



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 13, 2017 – 11:26 am 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 X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.7.2 (RC1), CSD as538be (2017)  
Xtriage (Phenix) : 1.9-1692  
EDS : trunk28620  
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)  
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) : recalc28949

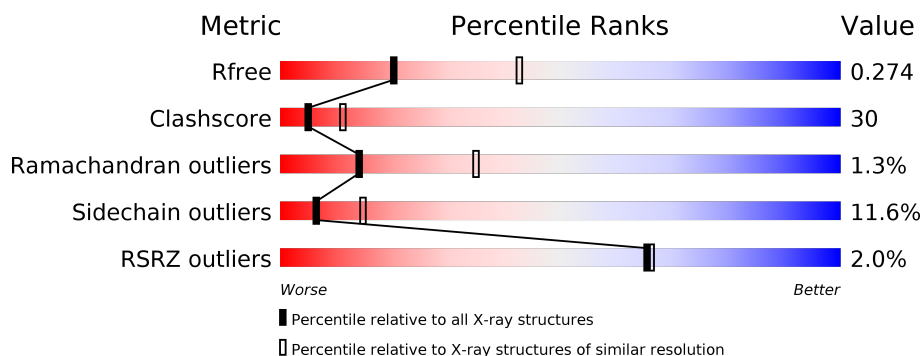
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 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))
$R_{free}$	100719	2259 (2.70-2.70)
Clashscore	112137	2590 (2.70-2.70)
Ramachandran outliers	110173	2550 (2.70-2.70)
Sidechain outliers	110143	2550 (2.70-2.70)
RSRZ outliers	101464	2275 (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  $\geq 3$ , 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	507	

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
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 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 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:

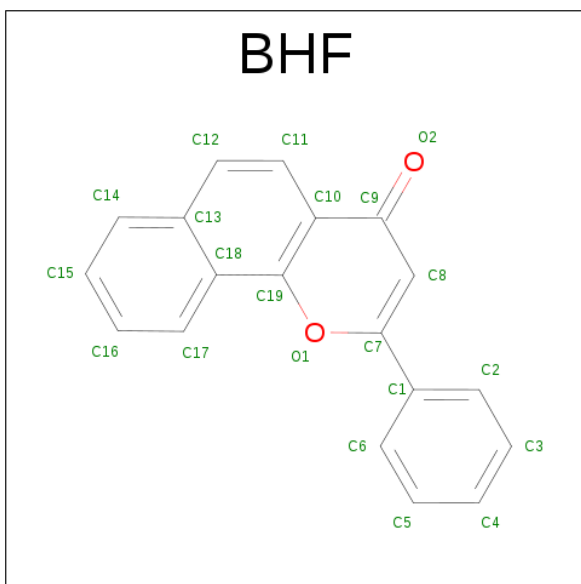
Chain	Residue	Modelled	Actual	Comment	Reference
A	41	MET	-	EXPRESSION TAG	UNP Q16678
A	42	ALA	-	EXPRESSION TAG	UNP Q16678
A	43	LYS	-	EXPRESSION TAG	UNP Q16678
A	44	LYS	-	EXPRESSION TAG	UNP Q16678
A	45	THR	-	EXPRESSION TAG	UNP Q16678
A	46	SER	-	EXPRESSION TAG	UNP Q16678
A	47	SER	-	EXPRESSION TAG	UNP Q16678
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	
			43	34	1	4	4	
							0	0

- 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	
			21	19	2	
					0	0

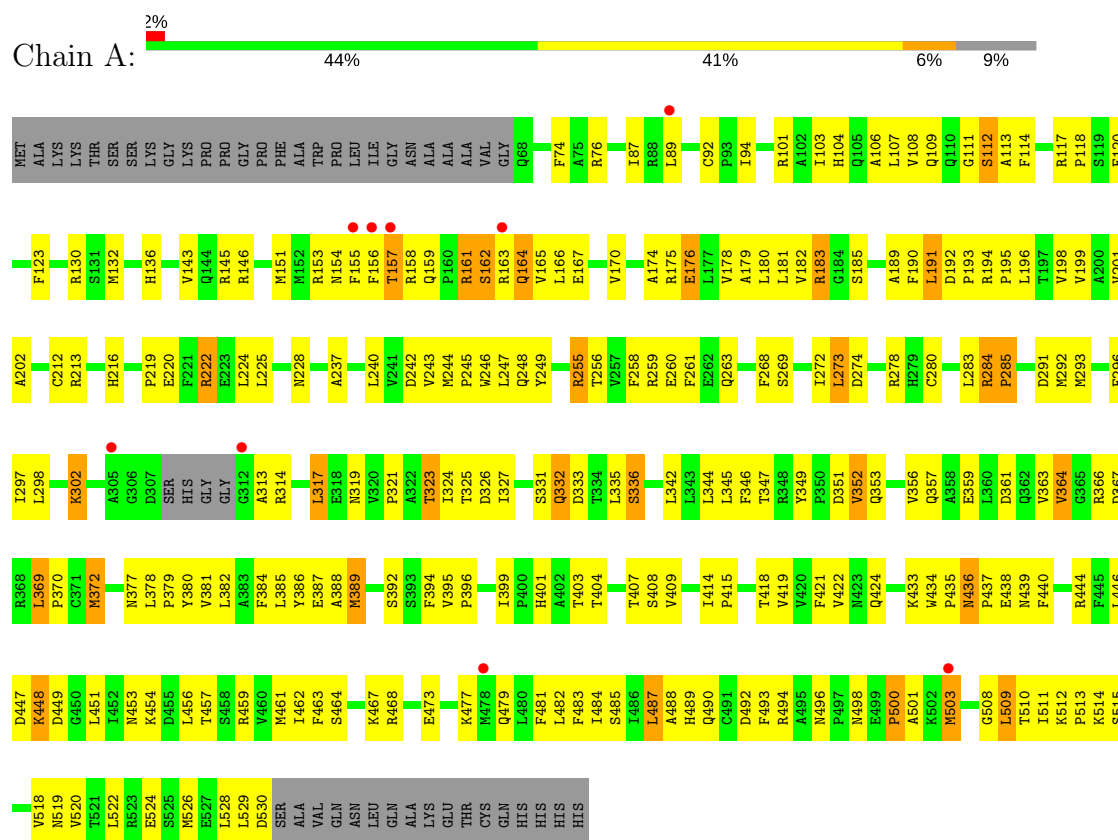
- Molecule 4 is water.

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 the various outlier classes 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



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	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
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.64 (at 2.69Å)	Xtriage
Refinement program	CNS 1.21	Depositor
R, $R_{free}$	0.235 , 0.278 0.230 , 0.274	Depositor DCC
$R_{free}$ test set	781 reflections (5.25%)	DCC
Wilson B-factor (Å <sup>2</sup> )	44.0	Xtriage
Anisotropy	0.774	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 49.6	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.92	EDS
Total number of atoms	3752	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	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.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: 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 Too-close contacts [i](#)

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	3655	0	3614	221	0
2	A	43	0	30	9	0

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

All (221) 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: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:101:ARG:HH11	1:A:101:ARG:HG3	1.56	0.69
1:A:342:LEU:HD13	1:A:484:ILE:HD12	1.75	0.69
1:A:359:GLU:OE1	1:A:381:VAL:HG23	1.93	0.69
1:A:433:LYS:HD2	1:A:434:TRP:NE1	2.08	0.69
1:A:388:ALA:O	1:A:392:SER:HB3	1.92	0.69
1:A:503:MET:HE1	1:A:514:LYS:O	1.93	0.69
1:A:435:PRO:O	1:A:437:PRO:CD	2.42	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:111:GLY:O	1:A:112:SER:HB2	1.94	0.67
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:130:ARG:HD3	1:A:319:ASN:OD1	1.97	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:434:TRP:HE3	1:A:444:ARG:HD3	1.61	0.65
1:A:387:GLU:HG3	1:A:440:PHE:CE1	2.32	0.64
1:A:347:THR:HG21	1:A:518:VAL:HG23	1.77	0.64
1:A:433:LYS:HD3	1:A:459:ARG:HG2	1.80	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:284:ARG:CB	1:A:285:PRO:CD	2.79	0.61
1:A:483:PHE:O	1:A:487:LEU:HD22	2.00	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:120:PHE:HB3	1:A:242:ASP:CG	2.22	0.59
1:A:272:ILE:HD13	1:A:324:ILE:HG23	1.83	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:284:ARG:HB3	1:A:285:PRO:HD2	1.85	0.58
1:A:369:LEU:CD2	1:A:369:LEU:H	2.16	0.58
1:A:436:ASN:H	1:A:444:ARG:HH12	1.51	0.58
1:A:114:PHE:CE1	1:A:407:THR:HG21	2.38	0.58
1:A:145:ARG:NH1	2:A:900:HEM:O2D	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:176:GLU:O	1:A:180:LEU:HG	2.05	0.55
1:A:353:GLN:O	1:A:357:GLN:HB2	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:189:ALA:O	1:A:519:ASN:ND2	2.31	0.54
1:A:361:ASP:OD1	1:A:366:ARG:NH2	2.41	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:156:PHE:O	1:A:157:THR:CB	2.58	0.51
1:A:326:ASP:HA	3:A:800:BHF:O2	2.10	0.51
1:A:219:PRO:HG2	1:A:220:GLU:H	1.75	0.51
1:A:389:MET:HG3	1:A:462:ILE:HD13	1.92	0.51
1:A:493:PHE:CD2	1:A:520:VAL:HG22	2.45	0.51
1:A:369:LEU:HD12	1:A:528:LEU:HD23	1.91	0.51
1:A:436:ASN:CB	1:A:444:ARG:NH2	2.45	0.51
1:A:151:MET:O	1:A:155:PHE:CB	2.58	0.51
1:A:161:ARG:HG3	1:A:164:GLN:HG3	1.93	0.51
1:A:366:ARG:HD2	1:A:524:GLU:OE2	2.11	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:327:ILE:HG23	2:A:900:HEM:HBC1	1.95	0.48
1:A:436:ASN:CA	1:A:444:ARG:HH22	2.26	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:272:ILE:CD1	1:A:325:THR:HA	2.45	0.47
1:A:225:LEU:HD22	1:A:332:GLN:HG3	1.97	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:258:PHE:O	1:A:261:PHE:HB3	2.16	0.46
1:A:117:ARG:HG2	1:A:468:ARG:NH2	2.31	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:194:ARG:HB2	1:A:195:PRO:HD3	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:292:MET:HE3	1:A:327:ILE:HG21	1.95	0.44
1:A:317:LEU:HA	1:A:317:LEU:HD23	1.77	0.44
1:A:490:GLN:HE21	1:A:528:LEU:HG	1.78	0.44
1:A:272:ILE:HG21	1:A:324:ILE:CG2	2.48	0.44
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:146:ARG:HD2	1:A:313:ALA:CB	2.48	0.44
1:A:446:LEU:HA	1:A:446:LEU:HD23	1.78	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:89:LEU:HB3	1:A:243:VAL:HG21	1.99	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:242:ASP:O	1:A:245:PRO:HD3	2.18	0.43
1:A:321:PRO:O	1:A:324:ILE:HG22	2.18	0.43
1:A:364:VAL:HG21	1:A:370:PRO:HG3	2.00	0.43
1:A:237:ALA:HB1	1:A:421:PHE:CD2	2.54	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:291:ASP:HB3	1:A:293:MET:H	1.83	0.43
1:A:363:VAL:HG21	1:A:377:ASN:O	2.19	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:319:ASN:O	1:A:323:THR:HG22	2.18	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: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:156:PHE:O	1:A:157:THR:CG2	2.68	0.42
1:A:331:SER:CB	2:A:900:HEM:HMC2	2.49	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: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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:414:ILE:HG22	1:A:418:THR:HG21	2.02	0.41
1:A:433:LYS:CB	1:A:433:LYS:NZ	2.84	0.41
1:A:192:ASP:C	1:A:192:ASP:OD1	2.59	0.41
1:A:237:ALA:HB1	1:A:421:PHE:CE2	2.54	0.41
1:A:182:VAL:HG21	1:A:522:LEU:HG	2.00	0.41
1:A:526:MET:CE	1:A:528:LEU:HD12	2.50	0.41
1:A:101:ARG:NH1	1:A:101:ARG:CG	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:193:PRO:O	1:A:196:LEU:HB2	2.21	0.41
1:A:399:ILE:O	1:A:401:HIS:HD2	2.03	0.41
1:A:404:THR:O	1:A:404:THR:HG22	2.21	0.41
1:A:500:PRO:HG2	1:A:501:ALA:H	1.85	0.41
1:A:199:VAL:HG21	1:A:222:ARG:HD3	2.02	0.41
1:A:202:ALA:CB	1:A:225:LEU:CD1	2.97	0.41
1:A:335:LEU:HD23	2:A:900:HEM:HAB	2.03	0.41
1:A:118:PRO:HG3	1:A:401:HIS:O	2.21	0.41
1:A:357:GLN:O	1:A:361:ASP:OD1	2.39	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: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:269:SER:O	1:A:273:LEU:HB2	2.21	0.40
1:A:274:ASP:OD1	1:A:278:ARG:NH2	2.53	0.40
1:A:464:SER:CB	2:A:900:HEM:HBA1	2.50	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 [i](#)

### 5.3.1 Protein backbone [i](#)

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%)	14	35

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	436	ASN
1	A	112	SER
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%)	6	15

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

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Mol	Chain	Res	Type
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
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 molecules in this entry.



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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

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	1.19	1 (4%)	26,34,34	1.07	1 (3%)
2	HEM	A	900	1	28,50,50	2.10	10 (35%)	17,82,82	1.54	4 (23%)

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	0/4/4/4
2	HEM	A	900	1	-	0/6/54/54	0/0/8/8

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	HEM	C3C-CAC	-5.93	1.35	1.47
3	A	800	BHF	C1-C7	-3.80	1.39	1.46
2	A	900	HEM	C1B-NB	-3.49	1.32	1.36
2	A	900	HEM	C4D-ND	-3.36	1.32	1.36
2	A	900	HEM	C3B-CAB	-2.95	1.42	1.47
2	A	900	HEM	C1C-NC	-2.75	1.33	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	900	HEM	C1D-ND	-2.31	1.31	1.36
2	A	900	HEM	C4A-NA	-2.29	1.31	1.36
2	A	900	HEM	CMC-C2C	2.12	1.56	1.51
2	A	900	HEM	CBB-CAB	2.73	1.48	1.28
2	A	900	HEM	CMA-C3A	2.90	1.57	1.51

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	800	BHF	C8-C9-C10	-2.57	120.21	123.05
2	A	900	HEM	C4C-C3C-C2C	-2.54	105.13	106.90
2	A	900	HEM	CAD-CBD-CGD	2.15	116.34	112.66
2	A	900	HEM	CBA-CAA-C2A	2.57	117.40	112.48
2	A	900	HEM	CMB-C2B-C3B	2.81	130.11	124.89

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	800	BHF	1	0
2	A	900	HEM	9	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 [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	459/507 (90%)	0.13	9 (1%) 65 66	27, 43, 68, 100	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	157	THR	3.8
1	A	163	ARG	3.5
1	A	312	GLY	3.2
1	A	305	ALA	3.0
1	A	89	LEU	2.2
1	A	155	PHE	2.1
1	A	156	PHE	2.1
1	A	503	MET	2.1
1	A	478	MET	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, 95<sup>th</sup> 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( $\text{\AA}^2$ )	Q<0.9
3	BHF	A	800	21/21	0.89	0.24	2.57	38,40,42,43	0
2	HEM	A	900	43/43	0.97	0.16	-0.45	22,27,37,41	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.