



Full wwPDB X-ray Structure Validation Report

Nov 6, 2014 – 07:17 PM EST

PDB ID : 4J6C
Title : The 1.9 Å crystal structure of CYP154C5 from *Nocardia farcinica* in complex with progesterone
Authors : Herzog, K.; Hoffmann, K.M.
Deposited on : 2013-02-11
Resolution : 1.90 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>

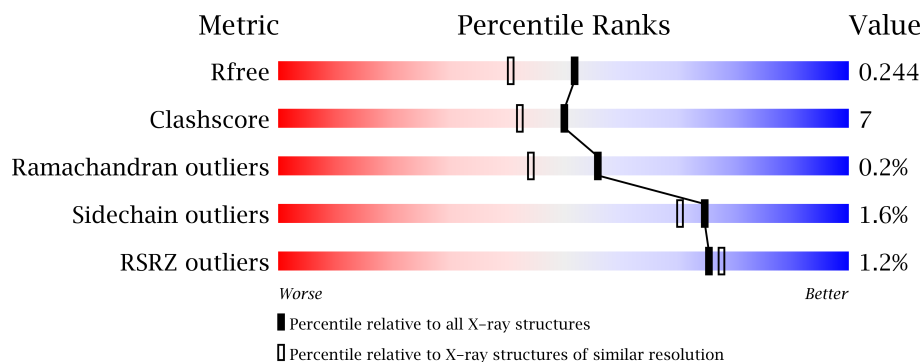
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.16 November 2013
Xtriage (Phenix) : dev-1439
EDS : stable24103
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.1.3
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable24103

1 Overall quality at a glance



The reported resolution of this entry is 1.90 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	3684 (1.90-1.90)
Clashscore	79885	4465 (1.90-1.90)
Ramachandran outliers	78287	4413 (1.90-1.90)
Sidechain outliers	78261	4414 (1.90-1.90)
RSRZ outliers	66119	3686 (1.90-1.90)

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	410	
1	B	410	

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

Mol	Type	Chain	Res	Geometry	Electron density
4	MG	A	503	-	X
5	FMT	B	507	-	X

2 Entry composition i

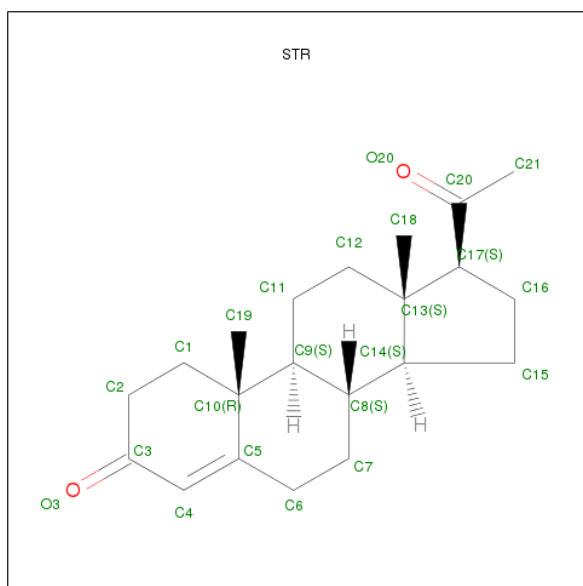
There are 7 unique types of molecules in this entry. The entry contains 7183 atoms, of which 0 are hydrogen and 0 are deuterium.

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 monooxygenase.

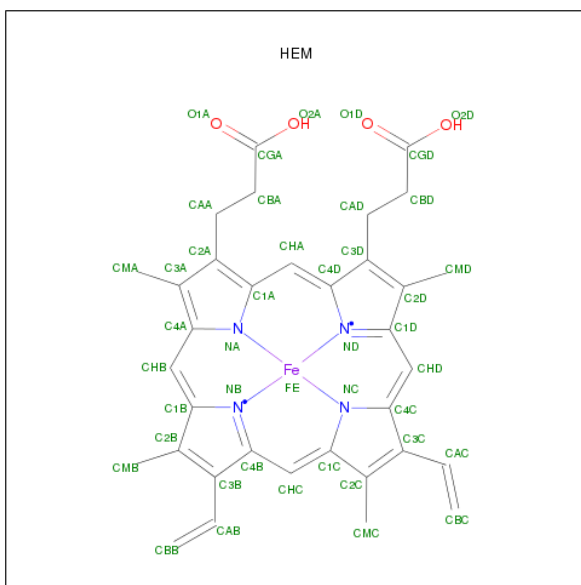
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	407	Total	C	N	O	S	0	6	0
			3181	2014	562	595	10			
1	B	406	Total	C	N	O	S	0	5	0
			3160	2004	555	591	10			

- Molecule 2 is PROGESTERONE (three-letter code: STR) (formula: $C_{21}H_{30}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			23	21	2		
2	B	1	Total	C	O	0	0
			23	21	2		

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).

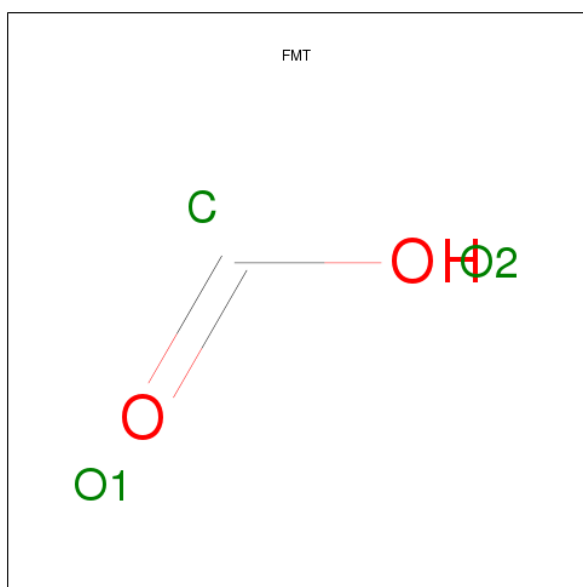


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
3	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Mg 1 1	0	0
4	A	1	Total Mg 1 1	0	0

- Molecule 5 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



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

- Molecule 6 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	1	Total K 1 1	0	0
6	A	1	Total K 1 1	0	0

- Molecule 7 is water.

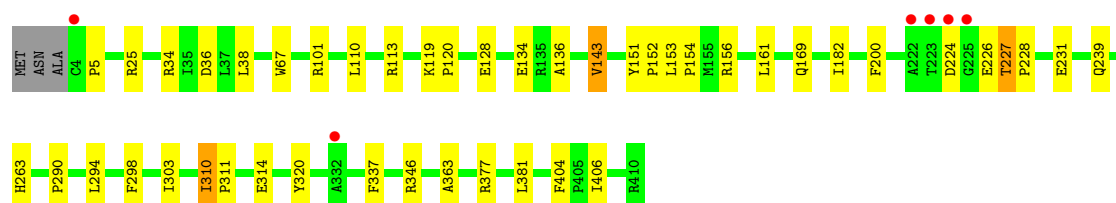
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	352	Total O 352 352	0	0
7	B	342	Total O 342 342	0	0

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.

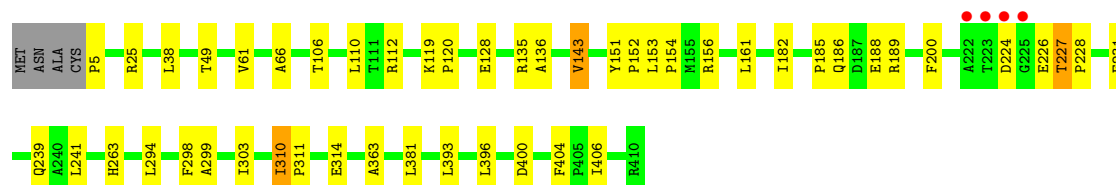
- Molecule 1: Cytochrome P450 monooxygenase

Chain A: 



- Molecule 1: Cytochrome P450 monooxygenase

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	103.00Å 103.00Å 217.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	82.55 – 1.90 42.01 – 1.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (82.55-1.90) 100.0 (42.01-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.67 (at 1.89Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.192 , 0.238 0.200 , 0.244	Depositor DCC
R_{free} test set	3448 reflections (5.35%)	DCC
Wilson B-factor (Å ²)	20.3	Xtriage
Anisotropy	0.063	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 39.4	EDS
Estimated twinning fraction	0.038 for h,-h-k,-l	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 67932 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7183	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 22.35 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 5.8016e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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, MG, FMT, K, STR

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	0.65	0/3246	0.48	1/4424 (0.0%)
1	B	0.67	1/3225 (0.0%)	0.46	1/4396 (0.0%)
All	All	0.66	1/6471 (0.0%)	0.47	2/8820 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	135	ARG	CZ-NH1	-9.02	1.21	1.33

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	101	ARG	NE-CZ-NH2	5.78	123.19	120.30
1	B	135	ARG	NE-CZ-NH2	5.38	122.99	120.30

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	3181	0	3131	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3160	0	3105	51	0
2	A	23	0	30	0	0
2	B	23	0	30	0	0
3	A	43	0	30	4	0
3	B	43	0	30	6	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	3	0	2	0	0
5	B	9	0	3	2	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
7	A	352	0	0	3	0
7	B	342	0	0	15	0
All	All	7183	0	6361	96	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 7.

All (96) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:226:GLU:HB3	1:B:227:THR:HA	1.21	1.16
1:B:226:GLU:CB	1:B:227:THR:HA	1.75	1.14
1:B:226:GLU:HB3	1:B:227:THR:CA	1.81	1.10
1:B:298:PHE:HB2	7:B:860:HOH:O	1.55	1.06
1:B:303:ILE:HB	1:B:310:ILE:HG23	1.39	1.03
1:A:226:GLU:HB3	1:A:227:THR:HA	1.38	1.00
1:A:226:GLU:CB	1:A:227:THR:HA	1.94	0.97
1:A:226:GLU:HB3	1:A:227:THR:CA	1.96	0.94
3:B:501:HEM:HBB2	3:B:501:HEM:HHC	1.53	0.89
1:B:61:VAL:HG22	7:B:860:HOH:O	1.72	0.89
1:B:303:ILE:HB	1:B:310:ILE:CG2	2.05	0.87
1:A:346:ARG:CG	1:A:346:ARG:NE	2.41	0.83
1:A:310:ILE:CD1	1:A:314:GLU:HB3	2.10	0.82
1:B:61:VAL:CG2	7:B:860:HOH:O	2.29	0.80
1:B:393:LEU:HD12	7:B:883:HOH:O	1.82	0.80
1:B:5:PRO:HB2	7:B:832:HOH:O	1.82	0.79
3:A:502:HEM:HBB2	3:A:502:HEM:HHC	1.63	0.79
1:A:25:ARG:NE	1:A:25:ARG:CG	2.47	0.78
1:B:226:GLU:HB3	1:B:227:THR:C	2.04	0.76
1:B:381:LEU:HD11	1:B:406:ILE:HD12	1.68	0.75
1:B:66:ALA:HB3	7:B:860:HOH:O	1.89	0.72
1:B:25:ARG:NE	1:B:25:ARG:CG	2.55	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:381:LEU:HD11	1:A:406:ILE:HD12	1.72	0.68
1:A:310:ILE:HD12	1:A:310:ILE:C	2.14	0.68
1:A:34:ARG:CG	1:A:34:ARG:NE	2.57	0.68
1:B:381:LEU:HD11	1:B:406:ILE:CD1	2.24	0.67
1:A:381:LEU:CD1	1:A:406:ILE:HD12	2.25	0.67
1:B:226:GLU:HB2	1:B:227:THR:HA	1.71	0.67
1:B:200:PHE:CD2	1:B:239:GLN:HG2	2.30	0.67
1:B:381:LEU:CD1	1:B:406:ILE:HD12	2.24	0.66
1:B:363:ALA:HB1	3:B:501:HEM:CBB	2.26	0.66
1:A:226:GLU:HB3	1:A:227:THR:C	2.15	0.65
1:A:381:LEU:HD11	1:A:406:ILE:CD1	2.26	0.65
1:A:200:PHE:CD2	1:A:239:GLN:HG2	2.32	0.64
1:B:186:GLN:HE22	1:B:189:ARG:HE	1.45	0.64
1:A:310:ILE:HD13	1:A:314:GLU:HB3	1.82	0.62
1:B:299:ALA:HB2	1:B:310:ILE:HD12	1.82	0.61
1:B:231[B]:GLU:HB2	7:B:884:HOH:O	2.00	0.61
3:B:501:HEM:HBB2	3:B:501:HEM:CHC	2.29	0.61
1:A:303:ILE:HB	1:A:310:ILE:HG23	1.81	0.61
1:B:393:LEU:CD1	7:B:883:HOH:O	2.43	0.60
1:A:363:ALA:HB1	3:A:502:HEM:CBB	2.33	0.59
1:A:153:LEU:HB3	1:A:154:PRO:HD3	1.84	0.58
1:B:310:ILE:HD11	1:B:314:GLU:HB3	1.85	0.58
1:B:66:ALA:CB	7:B:860:HOH:O	2.48	0.58
1:B:153:LEU:HB3	1:B:154:PRO:HD3	1.86	0.58
5:B:507:FMT:H	7:B:778:HOH:O	2.04	0.57
1:B:263:HIS:HE1	7:B:647:HOH:O	1.87	0.57
1:B:400:ASP:CG	7:B:883:HOH:O	2.43	0.56
1:B:110:LEU:HD21	1:B:161:LEU:HD21	1.87	0.56
1:B:310:ILE:HD13	1:B:310:ILE:C	2.27	0.55
1:A:320:TYR:CE1	7:A:743:HOH:O	2.54	0.55
1:A:226:GLU:HB2	1:A:227:THR:HA	1.87	0.55
3:B:501:HEM:HBC2	3:B:501:HEM:HMC2	1.89	0.55
1:B:186:GLN:NE2	1:B:189:ARG:HE	2.05	0.52
1:A:310:ILE:HD11	1:A:314:GLU:HB3	1.89	0.52
1:B:226:GLU:HB3	1:B:228:PRO:N	2.25	0.51
1:A:310:ILE:HD12	1:A:311:PRO:N	2.26	0.51
1:B:310:ILE:HD13	1:B:311:PRO:O	2.11	0.51
1:B:151:TYR:HB3	1:B:152:PRO:HD3	1.93	0.50
1:A:231:GLU:HB2	7:A:777:HOH:O	2.11	0.50
3:B:501:HEM:HBC2	3:B:501:HEM:CMC	2.42	0.50
1:A:5:PRO:HB2	7:A:742:HOH:O	2.12	0.50
1:A:128:GLU:CD	1:A:156[A]:ARG:HH12	2.16	0.48

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:38:LEU:HD11	1:A:182:ILE:HD11	1.96	0.48
1:A:110:LEU:HD21	1:A:161:LEU:HD21	1.94	0.48
1:B:299:ALA:CB	1:B:310:ILE:HD12	2.42	0.48
1:B:396:LEU:HB2	7:B:811:HOH:O	2.14	0.48
1:A:294:LEU:C	1:A:294:LEU:HD13	2.35	0.47
1:B:186:GLN:HE21	1:B:186:GLN:HA	1.80	0.46
1:B:49:THR:HG23	7:B:936:HOH:O	2.14	0.46
1:A:226:GLU:HB3	1:A:228:PRO:N	2.30	0.46
1:B:363:ALA:CB	3:B:501:HEM:CBB	2.93	0.46
1:A:151:TYR:HB3	1:A:152:PRO:HD3	1.96	0.46
1:B:310:ILE:CD1	1:B:314:GLU:HB3	2.45	0.46
1:B:182:ILE:HA	1:B:182:ILE:HD12	1.84	0.45
1:A:156[A]:ARG:HD3	1:A:169:GLN:HE22	1.81	0.45
1:B:106:THR:HG21	1:B:241:LEU:HD21	1.98	0.45
1:B:310:ILE:CD1	1:B:310:ILE:C	2.85	0.45
1:B:128:GLU:CD	1:B:156[A]:ARG:HH12	2.20	0.44
1:A:182:ILE:HD12	1:A:182:ILE:HA	1.87	0.44
1:B:38:LEU:HD11	1:B:182:ILE:HD11	1.99	0.44
5:B:507:FMT:C	7:B:778:HOH:O	2.64	0.44
3:A:502:HEM:HBB2	3:A:502:HEM:CHC	2.39	0.43
1:B:294:LEU:C	1:B:294:LEU:HD13	2.39	0.42
1:A:119:LYS:HB3	1:A:120:PRO:HD3	2.02	0.42
1:A:263:HIS:CE1	1:A:337:PHE:HB3	2.54	0.42
1:B:119:LYS:HB3	1:B:120:PRO:HD3	2.02	0.41
1:B:112:ARG:CG	1:B:112:ARG:NE	2.84	0.41
1:B:186:GLN:HA	1:B:186:GLN:NE2	2.34	0.41
1:B:136:ALA:HB1	1:B:143:VAL:HG13	2.03	0.41
1:A:134[B]:GLU:OE2	1:A:377:ARG:NH2	2.50	0.41
1:A:136:ALA:HB1	1:A:143:VAL:HG13	2.03	0.40
1:A:290:PRO:HG2	3:A:502:HEM:HAB	2.03	0.40
1:B:185:PRO:HB2	1:B:188[B]:GLU:HG3	2.03	0.40
1:A:67:TRP:CG	1:A:298:PHE:CE1	3.10	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	411/410 (100%)	400 (97%)	10 (2%)	1 (0%)	56	44
1	B	409/410 (100%)	398 (97%)	10 (2%)	1 (0%)	56	44
All	All	820/820 (100%)	798 (97%)	20 (2%)	2 (0%)	56	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	224	ASP
1	B	224	ASP

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	329/341 (96%)	323 (98%)	6 (2%)	71	66
1	B	325/341 (95%)	321 (99%)	4 (1%)	82	80
All	All	654/682 (96%)	644 (98%)	10 (2%)	75	73

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	ASP
1	A	113	ARG
1	A	143	VAL
1	A	227	THR
1	A	310	ILE
1	A	404	PHE
1	B	143	VAL
1	B	227	THR
1	B	310	ILE
1	B	404	PHE

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	53	GLN
1	A	169	GLN
1	A	392	ASN
1	B	53	GLN
1	B	186	GLN
1	B	263	HIS
1	B	392	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 12 ligands modelled in this entry, 4 are monoatomic - leaving 8 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$
2	STR	A	501	-	26,26,26	0.89	1 (3%)	42,42,42	1.26	4 (9%)
3	HEM	A	502	1	42,50,50	2.73	15 (35%)	27,82,82	2.54	6 (22%)
5	FMT	A	504	-	2,2,2	1.37	0	1,1,1	0.62	0
3	HEM	B	501	1	42,50,50	2.94	17 (40%)	27,82,82	2.58	6 (22%)
2	STR	B	503	-	26,26,26	0.80	0	42,42,42	1.14	4 (9%)
5	FMT	B	505	-	2,2,2	0.71	0	1,1,1	0.66	0
5	FMT	B	506	-	2,2,2	0.86	0	1,1,1	0.64	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	FMT	B	507	-	2,2,2	1.26	0	1,1,1	0.62	0

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	STR	A	501	-	-	0/4/62/62	0/4/4/4
3	HEM	A	502	1	-	0/14/114/114	0/0/8/8
5	FMT	A	504	-	-	0/0/0/0	0/0/0/0
3	HEM	B	501	1	-	0/14/114/114	0/0/8/8
2	STR	B	503	-	-	0/4/62/62	0/4/4/4
5	FMT	B	505	-	-	0/0/0/0	0/0/0/0
5	FMT	B	506	-	-	0/0/0/0	0/0/0/0
5	FMT	B	507	-	-	0/0/0/0	0/0/0/0

All (33) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	501	HEM	C1A-NA	7.28	1.48	1.36
3	A	502	HEM	C1A-NA	6.96	1.48	1.36
3	B	501	HEM	C4A-NA	6.58	1.47	1.36
3	B	501	HEM	CHC-C4B	6.07	1.47	1.39
3	B	501	HEM	C1C-NC	5.97	1.48	1.33
3	A	502	HEM	C1C-NC	5.67	1.47	1.33
3	A	502	HEM	C4A-NA	5.45	1.45	1.36
3	A	502	HEM	C4C-NC	5.33	1.46	1.33
3	B	501	HEM	C4A-C3A	5.32	1.52	1.43
3	B	501	HEM	C4C-NC	4.98	1.46	1.33
3	A	502	HEM	C4A-C3A	4.66	1.51	1.43
3	A	502	HEM	CHC-C4B	4.64	1.45	1.39
3	B	501	HEM	FE-NB	4.56	2.12	1.95
3	B	501	HEM	CMD-C2D	4.53	1.52	1.45
3	A	502	HEM	CMD-C2D	4.39	1.52	1.45
3	A	502	HEM	CMC-C2C	4.26	1.52	1.45
3	B	501	HEM	CMC-C2C	4.01	1.52	1.45
3	A	502	HEM	FE-NB	3.80	2.09	1.95
3	A	502	HEM	FE-ND	3.74	2.10	1.95
3	B	501	HEM	C1A-C2A	3.68	1.49	1.43
3	B	501	HEM	FE-ND	3.59	2.10	1.95
3	A	502	HEM	C1B-C2B	-3.44	1.42	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	502	HEM	C1A-C2A	3.14	1.48	1.43
3	A	502	HEM	C3B-CAB	2.77	1.49	1.40
3	B	501	HEM	CMB-C2B	2.76	1.50	1.45
3	B	501	HEM	C3B-CAB	2.73	1.49	1.40
3	B	501	HEM	C2A-C3A	2.71	1.45	1.37
3	A	502	HEM	C3C-CAC	2.65	1.49	1.40
3	B	501	HEM	C3C-CAC	2.41	1.48	1.40
3	B	501	HEM	C1C-C2C	2.36	1.49	1.43
3	B	501	HEM	O1A-CGA	2.19	1.29	1.22
3	A	502	HEM	C1C-C2C	2.08	1.49	1.43
2	A	501	STR	C13-C17	-2.04	1.52	1.56

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	501	HEM	C3A-C4A-NA	7.92	114.79	109.50
3	A	502	HEM	C3A-C4A-NA	7.68	114.63	109.50
3	B	501	HEM	C4A-NA-C1A	-7.39	98.73	107.93
3	A	502	HEM	C4A-NA-C1A	-7.24	98.92	107.93
3	B	501	HEM	C2A-C1A-NA	4.14	115.48	109.73
3	B	501	HEM	CBD-CAD-C3D	-3.99	105.84	114.51
3	A	502	HEM	C2A-C1A-NA	3.70	114.88	109.73
3	A	502	HEM	CBD-CAD-C3D	-3.58	106.74	114.51
2	A	501	STR	C6-C5-C10	3.40	120.49	116.69
2	A	501	STR	C21-C20-C17	2.75	121.38	117.60
2	B	503	STR	C6-C5-C10	2.63	119.63	116.69
3	B	501	HEM	C4A-C3A-C2A	-2.60	105.19	107.00
2	A	501	STR	C6-C5-C4	-2.58	117.67	120.91
2	B	503	STR	C6-C5-C4	-2.50	117.77	120.91
2	A	501	STR	O20-C20-C17	-2.19	118.62	121.78
3	B	501	HEM	CMC-C2C-C3C	2.18	129.42	124.26
3	A	502	HEM	CMC-C2C-C3C	2.15	129.35	124.26
3	A	502	HEM	C4A-C3A-C2A	-2.14	105.51	107.00
2	B	503	STR	C1-C10-C9	-2.10	105.81	108.61
2	B	503	STR	C5-C4-C3	-2.09	121.21	123.78

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	407/410 (99%)	0.05	6 (1%) 70 72	16, 27, 41, 82	0
1	B	406/410 (99%)	0.06	4 (0%) 79 81	17, 28, 41, 87	0
All	All	813/820 (99%)	0.06	10 (1%) 75 78	16, 27, 41, 87	0

All (10) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	223	THR	4.7
1	A	4	CYS	4.7
1	B	223	THR	3.8
1	A	225	GLY	3.7
1	A	222	ALA	3.6
1	B	225	GLY	3.4
1	B	224	ASP	3.1
1	A	224	ASP	2.6
1	B	222	ALA	2.1
1	A	332	ALA	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
5	FMT	B	507	3/3	0.18	3.09	26,26,27,32	0
4	MG	A	503	1/1	0.15	2.78	10,10,10,10	0
2	STR	A	501	23/23	0.15	1.92	15,17,20,24	0
4	MG	B	502	1/1	0.14	1.88	10,10,10,10	0
5	FMT	A	504	3/3	0.16	1.08	24,24,27,31	0
6	K	B	504	1/1	0.12	0.90	70,70,70,70	0
2	STR	B	503	23/23	0.14	0.67	15,18,21,26	0
3	HEM	A	502	43/43	0.12	0.62	14,17,21,26	0
3	HEM	B	501	43/43	0.12	0.35	17,19,22,26	0
6	K	A	505	1/1	0.10	-0.15	72,72,72,72	0
5	FMT	B	506	3/3	0.09	-1.19	40,40,46,46	0
5	FMT	B	505	3/3	0.11	-1.37	48,48,49,51	0

6.5 Other polymers ⓘ

There are no such residues in this entry.