



Full wwPDB X-ray Structure Validation Report

Mar 12, 2014 – 04:42 PM GMT

PDB ID : 4O7G
Title : Crystal Structure of Ascorbate-bound Cytochrome b561, crystal soaked in 1 M L-ascorbate for 40 minutes
Authors : Lu, P.; Ma, D.; Yan, C.; Gong, X.; Du, M.; Shi, Y.
Deposited on : 2013-12-24
Resolution : 2.21 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>

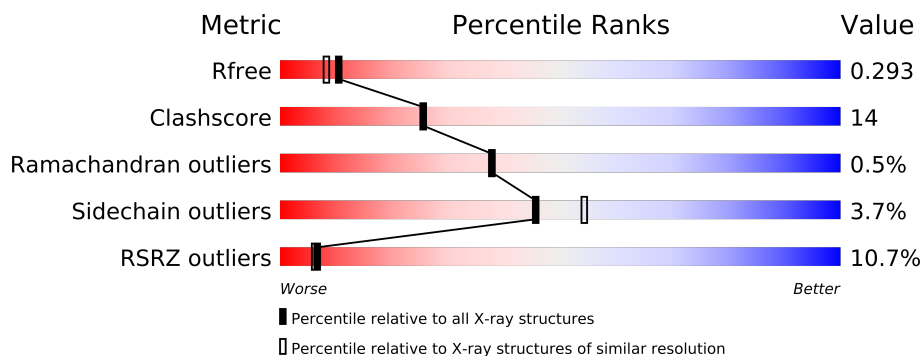
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1323
EDS : trunk22714
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) : trunk22714

1 Overall quality at a glance

The reported resolution of this entry is 2.21 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	3340 (2.24-2.20)
Clashscore	79885	4208 (2.24-2.20)
Ramachandran outliers	78287	4135 (2.24-2.20)
Sidechain outliers	78261	4136 (2.24-2.20)
RSRZ outliers	66119	3341 (2.24-2.20)

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	230	
1	B	230	

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 3575 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 Probable transmembrane ascorbate ferrireductase 2.

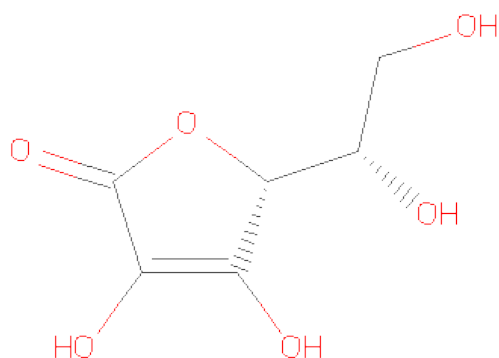
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	211	Total	C	N	O	S	0	0	0
			1652	1107	267	270	8			
1	B	212	Total	C	N	O	S	0	0	0
			1663	1116	268	271	8			

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



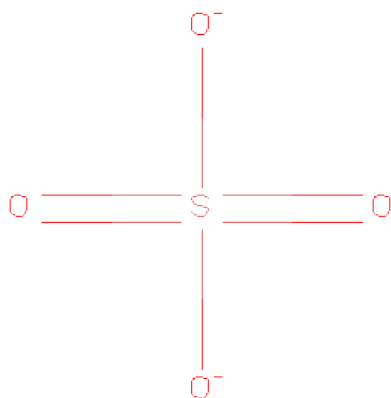
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is ASCORBIC ACID (three-letter code: ASC) (formula: $C_6H_8O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		
3	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 5 is water.

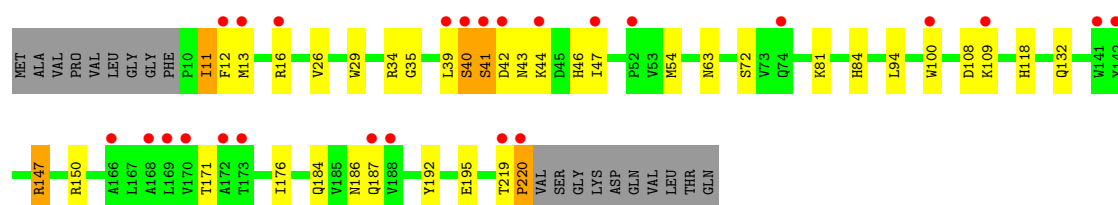
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	25	Total	O	0	0
			25	25		
5	B	22	Total	O	0	0
			22	22		

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 ($\text{RSRZ} > 2$). 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.

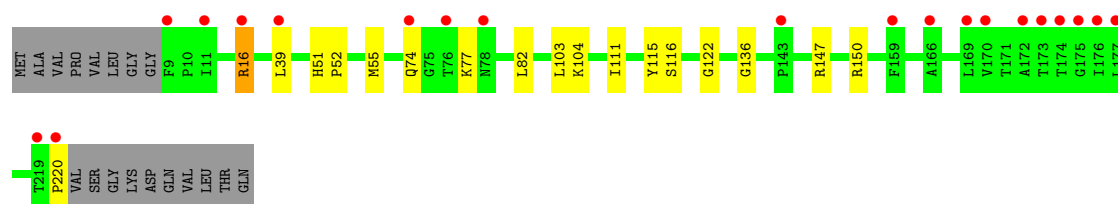
- Molecule 1: Probable transmembrane ascorbate ferrireductase 2

Chain A: 



- Molecule 1: Probable transmembrane ascorbate ferrireductase 2

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	73.19Å 108.70Å 111.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	28.08 – 2.21 28.08 – 2.21	Depositor EDS
% Data completeness (in resolution range)	98.3 (28.08-2.21) 98.3 (28.08-2.21)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.22Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7.3_928)	Depositor
R, R_{free}	0.202 , 0.269 0.242 , 0.293	Depositor DCC
R_{free} test set	1134 reflections (5.09%)	DCC
Wilson B-factor (Å ²)	43.1	Xtriage
Anisotropy	0.309	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 59.9	EDS
Estimated twinning fraction	0.014 for -h,-l,-k	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 22282 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3575	wwPDB-VP
Average B, all atoms (Å ²)	54.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.58% 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: ASC, 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	0.58	0/1696	0.53	2/2310 (0.1%)
1	B	0.62	1/1708 (0.1%)	0.51	0/2327
All	All	0.60	1/3404 (0.0%)	0.52	2/4637 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	220	PRO	N-CD	9.00	1.60	1.47

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	219	THR	C-N-CD	6.82	142.71	128.40
1	A	220	PRO	CA-N-CD	-5.36	104.00	111.50

There are no chirality outliers.

There are no planarity outliers.

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	1652	0	65	31	4

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1663	0	13	14	2
2	A	86	0	0	5	0
2	B	86	0	0	7	0
3	A	12	0	0	1	0
3	B	24	0	0	3	0
4	A	5	0	0	2	2
5	A	25	0	0	3	0
5	B	22	0	0	1	0
All	All	3575	0	78	49	4

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 14.

All (49) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:44:LYS:NZ	1:A:108:ASP:O	1.64	1.27
1:A:39:LEU:HD22	1:A:100:TRP:CZ3	1.74	1.23
1:A:39:LEU:HD22	1:A:100:TRP:HZ3	0.93	1.09
1:A:220:PRO:CB	5:A:425:HOH:O	2.00	1.09
1:A:39:LEU:CD2	1:A:100:TRP:HZ3	1.76	0.98
1:A:81:LYS:NZ	4:A:304:SO4:O4	1.97	0.96
1:B:77:LYS:NZ	3:B:303:ASC:O2	2.03	0.91
1:B:136:GLY:O	2:B:302:HEM:CBC	2.25	0.85
1:A:43:ASN:ND2	1:A:46:HIS:CD2	2.45	0.83
1:B:115:TYR:O	5:B:409:HOH:O	1.96	0.83
1:A:43:ASN:ND2	1:A:46:HIS:NE2	2.27	0.82
1:A:35:GLY:O	1:A:46:HIS:CE1	2.36	0.78
1:A:192:TYR:O	5:A:420:HOH:O	1.99	0.78
1:B:52:PRO:CB	2:B:301:HEM:CBB	2.67	0.72
1:A:43:ASN:CG	1:A:46:HIS:CD2	2.63	0.72
1:A:39:LEU:HB3	1:A:100:TRP:CH2	2.35	0.61
1:A:39:LEU:HB3	1:A:100:TRP:HH2	1.66	0.61
1:B:55:MET:CE	2:B:301:HEM:CAC	2.78	0.61
1:A:40:SER:HA	1:A:47:ILE:HD11	1.84	0.59
1:B:16:ARG:HH11	1:B:16:ARG:CB	2.16	0.58
1:A:54:MET:CE	1:A:94:LEU:CD2	2.80	0.58
1:A:44:LYS:HZ2	1:A:108:ASP:C	2.04	0.57
1:B:122:GLY:CA	2:B:301:HEM:CBC	2.82	0.56
1:A:184:GLN:OE1	5:A:408:HOH:O	2.18	0.55
1:A:150:ARG:NH1	4:A:304:SO4:O1	2.40	0.54
1:A:35:GLY:C	1:A:46:HIS:CE1	2.81	0.54
1:A:118:HIS:CE1	1:A:171:THR:O	2.65	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:44:LYS:NZ	1:A:108:ASP:C	2.56	0.50
1:A:29:TRP:NE1	1:A:195:GLU:OE2	2.44	0.49
3:A:303:ASC:O4	3:A:303:ASC:O6	2.30	0.49
1:B:116:SER:OG	2:B:301:HEM:O2A	2.30	0.49
1:B:103:LEU:CD2	2:B:301:HEM:O2D	2.61	0.49
1:A:39:LEU:O	1:A:40:SER:HB3	2.13	0.48
3:B:304:ASC:O4	3:B:304:ASC:O6	2.30	0.48
1:A:39:LEU:CD2	1:A:100:TRP:CZ3	2.64	0.48
2:A:302:HEM:CBB	2:A:302:HEM:CMB	2.92	0.47
2:A:301:HEM:CBC	2:A:301:HEM:CMC	2.93	0.47
1:A:12:PHE:CE1	1:A:72:SER:CB	3.00	0.45
1:A:176:ILE:CG1	2:A:301:HEM:CBC	2.96	0.44
1:B:150:ARG:NH1	3:B:303:ASC:O1	2.50	0.44
1:A:11:ILE:CG2	1:A:12:PHE:N	2.79	0.43
1:A:43:ASN:CB	1:A:46:HIS:CD2	3.03	0.42
1:B:51:HIS:CB	1:B:52:PRO:CD	2.98	0.42
1:B:16:ARG:CG	1:B:16:ARG:HH11	2.33	0.42
1:B:39:LEU:O	1:B:104:LYS:NZ	2.53	0.41
1:A:84:HIS:CD2	2:A:302:HEM:NC	2.88	0.41
1:A:186:ASN:O	1:A:187:GLN:CB	2.68	0.41
1:A:132:GLN:CG	2:A:302:HEM:CBB	2.98	0.40
1:B:55:MET:CE	2:B:301:HEM:C3C	3.03	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:147:ARG:NH1	4:A:304:SO4:O3[2_565]	1.23	0.97
1:A:42:ASP:OD1	1:B:147:ARG:CD[7_445]	2.08	0.12
1:A:42:ASP:CG	1:B:147:ARG:CD[7_445]	2.12	0.08
1:A:147:ARG:CZ	4:A:304:SO4:O3[2_565]	2.14	0.06

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	209/230 (91%)	198 (95%)	9 (4%)	2 (1%)	22	18
1	B	210/230 (91%)	206 (98%)	4 (2%)	0	100	100
All	All	419/460 (91%)	404 (96%)	13 (3%)	2 (0%)	38	38

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	40	SER
1	A	41	SER

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	176/191 (92%)	167 (95%)	9 (5%)	33	37
1	B	177/191 (93%)	173 (98%)	4 (2%)	63	74
All	All	353/382 (92%)	340 (96%)	13 (4%)	45	54

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ILE
1	A	13	MET
1	A	16	ARG
1	A	26	VAL
1	A	34	ARG
1	A	41	SER
1	A	63	ASN
1	A	109	LYS
1	A	147	ARG
1	B	16	ARG
1	B	74	GLN
1	B	82	LEU
1	B	111	ILE

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

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 ⓘ

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	$\# Z > 2$
2	HEM	A	301	1	50,50,50	2.73	15 (30%)	46,82,82	2.13	7 (15%)
2	HEM	A	302	1	50,50,50	2.74	12 (24%)	46,82,82	2.08	8 (17%)
3	ASC	A	303	-	12,12,12	1.93	1 (8%)	17,17,17	1.82	6 (35%)
4	SO4	A	304	-	4,4,4	0.31	0	6,6,6	0.09	0
2	HEM	B	301	1	50,50,50	2.59	15 (30%)	46,82,82	2.14	12 (26%)
2	HEM	B	302	1	50,50,50	2.73	14 (28%)	46,82,82	2.18	9 (19%)
3	ASC	B	303	-	12,12,12	1.93	1 (8%)	17,17,17	1.88	7 (41%)
3	ASC	B	304	-	12,12,12	1.93	1 (8%)	17,17,17	1.85	6 (35%)

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	301	1	-	0/14/114/114	0/0/8/8

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	302	1	-	0/14/114/114	0/0/8/8
3	ASC	A	303	-	-	0/6/22/22	0/1/1/1
4	SO4	A	304	-	-	0/0/0/0	0/0/0/0
2	HEM	B	301	1	-	0/14/114/114	0/0/8/8
2	HEM	B	302	1	-	0/14/114/114	0/0/8/8
3	ASC	B	303	-	-	0/6/22/22	0/1/1/1
3	ASC	B	304	-	-	0/6/22/22	0/1/1/1

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	302	HEM	C3B-C2B	-9.87	1.32	1.45
2	B	301	HEM	C3C-C2C	-8.95	1.34	1.45
2	B	302	HEM	C3D-C2D	8.28	1.54	1.44
2	A	302	HEM	C3C-C2C	-8.27	1.34	1.45
2	A	301	HEM	C3C-C2C	-8.24	1.34	1.45
2	A	301	HEM	C3B-C2B	-8.11	1.35	1.45
2	B	302	HEM	C3C-C2C	-8.06	1.35	1.45
2	A	301	HEM	C3D-C2D	8.04	1.53	1.44
2	B	301	HEM	C3B-C2B	-7.81	1.35	1.45
2	A	302	HEM	C3D-C2D	7.77	1.53	1.44
2	B	302	HEM	C3B-C2B	-7.75	1.35	1.45
2	B	301	HEM	C3D-C2D	6.67	1.52	1.44
3	A	303	ASC	O4-C1	6.53	1.45	1.35
3	B	304	ASC	O4-C1	6.50	1.45	1.35
3	B	303	ASC	O4-C1	6.48	1.45	1.35
2	B	302	HEM	C3C-CAC	4.97	1.56	1.40
2	B	302	HEM	C3B-CAB	4.82	1.56	1.40
2	A	301	HEM	C3C-CAC	4.74	1.55	1.40
2	A	301	HEM	C3B-CAB	4.61	1.55	1.40
2	B	302	HEM	CBB-CAB	4.58	1.55	1.29
2	A	301	HEM	CBC-CAC	4.53	1.54	1.29
2	B	301	HEM	CBB-CAB	4.51	1.54	1.29
2	A	301	HEM	CBB-CAB	4.51	1.54	1.29
2	A	302	HEM	CBB-CAB	4.51	1.54	1.29
2	B	301	HEM	C3B-CAB	4.49	1.55	1.40
2	A	302	HEM	CBC-CAC	4.49	1.54	1.29
2	B	302	HEM	CBC-CAC	4.44	1.54	1.29
2	B	301	HEM	CBC-CAC	4.38	1.54	1.29
2	A	302	HEM	C3B-CAB	4.28	1.54	1.40
2	B	301	HEM	C3C-CAC	4.24	1.54	1.40
2	A	302	HEM	C3C-CAC	4.00	1.53	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	302	HEM	FE-ND	3.95	2.12	1.96
2	B	302	HEM	FE-NB	3.60	2.10	1.96
2	B	302	HEM	FE-ND	3.54	2.10	1.96
2	B	301	HEM	FE-ND	3.43	2.09	1.96
2	A	301	HEM	FE-NB	3.22	2.09	1.96
2	A	301	HEM	FE-ND	3.13	2.08	1.96
2	A	301	HEM	FE-NC	3.00	2.07	1.95
2	A	301	HEM	CMC-C2C	2.54	1.55	1.47
2	B	301	HEM	CMD-C2D	2.51	1.55	1.47
2	A	302	HEM	CMC-C2C	2.50	1.55	1.47
2	B	301	HEM	CAA-C2A	2.50	1.56	1.52
2	A	302	HEM	CMB-C2B	2.47	1.55	1.47
2	B	301	HEM	CMC-C2C	2.46	1.55	1.47
2	B	302	HEM	CMC-C2C	2.44	1.55	1.47
2	A	301	HEM	CMB-C2B	2.44	1.55	1.47
2	A	301	HEM	CMD-C2D	2.44	1.55	1.47
2	B	302	HEM	CMB-C2B	2.37	1.55	1.47
2	A	302	HEM	CMD-C2D	2.35	1.55	1.47
2	B	302	HEM	CMD-C2D	2.34	1.54	1.47
2	B	301	HEM	FE-NC	2.32	2.05	1.95
2	B	301	HEM	CMB-C2B	2.23	1.54	1.47
2	A	301	HEM	CAA-C2A	2.20	1.55	1.52
2	A	302	HEM	C2D-C1D	-2.20	1.43	1.45
2	B	301	HEM	FE-NA	2.20	2.11	1.94
2	B	302	HEM	CHB-C1B	2.16	1.38	1.35
2	A	301	HEM	CHB-C1B	2.06	1.38	1.35
2	B	302	HEM	CHA-C4D	2.05	1.38	1.35
2	B	301	HEM	CHA-C4D	2.00	1.38	1.35

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	HEM	C3B-C4B-NB	-8.73	107.75	114.00
2	B	302	HEM	C3B-C4B-NB	-8.32	108.04	114.00
2	A	302	HEM	C3B-C4B-NB	-8.17	108.16	114.00
2	B	301	HEM	C3B-C4B-NB	-8.02	108.26	114.00
2	B	302	HEM	C4D-ND-C1D	6.79	112.15	105.11
2	B	301	HEM	C4D-ND-C1D	6.33	111.68	105.11
2	A	302	HEM	C4D-ND-C1D	6.12	111.46	105.11
2	A	301	HEM	C4D-ND-C1D	5.58	110.90	105.11
2	A	301	HEM	C3A-C4A-NA	-4.23	106.67	109.50
2	A	302	HEM	CBA-CAA-C2A	-4.18	105.67	112.63

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	HEM	CAD-CBD-CGD	-3.94	106.10	113.53
3	B	303	ASC	C4-O4-C1	-3.81	105.34	109.32
3	A	303	ASC	C4-O4-C1	-3.77	105.39	109.32
3	B	304	ASC	C4-O4-C1	-3.66	105.50	109.32
2	B	301	HEM	C2D-C1D-ND	-3.60	108.67	112.93
3	B	303	ASC	C5-C4-C3	-3.36	109.50	114.18
2	A	301	HEM	C1B-NB-C4B	3.34	108.58	105.11
2	B	302	HEM	CBA-CAA-C2A	-3.32	107.10	112.63
2	B	302	HEM	C2D-C1D-ND	-3.29	109.04	112.93
2	A	301	HEM	CAD-CBD-CGD	-3.27	107.36	113.53
2	A	301	HEM	C4A-C3A-C2A	3.22	109.23	107.00
2	B	302	HEM	C4A-CHB-C1B	-3.10	123.39	127.47
2	A	301	HEM	C2D-C1D-ND	-3.03	109.35	112.93
2	A	302	HEM	C2D-C1D-ND	-3.02	109.36	112.93
2	A	302	HEM	CBD-CAD-C3D	-2.97	107.88	114.37
3	B	304	ASC	C5-C4-C3	-2.96	110.06	114.18
2	B	301	HEM	C1B-NB-C4B	2.93	108.15	105.11
2	B	302	HEM	C1B-NB-C4B	2.90	108.12	105.11
3	B	304	ASC	O1-C1-C2	-2.79	125.67	129.45
3	A	303	ASC	O1-C1-C2	-2.75	125.72	129.45
2	B	301	HEM	CMA-C3A-C4A	-2.72	124.28	128.46
2	A	302	HEM	CMA-C3A-C4A	-2.71	124.30	128.46
3	B	304	ASC	O3-C3-C2	-2.69	124.28	132.36
3	A	303	ASC	O3-C3-C2	-2.67	124.33	132.36
2	B	301	HEM	C3A-C4A-NA	-2.66	107.72	109.50
3	B	304	ASC	O4-C1-O1	2.61	124.23	121.24
3	A	303	ASC	O4-C1-O1	2.55	124.16	121.24
2	B	302	HEM	CHD-C4C-NC	2.50	127.23	124.38
2	B	301	HEM	C4C-NC-C1C	2.49	108.04	105.51
3	A	303	ASC	C5-C4-C3	-2.45	110.77	114.18
3	B	303	ASC	O3-C3-C2	-2.44	125.03	132.36
3	B	303	ASC	O4-C1-O1	2.43	124.03	121.24
2	B	301	HEM	CHA-C4D-ND	2.40	127.66	124.28
2	B	301	HEM	CHC-C4B-NB	2.38	127.23	124.47
3	B	303	ASC	O1-C1-C2	-2.34	126.28	129.45
2	B	302	HEM	CBD-CAD-C3D	-2.27	109.41	114.37
3	B	303	ASC	O4-C4-C3	2.22	105.79	104.06
2	B	302	HEM	CMA-C3A-C4A	-2.22	125.05	128.46
2	B	301	HEM	CBD-CAD-C3D	-2.21	109.55	114.37
3	A	303	ASC	O3-C3-C4	2.19	124.31	118.05
2	A	302	HEM	C4A-CHB-C1B	-2.17	124.62	127.47
3	B	304	ASC	O3-C3-C4	2.17	124.25	118.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	301	HEM	C4A-C3A-C2A	2.09	108.45	107.00
3	B	303	ASC	O3-C3-C4	2.08	123.99	118.05
2	A	302	HEM	C1B-NB-C4B	2.06	107.25	105.11

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	211/230 (91%)	0.77	25 (11%) 5 5	25, 51, 78, 95	0
1	B	212/230 (92%)	0.51	20 (9%) 9 8	34, 51, 77, 96	0
All	All	423/460 (91%)	0.64	45 (10%) 6 6	25, 51, 78, 96	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	44	LYS	8.6
1	A	40	SER	8.3
1	A	142	TYR	7.9
1	A	39	LEU	7.8
1	A	12	PHE	6.7
1	A	100	TRP	6.6
1	B	220	PRO	6.0
1	B	74	GLN	5.5
1	A	42	ASP	5.2
1	A	47	ILE	4.9
1	A	13	MET	4.8
1	A	41	SER	4.3
1	A	109	LYS	3.8
1	B	174	THR	3.6
1	B	9	PHE	3.3
1	B	175	GLY	3.1
1	A	74	GLN	3.1
1	A	173	THR	3.1
1	A	219	THR	2.9
1	A	170	VAL	2.8
1	B	39	LEU	2.8
1	A	169	LEU	2.8
1	A	52	PRO	2.7
1	A	220	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	169	LEU	2.7
1	B	166	ALA	2.6
1	B	219	THR	2.6
1	B	173	THR	2.6
1	B	176	ILE	2.5
1	B	170	VAL	2.5
1	A	16	ARG	2.5
1	A	168	ALA	2.5
1	A	187	GLN	2.5
1	B	177	LEU	2.4
1	B	78	ASN	2.4
1	B	16	ARG	2.4
1	A	188	VAL	2.3
1	B	159	PHE	2.3
1	B	11	ILE	2.2
1	A	172	ALA	2.2
1	B	76	THR	2.2
1	B	143	PRO	2.2
1	B	172	ALA	2.1
1	A	166	ALA	2.1
1	A	141	TRP	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(Å ²)	Q<0.9
3	ASC	B	304	12/12	0.22	1.93	67,97,106,108	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	ASC	B	303	12/12	0.27	1.04	67,72,82,93	0
2	HEM	A	301	43/43	0.21	0.77	35,43,54,62	0
2	HEM	A	302	43/43	0.17	0.60	31,46,56,78	0
2	HEM	B	301	43/43	0.20	0.58	32,42,51,72	0
2	HEM	B	302	43/43	0.14	0.18	35,46,65,71	0
3	ASC	A	303	12/12	0.16	0.06	70,88,94,96	0
4	SO4	A	304	5/5	0.15	-0.02	79,89,99,103	0

6.5 Other polymers ⓘ

There are no such residues in this entry.