.03	1.38	1.36
4	D	147	HEM	CMA-C3A	2.03	1.55	1.51
4	D	147	HEM	CAA-C2A	2.11	1.55	1.52
4	A	142	HEM	CAA-C2A	2.14	1.55	1.52
4	C	142	HEM	C4C-NC	2.23	1.38	1.36
4	A	142	HEM	C1C-NC	2.32	1.38	1.36
4	A	142	HEM	FE-NC	2.40	2.05	1.95
4	C	142	HEM	FE-NC	2.42	2.05	1.95
4	B	147	HEM	C4C-NC	2.74	1.39	1.36
4	B	147	HEM	CAA-C2A	2.78	1.56	1.52
4	D	147	HEM	C1C-NC	3.34	1.40	1.36
4	B	147	HEM	C1C-NC	3.70	1.40	1.36
4	C	142	HEM	C1C-NC	3.75	1.40	1.36

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	147	HEM	CAA-C2A-C1A	-5.90	120.61	127.01
4	B	147	HEM	CAA-C2A-C1A	-4.57	122.04	127.01
4	C	142	HEM	CAA-C2A-C1A	-3.86	122.82	127.01
4	A	142	HEM	C3B-CAB-CBB	-3.45	119.16	124.46
4	C	142	HEM	C3C-CAC-CBC	-3.29	119.40	124.46
4	A	142	HEM	CAA-C2A-C1A	-3.17	123.57	127.01
4	A	142	HEM	CMA-C3A-C4A	-3.02	123.37	128.36
4	C	142	HEM	C3B-CAB-CBB	-2.99	119.87	124.46
4	B	147	HEM	C3B-CAB-CBB	-2.41	120.75	124.46
4	D	147	HEM	C3B-CAB-CBB	-2.34	120.86	124.46
4	B	147	HEM	CBD-CAD-C3D	-2.29	106.90	113.55
4	D	147	HEM	CBA-CAA-C2A	-2.19	108.61	112.53
4	C	142	HEM	C4B-CHC-C1C	-2.13	122.26	125.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	142	HEM	CMA-C3A-C4A	-2.09	124.90	128.36
4	B	147	HEM	CBA-CAA-C2A	-2.02	108.91	112.53
4	A	142	HEM	CMA-C3A-C2A	2.10	129.63	125.24
4	A	142	HEM	CAD-C3D-C4D	3.01	123.09	112.47
4	B	147	HEM	CAA-CBA-CGA	3.11	118.44	112.75
4	B	147	HEM	CAD-C3D-C4D	3.11	123.45	112.47
4	B	147	HEM	CMD-C2D-C3D	3.24	128.70	114.35
4	C	142	HEM	CAD-C3D-C4D	3.28	124.05	112.47
4	D	147	HEM	CAD-C3D-C4D	3.29	124.09	112.47
4	C	142	HEM	CMD-C2D-C3D	3.30	128.94	114.35
4	A	142	HEM	CMD-C2D-C3D	3.40	129.40	114.35
4	D	147	HEM	CMD-C2D-C3D	3.67	130.58	114.35
4	D	147	HEM	CAA-CBA-CGA	3.83	119.76	112.75
4	B	147	HEM	C2D-C3D-C4D	4.03	108.33	101.50
4	A	142	HEM	C2D-C3D-C4D	4.18	108.59	101.50
4	A	142	HEM	CMC-C2C-C3C	4.21	127.03	116.53
4	C	142	HEM	C2D-C3D-C4D	4.27	108.74	101.50
4	C	142	HEM	CMC-C2C-C3C	4.42	127.58	116.53
4	A	142	HEM	CMB-C2B-C3B	4.48	127.72	116.53
4	D	147	HEM	C2D-C3D-C4D	4.51	109.15	101.50
4	C	142	HEM	CMB-C2B-C3B	4.56	127.92	116.53
4	B	147	HEM	CMC-C2C-C3C	4.67	128.19	116.53
4	D	147	HEM	CAD-C3D-C2D	4.71	126.76	113.22
4	D	147	HEM	CMB-C2B-C3B	4.80	128.51	116.53
4	D	147	HEM	CMC-C2C-C3C	4.81	128.54	116.53
4	C	142	HEM	CAD-C3D-C2D	4.87	127.21	113.22
4	B	147	HEM	CMB-C2B-C3B	4.94	128.87	116.53
4	B	147	HEM	CAD-C3D-C2D	5.20	128.17	113.22
4	A	142	HEM	CAD-C3D-C2D	5.25	128.32	113.22

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	147	HEM	1	0
5	B	148	CMO	2	0
4	D	147	HEM	1	0
5	D	148	CMO	4	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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.