



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:24 AM GMT

PDB ID : 2CYM
Title : EFFECTS OF AMINO ACID SUBSTITUTION ON THREE-DIMENSIONAL STRUCTURE: AN X-RAY ANALYSIS OF CYTOCHROME C3 FROM DESULFOVIBRIO VULGARIS HILDENBOROUGH AT 2 ANGSTROMS RESOLUTION
Authors : Morimoto, Y.; Tani, T.; Okumura, H.; Higuchi, Y.; Yasuoka, N.
Deposited on : 1993-09-29
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 (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

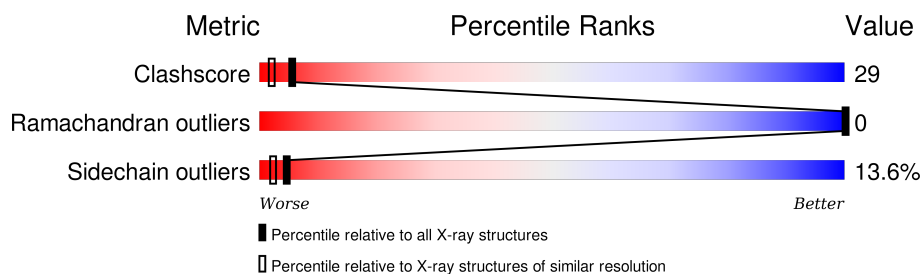
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (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	107	

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1031 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 CYTOCHROME C3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	107	Total	C	N	O	S	0	0	0
			810	495	152	152	11			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 3 is water.

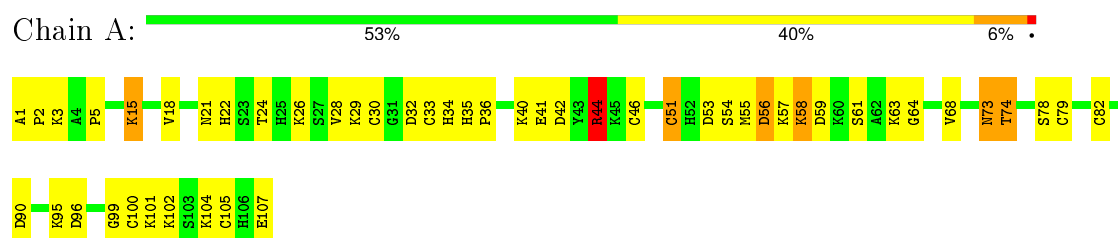
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	49	Total	O	0	0
			49	49		

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: CYTOCHROME C3



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 21	Depositor
Cell constants a, b, c, α , β , γ	43.17Å 62.91Å 41.17Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.00)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.212 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	1031	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM

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	1.03	0/827	1.62	14/1101 (1.3%)

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	2

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	90	ASP	CB-CG-OD2	9.44	126.80	118.30
1	A	107	GLU	OE1-CD-OE2	8.33	133.30	123.30
1	A	44	ARG	NE-CZ-NH1	7.88	124.24	120.30
1	A	96	ASP	CB-CG-OD1	-6.19	112.73	118.30
1	A	44	ARG	CA-CB-CG	6.01	126.63	113.40
1	A	56	ASP	CB-CG-OD1	5.88	123.59	118.30
1	A	53	ASP	CB-CG-OD1	5.87	123.58	118.30
1	A	44	ARG	NE-CZ-NH2	-5.72	117.44	120.30
1	A	56	ASP	CB-CG-OD2	-5.50	113.35	118.30
1	A	24	THR	CA-CB-OG1	-5.45	97.55	109.00
1	A	32	ASP	CB-CG-OD1	5.27	123.05	118.30
1	A	41	GLU	OE1-CD-OE2	-5.19	117.07	123.30
1	A	34	HIS	CA-CB-CG	-5.11	104.91	113.60
1	A	74	THR	O-C-N	-5.01	114.69	122.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	40	LYS	Mainchain
1	A	51	CYS	Mainchain

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	810	0	800	48	0
2	A	172	0	120	35	0
3	A	49	0	0	17	0
All	All	1031	0	920	56	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 29.

All (56) 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:105:CYS:SG	2:A:108:HEM:CAC	2.48	1.01
1:A:18:VAL:HG11	3:A:148:HOH:O	1.72	0.90
2:A:108:HEM:HAB	3:A:149:HOH:O	1.72	0.89
2:A:108:HEM:HAC	3:A:148:HOH:O	1.71	0.88
1:A:79:CYS:HG	2:A:111:HEM:CAB	1.88	0.87
1:A:1:ALA:HB1	3:A:130:HOH:O	1.76	0.85
1:A:2:PRO:HB3	3:A:131:HOH:O	1.80	0.82
1:A:51:CYS:SG	2:A:109:HEM:CAC	2.69	0.80
1:A:82:CYS:SG	2:A:111:HEM:CAC	2.69	0.80
1:A:100:CYS:SG	2:A:108:HEM:CAB	2.71	0.79
1:A:33:CYS:SG	2:A:110:HEM:CAC	2.73	0.76
1:A:46:CYS:SG	2:A:109:HEM:CAB	2.78	0.72
1:A:79:CYS:SG	2:A:111:HEM:CAB	2.80	0.69
1:A:79:CYS:SG	2:A:111:HEM:C3B	2.86	0.69
1:A:30:CYS:SG	2:A:110:HEM:CAB	2.81	0.69
1:A:28:VAL:HG11	3:A:143:HOH:O	1.93	0.68
1:A:79:CYS:HG	2:A:111:HEM:CBB	2.06	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:ASP:O	1:A:59:ASP:HB3	1.95	0.66
1:A:82:CYS:SG	2:A:111:HEM:C3C	2.90	0.65
2:A:111:HEM:HBC2	2:A:111:HEM:HMC2	1.80	0.63
1:A:79:CYS:SG	2:A:111:HEM:CBB	2.87	0.62
1:A:68:VAL:HA	3:A:115:HOH:O	1.99	0.62
2:A:108:HEM:HHC	3:A:149:HOH:O	1.99	0.61
1:A:105:CYS:SG	2:A:108:HEM:HAC	2.38	0.61
1:A:21:ASN:HB2	3:A:144:HOH:O	2.01	0.60
1:A:82:CYS:SG	2:A:111:HEM:CBC	2.90	0.60
2:A:111:HEM:CBC	3:A:143:HOH:O	2.51	0.57
1:A:46:CYS:SG	2:A:109:HEM:CBB	2.92	0.57
1:A:30:CYS:HG	2:A:110:HEM:CAB	2.20	0.55
1:A:74:THR:HG21	3:A:115:HOH:O	2.06	0.54
1:A:42:ASP:OD1	1:A:44:ARG:HG2	2.07	0.54
1:A:33:CYS:SG	2:A:110:HEM:CBC	2.98	0.52
1:A:51:CYS:SG	2:A:109:HEM:HAC	2.46	0.51
2:A:110:HEM:HBB2	2:A:110:HEM:HMB3	1.92	0.51
1:A:105:CYS:SG	2:A:108:HEM:C3C	3.05	0.50
1:A:33:CYS:SG	2:A:110:HEM:C3C	3.05	0.49
1:A:63:LYS:HD2	3:A:152:HOH:O	2.11	0.49
1:A:78:SER:HB3	3:A:115:HOH:O	2.13	0.48
1:A:105:CYS:SG	2:A:108:HEM:CBC	2.99	0.48
1:A:58:LYS:HD3	1:A:58:LYS:O	2.13	0.48
1:A:73:ASN:O	1:A:73:ASN:ND2	2.47	0.47
1:A:58:LYS:CD	1:A:58:LYS:O	2.63	0.47
1:A:35:HIS:HB2	1:A:36:PRO:CD	2.45	0.46
1:A:46:CYS:SG	2:A:109:HEM:C3B	3.10	0.45
1:A:54:SER:HB3	1:A:64:GLY:HA2	1.99	0.44
1:A:5:PRO:HD2	1:A:22:HIS:CE1	2.52	0.44
1:A:15:LYS:HE2	1:A:15:LYS:HB2	1.63	0.43
1:A:55:MET:HE3	3:A:153:HOH:O	2.18	0.43
1:A:99:GLY:C	3:A:149:HOH:O	2.58	0.42
1:A:5:PRO:HD3	2:A:110:HEM:C4A	2.54	0.42
1:A:30:CYS:SG	2:A:110:HEM:C3B	3.12	0.42
1:A:15:LYS:HD2	3:A:133:HOH:O	2.19	0.42
2:A:111:HEM:HMB1	3:A:148:HOH:O	2.20	0.42
1:A:104:LYS:HD3	2:A:111:HEM:HAA2	2.02	0.41
1:A:82:CYS:HG	2:A:111:HEM:CAC	2.24	0.41
2:A:110:HEM:CMB	2:A:110:HEM:HBB2	2.52	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	105/107 (98%)	98 (93%)	7 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	88/88 (100%)	76 (86%)	12 (14%)	5	2

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	15	LYS
1	A	26	LYS
1	A	29	LYS
1	A	44	ARG
1	A	57	LYS
1	A	58	LYS
1	A	61	SER
1	A	73	ASN
1	A	95	LYS
1	A	101	LYS
1	A	102	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

4 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	$\# Z > 2$
2	HEM	A	108	1	30,50,50	2.42	8 (26%)	24,82,82	3.13	15 (62%)
2	HEM	A	109	1	30,50,50	2.51	8 (26%)	24,82,82	3.15	12 (50%)
2	HEM	A	110	1	30,50,50	2.74	12 (40%)	24,82,82	2.72	12 (50%)
2	HEM	A	111	1	30,50,50	2.72	11 (36%)	24,82,82	2.76	10 (41%)

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
2	HEM	A	108	1	-	0/10/54/54	0/0/8/8
2	HEM	A	109	1	-	0/10/54/54	0/0/8/8

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	110	1	-	0/10/54/54	0/0/8/8
2	HEM	A	111	1	-	0/10/54/54	0/0/8/8

All (39) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	110	HEM	C3B-C4B	-7.66	1.45	1.51
2	A	109	HEM	C2D-C3D	-7.41	1.32	1.54
2	A	111	HEM	C3B-C4B	-7.31	1.45	1.51
2	A	111	HEM	C2D-C3D	-6.94	1.33	1.54
2	A	110	HEM	C2D-C3D	-6.52	1.35	1.54
2	A	108	HEM	C2D-C3D	-6.23	1.35	1.54
2	A	109	HEM	C3B-C4B	-6.14	1.46	1.51
2	A	110	HEM	C3D-C4D	-5.76	1.44	1.51
2	A	109	HEM	C3D-C4D	-5.67	1.44	1.51
2	A	111	HEM	C3D-C4D	-5.63	1.44	1.51
2	A	108	HEM	C3D-C4D	-5.46	1.44	1.51
2	A	108	HEM	C2C-C1C	-4.97	1.43	1.52
2	A	111	HEM	C2C-C1C	-4.76	1.43	1.52
2	A	108	HEM	C3B-C4B	-4.73	1.47	1.51
2	A	109	HEM	C2C-C1C	-3.71	1.45	1.52
2	A	110	HEM	C2C-C1C	-3.53	1.45	1.52
2	A	110	HEM	C2D-C1D	-2.51	1.43	1.51
2	A	109	HEM	C2D-C1D	-2.44	1.43	1.51
2	A	110	HEM	C2B-C1B	-2.40	1.44	1.51
2	A	111	HEM	C2D-C1D	-2.14	1.44	1.51
2	A	111	HEM	C2B-C1B	-2.12	1.44	1.51
2	A	108	HEM	C2B-C1B	-2.02	1.45	1.51
2	A	108	HEM	C3B-CAB	2.02	1.55	1.51
2	A	111	HEM	CMB-C2B	2.03	1.57	1.53
2	A	111	HEM	CMA-C3A	2.04	1.55	1.51
2	A	110	HEM	C4C-NC	2.15	1.38	1.36
2	A	109	HEM	CMA-C3A	2.17	1.56	1.51
2	A	111	HEM	FE-NC	2.20	2.04	1.95
2	A	109	HEM	C1C-NC	2.26	1.38	1.36
2	A	110	HEM	CMA-C3A	2.29	1.56	1.51
2	A	110	HEM	C1C-NC	2.33	1.38	1.36
2	A	108	HEM	C4C-NC	2.41	1.39	1.36
2	A	111	HEM	CAA-C2A	2.47	1.56	1.52
2	A	109	HEM	C4C-NC	3.00	1.39	1.36
2	A	110	HEM	CAA-C2A	3.23	1.57	1.52
2	A	110	HEM	C3B-CAB	3.29	1.57	1.51

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	111	HEM	C3B-CAB	3.36	1.57	1.51
2	A	110	HEM	C3C-CAC	4.11	1.59	1.51
2	A	108	HEM	C1C-NC	4.83	1.42	1.36

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	109	HEM	CAA-C2A-C1A	-6.40	120.06	127.01
2	A	111	HEM	CMA-C3A-C4A	-5.84	118.70	128.36
2	A	108	HEM	CMA-C3A-C4A	-5.58	119.14	128.36
2	A	111	HEM	CAA-C2A-C1A	-4.88	121.71	127.01
2	A	110	HEM	CAA-C2A-C1A	-4.16	122.49	127.01
2	A	110	HEM	C3B-CAB-CBB	-3.52	119.05	124.46
2	A	109	HEM	CBD-CAD-C3D	-3.40	103.65	113.55
2	A	110	HEM	CMA-C3A-C4A	-3.01	123.38	128.36
2	A	108	HEM	CAA-C2A-C1A	-2.93	123.82	127.01
2	A	111	HEM	C3C-CAC-CBC	-2.58	120.50	124.46
2	A	109	HEM	CBA-CAA-C2A	-2.40	108.23	112.53
2	A	110	HEM	C1D-CHD-C4C	-2.19	122.17	125.82
2	A	108	HEM	C4B-CHC-C1C	2.05	129.25	125.82
2	A	110	HEM	CAA-CBA-CGA	2.17	116.72	112.75
2	A	109	HEM	CAD-C3D-C4D	2.27	120.46	112.47
2	A	108	HEM	CBD-CAD-C3D	2.34	120.38	113.55
2	A	109	HEM	C3B-C4B-CHC	2.39	126.53	123.16
2	A	110	HEM	C3B-C4B-CHC	2.49	126.67	123.16
2	A	108	HEM	CBA-CAA-C2A	2.49	117.00	112.53
2	A	109	HEM	C2C-C1C-CHC	2.50	127.49	123.68
2	A	111	HEM	CMD-C2D-C3D	2.61	125.88	114.35
2	A	111	HEM	CAD-C3D-C4D	2.71	122.03	112.47
2	A	111	HEM	CMA-C3A-C2A	2.93	131.37	125.24
2	A	111	HEM	C3B-CAB-CBB	2.97	129.01	124.46
2	A	110	HEM	CAD-C3D-C4D	3.00	123.04	112.47
2	A	109	HEM	CMD-C2D-C3D	3.37	129.26	114.35
2	A	108	HEM	CAD-C3D-C4D	3.39	124.44	112.47
2	A	108	HEM	CMA-C3A-C2A	3.55	132.65	125.24
2	A	108	HEM	C2D-C3D-C4D	3.57	107.55	101.50
2	A	111	HEM	C2D-C3D-C4D	3.63	107.65	101.50
2	A	108	HEM	CMD-C2D-C3D	3.67	130.56	114.35
2	A	110	HEM	CMD-C2D-C3D	3.71	130.74	114.35
2	A	108	HEM	C3B-C4B-CHC	3.73	128.42	123.16
2	A	108	HEM	C2C-C1C-CHC	3.77	129.42	123.68
2	A	110	HEM	C2D-C3D-C4D	3.99	108.27	101.50

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	108	HEM	C3C-CAC-CBC	4.29	131.03	124.46
2	A	110	HEM	CMC-C2C-C3C	4.30	127.25	116.53
2	A	109	HEM	CMC-C2C-C3C	4.47	127.68	116.53
2	A	110	HEM	CMB-C2B-C3B	4.51	127.79	116.53
2	A	108	HEM	CMC-C2C-C3C	4.52	127.82	116.53
2	A	109	HEM	C2D-C3D-C4D	4.59	109.28	101.50
2	A	109	HEM	CMB-C2B-C3B	4.95	128.90	116.53
2	A	108	HEM	CMB-C2B-C3B	5.00	129.01	116.53
2	A	108	HEM	CAD-C3D-C2D	5.14	128.00	113.22
2	A	110	HEM	CAD-C3D-C2D	5.37	128.66	113.22
2	A	109	HEM	CAD-C3D-C2D	5.47	128.94	113.22
2	A	111	HEM	CMC-C2C-C3C	5.61	130.54	116.53
2	A	111	HEM	CAD-C3D-C2D	5.77	129.81	113.22
2	A	109	HEM	C3B-CAB-CBB	6.75	134.81	124.46

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	108	HEM	8	0
2	A	109	HEM	5	0
2	A	110	HEM	9	0
2	A	111	HEM	13	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.