



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 08:17 am GMT

PDB ID : 4O4P
Title : Structure of P450 BM3 A82F F87V in complex with S-omeprazol
Authors : Leys, D.
Deposited on : 2013-12-19
Resolution : 1.83 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 : 1.7.2 (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

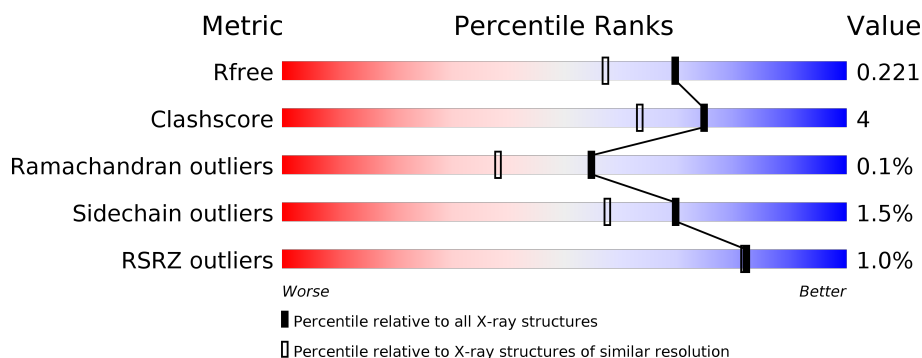
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.83 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	2964 (1.86-1.82)
Clashscore	112137	3197 (1.86-1.82)
Ramachandran outliers	110173	3164 (1.86-1.82)
Sidechain outliers	110143	3165 (1.86-1.82)
RSRZ outliers	101464	2973 (1.86-1.82)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	455	
1	B	455	

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	1C6	B	502	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7990 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

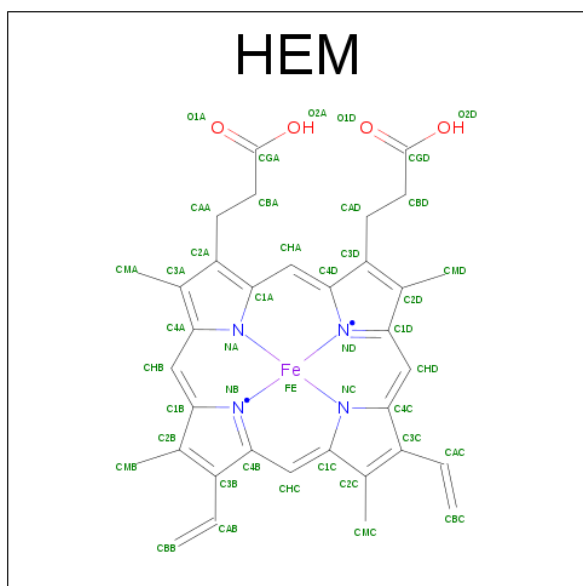
- Molecule 1 is a protein called Bifunctional P-450/NADPH-P450 reductase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	450	Total	C	N	O	S	0	6	0
			3671	2345	622	687	17			
1	B	446	Total	C	N	O	S	0	4	0
			3618	2312	613	676	17			

There are 4 discrepancies between the modelled and reference sequences:

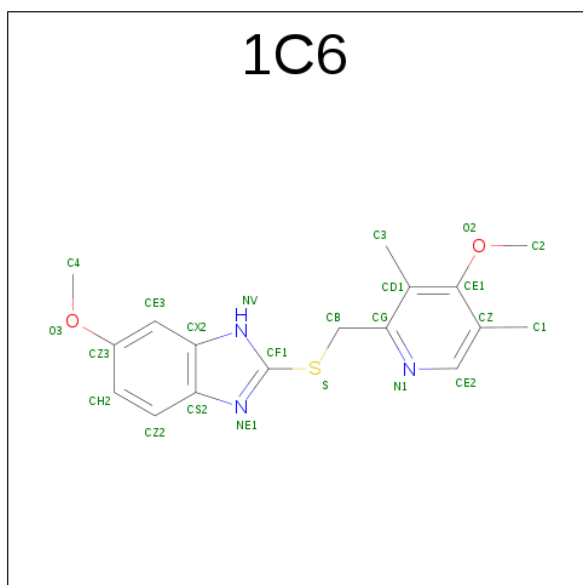
Chain	Residue	Modelled	Actual	Comment	Reference
A	82	PHE	ALA	ENGINEERED MUTATION	UNP P14779
A	87	VAL	PHE	ENGINEERED MUTATION	UNP P14779
B	82	PHE	ALA	ENGINEERED MUTATION	UNP P14779
B	87	VAL	PHE	ENGINEERED MUTATION	UNP P14779

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



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

- Molecule 3 is 6-METHOXY-2-[[[(4-METHOXY-3,5-DIMETHYLPYRIDIN-2-YL)METHYL]SULFANYL]-1H-BENZIMIDAZOLE (three-letter code: 1C6) (formula: C₁₇H₁₉N₃O₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			23	17	3	2	1		
3	B	1	Total	C	N	O	S	0	0
			23	17	3	2	1		

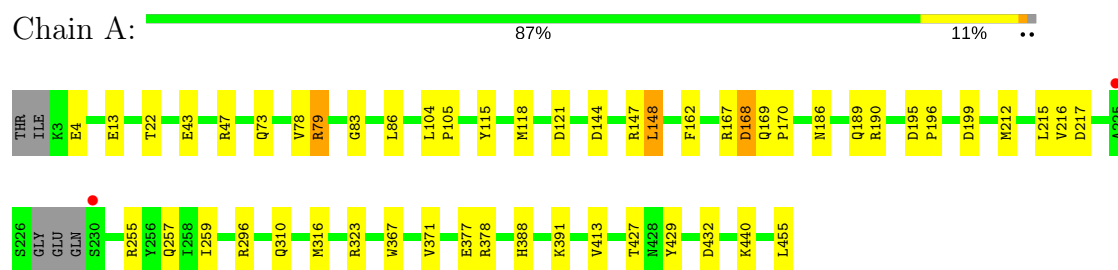
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	322	Total	O	0	0
			322	322		
4	B	247	Total	O	0	0
			247	247		

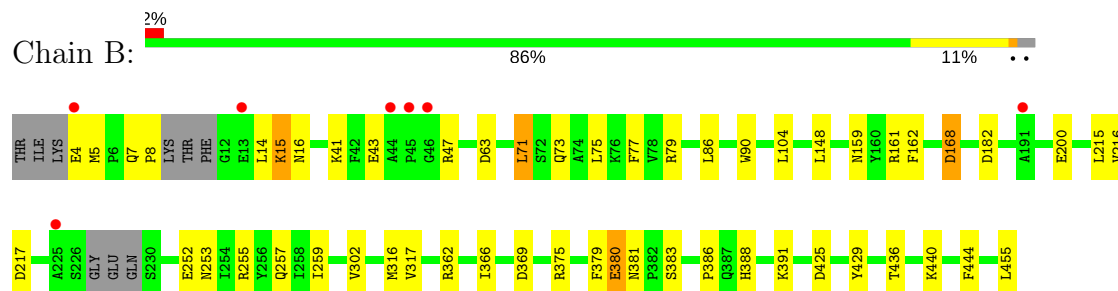
3 Residue-property plots

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: Bifunctional P-450/NADPH-P450 reductase



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.42Å 132.88Å 147.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	66.00 – 1.83 66.44 – 1.83	Depositor EDS
% Data completeness (in resolution range)	99.9 (66.00-1.83) 99.9 (66.44-1.83)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.36 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.186 , 0.221 0.188 , 0.221	Depositor DCC
R_{free} test set	5147 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	26.5	Xtriage
Anisotropy	0.021	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 39.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7990	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 11.76% 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, 0.333 respectively for untwinned datasets, and 0.375, 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, 1C6

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	1.19	3/3755 (0.1%)	1.04	17/5079 (0.3%)
1	B	1.09	7/3700 (0.2%)	0.95	7/5003 (0.1%)
All	All	1.14	10/7455 (0.1%)	1.00	24/10082 (0.2%)

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	161	ARG	CG-CD	-6.00	1.36	1.51
1	B	444	PHE	CE1-CZ	5.83	1.48	1.37
1	A	429	TYR	CE2-CZ	-5.64	1.31	1.38
1	B	429	TYR	CD1-CE1	5.50	1.47	1.39
1	B	302	VAL	CB-CG1	5.38	1.64	1.52
1	A	78	VAL	CB-CG2	5.37	1.64	1.52
1	B	379	PHE	CD2-CE2	-5.22	1.28	1.39
1	B	380	GLU	CB-CG	-5.07	1.42	1.52
1	A	43	GLU	CB-CG	-5.05	1.42	1.52
1	B	317	VAL	CB-CG1	5.05	1.63	1.52

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	432	ASP	CB-CG-OD1	10.68	127.91	118.30
1	A	323	ARG	NE-CZ-NH2	-9.29	115.65	120.30
1	A	296	ARG	NE-CZ-NH1	9.16	124.88	120.30
1	A	168[A]	ASP	CB-CG-OD1	-8.49	110.66	118.30
1	A	168[B]	ASP	CB-CG-OD1	-8.49	110.66	118.30
1	A	296	ARG	NE-CZ-NH2	-7.61	116.50	120.30
1	A	79	ARG	NE-CZ-NH1	6.85	123.72	120.30
1	A	190	ARG	NE-CZ-NH2	-6.84	116.88	120.30
1	A	147	ARG	NE-CZ-NH1	6.09	123.34	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	378	ARG	NE-CZ-NH1	5.76	123.18	120.30
1	A	144	ASP	CB-CG-OD1	5.75	123.47	118.30
1	B	168	ASP	CB-CG-OD1	-5.61	113.25	118.30
1	B	455	LEU	CA-CB-CG	5.58	128.14	115.30
1	B	375	ARG	NE-CZ-NH1	5.56	123.08	120.30
1	B	425	ASP	CB-CG-OD1	5.45	123.21	118.30
1	A	323	ARG	NE-CZ-NH1	5.45	123.03	120.30
1	A	199	ASP	CB-CG-OD1	5.35	123.11	118.30
1	A	455	LEU	CA-CB-CG	5.34	127.59	115.30
1	B	362	ARG	NE-CZ-NH1	5.25	122.93	120.30
1	B	63	ASP	CB-CG-OD2	-5.18	113.64	118.30
1	A	190	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	B	79	ARG	NE-CZ-NH1	5.12	122.86	120.30
1	A	121	ASP	CB-CG-OD1	5.09	122.89	118.30
1	A	199	ASP	CB-CG-OD2	-5.06	113.75	118.30

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	3671	0	3615	24	1
1	B	3618	0	3568	30	1
2	A	43	0	30	0	0
2	B	43	0	30	0	0
3	A	23	0	19	0	0
3	B	23	0	19	1	0
4	A	322	0	0	4	0
4	B	247	0	0	3	0
All	All	7990	0	7281	54	1

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 4.

All (54) 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:162:PHE:HE1	1:B:215:LEU:HD21	1.28	0.94
1:B:252:GLU:CD	4:B:736:HOH:O	2.08	0.91
1:B:162:PHE:CE1	1:B:215:LEU:HD21	2.08	0.87
1:B:216:VAL:HG21	1:B:259[B]:ILE:CG2	2.09	0.82
1:B:7:GLN:HE21	1:B:16:ASN:HD21	1.27	0.82
1:A:217:ASP:OD1	1:A:255:ARG:NH1	2.13	0.81
1:B:16:ASN:HD22	1:B:43:GLU:H	1.39	0.70
1:B:216:VAL:HG21	1:B:259[B]:ILE:HG22	1.73	0.70
1:B:7:GLN:HE21	1:B:16:ASN:ND2	1.89	0.69
1:A:388:HIS:HD2	1:A:391:LYS:HZ2	1.44	0.63
1:A:388:HIS:HD2	1:A:391:LYS:NZ	1.99	0.61
1:B:182[A]:ASP:OD1	1:B:436:THR:HG22	2.01	0.60
1:B:216:VAL:CG2	1:B:259[B]:ILE:HG22	2.32	0.59
1:B:47:ARG:CZ	1:B:73[A]:GLN:HG3	2.32	0.59
1:B:216:VAL:HG21	1:B:259[B]:ILE:HG23	1.84	0.59
1:A:427:THR:O	4:A:845:HOH:O	2.16	0.59
1:A:148:LEU:HD21	1:A:413:VAL:HG21	1.86	0.57
1:A:86:LEU:H	1:A:257:GLN:HE22	1.52	0.56
1:B:86:LEU:H	1:B:257:GLN:HE22	1.54	0.54
1:A:186:ASN:O	1:A:189:GLN:HG2	2.09	0.53
1:B:217:ASP:OD1	1:B:255:ARG:NH1	2.28	0.53
1:A:388:HIS:HE1	4:A:627:HOH:O	1.91	0.52
1:A:162:PHE:HE1	1:A:215:LEU:HD21	1.74	0.52
1:A:169:GLN:HB3	1:A:170:PRO:HD2	1.90	0.52
1:B:388:HIS:HD2	1:B:391:LYS:NZ	2.08	0.52
1:A:13:GLU:HG2	4:A:803:HOH:O	2.11	0.50
1:A:316:MET:CE	1:A:377:GLU:HA	2.43	0.49
1:A:216:VAL:HG21	1:A:259[B]:ILE:HG13	1.95	0.48
1:A:79:ARG:HG3	1:A:83:GLY:O	2.13	0.48
1:B:73[B]:GLN:NE2	1:B:77:PHE:CE1	2.82	0.47
1:A:310[B]:GLN:NE2	1:B:369:ASP:OD2	2.49	0.46
1:A:367:TRP:HB2	1:A:371:VAL:HG12	1.97	0.46
1:A:104:LEU:N	1:A:105:PRO:CD	2.79	0.46
1:A:47:ARG:NE	1:A:73[A]:GLN:HG3	2.31	0.46
1:B:366:ILE:HD12	1:B:386:PRO:HG2	1.98	0.46
1:A:167:ARG:HD3	4:A:911:HOH:O	2.14	0.46
1:B:316:MET:HE1	1:B:380:GLU:HG3	1.97	0.45
1:B:71:LEU:HD22	1:B:90:TRP:CE2	2.52	0.44
1:A:212:MET:HE3	1:A:259[A]:ILE:HD11	2.00	0.44
1:A:388:HIS:HA	1:A:391:LYS:HD3	2.00	0.44
1:A:162:PHE:CE1	1:A:215:LEU:HD21	2.51	0.43
1:B:200:GLU:N	4:B:724:HOH:O	2.40	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:GLN:HG3	1:B:8:PRO:HD2	2.02	0.42
1:B:200:GLU:HB3	4:B:724:HOH:O	2.18	0.42
1:B:388:HIS:HD2	1:B:391:LYS:HZ3	1.67	0.42
3:B:502:1C6:H7	3:B:502:1C6:H10	2.01	0.41
1:A:195:ASP:OD1	1:A:196:PRO:HD2	2.20	0.41
1:B:162:PHE:HE1	1:B:215:LEU:CD2	2.13	0.41
1:B:75:LEU:HA	1:B:75:LEU:HD23	1.91	0.41
1:A:115:TYR:HA	1:A:118:MET:CE	2.51	0.40
1:B:253:ASN:O	1:B:257:GLN:HG2	2.22	0.40
1:B:381:ASN:OD1	1:B:383:SER:HB2	2.21	0.40
1:B:73[B]:GLN:HG3	1:B:77:PHE:CE2	2.56	0.40
1:B:14:LEU:O	1:B:15:LYS:HB2	2.21	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168[B]:ASP:OD2	1:B:168:ASP:N[2_454]	2.05	0.15

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	452/455 (99%)	441 (98%)	11 (2%)	0	100	100
1	B	444/455 (98%)	429 (97%)	14 (3%)	1 (0%)	51	34
All	All	896/910 (98%)	870 (97%)	25 (3%)	1 (0%)	55	38

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	15	LYS

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	398/400 (100%)	394 (99%)	4 (1%)	80	73
1	B	392/400 (98%)	384 (98%)	8 (2%)	60	45
All	All	790/800 (99%)	778 (98%)	12 (2%)	70	57

All (12) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	4	GLU
1	A	22	THR
1	A	148	LEU
1	A	440	LYS
1	B	4	GLU
1	B	5	MET
1	B	41	LYS
1	B	71	LEU
1	B	104	LEU
1	B	148	LEU
1	B	159	ASN
1	B	440	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	95	ASN
1	A	125	GLN
1	A	159	ASN
1	A	201	ASN
1	A	236	HIS
1	A	253	ASN
1	A	257	GLN
1	A	319	ASN
1	A	388	HIS
1	A	403	GLN

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Mol	Chain	Res	Type
1	B	16	ASN
1	B	95	ASN
1	B	125	GLN
1	B	159	ASN
1	B	201	ASN
1	B	236	HIS
1	B	253	ASN
1	B	257	GLN
1	B	283	ASN
1	B	388	HIS
1	B	403	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](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	HEM	A	501	1	28,50,50	2.16	10 (35%)	17,82,82	1.81	4 (23%)
3	1C6	A	502	-	25,25,25	2.66	8 (32%)	25,35,35	2.13	7 (28%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	B	501	1	28,50,50	1.91	9 (32%)	17,82,82	2.08	6 (35%)
3	1C6	B	502	-	25,25,25	3.27	9 (36%)	25,35,35	2.34	9 (36%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	A	501	1	-	0/6/54/54	0/0/8/8
3	1C6	A	502	-	-	0/7/9/9	0/3/3/3
2	HEM	B	501	1	-	0/6/54/54	0/0/8/8
3	1C6	B	502	-	-	0/7/9/9	0/3/3/3

All (36) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	1C6	CB-S	-5.45	1.71	1.82
2	A	501	HEM	C3B-C2B	-4.60	1.34	1.40
2	A	501	HEM	C3C-C2C	-3.54	1.35	1.40
3	A	502	1C6	CB-S	-3.45	1.75	1.82
3	B	502	1C6	CF1-S	-2.47	1.71	1.75
3	A	502	1C6	CF1-S	-2.35	1.71	1.75
2	B	501	HEM	C3B-C2B	-2.13	1.37	1.40
3	A	502	1C6	CE2-CZ	-2.12	1.36	1.39
2	A	501	HEM	C4C-NC	2.04	1.39	1.36
3	A	502	1C6	CE3-CZ3	2.04	1.40	1.37
2	B	501	HEM	CAD-C3D	2.12	1.56	1.52
2	B	501	HEM	CMB-C2B	2.19	1.56	1.51
2	B	501	HEM	C1B-NB	2.21	1.39	1.36
2	A	501	HEM	C4A-NA	2.36	1.41	1.36
2	B	501	HEM	C3B-CAB	2.43	1.52	1.47
2	B	501	HEM	C3C-CAC	2.45	1.52	1.47
2	B	501	HEM	CAA-C2A	2.47	1.56	1.52
3	B	502	1C6	CS2-CX2	2.61	1.50	1.42
3	B	502	1C6	CB-CG	2.74	1.53	1.50
3	B	502	1C6	CF1-NV	2.86	1.39	1.34
2	A	501	HEM	C3C-CAC	2.88	1.53	1.47
3	B	502	1C6	CE3-CZ3	2.90	1.42	1.37
2	A	501	HEM	C3B-CAB	3.21	1.54	1.47
2	A	501	HEM	CMB-C2B	3.23	1.58	1.51
3	A	502	1C6	CB-CG	3.36	1.54	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C4D-ND	3.47	1.40	1.36
2	B	501	HEM	C3D-C2D	3.88	1.49	1.37
3	A	502	1C6	CE1-CD1	3.92	1.46	1.39
2	A	501	HEM	C1C-NC	3.92	1.41	1.36
2	A	501	HEM	C3D-C2D	4.05	1.49	1.37
2	B	501	HEM	C1C-NC	4.28	1.41	1.36
3	A	502	1C6	CE1-CZ	4.45	1.48	1.40
3	B	502	1C6	CE1-CZ	4.64	1.48	1.40
3	B	502	1C6	CE1-CD1	6.45	1.51	1.39
3	A	502	1C6	CD1-CG	9.13	1.48	1.39
3	B	502	1C6	CD1-CG	11.29	1.50	1.39

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	1C6	CZ-CE1-CD1	-6.65	116.28	121.53
3	B	502	1C6	CZ-CE1-CD1	-5.42	117.25	121.53
2	A	501	HEM	CBD-CAD-C3D	-4.65	103.60	112.47
2	B	501	HEM	CBD-CAD-C3D	-4.58	103.73	112.47
3	B	502	1C6	CZ-CE2-N1	-3.54	122.24	125.22
3	B	502	1C6	C3-CD1-CG	-3.36	119.65	122.67
3	A	502	1C6	CB-S-CF1	-3.05	99.16	102.48
3	B	502	1C6	CD1-CG-N1	-2.96	121.60	124.08
2	B	501	HEM	C3B-C4B-NB	-2.89	105.47	109.21
3	A	502	1C6	CD1-CG-N1	-2.73	121.79	124.08
3	A	502	1C6	CZ3-CE3-CX2	-2.47	117.58	119.13
3	B	502	1C6	C1-CZ-CE2	-2.46	116.65	120.21
2	A	501	HEM	C3B-C4B-NB	-2.38	106.13	109.21
3	A	502	1C6	O3-CZ3-CE3	-2.29	118.48	124.49
2	A	501	HEM	C3C-C4C-NC	-2.23	106.73	110.94
2	B	501	HEM	CMA-C3A-C4A	-2.16	125.14	128.46
3	A	502	1C6	C1-CZ-CE2	-2.13	117.13	120.21
2	B	501	HEM	CAD-CBD-CGD	-2.13	109.03	112.66
3	B	502	1C6	C3-CD1-CE1	2.11	124.50	121.30
2	B	501	HEM	C4C-C3C-C2C	2.48	108.63	106.90
3	B	502	1C6	C4-O3-CZ3	2.83	123.70	117.50
2	B	501	HEM	CMB-C2B-C3B	3.14	130.72	124.89
3	A	502	1C6	CE2-N1-CG	3.27	121.53	117.88
2	A	501	HEM	C4C-C3C-C2C	3.38	109.26	106.90
3	B	502	1C6	O2-CE1-CD1	3.78	124.09	118.84
3	B	502	1C6	CE2-N1-CG	4.74	123.18	117.88

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	502	1C6	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	450/455 (98%)	-0.15	2 (0%) 92 91	13, 23, 40, 55	0
1	B	446/455 (98%)	-0.07	7 (1%) 72 71	15, 28, 56, 74	0
All	All	896/910 (98%)	-0.11	9 (1%) 82 82	13, 25, 51, 74	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	13	GLU	3.4
1	B	44	ALA	2.8
1	B	191	ALA	2.7
1	A	225	ALA	2.7
1	B	46	GLY	2.6
1	B	225	ALA	2.5
1	B	4	GLU	2.5
1	B	45	PRO	2.0
1	A	230	SER	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	1C6	B	502	23/23	0.90	0.13	2.25	29,33,39,43	0
2	HEM	B	501	43/43	0.96	0.10	0.41	14,20,23,26	0
2	HEM	A	501	43/43	0.98	0.10	0.15	10,14,18,20	0
3	1C6	A	502	23/23	0.96	0.09	-0.80	13,19,26,27	0

6.5 Other polymers

There are no such residues in this entry.