



Full wwPDB X-ray Structure Validation Report

Feb 26, 2014 – 03:50 PM GMT

PDB ID : 3PM0
Title : Structural Characterization of the Complex between Alpha-Naphthoflavone and Human Cytochrome P450 1B1 (CYP1B1)
Authors : Wang, A.; Stout, C.D.; Johnson, E.F.
Deposited on : 2010-11-15
Resolution : 2.70 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

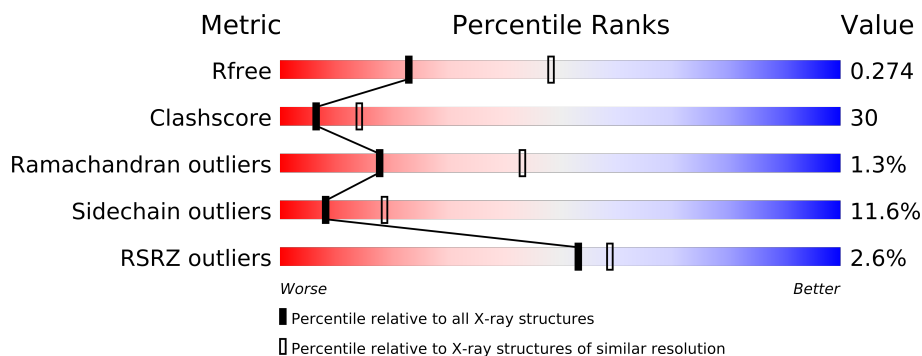
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.70 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	1557 (2.70-2.70)
Clashscore	79885	1939 (2.70-2.70)
Ramachandran outliers	78287	1905 (2.70-2.70)
Sidechain outliers	78261	1905 (2.70-2.70)
RSRZ outliers	66119	1559 (2.70-2.70)

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	507	

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	BHF	A	800	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3752 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 1B1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	459	3655	2325	657	653	20	0	0	0

There are 15 discrepancies between the modelled and reference sequences:

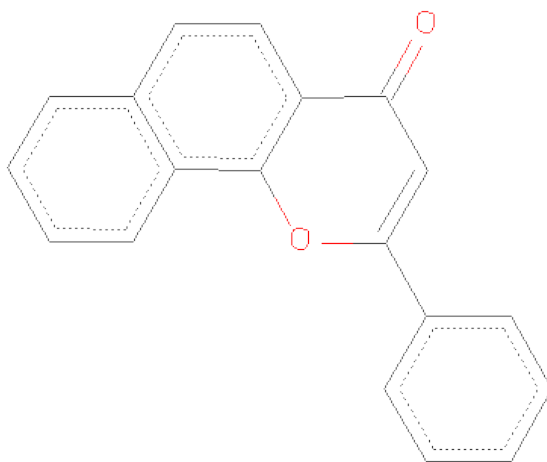
Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MET	-	EXPRESSION TAG	UNP Q16678
A	42	ALA	-	EXPRESSION TAG	UNP Q16678
A	43	LYS	-	EXPRESSION TAG	UNP Q16678
A	44	LYS	-	EXPRESSION TAG	UNP Q16678
A	45	THR	-	EXPRESSION TAG	UNP Q16678
A	46	SER	-	EXPRESSION TAG	UNP Q16678
A	47	SER	-	EXPRESSION TAG	UNP Q16678
A	48	LYS	-	EXPRESSION TAG	UNP Q16678
A	49	GLY	-	EXPRESSION TAG	UNP Q16678
A	50	LYS	-	EXPRESSION TAG	UNP Q16678
A	119	SER	ALA	SEE REMARK 999	UNP Q16678
A	544	HIS	-	EXPRESSION TAG	UNP Q16678
A	545	HIS	-	EXPRESSION TAG	UNP Q16678
A	546	HIS	-	EXPRESSION TAG	UNP Q16678
A	547	HIS	-	EXPRESSION TAG	UNP Q16678

- 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
			43	34	1	4	4	

- Molecule 3 is 2-PHENYL-4H-BENZO[H]CHROMEN-4-ONE (three-letter code: BHF) (formula: $C_{19}H_{12}O_2$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	C	O	0
			21	19	2	

- Molecule 4 is water.

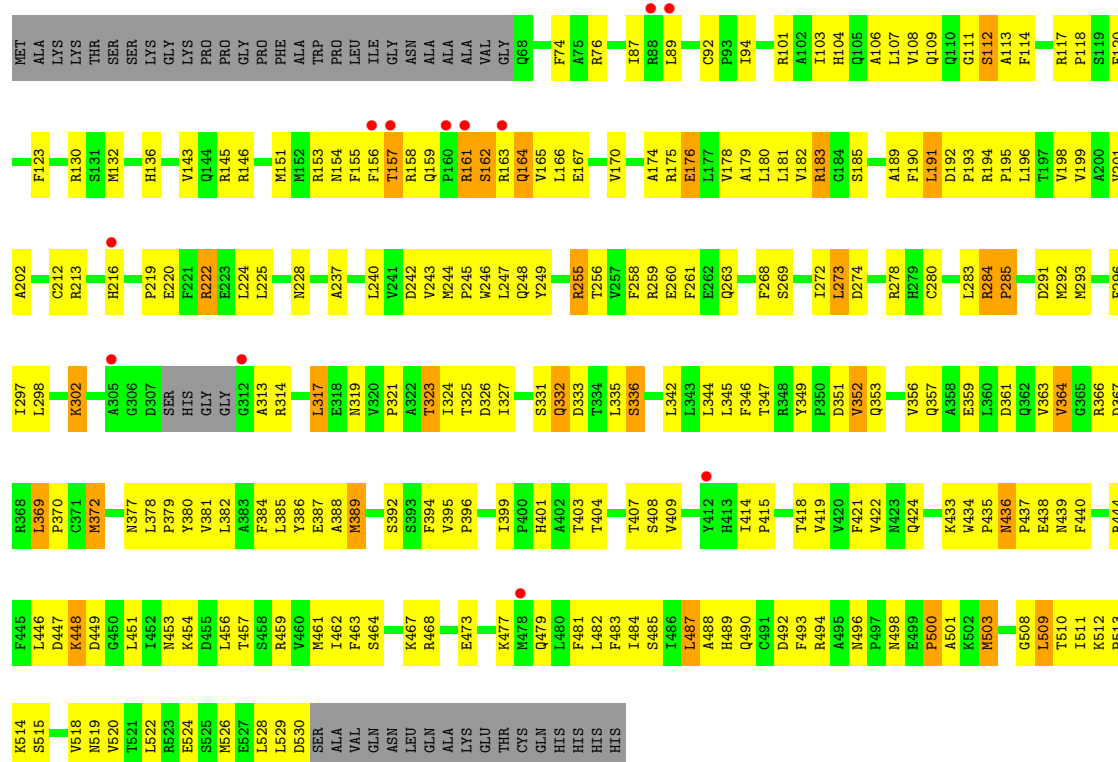
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	33	Total 33	O 33	0	0

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 1B1

Chain A: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	84.17Å 103.98Å 62.97Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	65.50 – 2.70 53.86 – 2.69	Depositor EDS
% Data completeness (in resolution range)	99.3 (65.50-2.70) 99.1 (53.86-2.69)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	0.11	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.64 (at 2.69Å)	Xtriage
Refinement program	CNS 1.21	Depositor
R, R_{free}	0.235 , 0.278 0.231 , 0.274	Depositor DCC
R_{free} test set	781 reflections (5.25%)	DCC
Wilson B-factor (Å ²)	44.0	Xtriage
Anisotropy	0.774	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 39.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 15708 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	3752	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.18% 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, BHF

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	0.50	0/3744	0.69	0/5072

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	494	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	3655	0	3614	221	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	43	0	30	9	0
3	A	21	0	12	1	0
4	A	33	0	0	0	0
All	All	3752	0	3656	221	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 30.

All (221) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:436:ASN:HB2	1:A:444:ARG:NH2	1.58	1.17
1:A:436:ASN:HB2	1:A:444:ARG:HH22	0.97	1.08
1:A:436:ASN:CB	1:A:444:ARG:HH22	1.69	1.06
1:A:453:ASN:ND2	1:A:456:LEU:HD23	1.73	1.02
1:A:454:LYS:HA	1:A:457:THR:HG22	1.36	1.02
1:A:292:MET:HE3	1:A:327:ILE:HD13	1.52	0.90
1:A:435:PRO:O	1:A:437:PRO:HD3	1.75	0.87
1:A:454:LYS:HA	1:A:457:THR:CG2	2.06	0.86
1:A:453:ASN:HD21	1:A:456:LEU:HD23	1.42	0.84
1:A:415:PRO:O	1:A:418:THR:HG22	1.78	0.83
1:A:157:THR:HG23	1:A:157:THR:O	1.79	0.82
1:A:155:PHE:HA	1:A:158:ARG:HD3	1.60	0.82
1:A:174:ALA:O	1:A:178:VAL:HG13	1.80	0.81
1:A:292:MET:HE1	1:A:327:ILE:HG21	1.65	0.79
1:A:162:SER:O	1:A:165:VAL:HG12	1.85	0.77
1:A:302:LYS:HA	1:A:302:LYS:HE3	1.67	0.77
1:A:395:VAL:HG11	2:A:900:HEM:HMA2	1.66	0.76
1:A:292:MET:CE	1:A:327:ILE:HG21	2.17	0.75
1:A:111:GLY:O	1:A:112:SER:CB	2.34	0.75
1:A:436:ASN:O	1:A:444:ARG:NH2	2.22	0.72
1:A:107:LEU:HD11	1:A:422:VAL:HG21	1.72	0.71
1:A:454:LYS:CA	1:A:457:THR:HG22	2.18	0.71
1:A:280:CYS:SG	1:A:297:ILE:HD13	2.32	0.70
1:A:438:GLU:OE2	1:A:438:GLU:HA	1.92	0.69
1:A:359:GLU:OE1	1:A:381:VAL:HG23	1.93	0.69
1:A:342:LEU:HD13	1:A:484:ILE:HD12	1.75	0.69
1:A:101:ARG:HH11	1:A:101:ARG:HG3	1.56	0.69
1:A:433:LYS:HD2	1:A:434:TRP:NE1	2.08	0.69
1:A:503:MET:HE1	1:A:514:LYS:O	1.93	0.69
1:A:388:ALA:O	1:A:392:SER:HB3	1.92	0.69
1:A:435:PRO:O	1:A:437:PRO:CD	2.42	0.68
1:A:111:GLY:O	1:A:112:SER:HB2	1.94	0.67

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:369:LEU:H	1:A:369:LEU:HD23	1.60	0.67
1:A:332:GLN:O	1:A:336:SER:HB2	1.94	0.67
1:A:284:ARG:CB	1:A:285:PRO:HD2	2.25	0.66
1:A:151:MET:O	1:A:155:PHE:HB2	1.96	0.66
1:A:490:GLN:NE2	1:A:528:LEU:HG	2.10	0.66
1:A:372:MET:CE	1:A:372:MET:HA	2.26	0.66
1:A:396:PRO:HD3	1:A:511:ILE:HG13	1.76	0.66
1:A:349:TYR:HB3	1:A:352:VAL:HG13	1.79	0.65
1:A:240:LEU:HD13	1:A:244:MET:HE1	1.77	0.65
1:A:434:TRP:HE3	1:A:444:ARG:HD3	1.61	0.65
1:A:372:MET:HE3	1:A:372:MET:HA	1.78	0.65
1:A:113:ALA:O	1:A:403:THR:HA	1.97	0.65
1:A:130:ARG:HD3	1:A:319:ASN:OD1	1.97	0.65
1:A:387:GLU:HG3	1:A:440:PHE:CE1	2.32	0.64
1:A:433:LYS:HD3	1:A:459:ARG:HG2	1.80	0.64
1:A:347:THR:HG21	1:A:518:VAL:HG23	1.77	0.64
1:A:319:ASN:O	1:A:323:THR:CG2	2.46	0.63
1:A:190:PHE:CE2	1:A:496:ASN:HA	2.33	0.63
1:A:378:LEU:HD12	1:A:481:PHE:HZ	1.64	0.63
1:A:434:TRP:CE3	1:A:444:ARG:HD3	2.33	0.63
1:A:319:ASN:O	1:A:323:THR:HG23	1.99	0.62
1:A:435:PRO:C	1:A:437:PRO:HD3	2.20	0.61
1:A:483:PHE:O	1:A:487:LEU:HD22	2.00	0.61
1:A:284:ARG:CB	1:A:285:PRO:CD	2.79	0.61
1:A:356:VAL:HG12	1:A:380:TYR:CE2	2.36	0.61
1:A:202:ALA:CB	1:A:225:LEU:HD11	2.31	0.60
1:A:222:ARG:HH11	1:A:222:ARG:CB	2.15	0.60
1:A:157:THR:CG2	1:A:157:THR:O	2.50	0.60
1:A:103:ILE:HG12	1:A:422:VAL:HG13	1.83	0.60
1:A:104:HIS:O	1:A:108:VAL:HB	2.02	0.60
1:A:272:ILE:HD13	1:A:324:ILE:HG23	1.83	0.59
1:A:120:PHE:HB3	1:A:242:ASP:CG	2.22	0.59
1:A:387:GLU:HG3	1:A:440:PHE:CD1	2.37	0.59
1:A:352:VAL:O	1:A:356:VAL:HG13	2.02	0.59
1:A:292:MET:HE3	1:A:327:ILE:CD1	2.29	0.59
1:A:436:ASN:H	1:A:444:ARG:HH12	1.51	0.58
1:A:369:LEU:CD2	1:A:369:LEU:H	2.16	0.58
1:A:284:ARG:HB3	1:A:285:PRO:HD2	1.85	0.58
1:A:145:ARG:NH1	2:A:900:HEM:O2D	2.36	0.58
1:A:114:PHE:CE1	1:A:407:THR:HG21	2.38	0.58
1:A:198:VAL:HA	1:A:336:SER:OG	2.04	0.58
1:A:154:ASN:O	1:A:158:ARG:HB3	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:331:SER:HB3	2:A:900:HEM:HMC2	1.86	0.57
1:A:378:LEU:N	1:A:379:PRO:HD3	2.20	0.56
1:A:240:LEU:HD13	1:A:244:MET:CE	2.36	0.56
1:A:372:MET:HE1	1:A:481:PHE:CD2	2.40	0.56
1:A:369:LEU:CD1	1:A:528:LEU:HD23	2.35	0.56
1:A:181:LEU:HD23	1:A:191:LEU:HD21	1.88	0.56
1:A:433:LYS:HD2	1:A:434:TRP:HE1	1.70	0.56
1:A:463:PHE:CE1	1:A:473:GLU:HG3	2.42	0.55
1:A:372:MET:CE	1:A:481:PHE:CD2	2.89	0.55
1:A:353:GLN:O	1:A:357:GLN:HB2	2.05	0.55
1:A:176:GLU:O	1:A:180:LEU:HG	2.05	0.55
1:A:156:PHE:O	1:A:157:THR:HG22	2.07	0.55
1:A:74:PHE:CE1	1:A:87:ILE:HD13	2.41	0.55
1:A:74:PHE:HE1	1:A:87:ILE:HD13	1.72	0.55
1:A:344:LEU:HD11	1:A:503:MET:HG3	1.89	0.55
1:A:498:ASN:O	1:A:500:PRO:HD3	2.07	0.54
1:A:146:ARG:HD2	1:A:313:ALA:HB2	1.89	0.54
1:A:361:ASP:OD1	1:A:366:ARG:NH2	2.41	0.54
1:A:189:ALA:O	1:A:519:ASN:ND2	2.31	0.54
1:A:94:ILE:HG23	1:A:419:VAL:HG12	1.91	0.53
1:A:342:LEU:HD22	1:A:384:PHE:CE2	2.44	0.53
1:A:248:GLN:HA	1:A:255:ARG:HG2	1.91	0.52
1:A:503:MET:HE1	1:A:515:SER:O	2.09	0.52
1:A:356:VAL:HG12	1:A:380:TYR:HE2	1.74	0.52
1:A:361:ASP:CG	1:A:366:ARG:NH2	2.62	0.52
1:A:326:ASP:HA	3:A:800:BHF:O2	2.10	0.51
1:A:156:PHE:O	1:A:157:THR:CB	2.58	0.51
1:A:219:PRO:HG2	1:A:220:GLU:H	1.75	0.51
1:A:493:PHE:CD2	1:A:520:VAL:HG22	2.45	0.51
1:A:389:MET:HG3	1:A:462:ILE:HD13	1.92	0.51
1:A:436:ASN:CB	1:A:444:ARG:NH2	2.45	0.51
1:A:369:LEU:HD12	1:A:528:LEU:HD23	1.91	0.51
1:A:151:MET:O	1:A:155:PHE:CB	2.58	0.51
1:A:366:ARG:HD2	1:A:524:GLU:OE2	2.11	0.51
1:A:161:ARG:HG3	1:A:164:GLN:HG3	1.93	0.51
1:A:143:VAL:HG13	1:A:313:ALA:HB1	1.93	0.51
1:A:228:ASN:HB3	1:A:333:ASP:OD1	2.11	0.51
1:A:179:ALA:O	1:A:183:ARG:HB2	2.12	0.50
1:A:255:ARG:CZ	1:A:259:ARG:NH1	2.74	0.50
1:A:201:VAL:HG11	1:A:335:LEU:HB2	1.92	0.50
1:A:132:MET:SD	1:A:145:ARG:HG3	2.52	0.50
1:A:274:ASP:OD1	1:A:278:ARG:NH1	2.46	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:395:VAL:HG11	2:A:900:HEM:CMA	2.38	0.49
1:A:447:ASP:CB	1:A:451:LEU:O	2.60	0.49
1:A:292:MET:HE2	1:A:296:PHE:HE2	1.78	0.49
1:A:394:PHE:O	1:A:510:THR:HA	2.13	0.48
1:A:284:ARG:HB3	1:A:285:PRO:CD	2.43	0.48
1:A:159:GLN:O	1:A:162:SER:HB2	2.14	0.48
1:A:246:TRP:HA	1:A:249:TYR:CE2	2.48	0.48
1:A:104:HIS:CE1	1:A:109:GLN:HG3	2.48	0.48
1:A:436:ASN:CA	1:A:444:ARG:HH22	2.26	0.48
1:A:327:ILE:HG23	2:A:900:HEM:HBC1	1.95	0.48
1:A:453:ASN:HD22	1:A:456:LEU:HD23	1.71	0.48
1:A:361:ASP:CG	1:A:366:ARG:HH21	2.18	0.47
1:A:130:ARG:CD	1:A:319:ASN:OD1	2.63	0.47
1:A:225:LEU:HD22	1:A:332:GLN:HG3	1.97	0.47
1:A:272:ILE:CD1	1:A:325:THR:HA	2.45	0.47
1:A:436:ASN:HB3	1:A:439:ASN:HB2	1.98	0.46
1:A:468:ARG:O	2:A:900:HEM:HBA2	2.15	0.46
1:A:387:GLU:HA	1:A:387:GLU:OE1	2.14	0.46
1:A:382:LEU:O	1:A:386:TYR:HD2	1.98	0.46
1:A:167:GLU:HG3	1:A:482:LEU:CD1	2.45	0.46
1:A:178:VAL:HG12	1:A:487:LEU:HD12	1.98	0.46
1:A:117:ARG:HG2	1:A:468:ARG:NH2	2.31	0.46
1:A:258:PHE:O	1:A:261:PHE:HB3	2.16	0.46
1:A:415:PRO:O	1:A:418:THR:CG2	2.57	0.46
1:A:199:VAL:HG11	1:A:216:HIS:HA	1.97	0.46
1:A:170:VAL:HG11	1:A:482:LEU:HB3	1.97	0.45
1:A:244:MET:CE	1:A:247:LEU:HD11	2.46	0.45
1:A:244:MET:HE3	1:A:247:LEU:HD11	1.98	0.45
1:A:463:PHE:CD1	1:A:473:GLU:HG3	2.51	0.45
1:A:194:ARG:HB2	1:A:195:PRO:HD3	1.97	0.45
1:A:156:PHE:O	1:A:157:THR:HB	2.16	0.45
1:A:132:MET:N	1:A:326:ASP:OD1	2.45	0.45
1:A:448:LYS:HE3	1:A:448:LYS:H	1.80	0.45
1:A:372:MET:HE3	1:A:481:PHE:CD2	2.51	0.45
1:A:94:ILE:HG21	1:A:421:PHE:CE1	2.52	0.45
1:A:384:PHE:CE2	1:A:484:ILE:CD1	3.00	0.45
1:A:111:GLY:O	1:A:112:SER:HB3	2.16	0.44
1:A:342:LEU:HD22	1:A:384:PHE:HE2	1.82	0.44
1:A:292:MET:HE3	1:A:327:ILE:HG21	1.95	0.44
1:A:490:GLN:HE21	1:A:528:LEU:HG	1.78	0.44
1:A:317:LEU:HA	1:A:317:LEU:HD23	1.77	0.44
1:A:272:ILE:HG21	1:A:324:ILE:CG2	2.48	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:364:VAL:HG11	1:A:370:PRO:HG3	2.00	0.44
1:A:101:ARG:HH11	1:A:101:ARG:CG	2.24	0.44
1:A:512:LYS:HA	1:A:513:PRO:HD3	1.90	0.44
1:A:453:ASN:HD21	1:A:456:LEU:CD2	2.22	0.44
1:A:190:PHE:CZ	1:A:496:ASN:HA	2.53	0.44
1:A:146:ARG:HD2	1:A:313:ALA:CB	2.48	0.44
1:A:89:LEU:HB3	1:A:243:VAL:HG21	1.99	0.44
1:A:446:LEU:HA	1:A:446:LEU:HD23	1.78	0.44
1:A:117:ARG:O	1:A:136:HIS:CE1	2.71	0.44
1:A:185:SER:OG	1:A:520:VAL:N	2.48	0.44
1:A:255:ARG:NE	1:A:259:ARG:CZ	2.81	0.44
1:A:485:SER:O	1:A:489:HIS:HB3	2.18	0.43
1:A:167:GLU:HG3	1:A:482:LEU:HD11	2.00	0.43
1:A:321:PRO:O	1:A:324:ILE:HG22	2.18	0.43
1:A:242:ASP:O	1:A:245:PRO:HD3	2.18	0.43
1:A:237:ALA:HB1	1:A:421:PHE:CD2	2.54	0.43
1:A:364:VAL:HG21	1:A:370:PRO:HG3	2.00	0.43
1:A:175:ARG:HE	1:A:529:LEU:CD2	2.31	0.43
1:A:477:LYS:C	1:A:479:GLN:H	2.22	0.43
1:A:396:PRO:HD2	1:A:509:LEU:O	2.18	0.43
1:A:363:VAL:HG21	1:A:377:ASN:O	2.19	0.43
1:A:291:ASP:HB3	1:A:293:MET:H	1.83	0.43
1:A:357:GLN:HG3	1:A:488:ALA:O	2.18	0.43
1:A:509:LEU:HA	1:A:509:LEU:HD12	1.82	0.42
1:A:384:PHE:CE2	1:A:484:ILE:HD11	2.54	0.42
1:A:424:GLN:OE1	1:A:461:MET:HG2	2.20	0.42
1:A:369:LEU:CD2	1:A:369:LEU:N	2.81	0.42
1:A:377:ASN:C	1:A:379:PRO:HD3	2.40	0.42
1:A:319:ASN:O	1:A:323:THR:HG22	2.18	0.42
1:A:202:ALA:HB1	1:A:225:LEU:HD11	2.02	0.42
1:A:225:LEU:HD22	1:A:332:GLN:CG	2.50	0.42
1:A:331:SER:CB	2:A:900:HEM:HMC2	2.49	0.42
1:A:156:PHE:O	1:A:157:THR:CG2	2.68	0.42
1:A:181:LEU:O	1:A:185:SER:HB2	2.20	0.42
1:A:118:PRO:HD2	1:A:123:PHE:HE1	1.85	0.42
1:A:414:ILE:HG22	1:A:418:THR:HG21	2.02	0.41
1:A:191:LEU:O	1:A:191:LEU:HD13	2.19	0.41
1:A:366:ARG:HA	1:A:366:ARG:NE	2.34	0.41
1:A:433:LYS:NZ	1:A:433:LYS:CB	2.84	0.41
1:A:237:ALA:HB1	1:A:421:PHE:CE2	2.54	0.41
1:A:192:ASP:C	1:A:192:ASP:OD1	2.59	0.41
1:A:182:VAL:HG21	1:A:522:LEU:HG	2.00	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:526:MET:CE	1:A:528:LEU:HD12	2.50	0.41
1:A:101:ARG:CG	1:A:101:ARG:NH1	2.83	0.41
1:A:222:ARG:HH11	1:A:222:ARG:HB2	1.84	0.41
1:A:256:THR:O	1:A:260:GLU:HG3	2.20	0.41
1:A:500:PRO:HG2	1:A:501:ALA:H	1.85	0.41
1:A:399:ILE:O	1:A:401:HIS:HD2	2.03	0.41
1:A:193:PRO:O	1:A:196:LEU:HB2	2.21	0.41
1:A:404:THR:O	1:A:404:THR:HG22	2.21	0.41
1:A:335:LEU:HD23	2:A:900:HEM:HAB	2.03	0.41
1:A:435:PRO:C	1:A:437:PRO:CD	2.88	0.41
1:A:384:PHE:HE2	1:A:484:ILE:CD1	2.34	0.41
1:A:202:ALA:CB	1:A:225:LEU:CD1	2.97	0.41
1:A:199:VAL:HG21	1:A:222:ARG:HD3	2.02	0.41
1:A:357:GLN:O	1:A:361:ASP:OD1	2.39	0.41
1:A:118:PRO:HG3	1:A:401:HIS:O	2.21	0.41
1:A:106:ALA:HB2	1:A:409:VAL:HG21	2.02	0.40
1:A:154:ASN:O	1:A:155:PHE:C	2.59	0.40
1:A:464:SER:CB	2:A:900:HEM:HBA1	2.50	0.40
1:A:274:ASP:OD1	1:A:278:ARG:NH2	2.53	0.40
1:A:269:SER:O	1:A:273:LEU:HB2	2.21	0.40
1:A:407:THR:OG1	1:A:408:SER:N	2.55	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	455/507 (90%)	418 (92%)	31 (7%)	6 (1%)	18	43

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	436	ASN
1	A	112	SER

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Mol	Chain	Res	Type
1	A	285	PRO
1	A	500	PRO
1	A	509	LEU
1	A	508	GLY

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	395/430 (92%)	349 (88%)	46 (12%)	8 18

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ARG
1	A	92	CYS
1	A	153	ARG
1	A	157	THR
1	A	161	ARG
1	A	162	SER
1	A	163	ARG
1	A	164	GLN
1	A	166	LEU
1	A	176	GLU
1	A	183	ARG
1	A	191	LEU
1	A	212	CYS
1	A	213	ARG
1	A	222	ARG
1	A	224	LEU
1	A	255	ARG
1	A	263	GLN
1	A	268	PHE
1	A	273	LEU
1	A	283	LEU
1	A	284	ARG
1	A	298	LEU
1	A	302	LYS

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Mol	Chain	Res	Type
1	A	314	ARG
1	A	317	LEU
1	A	323	THR
1	A	332	GLN
1	A	336	SER
1	A	345	LEU
1	A	346	PHE
1	A	351	ASP
1	A	352	VAL
1	A	364	VAL
1	A	367	ASP
1	A	369	LEU
1	A	372	MET
1	A	385	LEU
1	A	389	MET
1	A	448	LYS
1	A	449	ASP
1	A	467	LYS
1	A	487	LEU
1	A	492	ASP
1	A	503	MET
1	A	530	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	429	HIS
1	A	490	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 ⓘ

2 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$
3	BHF	A	800	-	24,24,24	1.42	1 (4%)	32,34,34	1.59	6 (18%)
2	HEM	A	900	1	49,50,50	4.11	22 (44%)	46,82,82	1.96	13 (28%)

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
3	BHF	A	800	-	-	0/4/4/4	0/1/4/4
2	HEM	A	900	1	-	0/14/114/114	0/0/8/8

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	HEM	C3D-C4D	-21.25	1.39	1.44
2	A	900	HEM	C2D-C1D	-12.65	1.41	1.44
2	A	900	HEM	C2B-C1B	-6.40	1.43	1.44
3	A	800	BHF	C1-C7	-5.25	1.39	1.46
2	A	900	HEM	C3D-C2D	-3.65	1.37	1.43
2	A	900	HEM	C1B-NB	-3.35	1.32	1.39
2	A	900	HEM	CBB-CAB	3.31	1.48	1.28
2	A	900	HEM	FE-NA	3.29	2.06	1.92
2	A	900	HEM	C4D-ND	-3.27	1.32	1.39
2	A	900	HEM	C1C-NC	-3.18	1.33	1.38
2	A	900	HEM	FE-NB	2.84	2.08	1.97
2	A	900	HEM	C4A-NA	-2.82	1.31	1.36
2	A	900	HEM	CMA-C3A	2.81	1.57	1.51
2	A	900	HEM	C1D-ND	-2.81	1.31	1.37
2	A	900	HEM	CMC-C2C	2.80	1.56	1.47
2	A	900	HEM	CHA-C4D	2.63	1.39	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	HEM	CMD-C2D	2.35	1.54	1.47
2	A	900	HEM	C4B-NB	-2.34	1.32	1.37
2	A	900	HEM	CHD-C4C	2.16	1.40	1.36
2	A	900	HEM	FE-NC	2.13	2.05	1.97
2	A	900	HEM	CBC-CAC	2.11	1.41	1.28
2	A	900	HEM	C4C-NC	-2.07	1.35	1.38
2	A	900	HEM	O2D-CGD	-2.05	1.23	1.30

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	900	HEM	CHD-C4C-NC	-4.77	120.59	124.73
2	A	900	HEM	C3B-C4B-NB	-4.18	111.01	114.00
3	A	800	BHF	C8-C7-C1	-4.14	122.76	127.13
3	A	800	BHF	O1-C7-C1	3.65	118.23	113.54
2	A	900	HEM	C4D-ND-C1D	3.65	108.90	105.16
2	A	900	HEM	C3A-C4A-NA	3.59	112.12	109.41
2	A	900	HEM	CHC-C1C-NC	-3.54	121.66	124.73
2	A	900	HEM	C2D-C1D-ND	-3.53	108.77	112.93
3	A	800	BHF	O1-C19-C18	3.13	119.86	116.13
2	A	900	HEM	CHC-C4B-NB	-3.04	122.06	124.58
3	A	800	BHF	C11-C10-C19	3.02	119.99	116.30
2	A	900	HEM	CBA-CAA-C2A	2.67	117.40	112.69
3	A	800	BHF	C8-C9-C10	-2.54	120.21	123.28
2	A	900	HEM	C4A-CHB-C1B	2.54	130.81	127.47
2	A	900	HEM	CHA-C1A-NA	-2.48	120.44	124.58
2	A	900	HEM	C1B-NB-C4B	2.37	107.58	105.16
3	A	800	BHF	C17-C18-C19	-2.13	119.99	123.78
2	A	900	HEM	O2A-CGA-CBA	2.12	121.72	114.22
2	A	900	HEM	C4B-CHC-C1C	2.09	132.07	126.57

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	459/507 (90%)	0.18	12 (2%) 53 59	27, 43, 68, 100	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	312	GLY	3.9
1	A	163	ARG	3.7
1	A	412	TYR	2.9
1	A	157	THR	2.8
1	A	161	ARG	2.5
1	A	89	LEU	2.4
1	A	216	HIS	2.3
1	A	160	PRO	2.2
1	A	478	MET	2.1
1	A	156	PHE	2.1
1	A	305	ALA	2.1
1	A	88	ARG	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
3	BHF	A	800	21/21	0.24	2.03	38,40,42,43	0
2	HEM	A	900	43/43	0.15	-0.59	22,27,37,41	0

6.5 Other polymers ⓘ

There are no such residues in this entry.