



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 5, 2026 – 02:22 PM UTC

PDB ID : 7SNM / pdb_00007snm
Title : Lanosterol-bound P450 domain of the CYP51-ferredoxin fusion protein from *Methylococcus capsulatus*
Authors : Lepesheva, G.I.; Hargrove, T.; Wawrzak, Z.
Deposited on : 2021-10-28
Resolution : 2.55 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

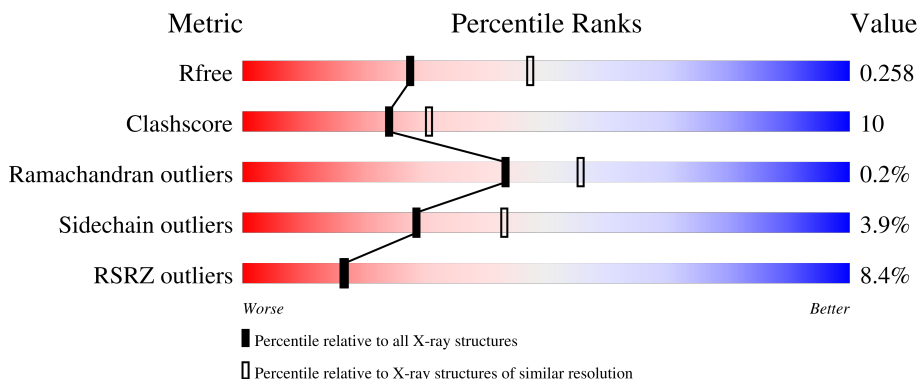
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.55 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}	180053	1091 (2.54-2.54)
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	551	
1	B	551	
1	C	551	
1	D	551	

2 Entry composition [i](#)

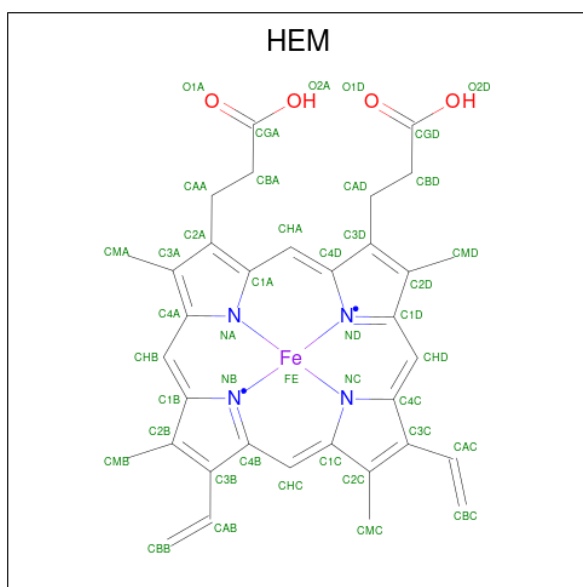
There are 4 unique types of molecules in this entry. The entry contains 14767 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 51.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	443	Total	C	N	O	S	0	0	0
			3574	2278	638	635	23			
1	B	443	Total	C	N	O	S	0	0	0
			3568	2275	635	635	23			
1	C	440	Total	C	N	O	S	0	0	0
			3530	2253	624	630	23			
1	D	439	Total	C	N	O	S	0	0	0
			3458	2212	602	622	22			

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



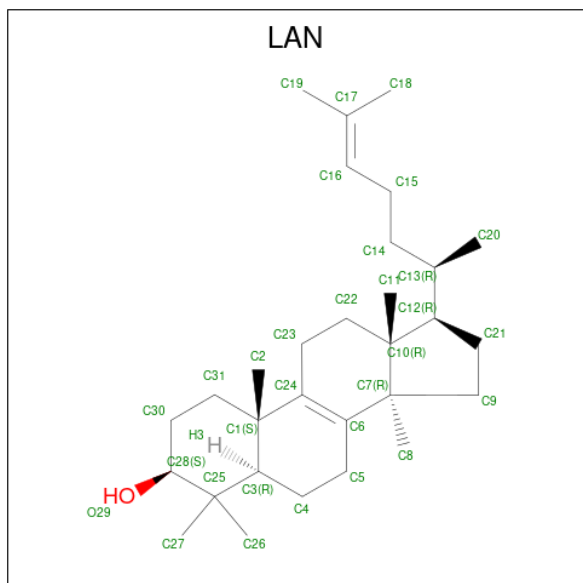
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		
2	B	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		

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

- Molecule 3 is LANOSTEROL (CCD ID: LAN) (formula: $C_{30}H_{50}O$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			31	30	1		
3	B	1	Total	C	O	0	0
			31	30	1		
3	C	1	Total	C	O	0	0
			31	30	1		
3	D	1	Total	C	O	0	0
			31	30	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	103	Total	O	0	0
			103	103		
4	B	95	Total	O	0	0
			95	95		
4	C	66	Total	O	0	0
			66	66		

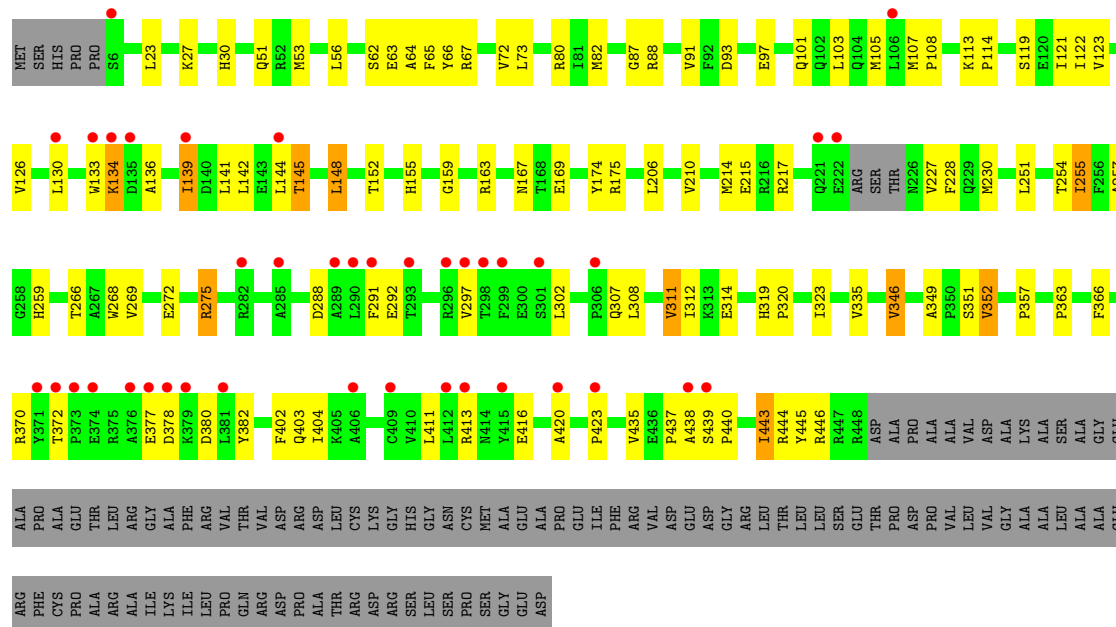
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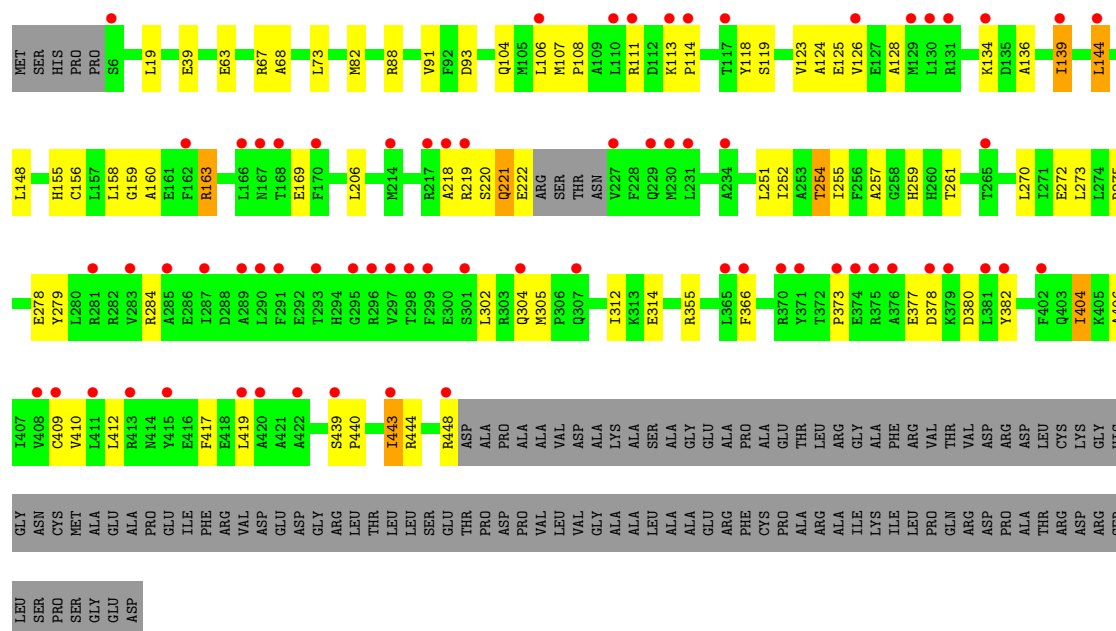
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	77	Total	O	0	0
			77	77		

ARG	VAL	ASP	GLU	ASP	GLY	ARG	LEU	THR	LEU	SER	LEU	VAL	GLY	ALA	ALA	LEU	ALA	ALA	GLU	ARG	PHE	CYS	PRO	ALA	ARG	ALA	ILE	LYS	ILE	LEU	PRO	GLN	ARG	ASP	PRO	ALA	THR	ARG	ASP	ARG	SER	LEU	SER	PRO	SER	GLY	GLU	ASP
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● Molecule 1: Cytochrome P450 51



● Molecule 1: Cytochrome P450 51



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	73.21Å 206.16Å 81.43Å 90.00° 114.09° 90.00°	Depositor
Resolution (Å)	74.45 – 2.55 74.45 – 2.55	Depositor EDS
% Data completeness (in resolution range)	99.0 (74.45-2.55) 99.0 (74.45-2.55)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.80 (at 2.55Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.204 , 0.246 0.208 , 0.258	Depositor DCC
R_{free} test set	3792 reflections (5.30%)	wwPDB-VP
Wilson B-factor (Å ²)	30.1	Xtriage
Anisotropy	0.312	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 60.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.042 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14767	wwPDB-VP
Average B, all atoms (Å ²)	59.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.76% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: LAN, 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	0.96	0/3657	1.31	4/4944 (0.1%)
1	B	0.96	0/3651	1.32	6/4937 (0.1%)
1	C	0.95	0/3612	1.31	3/4887 (0.1%)
1	D	0.96	0/3539	1.31	2/4795 (0.0%)
All	All	0.96	0/14459	1.31	15/19563 (0.1%)

There are no bond length outliers.

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	222	GLU	N-CA-C	-7.55	101.96	112.45
1	C	438	ALA	O-C-N	6.37	130.21	122.75
1	A	87	GLY	CA-C-O	-6.18	118.20	122.22
1	B	225	THR	CB-CA-C	-5.97	99.15	111.22
1	B	159	GLY	CA-C-O	-5.62	118.41	122.45

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	3574	0	3584	85	0
1	B	3568	0	3573	75	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3530	0	3514	74	0
1	D	3458	0	3396	49	0
2	A	43	0	30	2	0
2	B	43	0	30	4	0
2	C	43	0	30	1	0
2	D	43	0	30	2	0
3	A	31	0	50	2	0
3	B	31	0	50	5	0
3	C	31	0	50	4	0
3	D	31	0	50	3	0
4	A	103	0	0	1	0
4	B	95	0	0	0	0
4	C	66	0	0	0	0
4	D	77	0	0	0	0
All	All	14767	0	14387	288	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 10.

The worst 5 of 288 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:117:THR:HG23	1:B:373:PRO:HB3	1.43	1.00
1:A:370:ARG:HH22	1:A:384:TRP:NE1	1.69	0.89
1:B:225:THR:HG22	1:B:226:ASN:H	1.43	0.82
1:A:117:THR:CG2	1:B:373:PRO:HB3	2.11	0.81
1:B:149:THR:HA	1:B:152:THR:HG22	1.68	0.75

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	441/551 (80%)	412 (93%)	28 (6%)	1 (0%)	43	56
1	B	441/551 (80%)	415 (94%)	25 (6%)	1 (0%)	43	56
1	C	436/551 (79%)	409 (94%)	26 (6%)	1 (0%)	43	56
1	D	435/551 (79%)	412 (95%)	22 (5%)	1 (0%)	43	56
All	All	1753/2204 (80%)	1648 (94%)	101 (6%)	4 (0%)	43	56

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	91	VAL
1	B	91	VAL
1	C	91	VAL
1	D	91	VAL

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	381/466 (82%)	365 (96%)	16 (4%)	26	40
1	B	380/466 (82%)	370 (97%)	10 (3%)	40	59
1	C	374/466 (80%)	354 (95%)	20 (5%)	20	31
1	D	360/466 (77%)	348 (97%)	12 (3%)	33	50
All	All	1495/1864 (80%)	1437 (96%)	58 (4%)	28	43

5 of 58 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	119	SER
1	D	373	PRO
1	C	167	ASN
1	D	355	ARG
1	D	169	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such

sidechains are listed below:

Mol	Chain	Res	Type
1	D	336	GLN
1	C	164	HIS
1	B	221	GLN
1	B	75	GLN
1	C	30	HIS

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](#)

8 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	C	601	1	50,50,50	1.58	8 (16%)	67,82,82	1.75	15 (22%)
2	HEM	A	601	1	50,50,50	1.58	8 (16%)	67,82,82	1.68	13 (19%)
2	HEM	D	601	1	50,50,50	1.59	8 (16%)	67,82,82	1.76	14 (20%)
3	LAN	C	602	-	34,34,34	0.40	0	56,56,56	0.64	0
3	LAN	A	602	-	34,34,34	0.39	0	56,56,56	0.67	0
3	LAN	D	602	-	34,34,34	0.39	0	56,56,56	0.64	0
3	LAN	B	602	-	34,34,34	0.40	0	56,56,56	0.63	1 (1%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEM	B	601	1	50,50,50	1.57	8 (16%)	67,82,82	1.72	17 (25%)

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	C	601	1	-	3/14/54/54	-
2	HEM	A	601	1	-	3/14/54/54	-
2	HEM	D	601	1	-	1/14/54/54	-
3	LAN	C	602	-	-	0/10/82/82	0/4/4/4
3	LAN	A	602	-	-	3/10/82/82	0/4/4/4
3	LAN	D	602	-	-	0/10/82/82	0/4/4/4
3	LAN	B	602	-	-	1/10/82/82	0/4/4/4
2	HEM	B	601	1	-	2/14/54/54	-

The worst 5 of 32 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	601	HEM	FE-NB	5.10	2.10	1.94
2	A	601	HEM	FE-NB	5.05	2.10	1.94
2	D	601	HEM	FE-NB	5.01	2.10	1.94
2	C	601	HEM	FE-NB	5.00	2.10	1.94
2	C	601	HEM	FE-NC	4.26	2.09	1.95

The worst 5 of 60 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	601	HEM	CHC-C4B-NB	5.88	130.75	124.42
2	D	601	HEM	CHC-C4B-NB	5.43	130.27	124.42
2	C	601	HEM	CHC-C4B-NB	5.37	130.20	124.42
2	A	601	HEM	CHC-C4B-NB	5.27	130.09	124.42
2	D	601	HEM	CHD-C1D-ND	5.15	129.97	124.42

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	602	LAN	C15-C16-C17-C19

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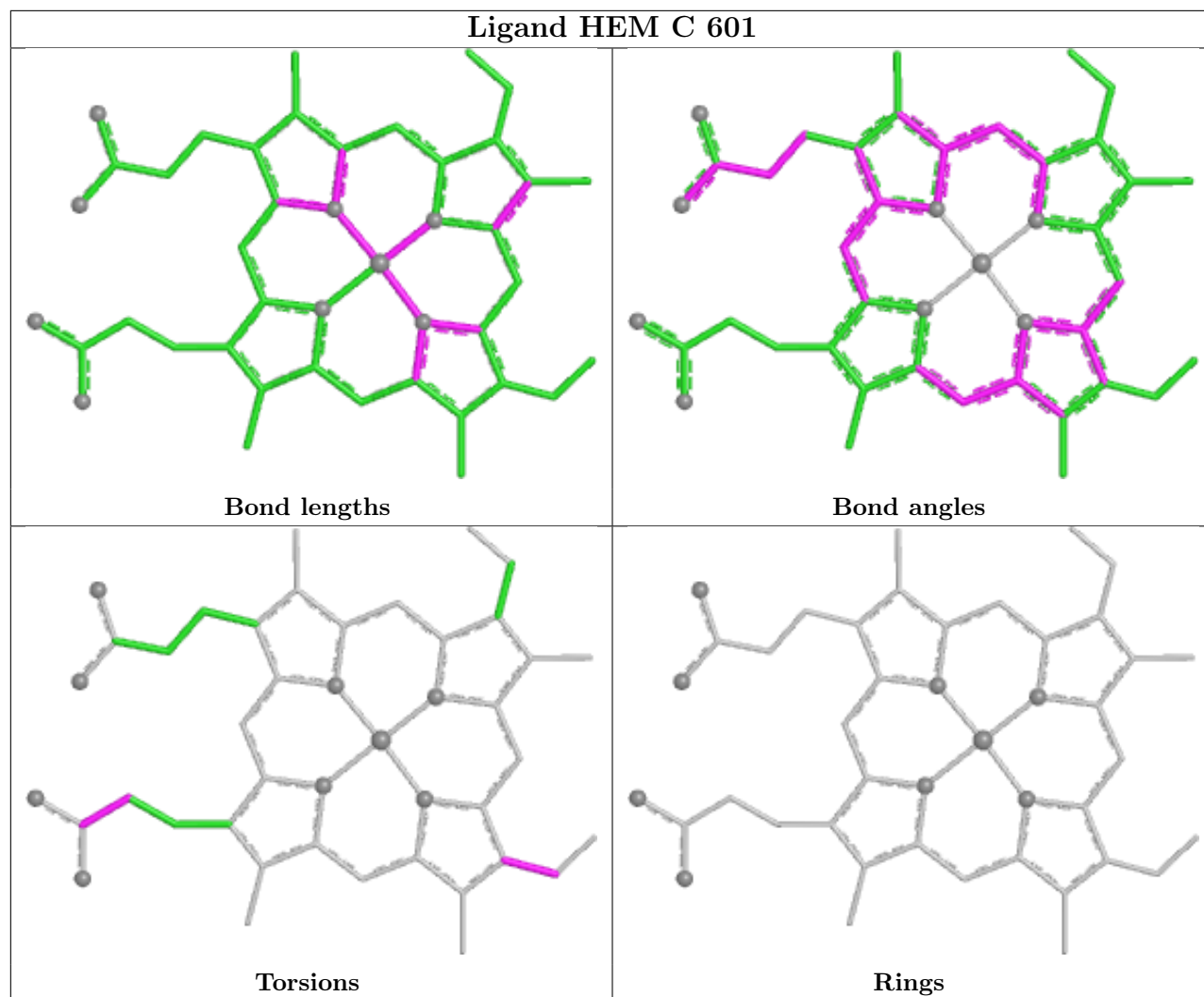
Mol	Chain	Res	Type	Atoms
3	A	602	LAN	C15-C16-C17-C18
3	B	602	LAN	C20-C13-C14-C15
3	A	602	LAN	C20-C13-C14-C15
2	B	601	HEM	CAD-CBD-CGD-O1D

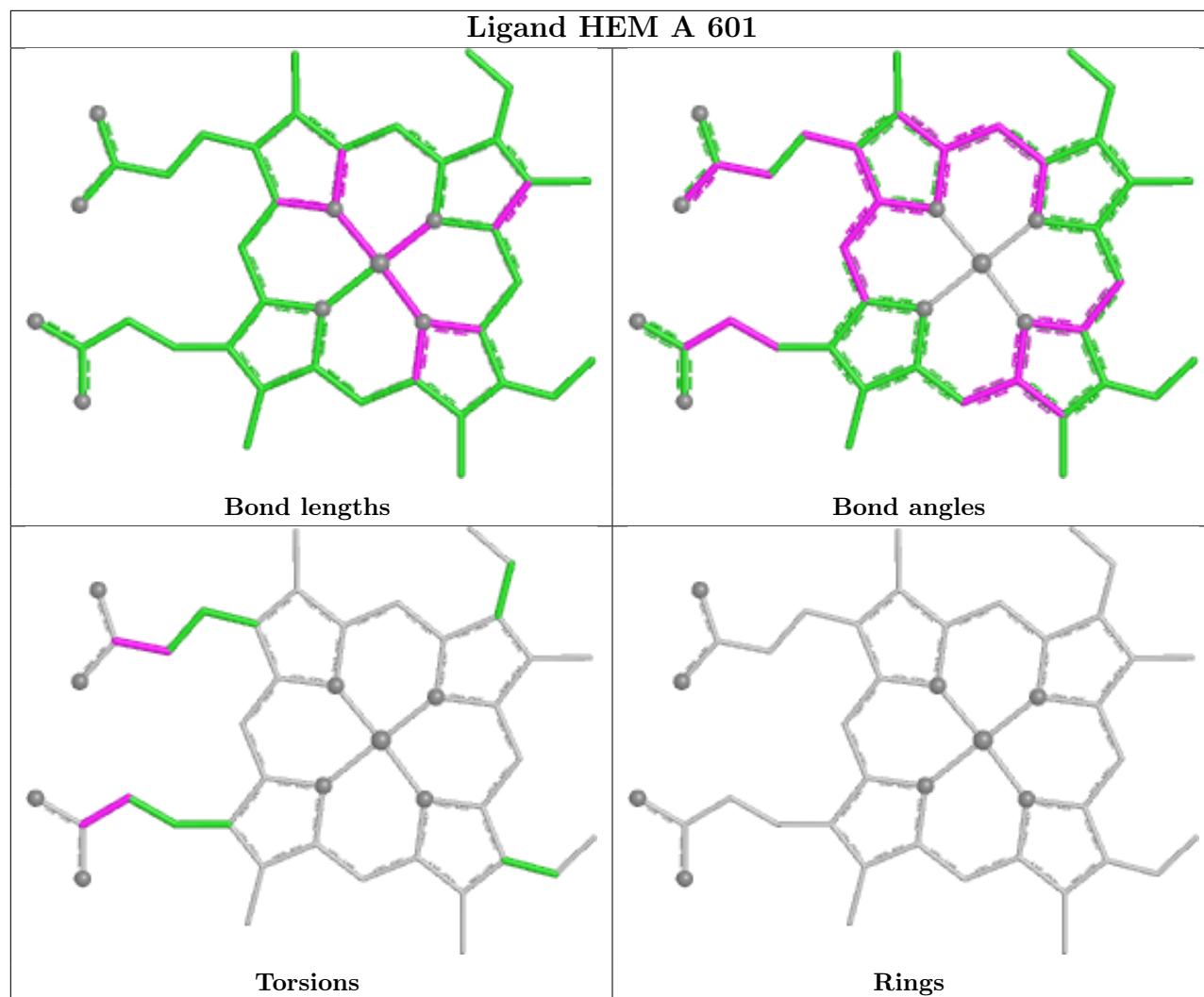
There are no ring outliers.

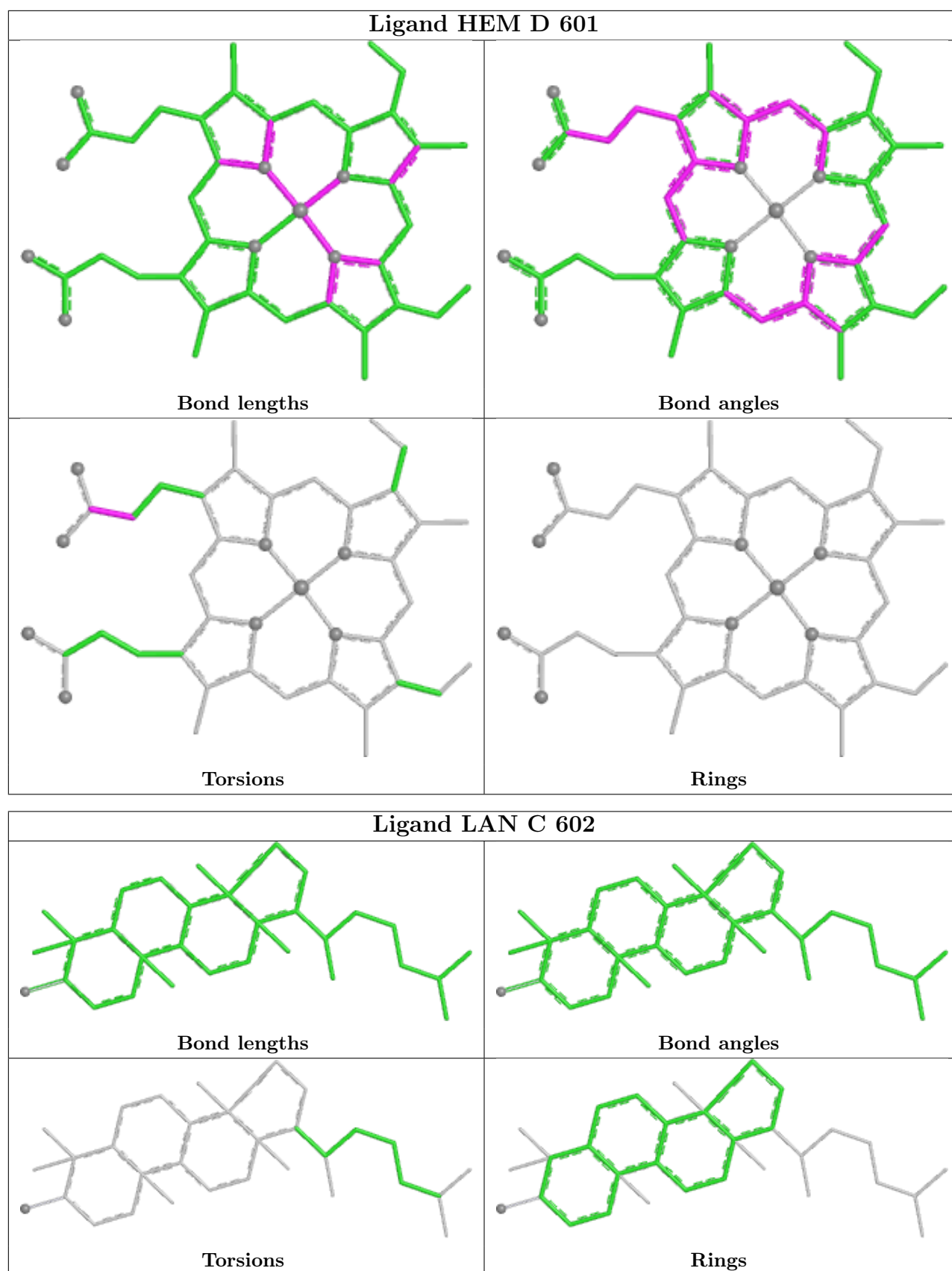
8 monomers are involved in 22 short contacts:

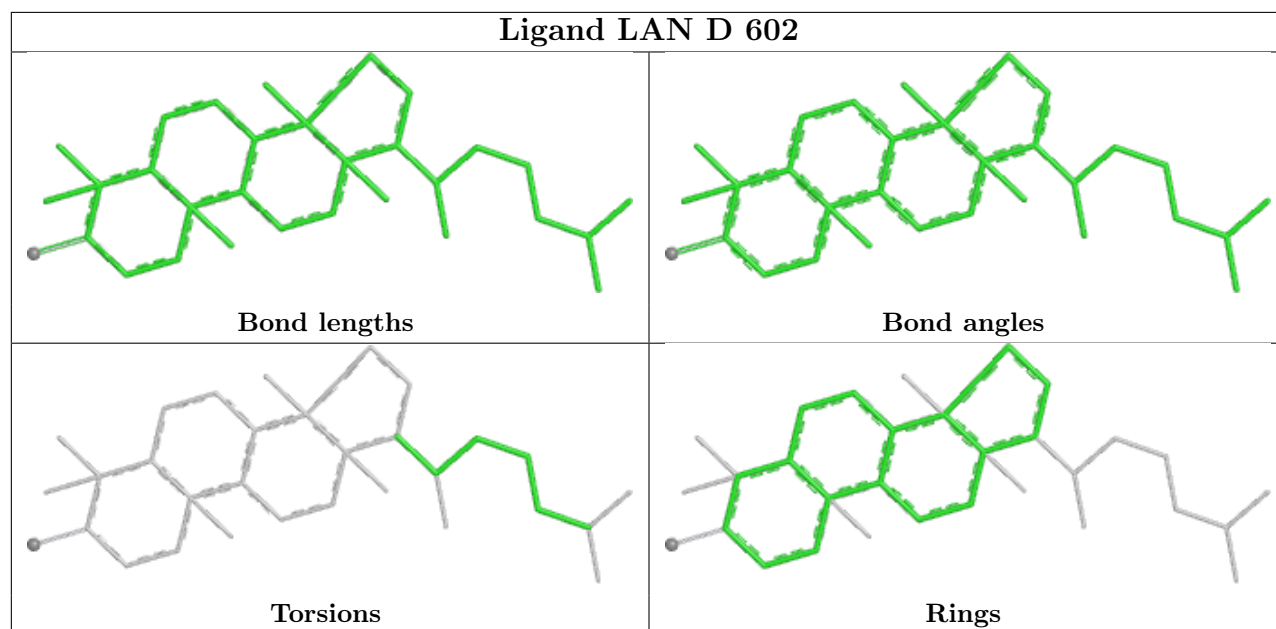
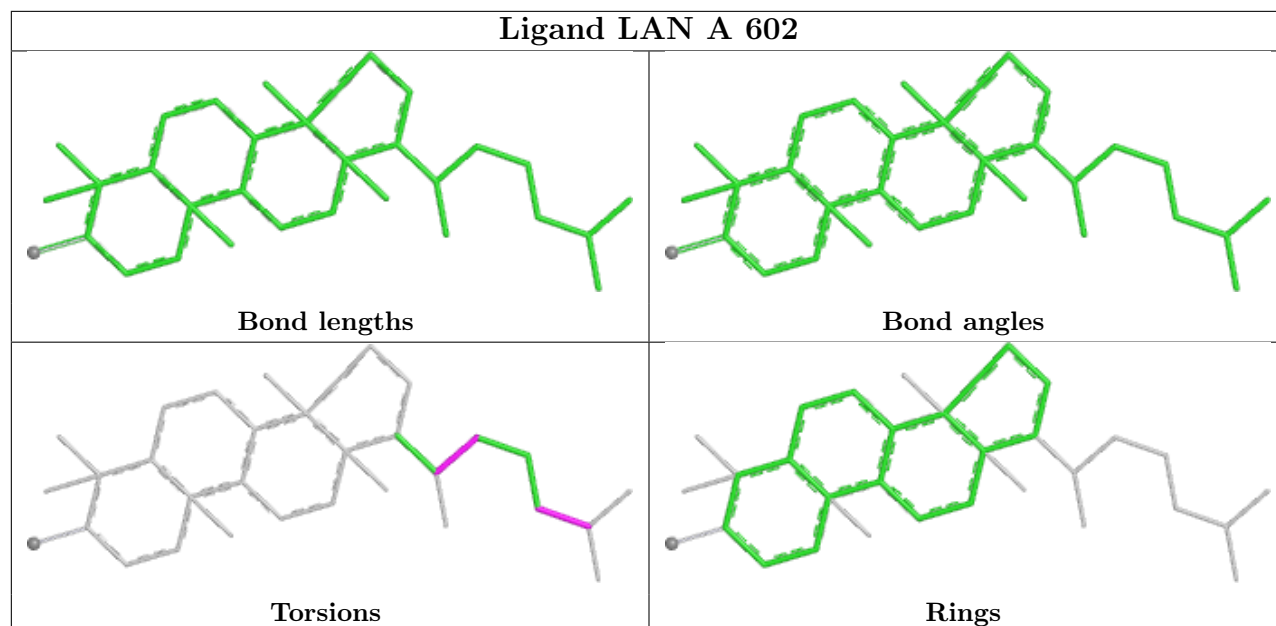
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	601	HEM	1	0
2	A	601	HEM	2	0
2	D	601	HEM	2	0
3	C	602	LAN	4	0
3	A	602	LAN	2	0
3	D	602	LAN	3	0
3	B	602	LAN	5	0
2	B	601	HEM	4	0

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.

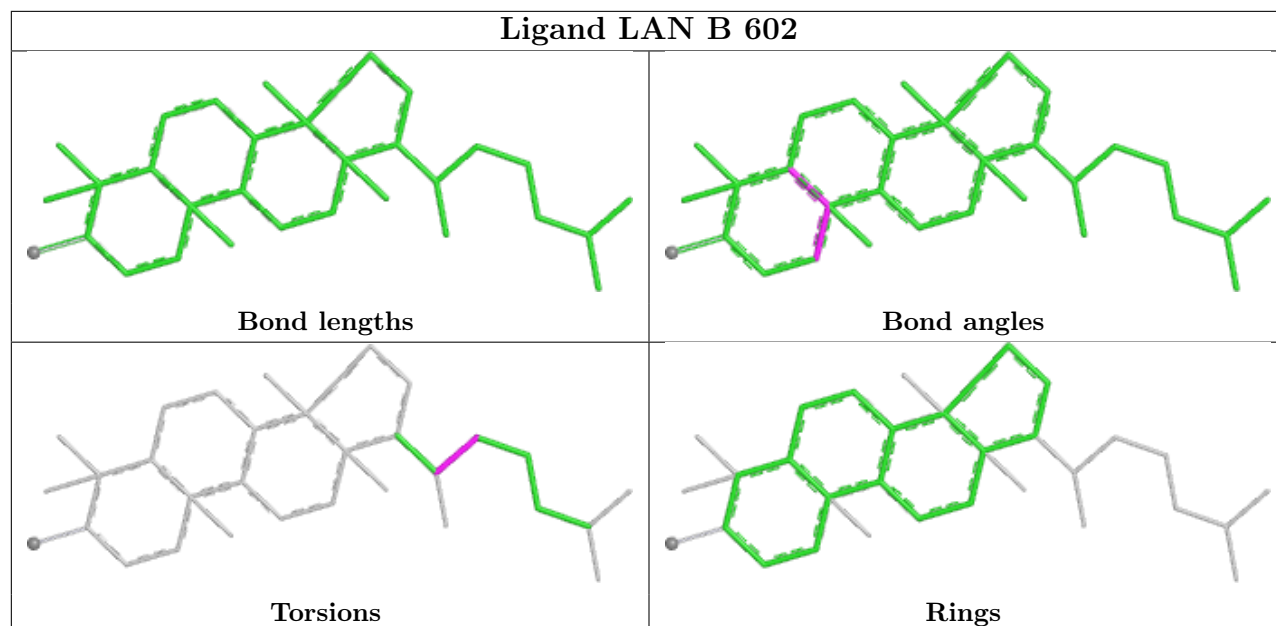




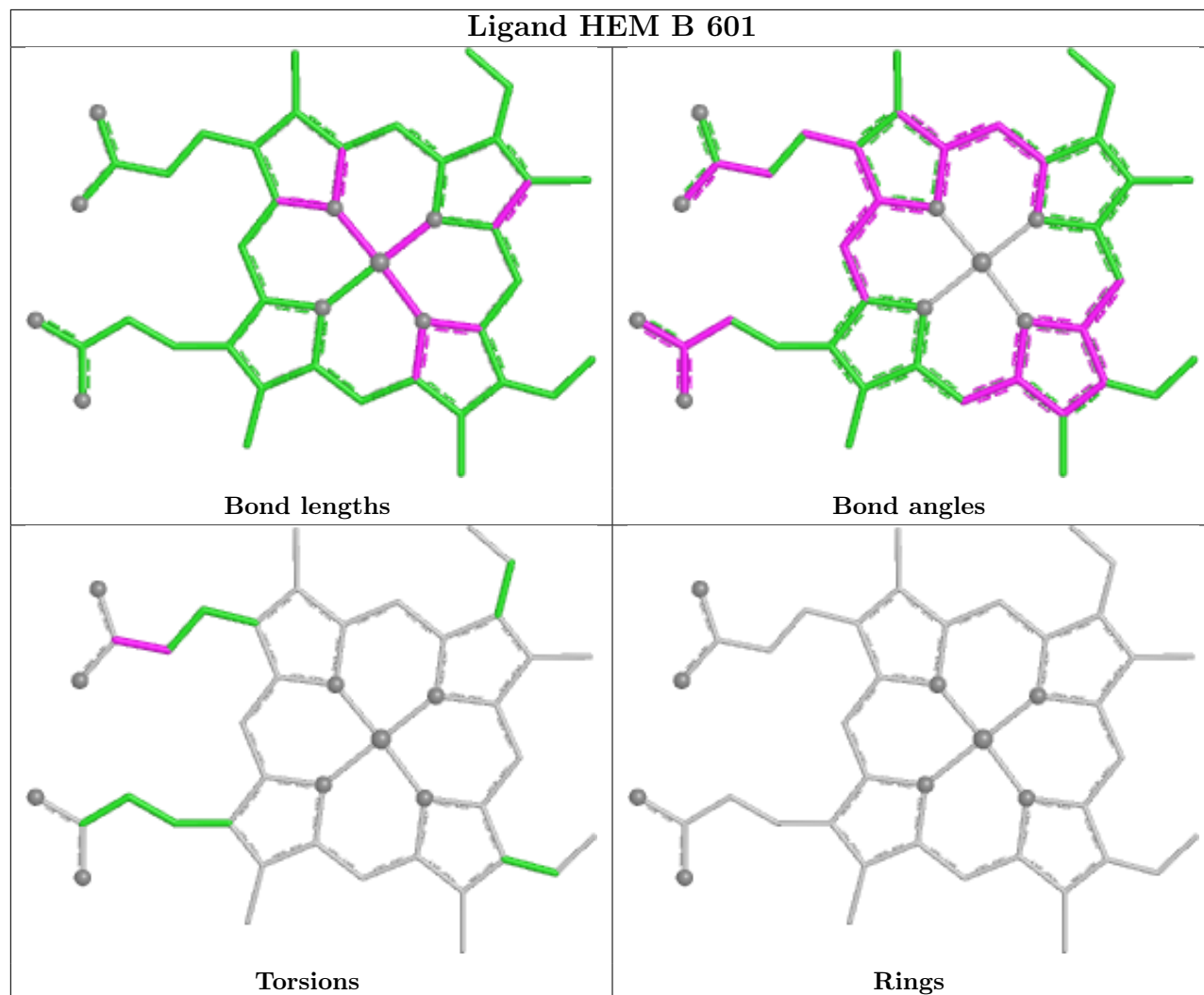




Ligand LAN B 602



Ligand HEM B 601



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	443/551 (80%)	0.04	16 (3%)	46 47	13, 38, 97, 138	0
1	B	443/551 (80%)	0.18	23 (5%)	33 32	12, 41, 104, 193	0
1	C	440/551 (79%)	0.50	40 (9%)	15 14	16, 55, 144, 172	0
1	D	439/551 (79%)	0.81	69 (15%)	5 4	15, 71, 153, 204	0
All	All	1765/2204 (80%)	0.38	148 (8%)	17 16	12, 51, 132, 204	0

The worst 5 of 148 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	371	TYR	5.1
1	D	227	VAL	4.9
1	D	297	VAL	4.8
1	B	381	LEU	4.8
1	C	371	TYR	4.6

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](#)

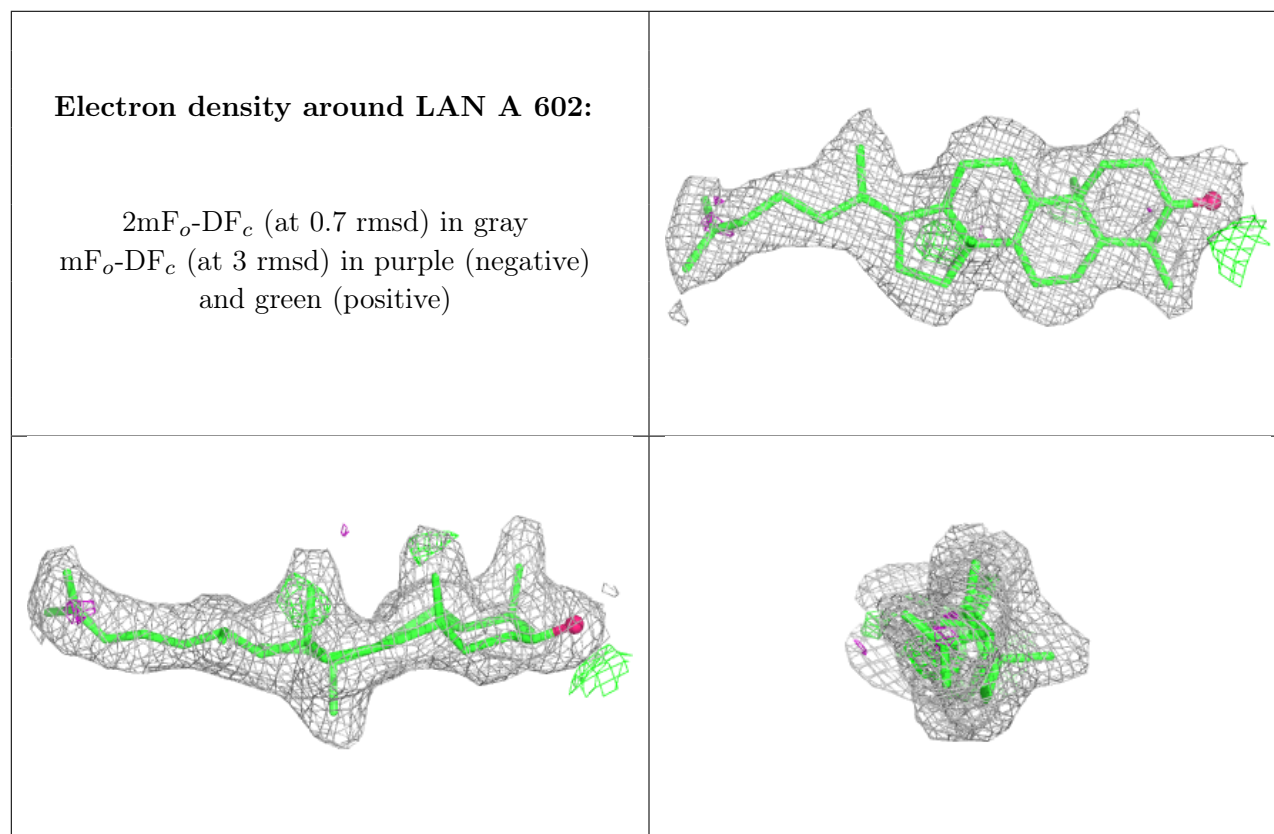
There are no oligosaccharides in this entry.

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.

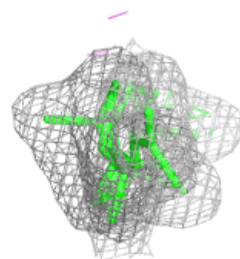
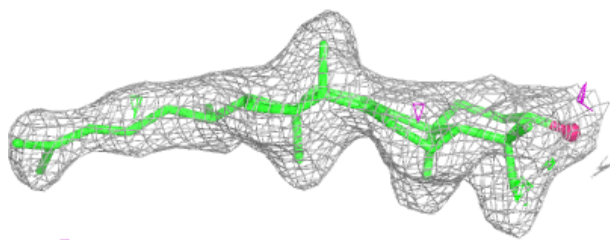
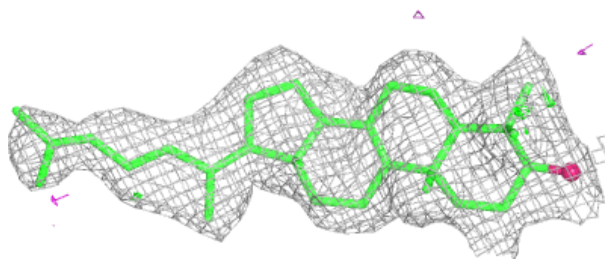
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	LAN	A	602	31/31	0.94	0.08	12,12,19,21	0
3	LAN	C	602	31/31	0.94	0.10	21,26,50,65	0
3	LAN	D	602	31/31	0.94	0.10	24,31,51,69	0
3	LAN	B	602	31/31	0.96	0.08	11,12,60,67	0
2	HEM	D	601	43/43	0.97	0.08	32,40,49,56	0
2	HEM	C	601	43/43	0.98	0.07	27,35,40,41	0
2	HEM	A	601	43/43	0.98	0.06	12,13,16,17	0
2	HEM	B	601	43/43	0.98	0.06	12,15,21,27	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.

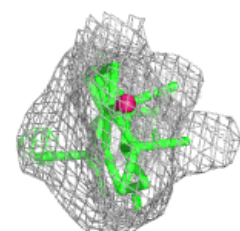
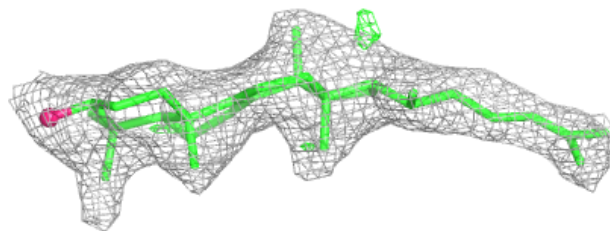
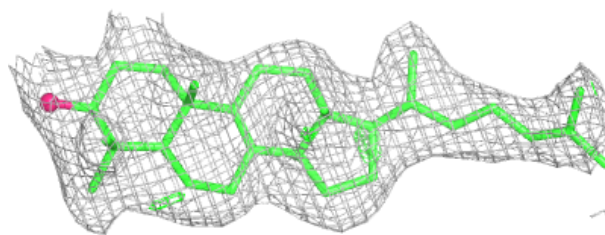


Electron density around LAN C 602:

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

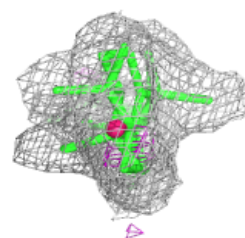
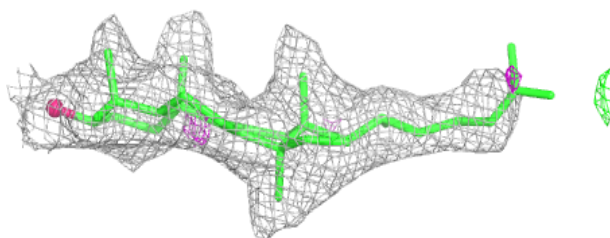
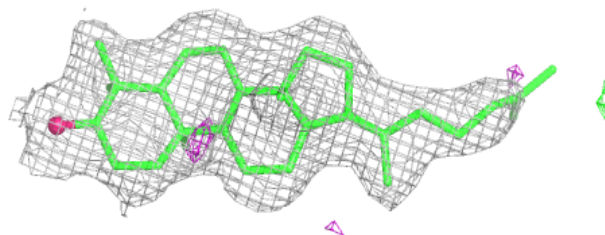
**Electron density around LAN D 602:**

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

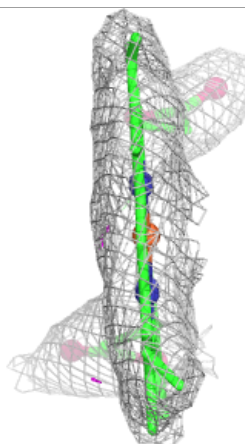
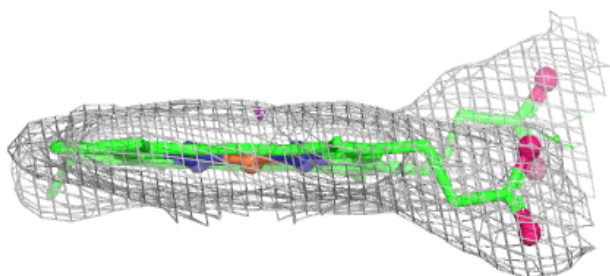
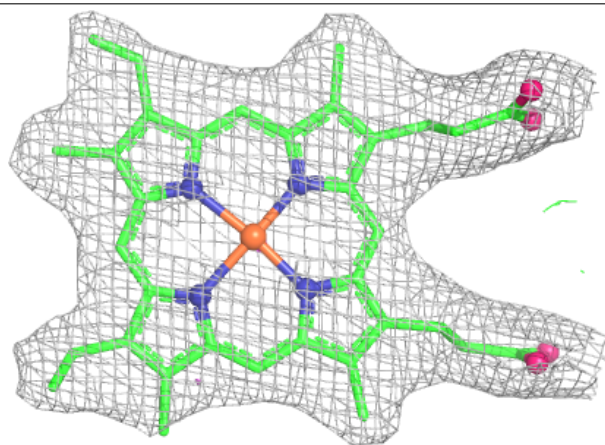


Electron density around LAN B 602:

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

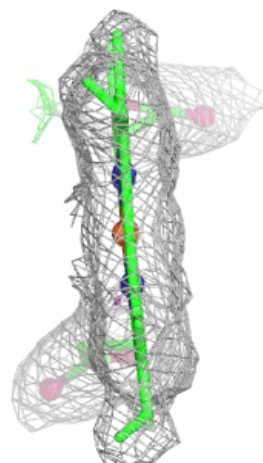
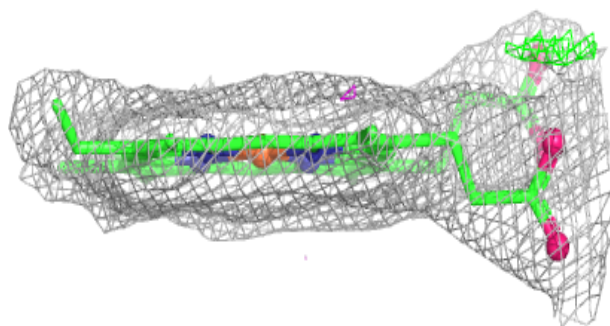
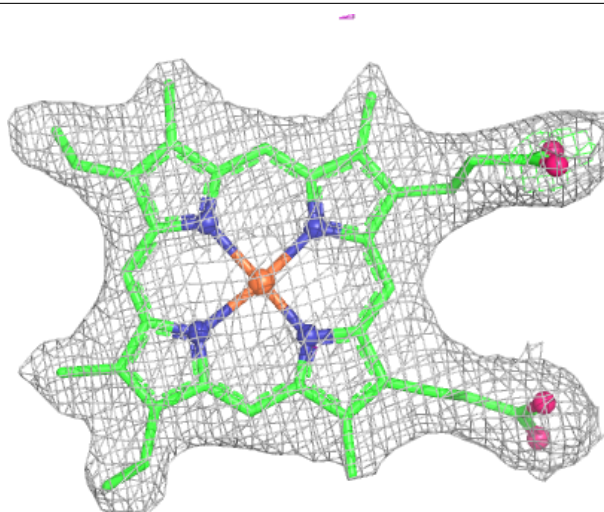
**Electron density around HEM D 601:**

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



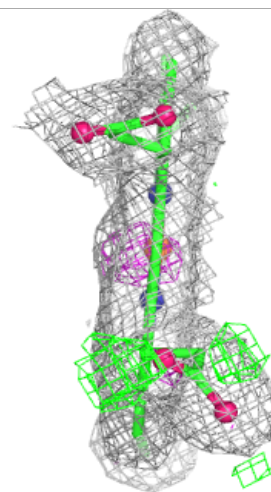
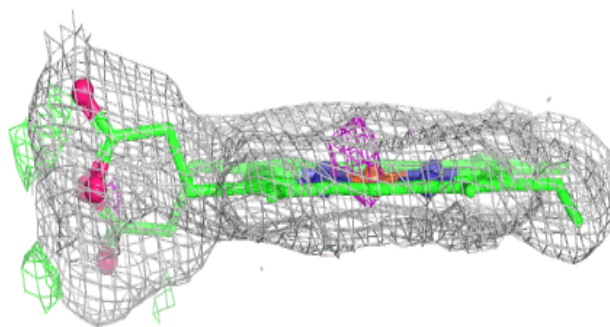
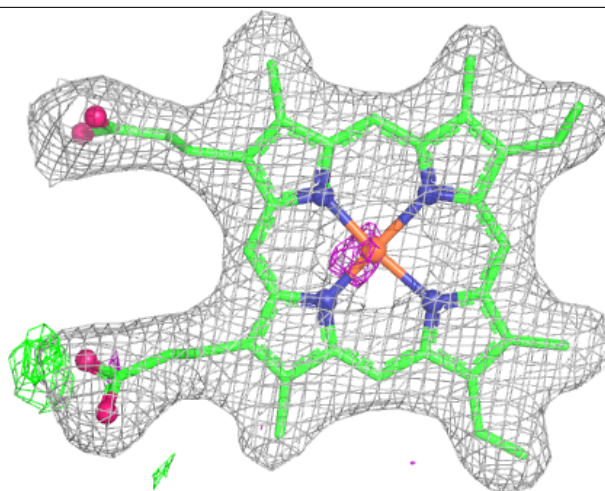
Electron density around HEM C 601:

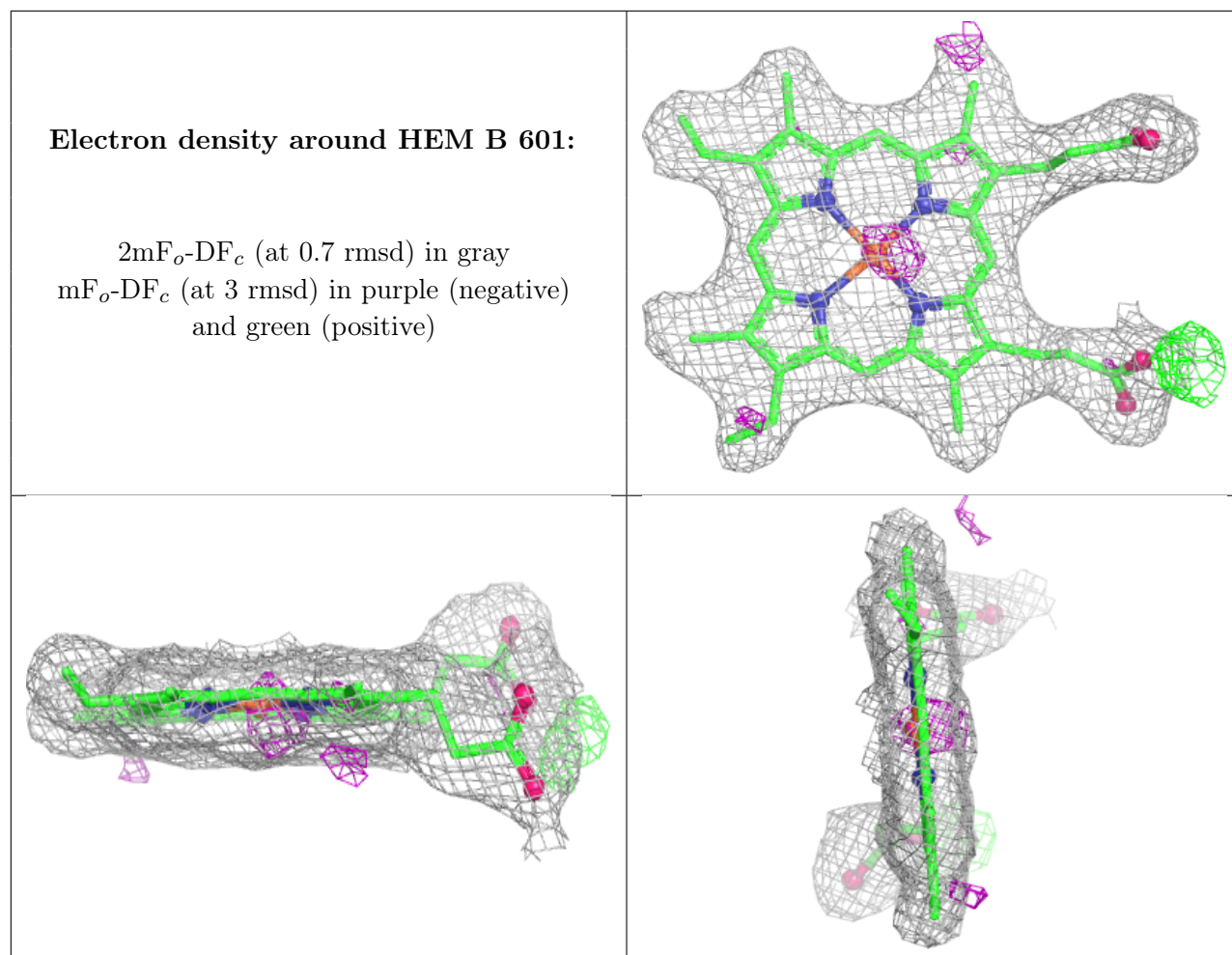
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 601:

$2mF_o-DF_c$ (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.