



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 05:11 PM GMT

PDB ID : 4HGH
Title : Crystal structure of P450 BM3 5F5 heme domain variant complexed with styrene (dataset I)
Authors : Shehzad, A.; Panneerselvam, S.; Bocola, M.; Mueller-Dieckmann, J.; Wilmanns, M.; Schwaneberg, U.
Deposited on : 2012-10-08
Resolution : 1.40 Å(reported)

This is a Full wwPDB X-ray Structure 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/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

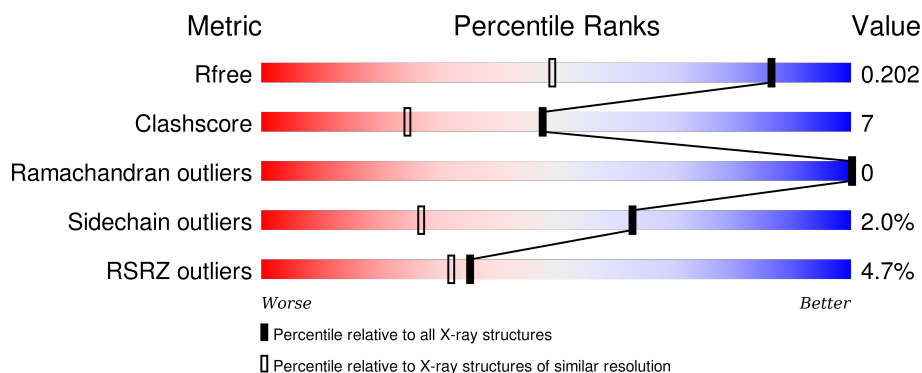
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.40 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}	91344	1199 (1.40-1.40)
Clashscore	102246	1295 (1.40-1.40)
Ramachandran outliers	100387	1259 (1.40-1.40)
Sidechain outliers	100360	1258 (1.40-1.40)
RSRZ outliers	91569	1198 (1.40-1.40)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	455	<div> <div>4%</div> <div>88%</div> <div>11%</div> </div>
1	B	455	<div> <div>6%</div> <div>84%</div> <div>15%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	GOL	A	503	-	-	-	X
4	GOL	A	506	-	-	-	X
5	MES	A	504	-	-	-	X
6	PEG	A	505	-	-	X	X

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 8824 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bifunctional P-450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	455	Total	C	N	O	S	0	28	0
			3817	2458	636	704	19			
1	B	455	Total	C	N	O	S	0	33	0
			3851	2480	642	710	19			

There are 4 discrepancies between the modelled and reference sequences:

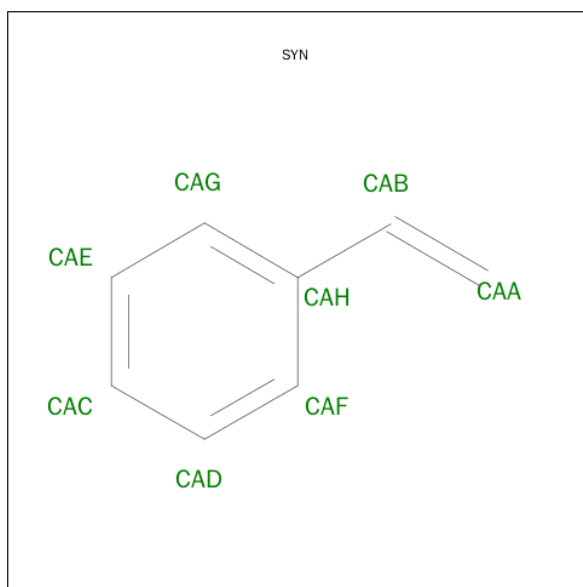
Chain	Residue	Modelled	Actual	Comment	Reference
A	87	ALA	PHE	ENGINEERED MUTATION	UNP P14779
A	235	ALA	THR	ENGINEERED MUTATION	UNP P14779
B	87	ALA	PHE	ENGINEERED MUTATION	UNP P14779
B	235	ALA	THR	ENGINEERED MUTATION	UNP P14779

- 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	B	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is ETHENYLBENZENE (three-letter code: SYN) (formula: C_8H_8).



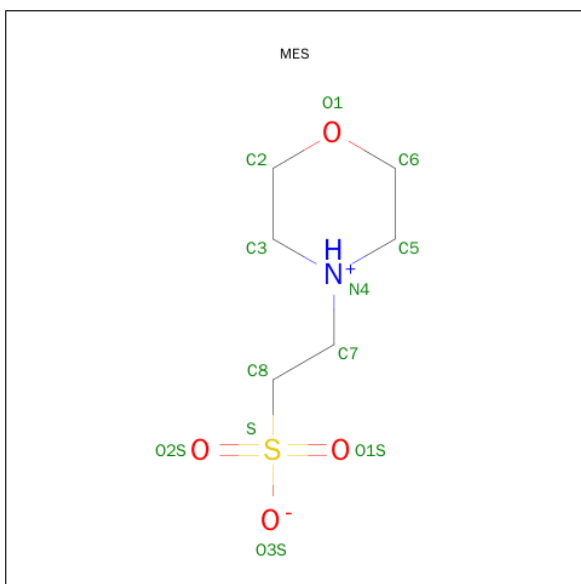
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C	0	0
			8	8		
3	B	1	Total	C	0	0
			8	8		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 5 is 2-(N-MORPHOLINO)-ETHANESULFONIC ACID (three-letter code: MES) (formula: $C_6H_{13}NO_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	N	O	S	0	0
			12	6	1	4	1		

- Molecule 6 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			7	4	3		

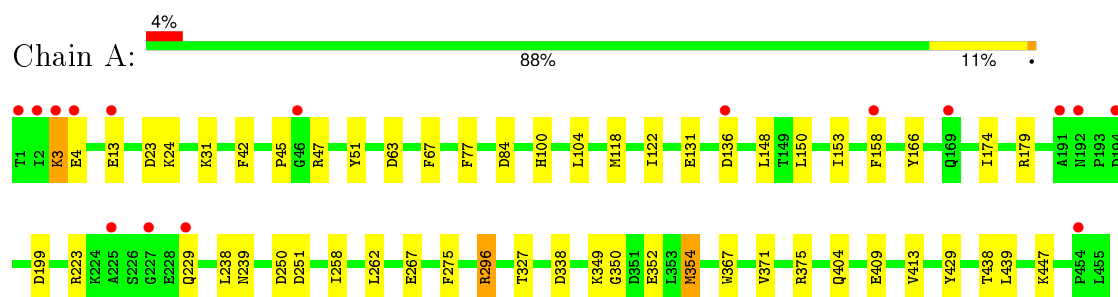
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	536	Total	O	0	0
			536	536		
7	B	475	Total	O	0	0
			475	475		

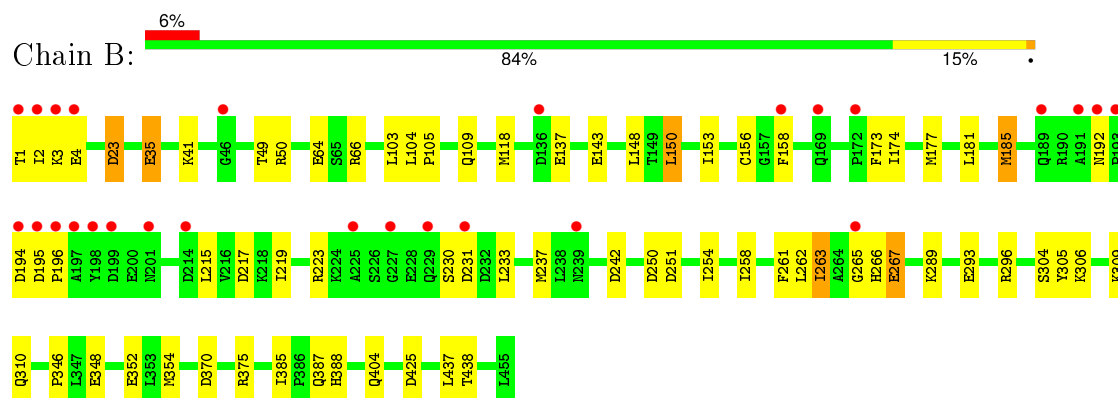
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($\text{RSRZ} > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Bifunctional P-450/NADPH-P450 reductase



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4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	59.09 Å 149.27 Å 65.16 Å 90.00° 98.38° 90.00°	Depositor
Resolution (Å)	19.87 – 1.40 19.87 – 1.40	Depositor EDS
% Data completeness (in resolution range)	96.3 (19.87-1.40) 96.4 (19.87-1.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.54 (at 1.40 Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.190 , 0.204 0.188 , 0.202	Depositor DCC
R_{free} test set	818 reflections (0.39%)	DCC
Wilson B-factor (Å ²)	12.3	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 56.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 210565 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	8824	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.54% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

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: GOL, PEG, SYN, MES, HEM

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	1.29	8/3978 (0.2%)	1.17	22/5373 (0.4%)
1	B	1.29	13/4016 (0.3%)	1.13	14/5425 (0.3%)
All	All	1.29	21/7994 (0.3%)	1.15	36/10798 (0.3%)

All (21) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	64	GLU	CD-OE2	-8.58	1.16	1.25
1	A	409[A]	GLU	CG-CD	-8.32	1.39	1.51
1	A	409[B]	GLU	CG-CD	-8.32	1.39	1.51
1	B	137	GLU	CD-OE1	-6.99	1.18	1.25
1	B	143	GLU	CD-OE2	-6.11	1.19	1.25
1	B	35	GLU	CG-CD	5.92	1.60	1.51
1	B	137	GLU	CD-OE2	-5.86	1.19	1.25
1	B	293	GLU	CD-OE1	-5.84	1.19	1.25
1	B	348	GLU	CD-OE2	-5.75	1.19	1.25
1	A	352	GLU	CD-OE2	5.72	1.31	1.25
1	B	156	CYS	C-O	5.54	1.33	1.23
1	B	143	GLU	CD-OE1	-5.53	1.19	1.25
1	A	375	ARG	CZ-NH1	5.47	1.40	1.33
1	B	352	GLU	CG-CD	-5.41	1.43	1.51
1	A	166	TYR	CG-CD2	5.37	1.46	1.39
1	A	131	GLU	CD-OE2	5.30	1.31	1.25
1	A	429	TYR	CG-CD1	5.18	1.45	1.39
1	A	275	PHE	CG-CD2	5.03	1.46	1.38
1	B	304	SER	CB-OG	5.02	1.48	1.42
1	B	310[A]	GLN	CG-CD	5.00	1.62	1.51
1	B	310[B]	GLN	CG-CD	5.00	1.62	1.51

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	296[A]	ARG	NE-CZ-NH1	7.47	124.03	120.30
1	A	296[B]	ARG	NE-CZ-NH1	7.47	124.03	120.30
1	A	84	ASP	CB-CG-OD2	-6.72	112.25	118.30
1	B	251	ASP	CB-CG-OD2	6.35	124.02	118.30
1	B	50	ARG	NE-CZ-NH2	-5.95	117.32	120.30
1	B	296	ARG	NE-CZ-NH1	5.93	123.26	120.30
1	B	263[A]	ILE	CB-CA-C	-5.91	99.77	111.60
1	B	263[B]	ILE	CB-CA-C	-5.91	99.77	111.60
1	B	23	ASP	CB-CG-OD1	5.86	123.57	118.30
1	A	223	ARG	NE-CZ-NH1	5.79	123.19	120.30
1	A	354[A]	MET	CG-SD-CE	5.78	109.45	100.20
1	A	354[B]	MET	CG-SD-CE	5.78	109.45	100.20
1	A	250	ASP	CB-CG-OD1	5.77	123.49	118.30
1	A	179	ARG	NE-CZ-NH2	-5.71	117.45	120.30
1	B	217	ASP	CB-CG-OD1	5.69	123.42	118.30
1	B	223	ARG	NE-CZ-NH1	5.67	123.14	120.30
1	A	77	PHE	CB-CG-CD2	-5.64	116.85	120.80
1	A	199	ASP	CB-CG-OD1	5.63	123.37	118.30
1	A	63	ASP	CB-CG-OD2	-5.62	113.24	118.30
1	A	338	ASP	CB-CG-OD1	5.56	123.30	118.30
1	A	239	ASN	CB-CA-C	-5.53	99.33	110.40
1	B	425	ASP	CB-CG-OD2	-5.49	113.36	118.30
1	A	375	ARG	NE-CZ-NH2	-5.49	117.56	120.30
1	B	237	MET	CG-SD-CE	5.44	108.90	100.20
1	B	23	ASP	CB-CG-OD2	-5.44	113.41	118.30
1	B	242	ASP	CB-CG-OD2	-5.27	113.56	118.30
1	A	250	ASP	CB-CG-OD2	-5.20	113.62	118.30
1	A	63	ASP	CB-CG-OD1	5.19	122.97	118.30
1	B	66	ARG	NE-CZ-NH2	-5.19	117.71	120.30
1	A	199	ASP	CB-CG-OD2	-5.18	113.64	118.30
1	A	47	ARG	NE-CZ-NH2	5.17	122.89	120.30
1	A	251	ASP	CB-CG-OD1	5.16	122.95	118.30
1	A	42	PHE	CB-CG-CD2	-5.15	117.19	120.80
1	A	67	PHE	CB-CG-CD1	-5.10	117.23	120.80
1	A	238	LEU	CB-CG-CD2	5.09	119.66	111.00
1	B	305	TYR	CB-CG-CD1	-5.04	117.97	121.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3817	0	3879	31	0
1	B	3851	0	3919	81	0
2	A	43	0	30	2	0
2	B	43	0	30	2	0
3	A	8	0	8	0	0
3	B	8	0	8	0	0
4	A	12	0	16	1	0
5	A	12	0	12	0	0
5	B	12	0	12	1	0
6	A	7	0	10	4	0
7	A	536	0	0	11	0
7	B	475	0	0	13	0
All	All	8824	0	7924	117	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (117) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150[A]:LEU:HD11	1:B:174:ILE:CD1	1.30	1.58
1:B:150[A]:LEU:CD1	1:B:174:ILE:CD1	1.99	1.39
1:B:150[A]:LEU:CD1	1:B:174:ILE:HD13	1.56	1.35
1:B:309[A]:LYS:NZ	7:B:968:HOH:O	1.61	1.33
1:B:158[A]:PHE:CE1	1:B:258:ILE:HG12	1.78	1.17
1:B:150[A]:LEU:CD1	1:B:174:ILE:HD11	1.75	1.14
1:B:375[B]:ARG:NH2	5:B:503:MES:O2S	1.84	1.09
1:B:261:PHE:O	1:B:265[B]:GLY:N	1.87	1.07
1:B:150[A]:LEU:HD11	1:B:174:ILE:HD11	1.25	1.06
1:B:158[B]:PHE:CE2	1:B:262[B]:LEU:HD21	1.91	1.04
1:B:158[B]:PHE:HE2	1:B:262[B]:LEU:HD21	1.21	1.03
1:B:150[A]:LEU:HD12	1:B:174:ILE:HD13	1.46	0.98
1:B:150[A]:LEU:HD11	1:B:174:ILE:CG1	1.94	0.98
1:B:150[B]:LEU:HD13	1:B:266[B]:HIS:CE1	1.98	0.98
1:B:267[B]:GLU:CG	1:B:438[B]:THR:HG21	1.95	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:267[B]:GLU:HG2	1:B:438[B]:THR:HG21	1.50	0.93
1:B:150[A]:LEU:HD11	1:B:174:ILE:HD13	1.18	0.93
1:B:150[B]:LEU:CD1	1:B:266[B]:HIS:CE1	2.53	0.92
1:B:158[A]:PHE:HE1	1:B:258:ILE:HG12	1.31	0.88
1:B:150[B]:LEU:HD13	1:B:266[B]:HIS:CG	2.08	0.87
1:A:267[B]:GLU:OE2	7:A:1021:HOH:O	1.95	0.85
1:B:370:ASP:OD2	1:B:375[A]:ARG:NH1	2.10	0.84
1:B:150[B]:LEU:HD13	1:B:266[B]:HIS:ND1	1.91	0.84
1:B:150[A]:LEU:HD13	1:B:174:ILE:HD11	1.62	0.81
1:B:233[B]:LEU:HD21	1:B:261:PHE:CD2	2.15	0.80
1:A:118[B]:MET:SD	7:A:866:HOH:O	2.40	0.79
1:A:3:LYS:HE2	7:A:1004:HOH:O	1.81	0.79
1:A:327:THR:O	1:A:438[B]:THR:HG23	1.87	0.75
1:A:438[B]:THR:HG22	1:A:439:LEU:N	2.01	0.75
1:B:185:MET:HE1	1:B:437:LEU:HD23	1.69	0.74
1:B:150[B]:LEU:HD13	1:B:266[B]:HIS:CD2	2.22	0.74
1:B:375[A]:ARG:NE	7:B:877:HOH:O	2.13	0.73
1:B:118[A]:MET:SD	7:B:802:HOH:O	2.46	0.73
1:A:136:ASP:O	1:A:447:LYS:NZ	2.24	0.71
1:B:261:PHE:O	1:B:265[B]:GLY:CA	2.39	0.69
1:B:267[B]:GLU:HG2	1:B:438[B]:THR:CG2	2.21	0.69
1:B:158[B]:PHE:HE2	1:B:262[B]:LEU:CD2	2.01	0.68
1:B:150[B]:LEU:HD11	1:B:266[B]:HIS:CE1	2.29	0.67
1:B:177:MET:SD	1:B:263[B]:ILE:HG12	2.35	0.67
1:A:3:LYS:CE	7:A:1004:HOH:O	2.41	0.66
1:B:233[B]:LEU:HD21	1:B:261:PHE:CE2	2.32	0.65
1:B:153[A]:ILE:HG13	7:B:812:HOH:O	1.97	0.64
1:B:2:ILE:HD13	1:B:346[A]:PRO:HG3	1.81	0.62
1:B:150[A]:LEU:HD11	1:B:174:ILE:HG12	1.80	0.62
1:B:267[B]:GLU:HG3	1:B:438[B]:THR:HG21	1.77	0.61
1:A:349:LYS:HE2	6:A:505:PEG:H32	1.83	0.60
1:B:266[A]:HIS:CE1	1:B:267[A]:GLU:OE2	2.54	0.59
1:A:23[A]:ASP:OD2	7:A:1061:HOH:O	2.15	0.59
1:B:150[B]:LEU:HD13	1:B:266[B]:HIS:NE2	2.18	0.59
1:A:153[A]:ILE:HG13	7:A:978:HOH:O	2.03	0.58
1:A:350:GLY:HA2	6:A:505:PEG:H41	1.84	0.58
1:A:158[B]:PHE:CE1	1:A:258:ILE:HG12	2.39	0.57
6:A:505:PEG:H11	7:A:768:HOH:O	2.03	0.56
1:B:109:GLN:NE2	1:B:404[B]:GLN:HG3	2.20	0.56
1:A:438[B]:THR:HG22	1:A:439:LEU:H	1.68	0.56
1:B:177:MET:CE	1:B:266[B]:HIS:HE1	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:261:PHE:O	1:B:265[B]:GLY:HA3	2.05	0.56
1:B:375[A]:ARG:CG	7:B:877:HOH:O	2.54	0.55
1:B:49[A]:THR:CG2	1:B:354[A]:MET:HG2	2.36	0.55
1:B:385:ILE:HD12	7:B:957:HOH:O	2.07	0.55
1:A:150[A]:LEU:HD13	1:A:174:ILE:HD13	1.89	0.55
1:B:103:LEU:HD22	1:B:233[B]:LEU:CD1	2.37	0.55
1:B:158[B]:PHE:CZ	1:B:262[B]:LEU:HD21	2.39	0.54
1:B:185:MET:CE	1:B:437:LEU:HD23	2.37	0.54
1:B:23:ASP:OD2	7:B:1004:HOH:O	2.19	0.52
1:B:375[A]:ARG:HG2	7:B:877:HOH:O	2.09	0.52
1:A:350:GLY:CA	6:A:505:PEG:H41	2.39	0.52
1:A:438[B]:THR:CG2	1:A:439:LEU:H	2.22	0.52
1:B:150[B]:LEU:CD1	1:B:266[B]:HIS:ND1	2.68	0.52
1:B:387:GLN:HG2	1:B:388:HIS:CD2	2.47	0.50
1:A:438[B]:THR:CG2	1:A:439:LEU:N	2.67	0.50
1:B:181:LEU:HD11	1:B:263[B]:ILE:HG23	1.92	0.50
1:B:185:MET:HE2	1:B:437:LEU:HA	1.94	0.49
1:B:267[A]:GLU:HG3	7:B:963:HOH:O	2.11	0.49
1:A:148[B]:LEU:HD11	1:A:413:VAL:HG21	1.94	0.49
1:B:263[A]:ILE:HG22	1:B:263[A]:ILE:O	2.11	0.49
1:A:296[A]:ARG:NH1	7:A:995:HOH:O	2.40	0.49
1:B:375[A]:ARG:CD	7:B:877:HOH:O	2.60	0.49
1:B:262[B]:LEU:O	1:B:266[B]:HIS:ND1	2.39	0.48
1:B:177:MET:HE1	1:B:266[B]:HIS:HE1	1.78	0.48
1:A:100:HIS:HE1	4:A:506:GOL:O2	1.96	0.48
1:B:258:ILE:O	1:B:262[B]:LEU:HG	2.14	0.47
1:A:51:TYR:CE2	1:A:354[B]:MET:HG2	2.49	0.47
1:A:31[A]:LYS:HE2	7:A:746:HOH:O	2.15	0.47
1:B:173:PHE:CE2	1:B:262[B]:LEU:HD13	2.50	0.47
2:A:501:HEM:HBC2	2:A:501:HEM:HMC2	1.97	0.46
2:B:501:HEM:HBC2	2:B:501:HEM:HMC2	1.96	0.46
1:A:153[B]:ILE:HG23	1:A:262:LEU:HD23	1.96	0.46
1:B:195:ASP:HA	1:B:196:PRO:HD3	1.83	0.45
1:B:103:LEU:HD22	1:B:233[B]:LEU:HD11	1.98	0.45
2:A:501:HEM:HBC2	2:A:501:HEM:CMC	2.47	0.45
1:B:173:PHE:HB2	1:B:215:LEU:HD22	1.98	0.44
1:B:309[B]:LYS:HD2	1:B:309[B]:LYS:HA	1.74	0.44
1:B:250:ASP:O	1:B:254:ILE:HG13	2.17	0.44
1:A:150[B]:LEU:HG	1:A:174:ILE:HD13	1.99	0.44
1:B:1:THR:HG22	1:B:2:ILE:N	2.32	0.44
1:A:23[B]:ASP:HB3	1:A:24:LYS:HG3	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:233[B]:LEU:HD11	1:B:261:PHE:CE2	2.53	0.44
1:B:35:GLU:HG3	7:B:1030:HOH:O	2.17	0.44
2:B:501:HEM:HBC2	2:B:501:HEM:CMC	2.48	0.43
1:B:306:LYS:HB2	1:B:306:LYS:HE2	1.76	0.43
1:B:177:MET:SD	1:B:263[B]:ILE:CG1	3.05	0.43
1:B:267[A]:GLU:HG3	7:B:1074:HOH:O	2.18	0.43
1:B:289[A]:LYS:HG3	7:B:824:HOH:O	2.18	0.43
1:A:404[B]:GLN:NE2	7:A:981:HOH:O	2.50	0.43
1:A:31[A]:LYS:HG3	7:A:746:HOH:O	2.19	0.42
1:B:266[A]:HIS:CE1	1:B:267[A]:GLU:HG2	2.54	0.42
1:B:354[B]:MET:HE3	1:B:354[B]:MET:HB2	1.65	0.42
1:B:103:LEU:CD2	1:B:233[B]:LEU:HD12	2.50	0.42
1:B:2:ILE:HD13	1:B:346[B]:PRO:HG2	2.02	0.42
1:B:266[A]:HIS:HE1	1:B:267[A]:GLU:OE2	2.03	0.41
1:B:104:LEU:HB3	1:B:105:PRO:HD3	2.03	0.41
1:A:150[B]:LEU:HG	1:A:174:ILE:CD1	2.51	0.41
1:A:367:TRP:HB2	1:A:371:VAL:HG12	2.02	0.41
1:B:215:LEU:O	1:B:219:ILE:HG12	2.21	0.41
1:A:104[B]:LEU:HD23	1:A:104[B]:LEU:HA	1.78	0.41
1:A:122:ILE:HG22	1:A:148[B]:LEU:HD23	2.03	0.41

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	481/455 (106%)	466 (97%)	15 (3%)	0	100	100
1	B	486/455 (107%)	468 (96%)	18 (4%)	0	100	100
All	All	967/910 (106%)	934 (97%)	33 (3%)	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	424/397 (107%)	419 (99%)	5 (1%)	78	52
1	B	428/397 (108%)	415 (97%)	13 (3%)	48	12
All	All	852/794 (107%)	834 (98%)	18 (2%)	63	24

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	LYS
1	A	4	GLU
1	A	13	GLU
1	A	45	PRO
1	A	229	GLN
1	B	3	LYS
1	B	4	GLU
1	B	41	LYS
1	B	148	LEU
1	B	150[A]	LEU
1	B	150[B]	LEU
1	B	185	MET
1	B	192	ASN
1	B	194	ASP
1	B	230	SER
1	B	231	ASP
1	B	267[A]	GLU
1	B	267[B]	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	169	GLN
1	A	239	ASN
1	A	403	GLN
1	B	169	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

9 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	501	1,7	30,50,50	2.61	10 (33%)	24,82,82	3.01	17 (70%)
3	SYN	A	502	-	8,8,8	1.66	2 (25%)	9,9,9	0.86	0
4	GOL	A	503	-	5,5,5	0.59	0	5,5,5	1.24	1 (20%)
5	MES	A	504	-	11,12,12	0.68	0	14,16,16	2.24	5 (35%)
6	PEG	A	505	-	6,6,6	1.02	0	5,5,5	1.71	2 (40%)
4	GOL	A	506	-	5,5,5	0.59	0	5,5,5	0.98	0
2	HEM	B	501	1,7	30,50,50	2.82	11 (36%)	24,82,82	2.98	14 (58%)
3	SYN	B	502	-	8,8,8	1.73	4 (50%)	9,9,9	2.35	3 (33%)
5	MES	B	503	-	11,12,12	0.85	0	14,16,16	3.27	7 (50%)

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	501	1,7	-	0/10/54/54	0/0/8/8
3	SYN	A	502	-	-	0/2/2/2	0/1/1/1
4	GOL	A	503	-	-	0/4/4/4	0/0/0/0
5	MES	A	504	-	-	0/6/14/14	0/1/1/1
6	PEG	A	505	-	-	0/4/4/4	0/0/0/0
4	GOL	A	506	-	-	0/4/4/4	0/0/0/0
2	HEM	B	501	1,7	-	0/10/54/54	0/0/8/8
3	SYN	B	502	-	-	0/2/2/2	0/1/1/1
5	MES	B	503	-	-	0/6/14/14	0/1/1/1

All (27) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	C3B-C4B	-9.08	1.43	1.51
2	B	501	HEM	C3D-C4D	-7.96	1.41	1.51
2	A	501	HEM	C3B-C4B	-7.83	1.44	1.51
2	A	501	HEM	C3D-C4D	-6.04	1.43	1.51
2	A	501	HEM	C2C-C1C	-4.21	1.44	1.52
2	B	501	HEM	C2C-C1C	-3.87	1.45	1.52
2	A	501	HEM	C2D-C1D	-3.25	1.41	1.51
2	B	501	HEM	C2D-C1D	-2.72	1.43	1.51
2	A	501	HEM	C2B-C1B	-2.66	1.43	1.51
2	A	501	HEM	C2D-C3D	-2.45	1.47	1.54
2	B	501	HEM	C2B-C1B	-2.27	1.44	1.51
2	B	501	HEM	C2D-C3D	-2.08	1.48	1.54
3	B	502	SYN	CAH-CAB	-2.05	1.41	1.49
2	B	501	HEM	C4C-NC	-2.01	1.33	1.36
3	B	502	SYN	CAE-CAG	2.23	1.43	1.38
3	A	502	SYN	CAF-CAH	2.32	1.43	1.39
3	B	502	SYN	CAG-CAH	2.43	1.44	1.39
3	B	502	SYN	CAA-CAB	2.61	1.47	1.28
2	B	501	HEM	CAA-C2A	2.72	1.56	1.52
2	A	501	HEM	FE-ND	2.72	2.11	1.97
2	A	501	HEM	C1C-NC	2.74	1.39	1.36
2	B	501	HEM	FE-NB	2.86	2.12	1.97
2	B	501	HEM	FE-ND	3.05	2.13	1.97
2	A	501	HEM	CAA-C2A	3.29	1.57	1.52
3	A	502	SYN	CAA-CAB	3.31	1.52	1.28
2	A	501	HEM	FE-NC	4.46	2.13	1.95
2	B	501	HEM	FE-NC	4.89	2.15	1.95

All (49) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	HEM	CMA-C3A-C4A	-4.37	121.13	128.36
2	B	501	HEM	C3B-C4B-NB	-3.37	105.18	111.63
2	A	501	HEM	CAA-C2A-C1A	-3.17	123.56	127.01
3	B	502	SYN	CAD-CAF-CAH	-3.12	116.67	120.64
2	B	501	HEM	CMA-C3A-C4A	-3.12	123.21	128.36
2	A	501	HEM	C3B-C4B-NB	-2.49	106.87	111.63
5	B	503	MES	O1-C6-C5	-2.39	106.36	111.84
2	A	501	HEM	C2C-C1C-NC	-2.35	106.24	110.21
4	A	503	GOL	C3-C2-C1	-2.30	102.11	111.12
5	B	503	MES	O3S-S-O2S	-2.26	106.35	111.61
2	B	501	HEM	C3C-CAC-CBC	-2.20	121.09	124.46
5	A	504	MES	O1-C2-C3	-2.05	107.13	111.84
3	B	502	SYN	CAE-CAG-CAH	2.01	123.20	120.64
2	A	501	HEM	C3B-CAB-CBB	2.04	127.59	124.46
2	A	501	HEM	CMA-C3A-C2A	2.05	129.53	125.24
5	B	503	MES	C7-N4-C5	2.16	116.79	111.27
5	A	504	MES	C7-N4-C3	2.29	117.14	111.27
6	A	505	PEG	O2-C3-C4	2.36	121.32	110.43
2	A	501	HEM	CMD-C2D-C3D	2.37	124.83	114.35
2	B	501	HEM	C2C-C1C-CHC	2.41	127.34	123.68
2	A	501	HEM	C1D-CHD-C4C	2.45	129.91	125.82
6	A	505	PEG	O2-C2-C1	2.57	122.26	110.43
2	B	501	HEM	CAA-CBA-CGA	2.58	117.48	112.75
2	A	501	HEM	CAA-CBA-CGA	2.70	117.69	112.75
2	B	501	HEM	C4B-CHC-C1C	2.79	130.49	125.82
2	A	501	HEM	C2D-C3D-C4D	2.90	106.41	101.50
2	A	501	HEM	C4B-CHC-C1C	2.93	130.72	125.82
2	B	501	HEM	CMD-C2D-C3D	2.99	127.57	114.35
5	A	504	MES	C2-C3-N4	3.05	114.75	110.12
5	B	503	MES	C7-C8-S	3.12	122.18	112.51
2	B	501	HEM	C2D-C3D-C4D	3.29	107.08	101.50
2	B	501	HEM	C1D-CHD-C4C	3.44	131.57	125.82
5	B	503	MES	C7-N4-C3	3.65	120.64	111.27
2	B	501	HEM	C3B-C4B-CHC	3.69	128.36	123.16
2	A	501	HEM	CAD-C3D-C4D	3.89	126.20	112.47
2	A	501	HEM	C2C-C1C-CHC	4.15	129.99	123.68
2	B	501	HEM	CAD-C3D-C2D	4.15	125.16	113.22
2	A	501	HEM	C3B-C4B-CHC	4.33	129.26	123.16
2	B	501	HEM	CAD-C3D-C4D	4.34	127.77	112.47
2	A	501	HEM	CMB-C2B-C3B	4.61	128.04	116.53
5	A	504	MES	C5-N4-C3	4.75	119.19	108.90
5	A	504	MES	C7-N4-C5	4.84	123.67	111.27
2	A	501	HEM	CAD-C3D-C2D	4.91	127.34	113.22

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	SYN	CAC-CAD-CAF	5.01	127.52	120.19
2	A	501	HEM	CMC-C2C-C3C	5.21	129.53	116.53
2	B	501	HEM	CMB-C2B-C3B	5.94	131.37	116.53
2	B	501	HEM	CMC-C2C-C3C	6.25	132.12	116.53
5	B	503	MES	C5-N4-C3	6.31	122.57	108.90
5	B	503	MES	O1S-S-C8	8.06	113.78	106.91

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	HEM	2	0
6	A	505	PEG	4	0
4	A	506	GOL	1	0
2	B	501	HEM	2	0
5	B	503	MES	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	455/455 (100%)	-0.13	16 (3%)	48	44	7, 13, 29, 107	0
1	B	455/455 (100%)	0.16	27 (5%)	26	23	7, 14, 51, 142	0
All	All	910/910 (100%)	0.02	43 (4%)	35	32	7, 14, 41, 142	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	191	ALA	9.6
1	A	1	THR	8.0
1	B	1	THR	7.5
1	B	2	ILE	6.1
1	B	193	PRO	5.2
1	B	198	TYR	5.1
1	B	192	ASN	5.1
1	B	196	PRO	4.9
1	B	194	ASP	4.6
1	A	2	ILE	4.6
1	B	199	ASP	4.6
1	B	136	ASP	4.5
1	B	3	LYS	4.4
1	A	191	ALA	4.3
1	B	4	GLU	4.1
1	B	195	ASP	4.1
1	B	197	ALA	4.0
1	B	225	ALA	3.7
1	B	201	ASN	3.7
1	B	229	GLN	3.4
1	B	265[A]	GLY	3.4
1	B	169	GLN	3.4
1	B	231	ASP	3.3
1	B	227	GLY	3.1

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Mol	Chain	Res	Type	RSRZ
1	A	4	GLU	3.0
1	A	225	ALA	2.9
1	A	3	LYS	2.8
1	B	158[A]	PHE	2.8
1	B	239	ASN	2.5
1	A	136	ASP	2.5
1	A	46	GLY	2.5
1	A	229	GLN	2.4
1	B	214	ASP	2.4
1	B	46	GLY	2.4
1	A	454	PRO	2.4
1	A	169	GLN	2.3
1	B	172	PRO	2.3
1	A	194	ASP	2.2
1	B	189	GLN	2.2
1	A	227	GLY	2.2
1	A	158[A]	PHE	2.2
1	A	192	ASN	2.1
1	A	13	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	GOL	A	503	6/6	0.94	0.15	11.79	21,34,45,47	0
5	MES	A	504	12/12	0.43	0.26	11.40	30,43,80,123	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
6	PEG	A	505	7/7	0.83	0.17	4.49	29,32,38,41	0
4	GOL	A	506	6/6	0.82	0.15	3.75	26,28,30,41	0
5	MES	B	503	12/12	0.72	0.15	1.67	34,45,71,80	0
3	SYN	B	502	8/8	0.87	0.12	0.98	26,31,41,43	0
3	SYN	A	502	8/8	0.84	0.10	0.34	23,26,29,30	0
2	HEM	B	501	43/43	0.98	0.07	-0.47	6,8,10,22	0
2	HEM	A	501	43/43	0.99	0.06	-0.61	5,7,10,17	0

6.5 Other polymers

There are no such residues in this entry.