



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:03 PM GMT

PDB ID : 4KEY
Title : Structure of P450 BM3 A82F F87V in complex with omeprazole
Authors : Leys, D.
Deposited on : 2013-04-26
Resolution : 2.05 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

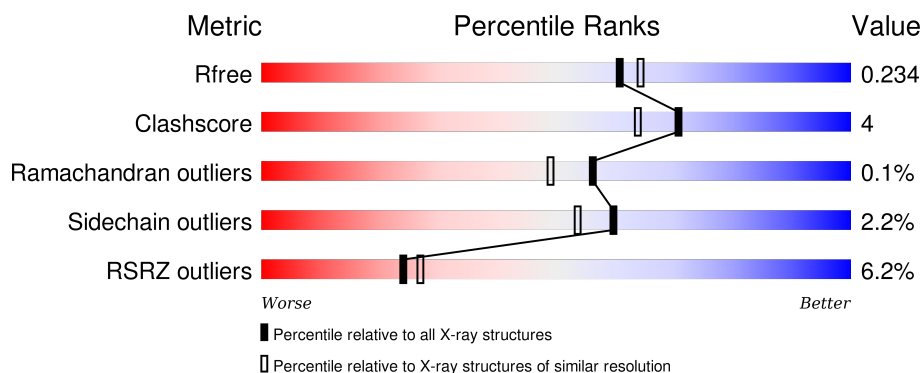
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.05 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}	91344	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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	 12% 88% 9% ..
1	B	455	 12% 82% 13% ..

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	A	502	-	-	-	X
3	1C6	B	502	-	-	-	X

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 7764 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	5	0
			3673	2344	623	689	17			
1	B	442	Total	C	N	O	S	0	0	0
			3553	2274	604	658	17			

There are 4 discrepancies between the modelled and reference sequences:

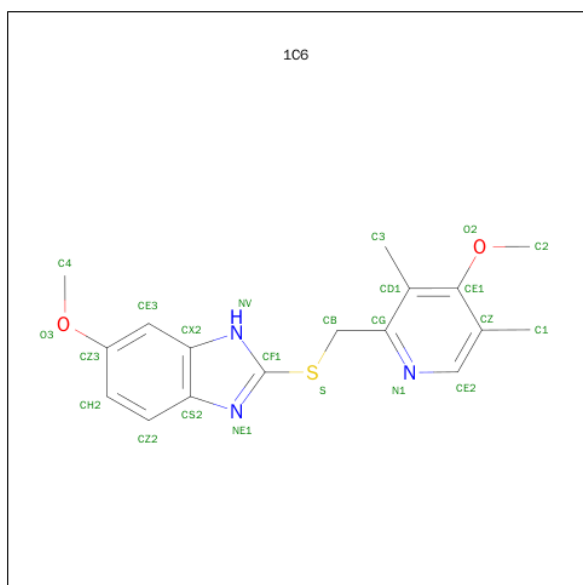
Chain	Residue	Modelled	Actual	Comment	Reference
A	82	PHE	ALA	ENGINEERED MUTATION	UNP P14779
A	87	VAL	PHE	ENGINEERED MUTATION	UNP P14779
B	82	PHE	ALA	ENGINEERED MUTATION	UNP P14779
B	87	VAL	PHE	ENGINEERED MUTATION	UNP P14779

- 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.

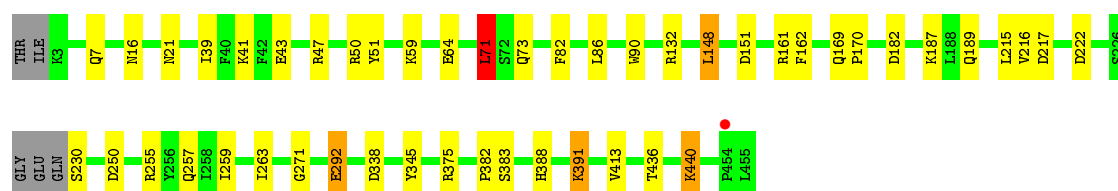
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	264	Total	O	0	0
			264	264		
4	B	142	Total	O	0	0
			142	142		

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.

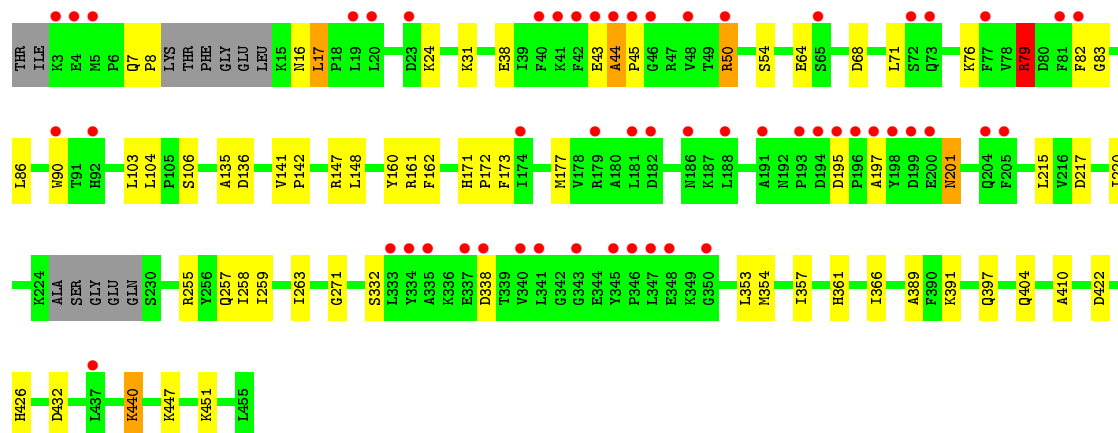
- Molecule 1: Bifunctional P-450/NADPH-P450 reductase

Chain A: 



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

Chain B: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.31Å 130.66Å 145.98Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.95 – 2.05 54.95 – 2.05	Depositor EDS
% Data completeness (in resolution range)	100.0 (54.95-2.05) 100.0 (54.95-2.05)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 2.05Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.187 , 0.235 0.191 , 0.234	Depositor DCC
R_{free} test set	3640 reflections (5.33%)	DCC
Wilson B-factor (Å ²)	32.4	Xtriage
Anisotropy	0.054	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 44.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 71956 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7764	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 7.52% 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 [i](#)

5.1 Standard geometry [i](#)

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	1.10	4/3757 (0.1%)	0.94	9/5080 (0.2%)
1	B	0.95	2/3635 (0.1%)	0.85	6/4918 (0.1%)
All	All	1.03	6/7392 (0.1%)	0.89	15/9998 (0.2%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	292	GLU	CG-CD	6.68	1.61	1.51
1	A	383	SER	CB-OG	-6.00	1.34	1.42
1	B	404	GLN	CG-CD	5.86	1.64	1.51
1	A	64	GLU	CD-OE2	-5.41	1.19	1.25
1	B	410	ALA	CA-CB	-5.33	1.41	1.52
1	A	345	TYR	CE1-CZ	-5.12	1.31	1.38

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	147	ARG	NE-CZ-NH1	8.08	124.34	120.30
1	A	132	ARG	NE-CZ-NH1	6.65	123.62	120.30
1	B	147	ARG	NE-CZ-NH2	-6.44	117.08	120.30
1	A	391	LYS	CD-CE-NZ	-5.83	98.28	111.70
1	A	383	SER	CB-CA-C	-5.82	99.05	110.10
1	A	71	LEU	CB-CG-CD1	5.72	120.72	111.00
1	B	50	ARG	NE-CZ-NH1	5.55	123.08	120.30
1	B	68	ASP	CB-CG-OD1	5.40	123.16	118.30
1	A	161	ARG	NE-CZ-NH2	-5.40	117.60	120.30
1	A	50	ARG	NE-CZ-NH2	-5.31	117.64	120.30
1	A	250	ASP	CB-CG-OD2	-5.31	113.52	118.30
1	A	375	ARG	NE-CZ-NH2	-5.29	117.65	120.30
1	B	391	LYS	CD-CE-NZ	-5.23	99.66	111.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	79	ARG	NE-CZ-NH1	5.22	122.91	120.30
1	A	132	ARG	NE-CZ-NH2	-5.12	117.74	120.30

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	3673	0	3622	26	0
1	B	3553	0	3504	37	0
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	264	0	0	3	0
4	B	142	0	0	2	0
All	All	7764	0	7224	64	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 4.

All (64) 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:16:ASN:HD22	1:A:43:GLU:H	1.23	0.84
1:A:7:GLN:HE21	1:A:16:ASN:HD21	1.38	0.71
1:B:16:ASN:HD22	1:B:43:GLU:H	1.38	0.71
1:B:50:ARG:HB2	1:B:353:LEU:HD23	1.72	0.70
1:A:47:ARG:NH2	1:A:73[A]:GLN:HG3	2.07	0.69
1:A:338:ASP:OD1	4:A:804:HOH:O	2.09	0.69
1:A:7:GLN:HE21	1:A:16:ASN:ND2	1.92	0.67
1:A:71:LEU:HD22	1:A:90:TRP:CE2	2.29	0.67
1:B:217:ASP:OD1	1:B:255:ARG:NH1	2.22	0.66
1:A:148:LEU:HD21	1:A:413:VAL:HG21	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:86:LEU:H	1:B:257:GLN:HE22	1.46	0.64
1:B:79:ARG:HD3	1:B:83:GLY:O	1.98	0.64
1:B:7:GLN:HG3	1:B:8:PRO:HD2	1.81	0.62
1:B:44:ALA:HB1	1:B:45:PRO:CD	2.28	0.62
1:A:217:ASP:OD1	1:A:255:ARG:NH1	2.30	0.61
1:B:82:PHE:HE1	1:B:263:ILE:HD12	1.66	0.61
1:A:162:PHE:CE1	1:A:215:LEU:HD21	2.35	0.61
1:A:47:ARG:CZ	1:A:73[A]:GLN:HG3	2.32	0.60
1:B:17:LEU:HD13	1:B:45:PRO:HD3	1.84	0.60
1:A:292:GLU:OE1	4:A:767:HOH:O	2.17	0.59
1:A:86:LEU:H	1:A:257:GLN:HE22	1.51	0.58
1:B:162:PHE:CE1	1:B:215:LEU:HD21	2.38	0.58
3:B:502:1C6:H10	4:B:732:HOH:O	2.03	0.57
1:B:271:GLY:HA2	1:B:440:LYS:HG3	1.86	0.57
1:A:388:HIS:HD2	1:A:391:LYS:NZ	2.03	0.56
1:B:201:ASN:N	1:B:201:ASN:HD22	2.04	0.55
1:A:82:PHE:HE1	1:A:263:ILE:HD12	1.71	0.55
1:B:332:SER:HB2	1:B:354:MET:SD	2.47	0.54
1:A:271:GLY:HA2	1:A:440:LYS:HG3	1.87	0.54
1:B:366:ILE:HG21	1:B:389:ALA:HB1	1.90	0.54
1:B:24:LYS:NZ	1:B:432:ASP:OD1	2.32	0.53
1:B:135:ALA:O	1:B:136:ASP:CB	2.54	0.53
1:B:422:ASP:OD2	1:B:451:LYS:NZ	2.38	0.53
1:B:82:PHE:CE1	1:B:263:ILE:HD12	2.46	0.51
1:B:162:PHE:HE1	1:B:215:LEU:HD21	1.75	0.51
1:B:177:MET:SD	1:B:263:ILE:HG23	2.52	0.50
1:A:182[A]:ASP:OD1	1:A:436:THR:HG22	2.11	0.50
1:B:201:ASN:H	1:B:201:ASN:HD22	1.60	0.49
1:B:141:VAL:HB	1:B:142:PRO:HD3	1.93	0.49
1:A:21:ASN:OD1	1:A:189:GLN:OE1	2.30	0.49
1:A:59:LYS:NZ	4:A:834:HOH:O	2.46	0.49
1:A:216:VAL:HG21	1:A:259:ILE:HG13	1.97	0.47
1:B:38:GLU:HB2	1:B:54:SER:HB3	1.97	0.47
1:B:103:LEU:O	1:B:106:SER:HB2	2.15	0.47
1:B:44:ALA:HB1	1:B:45:PRO:HD2	1.97	0.46
1:B:71:LEU:HD22	1:B:90:TRP:CE2	2.51	0.46
1:A:388:HIS:HA	1:A:391:LYS:HD3	1.97	0.45
1:A:71:LEU:HD22	1:A:90:TRP:NE1	2.32	0.45
1:A:39:ILE:HA	1:A:51:TYR:O	2.18	0.44
1:A:151:ASP:OD1	1:A:162:PHE:HB2	2.18	0.44
1:B:422:ASP:CG	1:B:451:LYS:NZ	2.71	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLN:HB3	1:A:170:PRO:HD2	2.01	0.43
1:B:357:ILE:HG22	1:B:361:HIS:CE1	2.53	0.43
1:B:220:ILE:HD11	1:B:258:ILE:HD11	2.02	0.41
1:B:161:ARG:HA	4:B:738:HOH:O	2.20	0.41
1:B:160:TYR:CE1	1:B:215:LEU:HD11	2.55	0.41
1:B:64:GLU:OE2	1:B:397:GLN:HG2	2.21	0.41
1:A:162:PHE:HE1	1:A:215:LEU:HD21	1.82	0.41
1:A:162:PHE:CZ	1:A:215:LEU:HD21	2.55	0.41
1:B:426:HIS:CD2	1:B:447:LYS:HE3	2.55	0.41
1:B:76:LYS:O	1:B:79:ARG:HB2	2.21	0.41
1:B:173:PHE:CD1	1:B:215:LEU:CD2	3.03	0.41
1:B:171:HIS:CG	1:B:172:PRO:HD2	2.55	0.41
1:B:195:ASP:OD1	1:B:197:ALA:HB3	2.20	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	451/455 (99%)	438 (97%)	13 (3%)	0	100	100
1	B	436/455 (96%)	424 (97%)	11 (2%)	1 (0%)	52	43
All	All	887/910 (98%)	862 (97%)	24 (3%)	1 (0%)	56	49

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	44	ALA

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	400/400 (100%)	392 (98%)	8 (2%)	63	58
1	B	383/400 (96%)	374 (98%)	9 (2%)	58	51
All	All	783/800 (98%)	766 (98%)	17 (2%)	60	53

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	41	LYS
1	A	71	LEU
1	A	148	LEU
1	A	187	LYS
1	A	222	ASP
1	A	230	SER
1	A	382	PRO
1	A	440	LYS
1	B	17	LEU
1	B	31	LYS
1	B	79	ARG
1	B	104	LEU
1	B	148	LEU
1	B	201	ASN
1	B	259	ILE
1	B	338	ASP
1	B	440	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	16	ASN
1	A	95	ASN
1	A	109	GLN
1	A	159	ASN
1	A	189	GLN

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Mol	Chain	Res	Type
1	A	236	HIS
1	A	253	ASN
1	A	257	GLN
1	A	283	ASN
1	A	319	ASN
1	A	388	HIS
1	A	403	GLN
1	B	16	ASN
1	B	95	ASN
1	B	109	GLN
1	B	125	GLN
1	B	134	ASN
1	B	159	ASN
1	B	201	ASN
1	B	253	ASN
1	B	257	GLN
1	B	310	GLN
1	B	388	HIS
1	B	403	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

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	30,50,50	2.41	9 (30%)	24,82,82	2.60	11 (45%)
3	1C6	A	502	-	25,25,25	3.27	9 (36%)	24,35,35	2.25	8 (33%)
2	HEM	B	501	1	30,50,50	2.06	7 (23%)	24,82,82	2.74	12 (50%)
3	1C6	B	502	-	25,25,25	3.80	10 (40%)	24,35,35	3.11	8 (33%)

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	-	0/10/54/54	0/0/8/8
3	1C6	A	502	-	-	0/7/9/9	0/3/3/3
2	HEM	B	501	1	-	0/10/54/54	0/0/8/8
3	1C6	B	502	-	-	0/7/9/9	0/3/3/3

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	501	HEM	C3B-C4B	-8.58	1.44	1.51
3	B	502	1C6	CB-S	-6.41	1.69	1.82
2	B	501	HEM	C3B-C4B	-5.02	1.47	1.51
2	B	501	HEM	C3D-C4D	-4.41	1.45	1.51
2	A	501	HEM	C2C-C1C	-4.34	1.44	1.52
3	A	502	1C6	CB-S	-4.21	1.73	1.82
2	B	501	HEM	C2C-C1C	-4.11	1.44	1.52
2	A	501	HEM	C3D-C4D	-3.85	1.46	1.51
3	B	502	1C6	CF1-S	-3.23	1.69	1.75
2	A	501	HEM	C2D-C3D	-2.96	1.45	1.54
3	A	502	1C6	CF1-S	-2.69	1.70	1.75
2	A	501	HEM	C2B-C1B	-2.38	1.44	1.51
2	B	501	HEM	C2D-C1D	-2.16	1.44	1.51
3	A	502	1C6	O2-CE1	-2.09	1.34	1.38
3	A	502	1C6	CS2-CX2	2.05	1.48	1.42
3	B	502	1C6	CF1-NV	2.14	1.39	1.34
2	A	501	HEM	CHD-C4C	2.19	1.41	1.36
3	B	502	1C6	CB-CG	2.29	1.53	1.50
3	B	502	1C6	CG-N1	2.34	1.37	1.34
2	A	501	HEM	FE-ND	2.37	2.10	1.97

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	502	1C6	CS2-CX2	2.54	1.50	1.42
2	B	501	HEM	FE-NB	2.63	2.11	1.97
2	A	501	HEM	FE-NC	2.83	2.07	1.95
2	B	501	HEM	FE-ND	3.09	2.13	1.97
3	A	502	1C6	CG-N1	3.22	1.38	1.34
3	A	502	1C6	CE3-CZ3	3.39	1.43	1.37
2	A	501	HEM	C1C-NC	3.70	1.40	1.36
2	B	501	HEM	FE-NC	3.73	2.10	1.95
3	B	502	1C6	CE3-CZ3	3.73	1.43	1.37
3	A	502	1C6	CE1-CD1	3.99	1.47	1.39
3	A	502	1C6	CE1-CZ	4.45	1.48	1.40
3	B	502	1C6	CE1-CZ	5.95	1.51	1.40
3	B	502	1C6	CE1-CD1	6.06	1.51	1.39
3	A	502	1C6	CD1-CG	12.86	1.50	1.39
3	B	502	1C6	CD1-CG	13.86	1.51	1.39

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	502	1C6	CD1-CG-N1	-8.61	119.71	124.18
3	B	502	1C6	CZ-CE1-CD1	-5.84	117.24	121.57
3	A	502	1C6	CD1-CG-N1	-5.18	121.50	124.18
3	A	502	1C6	CB-S-CF1	-4.90	97.34	102.57
2	B	501	HEM	CMA-C3A-C4A	-3.14	123.17	128.36
2	A	501	HEM	CBD-CAD-C3D	-2.80	105.40	113.55
2	B	501	HEM	C3B-C4B-NB	-2.74	106.38	111.63
2	A	501	HEM	CMA-C3A-C4A	-2.67	123.94	128.36
3	B	502	1C6	CZ-CE2-N1	-2.47	123.02	125.23
2	B	501	HEM	CAA-C2A-C1A	-2.29	124.52	127.01
3	A	502	1C6	CZ-CE1-CD1	-2.28	119.87	121.57
3	A	502	1C6	CZ-CE2-N1	-2.14	123.32	125.23
2	A	501	HEM	C3B-C4B-NB	-2.12	107.57	111.63
2	B	501	HEM	CBD-CAD-C3D	-2.09	107.48	113.55
3	A	502	1C6	CH2-CZ3-CE3	2.07	123.81	120.84
2	B	501	HEM	C4B-CHC-C1C	2.33	129.72	125.82
3	B	502	1C6	C4-O3-CZ3	2.41	123.15	117.51
3	B	502	1C6	CZ3-CE3-CX2	2.41	120.57	119.14
3	B	502	1C6	CE3-CX2-NV	2.52	138.18	130.70
2	B	501	HEM	C3B-C4B-CHC	2.62	126.86	123.16
2	A	501	HEM	C3B-C4B-CHC	2.80	127.10	123.16
3	A	502	1C6	CE2-N1-CG	2.88	121.25	117.83
2	A	501	HEM	C2C-C1C-CHC	2.91	128.11	123.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	502	1C6	C4-O3-CZ3	3.05	124.66	117.51
2	A	501	HEM	C2D-C3D-C4D	3.36	107.19	101.50
2	A	501	HEM	CMD-C2D-C3D	3.43	129.54	114.35
2	B	501	HEM	CMD-C2D-C3D	3.48	129.75	114.35
3	A	502	1C6	C3-CD1-CG	3.72	126.26	122.61
2	B	501	HEM	CAD-C3D-C4D	3.79	125.84	112.47
2	A	501	HEM	CAD-C3D-C2D	3.94	124.54	113.22
2	B	501	HEM	CAD-C3D-C2D	4.17	125.19	113.22
2	B	501	HEM	C2D-C3D-C4D	4.39	108.94	101.50
2	A	501	HEM	CAD-C3D-C4D	4.47	128.25	112.47
2	A	501	HEM	CMB-C2B-C3B	4.65	128.13	116.53
3	B	502	1C6	CE2-N1-CG	5.05	123.84	117.83
2	B	501	HEM	CMB-C2B-C3B	5.12	129.31	116.53
2	A	501	HEM	CMC-C2C-C3C	5.28	129.71	116.53
2	B	501	HEM	CMC-C2C-C3C	6.01	131.54	116.53
3	B	502	1C6	CB-S-CF1	7.46	110.53	102.57

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

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	450/455 (98%)	-0.10	1 (0%) 95 96	17, 27, 46, 63	0
1	B	442/455 (97%)	0.51	54 (12%) 5 6	22, 42, 81, 100	0
All	All	892/910 (98%)	0.20	55 (6%) 24 27	17, 33, 72, 100	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	46	GLY	6.9
1	B	193	PRO	5.7
1	B	81	PHE	5.5
1	B	194	ASP	5.3
1	B	3	LYS	5.2
1	B	191	ALA	5.0
1	B	19	LEU	4.7
1	B	188	LEU	4.7
1	B	198	TYR	4.6
1	B	199	ASP	4.6
1	B	44	ALA	4.1
1	B	181	LEU	4.1
1	B	205	PHE	3.9
1	B	197	ALA	3.4
1	B	437	LEU	3.3
1	B	45	PRO	3.3
1	B	4	GLU	3.3
1	B	347	LEU	3.3
1	B	174	ILE	3.2
1	B	340	VAL	3.2
1	B	338	ASP	3.1
1	B	82	PHE	3.1
1	B	343	GLY	3.1
1	B	42	PHE	2.9

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Mol	Chain	Res	Type	RSRZ
1	B	204	GLN	2.9
1	B	348	GLU	2.9
1	B	196	PRO	2.8
1	B	77	PHE	2.8
1	B	48	VAL	2.8
1	B	5	MET	2.8
1	B	333	LEU	2.8
1	B	346	PRO	2.7
1	B	345	TYR	2.7
1	B	65	SER	2.6
1	B	90	TRP	2.6
1	B	40	PHE	2.5
1	B	50	ARG	2.5
1	B	337	GLU	2.5
1	B	200	GLU	2.5
1	A	454	PRO	2.4
1	B	182	ASP	2.3
1	B	341	LEU	2.3
1	B	92	HIS	2.2
1	B	72	SER	2.2
1	B	43	GLU	2.2
1	B	41	LYS	2.2
1	B	186	ASN	2.2
1	B	350	GLY	2.2
1	B	20	LEU	2.2
1	B	73	GLN	2.1
1	B	179	ARG	2.1
1	B	195	ASP	2.1
1	B	23	ASP	2.1
1	B	335	ALA	2.1
1	B	334	TYR	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	A	502	23/23	0.91	0.15	3.51	28,33,40,42	0
3	1C6	B	502	23/23	0.78	0.34	2.57	65,76,83,84	0
2	HEM	B	501	43/43	0.96	0.11	-0.29	22,33,39,40	0
2	HEM	A	501	43/43	0.98	0.11	-0.74	12,17,20,22	0

6.5 Other polymers [i](#)

There are no such residues in this entry.