



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 26, 2014 – 02:24 PM GMT

PDB ID : 1VYD
Title : Crystal structure of cytochrome c2 mutant G95E
Authors : Dumortier, C.; Fitch, J.; Van Petegem, F.; Vermeulen, W.; Meyer, T.E.; Van Beeumen, J.J.; Cusanovich, M.A.
Deposited on : 2004-04-27
Resolution : 2.30 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

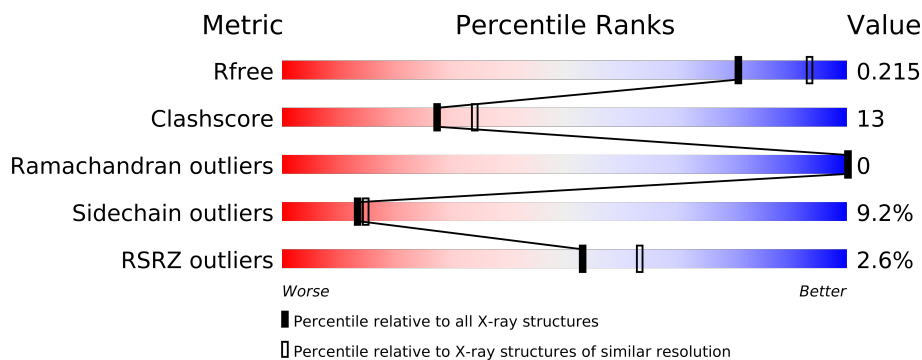
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.30 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))
R_{free}	66092	2929 (2.30-2.30)
Clashscore	79885	3679 (2.30-2.30)
Ramachandran outliers	78287	3642 (2.30-2.30)
Sidechain outliers	78261	3641 (2.30-2.30)
RSRZ outliers	66119	2930 (2.30-2.30)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	116	
1	B	116	

2 Entry composition i

There are 3 unique types of molecules in this entry. The entry contains 1904 atoms, of which 0 are hydrogen and 0 are deuterium.

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 C2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	116	Total	C	N	O	S	0	0	0
			868	555	141	169	3			
1	B	116	Total	C	N	O	S	0	0	0
			868	555	141	169	3			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	95	GLU	GLY	ENGINEERED MUTATION	UNP P00094
B	95	GLU	GLY	ENGINEERED MUTATION	UNP P00094

- 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		

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

- Molecule 3 is water.

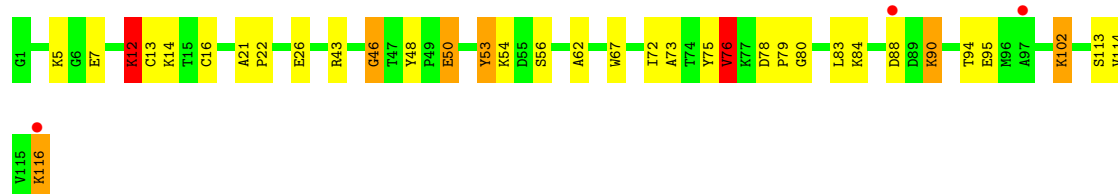
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	39	Total	O	0	0
			39	39		
3	B	43	Total	O	0	0
			43	43		

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.

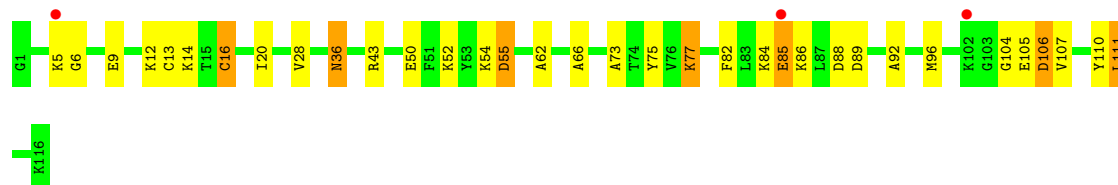
• Molecule 1: CYTOCHROME C2

Chain A: 



• Molecule 1: CYTOCHROME C2

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	100.23Å 100.23Å 162.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	10.00 – 2.30 9.93 – 2.30	Depositor EDS
% Data completeness (in resolution range)	88.0 (10.00-2.30) 88.9 (9.93-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.34 (at 2.31Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.198 , 0.242 0.178 , 0.215	Depositor DCC
R_{free} test set	601 reflections (5.07%)	DCC
Wilson B-factor (Å ²)	36.0	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 50.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 12459 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	1904	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.98% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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	0.88	0/884	1.88	18/1188 (1.5%)
1	B	0.84	0/884	1.78	12/1188 (1.0%)
All	All	0.86	0/1768	1.83	30/2376 (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	3
1	B	0	3
All	All	0	6

There are no bond length outliers.

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	43	ARG	NE-CZ-NH2	-15.22	112.69	120.30
1	A	43	ARG	NE-CZ-NH2	14.28	127.44	120.30
1	B	55	ASP	CB-CG-OD2	12.80	129.82	118.30
1	A	78	ASP	CB-CG-OD2	11.28	128.45	118.30
1	A	43	ARG	NH1-CZ-NH2	-10.29	108.08	119.40
1	B	50	GLU	OE1-CD-OE2	-9.56	111.83	123.30
1	A	43	ARG	NE-CZ-NH1	8.35	124.47	120.30
1	A	48	TYR	CB-CG-CD2	-7.57	116.46	121.00
1	A	43	ARG	CD-NE-CZ	6.76	133.07	123.60
1	A	53	TYR	CB-CG-CD2	-6.72	116.97	121.00
1	A	78	ASP	CB-CG-OD1	-6.46	112.49	118.30
1	A	73	ALA	CB-CA-C	-5.99	101.12	110.10

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	89	ASP	CB-CG-OD2	5.95	123.65	118.30
1	A	26	GLU	OE1-CD-OE2	-5.81	116.33	123.30
1	A	76	VAL	N-CA-CB	-5.78	98.78	111.50
1	A	7	GLU	OE1-CD-OE2	-5.69	116.48	123.30
1	B	28	VAL	CA-CB-CG2	5.68	119.42	110.90
1	B	106	ASP	CB-CG-OD1	-5.64	113.22	118.30
1	B	110	TYR	CB-CG-CD2	-5.63	117.62	121.00
1	A	50	GLU	O-C-N	-5.56	113.81	122.70
1	A	12	LYS	CG-CD-CE	5.50	128.41	111.90
1	B	110	TYR	CB-CG-CD1	5.42	124.25	121.00
1	B	88	ASP	CB-CG-OD2	5.39	123.15	118.30
1	B	43	ARG	NH1-CZ-NH2	5.38	125.32	119.40
1	B	111	LEU	CA-CB-CG	5.24	127.34	115.30
1	A	76	VAL	O-C-N	-5.21	114.37	122.70
1	B	75	TYR	CB-CG-CD2	-5.20	117.88	121.00
1	A	76	VAL	CA-CB-CG1	5.19	118.68	110.90
1	A	56	SER	CB-CA-C	-5.16	100.30	110.10
1	A	102	LYS	CA-CB-CG	5.02	124.45	113.40

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	12	LYS	Mainchain
1	A	46	GLY	Mainchain
1	A	67	TRP	Mainchain
1	B	16	CYS	Mainchain
1	B	55	ASP	Mainchain
1	B	66	ALA	Mainchain

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	868	0	876	23	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	868	0	876	25	0
2	A	43	0	30	8	0
2	B	43	0	30	11	0
3	A	39	0	0	2	0
3	B	43	0	0	0	0
All	All	1904	0	1812	47	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 13.

All (47) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:13:CYS:SG	2:A:1117:HEM:CAB	2.31	1.19
1:A:16:CYS:SG	2:A:1117:HEM:CAC	2.39	1.10
1:A:13:CYS:HG	2:A:1117:HEM:CAB	1.64	1.08
1:B:16:CYS:SG	2:B:1117:HEM:CAC	2.49	1.01
1:B:13:CYS:SG	2:B:1117:HEM:CAB	2.51	0.98
1:A:13:CYS:SG	2:A:1117:HEM:HAB	2.02	0.96
1:A:54:LYS:HG3	1:A:95:GLU:HG2	1.58	0.86
1:A:16:CYS:SG	2:A:1117:HEM:HAC	2.25	0.76
1:B:20:ILE:H	1:B:36:ASN:HD21	1.33	0.75
1:B:13:CYS:HG	2:B:1117:HEM:CAB	2.04	0.71
2:B:1117:HEM:HMB1	2:B:1117:HEM:HBB2	1.73	0.69
1:B:20:ILE:H	1:B:36:ASN:ND2	1.91	0.67
1:A:113:SER:HA	1:A:116:LYS:HD2	1.85	0.58
1:B:16:CYS:SG	2:B:1117:HEM:C3C	2.98	0.56
1:A:95:GLU:HG3	3:A:2039:HOH:O	2.06	0.55
1:B:85:GLU:HG3	1:B:86:LYS:N	2.23	0.54
1:B:84:LYS:HG2	1:B:92:ALA:HB3	1.89	0.54
1:B:5:LYS:HD3	1:B:106:ASP:OD1	2.09	0.53
1:A:116:LYS:HA	3:A:2037:HOH:O	2.08	0.53
1:A:16:CYS:SG	2:A:1117:HEM:CBC	2.96	0.52
1:A:16:CYS:SG	2:A:1117:HEM:C3C	3.03	0.51
1:B:13:CYS:SG	2:B:1117:HEM:C3B	3.04	0.51
1:B:13:CYS:SG	2:B:1117:HEM:CBB	2.98	0.51
1:A:50:GLU:HA	1:B:62:ALA:O	2.13	0.49
1:B:16:CYS:SG	2:B:1117:HEM:CBC	3.01	0.48
1:A:72:ILE:O	1:A:76:VAL:HB	2.14	0.48
1:A:54:LYS:HG3	1:A:95:GLU:CG	2.36	0.47
1:B:9:GLU:OE1	1:B:12:LYS:NZ	2.41	0.46
1:B:16:CYS:SG	2:B:1117:HEM:HAC	2.51	0.45
1:B:6:GLY:HA3	1:B:106:ASP:O	2.17	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:73:ALA:O	1:B:77:LYS:HD2	2.17	0.45
1:A:62:ALA:HB3	1:B:52:LYS:HE2	1.98	0.44
1:A:21:ALA:HB1	1:A:22:PRO:HD2	2.01	0.43
1:A:75:TYR:CE1	1:A:79:PRO:HB3	2.54	0.43
1:B:96:MET:HG3	2:B:1117:HEM:C4C	2.54	0.43
1:B:104:GLY:O	1:B:107:VAL:HG12	2.19	0.42
1:B:13:CYS:SG	2:B:1117:HEM:HAB	2.54	0.42
1:B:82:PHE:O	1:B:85:GLU:HG3	2.19	0.42
1:A:46:GLY:HA2	1:A:53:TYR:CZ	2.54	0.42
1:A:13:CYS:SG	2:A:1117:HEM:C3B	3.03	0.42
1:A:80:GLY:O	1:A:84:LYS:HG3	2.20	0.42
1:B:77:LYS:HZ1	1:B:104:GLY:H	1.67	0.42
1:A:113:SER:CA	1:A:116:LYS:HD2	2.50	0.41
1:A:83:LEU:HD12	1:A:94:THR:HB	2.01	0.41
1:B:84:LYS:HE3	1:B:92:ALA:O	2.21	0.41
1:B:85:GLU:HG3	1:B:86:LYS:H	1.85	0.41
1:A:90:LYS:HB2	1:A:90:LYS:HE3	1.96	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	114/116 (98%)	111 (97%)	3 (3%)	0	100	100
1	B	114/116 (98%)	113 (99%)	1 (1%)	0	100	100
All	All	228/232 (98%)	224 (98%)	4 (2%)	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	87/87 (100%)	78 (90%)	9 (10%)	10	11
1	B	87/87 (100%)	80 (92%)	7 (8%)	17	20
All	All	174/174 (100%)	158 (91%)	16 (9%)	13	15

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LYS
1	A	12	LYS
1	A	14	LYS
1	A	76	VAL
1	A	88	ASP
1	A	90	LYS
1	A	102	LYS
1	A	114	VAL
1	A	116	LYS
1	B	14	LYS
1	B	36	ASN
1	B	54	LYS
1	B	77	LYS
1	B	85	GLU
1	B	105	GLU
1	B	111	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	36	ASN

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

2 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	1117	1	49,50,50	2.16	11 (22%)	46,82,82	2.91	17 (36%)
2	HEM	B	1117	1	49,50,50	3.10	17 (34%)	46,82,82	2.56	15 (32%)

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	1117	1	-	0/14/114/114	0/0/8/8
2	HEM	B	1117	1	-	0/14/114/114	0/0/8/8

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1117	HEM	C2D-C1D	13.67	1.48	1.44
2	A	1117	HEM	C3C-C2C	-6.59	1.32	1.43
2	B	1117	HEM	C3D-C2D	-5.95	1.33	1.43
2	B	1117	HEM	C3C-C2C	-5.85	1.33	1.43
2	A	1117	HEM	C3B-C2B	-5.33	1.34	1.43
2	B	1117	HEM	C3C-CAC	5.24	1.56	1.40
2	B	1117	HEM	C3B-C2B	-5.23	1.34	1.43
2	A	1117	HEM	C3D-C2D	-4.98	1.35	1.43
2	B	1117	HEM	C3B-CAB	4.93	1.55	1.40
2	A	1117	HEM	C3B-CAB	4.75	1.55	1.40
2	A	1117	HEM	C3C-CAC	4.56	1.54	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1117	HEM	CHA-C4D	4.09	1.41	1.35
2	B	1117	HEM	C3D-C4D	-3.84	1.43	1.44
2	B	1117	HEM	C2B-C1B	-3.74	1.43	1.44
2	B	1117	HEM	C4A-C3A	3.56	1.44	1.40
2	B	1117	HEM	CMC-C2C	3.49	1.58	1.47
2	A	1117	HEM	CMB-C2B	3.15	1.57	1.47
2	A	1117	HEM	CAA-C2A	2.99	1.57	1.52
2	B	1117	HEM	CMD-C2D	2.81	1.56	1.47
2	A	1117	HEM	C4A-C3A	2.70	1.43	1.40
2	B	1117	HEM	FE-NA	2.53	2.03	1.92
2	A	1117	HEM	CHA-C4D	2.46	1.39	1.35
2	B	1117	HEM	CHB-C1B	2.43	1.39	1.35
2	A	1117	HEM	CMC-C2C	2.26	1.54	1.47
2	B	1117	HEM	CMB-C2B	2.25	1.54	1.47
2	B	1117	HEM	CHD-C4C	2.25	1.40	1.36
2	B	1117	HEM	O1A-CGA	2.23	1.30	1.22
2	A	1117	HEM	CMD-C2D	2.18	1.54	1.47

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1117	HEM	C3B-C4B-NB	-9.04	107.53	114.00
2	B	1117	HEM	C3B-C4B-NB	-8.23	108.11	114.00
2	A	1117	HEM	CMA-C3A-C4A	-7.42	117.20	128.62
2	A	1117	HEM	C4A-CHB-C1B	-6.79	118.54	127.47
2	A	1117	HEM	CHC-C4B-NB	6.74	130.18	124.58
2	B	1117	HEM	CHD-C4C-NC	5.82	129.79	124.73
2	B	1117	HEM	CAD-CBD-CGD	5.64	131.07	113.48
2	B	1117	HEM	CHD-C1D-ND	5.50	129.15	124.58
2	A	1117	HEM	CMA-C3A-C2A	5.27	134.88	124.94
2	B	1117	HEM	CMA-C3A-C4A	-4.00	122.47	128.62
2	A	1117	HEM	CBD-CAD-C3D	-3.68	106.34	114.37
2	B	1117	HEM	C1D-CHD-C4C	-3.67	116.91	126.57
2	A	1117	HEM	C1A-CHA-C4D	-3.59	122.74	127.47
2	A	1117	HEM	CHB-C4A-NA	3.59	130.57	124.58
2	B	1117	HEM	C4A-CHB-C1B	-3.58	122.76	127.47
2	B	1117	HEM	CHC-C4B-NB	3.21	127.25	124.58
2	B	1117	HEM	C1A-CHA-C4D	-3.13	123.35	127.47
2	A	1117	HEM	O1D-CGD-CBD	-3.04	112.56	123.03
2	B	1117	HEM	C4C-NC-C1C	3.03	108.69	105.53
2	A	1117	HEM	C2D-C1D-ND	-2.93	109.47	112.93
2	A	1117	HEM	C3A-C4A-CHB	-2.82	120.64	126.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	1117	HEM	CMA-C3A-C2A	2.79	130.21	124.94
2	A	1117	HEM	CHB-C1B-NB	2.61	127.90	124.31
2	B	1117	HEM	CBA-CAA-C2A	-2.45	108.38	112.69
2	A	1117	HEM	C4B-CHC-C1C	-2.43	120.17	126.57
2	A	1117	HEM	CHC-C1C-NC	2.32	126.74	124.73
2	B	1117	HEM	C2D-C1D-ND	-2.31	110.20	112.93
2	A	1117	HEM	O2D-CGD-CBD	2.24	122.14	114.22
2	A	1117	HEM	CBA-CAA-C2A	-2.20	108.82	112.69
2	A	1117	HEM	CMC-C2C-C3C	2.17	131.27	126.16
2	B	1117	HEM	CHC-C1C-NC	2.05	126.52	124.73
2	B	1117	HEM	C3A-C4A-NA	-2.01	107.89	109.41

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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 ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	116/116 (100%)	-0.45	3 (2%) 53 63	20, 34, 51, 60	0
1	B	116/116 (100%)	-0.33	3 (2%) 53 63	22, 38, 59, 63	0
All	All	232/232 (100%)	-0.39	6 (2%) 53 63	20, 35, 57, 63	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	116	LYS	3.9
1	B	102	LYS	2.7
1	A	88	ASP	2.6
1	A	97	ALA	2.3
1	B	85	GLU	2.3
1	B	5	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors

of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	HEM	B	1117	43/43	0.09	-0.20	20,26,34,39	0
2	HEM	A	1117	43/43	0.08	-0.29	21,23,29,32	0

6.5 Other polymers ⓘ

There are no such residues in this entry.