



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 10:18 AM GMT

PDB ID : 3HQ8
Title : CcpA from *G. sulfurreducens* S134P/V135K variant
Authors : Hoffmann, M.; Seidel, J.; Einsle, O.
Deposited on : 2009-06-05
Resolution : 2.40 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>

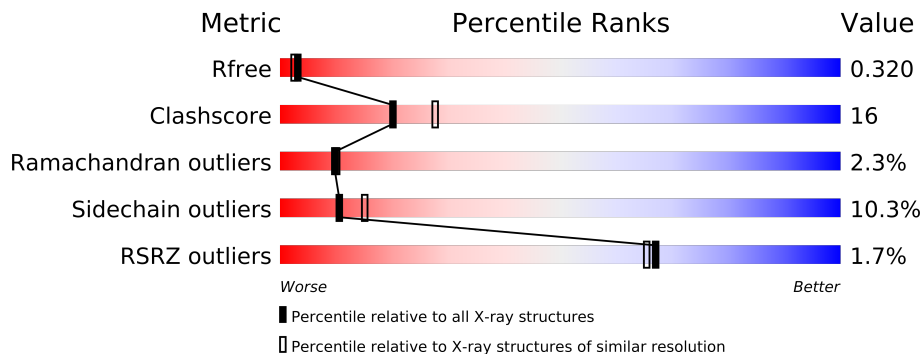
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.40 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	2207 (2.40-2.40)
Clashscore	79885	2789 (2.40-2.40)
Ramachandran outliers	78287	2736 (2.40-2.40)
Sidechain outliers	78261	2737 (2.40-2.40)
RSRZ outliers	66119	2210 (2.40-2.40)

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	345	
1	B	345	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	CA	A	402	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 5094 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.

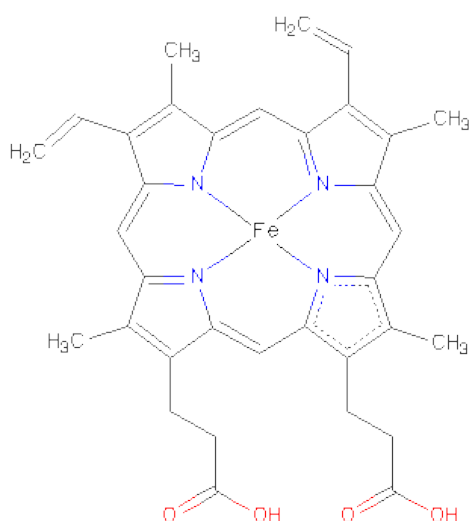
- Molecule 1 is a protein called Cytochrome c551 peroxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	315	Total	C	N	O	S	0	0	0
			2385	1525	408	443	9			
1	B	313	Total	C	N	O	S	0	1	0
			2381	1521	408	443	9			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	134	PRO	SER	ENGINEERED	UNP Q749D0
A	135	LYS	VAL	ENGINEERED	UNP Q749D0
B	134	PRO	SER	ENGINEERED	UNP Q749D0
B	135	LYS	VAL	ENGINEERED	UNP Q749D0

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

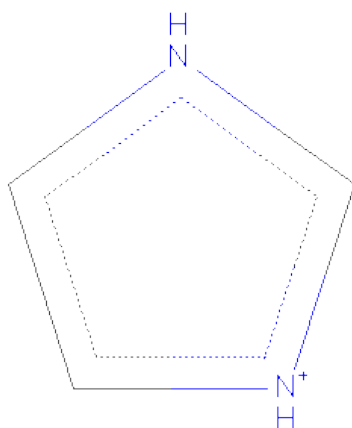


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Ca	0	0
			1	1		
3	A	1	Total	Ca	0	0
			1	1		

- Molecule 4 is IMIDAZOLE (three-letter code: IMD) (formula: C₃H₅N₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	N	0	0
			5	3	2		
4	B	1	Total	C	N	0	0
			5	3	2		

- Molecule 5 is water.

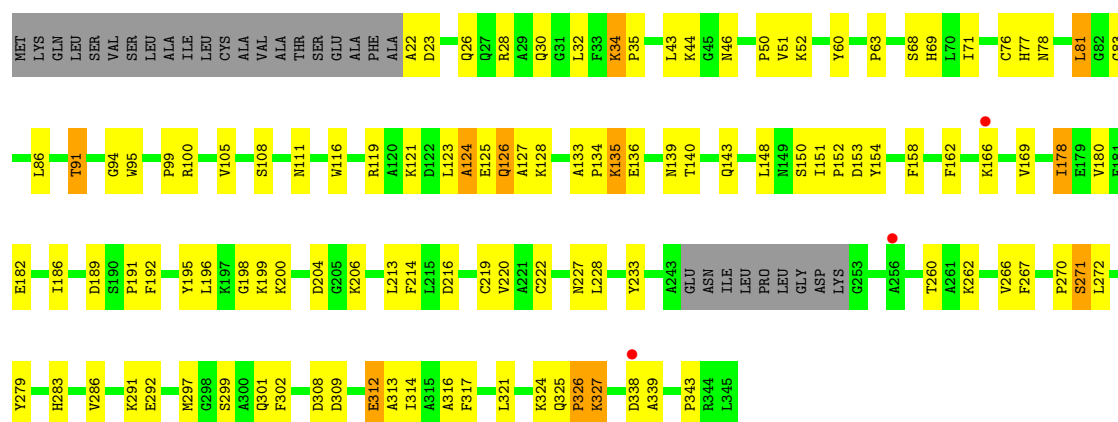
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	70	Total 70	O 70	0	0
5	B	74	Total 74	O 74	0	0

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.

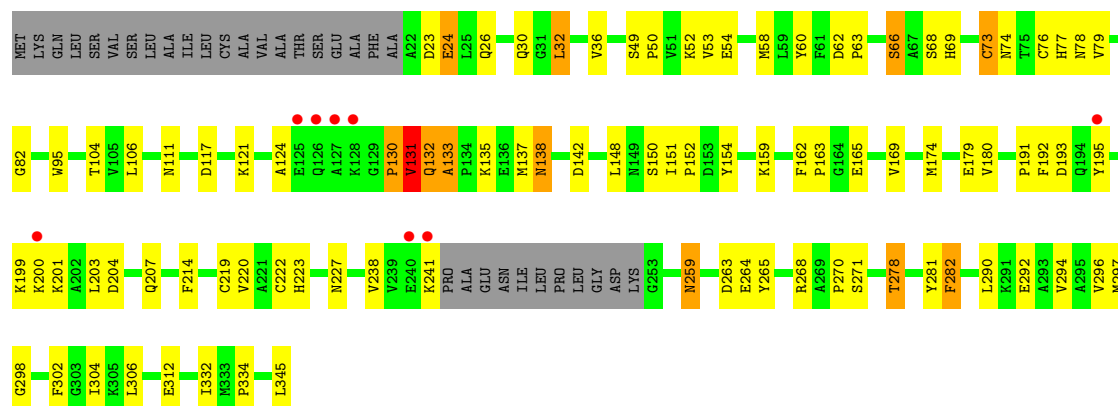
• Molecule 1: Cytochrome c551 peroxidase

Chain A: 



• Molecule 1: Cytochrome c551 peroxidase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	56.16Å 89.51Å 75.11Å 90.00° 107.57° 90.00°	Depositor
Resolution (Å)	44.77 – 2.40 44.23 – 2.40	Depositor EDS
% Data completeness (in resolution range)	98.8 (44.77-2.40) 98.7 (44.23-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.17	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.229 , 0.324 0.226 , 0.320	Depositor DCC
R_{free} test set	1370 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	38.7	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 25.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	0 of 27472 reflections	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	5094	wwPDB-VP
Average B, all atoms (Å ²)	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.13% 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, CA, IMD

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	0.64	0/2444	0.75	1/3315 (0.0%)
1	B	0.60	0/2439	0.72	0/3307
All	All	0.62	0/4883	0.74	1/6622 (0.0%)

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	1
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	81	LEU	CA-CB-CG	5.87	128.81	115.30

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	126	GLN	Peptide
1	B	130	PRO	Peptide

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	2385	0	2388	89	0
1	B	2381	0	2382	71	0
2	A	86	0	60	22	0
2	B	86	0	60	23	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	5	0	4	0	0
4	B	5	0	4	0	0
5	A	70	0	0	3	0
5	B	74	0	0	0	0
All	All	5094	0	4898	157	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 16.

All (157) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:219:CYS:SG	2:B:401:HEM:HAB	1.56	1.43
1:A:219:CYS:SG	2:A:401:HEM:HAB	1.63	1.37
1:A:76:CYS:SG	2:A:400:HEM:CAC	2.12	1.36
1:A:219:CYS:SG	2:A:401:HEM:CAB	2.22	1.28
1:A:219:CYS:HG	2:A:401:HEM:CAB	1.47	1.27
1:B:219:CYS:SG	2:B:401:HEM:CAB	2.23	1.25
1:B:76:CYS:SG	2:B:400:HEM:CBC	2.31	1.18
1:A:222:CYS:SG	2:A:401:HEM:CAC	2.32	1.16
1:A:76:CYS:SG	2:A:400:HEM:HAC	1.89	1.08
1:B:73:CYS:SG	2:B:400:HEM:CAB	2.42	1.06
1:B:76:CYS:SG	2:B:400:HEM:CAC	2.47	1.02
1:B:222:CYS:SG	2:B:401:HEM:CAC	2.47	1.01
1:B:222:CYS:HG	2:B:401:HEM:CAC	1.73	1.00
1:B:219:CYS:HG	2:B:401:HEM:CAB	1.70	1.00
1:A:134:PRO:O	1:A:135:LYS:HB2	1.58	0.99
1:A:222:CYS:SG	2:A:401:HEM:HAC	2.12	0.89
1:B:131:VAL:HA	1:B:132:GLN:HB2	1.52	0.89

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:127:ALA:HB1	1:A:128:LYS:HA	1.56	0.86
1:B:73:CYS:SG	2:B:400:HEM:HAB	2.16	0.86
1:B:297:MET:HG2	2:B:401:HEM:C4B	2.17	0.79
1:A:76:CYS:SG	2:A:400:HEM:CBC	2.70	0.79
1:A:150:SER:HB2	1:A:339:ALA:O	1.85	0.76
1:B:298:GLY:HA3	1:B:306:LEU:HD12	1.68	0.74
1:A:76:CYS:HG	2:A:400:HEM:HAC	1.51	0.73
1:A:108:SER:HA	1:A:111:ASN:ND2	2.05	0.72
1:A:206:LYS:NZ	1:A:312:GLU:HG2	2.05	0.70
1:A:140:THR:H	1:A:143:GLN:NE2	1.89	0.69
1:B:117:ASP:HB2	1:B:265:TYR:HE2	1.59	0.66
1:A:297:MET:HE3	2:A:401:HEM:C4D	2.30	0.66
1:A:95:TRP:HB2	1:B:95:TRP:HB2	1.78	0.66
1:B:60:TYR:CE1	1:B:73:CYS:HB2	2.31	0.65
1:A:108:SER:HA	1:A:111:ASN:HD21	1.62	0.64
1:A:127:ALA:CB	1:A:128:LYS:HA	2.26	0.64
1:A:68:SER:O	1:A:69:HIS:HB2	1.98	0.64
1:B:222:CYS:SG	2:B:401:HEM:C3C	2.89	0.63
1:A:136:GLU:HB2	2:A:400:HEM:HBC2	1.80	0.63
1:A:313:ALA:O	1:A:316:ALA:HB3	1.99	0.62
1:A:63:PRO:O	1:A:69:HIS:HA	1.99	0.62
1:B:292:GLU:O	1:B:296:VAL:HG23	1.99	0.62
1:B:151:ILE:HG21	1:B:334:PRO:HG2	1.83	0.61
1:B:214:PHE:O	1:B:219:CYS:HB2	2.01	0.61
1:A:52:LYS:HE2	1:A:162:PHE:CE1	2.35	0.61
1:A:22:ALA:HB1	5:A:347:HOH:O	2.01	0.61
1:A:297:MET:HE3	2:A:401:HEM:CHA	2.31	0.60
1:A:76:CYS:SG	2:A:400:HEM:C3C	2.92	0.59
1:A:297:MET:CE	2:A:401:HEM:C4D	2.85	0.59
1:A:228:LEU:HD22	1:A:321:LEU:HD11	1.85	0.58
1:A:178:ILE:O	1:A:182:GLU:HG3	2.03	0.58
1:B:58:MET:O	1:B:62:ASP:HB2	2.02	0.58
1:A:105:VAL:O	1:A:105:VAL:HG22	2.03	0.58
1:A:206:LYS:HZ1	1:A:312:GLU:HG2	1.66	0.58
1:B:282:PHE:CD1	2:B:401:HEM:HBD1	2.39	0.57
1:A:52:LYS:HE2	1:A:162:PHE:CZ	2.40	0.57
1:B:53:VAL:HG22	1:B:180:VAL:HG12	1.86	0.57
1:B:290:LEU:HG	1:B:290:LEU:O	2.03	0.57
1:A:222:CYS:SG	2:A:401:HEM:CBC	2.91	0.57
1:A:60:TYR:OH	1:A:105:VAL:HG11	2.05	0.57
1:A:299:SER:HA	5:A:390:HOH:O	2.05	0.57
1:B:73:CYS:HB3	2:B:400:HEM:C3B	2.41	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:228:LEU:HD22	1:A:321:LEU:CD1	2.37	0.55
1:B:204:ASP:OD1	1:B:207:GLN:HG3	2.07	0.55
1:B:297:MET:HG2	2:B:401:HEM:C3B	2.42	0.54
1:A:69:HIS:CD2	1:A:343:PRO:HG3	2.43	0.54
1:A:28:ARG:O	1:A:32:LEU:HD12	2.07	0.54
1:A:297:MET:O	1:A:299:SER:N	2.41	0.53
1:A:126:GLN:HB3	1:A:127:ALA:CA	2.39	0.53
1:A:192:PHE:CD1	1:A:317:PHE:HE1	2.26	0.53
1:A:169:VAL:HG12	1:A:169:VAL:O	2.08	0.53
1:A:28:ARG:NH2	1:A:216:ASP:OD1	2.33	0.52
1:B:203:LEU:HB3	1:B:207:GLN:HB2	1.91	0.52
1:A:23:ASP:HB2	1:A:200:LYS:NZ	2.25	0.52
1:A:127:ALA:CB	1:A:128:LYS:CA	2.88	0.52
1:B:117:ASP:HB2	1:B:265:TYR:CE2	2.44	0.52
1:A:91:THR:HG21	1:B:345:LEU:HD22	1.91	0.52
1:B:238:VAL:HG21	1:B:282:PHE:CE1	2.45	0.52
1:A:139:ASN:OD1	1:A:143:GLN:NE2	2.43	0.52
1:A:134:PRO:O	1:A:135:LYS:CB	2.42	0.51
1:B:137:MET:HE3	2:B:400:HEM:HHC	1.92	0.51
1:A:297:MET:HG3	2:A:401:HEM:C4A	2.45	0.51
1:B:26:GLN:HE21	1:B:30:GLN:NE2	2.09	0.51
1:A:195:TYR:HA	1:A:199:LYS:O	2.10	0.51
1:B:26:GLN:HE21	1:B:30:GLN:HE22	1.59	0.50
1:B:282:PHE:HD1	2:B:401:HEM:HBD1	1.76	0.50
1:B:222:CYS:SG	2:B:401:HEM:HAC	2.46	0.50
1:A:126:GLN:HB3	1:A:127:ALA:HA	1.94	0.50
1:B:302:PHE:HB2	1:B:304:ILE:HG13	1.94	0.50
1:B:223:HIS:HE1	1:B:270:PRO:HD2	1.77	0.50
1:A:214:PHE:O	1:A:219:CYS:HB2	2.11	0.50
1:A:99:PRO:O	1:A:100:ARG:HD3	2.12	0.50
1:B:214:PHE:CE1	1:B:219:CYS:HB3	2.47	0.49
1:B:73:CYS:SG	2:B:400:HEM:C3B	3.03	0.49
1:A:46:ASN:ND2	1:A:180:VAL:HG23	2.28	0.49
1:A:325:GLN:HB2	1:B:332:ILE:HD13	1.94	0.49
1:B:82:GLY:HA2	1:B:106:LEU:HD11	1.95	0.49
1:A:222:CYS:SG	2:A:401:HEM:C3C	3.03	0.49
1:A:23:ASP:HB2	1:A:200:LYS:HZ3	1.78	0.48
1:A:204:ASP:C	1:A:204:ASP:OD1	2.51	0.48
1:B:270:PRO:HG2	2:B:401:HEM:HBA1	1.95	0.48
1:B:77:HIS:CD2	2:B:400:HEM:NB	2.81	0.48
1:B:150:SER:O	1:B:152:PRO:HD3	2.13	0.47
1:A:78:ASN:HB3	1:A:81:LEU:HB2	1.95	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:119:ARG:HE	1:A:121:LYS:HE3	1.79	0.47
1:B:60:TYR:CD1	1:B:73:CYS:HB2	2.50	0.47
1:A:270:PRO:HD3	2:A:401:HEM:C3D	2.50	0.47
1:B:76:CYS:SG	2:B:400:HEM:C3C	3.07	0.46
1:B:204:ASP:OD1	1:B:204:ASP:C	2.53	0.46
1:B:52:LYS:HE2	1:B:162:PHE:CE1	2.51	0.46
1:A:26:GLN:O	1:A:30:GLN:HG3	2.15	0.46
1:B:259[B]:ASN:ND2	1:B:259[B]:ASN:C	2.68	0.46
1:B:297:MET:HG2	2:B:401:HEM:NB	2.30	0.46
1:B:32:LEU:HG	1:B:220:VAL:HG21	1.98	0.46
1:A:116:TRP:CE3	1:A:283:HIS:HB3	2.51	0.46
1:A:297:MET:HG3	2:A:401:HEM:NA	2.32	0.45
1:B:74:ASN:OD1	1:B:78:ASN:HA	2.15	0.45
1:B:24:GLU:HG2	1:B:24:GLU:O	2.17	0.45
1:B:54:GLU:O	1:B:58:MET:HG3	2.17	0.45
1:B:192:PHE:O	1:B:195:TYR:HB3	2.17	0.45
1:A:123:LEU:HA	1:A:126:GLN:NE2	2.32	0.44
1:B:131:VAL:CA	1:B:132:GLN:HB2	2.36	0.44
1:A:260:THR:HG21	1:A:267:PHE:HE2	1.83	0.44
1:B:104:THR:HB	1:B:278:THR:HG21	2.00	0.44
1:B:238:VAL:HG21	1:B:282:PHE:CZ	2.53	0.44
1:A:227:ASN:HB3	1:A:271:SER:HB2	1.99	0.44
1:A:34:LYS:HD3	1:A:35:PRO:HD2	1.99	0.44
1:A:123:LEU:O	1:A:124:ALA:C	2.56	0.43
1:A:151:ILE:HA	1:A:152:PRO:HD2	1.81	0.43
1:A:327:LYS:HA	1:A:327:LYS:HD2	1.94	0.43
1:A:94:GLY:O	1:A:95:TRP:CB	2.67	0.42
1:B:259[B]:ASN:C	1:B:259[B]:ASN:HD22	2.22	0.42
1:B:63:PRO:O	1:B:66:SER:OG	2.37	0.42
1:A:63:PRO:HA	1:A:71:ILE:O	2.19	0.42
1:A:272:LEU:HA	1:A:272:LEU:HD23	1.66	0.42
1:B:132:GLN:HG3	1:B:133:ALA:N	2.34	0.42
1:A:192:PHE:CE1	1:A:317:PHE:HE1	2.36	0.42
1:B:111:ASN:OD1	1:B:227:ASN:ND2	2.34	0.42
1:A:213:LEU:HB3	1:A:314:ILE:HD11	2.02	0.42
1:B:138:ASN:O	1:B:138:ASN:ND2	2.43	0.42
1:A:272:LEU:HD11	2:A:401:HEM:CHB	2.50	0.42
1:A:220:VAL:HG22	1:A:220:VAL:O	2.20	0.42
1:B:49:SER:O	1:B:50:PRO:C	2.58	0.41
1:A:68:SER:O	1:A:343:PRO:HB3	2.20	0.41
1:A:233:TYR:HB3	1:A:266:VAL:HB	2.02	0.41
1:B:159:LYS:O	1:B:163:PRO:HA	2.20	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:270:PRO:HG2	2:A:401:HEM:CBA	2.51	0.41
1:B:76:CYS:SG	2:B:400:HEM:HBC2	2.47	0.41
1:B:281:TYR:O	1:B:282:PHE:HB2	2.21	0.41
1:A:297:MET:HB3	2:A:401:HEM:C4B	2.56	0.41
1:A:133:ALA:HB1	1:A:136:GLU:HG2	2.03	0.41
1:A:192:PHE:CE2	1:A:196:LEU:HD11	2.56	0.41
1:A:148:LEU:HD22	1:A:158:PHE:HE2	1.86	0.41
1:B:36:VAL:HG23	1:B:193:ASP:OD2	2.21	0.41
1:A:77:HIS:HA	1:A:83:GLY:O	2.20	0.41
1:A:325:GLN:HA	1:A:326:PRO:HD3	1.94	0.41
1:A:200:LYS:HE3	5:A:348:HOH:O	2.21	0.40
1:B:68:SER:O	1:B:69:HIS:HB2	2.21	0.40
1:B:290:LEU:O	1:B:294:VAL:HG23	2.21	0.40
1:B:263:ASP:O	1:B:264:GLU:HB2	2.22	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	311/345 (90%)	271 (87%)	34 (11%)	6 (2%)	12	14
1	B	310/345 (90%)	277 (89%)	25 (8%)	8 (3%)	8	8
All	All	621/690 (90%)	548 (88%)	59 (10%)	14 (2%)	10	10

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	124	ALA
1	A	135	LYS
1	B	130	PRO
1	B	131	VAL
1	B	132	GLN
1	A	198	GLY

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Mol	Chain	Res	Type
1	B	124	ALA
1	B	169	VAL
1	B	282	PHE
1	B	191	PRO
1	A	191	PRO
1	A	50	PRO
1	A	279	TYR
1	B	133	ALA

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	252/276 (91%)	225 (89%)	27 (11%)	10	13
1	B	252/276 (91%)	226 (90%)	26 (10%)	10	15
All	All	504/552 (91%)	451 (90%)	53 (10%)	10	14

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

Mol	Chain	Res	Type
1	A	34	LYS
1	A	43	LEU
1	A	44	LYS
1	A	51	VAL
1	A	86	LEU
1	A	91	THR
1	A	125	GLU
1	A	153	ASP
1	A	154	TYR
1	A	166	LYS
1	A	178	ILE
1	A	186	ILE
1	A	189	ASP
1	A	262	LYS
1	A	271	SER
1	A	286	VAL
1	A	291	LYS

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Mol	Chain	Res	Type
1	A	292	GLU
1	A	301	GLN
1	A	302	PHE
1	A	308	ASP
1	A	309	ASP
1	A	312	GLU
1	A	324	LYS
1	A	326	PRO
1	A	327	LYS
1	A	338	ASP
1	B	23	ASP
1	B	24	GLU
1	B	32	LEU
1	B	66	SER
1	B	73	CYS
1	B	79	VAL
1	B	121	LYS
1	B	131	VAL
1	B	135	LYS
1	B	138	ASN
1	B	142	ASP
1	B	148	LEU
1	B	154	TYR
1	B	165	GLU
1	B	174	MET
1	B	179	GLU
1	B	199	LYS
1	B	200	LYS
1	B	201	LYS
1	B	241	LYS
1	B	259[A]	ASN
1	B	259[B]	ASN
1	B	268	ARG
1	B	271	SER
1	B	278	THR
1	B	312	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	69	HIS
1	A	107	ASN

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Mol	Chain	Res	Type
1	A	139	ASN
1	A	143	GLN
1	A	301	GLN
1	B	26	GLN
1	B	107	ASN
1	B	173	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 ⓘ

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 for Mogul analysis.

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	IMD	A	346	2	5,5,5	1.30	1 (20%)	5,5,5	0.60	0
2	HEM	A	400	1,4	49,50,50	2.25	15 (30%)	46,82,82	2.38	13 (28%)
2	HEM	A	401	1	49,50,50	2.76	15 (30%)	46,82,82	2.30	12 (26%)
4	IMD	B	346	2	5,5,5	1.32	1 (20%)	5,5,5	0.20	0
2	HEM	B	400	1,4	49,50,50	2.34	13 (26%)	46,82,82	1.95	11 (23%)
2	HEM	B	401	1	49,50,50	2.37	14 (28%)	46,82,82	1.97	11 (23%)

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	IMD	A	346	2	-	0/0/0/0	0/1/1/1
2	HEM	A	400	1,4	-	0/14/114/114	0/0/8/8
2	HEM	A	401	1	-	0/14/114/114	0/0/8/8
4	IMD	B	346	2	-	0/0/0/0	0/1/1/1
2	HEM	B	400	1,4	-	0/14/114/114	0/0/8/8
2	HEM	B	401	1	-	0/14/114/114	0/0/8/8

All (59) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HEM	C3D-C4D	9.25	1.46	1.44
2	A	401	HEM	C3D-C2D	6.28	1.54	1.43
2	B	400	HEM	C3D-C2D	5.86	1.54	1.43
2	B	401	HEM	C3B-C2B	-5.70	1.33	1.43
2	B	401	HEM	C3D-C4D	5.66	1.46	1.44
2	B	400	HEM	C3B-C2B	-5.49	1.34	1.43
2	A	400	HEM	C3D-C2D	5.48	1.53	1.43
2	A	401	HEM	C3C-C2C	-5.37	1.34	1.43
2	B	400	HEM	C3B-CAB	5.34	1.57	1.40
2	A	400	HEM	C3C-C2C	-5.34	1.34	1.43
2	B	400	HEM	C2B-C1B	5.34	1.45	1.44
2	A	401	HEM	C3B-C2B	-5.30	1.34	1.43
2	A	401	HEM	C2D-C1D	-5.30	1.43	1.44
2	B	401	HEM	C3D-C2D	5.16	1.52	1.43
2	B	400	HEM	C3C-CAC	5.11	1.56	1.40
2	A	400	HEM	C3B-C2B	-5.09	1.34	1.43
2	A	401	HEM	FE-NA	5.08	2.14	1.92
2	B	401	HEM	C3C-CAC	5.07	1.56	1.40
2	A	400	HEM	C3C-CAC	5.06	1.56	1.40
2	B	401	HEM	C3C-C2C	-5.00	1.35	1.43
2	A	401	HEM	C3C-CAC	4.95	1.56	1.40
2	A	400	HEM	C3B-CAB	4.92	1.55	1.40
2	A	401	HEM	C3B-CAB	4.77	1.55	1.40
2	B	400	HEM	C4A-C3A	4.63	1.46	1.40
2	B	401	HEM	FE-ND	4.61	2.14	1.97
2	B	401	HEM	C4A-C3A	4.43	1.45	1.40
2	B	401	HEM	C3B-CAB	4.41	1.54	1.40
2	B	400	HEM	C3C-C2C	-4.32	1.36	1.43
2	A	400	HEM	C4A-C3A	4.17	1.45	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	HEM	C2D-C1D	4.04	1.45	1.44
2	A	400	HEM	FE-ND	3.62	2.11	1.97
2	B	400	HEM	FE-ND	3.45	2.10	1.97
2	A	400	HEM	CMB-C2B	3.24	1.57	1.47
2	B	400	HEM	FE-NA	3.19	2.06	1.92
2	A	401	HEM	C4A-C3A	3.11	1.44	1.40
2	A	401	HEM	CHA-C4D	2.80	1.39	1.35
2	A	400	HEM	CHA-C4D	2.67	1.39	1.35
2	B	400	HEM	CMB-C2B	2.63	1.55	1.47
2	A	401	HEM	CMB-C2B	2.60	1.55	1.47
2	A	401	HEM	CMC-C2C	2.59	1.55	1.47
2	A	401	HEM	CMD-C2D	2.59	1.55	1.47
2	B	401	HEM	CMD-C2D	2.57	1.55	1.47
2	B	400	HEM	CMD-C2D	2.45	1.55	1.47
2	B	400	HEM	CMC-C2C	2.43	1.55	1.47
2	A	400	HEM	C3D-C4D	2.43	1.45	1.44
2	A	400	HEM	C2D-C1D	-2.40	1.44	1.44
2	B	401	HEM	CMB-C2B	2.40	1.54	1.47
2	A	400	HEM	C4A-NA	2.39	1.41	1.36
2	B	401	HEM	CMC-C2C	2.36	1.54	1.47
2	A	400	HEM	CMA-C3A	2.30	1.56	1.51
4	B	346	IMD	C2-N3	2.29	1.36	1.31
2	B	400	HEM	FE-NC	-2.24	1.89	1.97
2	A	401	HEM	C3B-C4B	2.23	1.47	1.44
2	A	401	HEM	CMA-C3A	2.18	1.56	1.51
2	A	400	HEM	CMD-C2D	2.16	1.54	1.47
2	B	401	HEM	FE-NB	2.15	2.05	1.97
4	A	346	IMD	C2-N3	2.15	1.36	1.31
2	A	400	HEM	CMC-C2C	2.11	1.53	1.47
2	B	401	HEM	CAA-C2A	2.11	1.55	1.52

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	HEM	C3B-C4B-NB	-8.22	108.12	114.00
2	A	400	HEM	C3B-C4B-NB	-8.10	108.20	114.00
2	A	401	HEM	C3B-C4B-NB	-8.02	108.26	114.00
2	B	400	HEM	C3B-C4B-NB	-6.93	109.04	114.00
2	A	401	HEM	CBD-CAD-C3D	-6.88	99.36	114.37
2	A	400	HEM	C4D-ND-C1D	6.52	111.83	105.16
2	A	401	HEM	C4D-ND-C1D	6.08	111.39	105.16
2	B	400	HEM	C4D-ND-C1D	5.50	110.79	105.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	400	HEM	C1A-CHA-C4D	-4.88	121.05	127.47
2	A	400	HEM	CBD-CAD-C3D	-4.38	104.81	114.37
2	B	401	HEM	C4D-ND-C1D	4.36	109.62	105.16
2	B	401	HEM	CBA-CAA-C2A	-4.05	105.55	112.69
2	A	400	HEM	CHD-C4C-NC	3.91	128.13	124.73
2	A	401	HEM	C4A-CHB-C1B	-3.68	122.62	127.47
2	B	400	HEM	C1A-CHA-C4D	-3.49	122.88	127.47
2	A	400	HEM	CAA-CBA-CGA	-3.49	102.25	113.47
2	A	401	HEM	CHD-C1D-ND	3.37	127.38	124.58
2	B	401	HEM	C1B-NB-C4B	3.19	108.42	105.16
2	A	400	HEM	C4B-CHC-C1C	-3.16	118.25	126.57
2	B	401	HEM	CAD-C3D-C4D	3.02	129.95	124.53
2	B	400	HEM	C2D-C1D-ND	-2.97	109.43	112.93
2	B	400	HEM	C1B-NB-C4B	2.77	108.00	105.16
2	A	400	HEM	C2D-C1D-ND	-2.74	109.69	112.93
2	B	401	HEM	C1A-CHA-C4D	-2.74	123.87	127.47
2	A	400	HEM	C1B-NB-C4B	2.72	107.95	105.16
2	A	401	HEM	CBA-CAA-C2A	-2.66	108.00	112.69
2	B	400	HEM	CHD-C4C-NC	2.58	126.97	124.73
2	A	400	HEM	CMA-C3A-C4A	-2.53	124.73	128.62
2	B	400	HEM	CMA-C3A-C4A	-2.53	124.74	128.62
2	B	401	HEM	C4A-C3A-C2A	2.42	108.68	107.00
2	A	400	HEM	C1A-C2A-C3A	2.40	109.41	106.92
2	A	400	HEM	CBA-CAA-C2A	-2.39	108.49	112.69
2	A	401	HEM	C2D-C1D-ND	-2.35	110.16	112.93
2	A	401	HEM	CMA-C3A-C4A	-2.34	125.02	128.62
2	B	400	HEM	C4C-NC-C1C	2.29	107.92	105.53
2	A	400	HEM	CAD-C3D-C4D	2.26	128.59	124.53
2	A	401	HEM	C1A-C2A-C3A	2.26	109.26	106.92
2	B	401	HEM	C3A-C4A-NA	-2.21	107.75	109.41
2	B	400	HEM	CAD-C3D-C4D	2.19	128.46	124.53
2	A	401	HEM	C2A-C1A-NA	-2.14	106.77	109.73
2	B	400	HEM	CBD-CAD-C3D	-2.12	109.75	114.37
2	A	401	HEM	CAA-C2A-C3A	-2.11	122.98	129.00
2	B	401	HEM	O1A-CGA-CBA	-2.07	115.90	123.03
2	B	401	HEM	C2D-C1D-ND	-2.06	110.50	112.93
2	B	400	HEM	C4A-CHB-C1B	-2.06	124.77	127.47
2	B	401	HEM	CMA-C3A-C4A	-2.05	125.46	128.62
2	A	401	HEM	CMD-C2D-C3D	2.01	130.15	125.60

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	315/345 (91%)	-0.01	3 (0%) 79 79	19, 40, 59, 74	0
1	B	313/345 (90%)	0.14	8 (2%) 53 51	22, 40, 72, 83	0
All	All	628/690 (91%)	0.06	11 (1%) 67 63	19, 40, 66, 83	0

All (11) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	127	ALA	5.5
1	B	128	LYS	4.4
1	B	195	TYR	4.2
1	B	241	LYS	3.6
1	B	126	GLN	2.6
1	A	256	ALA	2.4
1	B	125	GLU	2.3
1	B	240	GLU	2.2
1	A	338	ASP	2.1
1	A	166	LYS	2.1
1	B	200	LYS	2.1

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	CA	A	402	1/1	0.18	2.16	31,31,31,31	0
3	CA	B	402	1/1	0.22	1.86	37,37,37,37	0
2	HEM	B	400	43/43	0.18	0.59	24,34,43,48	0
2	HEM	B	401	43/43	0.16	0.31	27,38,40,40	0
2	HEM	A	400	43/43	0.16	0.12	15,25,31,33	0
4	IMD	A	346	5/5	0.15	0.05	26,26,27,27	0
2	HEM	A	401	43/43	0.14	-0.43	20,31,33,33	0
4	IMD	B	346	5/5	0.12	-1.38	40,40,42,42	0

6.5 Other polymers ⓘ

There are no such residues in this entry.