



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 02:57 PM GMT

PDB ID : 4KTF
Title : Crystal structure of Mycobacterium tuberculosis CYP121 in complex with 4, 4'-(3-amino-1H-pyrazole-4,5-diyl)diphenol
Authors : Hudson, S.A.
Deposited on : 2013-05-20
Resolution : 1.35 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>

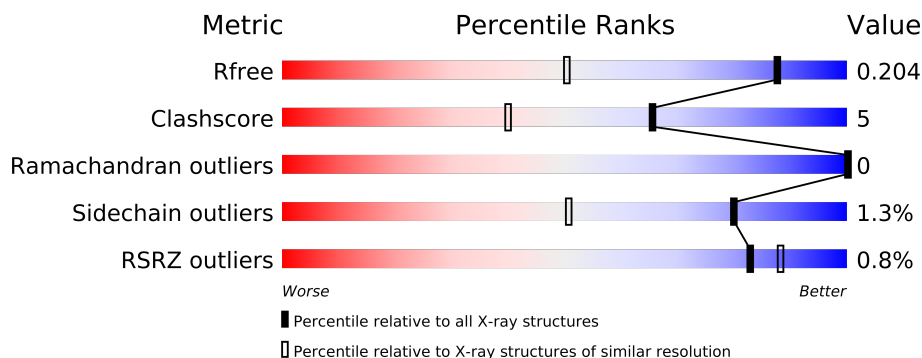
The following versions of software and data (see [references](#)) were used in the production of this report:

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 1.35 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	1519 (1.40-1.32)
Clashscore	79885	1707 (1.40-1.32)
Ramachandran outliers	78287	1662 (1.40-1.32)
Sidechain outliers	78261	1661 (1.40-1.32)
RSRZ outliers	66119	1519 (1.40-1.32)

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	395	

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	SO4	A	403	-	X
3	SO4	A	405	-	X

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 3739 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 P450 121.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	393	3068	1963	533	561	11	0	25	1

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



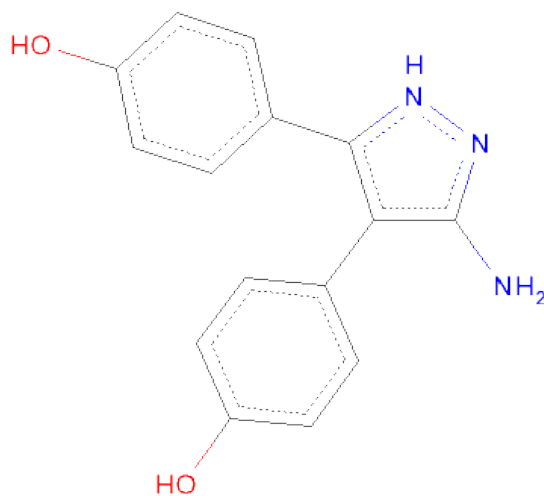
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



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

- Molecule 4 is 4,4'-(3-AMINO-1H-PYRAZOLE-4,5-DIYL)DIPHENOL (three-letter code: 1TM) (formula: C₁₅H₁₃N₃O₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			20	15	3	2		

- Molecule 5 is water.

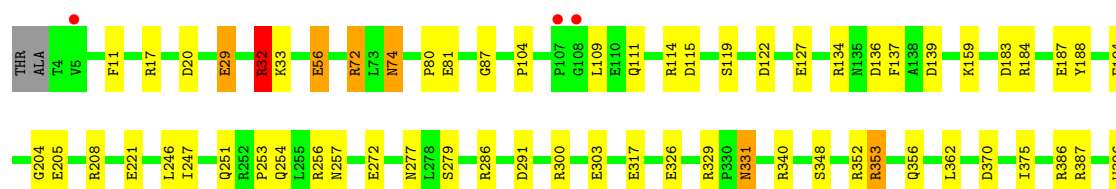
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	588	Total	O	0	0
			588	588		

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 P450 121

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 65 2 2	Depositor
Cell constants a, b, c, α , β , γ	78.20Å 78.20Å 263.19Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	35.71 – 1.35 35.71 – 1.35	Depositor EDS
% Data completeness (in resolution range)	91.3 (35.71-1.35) 91.4 (35.71-1.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.07 (at 1.35Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.166 , 0.195 0.179 , 0.204	Depositor DCC
R_{free} test set	4818 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	11.3	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 33.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Outliers	0 of 96125 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	3739	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.77% 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, 1TM, 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	2.26	18/3244 (0.6%)	1.46	41/4412 (0.9%)

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

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	194[A]	GLU	CG-CD	69.84	2.56	1.51
1	A	194[B]	GLU	CG-CD	69.84	2.56	1.51
1	A	81	GLU	CD-OE2	9.23	1.35	1.25
1	A	32	ARG	CZ-NH1	8.80	1.44	1.33
1	A	396	TRP	CG-CD1	7.56	1.47	1.36
1	A	81	GLU	CD-OE1	7.52	1.33	1.25
1	A	272	GLU	CD-OE2	6.70	1.33	1.25
1	A	87	GLY	N-CA	-6.63	1.36	1.46
1	A	279[A]	SER	CB-OG	-6.37	1.33	1.42
1	A	279[B]	SER	CB-OG	-6.37	1.33	1.42
1	A	56	GLU	CD-OE2	6.29	1.32	1.25
1	A	205	GLU	CD-OE1	-6.28	1.18	1.25
1	A	256	ARG	CZ-NH2	5.92	1.40	1.33
1	A	204	GLY	N-CA	5.90	1.54	1.46
1	A	326	GLU	CD-OE2	-5.75	1.19	1.25
1	A	29	GLU	CD-OE2	5.42	1.31	1.25
1	A	221	GLU	CD-OE2	5.22	1.31	1.25
1	A	253	PRO	N-CA	-5.16	1.38	1.47

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	32	ARG	NE-CZ-NH1	19.66	130.13	120.30
1	A	32	ARG	NE-CZ-NH2	-17.19	111.71	120.30
1	A	352	ARG	NE-CZ-NH1	-10.51	115.04	120.30
1	A	370	ASP	CB-CG-OD2	-9.64	109.62	118.30
1	A	122	ASP	CB-CG-OD1	9.43	126.78	118.30
1	A	208	ARG	NE-CZ-NH1	-9.30	115.65	120.30
1	A	136	ASP	CB-CG-OD2	8.57	126.01	118.30
1	A	329	ARG	NE-CZ-NH2	8.23	124.41	120.30
1	A	184	ARG	NE-CZ-NH2	-8.12	116.24	120.30
1	A	194[A]	GLU	CG-CD-OE2	8.10	134.49	118.30
1	A	194[B]	GLU	CG-CD-OE2	8.10	134.49	118.30
1	A	256	ARG	NE-CZ-NH1	-8.08	116.26	120.30
1	A	137	PHE	CB-CG-CD2	7.61	126.12	120.80
1	A	286	ARG	NE-CZ-NH1	7.32	123.96	120.30
1	A	56	GLU	OE1-CD-OE2	7.22	131.96	123.30
1	A	353	ARG	NE-CZ-NH1	7.21	123.91	120.30
1	A	194[A]	GLU	CB-CG-CD	-7.06	95.14	114.20
1	A	194[B]	GLU	CB-CG-CD	-7.06	95.14	114.20
1	A	291	ASP	CB-CG-OD1	6.69	124.32	118.30
1	A	188	TYR	CG-CD1-CE1	-6.68	115.95	121.30
1	A	139	ASP	CB-CG-OD2	-6.67	112.29	118.30
1	A	72	ARG	NE-CZ-NH2	6.58	123.59	120.30
1	A	139	ASP	CB-CG-OD1	6.55	124.19	118.30
1	A	115	ASP	CB-CG-OD2	-6.53	112.43	118.30
1	A	256	ARG	NE-CZ-NH2	6.46	123.53	120.30
1	A	303	GLU	OE1-CD-OE2	-6.37	115.66	123.30
1	A	137	PHE	CB-CG-CD1	-6.23	116.44	120.80
1	A	329	ARG	NH1-CZ-NH2	-6.13	112.66	119.40
1	A	11	PHE	CB-CG-CD1	5.89	124.92	120.80
1	A	11	PHE	CB-CG-CD2	-5.82	116.72	120.80
1	A	194[A]	GLU	CG-CD-OE1	-5.71	106.87	118.30
1	A	194[B]	GLU	CG-CD-OE1	-5.71	106.87	118.30
1	A	183	ASP	CB-CG-OD1	5.69	123.42	118.30
1	A	291	ASP	CB-CG-OD2	-5.67	113.19	118.30
1	A	56	GLU	CG-CD-OE2	-5.66	106.98	118.30
1	A	20	ASP	CB-CG-OD1	5.62	123.36	118.30
1	A	348	SER	N-CA-CB	5.44	118.66	110.50
1	A	387	ARG	CD-NE-CZ	5.10	130.75	123.60
1	A	370	ASP	CB-CG-OD1	5.06	122.86	118.30
1	A	134[A]	ARG	NE-CZ-NH2	5.06	122.83	120.30
1	A	134[B]	ARG	NE-CZ-NH2	5.06	122.83	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	72	ARG	Sidechain

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	3068	0	3083	32	0
2	A	43	0	30	1	0
3	A	20	0	0	0	0
4	A	20	0	12	0	0
5	A	588	0	0	13	2
All	All	3739	0	3125	33	2

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 5.

All (33) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:17:ARG:NE	5:A:1042:HOH:O	2.12	0.82
1:A:251[B]:GLN:NE2	1:A:375:ILE:HD13	1.95	0.81
1:A:353:ARG:HH11	1:A:356:GLN:HE21	1.26	0.80
1:A:251[B]:GLN:NE2	1:A:375:ILE:CD1	2.45	0.79
1:A:127[A]:GLU:CG	5:A:1043:HOH:O	2.34	0.75
1:A:127[B]:GLU:CG	5:A:1043:HOH:O	2.38	0.71
1:A:29:GLU:HG2	1:A:32:ARG:HB3	1.73	0.70
1:A:187[A]:GLU:OE2	5:A:1027:HOH:O	2.11	0.69
1:A:254:GLN:CG	5:A:981:HOH:O	2.41	0.68
1:A:74:ASN:HD22	1:A:74:ASN:C	2.02	0.63
1:A:254:GLN:HG3	5:A:981:HOH:O	1.99	0.62
1:A:353:ARG:HH11	1:A:356:GLN:NE2	1.96	0.61
1:A:300:ARG:CG	5:A:1079:HOH:O	2.48	0.60
1:A:111:GLN:NE2	1:A:114[B]:ARG:HH21	1.98	0.60
1:A:254:GLN:HG2	5:A:981:HOH:O	2.03	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:247:ILE:O	1:A:251[B]:GLN:HG3	2.09	0.53
1:A:331:ASN:H	1:A:331:ASN:HD22	1.56	0.52
2:A:401:HEM:HMC1	2:A:401:HEM:HBC2	1.93	0.51
1:A:257:ASN:ND2	5:A:1053:HOH:O	2.22	0.49
1:A:29:GLU:CG	1:A:32:ARG:HB3	2.43	0.48
1:A:317:GLU:HG3	5:A:1066:HOH:O	2.12	0.48
1:A:159[B]:LYS:CG	5:A:1026:HOH:O	2.63	0.46
1:A:56:GLU:OE2	1:A:340[B]:ARG:HD2	2.16	0.46
1:A:246:LEU:HD22	1:A:362[B]:LEU:HD22	1.99	0.45
1:A:104:PRO:HA	1:A:109:LEU:HD23	1.98	0.44
1:A:251[B]:GLN:HE22	1:A:375:ILE:HD13	1.76	0.44
1:A:331:ASN:N	1:A:331:ASN:HD22	2.16	0.41
1:A:33[B]:LYS:CD	5:A:877:HOH:O	2.68	0.41
1:A:246:LEU:HD22	1:A:362[B]:LEU:CD2	2.50	0.41
1:A:80[A]:PRO:HG3	5:A:947:HOH:O	2.21	0.41

All (2) 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)
5:A:849:HOH:O	5:A:849:HOH:O[10_665]	1.48	0.72
5:A:814:HOH:O	5:A:953:HOH:O[10_555]	1.96	0.24

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	417/395 (106%)	409 (98%)	8 (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	330/325 (102%)	326 (99%)	4 (1%)	82	54

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	32	ARG
1	A	74	ASN
1	A	277	ASN
1	A	331	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	111	GLN
1	A	154	GLN
1	A	331	ASN
1	A	342	GLN
1	A	356	GLN
1	A	385	GLN

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 ⓘ

6 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	401	1,5	49,50,50	4.23	20 (40%)	46,82,82	2.82	9 (19%)
3	SO4	A	402	-	4,4,4	2.54	2 (50%)	6,6,6	2.23	2 (33%)
3	SO4	A	403	-	4,4,4	1.06	0	6,6,6	0.35	0
3	SO4	A	404	-	4,4,4	1.75	2 (50%)	6,6,6	2.37	2 (33%)
3	SO4	A	405	-	4,4,4	1.81	1 (25%)	6,6,6	0.84	0
4	1TM	A	406	-	22,22,22	1.83	6 (27%)	30,31,31	2.08	8 (26%)

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	401	1,5	-	0/14/114/114	0/0/8/8
3	SO4	A	402	-	-	0/0/0/0	0/0/0/0
3	SO4	A	403	-	-	0/0/0/0	0/0/0/0
3	SO4	A	404	-	-	0/0/0/0	0/0/0/0
3	SO4	A	405	-	-	0/0/0/0	0/0/0/0
4	1TM	A	406	-	-	0/8/8/8	0/3/3/3

All (31) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HEM	C2D-C1D	18.74	1.49	1.44
2	A	401	HEM	C3D-C4D	17.15	1.48	1.44
2	A	401	HEM	C3C-C2C	-5.29	1.34	1.43
2	A	401	HEM	CHB-C1B	4.69	1.42	1.35
2	A	401	HEM	C3B-C2B	-4.30	1.36	1.43
2	A	401	HEM	C3D-C2D	-3.90	1.36	1.43
2	A	401	HEM	C4A-C3A	3.84	1.45	1.40
4	A	406	1TM	CAK-NAM	3.83	1.44	1.34
3	A	402	SO4	O2-S	3.79	1.59	1.47
2	A	401	HEM	C2B-C1B	3.77	1.45	1.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	HEM	C4C-NC	3.44	1.42	1.38
3	A	402	SO4	O1-S	3.34	1.58	1.47
2	A	401	HEM	C3B-CAB	3.34	1.50	1.40
2	A	401	HEM	FE-NB	3.16	2.09	1.97
4	A	406	1TM	CAS-CAR	3.06	1.44	1.38
2	A	401	HEM	CHA-C4D	3.04	1.40	1.35
2	A	401	HEM	C1B-NB	-2.94	1.33	1.39
4	A	406	1TM	CAC-CAB	-2.93	1.43	1.47
2	A	401	HEM	C4B-NB	-2.85	1.31	1.37
2	A	401	HEM	C4A-NA	2.80	1.42	1.36
4	A	406	1TM	CAO-CAP	2.74	1.44	1.38
3	A	405	SO4	O1-S	2.71	1.55	1.47
2	A	401	HEM	C3C-CAC	2.68	1.48	1.40
2	A	401	HEM	C4A-CHB	-2.60	1.32	1.39
4	A	406	1TM	CAJ-CAN	-2.51	1.45	1.50
2	A	401	HEM	FE-NA	2.51	2.03	1.92
3	A	404	SO4	O4-S	2.28	1.55	1.47
4	A	406	1TM	CAG-CAF	2.15	1.43	1.38
3	A	404	SO4	O1-S	2.05	1.53	1.47
2	A	401	HEM	O1D-CGD	2.04	1.29	1.22
2	A	401	HEM	C3B-C4B	2.00	1.46	1.44

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	HEM	C3B-C4B-NB	-12.71	104.91	114.00
2	A	401	HEM	C1B-NB-C4B	9.16	114.54	105.16
2	A	401	HEM	CHC-C4B-NB	6.85	130.27	124.58
3	A	404	SO4	O4-S-O3	5.33	131.59	109.08
4	A	406	1TM	CAJ-CAB-CAC	4.48	138.78	129.27
4	A	406	1TM	CAH-CAC-CAB	4.38	123.90	120.65
3	A	402	SO4	O4-S-O3	4.31	127.28	109.08
4	A	406	1TM	CAB-NAA-NAL	4.24	116.66	111.18
4	A	406	1TM	CAC-CAB-NAA	-4.00	116.08	121.52
4	A	406	1TM	CAJ-CAB-NAA	-3.44	104.89	106.71
4	A	406	1TM	CAN-CAJ-CAB	3.02	132.18	126.71
2	A	401	HEM	C1A-CHA-C4D	-2.91	123.64	127.47
2	A	401	HEM	CHC-C1C-NC	-2.76	122.33	124.73
3	A	402	SO4	O2-S-O1	-2.75	100.45	109.53
4	A	406	1TM	CAS-CAN-CAJ	2.69	125.27	120.77
2	A	401	HEM	CHB-C1B-NB	2.59	127.87	124.31
4	A	406	1TM	CAR-CAS-CAN	2.20	124.52	121.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	HEM	CMA-C3A-C4A	-2.05	125.46	128.62
3	A	404	SO4	O4-S-O2	-2.05	91.54	110.12
2	A	401	HEM	CHD-C1D-ND	-2.04	122.89	124.58
2	A	401	HEM	O2A-CGA-O1A	-2.03	118.13	123.30

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	393/395 (99%)	-0.17	3 (0%) 83 88	5, 10, 22, 33	2 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	107	PRO	4.0
1	A	5	VAL	2.6
1	A	108	GLY	2.4

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	SO4	A	403	5/5	0.25	7.33	27,33,38,40	0
3	SO4	A	405	5/5	0.16	6.11	16,25,32,32	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	SO4	A	402	5/5	0.10	1.69	21,23,27,34	0
3	SO4	A	404	5/5	0.09	0.73	15,16,22,25	0
2	HEM	A	401	43/43	0.07	-0.35	5,8,11,13	0
4	1TM	A	406	20/20	0.05	-1.20	6,8,10,10	0

6.5 Other polymers ⓘ

There are no such residues in this entry.