



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2025 – 12:36 AM EST

PDB ID : 9CJG
Title : P450-G9 from *Actinokineospora terrae*, a non-canonical, serine-ligated cytochrome P450 in the ligand-free, closed conformation
Authors : Ireland, K.A.; Davis, K.M.
Deposited on : 2024-07-06
Resolution : 1.85 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

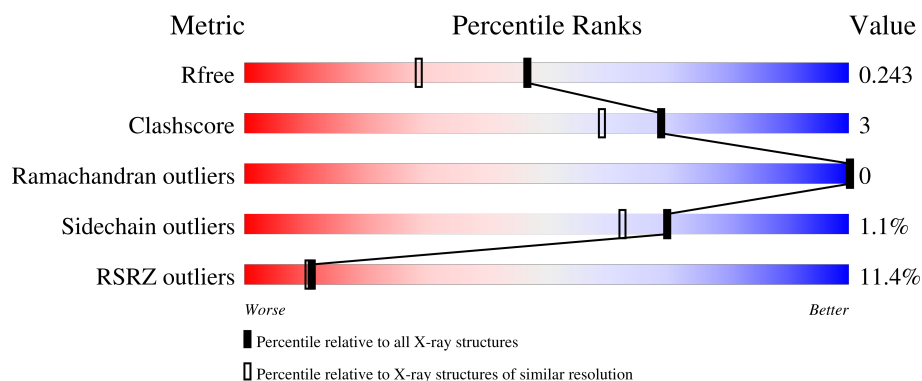
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.85 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}	164625	3097 (1.86-1.86)
Clashscore	180529	3359 (1.86-1.86)
Ramachandran outliers	177936	3335 (1.86-1.86)
Sidechain outliers	177891	3335 (1.86-1.86)
RSRZ outliers	164620	3097 (1.86-1.86)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	411	<div> <div>3%</div> <div>88%</div> <div>6%</div> <div>6%</div> </div>
1	B	411	<div> <div>18%</div> <div>85%</div> <div>9%</div> <div>7%</div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6282 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-G9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	388	Total	C	N	O	S	0	0	0
			2911	1853	527	529	2			
1	B	384	Total	C	N	O	S	0	0	0
			2780	1781	498	499	2			

There are 40 discrepancies between the modelled and reference sequences:

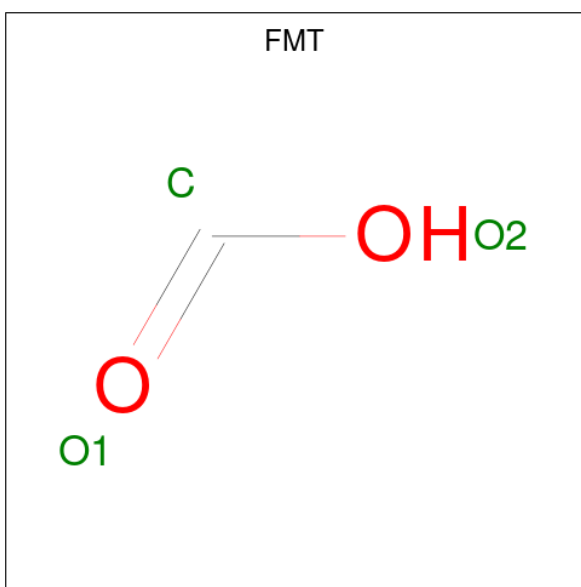
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A1H9KM06
A	-18	GLY	-	expression tag	UNP A0A1H9KM06
A	-17	SER	-	expression tag	UNP A0A1H9KM06
A	-16	SER	-	expression tag	UNP A0A1H9KM06
A	-15	HIS	-	expression tag	UNP A0A1H9KM06
A	-14	HIS	-	expression tag	UNP A0A1H9KM06
A	-13	HIS	-	expression tag	UNP A0A1H9KM06
A	-12	HIS	-	expression tag	UNP A0A1H9KM06
A	-11	HIS	-	expression tag	UNP A0A1H9KM06
A	-10	HIS	-	expression tag	UNP A0A1H9KM06
A	-9	SER	-	expression tag	UNP A0A1H9KM06
A	-8	SER	-	expression tag	UNP A0A1H9KM06
A	-7	GLY	-	expression tag	UNP A0A1H9KM06
A	-6	LEU	-	expression tag	UNP A0A1H9KM06
A	-5	VAL	-	expression tag	UNP A0A1H9KM06
A	-4	PRO	-	expression tag	UNP A0A1H9KM06
A	-3	ARG	-	expression tag	UNP A0A1H9KM06
A	-2	GLY	-	expression tag	UNP A0A1H9KM06
A	-1	SER	-	expression tag	UNP A0A1H9KM06
A	0	HIS	-	expression tag	UNP A0A1H9KM06
B	-19	MET	-	initiating methionine	UNP A0A1H9KM06
B	-18	GLY	-	expression tag	UNP A0A1H9KM06
B	-17	SER	-	expression tag	UNP A0A1H9KM06
B	-16	SER	-	expression tag	UNP A0A1H9KM06
B	-15	HIS	-	expression tag	UNP A0A1H9KM06

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP A0A1H9KM06
B	-13	HIS	-	expression tag	UNP A0A1H9KM06
B	-12	HIS	-	expression tag	UNP A0A1H9KM06
B	-11	HIS	-	expression tag	UNP A0A1H9KM06
B	-10	HIS	-	expression tag	UNP A0A1H9KM06
B	-9	SER	-	expression tag	UNP A0A1H9KM06
B	-8	SER	-	expression tag	UNP A0A1H9KM06
B	-7	GLY	-	expression tag	UNP A0A1H9KM06
B	-6	LEU	-	expression tag	UNP A0A1H9KM06
B	-5	VAL	-	expression tag	UNP A0A1H9KM06
B	-4	PRO	-	expression tag	UNP A0A1H9KM06
B	-3	ARG	-	expression tag	UNP A0A1H9KM06
B	-2	GLY	-	expression tag	UNP A0A1H9KM06
B	-1	SER	-	expression tag	UNP A0A1H9KM06
B	0	HIS	-	expression tag	UNP A0A1H9KM06

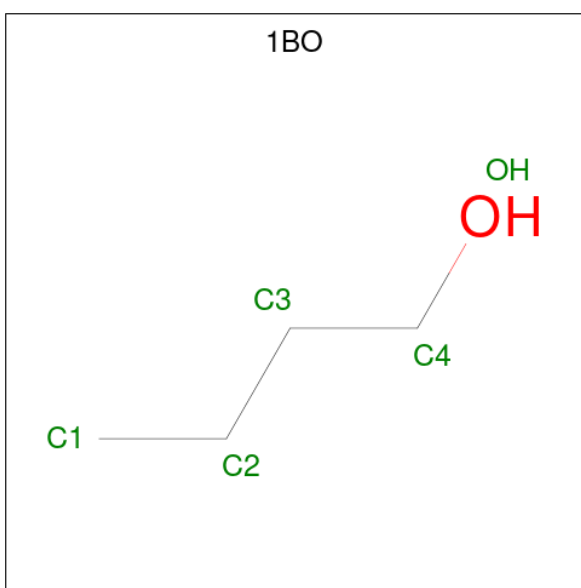
- # HEM

- Molecule 3 is FORMIC ACID (three-letter code: FMT) (formula: CH_2O_2).



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

- Molecule 4 is 1-BUTANOL (three-letter code: 1BO) (formula: $C_4H_{10}O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			5	4	1		
4	A	1	Total	C	O	0	0
			5	4	1		

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	342	Total 342	O 342	0	0
5	B	147	Total 147	O 147	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.13Å 59.65Å 233.80Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	57.80 – 1.85 57.80 – 1.85	Depositor EDS
% Data completeness (in resolution range)	99.3 (57.80-1.85) 99.3 (57.80-1.85)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 1.86Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.214 , 0.242 0.216 , 0.243	Depositor DCC
R_{free} test set	62612 reflections (3.12%)	wwPDB-VP
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.540	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 48.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6282	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.65% 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: FMT, IBO, 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	0.31	0/2982	0.54	0/4078
1	B	0.39	0/2847	0.59	0/3894
All	All	0.35	0/5829	0.56	0/7972

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	2911	0	2871	15	0
1	B	2780	0	2675	24	0
2	A	43	0	30	3	0
2	B	43	0	30	6	0
3	A	6	0	2	0	0
4	A	10	0	20	0	0
5	A	342	0	0	0	0
5	B	147	0	0	0	0
All	All	6282	0	5628	39	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 3.

All (39) 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:135:ALA:HA	1:B:236:VAL:HG22	1.84	0.58
1:A:135:ALA:HA	1:A:236:VAL:HG22	1.84	0.58
1:B:366:ALA:HB3	1:B:387:PRO:HB2	1.85	0.57
1:B:129:LEU:HD23	1:B:386:LEU:HD23	1.88	0.55
1:B:129:LEU:HD21	1:B:243:VAL:HG21	1.88	0.54
1:A:231:GLY:HA2	2:A:401:HEM:HAB	1.89	0.53
1:B:263:LEU:HB3	1:B:264:PRO:HD3	1.90	0.53
1:B:343:GLY:HA3	2:B:401:HEM:C3B	2.44	0.53
1:A:153:ARG:O	1:A:157:ARG:HG3	2.10	0.52
1:A:155:ARG:HG3	1:A:156:PHE:N	2.26	0.50
1:B:341:SER:HB2	2:B:401:HEM:ND	2.27	0.49
1:B:142:ILE:HD11	1:B:227:ALA:HB1	1.94	0.49
1:A:137:LEU:HB3	1:A:138:PRO:HD3	1.96	0.47
1:B:139:LEU:HD21	1:B:157:ARG:HG2	1.96	0.47
1:A:230:ALA:HB1	2:A:401:HEM:C4B	2.51	0.46
1:B:73:PRO:HD2	1:B:226:LEU:HD21	1.97	0.45
1:A:166:PRO:HB3	1:A:170:GLN:NE2	2.31	0.45
1:A:72:PRO:HA	1:A:73:PRO:HD3	1.80	0.44
1:B:343:GLY:HA3	2:B:401:HEM:C2B	2.53	0.44
1:B:139:LEU:CD2	1:B:157:ARG:HG2	2.48	0.44
1:B:226:LEU:HD23	1:B:226:LEU:HA	1.82	0.43
1:B:129:LEU:CD2	1:B:386:LEU:HD23	2.48	0.42
1:B:130:ILE:HA	1:B:134:ALA:HB3	2.01	0.42
1:A:56:SER:HB2	1:A:83:ASP:OD2	2.19	0.42
1:B:231:GLY:HA2	2:B:401:HEM:CBB	2.50	0.42
1:A:341:SER:HB2	2:A:401:HEM:ND	2.35	0.42
1:A:164:LEU:C	1:A:166:PRO:HD3	2.41	0.41
1:A:142:ILE:HD12	1:A:142:ILE:HA	1.92	0.41
1:B:25:GLY:HA3	1:B:42:ARG:HG3	2.02	0.41
1:B:277:LEU:HD23	2:B:401:HEM:HMD3	2.02	0.41
1:B:17:TYR:CG	1:B:308:SER:HB2	2.56	0.41
1:A:296:VAL:HA	1:A:297:PRO:HD3	1.97	0.41
1:B:164:LEU:C	1:B:166:PRO:HD3	2.41	0.41
1:A:202:LEU:HD12	1:A:202:LEU:HA	1.88	0.41
1:B:72:PRO:HA	1:B:73:PRO:HD3	1.85	0.41
1:B:90:LEU:HD23	1:B:90:LEU:HA	1.93	0.41
1:B:43:TYR:HA	1:B:309:ALA:HB1	2.02	0.40
1:B:230:ALA:HB1	2:B:401:HEM:NB	2.35	0.40
1:A:229:TRP:CE3	1:A:230:ALA:HB2	2.56	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	386/411 (94%)	374 (97%)	12 (3%)	0	100	100
1	B	381/411 (93%)	364 (96%)	17 (4%)	0	100	100
All	All	767/822 (93%)	738 (96%)	29 (4%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	286/320 (89%)	283 (99%)	3 (1%)	73	67
1	B	254/320 (79%)	251 (99%)	3 (1%)	67	59
All	All	540/640 (84%)	534 (99%)	6 (1%)	70	62

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	ASP
1	A	98	PHE
1	A	282	ARG
1	B	76	ASP
1	B	98	PHE

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Mol	Chain	Res	Type
1	B	282	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

6 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	B	401	1	42,50,50	1.32	6 (14%)	46,82,82	1.78	11 (23%)
3	FMT	A	405	-	2,2,2	0.74	0	1,1,1	0.26	0
3	FMT	A	402	-	2,2,2	0.73	0	1,1,1	0.21	0
4	1BO	A	404	-	4,4,4	0.12	0	3,3,3	0.09	0
2	HEM	A	401	1	42,50,50	1.30	5 (11%)	46,82,82	1.85	11 (23%)
4	1BO	A	403	-	4,4,4	0.35	0	3,3,3	0.45	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	401	1	-	2/12/54/54	-
4	1BO	A	403	-	-	0/2/2/2	-
2	HEM	B	401	1	-	1/12/54/54	-
4	1BO	A	404	-	-	2/2/2/2	-

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	401	HEM	C1B-NB	-3.41	1.34	1.40
2	A	401	HEM	C1B-NB	-3.30	1.34	1.40
2	B	401	HEM	C4D-ND	-3.19	1.34	1.40
2	A	401	HEM	C4D-ND	-3.18	1.34	1.40
2	B	401	HEM	C1D-ND	-2.38	1.34	1.38
2	A	401	HEM	FE-NB	2.38	2.11	1.98
2	B	401	HEM	FE-NB	2.35	2.11	1.98
2	B	401	HEM	CHB-C1B	2.32	1.40	1.34
2	A	401	HEM	C1D-ND	-2.25	1.34	1.38
2	A	401	HEM	CHB-C1B	2.21	1.40	1.34
2	B	401	HEM	C4B-NB	-2.19	1.34	1.38

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	HEM	CHC-C4B-NB	5.17	130.00	124.44
2	A	401	HEM	CHC-C4B-NB	5.03	129.85	124.44
2	A	401	HEM	CHD-C1D-ND	4.15	128.91	124.44
2	B	401	HEM	CHD-C1D-ND	3.96	128.70	124.44
2	A	401	HEM	CHA-C4D-ND	3.75	129.02	124.37
2	A	401	HEM	CHB-C1B-NB	3.60	128.84	124.37
2	B	401	HEM	CHA-C4D-ND	3.58	128.81	124.37
2	A	401	HEM	C1B-NB-C4B	3.25	109.05	105.21
2	B	401	HEM	CHB-C1B-NB	3.22	128.36	124.37
2	B	401	HEM	C1B-NB-C4B	3.16	108.95	105.21
2	A	401	HEM	CBA-CAA-C2A	-3.02	107.46	112.54
2	B	401	HEM	CBA-CAA-C2A	-2.91	107.65	112.54
2	A	401	HEM	C3B-C4B-NB	-2.52	107.66	109.47
2	A	401	HEM	CHD-C1D-C2D	-2.46	121.14	125.03
2	A	401	HEM	CHA-C4D-C3D	-2.45	120.70	125.23
2	B	401	HEM	CHD-C1D-C2D	-2.45	121.17	125.03
2	B	401	HEM	CHA-C4D-C3D	-2.26	121.06	125.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	401	HEM	C4B-C3B-C2B	-2.12	105.33	107.28
2	B	401	HEM	O2A-CGA-CBA	2.06	120.51	114.00
2	B	401	HEM	CAD-CBD-CGD	-2.05	108.23	113.67
2	A	401	HEM	CAD-CBD-CGD	-2.04	108.26	113.67
2	A	401	HEM	O2A-CGA-CBA	2.02	120.40	114.00

There are no chirality outliers.

All (5) torsion outliers are listed below:

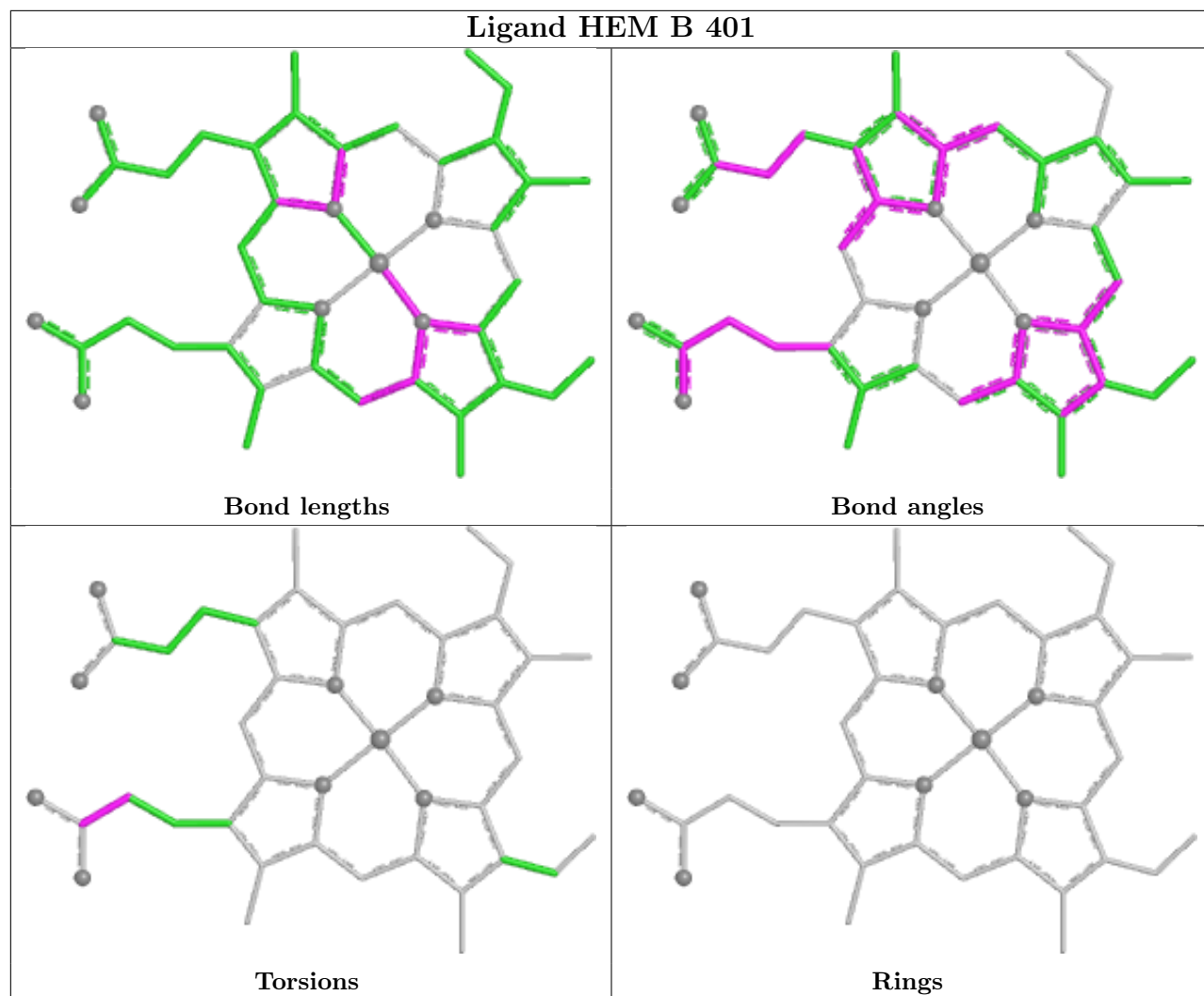
Mol	Chain	Res	Type	Atoms
4	A	404	1BO	C1-C2-C3-C4
2	A	401	HEM	C2B-C3B-CAB-CBB
4	A	404	1BO	C2-C3-C4-OH
2	A	401	HEM	C4B-C3B-CAB-CBB
2	B	401	HEM	CAA-CBA-CGA-O1A

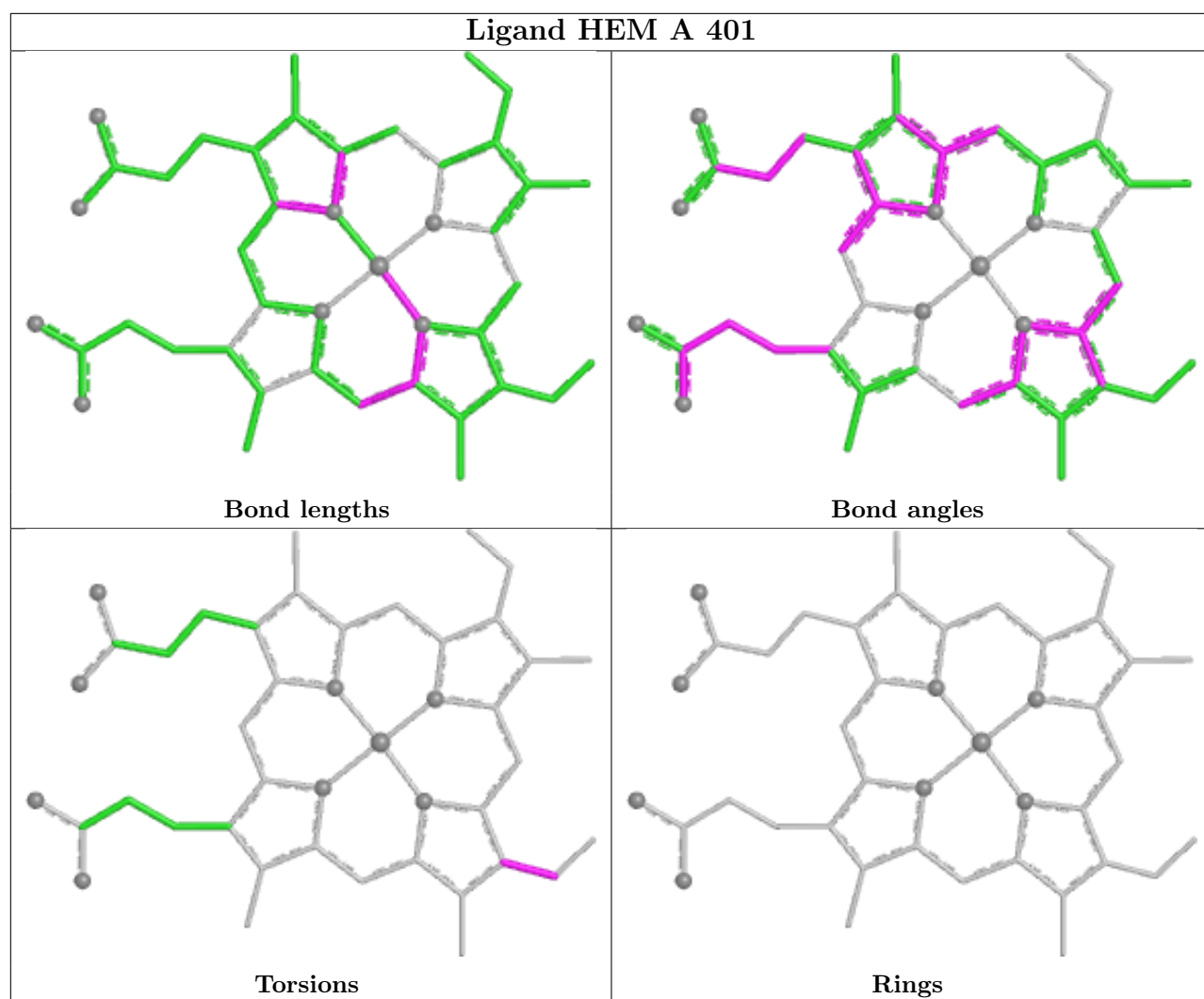
There are no ring outliers.

2 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	401	HEM	6	0
2	A	401	HEM	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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	388/411 (94%)	0.04	13 (3%)	48	51	10, 22, 40, 66	0
1	B	384/411 (93%)	1.11	75 (19%)	4	3	18, 36, 57, 75	0
All	All	772/822 (93%)	0.57	88 (11%)	11	10	10, 28, 54, 75	0

All (88) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	328	PRO	4.8
1	B	168	PRO	4.8
1	B	210	ASP	4.1
1	B	14	TYR	3.7
1	B	256	LEU	3.7
1	B	327	ASP	3.5
1	B	99	THR	3.5
1	A	391	GLY	3.5
1	B	100	PRO	3.4
1	A	5	PHE	3.4
1	B	330	PRO	3.3
1	B	370	ALA	3.3
1	B	19	ARG	3.3
1	B	23	GLY	3.3
1	B	315	VAL	3.2
1	B	262	LEU	3.2
1	B	261	ALA	3.2
1	B	263	LEU	3.2
1	B	312	ASP	3.2
1	B	16	ILE	3.2
1	B	254	ALA	3.2
1	A	75	LEU	3.1
1	B	26	ILE	3.1
1	B	17	TYR	3.1

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Mol	Chain	Res	Type	RSRZ
1	B	8	ALA	3.1
1	A	169	ASP	2.9
1	B	18	ALA	2.8
1	B	374	TRP	2.8
1	A	4	LEU	2.7
1	B	270	LEU	2.7
1	B	40	VAL	2.7
1	B	259	ASP	2.7
1	B	293	GLY	2.7
1	B	322	LEU	2.7
1	B	9	PHE	2.6
1	B	39	VAL	2.6
1	B	294	VAL	2.6
1	B	258	SER	2.6
1	B	365	ALA	2.6
1	B	159	TRP	2.6
1	B	291	ILE	2.6
1	B	20	LEU	2.6
1	B	56	SER	2.5
1	B	253	LEU	2.5
1	B	15	PRO	2.5
1	B	255	ASP	2.5
1	B	321	THR	2.5
1	B	27	ARG	2.5
1	B	356	SER	2.5
1	B	324	LEU	2.5
1	A	168	PRO	2.4
1	B	314	THR	2.4
1	B	332	LEU	2.4
1	B	386	LEU	2.4
1	A	195	ALA	2.4
1	B	22	ALA	2.4
1	B	319	PRO	2.4
1	B	98	PHE	2.4
1	A	211	ASP	2.3
1	B	104	HIS	2.3
1	B	238	LEU	2.3
1	B	75	LEU	2.3
1	B	318	LEU	2.3
1	B	264	PRO	2.2
1	A	212	GLU	2.2
1	B	11	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	273	PHE	2.2
1	A	124	GLY	2.2
1	B	212	GLU	2.2
1	A	159	TRP	2.2
1	B	389	LEU	2.2
1	B	103	VAL	2.1
1	A	199	HIS	2.1
1	B	292	ALA	2.1
1	B	316	PHE	2.1
1	A	155	ARG	2.1
1	B	360	ALA	2.1
1	B	31	THR	2.1
1	B	313	GLU	2.1
1	B	124	GLY	2.1
1	B	391	GLY	2.1
1	B	33	ASP	2.0
1	B	208	VAL	2.0
1	B	367	VAL	2.0
1	B	169	ASP	2.0
1	B	317	ASP	2.0
1	B	13	PRO	2.0
1	B	289	LEU	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 monosaccharides 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. 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.

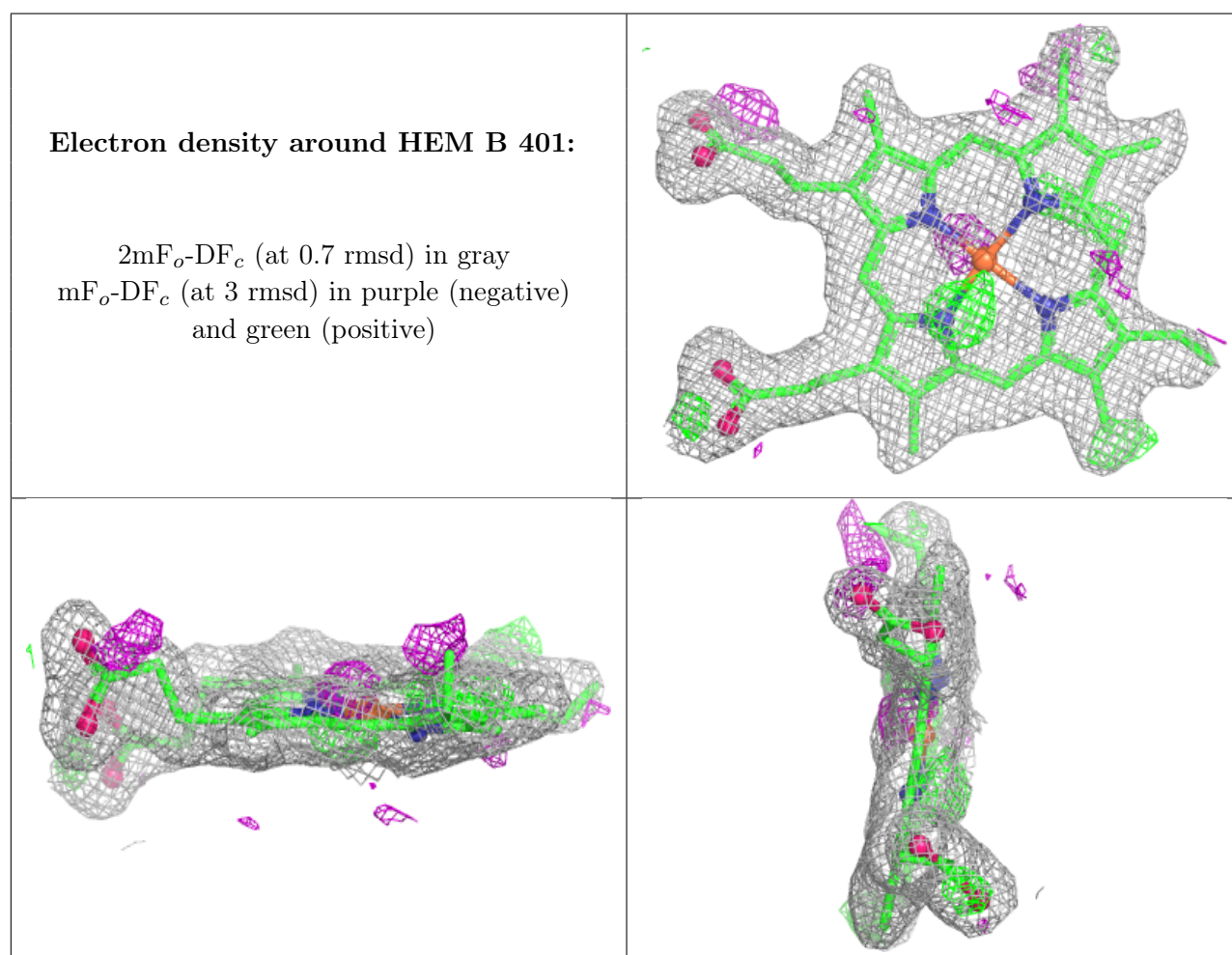
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	1BO	A	404	5/5	0.77	0.15	28,29,40,47	0

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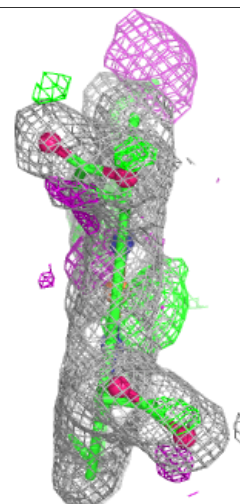
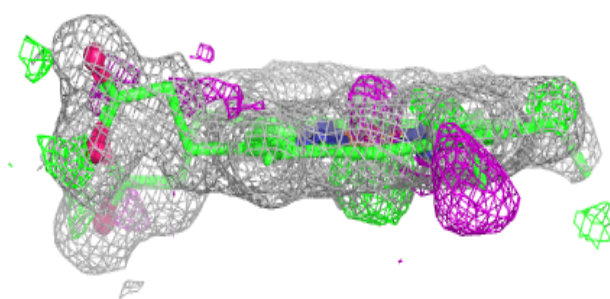
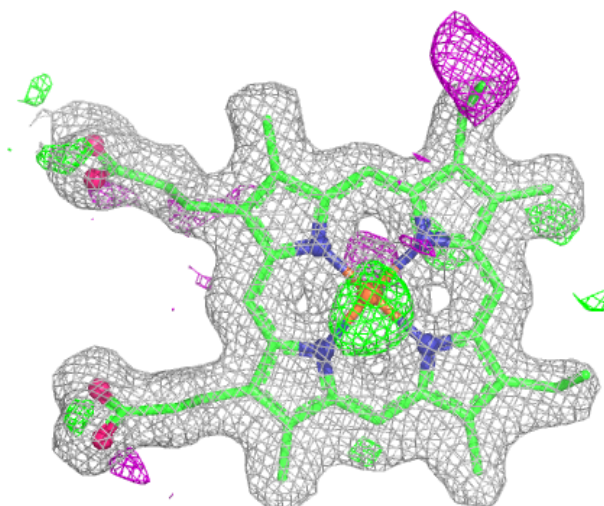
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	FMT	A	405	3/3	0.79	0.16	34,34,34,35	0
4	1BO	A	403	5/5	0.80	0.14	26,31,35,38	0
3	FMT	A	402	3/3	0.90	0.10	23,23,30,37	0
2	HEM	B	401	43/43	0.91	0.13	16,28,34,44	0
2	HEM	A	401	43/43	0.94	0.11	7,16,23,32	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around HEM A 401:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers ⓘ

There are no such residues in this entry.