



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 5, 2026 – 08:26 PM EDT

PDB ID : 3N6F / pdb_00003n6f
Title : Structure of endothelial nitric oxide synthase N368D single mutant heme domain complexed with 6,6'-(2,2'-(pyridine-3,5-diyl)bis(ethane-2,1-diyl))bis(4-methylpyridin-2-amine)
Authors : Delker, S.L.; Li, H.; Poulos, T.L.
Deposited on : 2010-05-25
Resolution : 2.18 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	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

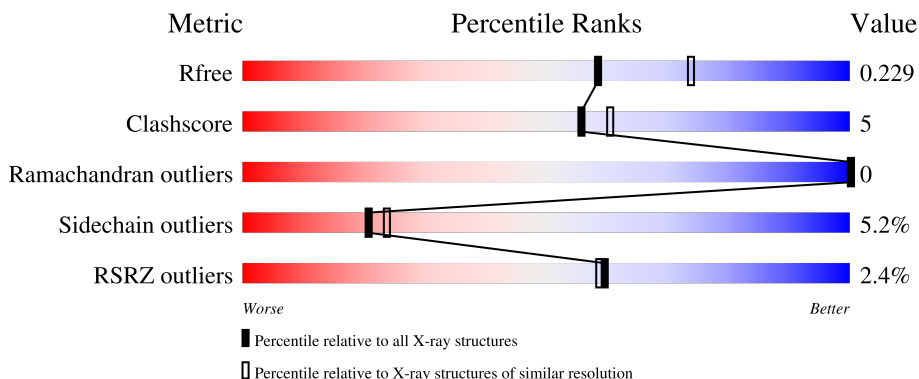
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.18 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	8975 (2.20-2.16)
Clashscore	190562	9786 (2.20-2.16)
Ramachandran outliers	187476	9664 (2.20-2.16)
Sidechain outliers	187428	9664 (2.20-2.16)
RSRZ outliers	180081	8979 (2.20-2.16)

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	444	
1	B	444	

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 6910 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.

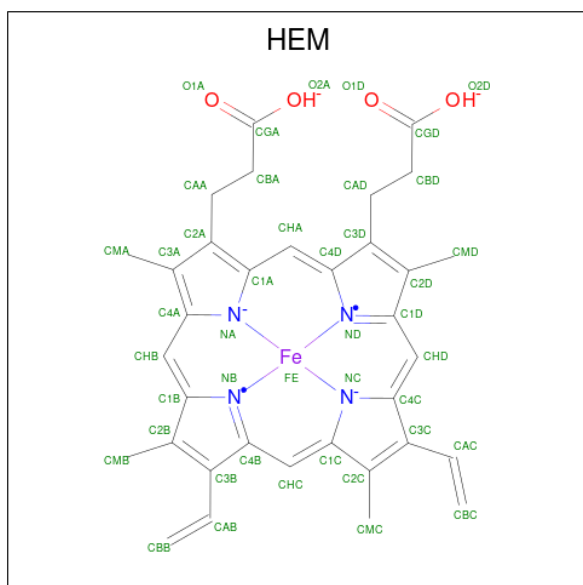
- Molecule 1 is a protein called 'Nitric oxide synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			3213	2042	566	589	16			
1	B	403	Total	C	N	O	S	0	0	0
			3209	2040	565	588	16			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ARG	CYS	SEE REMARK 999	UNP P29476
A	368	ASP	ASN	engineered mutation	UNP P29476
B	100	ARG	CYS	SEE REMARK 999	UNP P29476
B	368	ASP	ASN	engineered mutation	UNP P29476

- 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 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



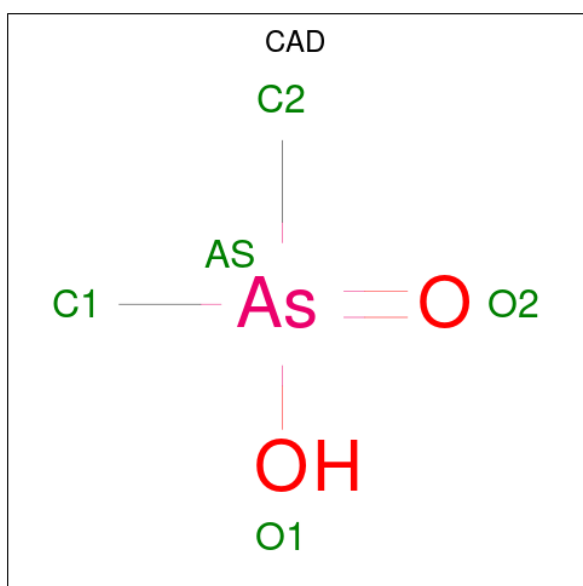
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is CACODYLIC ACID (CCD ID: CAD) (formula: $C_2H_7AsO_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	As	C	0	0
			3	1	2		
5	B	1	Total	As	C	0	0
			3	1	2		

-
- The chemical structure of H4B (Hydroxybenzoylserine) is shown. The molecule consists of a benzene ring substituted with an amino group (NH₂) at the 2-position and a carboxylic acid group (COOH) at the 4-position. The benzene ring is connected to a serine side chain at the 1-position. The serine side chain consists of a chiral center (C6(R)) bonded to a hydroxyl group (OH) and a hydrogen atom (H). The chiral center is also bonded to a methylene group (C9(R)), which is further bonded to a carboxylic acid group (COOH). The carboxylic acid group is shown with a red oxygen atom (O4) and a hydrogen atom (H). The serine side chain is also bonded to a hydrogen atom (H) at the C6(R) position. The atom labels are: N8, C7, C6(R), C9(R), C10(S), C11, O9, HO, O10, N5, N1, C8A, C4A, C4, O4, N3, N2, NH₂.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total 17	C 9	N 5	O 3	0	0
6	B	1	Total 17	C 9	N 5	O 3	0	0

-
- Chemical structure of compound XFJ, a bis-quinoline derivative. The structure consists of two quinoline rings connected by a 1,4-bis(methylene) linker. The top quinoline ring has an amino group (NH₂) at position 8. The bottom quinoline ring has an amino group (NH₂) at position 6. Carbon atoms are labeled C01 through C18, and nitrogen atoms are labeled N01 and N02.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	N	0	0
			26	21	5		
7	B	1	Total	C	N	0	0
			26	21	5		

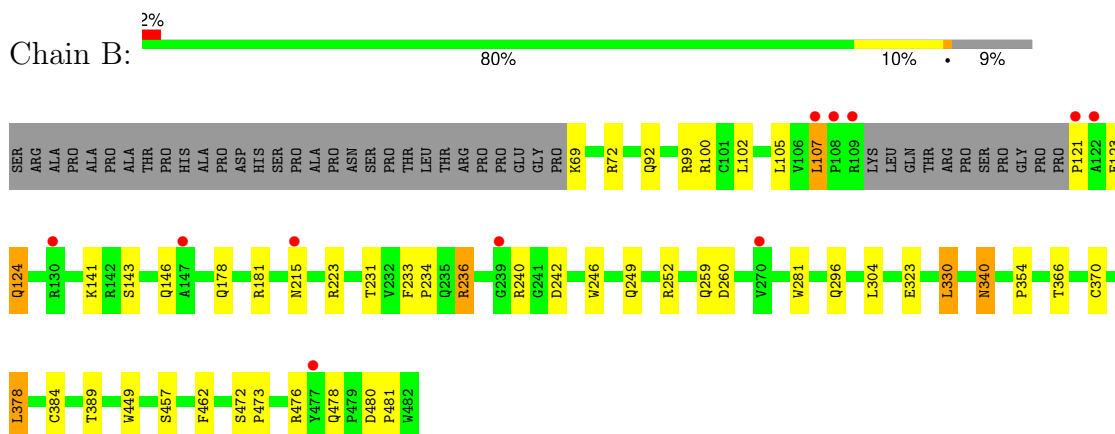
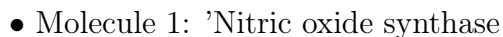
- Molecule 8 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	1	Total	Zn	0	0
			1	1		

- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	153	Total	O	0	0
			153	153		
9	B	136	Total	O	0	0
			136	136		

- Molecule 1: 'Nitric oxide synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.99Å 106.72Å 157.18Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	39.30 – 2.18 39.30 – 2.18	Depositor EDS
% Data completeness (in resolution range)	99.0 (39.30-2.18) 99.0 (39.30-2.18)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.18Å)	Xtriage
Refinement program	REFMAC 5.5.0089, CNS	Depositor
R, R_{free}	0.178 , 0.218 0.191 , 0.229	Depositor DCC
R_{free} test set	2557 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	40.4	Xtriage
Anisotropy	0.151	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 28.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6910	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.82% 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: ZN, H4B, HEM, CAD, ACT, XFJ, GOL

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	0.65	0/3302	0.81	0/4497
1	B	0.65	0/3298	0.86	3/4491 (0.1%)
All	All	0.65	0/6600	0.83	3/8988 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	240	ARG	N-CA-C	6.70	118.52	107.73
1	B	107	LEU	CA-C-N	-5.79	113.22	120.51
1	B	107	LEU	C-N-CA	-5.79	113.22	120.51

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	3213	0	3118	26	0
1	B	3209	0	3116	31	0
2	A	43	0	30	0	0
2	B	43	0	30	2	0
3	A	4	0	3	0	0
3	B	4	0	3	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	6	0	8	0	0
4	B	6	0	8	0	0
5	A	3	0	0	0	0
5	B	3	0	0	1	0
6	A	17	0	15	1	0
6	B	17	0	15	1	0
7	A	26	0	25	0	0
7	B	26	0	25	0	0
8	A	1	0	0	0	0
9	A	153	0	0	2	0
9	B	136	0	0	5	0
All	All	6910	0	6396	59	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 5.

The worst 5 of 59 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:236:ARG:HD3	1:B:242:ASP:CG	1.99	0.87
1:B:215:ASN:HB2	9:B:1247:HOH:O	1.74	0.86
1:B:236:ARG:HD3	1:B:242:ASP:OD1	1.83	0.77
1:B:121:PRO:HD2	1:B:124:GLN:HB2	1.72	0.72
1:B:72:ARG:NH1	9:B:1163:HOH:O	2.26	0.68

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	400/444 (90%)	392 (98%)	8 (2%)	0	100 100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	399/444 (90%)	393 (98%)	6 (2%)	0	100	100
All	All	799/888 (90%)	785 (98%)	14 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	343/377 (91%)	325 (95%)	18 (5%)	21	24
1	B	343/377 (91%)	325 (95%)	18 (5%)	21	24
All	All	686/754 (91%)	650 (95%)	36 (5%)	21	24

5 of 36 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	223	ARG
1	B	389	THR
1	B	236	ARG
1	B	330	LEU
1	A	311	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	296	GLN
1	B	405	ASN
1	B	478	GLN
1	B	376	ASN
1	B	146	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 13 ligands modelled in this entry, 1 is monoatomic - leaving 12 for Mogul analysis.

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	$\# Z > 2$
5	CAD	A	950	1	0,2,4	-	-	0,1,6	-	-
4	GOL	B	880	-	5,5,5	0.37	0	5,5,5	0.58	0
2	HEM	B	500	1	50,50,50	2.11	10 (20%)	67,82,82	1.50	10 (14%)
4	GOL	A	880	-	5,5,5	0.32	0	5,5,5	0.35	0
6	H4B	B	600	-	17,18,18	0.97	2 (11%)	14,26,26	1.94	5 (35%)
2	HEM	A	500	1	50,50,50	1.86	10 (20%)	67,82,82	1.32	7 (10%)
3	ACT	A	860	-	3,3,3	0.88	0	3,3,3	0.58	0
7	XFJ	A	800	-	28,28,28	0.65	0	36,38,38	1.93	8 (22%)
7	XFJ	B	800	-	28,28,28	0.51	0	36,38,38	2.10	10 (27%)
6	H4B	A	600	-	17,18,18	0.96	0	14,26,26	1.83	4 (28%)
3	ACT	B	860	-	3,3,3	0.81	0	3,3,3	0.73	0
5	CAD	B	950	1	0,2,4	-	-	0,1,6	-	-

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
4	GOL	B	880	-	-	2/4/4/4	-
2	HEM	B	500	1	-	0/14/54/54	-
4	GOL	A	880	-	-	0/4/4/4	-
2	HEM	A	500	1	-	0/14/54/54	-
7	XFJ	A	800	-	-	2/10/10/10	0/3/3/3
7	XFJ	B	800	-	-	1/10/10/10	0/3/3/3
6	H4B	A	600	-	-	0/8/17/17	0/2/2/2
6	H4B	B	600	-	-	0/8/17/17	0/2/2/2

The worst 5 of 22 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	500	HEM	C3D-C2D	7.51	1.53	1.36
2	A	500	HEM	C3D-C2D	6.65	1.51	1.36
2	B	500	HEM	FE-NA	5.22	2.12	1.95
2	B	500	HEM	FE-NC	5.15	2.12	1.95
2	A	500	HEM	FE-NA	4.76	2.10	1.95

The worst 5 of 44 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	800	XFJ	C22-N21-C26	6.72	123.10	118.07
7	A	800	XFJ	C02-N01-C06	5.95	122.52	118.07
7	A	800	XFJ	C22-N21-C26	5.21	121.96	118.07
7	B	800	XFJ	C02-N01-C06	4.93	121.76	118.07
6	B	600	H4B	C2-N1-C8A	4.41	121.14	113.36

There are no chirality outliers.

All (5) torsion outliers are listed below:

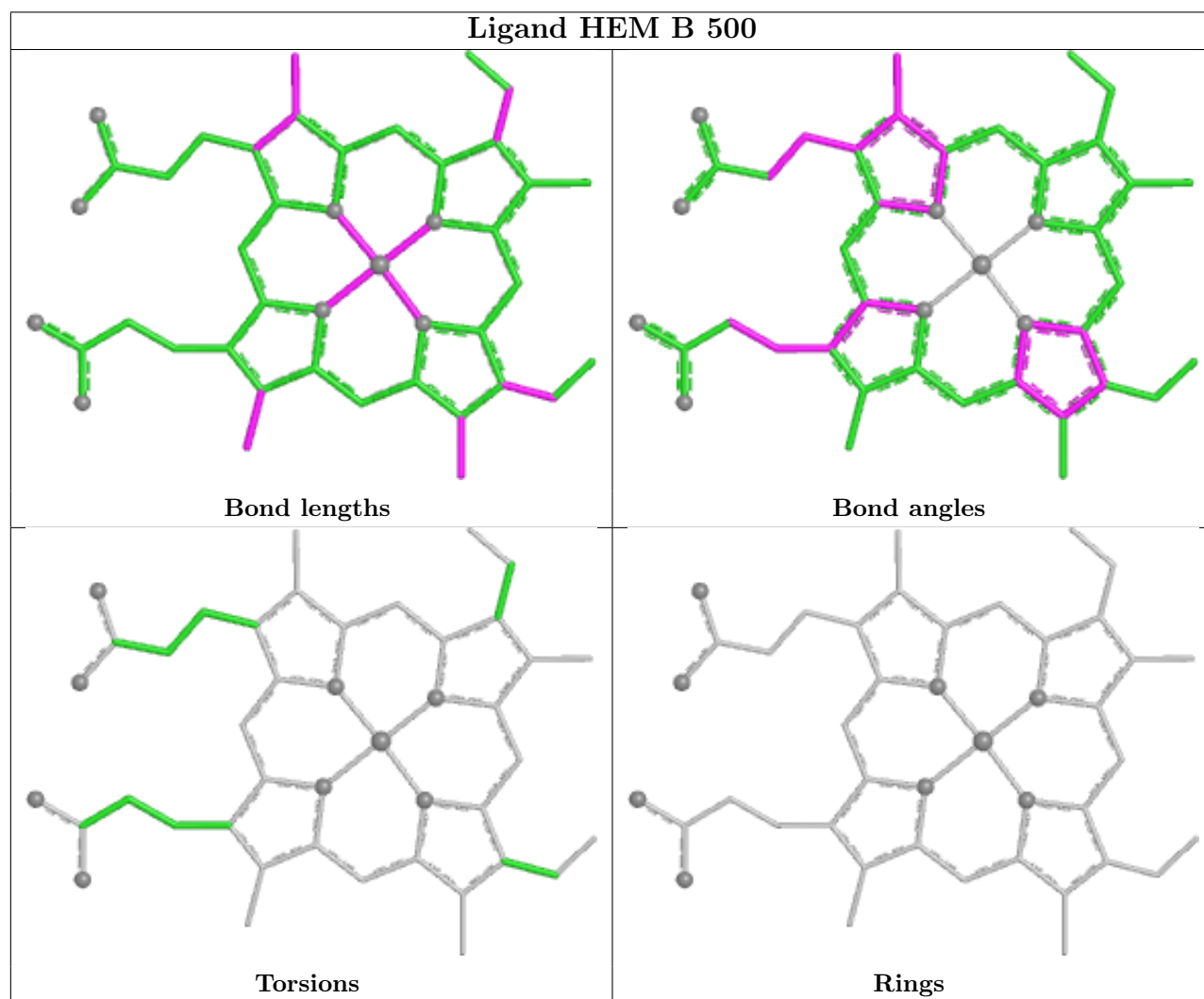
Mol	Chain	Res	Type	Atoms
4	B	880	GOL	C1-C2-C3-O3
4	B	880	GOL	O2-C2-C3-O3
7	A	800	XFJ	C05-C06-C08-C09
7	A	800	XFJ	N01-C06-C08-C09
7	B	800	XFJ	N01-C06-C08-C09

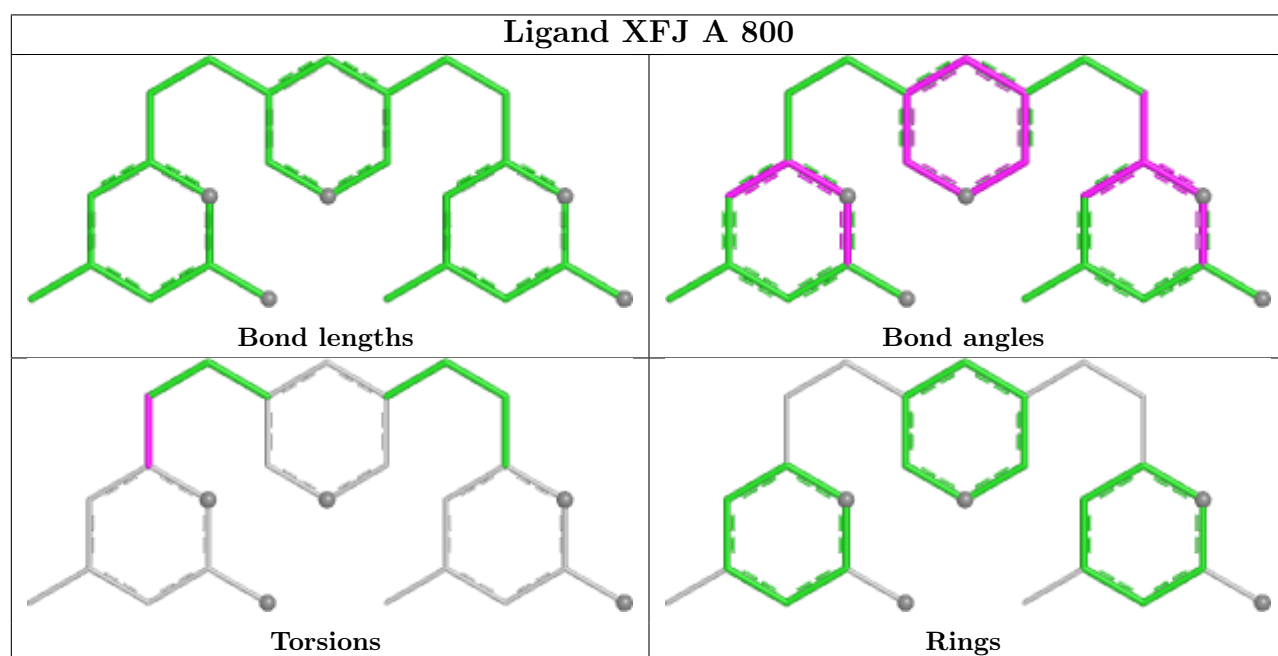
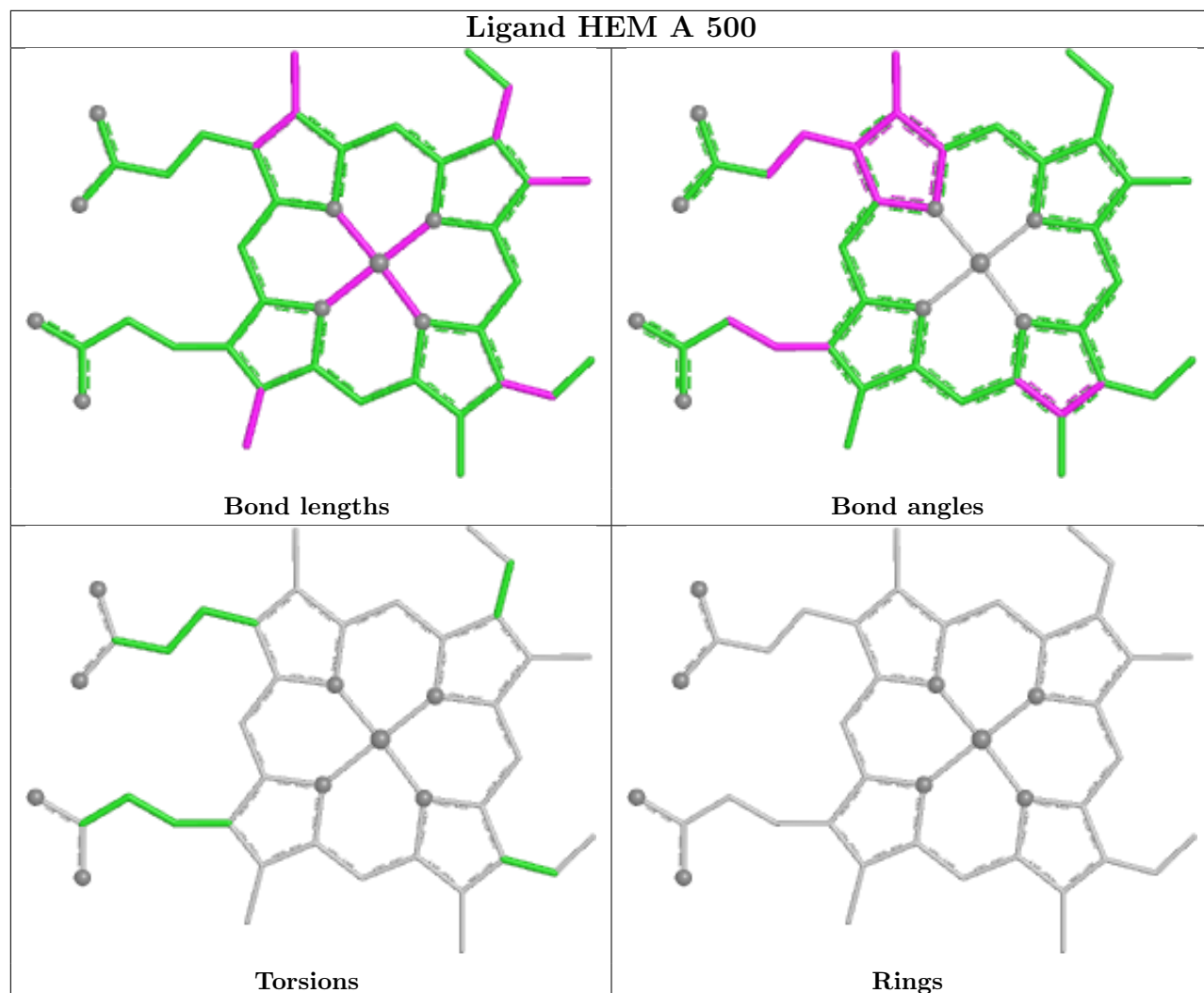
There are no ring outliers.

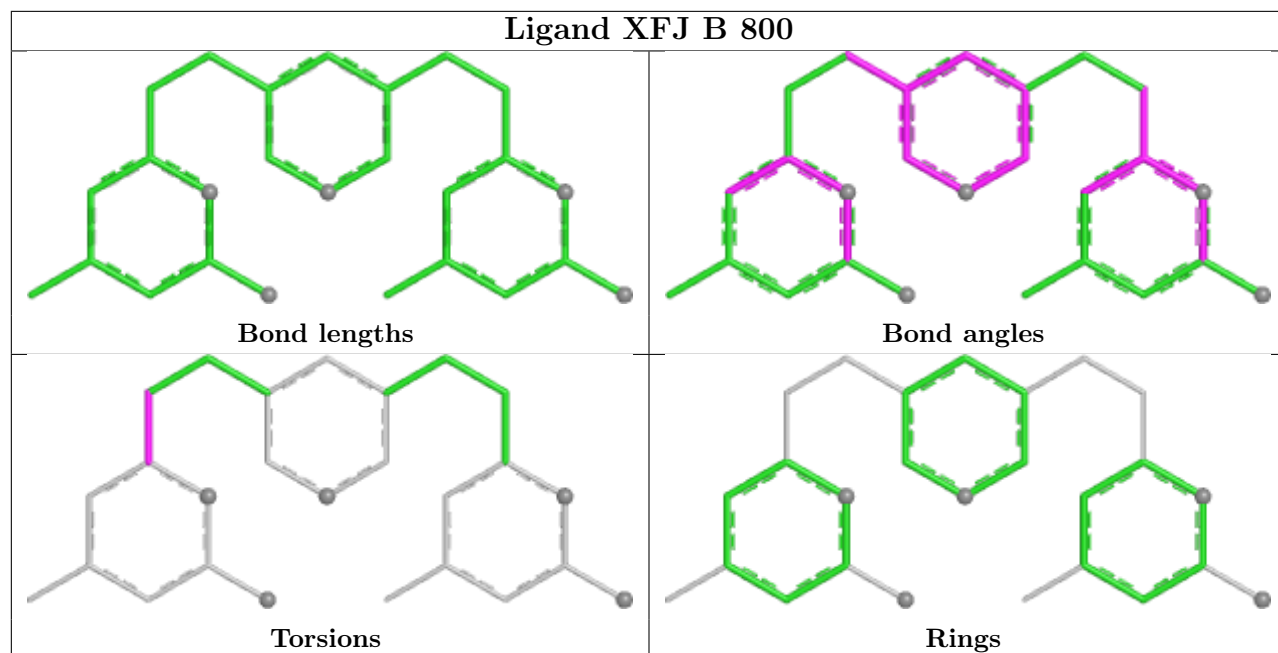
4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	500	HEM	2	0
6	B	600	H4B	1	0
6	A	600	H4B	1	0
5	B	950	CAD	1	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 [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	404/444 (90%)	-0.01	8 (1%) 65 64	28, 41, 67, 81	0
1	B	403/444 (90%)	0.04	11 (2%) 56 55	28, 44, 67, 88	0
All	All	807/888 (90%)	0.01	19 (2%) 59 59	28, 42, 67, 88	0

The worst 5 of 19 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	122	ALA	5.4
1	B	108	PRO	4.6
1	A	239	GLY	4.0
1	B	477	TYR	3.3
1	B	121	PRO	3.0

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 oligosaccharides in this entry.

6.4 Ligands [i](#)

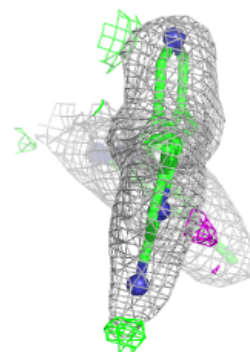
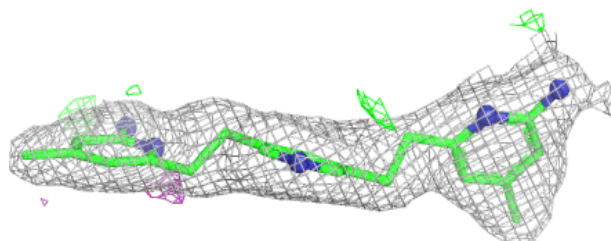
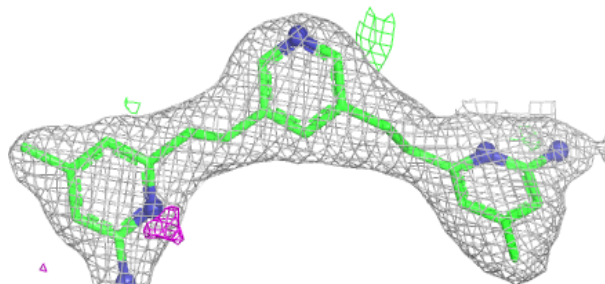
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists 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
4	GOL	A	880	6/6	0.81	0.14	60,61,63,63	0
4	GOL	B	880	6/6	0.86	0.17	58,58,58,60	0
7	XFJ	B	800	26/26	0.93	0.09	34,44,49,49	0
3	ACT	B	860	4/4	0.95	0.12	55,55,55,56	0
5	CAD	B	950	3/5	0.95	0.14	69,69,69,70	0
7	XFJ	A	800	26/26	0.95	0.08	28,39,42,43	0
3	ACT	A	860	4/4	0.95	0.09	45,45,45,47	0
6	H4B	A	600	17/17	0.97	0.05	30,34,36,36	0
6	H4B	B	600	17/17	0.97	0.05	30,32,35,35	0
5	CAD	A	950	3/5	0.98	0.10	65,65,65,65	0
2	HEM	A	500	43/43	0.98	0.06	27,29,40,44	0
2	HEM	B	500	43/43	0.99	0.04	26,30,36,39	0
8	ZN	A	900	1/1	1.00	0.01	34,34,34,34	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.

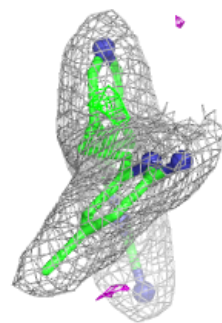
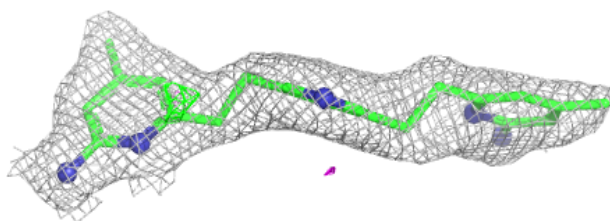
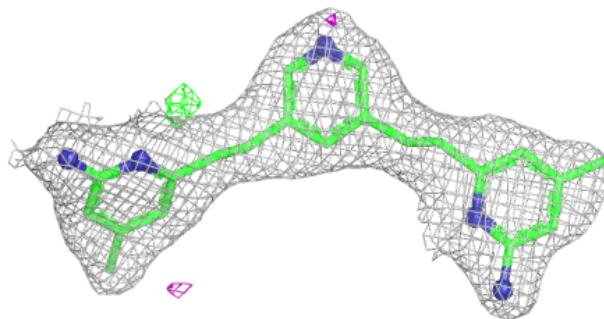
Electron density around XFJ B 800:

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



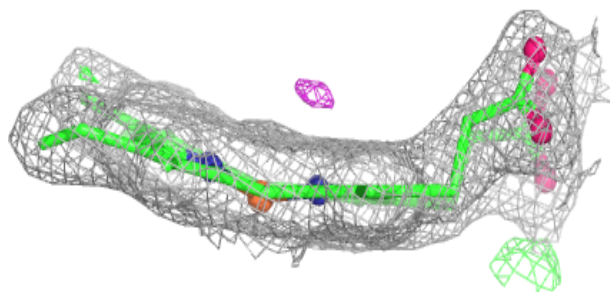
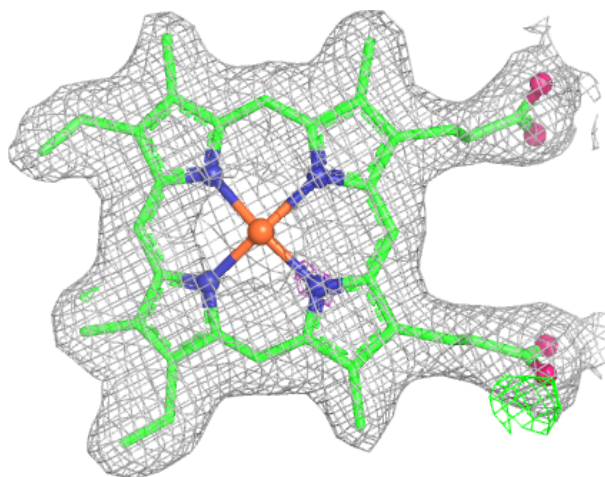
Electron density around XFJ A 800:

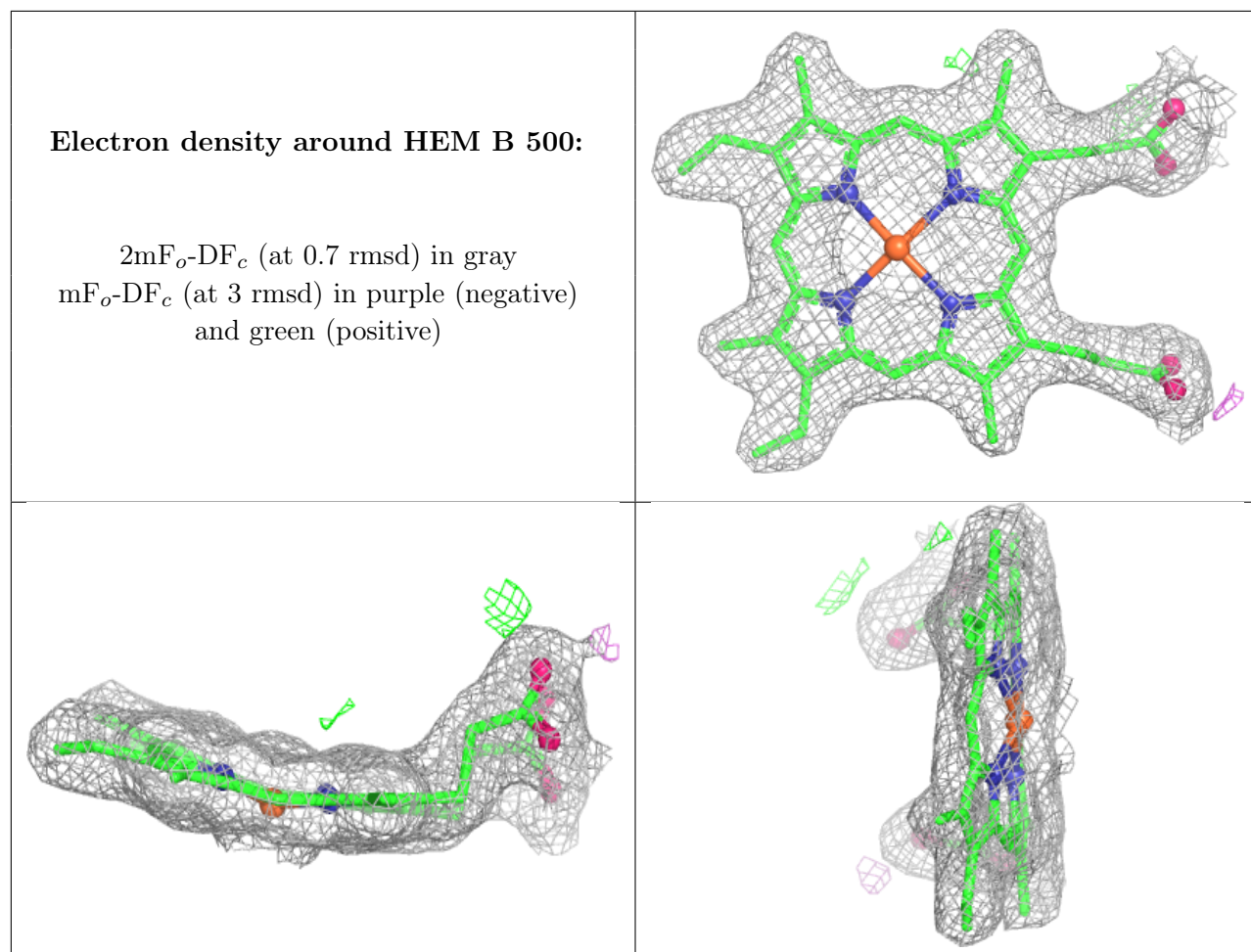
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEM A 500:

$2mF_o-DF_c$ (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.