



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 03:33 AM GMT

PDB ID : 3S3B
Title : Structure of Thermus thermophilus cytochrome ba3 oxidase 240s after Xe de-pressurization
Authors : Luna, V.M.; Fee, J.A.; Deniz, A.A.; Stout, C.D.
Deposited on : 2011-05-18
Resolution : 3.30 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>

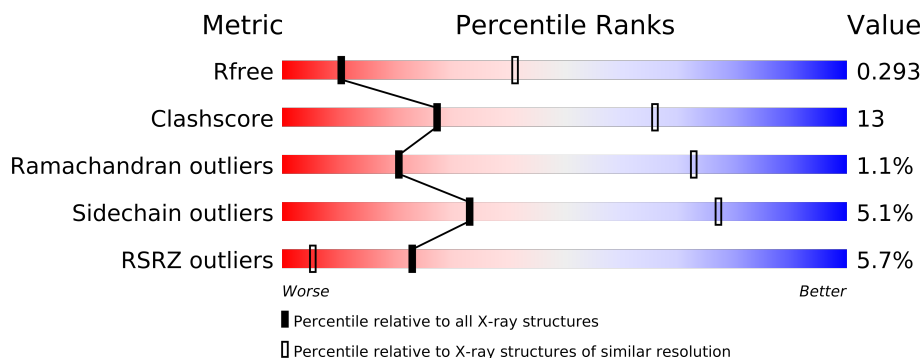
The following versions of software and data (see [references](#)) were used in the production of this report:

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 3.30 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	1341 (3.40-3.20)
Clashscore	79885	1696 (3.40-3.20)
Ramachandran outliers	78287	1664 (3.40-3.20)
Sidechain outliers	78261	1662 (3.40-3.20)
RSRZ outliers	66119	1342 (3.40-3.20)

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	568	
2	B	166	
3	C	33	

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

Mol	Type	Chain	Res	Geometry	Electron density
8	XE	A	563	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 6063 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 c oxidase subunit 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	555	Total	C	N	O	S	0	0	0
			4392	2977	705	694	16			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-5	MET	-	EXPRESSION TAG	UNP Q5SJ79
A	-4	HIS	-	EXPRESSION TAG	UNP Q5SJ79
A	-3	HIS	-	EXPRESSION TAG	UNP Q5SJ79
A	-2	HIS	-	EXPRESSION TAG	UNP Q5SJ79
A	-1	HIS	-	EXPRESSION TAG	UNP Q5SJ79
A	0	HIS	-	EXPRESSION TAG	UNP Q5SJ79
A	1	HIS	-	EXPRESSION TAG	UNP Q5SJ79

- Molecule 2 is a protein called Cytochrome c oxidase subunit 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	166	Total	C	N	O	S	0	0	0
			1298	844	216	234	4			

- Molecule 3 is a protein called Cytochrome c oxidase polypeptide 2A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	33	Total	C	N	O	0	0	0
			259	179	39	41			

- Molecule 4 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

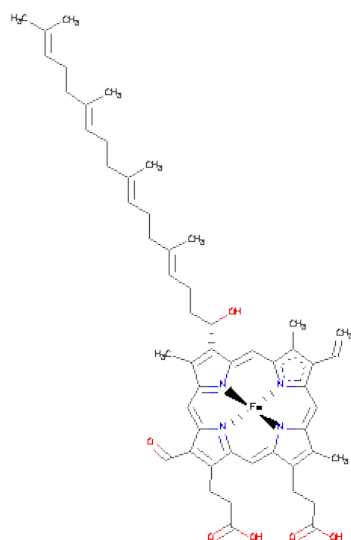
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cu	0	0
			1	1		

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



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

- Molecule 6 is HEME-AS (three-letter code: HAS) (formula: $C_{54}H_{64}FeN_4O_6$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
6	A	1	65	54	1	4	6	0	0

- Molecule 7 is DINUCLEAR COPPER ION (three-letter code: CUA) (formula: Cu₂).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	B	1	Total	Cu	0	0
			2	2		

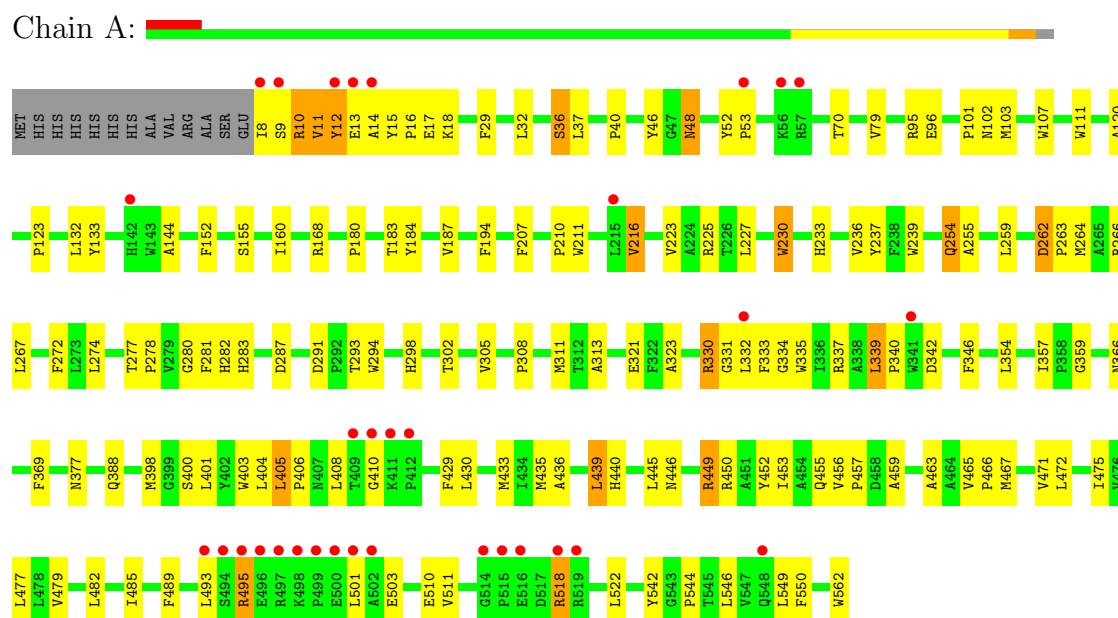
- Molecule 8 is XENON (three-letter code: XE) (formula: Xe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	3	Total	Xe	0	0
			3	3		

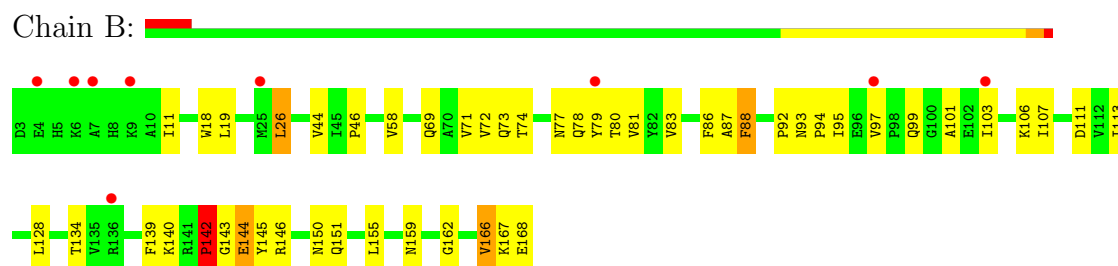
3 Residue-property plots

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 c oxidase subunit 1



• Molecule 2: Cytochrome c oxidase subunit 2



• Molecule 3: Cytochrome c oxidase polypeptide 2A



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	109.92Å 109.92Å 170.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	92.32 – 3.30 50.39 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.8 (92.32-3.30) 99.8 (50.39-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.15	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.55 (at 3.33Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.257 , 0.299 0.257 , 0.293	Depositor DCC
R_{free} test set	822 reflections (5.31%)	DCC
Wilson B-factor (Å ²)	86.9	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	1 of 16293 reflections (0.006%)	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	6063	wwPDB-VP
Average B, all atoms (Å ²)	86.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.30% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, HAS, CUA, CU, XE

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	0.70	2/4549 (0.0%)	0.71	2/6243 (0.0%)
2	B	0.81	2/1335 (0.1%)	0.74	0/1822
3	C	0.74	0/265	0.66	0/359
All	All	0.73	4/6149 (0.1%)	0.71	2/8424 (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
2	B	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	144	GLU	CB-CG	6.75	1.65	1.52
2	B	144	GLU	N-CA	6.11	1.58	1.46
1	A	510	GLU	CG-CD	-5.78	1.43	1.51
1	A	12	TYR	CE2-CZ	5.23	1.45	1.38

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	10	ARG	NE-CZ-NH1	-8.43	116.08	120.30
1	A	510	GLU	OE1-CD-OE2	5.53	129.93	123.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	142	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	4392	0	4505	119	28
2	B	1298	0	1280	33	28
3	C	259	0	279	10	0
4	A	1	0	0	0	0
5	A	43	0	30	5	0
6	A	65	0	62	9	0
7	B	2	0	0	0	0
8	A	3	0	0	0	0
All	All	6063	0	6156	157	28

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

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:233:HIS:NE2	1:A:237:TYR:HE2	0.91	1.38
1:A:233:HIS:NE2	1:A:237:TYR:CE2	1.79	1.37
1:A:233:HIS:CD2	1:A:237:TYR:HE2	1.80	0.98
1:A:262:ASP:HB2	1:A:511:VAL:HG11	1.47	0.93
1:A:15:TYR:HB2	1:A:18:LYS:HD2	1.58	0.86
1:A:313:ALA:HB2	6:A:801:HAS:H273	1.58	0.86
3:C:2:GLU:HG3	3:C:3:GLU:H	1.44	0.81
1:A:233:HIS:CD2	1:A:237:TYR:CE2	2.63	0.80
5:A:800:HEM:HBC2	5:A:800:HEM:HMC2	1.62	0.80
1:A:335:TRP:O	1:A:339:LEU:HD22	1.83	0.79
2:B:97:VAL:HG23	2:B:166:VAL:HG12	1.65	0.78
1:A:233:HIS:CE1	1:A:237:TYR:CE2	2.73	0.77
1:A:9:SER:HB3	1:A:12:TYR:CD2	2.22	0.75
1:A:388:GLN:CB	6:A:801:HAS:HAC	2.19	0.72
1:A:357:ILE:HG23	3:C:15:LEU:HD12	1.74	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:B:145:TYR:HE1	2:B:166:VAL:HG21	1.56	0.70
1:A:388:GLN:HB3	6:A:801:HAS:CAC	2.22	0.69
1:A:29:PHE:CE1	1:A:401:LEU:HD11	2.28	0.69
1:A:435:MET:HG2	1:A:439:LEU:HD22	1.77	0.67
1:A:465:VAL:HG23	1:A:466:PRO:HD3	1.75	0.67
1:A:410:GLY:HA2	1:A:501:LEU:HD12	1.77	0.66
2:B:145:TYR:CE1	2:B:166:VAL:HG21	2.31	0.66
1:A:388:GLN:HB3	6:A:801:HAS:HAC	1.79	0.64
1:A:401:LEU:HG	1:A:405:LEU:HD22	1.79	0.63
5:A:800:HEM:HBC2	5:A:800:HEM:CMC	2.28	0.63
1:A:449:ARG:HH12	6:A:801:HAS:CGA	2.12	0.63
1:A:518:ARG:HG2	1:A:518:ARG:HH21	1.66	0.60
1:A:388:GLN:HB2	6:A:801:HAS:HAC	1.83	0.58
1:A:377:ASN:HB3	2:B:150:ASN:O	2.04	0.58
1:A:410:GLY:C	1:A:501:LEU:HD11	2.23	0.58
1:A:482:LEU:HA	1:A:485:ILE:HD12	1.86	0.58
2:B:69:GLN:HA	2:B:69:GLN:NE2	2.18	0.57
1:A:357:ILE:HG23	3:C:15:LEU:CD1	2.34	0.57
1:A:223:VAL:HG12	1:A:549:LEU:HB3	1.85	0.57
3:C:2:GLU:CG	3:C:3:GLU:H	2.16	0.57
2:B:97:VAL:HG23	2:B:166:VAL:CG1	2.34	0.57
1:A:467:MET:O	1:A:471:VAL:HG23	2.05	0.57
1:A:332:LEU:HD12	3:C:6:LYS:HB3	1.86	0.57
1:A:489:PHE:HB3	1:A:493:LEU:HD22	1.87	0.57
1:A:332:LEU:HD23	1:A:333:PHE:CE2	2.40	0.56
2:B:73:GLN:HB2	2:B:79:TYR:CE1	2.40	0.56
2:B:71:VAL:HG13	2:B:81:VAL:HG22	1.87	0.56
3:C:2:GLU:HG3	3:C:3:GLU:N	2.17	0.56
2:B:143:GLY:N	2:B:166:VAL:HG23	2.21	0.56
1:A:302:THR:O	1:A:305:VAL:HG12	2.06	0.55
1:A:410:GLY:HA2	1:A:501:LEU:CD1	2.37	0.55
1:A:436:ALA:O	1:A:440:HIS:ND1	2.32	0.55
1:A:463:ALA:O	1:A:467:MET:HG3	2.06	0.54
1:A:11:VAL:HG22	1:A:503:GLU:CG	2.36	0.54
1:A:48:ASN:HB2	1:A:467:MET:SD	2.47	0.54
1:A:32:LEU:O	1:A:36:SER:HB2	2.08	0.53
1:A:183:THR:O	1:A:187:VAL:HG23	2.09	0.53
1:A:207:PHE:O	1:A:211:TRP:HB2	2.09	0.53
1:A:398:MET:O	1:A:401:LEU:HB2	2.09	0.52
1:A:29:PHE:CZ	1:A:401:LEU:HD11	2.44	0.52
2:B:18:TRP:CE3	3:C:12:ILE:HD12	2.45	0.52
1:A:282:HIS:CD2	1:A:283:HIS:CD2	2.97	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:429:PHE:O	1:A:433:MET:HG2	2.09	0.51
1:A:225:ARG:HD3	1:A:287:ASP:OD1	2.11	0.51
1:A:37:LEU:O	1:A:40:PRO:HD2	2.11	0.50
2:B:74:THR:HG21	2:B:80:THR:OG1	2.11	0.50
1:A:294:TRP:CZ2	1:A:544:PRO:HG2	2.47	0.50
2:B:139:PHE:CG	2:B:166:VAL:HG11	2.46	0.50
1:A:330:ARG:HG3	1:A:334:GLY:CA	2.42	0.50
1:A:262:ASP:HB3	1:A:263:PRO:HD3	1.94	0.49
1:A:282:HIS:NE2	1:A:283:HIS:CD2	2.80	0.49
1:A:46:TYR:HA	1:A:453:ILE:HD11	1.94	0.49
1:A:227:LEU:O	1:A:230:TRP:HB3	2.12	0.49
1:A:29:PHE:CZ	1:A:401:LEU:CD1	2.96	0.49
1:A:410:GLY:O	1:A:501:LEU:HD11	2.13	0.49
1:A:281:PHE:H	1:A:298:HIS:CD2	2.31	0.48
1:A:233:HIS:O	1:A:236:VAL:HG22	2.14	0.48
2:B:145:TYR:CE1	2:B:166:VAL:CG2	2.96	0.48
1:A:9:SER:O	1:A:12:TYR:N	2.46	0.48
5:A:800:HEM:HMC2	5:A:800:HEM:CBC	2.41	0.47
1:A:518:ARG:HG2	1:A:518:ARG:NH2	2.28	0.47
1:A:10:ARG:O	1:A:13:GLU:N	2.47	0.47
1:A:456:VAL:HG23	1:A:456:VAL:O	2.14	0.47
1:A:12:TYR:O	1:A:16:PRO:N	2.47	0.47
1:A:562:TRP:HA	2:B:155:LEU:HG	1.96	0.47
1:A:403:TRP:C	1:A:406:PRO:HD2	2.35	0.47
1:A:48:ASN:HD21	1:A:457:PRO:HA	1.79	0.47
1:A:236:VAL:HG12	1:A:239:TRP:CZ3	2.49	0.47
1:A:79:VAL:HA	1:A:152:PHE:CZ	2.49	0.46
1:A:495:ARG:HA	1:A:495:ARG:HD3	1.79	0.46
1:A:259:LEU:HD23	1:A:511:VAL:HG22	1.98	0.46
1:A:465:VAL:HG23	1:A:466:PRO:CD	2.43	0.46
1:A:280:GLY:HA3	1:A:542:TYR:OH	2.16	0.46
1:A:120:ALA:O	1:A:123:PRO:HD2	2.16	0.46
3:C:13:LEU:HD22	3:C:13:LEU:HA	1.76	0.46
2:B:69:GLN:HA	2:B:69:GLN:HE21	1.81	0.45
2:B:83:VAL:HB	2:B:107:ILE:HG23	1.99	0.45
1:A:291:ASP:OD2	1:A:293:THR:HB	2.15	0.45
1:A:254:GLN:HE21	1:A:254:GLN:HA	1.81	0.45
2:B:145:TYR:HE1	2:B:166:VAL:CG2	2.25	0.45
2:B:93:ASN:HD21	2:B:162:GLY:HA2	1.81	0.45
2:B:92:PRO:O	2:B:95:ILE:HG13	2.16	0.45
1:A:330:ARG:HG3	1:A:334:GLY:HA3	1.99	0.45
1:A:401:LEU:O	1:A:405:LEU:HB2	2.17	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:400:SER:HA	1:A:403:TRP:NE1	2.32	0.45
1:A:96:GLU:OE2	1:A:180:PRO:HB2	2.17	0.44
2:B:113:ILE:HG22	2:B:151:GLN:HG2	1.98	0.44
1:A:465:VAL:CG2	1:A:466:PRO:HD3	2.44	0.44
2:B:101:ALA:O	2:B:103:ILE:HD12	2.16	0.44
1:A:321:GLU:HA	1:A:335:TRP:CE3	2.52	0.44
1:A:404:LEU:HG	1:A:408:LEU:HD11	1.99	0.44
1:A:359:GLY:HA3	1:A:388:GLN:NE2	2.33	0.44
1:A:255:ALA:O	1:A:323:ALA:HB1	2.18	0.44
1:A:107:TRP:O	1:A:111:TRP:HD1	2.01	0.44
1:A:337:ARG:HA	1:A:337:ARG:HD2	1.73	0.44
2:B:106:LYS:HG2	2:B:134:THR:HG23	1.99	0.44
3:C:2:GLU:CG	3:C:3:GLU:N	2.80	0.43
1:A:445:LEU:O	1:A:446:ASN:HB2	2.17	0.43
1:A:101:PRO:O	1:A:103:MET:N	2.52	0.43
1:A:449:ARG:O	1:A:450:ARG:HB2	2.18	0.43
1:A:184:TYR:CD2	1:A:266:ARG:HG2	2.53	0.43
1:A:52:TYR:N	1:A:53:PRO:HD2	2.33	0.43
2:B:86:PHE:O	2:B:88:PHE:N	2.51	0.43
1:A:459:ALA:O	2:B:146:ARG:NH1	2.24	0.43
1:A:477:LEU:HD12	5:A:800:HEM:HMB3	2.00	0.43
1:A:15:TYR:HA	1:A:17:GLU:OE1	2.19	0.43
1:A:267:LEU:HB3	1:A:311:MET:HE1	2.00	0.43
2:B:11:ILE:HG13	3:C:5:PRO:HG3	2.00	0.43
1:A:12:TYR:O	1:A:16:PRO:CA	2.67	0.43
1:A:281:PHE:H	1:A:298:HIS:HD2	1.65	0.43
1:A:123:PRO:HG3	1:A:144:ALA:HB3	2.01	0.43
1:A:230:TRP:C	1:A:230:TRP:CD1	2.92	0.43
1:A:264:MET:HA	1:A:264:MET:CE	2.49	0.43
2:B:103:ILE:N	2:B:103:ILE:HD12	2.34	0.42
1:A:477:LEU:CD1	5:A:800:HEM:HMB3	2.48	0.42
6:A:801:HAS:HHC	6:A:801:HAS:H122	2.00	0.42
1:A:489:PHE:O	1:A:493:LEU:HB2	2.20	0.42
1:A:339:LEU:HB3	1:A:346:PHE:CZ	2.54	0.42
2:B:74:THR:HG23	2:B:78:GLN:OE1	2.19	0.42
1:A:152:PHE:O	1:A:155:SER:HB3	2.20	0.42
2:B:93:ASN:ND2	2:B:162:GLY:HA2	2.35	0.42
1:A:160:ILE:HD13	1:A:194:PHE:HB2	2.02	0.42
1:A:70:THR:OG1	1:A:132:LEU:HD23	2.21	0.41
1:A:272:PHE:CZ	1:A:308:PRO:HB2	2.55	0.41
2:B:93:ASN:HA	2:B:94:PRO:HA	1.82	0.41
1:A:455:GLN:C	1:A:457:PRO:HD3	2.40	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:11:VAL:HG22	1:A:503:GLU:HG2	2.02	0.41
1:A:48:ASN:HA	1:A:48:ASN:HD22	1.69	0.41
1:A:330:ARG:HB2	1:A:331:GLY:H	1.69	0.41
1:A:546:LEU:O	1:A:550:PHE:HD1	2.03	0.41
1:A:403:TRP:O	1:A:406:PRO:HD2	2.21	0.41
2:B:97:VAL:O	2:B:166:VAL:HA	2.19	0.41
1:A:277:THR:N	1:A:278:PRO:CD	2.84	0.41
2:B:44:VAL:O	2:B:46:PRO:HD3	2.21	0.41
1:A:282:HIS:CD2	6:A:801:HAS:OMD	2.74	0.41
1:A:475:ILE:O	1:A:479:VAL:HG23	2.20	0.41
1:A:210:PRO:HB2	1:A:216:VAL:HG22	2.02	0.40
1:A:340:PRO:C	1:A:342:ASP:H	2.25	0.40
1:A:95:ARG:HD2	1:A:95:ARG:HH21	1.78	0.40
1:A:366:ASN:HB3	6:A:801:HAS:HMD	2.04	0.40
1:A:377:ASN:HB3	2:B:150:ASN:HB2	2.02	0.40
2:B:26:LEU:HA	2:B:26:LEU:HD12	1.98	0.40

All (28) 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:10:ARG:N	2:B:142:PRO:CA[3_444]	1.02	1.18
1:A:10:ARG:N	2:B:142:PRO:C[3_444]	1.30	0.90
1:A:10:ARG:N	2:B:142:PRO:CB[3_444]	1.39	0.81
1:A:13:GLU:OE1	2:B:166:VAL:O[3_444]	1.46	0.74
1:A:13:GLU:OE1	2:B:166:VAL:C[3_444]	1.51	0.69
1:A:10:ARG:CA	2:B:142:PRO:CA[3_444]	1.69	0.51
1:A:10:ARG:NH1	2:B:139:PHE:C[3_444]	1.71	0.49
1:A:8:ILE:N	2:B:144:GLU:N[3_444]	1.72	0.48
1:A:11:VAL:N	2:B:142:PRO:CB[3_444]	1.81	0.39
1:A:10:ARG:N	2:B:143:GLY:N[3_444]	1.83	0.37
1:A:10:ARG:NH1	2:B:139:PHE:CB[3_444]	1.83	0.37
1:A:9:SER:C	2:B:142:PRO:C[3_444]	1.90	0.30
1:A:9:SER:C	2:B:142:PRO:CB[3_444]	1.90	0.30
1:A:10:ARG:CA	2:B:142:PRO:CB[3_444]	1.91	0.29
1:A:11:VAL:O	2:B:168:GLU:OE1[3_444]	1.91	0.29
1:A:10:ARG:CD	2:B:99:GLN:CB[3_444]	1.97	0.23
1:A:10:ARG:C	2:B:142:PRO:CB[3_444]	1.98	0.22
1:A:9:SER:C	2:B:143:GLY:N[3_444]	2.02	0.18
1:A:10:ARG:NH1	2:B:139:PHE:CA[3_444]	2.03	0.17
1:A:10:ARG:NH1	2:B:140:LYS:N[3_444]	2.04	0.16
1:A:10:ARG:CB	2:B:142:PRO:CA[3_444]	2.06	0.14

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:10:ARG:CZ	2:B:139:PHE:CB[3_444]	2.06	0.14
1:A:10:ARG:NH2	2:B:139:PHE:CB[3_444]	2.07	0.13
1:A:15:TYR:CD2	2:B:168:GLU:OE2[3_444]	2.07	0.13
1:A:15:TYR:N	2:B:168:GLU:OE1[3_444]	2.09	0.11
1:A:8:ILE:N	2:B:144:GLU:CA[3_444]	2.15	0.05
1:A:14:ALA:N	2:B:168:GLU:CB[3_444]	2.16	0.04
1:A:13:GLU:OE1	2:B:167:LYS:N[3_444]	2.18	0.02

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	553/568 (97%)	508 (92%)	41 (7%)	4 (1%)	30	81
2	B	164/166 (99%)	148 (90%)	13 (8%)	3 (2%)	13	61
3	C	31/33 (94%)	26 (84%)	4 (13%)	1 (3%)	6	43
All	All	748/767 (98%)	682 (91%)	58 (8%)	8 (1%)	21	73

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	330	ARG
1	A	11	VAL
1	A	518	ARG
1	A	102	ASN
2	B	87	ALA
2	B	142	PRO
2	B	88	PHE
3	C	4	LYS

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	451/462 (98%)	431 (96%)	20 (4%)	39	81
2	B	136/136 (100%)	127 (93%)	9 (7%)	24	68
3	C	26/26 (100%)	24 (92%)	2 (8%)	18	60
All	All	613/624 (98%)	582 (95%)	31 (5%)	33	77

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	SER
1	A	48	ASN
1	A	133	TYR
1	A	168	ARG
1	A	216	VAL
1	A	230	TRP
1	A	254	GLN
1	A	262	ASP
1	A	274	LEU
1	A	339	LEU
1	A	354	LEU
1	A	369	PHE
1	A	405	LEU
1	A	430	LEU
1	A	439	LEU
1	A	449	ARG
1	A	452	TYR
1	A	472	LEU
1	A	495	ARG
1	A	522	LEU
2	B	19	LEU
2	B	26	LEU
2	B	58	VAL
2	B	72	VAL
2	B	77	ASN
2	B	111	ASP
2	B	128	LEU
2	B	159	ASN
2	B	166	VAL
3	C	3	GLU
3	C	13	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	42	GLN
1	A	48	ASN
1	A	76	ASN
1	A	254	GLN
1	A	298	HIS
1	A	388	GLN
2	B	69	GLN
2	B	77	ASN
2	B	122	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 7 ligands modelled in this entry, 4 are monoatomic - leaving 3 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	$\# Z > 2$
5	HEM	A	800	1	49,50,50	2.20	13 (26%)	46,82,82	1.89	8 (17%)
6	HAS	A	801	1	72,72,72	2.15	21 (29%)	86,109,109	3.57	32 (37%)
7	CUA	B	802	2	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
5	HEM	A	800	1	-	0/14/114/114	0/0/8/8
6	HAS	A	801	1	-	0/36/82/82	0/0/8/8
7	CUA	B	802	2	-	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)
6	A	801	HAS	C4A-C3A	6.77	1.48	1.40
5	A	800	HEM	C3C-C2C	-6.05	1.33	1.43
6	A	801	HAS	C4C-C3C	5.93	1.49	1.41
6	A	801	HAS	C1C-C2C	5.92	1.47	1.40
5	A	800	HEM	C3B-C2B	-5.43	1.34	1.43
5	A	800	HEM	C3D-C2D	4.81	1.52	1.43
5	A	800	HEM	C3B-CAB	4.55	1.54	1.40
5	A	800	HEM	C3C-CAC	4.53	1.54	1.40
5	A	800	HEM	C4A-C3A	4.47	1.45	1.40
6	A	801	HAS	FE-NA	4.28	2.10	1.92
6	A	801	HAS	FE-NB	3.94	2.09	1.92
6	A	801	HAS	C3B-C11	-3.91	1.48	1.52
6	A	801	HAS	FE-ND	3.87	2.09	1.92
6	A	801	HAS	FE-NC	3.86	2.09	1.92
5	A	800	HEM	C2D-C1D	3.59	1.45	1.44
5	A	800	HEM	FE-NA	3.41	2.07	1.92
6	A	801	HAS	C4C-CHD	3.34	1.48	1.38
6	A	801	HAS	C4A-CHD	3.30	1.48	1.39
6	A	801	HAS	C1A-C2A	3.25	1.49	1.43
6	A	801	HAS	C4D-CHA	3.11	1.47	1.38
6	A	801	HAS	C1B-CHB	3.10	1.47	1.38
6	A	801	HAS	C1A-CHA	3.09	1.48	1.39
6	A	801	HAS	C2A-C3A	3.07	1.46	1.37
5	A	800	HEM	C3D-C4D	-3.01	1.43	1.44
6	A	801	HAS	C4B-CHC	2.85	1.47	1.38
6	A	801	HAS	C3C-C2C	2.80	1.46	1.41
6	A	801	HAS	C1D-CHB	2.76	1.47	1.39
6	A	801	HAS	C1C-CHC	2.71	1.47	1.39
5	A	800	HEM	C2B-C1B	2.60	1.45	1.44
5	A	800	HEM	CMC-C2C	2.46	1.55	1.47
6	A	801	HAS	C2D-C3D	2.35	1.45	1.41
5	A	800	HEM	CMD-C2D	2.23	1.54	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	800	HEM	CMB-C2B	2.17	1.54	1.47
6	A	801	HAS	C2D-C1D	-2.06	1.36	1.40

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	801	HAS	C4C-C3C-C2C	-18.76	93.76	106.87
5	A	800	HEM	C3B-C4B-NB	-8.00	108.28	114.00
6	A	801	HAS	C4D-C3D-C2D	-7.89	100.25	106.80
6	A	801	HAS	C3D-C4D-ND	7.76	115.46	109.64
6	A	801	HAS	C3A-C4A-NA	7.72	115.24	109.41
6	A	801	HAS	C3C-C2C-C1C	7.36	111.31	107.00
6	A	801	HAS	C3C-C4C-NC	7.34	122.12	111.52
6	A	801	HAS	C2B-C1B-NB	6.53	114.34	109.41
6	A	801	HAS	C4A-C3A-C2A	-6.16	102.71	107.00
5	A	800	HEM	C4D-ND-C1D	5.18	110.47	105.16
6	A	801	HAS	C3B-C4B-NB	5.10	114.84	109.90
6	A	801	HAS	C3C-CAC-CBC	-4.96	115.67	125.95
6	A	801	HAS	CMC-C2C-C3C	-4.65	117.64	124.97
6	A	801	HAS	C2C-C3C-CAC	-4.58	117.94	127.33
6	A	801	HAS	C25-C23-C24	4.28	121.89	115.39
6	A	801	HAS	OMD-CMD-C2D	-4.20	116.25	125.32
6	A	801	HAS	C4C-NC-C1C	-4.12	101.33	106.76
6	A	801	HAS	C2A-C1A-NA	4.05	115.36	109.73
6	A	801	HAS	C1A-C2A-C3A	-3.99	102.79	106.92
6	A	801	HAS	C4B-C3B-C2B	-3.83	104.19	106.87
6	A	801	HAS	C1B-C2B-C3B	-3.80	104.78	107.00
6	A	801	HAS	CAD-C3D-C4D	3.72	131.39	124.67
6	A	801	HAS	C26-C15-C16	3.39	120.54	115.39
5	A	800	HEM	C2D-C1D-ND	-3.38	108.94	112.93
6	A	801	HAS	C2D-C1D-ND	3.25	111.80	108.64
5	A	800	HEM	C4A-C3A-C2A	3.20	109.22	107.00
6	A	801	HAS	C21-C22-C23	-3.04	121.23	127.80
6	A	801	HAS	C4B-NB-C1B	-2.81	101.37	105.58
5	A	800	HEM	CAD-CBD-CGD	-2.70	105.06	113.48
6	A	801	HAS	C32-C30-C31	2.58	121.21	114.62
6	A	801	HAS	C4B-C3B-C11	2.51	129.19	124.67
6	A	801	HAS	C27-C19-C20	2.50	119.19	115.39
6	A	801	HAS	C4A-NA-C1A	-2.45	103.54	106.76
6	A	801	HAS	C3A-C4A-CHD	-2.45	121.36	126.00
6	A	801	HAS	C13-C12-C11	-2.44	110.13	114.06
5	A	800	HEM	C1B-NB-C4B	2.32	107.53	105.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	801	HAS	C4D-ND-C1D	-2.20	103.87	106.76
5	A	800	HEM	O1D-CGD-CBD	-2.08	115.88	123.03
5	A	800	HEM	CBA-CAA-C2A	-2.07	109.04	112.69
6	A	801	HAS	O2D-CGD-CBD	2.03	121.41	114.22

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	555/568 (97%)	0.32	32 (5%) 22 5	86, 86, 86, 86	0
2	B	166/166 (100%)	0.56	9 (5%) 25 6	86, 86, 86, 86	0
3	C	33/33 (100%)	0.29	2 (6%) 21 5	86, 86, 86, 86	0
All	All	754/767 (98%)	0.37	43 (5%) 23 5	86, 86, 86, 86	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	498	LYS	6.2
1	A	519	ARG	5.2
1	A	497	ARG	4.8
1	A	496	GLU	4.7
1	A	516	GLU	4.5
1	A	8	ILE	4.4
1	A	495	ARG	4.4
1	A	502	ALA	4.3
2	B	7	ALA	3.8
1	A	515	PRO	3.5
1	A	9	SER	3.5
1	A	411	LYS	3.4
1	A	500	GLU	3.3
1	A	14	ALA	3.2
1	A	501	LEU	3.2
1	A	518	ARG	3.2
1	A	499	PRO	3.2
1	A	12	TYR	3.2
2	B	136	ARG	3.1
2	B	79	TYR	3.0
1	A	57	ARG	2.9
3	C	2	GLU	2.9
1	A	13	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	6	LYS	2.9
2	B	103	ILE	2.8
1	A	409	THR	2.8
3	C	3	GLU	2.7
1	A	514	GLY	2.7
2	B	4	GLU	2.6
1	A	410	GLY	2.6
1	A	493	LEU	2.6
1	A	412	PRO	2.6
1	A	548	GLN	2.5
2	B	25	MET	2.4
2	B	9	LYS	2.3
2	B	97	VAL	2.3
1	A	56	LYS	2.3
1	A	494	SER	2.3
1	A	341	TRP	2.3
1	A	142	HIS	2.2
1	A	215	LEU	2.1
1	A	332	LEU	2.1
1	A	53	PRO	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
8	XE	A	563	1/1	0.54	1.33	86,86,86,86	1

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	HAS	A	801	65/65	0.46	1.28	86,86,86,86	0
5	HEM	A	800	43/43	0.27	-0.08	86,86,86,86	0
8	XE	A	565	1/1	0.21	-0.22	86,86,86,86	1
8	XE	A	564	1/1	0.15	-1.17	86,86,86,86	1
7	CUA	B	802	2/2	0.14	-1.51	86,86,86,86	0
4	CU	A	803	1/1	0.30	-	86,86,86,86	0

6.5 Other polymers ⓘ

There are no such residues in this entry.