



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:07 AM GMT

PDB ID : 2FRZ
Title : Crystal Structure of Cytochrome P450cam mutant (F87W/Y96F/V247L/C334A)
Authors : Rao, Z.; Wong, L.L.; Xu, F.; Bell, S.G.
Deposited on : 2006-01-20
Resolution : 2.10 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.

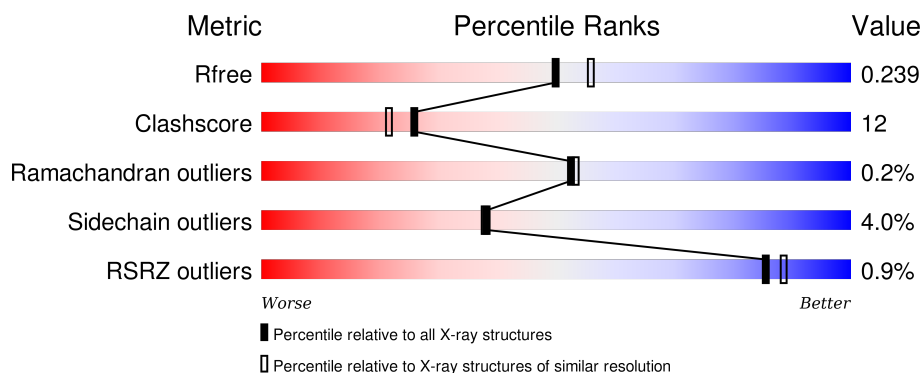
The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

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X-RAY DIFFRACTION

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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	414	<div> <div></div> <div>72%</div> <div>23%</div> <div>..</div> </div>
1	B	414	<div> <div></div> <div>73%</div> <div>22%</div> <div>..</div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7027 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-cam.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	405	Total	C	N	O	S	0	0	0
			3209	2036	561	595	17			
1	B	405	Total	C	N	O	S	0	0	0
			3209	2036	561	595	17			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	87	TRP	PHE	ENGINEERED	UNP P00183
A	96	PHE	TYR	ENGINEERED	UNP P00183
A	247	LEU	VAL	ENGINEERED	UNP P00183
A	334	ALA	CYS	ENGINEERED	UNP P00183
B	87	TRP	PHE	ENGINEERED	UNP P00183
B	96	PHE	TYR	ENGINEERED	UNP P00183
B	247	LEU	VAL	ENGINEERED	UNP P00183
B	334	ALA	CYS	ENGINEERED	UNP P00183

- Molecule 2 is POTASSIUM ION (three-letter code: K) (formula: K).

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

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

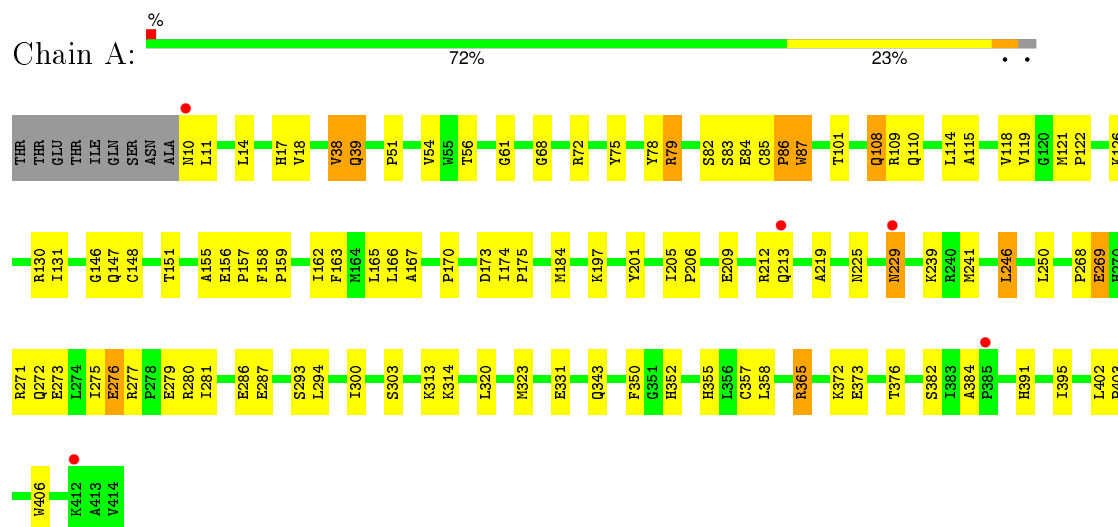
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	268	Total	O	0	0
			268	268		
4	B	253	Total	O	0	0
			253	253		

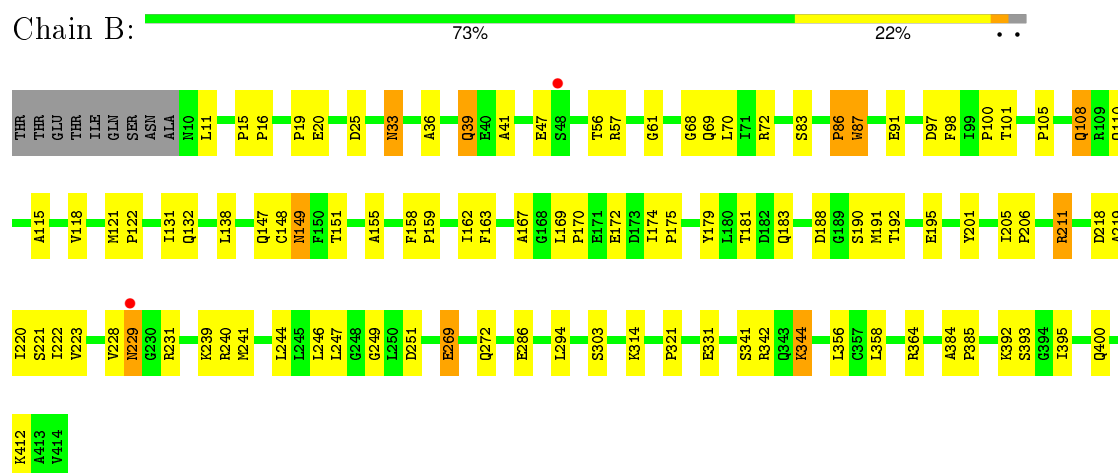
3 Residue-property plots

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-cam



• Molecule 1: Cytochrome P450-cam



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.82Å 62.11Å 94.94Å 90.00° 90.46° 90.00°	Depositor
Resolution (Å)	50.00 – 2.10 45.49 – 2.09	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.10) 97.2 (45.49-2.09)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.18 (at 2.10Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.190 , 0.243 0.189 , 0.239	Depositor DCC
R_{free} test set	4602 reflections (11.48%)	DCC
Wilson B-factor (Å ²)	21.2	Xtriage
Anisotropy	0.212	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 55.8	EDS
Estimated twinning fraction	0.028 for h,-k,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 45964 reflections	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	7027	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.78% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, K

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	0.32	0/3289	0.57	0/4470
1	B	0.32	0/3289	0.56	0/4470
All	All	0.32	0/6578	0.57	0/8940

There are no bond length outliers.

There are no bond angle outliers.

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	3209	0	3159	82	0
1	B	3209	0	3159	78	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	43	0	30	2	0
3	B	43	0	30	2	0
4	A	268	0	0	4	0
4	B	253	0	0	4	0
All	All	7027	0	6378	160	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 12.

All (160) 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:110:GLN:HB3	1:B:228:VAL:HG23	1.55	0.87
1:B:344:LYS:HA	1:B:344:LYS:HE3	1.53	0.87
1:B:223:VAL:HG22	1:B:241:MET:HE2	1.59	0.85
1:A:38:VAL:HG22	1:A:391:HIS:HB3	1.60	0.83
1:B:158:PHE:HB3	1:B:159:PRO:HD3	1.61	0.82
1:B:163:PHE:HE2	1:B:246:LEU:HD23	1.46	0.80
1:B:181:THR:HG23	1:B:247:LEU:HD12	1.70	0.73
1:B:97:ASP:O	1:B:240:ARG:HD2	1.90	0.72
1:B:181:THR:CG2	1:B:251:ASP:HB2	2.20	0.71
1:A:268:PRO:O	1:A:272:GLN:HG2	1.89	0.71
1:A:163:PHE:HE2	1:A:246:LEU:HD22	1.56	0.71
1:B:269:GLU:CD	1:B:269:GLU:H	1.91	0.70
1:A:131:ILE:HG12	1:A:162:ILE:HD13	1.74	0.69
1:B:163:PHE:CE2	1:B:246:LEU:HD23	2.29	0.68
1:B:392:LYS:HB2	1:B:400:GLN:HE21	1.59	0.67
1:A:365:ARG:HA	1:A:365:ARG:HE	1.59	0.67
1:B:211:ARG:HG2	1:B:221:SER:OG	1.96	0.66
1:B:181:THR:HG21	1:B:251:ASP:HB2	1.75	0.66
1:B:110:GLN:HG3	1:B:229:ASN:HB2	1.76	0.65
1:B:272:GLN:HG2	4:B:2586:HOH:O	1.96	0.64
1:A:51:PRO:HG2	1:A:54:VAL:HG12	1.78	0.63
1:A:209:GLU:HG2	1:A:213:GLN:HE21	1.63	0.63
1:A:212:ARG:HA	1:A:225:ASN:HD21	1.62	0.63
1:B:105:PRO:HG3	1:B:108:GLN:NE2	2.13	0.62
1:A:79:ARG:HD3	4:A:1494:HOH:O	1.99	0.62
1:A:303:SER:HA	1:A:314:LYS:HB2	1.81	0.61
1:B:68:GLY:HA3	1:B:331:GLU:OE2	2.02	0.60
1:A:130:ARG:HH11	1:A:165:LEU:HD13	1.67	0.60
1:B:151:THR:HA	1:B:155:ALA:HB3	1.83	0.59
1:B:91:GLU:CD	1:B:91:GLU:H	2.05	0.59
1:B:110:GLN:HG2	4:B:2538:HOH:O	2.02	0.59
1:B:218:ASP:O	1:B:222:ILE:HG12	2.02	0.59
1:B:249:GLY:N	3:B:417:HEM:HBC2	2.17	0.59
1:A:294:LEU:H	1:A:294:LEU:HD23	1.67	0.59
1:B:56:THR:O	1:B:61:GLY:HA2	2.04	0.58
1:B:11:LEU:HD12	1:B:57:ARG:HB2	1.85	0.58
1:B:170:PRO:HB2	1:B:172:GLU:OE1	2.03	0.58
1:B:223:VAL:CG2	1:B:241:MET:HE2	2.30	0.58
1:B:172:GLU:CD	1:B:172:GLU:H	2.07	0.58
1:B:131:ILE:HG12	1:B:162:ILE:HD13	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:MET:CE	1:A:197:LYS:HA	2.35	0.57
1:B:286:GLU:OE2	1:B:364:ARG:NE	2.33	0.57
1:A:229:ASN:CG	1:A:229:ASN:O	2.43	0.56
1:B:201:TYR:O	1:B:205:ILE:HG13	2.06	0.56
1:A:273:GLU:HB3	1:A:280:ARG:NH2	2.20	0.56
1:A:83:SER:HB3	1:A:101:THR:O	2.05	0.56
1:A:38:VAL:HG22	1:A:391:HIS:CB	2.32	0.56
1:A:39:GLN:NE2	1:A:39:GLN:H	2.04	0.55
1:A:269:GLU:H	1:A:269:GLU:CD	2.07	0.55
1:B:294:LEU:HD23	1:B:294:LEU:H	1.71	0.55
1:B:223:VAL:HG22	1:B:241:MET:CE	2.34	0.55
1:B:87:TRP:CZ2	1:B:395:ILE:HG21	2.42	0.54
1:B:158:PHE:CB	1:B:159:PRO:HD3	2.37	0.54
1:B:201:TYR:HB3	1:B:239:LYS:HD2	1.89	0.54
1:B:110:GLN:CB	1:B:228:VAL:HG23	2.34	0.53
1:A:17:HIS:CD2	1:A:313:LYS:HG2	2.42	0.53
1:A:205:ILE:HB	1:A:206:PRO:HD3	1.90	0.53
1:A:110:GLN:NE2	1:A:229:ASN:HB3	2.23	0.53
1:B:174:ILE:HB	1:B:175:PRO:HD3	1.91	0.53
1:B:344:LYS:CA	1:B:344:LYS:HE3	2.33	0.53
1:B:110:GLN:CG	1:B:229:ASN:HB2	2.39	0.52
1:A:130:ARG:NH1	1:A:165:LEU:HD13	2.23	0.52
1:B:228:VAL:O	1:B:231:ARG:HG3	2.09	0.52
1:B:83:SER:HB3	1:B:101:THR:O	2.09	0.52
1:A:294:LEU:HD23	1:A:294:LEU:N	2.23	0.52
1:A:72:ARG:HH12	1:A:331:GLU:CD	2.11	0.52
1:A:163:PHE:CE2	1:A:246:LEU:HD22	2.42	0.52
1:A:10:ASN:CG	1:A:11:LEU:H	2.13	0.51
1:B:121:MET:HB3	1:B:122:PRO:HD3	1.92	0.51
1:A:174:ILE:HB	1:A:175:PRO:HD3	1.93	0.51
1:A:151:THR:HA	1:A:155:ALA:HB3	1.93	0.51
1:A:83:SER:O	1:A:86:PRO:HD3	2.10	0.50
1:A:121:MET:HB3	1:A:122:PRO:HD3	1.92	0.50
1:A:158:PHE:HB3	1:A:159:PRO:CD	2.41	0.50
1:A:276:GLU:C	1:A:277:ARG:HD2	2.31	0.50
3:A:417:HEM:HBC2	3:A:417:HEM:HMC2	1.94	0.50
1:B:36:ALA:HB1	4:B:2596:HOH:O	2.12	0.50
1:A:110:GLN:HE21	1:A:229:ASN:HB3	1.77	0.50
1:B:97:ASP:OD1	1:B:240:ARG:HD3	2.12	0.49
1:B:392:LYS:HD3	1:B:393:SER:N	2.27	0.49
1:A:170:PRO:HG2	1:A:173:ASP:OD1	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:LEU:HD13	1:A:250:LEU:HD13	1.95	0.49
1:B:181:THR:HG23	1:B:247:LEU:CD1	2.42	0.49
1:B:341:SER:O	1:B:342:ARG:C	2.52	0.49
1:B:205:ILE:HB	1:B:206:PRO:HD3	1.95	0.48
1:B:83:SER:O	1:B:86:PRO:HD3	2.13	0.48
1:B:110:GLN:HB3	1:B:228:VAL:CG2	2.34	0.48
1:A:201:TYR:HB3	1:A:239:LYS:HD2	1.95	0.48
1:A:271:ARG:O	1:A:275:ILE:HG13	2.14	0.48
1:A:82:SER:OG	1:A:84:GLU:HB2	2.13	0.48
1:B:11:LEU:N	1:B:11:LEU:HD22	2.29	0.47
1:B:294:LEU:N	1:B:294:LEU:HD23	2.28	0.47
1:A:68:GLY:HA3	1:A:331:GLU:OE2	2.14	0.47
1:A:126:LYS:HB2	4:A:1565:HOH:O	2.13	0.47
1:B:412:LYS:CB	1:B:412:LYS:NZ	2.77	0.47
1:B:179:TYR:O	1:B:183:GLN:HG2	2.14	0.47
1:A:201:TYR:O	1:A:205:ILE:HG13	2.14	0.47
1:B:269:GLU:CD	1:B:269:GLU:N	2.65	0.47
1:B:98:PHE:HB3	1:B:244:LEU:HB2	1.96	0.46
1:A:115:ALA:HB3	1:A:358:LEU:HD13	1.97	0.46
1:A:350:PHE:HB3	1:A:357:CYS:HB3	1.98	0.46
1:B:169:LEU:HD21	1:B:220:ILE:HD12	1.97	0.46
1:A:373:GLU:O	1:A:376:THR:HG22	2.16	0.46
1:A:281:ILE:HD12	1:A:372:LYS:HD3	1.97	0.46
1:B:70:LEU:HD12	1:B:70:LEU:N	2.31	0.46
1:B:191:MET:HB3	1:B:195:GLU:HB2	1.98	0.45
1:B:192:THR:OG1	1:B:195:GLU:HG3	2.16	0.45
1:A:355:HIS:HD1	3:A:417:HEM:CGD	2.30	0.45
1:A:268:PRO:HD2	1:A:269:GLU:OE2	2.16	0.45
1:B:118:VAL:CG2	1:B:219:ALA:HB1	2.47	0.45
1:B:39:GLN:NE2	1:B:39:GLN:H	2.14	0.45
1:B:303:SER:HA	1:B:314:LYS:HB2	1.98	0.44
1:A:108:GLN:HE21	1:A:108:GLN:HB3	1.59	0.44
1:B:118:VAL:HG21	1:B:219:ALA:HB1	1.99	0.44
1:A:115:ALA:O	1:A:118:VAL:HG22	2.17	0.44
1:A:286:GLU:HG2	4:A:1488:HOH:O	2.17	0.44
1:A:146:GLY:HA2	1:A:406:TRP:NE1	2.33	0.44
1:A:56:THR:O	1:A:61:GLY:HA2	2.18	0.44
1:B:108:GLN:HB3	1:B:108:GLN:HE21	1.56	0.44
1:A:294:LEU:CD2	1:A:294:LEU:N	2.81	0.44
1:B:167:ALA:O	1:B:211:ARG:NH2	2.51	0.43
1:A:72:ARG:HD3	1:A:352:HIS:CE1	2.53	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:149:ASN:HD22	1:B:149:ASN:C	2.22	0.43
1:A:114:LEU:HD23	1:A:241:MET:CE	2.48	0.43
1:B:33:ASN:HB3	1:B:41:ALA:HA	1.98	0.43
1:A:87:TRP:CZ2	1:A:395:ILE:HG21	2.53	0.43
1:A:293:SER:OG	1:A:323:MET:HA	2.19	0.43
1:A:85:CYS:HB3	1:A:300:ILE:HB	2.01	0.43
1:A:130:ARG:HE	1:A:165:LEU:HD13	1.84	0.43
1:A:156:GLU:N	1:A:157:PRO:HD2	2.34	0.43
1:A:72:ARG:NH1	1:A:331:GLU:OE2	2.40	0.43
1:A:118:VAL:HG23	1:A:119:VAL:HG13	1.99	0.42
1:A:78:TYR:CE1	1:A:79:ARG:HD2	2.54	0.42
1:A:167:ALA:HA	1:A:219:ALA:HB3	2.01	0.42
1:B:15:PRO:HA	1:B:16:PRO:HD3	1.93	0.42
1:A:382:SER:HB2	4:A:1547:HOH:O	2.19	0.42
1:B:147:GLN:O	1:B:148:CYS:HB3	2.19	0.42
1:A:10:ASN:ND2	1:A:11:LEU:H	2.18	0.42
1:A:277:ARG:HB3	1:A:279:GLU:OE1	2.20	0.42
1:B:115:ALA:HB3	1:B:358:LEU:HD13	2.00	0.42
1:B:19:PRO:HD3	4:B:2437:HOH:O	2.19	0.42
1:A:343:GLN:HA	1:A:343:GLN:HE21	1.85	0.42
1:A:384:ALA:HB3	1:A:403:PRO:HB2	2.02	0.41
1:B:344:LYS:CE	1:B:344:LYS:HA	2.38	0.41
1:A:130:ARG:HE	1:A:165:LEU:CD1	2.34	0.41
1:A:87:TRP:CH2	1:A:395:ILE:HG21	2.54	0.41
1:A:114:LEU:HD23	1:A:241:MET:HE1	2.02	0.41
1:A:165:LEU:HD11	1:B:20:GLU:OE1	2.20	0.41
1:A:14:LEU:HD11	1:A:18:VAL:CG1	2.51	0.41
1:A:287:GLU:HA	1:A:287:GLU:OE1	2.21	0.41
1:A:165:LEU:HD23	1:A:166:LEU:N	2.36	0.41
1:B:100:PRO:HG2	3:B:417:HEM:CGD	2.51	0.41
1:A:365:ARG:HA	1:A:365:ARG:NE	2.31	0.40
1:A:343:GLN:NE2	1:A:343:GLN:HA	2.36	0.40
1:A:272:GLN:HA	1:A:272:GLN:NE2	2.37	0.40
1:A:75:TYR:CZ	1:A:320:LEU:HB2	2.56	0.40
1:A:147:GLN:O	1:A:148:CYS:HB3	2.21	0.40
1:B:188:ASP:OD1	1:B:190:SER:OG	2.35	0.40
1:B:384:ALA:HA	1:B:385:PRO:HD3	1.98	0.40
1:B:72:ARG:NH1	1:B:331:GLU:OE2	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	403/414 (97%)	386 (96%)	17 (4%)	0	100	100
1	B	403/414 (97%)	383 (95%)	18 (4%)	2 (0%)	34	30
All	All	806/828 (97%)	769 (95%)	35 (4%)	2 (0%)	52	53

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	25	ASP
1	B	321	PRO

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	349/357 (98%)	336 (96%)	13 (4%)	41	41
1	B	349/357 (98%)	334 (96%)	15 (4%)	35	34
All	All	698/714 (98%)	670 (96%)	28 (4%)	38	38

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	38	VAL
1	A	39	GLN
1	A	79	ARG
1	A	86	PRO

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Mol	Chain	Res	Type
1	A	87	TRP
1	A	108	GLN
1	A	109	ARG
1	A	229	ASN
1	A	246	LEU
1	A	269	GLU
1	A	276	GLU
1	A	365	ARG
1	A	402	LEU
1	B	33	ASN
1	B	39	GLN
1	B	47	GLU
1	B	69	GLN
1	B	86	PRO
1	B	87	TRP
1	B	108	GLN
1	B	132	GLN
1	B	138	LEU
1	B	149	ASN
1	B	211	ARG
1	B	229	ASN
1	B	269	GLU
1	B	344	LYS
1	B	356	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	10	ASN
1	A	30	ASN
1	A	39	GLN
1	A	46	GLN
1	A	108	GLN
1	A	110	GLN
1	A	117	GLN
1	A	129	ASN
1	A	132	GLN
1	A	213	GLN
1	A	225	ASN
1	A	272	GLN
1	A	311	GLN
1	A	343	GLN

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Mol	Chain	Res	Type
1	A	400	GLN
1	B	33	ASN
1	B	39	GLN
1	B	46	GLN
1	B	108	GLN
1	B	132	GLN
1	B	149	ASN
1	B	210	GLN
1	B	225	ASN
1	B	229	ASN
1	B	343	GLN
1	B	390	GLN
1	B	400	GLN

5.3.3 RNA [i](#)

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](#)

Of 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 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	$\# Z > 2$
3	HEM	A	417	1,4	30,50,50	2.45	10 (33%)	24,82,82	2.73	11 (45%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	HEM	B	417	1,4	30,50,50	2.41	10 (33%)	24,82,82	2.79	11 (45%)

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	HEM	A	417	1,4	-	0/10/54/54	0/0/8/8
3	HEM	B	417	1,4	-	0/10/54/54	0/0/8/8

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	417	HEM	C2D-C3D	-6.78	1.34	1.54
3	B	417	HEM	C2D-C3D	-6.53	1.34	1.54
3	B	417	HEM	C3B-C4B	-4.37	1.48	1.51
3	A	417	HEM	C3B-C4B	-4.30	1.48	1.51
3	B	417	HEM	C2C-C1C	-4.01	1.45	1.52
3	B	417	HEM	C3D-C4D	-3.73	1.46	1.51
3	A	417	HEM	C3D-C4D	-3.63	1.46	1.51
3	A	417	HEM	C2C-C1C	-3.46	1.46	1.52
3	B	417	HEM	FE-ND	2.13	2.08	1.97
3	A	417	HEM	FE-ND	2.18	2.09	1.97
3	B	417	HEM	FE-NC	2.40	2.05	1.95
3	A	417	HEM	FE-NC	2.44	2.05	1.95
3	A	417	HEM	CAA-C2A	2.93	1.57	1.52
3	B	417	HEM	CAA-C2A	3.10	1.57	1.52
3	B	417	HEM	C1C-NC	3.32	1.40	1.36
3	A	417	HEM	C1C-NC	3.36	1.40	1.36
3	A	417	HEM	C3B-CAB	3.55	1.58	1.51
3	B	417	HEM	C3B-CAB	3.84	1.58	1.51
3	B	417	HEM	C3C-CAC	3.95	1.58	1.51
3	A	417	HEM	C3C-CAC	5.14	1.61	1.51

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	417	HEM	C3C-CAC-CBC	-4.44	117.64	124.46
3	B	417	HEM	CMA-C3A-C4A	-3.14	123.17	128.36
3	A	417	HEM	CMA-C3A-C4A	-3.05	123.32	128.36
3	A	417	HEM	C3C-CAC-CBC	-2.90	120.01	124.46

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	417	HEM	CBD-CAD-C3D	2.07	119.58	113.55
3	A	417	HEM	CBD-CAD-C3D	2.08	119.60	113.55
3	A	417	HEM	CMA-C3A-C2A	2.27	129.99	125.24
3	B	417	HEM	CMA-C3A-C2A	2.38	130.22	125.24
3	A	417	HEM	CAD-C3D-C4D	2.83	122.45	112.47
3	B	417	HEM	CAD-C3D-C4D	2.86	122.57	112.47
3	A	417	HEM	CBA-CAA-C2A	2.95	117.82	112.53
3	B	417	HEM	CBA-CAA-C2A	2.95	117.82	112.53
3	A	417	HEM	CMD-C2D-C3D	3.34	129.11	114.35
3	B	417	HEM	CMD-C2D-C3D	3.36	129.20	114.35
3	A	417	HEM	C2D-C3D-C4D	4.58	109.27	101.50
3	B	417	HEM	C2D-C3D-C4D	4.60	109.30	101.50
3	B	417	HEM	CMC-C2C-C3C	4.62	128.05	116.53
3	A	417	HEM	CMC-C2C-C3C	5.03	129.09	116.53
3	B	417	HEM	CAD-C3D-C2D	5.18	128.12	113.22
3	A	417	HEM	CAD-C3D-C2D	5.24	128.28	113.22
3	A	417	HEM	CMB-C2B-C3B	5.87	131.18	116.53
3	B	417	HEM	CMB-C2B-C3B	5.90	131.26	116.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	417	HEM	2	0
3	B	417	HEM	2	0

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 [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, 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	405/414 (97%)	-0.28	5 (1%) 81 85	12, 20, 37, 59	0
1	B	405/414 (97%)	-0.25	2 (0%) 91 93	13, 22, 38, 54	0
All	All	810/828 (97%)	-0.27	7 (0%) 85 88	12, 21, 38, 59	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	213	GLN	3.5
1	B	48	SER	3.3
1	B	229	ASN	3.1
1	A	10	ASN	2.9
1	A	385	PRO	2.7
1	A	229	ASN	2.6
1	A	412	LYS	2.1

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(\AA^2)	Q<0.9
3	HEM	A	417	43/43	0.97	0.11	0.83	11,14,17,20	0
3	HEM	B	417	43/43	0.96	0.11	0.59	12,16,19,25	0
2	K	B	2418	1/1	0.99	0.07	-0.74	17,17,17,17	0
2	K	A	1418	1/1	0.99	0.05	-2.30	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.