Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:55 PM GMT

PDB ID	:	1R9O
Title	:	Crystal Structure of P4502C9 with Flurbiprofen bound
Authors	:	Wester, M.R.; Yano, J.K.; Schoch, G.A.; Griffin, K.J.; Stout, C.D.; Johnson, E.F.
Deposited on	:	2003-10-30
Resolution	:	2.00 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 ①) were used in the production of this report:

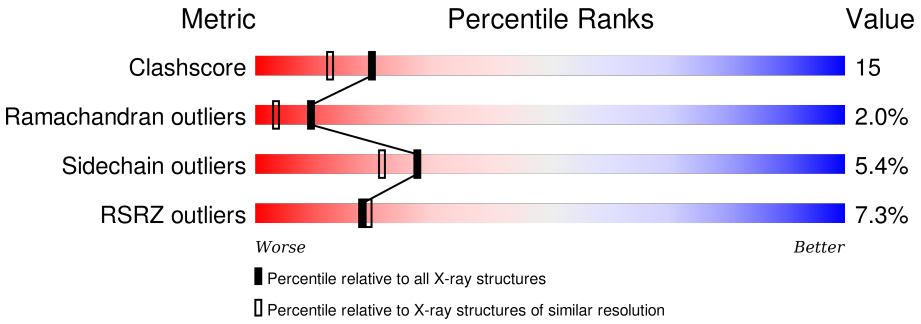
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)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

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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))
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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 <=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	477	7% 71% 20% • 5%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 3972 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.

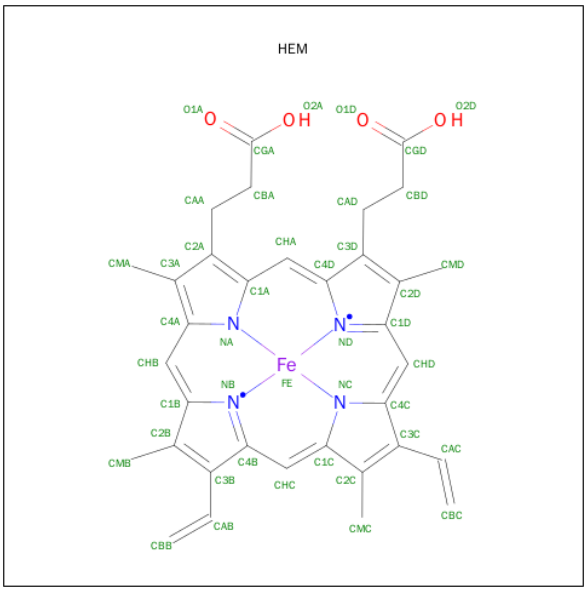
- Molecule 1 is a protein called Cytochrome P450 2C9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	455	3650	2356	615	656	23	4	0	0

There are 10 discrepancies between the modelled and reference sequences:

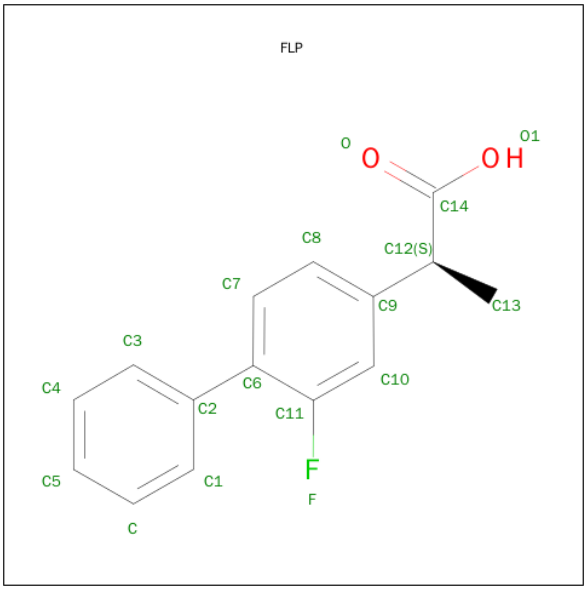
Chain	Residue	Modelled	Actual	Comment	Reference
A	18	MET	-	ENGINEERED	UNP P11712
A	19	ALA	-	ENGINEERED	UNP P11712
A	20	LYS	-	ENGINEERED	UNP P11712
A	21	LYS	-	ENGINEERED	UNP P11712
A	22	THR	-	ENGINEERED	UNP P11712
A	490	ILE	VAL	ENGINEERED	UNP P11712
A	491	HIS	-	EXPRESSION TAG	UNP P11712
A	492	HIS	-	EXPRESSION TAG	UNP P11712
A	493	HIS	-	EXPRESSION TAG	UNP P11712
A	494	HIS	-	EXPRESSION TAG	UNP P11712

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



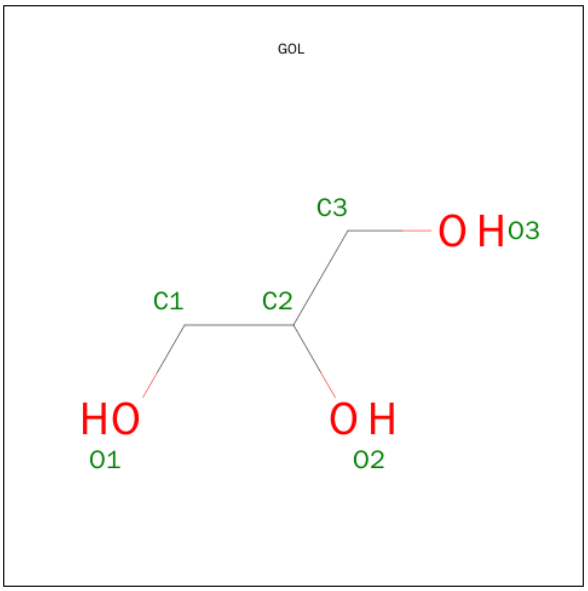
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is FLURBIPROFEN (three-letter code: FLP) (formula: $C_{15}H_{13}FO_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	F	O	0	0
			18	15	1	2		

- 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		

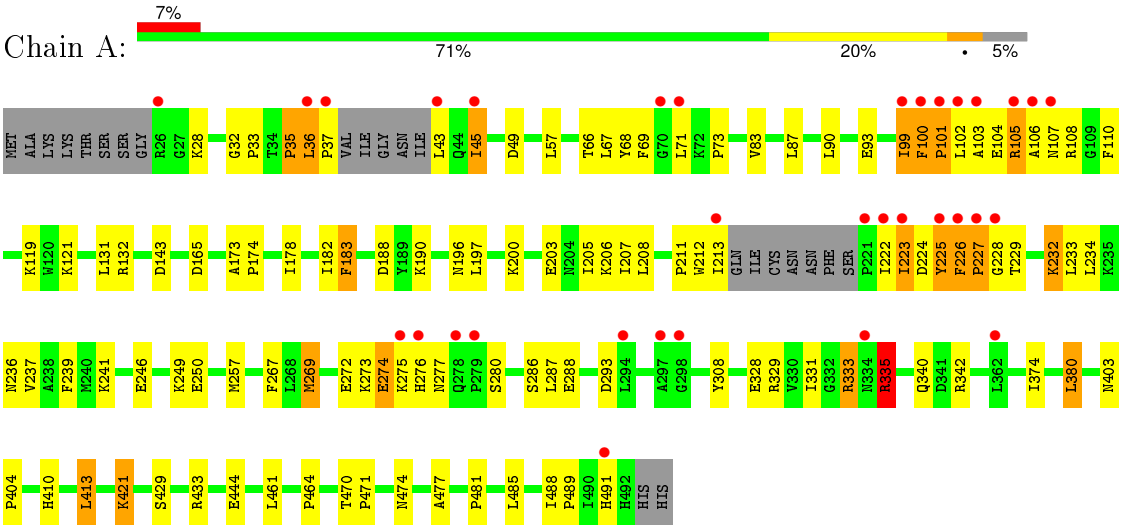
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	255	Total	O	0	0
			255	255		

3 Residue-property plots (i)

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 2C9



4 Data and refinement statistics (i)

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α, β, γ	91.05Å 91.05Å 169.48Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 2.00 31.14 – 2.00	Depositor EDS
% Data completeness (in resolution range)	99.8 (50.00-2.00) 100.0 (31.14-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	2.23 (at 2.00Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.191 , 0.236 0.192 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	29.4	Xtriage
Anisotropy	0.011	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 57.9	EDS
Estimated twinning fraction	0.032 for h,-h-k,-l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Outliers	0 of 35396 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	3972	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.22% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: FLP, GOL, 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	0.89	2/3739 (0.1%)	0.91	7/5051 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	83	VAL	CB-CG1	5.42	1.64	1.52
1	A	444	GLU	CD-OE2	-5.32	1.19	1.25

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	433	ARG	NE-CZ-NH1	-11.08	114.76	120.30
1	A	433	ARG	NE-CZ-NH2	7.05	123.83	120.30
1	A	132	ARG	NE-CZ-NH2	-6.77	116.91	120.30
1	A	183	PHE	CB-CA-C	-6.10	98.21	110.40
1	A	335	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	A	165	ASP	N-CA-C	-5.76	95.45	111.00
1	A	228	GLY	N-CA-C	-5.03	100.53	113.10

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	3650	0	3689	111	0
2	A	43	0	30	0	0
3	A	18	0	12	2	0
4	A	6	0	8	2	0
5	A	255	0	0	7	0
All	All	3972	0	3739	111	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 15.

All (111) 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:232:LYS:HE2	1:A:236:ASN:ND2	1.59	1.14
1:A:232:LYS:HE2	1:A:236:ASN:CG	1.87	0.94
1:A:35:PRO:C	1:A:37:PRO:HD3	1.86	0.94
1:A:232:LYS:HE2	1:A:236:ASN:HD21	1.34	0.93
1:A:333:ARG:NE	1:A:333:ARG:H	1.72	0.88
1:A:208:LEU:HD21	1:A:233:LEU:HD21	1.56	0.88
1:A:226:PHE:H	1:A:227:PRO:HD2	1.35	0.87
1:A:226:PHE:H	1:A:227:PRO:CD	2.01	0.73
1:A:99:ILE:O	1:A:100:PHE:HB2	1.88	0.73
1:A:274:GLU:HG2	1:A:280:SER:OG	1.89	0.71
1:A:93:GLU:HG2	1:A:374:ILE:HD12	1.74	0.70
1:A:226:PHE:CD1	1:A:227:PRO:HD3	2.30	0.66
1:A:208:LEU:CD2	1:A:233:LEU:HD21	2.26	0.66
1:A:68:TYR:CE2	1:A:73:PRO:HB3	2.32	0.64
1:A:173:ALA:HB3	1:A:174:PRO:HD3	1.78	0.64
1:A:203:GLU:OE1	1:A:232:LYS:HD3	1.98	0.63
1:A:93:GLU:CB	1:A:374:ILE:HD12	2.28	0.63
1:A:226:PHE:O	1:A:227:PRO:C	2.37	0.63
1:A:196:ASN:O	1:A:200:LYS:HG2	2.00	0.62
1:A:105:ARG:HA	1:A:105:ARG:HE	1.65	0.61
1:A:105:ARG:HA	1:A:105:ARG:NE	2.16	0.61
1:A:232:LYS:HE2	1:A:236:ASN:OD1	2.00	0.61
1:A:90:LEU:HD13	1:A:374:ILE:HD11	1.83	0.61
1:A:93:GLU:CG	1:A:374:ILE:HD12	2.31	0.61
1:A:223:ILE:HG13	1:A:224:ASP:H	1.67	0.60
1:A:105:ARG:C	1:A:107:ASN:N	2.55	0.57
1:A:36:LEU:N	1:A:37:PRO:HD3	2.19	0.57
1:A:225:TYR:CD2	1:A:229:THR:HG21	2.39	0.57
1:A:213:ILE:HD13	1:A:224:ASP:OD2	2.05	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:ILE:HG13	1:A:223:ILE:N	2.21	0.56
1:A:106:ALA:HB2	1:A:234:LEU:HD23	1.88	0.55
1:A:491:HIS:HD2	5:A:758:HOH:O	1.90	0.55
1:A:178:ILE:HG22	1:A:182:ILE:HD12	1.88	0.54
1:A:331:ILE:HG23	1:A:335:ARG:CZ	2.38	0.54
1:A:257:MET:SD	1:A:269:MET:HG2	2.47	0.54
1:A:477:ALA:CB	4:A:502:GOL:H32	2.38	0.54
1:A:222:ILE:HG13	1:A:223:ILE:H	1.73	0.54
1:A:106:ALA:HB2	1:A:234:LEU:CD2	2.38	0.54
1:A:237:VAL:O	1:A:241:LYS:HG3	2.08	0.53
1:A:246:GLU:O	1:A:250:GLU:HG3	2.09	0.53
1:A:36:LEU:N	1:A:37:PRO:CD	2.73	0.52
1:A:102:LEU:HG	1:A:103:ALA:H	1.74	0.52
1:A:223:ILE:HG23	1:A:224:ASP:N	2.23	0.51
1:A:232:LYS:CE	1:A:236:ASN:OD1	2.57	0.51
1:A:410:HIS:HA	1:A:413:LEU:HB2	1.91	0.51
1:A:105:ARG:C	1:A:107:ASN:H	2.14	0.51
1:A:275:LYS:O	1:A:276:HIS:HB2	2.09	0.51
1:A:249:LYS:HE2	5:A:832:HOH:O	2.10	0.50
1:A:421:LYS:HD3	5:A:753:HOH:O	2.12	0.49
1:A:232:LYS:CE	1:A:236:ASN:CG	2.71	0.49
1:A:485:LEU:N	1:A:485:LEU:HD23	2.27	0.49
1:A:477:ALA:HB2	4:A:502:GOL:H32	1.93	0.49
1:A:222:ILE:O	1:A:223:ILE:O	2.32	0.48
1:A:107:ASN:OD1	1:A:110:PHE:HD1	1.97	0.48
1:A:211:PRO:O	1:A:224:ASP:HB2	2.13	0.48
1:A:45:ILE:HG13	1:A:49:ASP:HB2	1.94	0.48
1:A:208:LEU:HG	1:A:233:LEU:HD11	1.95	0.47
1:A:410:HIS:CE1	1:A:413:LEU:HD23	2.49	0.47
1:A:69:PHE:CD1	1:A:69:PHE:N	2.82	0.47
1:A:293:ASP:HA	3:A:501:FLP:H8	1.96	0.47
1:A:257:MET:SD	1:A:269:MET:CG	3.04	0.46
1:A:226:PHE:CD1	1:A:227:PRO:CD	2.98	0.46
1:A:90:LEU:HD13	1:A:374:ILE:CD1	2.46	0.46
1:A:102:LEU:HD22	5:A:843:HOH:O	2.16	0.46
1:A:267:PHE:CG	1:A:287:LEU:HD13	2.51	0.46
1:A:470:THR:HA	1:A:471:PRO:HD3	1.85	0.45
1:A:226:PHE:N	1:A:227:PRO:CD	2.72	0.45
1:A:223:ILE:HG13	1:A:224:ASP:N	2.32	0.45
1:A:102:LEU:HB3	1:A:104:GLU:OE2	2.16	0.45
1:A:211:PRO:O	1:A:224:ASP:CB	2.65	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:ASN:ND2	5:A:828:HOH:O	2.49	0.45
1:A:241:LYS:HB3	1:A:288:GLU:OE2	2.17	0.44
1:A:329:ARG:HH11	1:A:329:ARG:HG3	1.82	0.44
1:A:43:LEU:N	1:A:43:LEU:HD12	2.33	0.44
1:A:35:PRO:O	1:A:36:LEU:HB2	2.17	0.44
1:A:28:LYS:HB3	1:A:380:LEU:HD12	2.00	0.44
1:A:207:ILE:HG21	1:A:233:LEU:HD13	1.99	0.43
1:A:100:PHE:CD2	1:A:101:PRO:HD2	2.53	0.43
1:A:470:THR:O	1:A:481:PRO:HD3	2.19	0.43
1:A:32:GLY:HA3	1:A:66:THR:O	2.19	0.43
1:A:200:LYS:HD2	1:A:236:ASN:OD1	2.18	0.43
1:A:269:MET:O	1:A:273:LYS:HD3	2.18	0.43
1:A:107:ASN:ND2	1:A:110:PHE:CD1	2.86	0.43
1:A:104:GLU:CB	1:A:108:ARG:HG2	2.48	0.43
1:A:272:GLU:O	1:A:275:LYS:HG2	2.18	0.43
1:A:342:ARG:HD2	5:A:737:HOH:O	2.19	0.43
1:A:461:LEU:HD22	1:A:461:LEU:N	2.34	0.43
1:A:35:PRO:O	1:A:37:PRO:HD3	2.19	0.42
1:A:188:ASP:OD1	1:A:190:LYS:HB2	2.19	0.42
1:A:57:LEU:HD13	1:A:67:LEU:HD21	2.01	0.42
1:A:108:ARG:HH21	3:A:501:FLP:H12	1.84	0.42
1:A:45:ILE:HG23	1:A:212:TRP:HB2	2.02	0.42
1:A:100:PHE:HD2	1:A:101:PRO:HD2	1.84	0.42
1:A:328:GLU:OE2	1:A:333:ARG:NH1	2.53	0.42
1:A:488:ILE:O	1:A:489:PRO:C	2.58	0.42
1:A:340:GLN:HG3	5:A:699:HOH:O	2.19	0.41
1:A:239:PHE:C	1:A:239:PHE:CD1	2.94	0.41
1:A:333:ARG:H	1:A:333:ARG:HE	1.60	0.41
1:A:331:ILE:HG22	1:A:331:ILE:O	2.21	0.41
1:A:104:GLU:HB3	1:A:108:ARG:HG2	2.02	0.41
1:A:107:ASN:OD1	1:A:110:PHE:CD1	2.74	0.41
1:A:206:LYS:HG3	1:A:206:LYS:O	2.18	0.41
1:A:33:PRO:HB2	1:A:37:PRO:HG2	2.02	0.41
1:A:333:ARG:CD	1:A:333:ARG:H	2.34	0.41
1:A:28:LYS:CB	1:A:380:LEU:HD12	2.51	0.41
1:A:403:ASN:N	1:A:404:PRO:CD	2.83	0.41
1:A:121:LYS:HB2	1:A:121:LYS:HE2	1.73	0.41
1:A:104:GLU:O	1:A:107:ASN:N	2.42	0.40
1:A:474:ASN:HB3	1:A:477:ALA:O	2.21	0.40
1:A:99:ILE:CG1	1:A:100:PHE:N	2.85	0.40
1:A:104:GLU:C	1:A:106:ALA:N	2.75	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	449/477 (94%)	419 (93%)	21 (5%)	9 (2%)	9 3

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	35	PRO
1	A	36	LEU
1	A	223	ILE
1	A	226	PHE
1	A	227	PRO
1	A	429	SER
1	A	99	ILE
1	A	101	PRO
1	A	100	PHE

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	411/430 (96%)	389 (95%)	22 (5%)	27 21

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	ILE
1	A	71	LEU
1	A	87	LEU
1	A	105	ARG
1	A	119	LYS
1	A	131	LEU
1	A	143	ASP
1	A	183	PHE
1	A	197	LEU
1	A	205	ILE
1	A	225	TYR
1	A	232	LYS
1	A	269	MET
1	A	274	GLU
1	A	286	SER
1	A	308	TYR
1	A	333	ARG
1	A	335	ARG
1	A	380	LEU
1	A	413	LEU
1	A	421	LYS
1	A	464	PRO

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	316	HIS
1	A	398	ASN

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)

3 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	500	1,5	30,50,50	3.26	13 (43%)	24,82,82	2.58	8 (33%)
3	FLP	A	501	-	16,19,19	3.86	15 (93%)	21,26,26	1.18	2 (9%)
4	GOL	A	502	-	5,5,5	1.17	0	5,5,5	0.67	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	500	1,5	-	0/10/54/54	0/0/8/8
3	FLP	A	501	-	-	0/8/12/12	0/2/2/2
4	GOL	A	502	-	-	0/4/4/4	0/0/0/0

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	500	HEM	C3B-C4B	-7.93	1.44	1.51
2	A	500	HEM	C3C-CAC	-6.70	1.38	1.51
2	A	500	HEM	C2D-C3D	-6.56	1.34	1.54
2	A	500	HEM	C2C-C1C	-6.28	1.40	1.52
2	A	500	HEM	C3D-C4D	-5.82	1.44	1.51
2	A	500	HEM	C2D-C1D	-3.63	1.40	1.51
2	A	500	HEM	C3B-CAB	-3.46	1.44	1.51
2	A	500	HEM	C1C-NC	-3.34	1.31	1.36
3	A	501	FLP	C6-C2	-3.24	1.44	1.49
2	A	500	HEM	C2B-C1B	-2.78	1.42	1.51
3	A	501	FLP	C13-C12	2.06	1.57	1.53
3	A	501	FLP	C4-C3	2.10	1.43	1.38
3	A	501	FLP	C5-C	2.46	1.44	1.38
3	A	501	FLP	C5-C4	2.49	1.44	1.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	501	FLP	C9-C12	2.57	1.57	1.52
2	A	500	HEM	FE-NB	2.67	2.11	1.97
3	A	501	FLP	C8-C7	2.86	1.43	1.38
2	A	500	HEM	CBB-CAB	2.87	1.45	1.29
3	A	501	FLP	C10-C9	3.01	1.44	1.39
2	A	500	HEM	CMA-C3A	3.33	1.58	1.51
3	A	501	FLP	C-C1	3.36	1.45	1.38
2	A	500	HEM	FE-NC	3.55	2.09	1.95
3	A	501	FLP	C10-C11	3.71	1.44	1.37
3	A	501	FLP	C1-C2	3.91	1.47	1.39
3	A	501	FLP	C8-C9	4.25	1.46	1.39
3	A	501	FLP	C3-C2	4.67	1.49	1.39
3	A	501	FLP	C7-C6	5.93	1.50	1.39
3	A	501	FLP	C6-C11	8.08	1.49	1.38

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	501	FLP	C10-C11-C6	-2.75	120.78	124.24
2	A	500	HEM	C1D-CHD-C4C	2.40	129.83	125.82
2	A	500	HEM	CMD-C2D-C3D	2.60	125.87	114.35
2	A	500	HEM	C3C-CAC-CBC	2.69	128.58	124.46
3	A	501	FLP	F-C11-C6	3.01	123.29	118.80
2	A	500	HEM	C3B-CAB-CBB	3.07	129.16	124.46
2	A	500	HEM	CAD-C3D-C4D	3.41	124.50	112.47
2	A	500	HEM	CMC-C2C-C3C	4.58	127.97	116.53
2	A	500	HEM	CAD-C3D-C2D	6.00	130.47	113.22
2	A	500	HEM	CMB-C2B-C3B	6.21	132.02	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	501	FLP	2	0
4	A	502	GOL	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

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, 95^{th} 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	455/477 (95%)	0.12	33 (7%)	18	19	17, 34, 63, 77	1 (0%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	228	GLY	7.6
1	A	221	PRO	7.1
1	A	226	PHE	7.0
1	A	36	LEU	6.6
1	A	105	ARG	5.8
1	A	225	TYR	5.6
1	A	37	PRO	5.4
1	A	100	PHE	5.3
1	A	71	LEU	5.1
1	A	102	LEU	4.9
1	A	99	ILE	4.8
1	A	227	PRO	4.3
1	A	45	ILE	4.3
1	A	276	HIS	3.9
1	A	222	ILE	3.9
1	A	106	ALA	3.8
1	A	101	PRO	3.7
1	A	223	ILE	3.7
1	A	43	LEU	3.5
1	A	107	ASN	3.4
1	A	213	ILE	2.7
1	A	334	ASN	2.7
1	A	103	ALA	2.6
1	A	294	LEU	2.5
1	A	297	ALA	2.4
1	A	26	ARG	2.4
1	A	278	GLN	2.2

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Mol	Chain	Res	Type	RSRZ
1	A	275	LYS	2.2
1	A	279	PRO	2.1
1	A	362	LEU	2.1
1	A	491	HIS	2.1
1	A	298	GLY	2.1
1	A	70	GLY	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^{th} 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	502	6/6	0.68	0.24	1.98	48,49,50,50	0
3	FLP	A	501	18/18	0.84	0.17	0.55	43,49,51,55	0
2	HEM	A	500	43/43	0.98	0.17	0.01	13,18,24,25	0

6.5 Other polymers (i)

There are no such residues in this entry.