



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 6, 2023 – 10:20 PM EDT

PDB ID : 1JBQ
Title : STRUCTURE OF HUMAN CYSTATHIONINE BETA-SYNTASE: A
UNIQUE PYRIDOXAL 5'-PHOSPHATE DEPENDENT HEMEPROTEIN
Authors : Meier, M.; Janosik, M.; Kery, V.; Kraus, J.P.; Burkhard, P.
Deposited on : 2001-06-06
Resolution : 2.60 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : 2.35
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

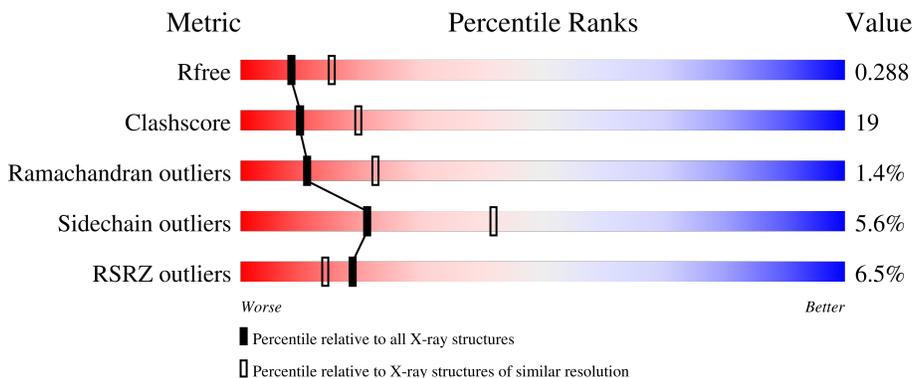
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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}	130704	3163 (2.60-2.60)
Clashscore	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)
RSRZ outliers	127900	3104 (2.60-2.60)

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	435	 3% 51% 26% 20%
1	B	435	 8% 51% 26% 20%
1	C	435	 6% 50% 27% 20%
1	D	435	 6% 50% 27% 20%
1	E	435	 4% 52% 25% 20%

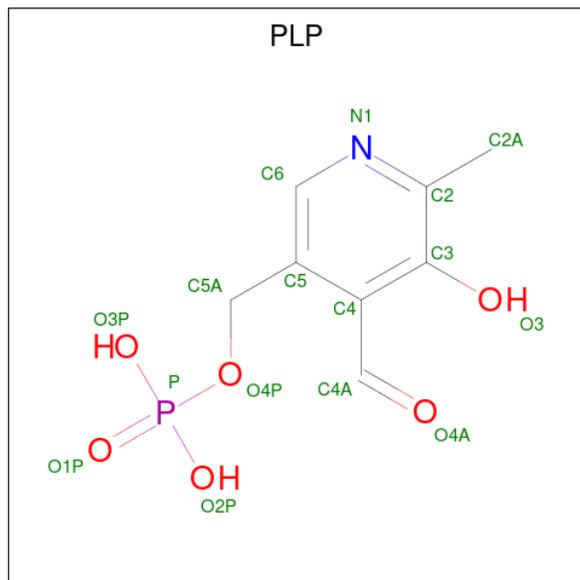
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Mol	Chain	Length	Quality of chain
1	F	435	 <p>A horizontal bar chart representing the quality of chain. The bar is divided into four segments: a small red segment (4%), a large green segment (53%), a yellow segment (25%), and a grey segment (20%).</p>

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		
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	E	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	F	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is PYRIDOXAL-5'-PHOSPHATE (three-letter code: PLP) (formula: C₈H₁₀NO₆P).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
3	A	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	B	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	C	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	D	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	E	1	Total	C	N	O	P	0	0
			15	8	1	5	1		
3	F	1	Total	C	N	O	P	0	0
			15	8	1	5	1		

- Molecule 4 is water.

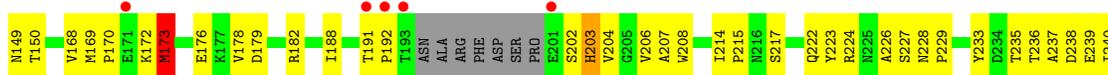
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	22	Total O 22 22	0	0
4	B	30	Total O 30 30	0	0
4	C	25	Total O 25 25	0	0
4	D	24	Total O 24 24	0	0
4	E	31	Total O 31 31	0	0
4	F	22	Total O 22 22	0	0



● Molecule 1: CYSTATHIONINE BETA-SYNTASE

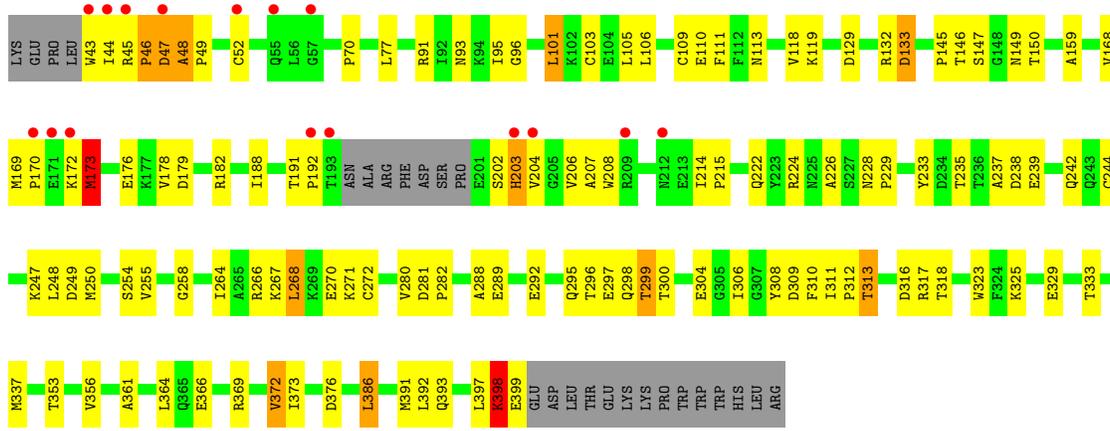


● Molecule 1: CYSTATHIONINE BETA-SYNTASE

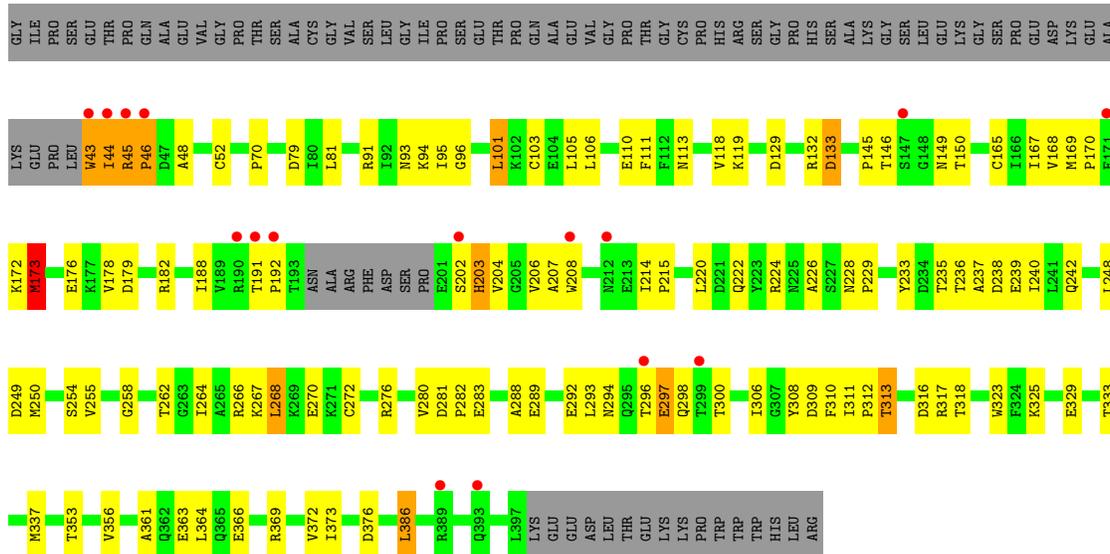


● Molecule 1: CYSTATHIONINE BETA-SYNTASE





● Molecule 1: CYSTATHIONINE BETA-SYNTASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	144.52Å 144.52Å 108.16Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.34 – 2.60 43.34 – 2.60	Depositor EDS
% Data completeness (in resolution range)	82.3 (43.34-2.60) 82.3 (43.34-2.60)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.49 (at 2.61Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.257 , 0.296 0.250 , 0.288	Depositor DCC
R_{free} test set	3219 reflections (4.83%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtrriage
Anisotropy	0.110	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 47.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	0.000 for -h,-k,l 0.036 for h,-h-k,-l 0.000 for -k,-h,-l	Xtrriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	16522	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 43.39 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.7767e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: PLP, 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	0.39	0/2715	0.67	1/3672 (0.0%)
1	B	0.38	0/2715	0.65	0/3672
1	C	0.39	0/2715	0.65	0/3672
1	D	0.37	0/2715	0.66	1/3672 (0.0%)
1	E	0.38	0/2733	0.67	0/3695
1	F	0.38	0/2715	0.66	0/3672
All	All	0.38	0/16308	0.66	2/22055 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	44	ILE	N-CA-C	-8.54	87.93	111.00
1	D	44	ILE	N-CA-C	-7.01	92.06	111.00

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	2667	0	2691	108	0
1	B	2667	0	2691	103	0
1	C	2667	0	2691	108	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2667	0	2691	114	0
1	E	2685	0	2710	96	0
1	F	2667	0	2691	107	0
2	A	43	0	30	4	0
2	B	43	0	30	2	0
2	C	43	0	30	2	0
2	D	43	0	30	3	0
2	E	43	0	30	2	0
2	F	43	0	30	3	0
3	A	15	0	7	1	0
3	B	15	0	7	1	0
3	C	15	0	7	1	0
3	D	15	0	7	1	0
3	E	15	0	7	1	0
3	F	15	0	7	1	0
4	A	22	0	0	4	0
4	B	30	0	0	2	0
4	C	25	0	0	6	0
4	D	24	0	0	7	0
4	E	31	0	0	2	0
4	F	22	0	0	0	0
All	All	16522	0	16387	628	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 19.

The worst 5 of 628 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:E:296:THR:HG22	1:E:298:GLN:H	1.30	0.94
1:A:296:THR:HG22	1:A:298:GLN:H	1.33	0.93
1:C:296:THR:HG22	1:C:298:GLN:H	1.32	0.92
1:D:296:THR:HG22	1:D:298:GLN:H	1.33	0.92
1:B:296:THR:HG22	1:B:298:GLN:H	1.33	0.91

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	344/435 (79%)	325 (94%)	16 (5%)	3 (1%)	17	35
1	B	344/435 (79%)	319 (93%)	19 (6%)	6 (2%)	9	18
1	C	344/435 (79%)	323 (94%)	18 (5%)	3 (1%)	17	35
1	D	344/435 (79%)	319 (93%)	19 (6%)	6 (2%)	9	18
1	E	346/435 (80%)	321 (93%)	19 (6%)	6 (2%)	9	18
1	F	344/435 (79%)	322 (94%)	17 (5%)	5 (2%)	10	21
All	All	2066/2610 (79%)	1929 (93%)	108 (5%)	29 (1%)	11	22

5 of 29 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	50	SER
1	B	146	THR
1	B	173	MET
1	C	173	MET
1	D	45	ARG

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	288/362 (80%)	272 (94%)	16 (6%)	21	42
1	B	288/362 (80%)	273 (95%)	15 (5%)	23	46
1	C	288/362 (80%)	272 (94%)	16 (6%)	21	42

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	D	288/362 (80%)	272 (94%)	16 (6%)	21	42
1	E	290/362 (80%)	272 (94%)	18 (6%)	18	37
1	F	288/362 (80%)	272 (94%)	16 (6%)	21	42
All	All	1730/2172 (80%)	1633 (94%)	97 (6%)	21	42

5 of 97 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	D	299	THR
1	E	208	TRP
1	D	313	THR
1	E	101	LEU
1	E	309	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 25 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	341	GLN
1	E	242	GLN
1	F	393	GLN
1	E	222	GLN
1	E	393	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry

12 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
3	PLP	F	1001	1	15,15,16	1.28	2 (13%)	20,22,23	1.13	2 (10%)
3	PLP	C	1001	1	15,15,16	1.21	0	20,22,23	1.11	2 (10%)
2	HEM	B	501	1	41,50,50	1.75	7 (17%)	45,82,82	1.32	7 (15%)
3	PLP	E	1001	1	15,15,16	1.51	1 (6%)	20,22,23	1.10	1 (5%)
2	HEM	D	501	1	41,50,50	1.76	7 (17%)	45,82,82	1.33	7 (15%)
2	HEM	C	501	1	41,50,50	1.75	9 (21%)	45,82,82	1.35	7 (15%)
2	HEM	E	501	1	41,50,50	1.75	8 (19%)	45,82,82	1.38	6 (13%)
2	HEM	A	501	1	41,50,50	1.79	9 (21%)	45,82,82	1.34	7 (15%)
3	PLP	B	1001	1	15,15,16	1.24	1 (6%)	20,22,23	1.11	1 (5%)
2	HEM	F	501	1	41,50,50	1.76	8 (19%)	45,82,82	1.38	6 (13%)
3	PLP	D	1001	1	15,15,16	1.26	1 (6%)	20,22,23	1.11	1 (5%)
3	PLP	A	1001	1	15,15,16	1.29	2 (13%)	20,22,23	1.16	1 (5%)

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
3	PLP	F	1001	1	-	0/6/6/8	0/1/1/1
3	PLP	C	1001	1	-	0/6/6/8	0/1/1/1
2	HEM	B	501	1	-	4/12/54/54	-
3	PLP	E	1001	1	-	0/6/6/8	0/1/1/1
2	HEM	D	501	1	-	4/12/54/54	-
2	HEM	C	501	1	-	4/12/54/54	-
2	HEM	E	501	1	-	4/12/54/54	-
2	HEM	A	501	1	-	4/12/54/54	-
3	PLP	B	1001	1	-	0/6/6/8	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	F	501	1	-	4/12/54/54	-
3	PLP	D	1001	1	-	0/6/6/8	0/1/1/1
3	PLP	A	1001	1	-	0/6/6/8	0/1/1/1

The worst 5 of 55 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	CBB-CAB	4.97	1.54	1.30
2	D	501	HEM	CBB-CAB	4.95	1.54	1.30
2	A	501	HEM	CBB-CAB	4.91	1.54	1.30
2	C	501	HEM	CBB-CAB	4.89	1.54	1.30
2	E	501	HEM	CBB-CAB	4.85	1.54	1.30

The worst 5 of 48 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	501	HEM	C4C-CHD-C1D	4.23	128.13	122.56
2	E	501	HEM	C4C-CHD-C1D	4.08	127.95	122.56
2	A	501	HEM	C4C-CHD-C1D	3.98	127.82	122.56
2	D	501	HEM	C4C-CHD-C1D	3.98	127.81	122.56
2	B	501	HEM	C4C-CHD-C1D	3.89	127.70	122.56

There are no chirality outliers.

5 of 24 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	HEM	C2B-C3B-CAB-CBB
2	B	501	HEM	C2B-C3B-CAB-CBB
2	C	501	HEM	C2B-C3B-CAB-CBB
2	D	501	HEM	C2B-C3B-CAB-CBB
2	E	501	HEM	C2B-C3B-CAB-CBB

There are no ring outliers.

12 monomers are involved in 22 short contacts:

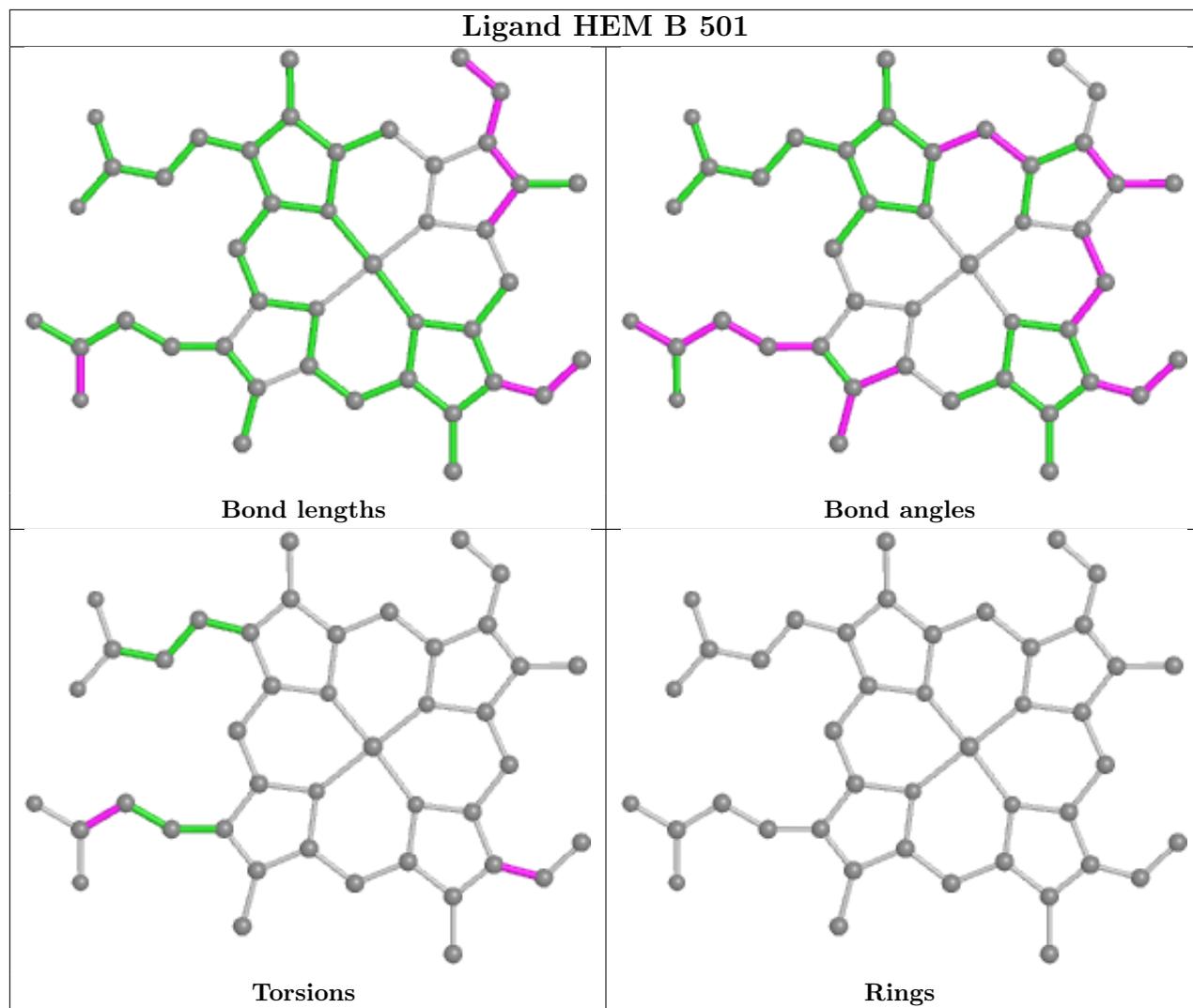
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	F	1001	PLP	1	0
3	C	1001	PLP	1	0
2	B	501	HEM	2	0
3	E	1001	PLP	1	0
2	D	501	HEM	3	0

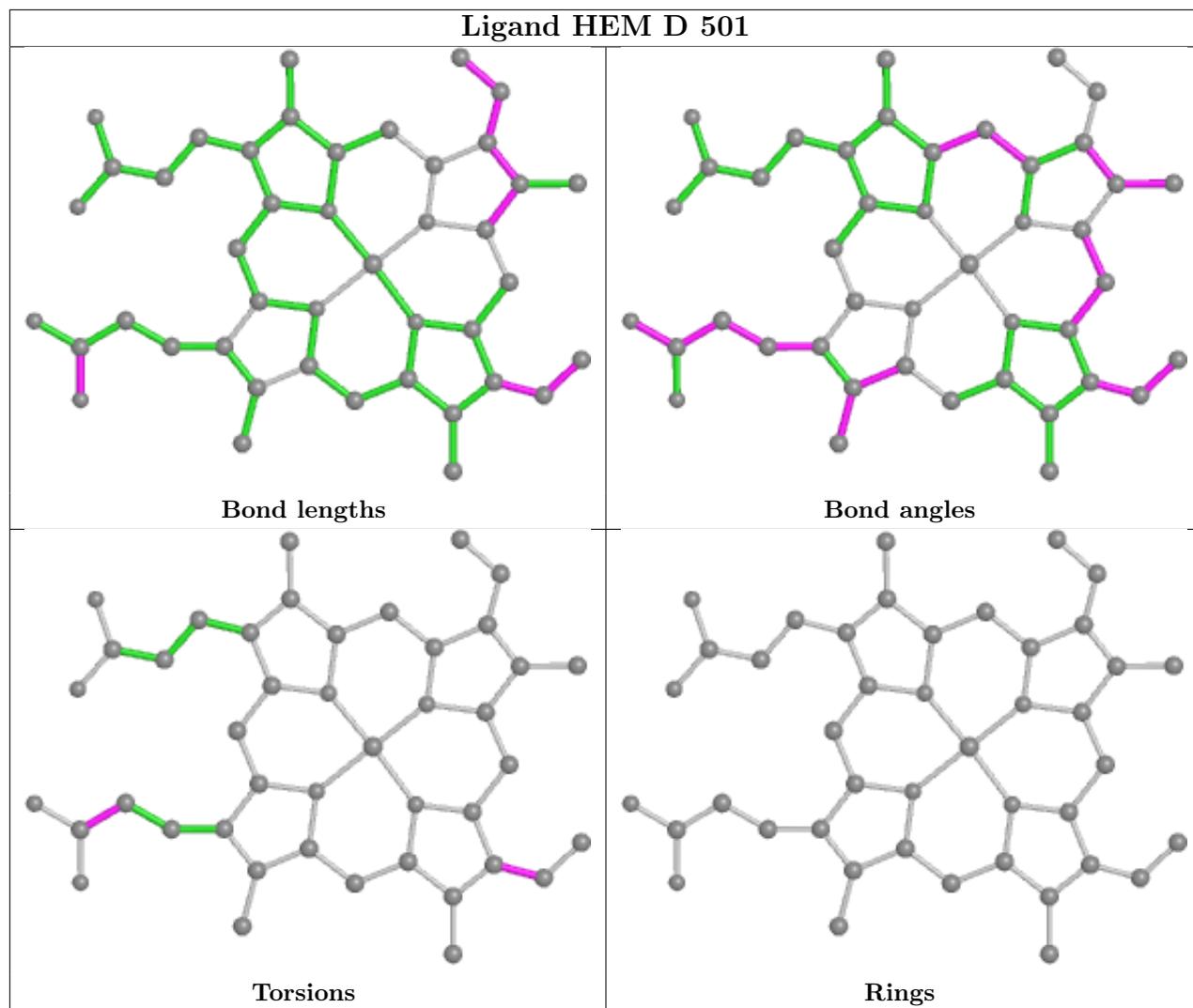
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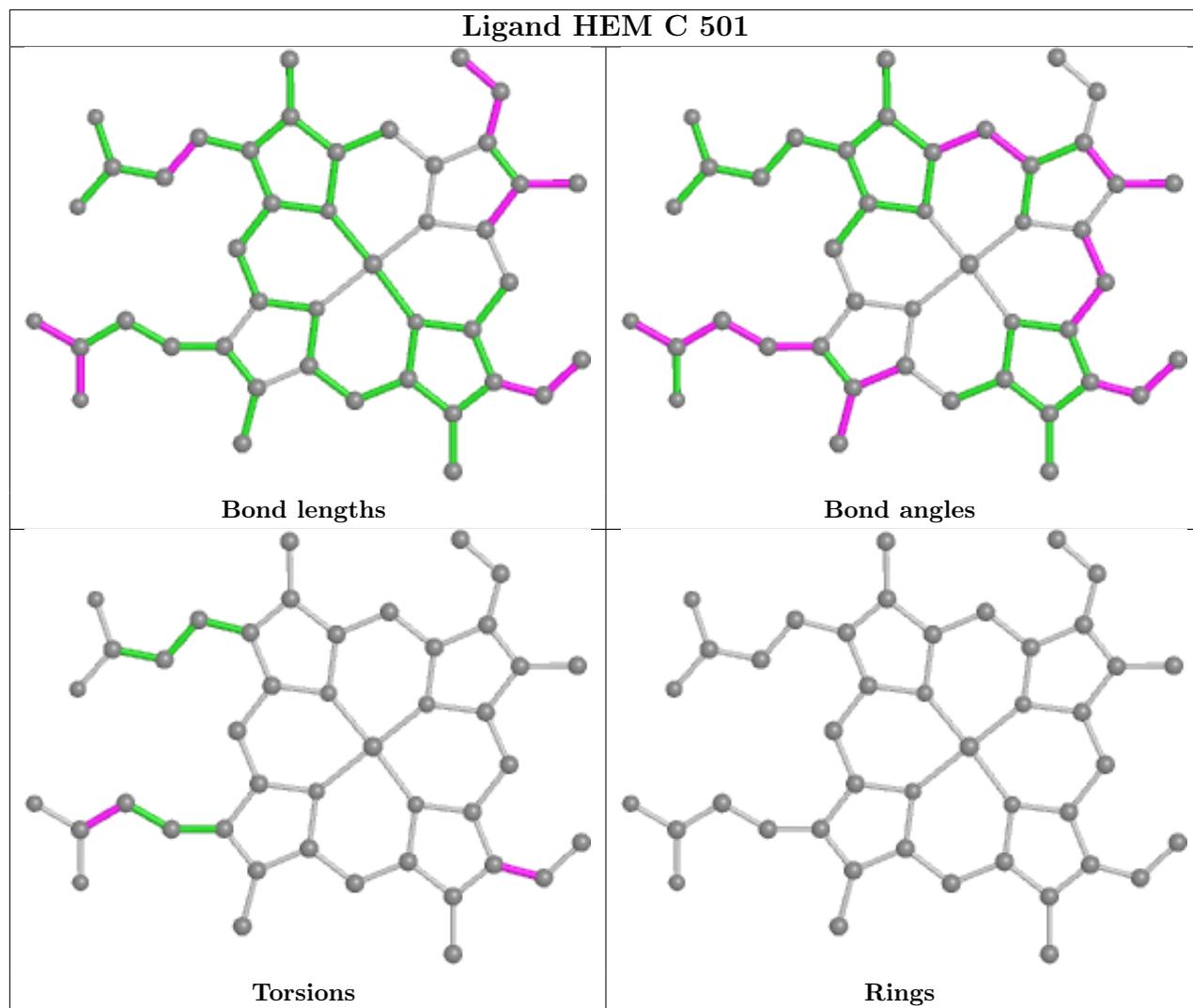
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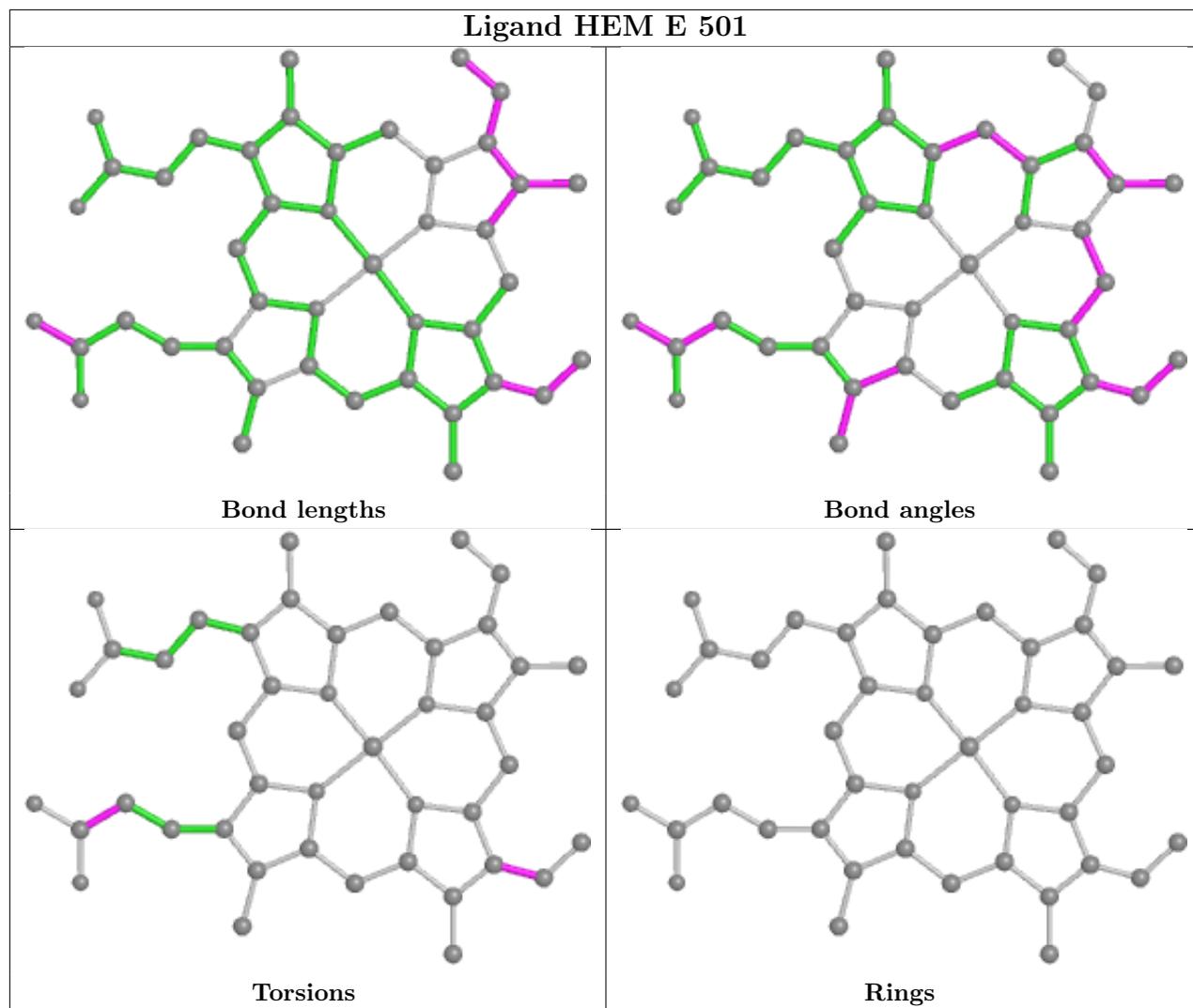
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	501	HEM	2	0
2	E	501	HEM	2	0
2	A	501	HEM	4	0
3	B	1001	PLP	1	0
2	F	501	HEM	3	0
3	D	1001	PLP	1	0
3	A	1001	PLP	1	0

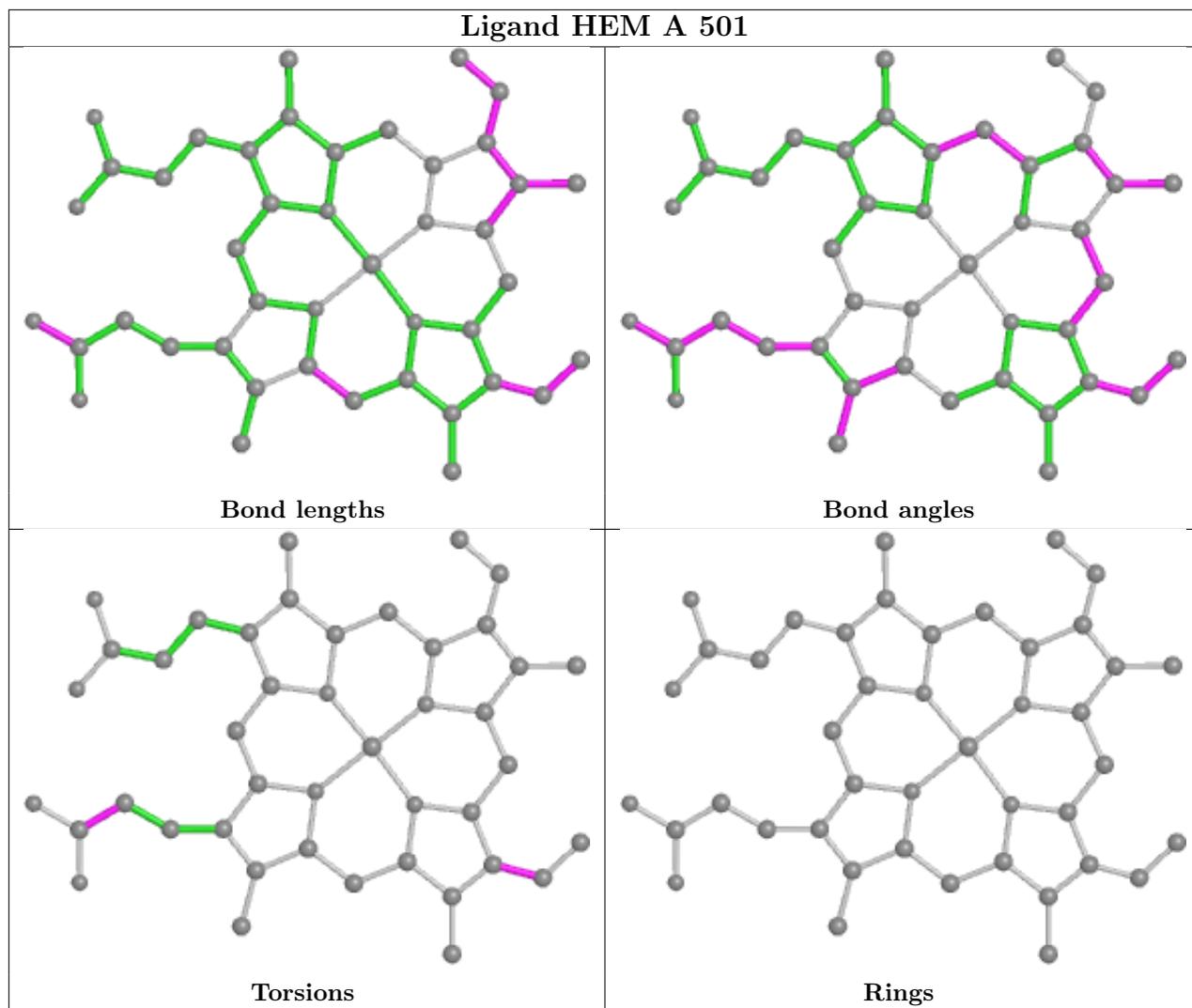
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

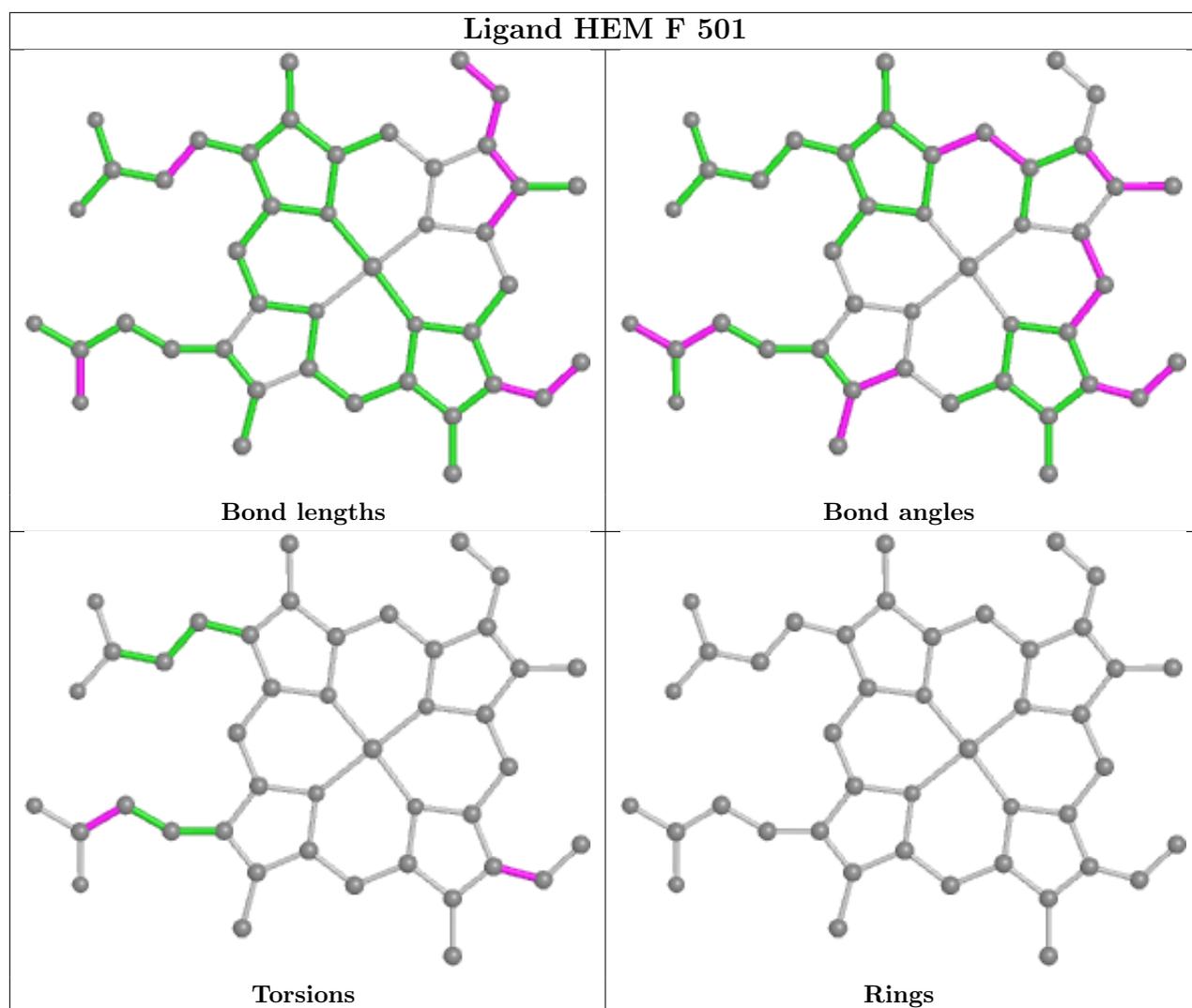












5.7 Other polymers [i](#)

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	348/435 (80%)	0.21	15 (4%) 35 28	12, 32, 66, 85	0
1	B	348/435 (80%)	0.61	36 (10%) 6 4	9, 35, 75, 94	0
1	C	348/435 (80%)	0.44	26 (7%) 14 10	15, 36, 72, 87	0
1	D	348/435 (80%)	0.46	27 (7%) 13 9	17, 39, 76, 92	0
1	E	350/435 (80%)	0.15	16 (4%) 32 26	9, 27, 67, 90	0
1	F	348/435 (80%)	0.08	16 (4%) 32 26	9, 26, 67, 89	0
All	All	2090/2610 (80%)	0.32	136 (6%) 18 14	9, 33, 72, 94	0

The worst 5 of 136 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	193	THR	9.7
1	E	193	THR	6.7
1	B	43	TRP	6.3
1	F	147	SER	6.3
1	D	392	LEU	6.1

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

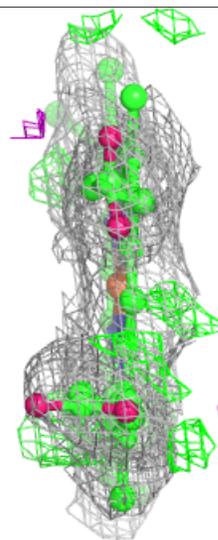
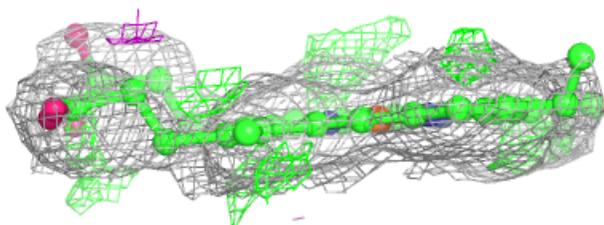
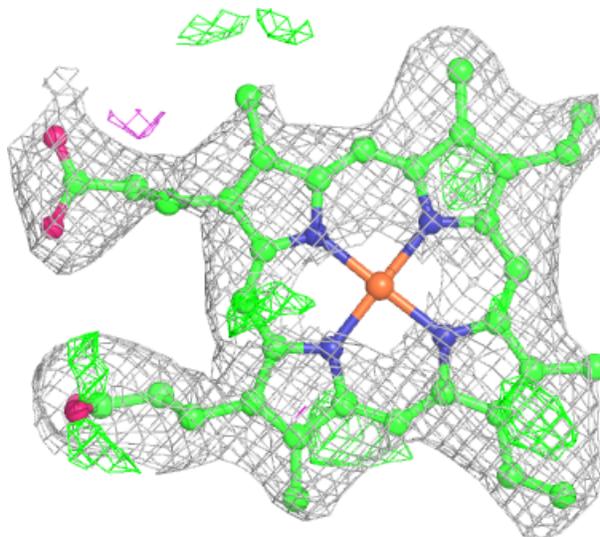
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	HEM	B	501	43/43	0.89	0.20	17,38,45,48	0
2	HEM	C	501	43/43	0.90	0.26	17,37,49,61	0
2	HEM	A	501	43/43	0.91	0.18	5,31,43,52	0
2	HEM	D	501	43/43	0.91	0.26	17,45,54,64	0
2	HEM	F	501	43/43	0.94	0.17	5,25,40,51	0
3	PLP	B	1001	15/16	0.94	0.18	2,15,25,26	0
2	HEM	E	501	43/43	0.95	0.15	8,27,42,48	0
3	PLP	C	1001	15/16	0.95	0.17	11,22,25,28	0
3	PLP	D	1001	15/16	0.95	0.15	15,22,31,32	0
3	PLP	A	1001	15/16	0.96	0.14	2,19,29,30	0
3	PLP	E	1001	15/16	0.97	0.16	2,12,18,23	0
3	PLP	F	1001	15/16	0.97	0.15	4,11,17,25	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.

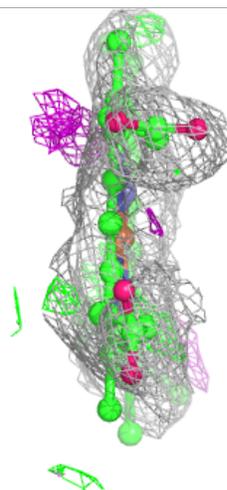
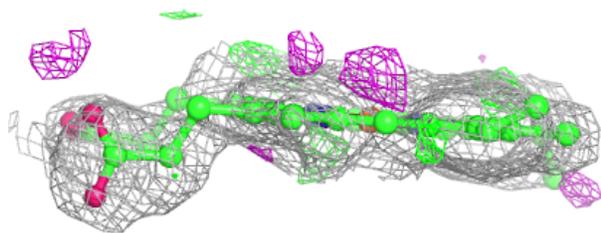
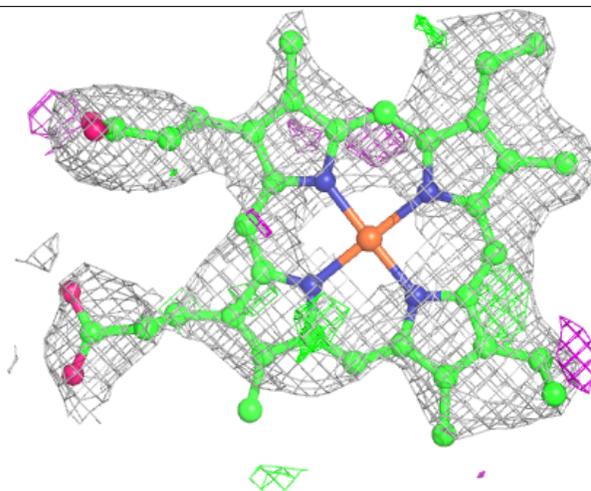
Electron density around HEM B 501:

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



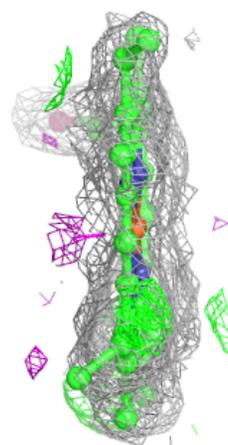
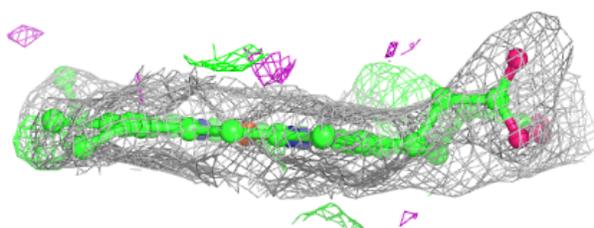
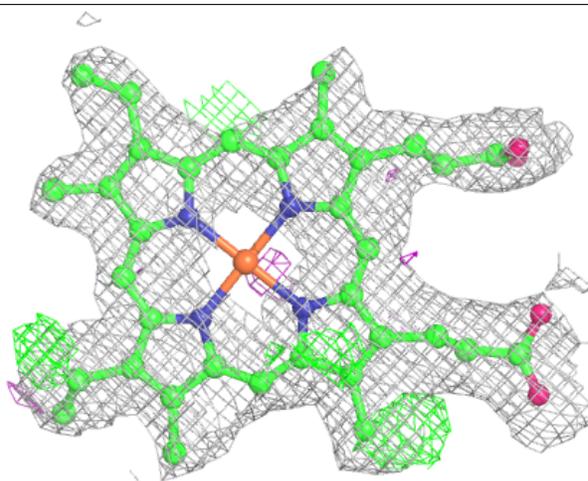
Electron density around HEM C 501:

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



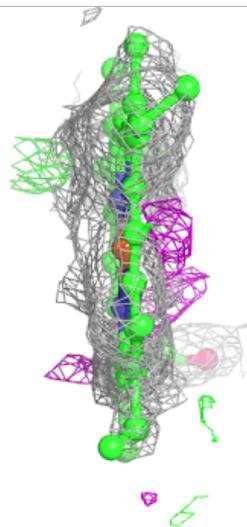
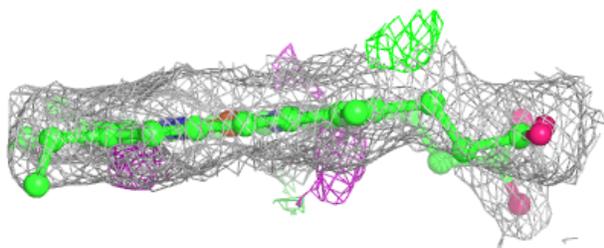
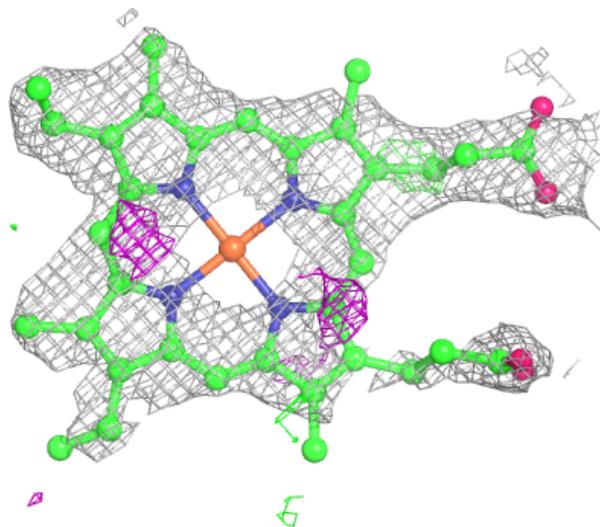
Electron density around HEM A 501:

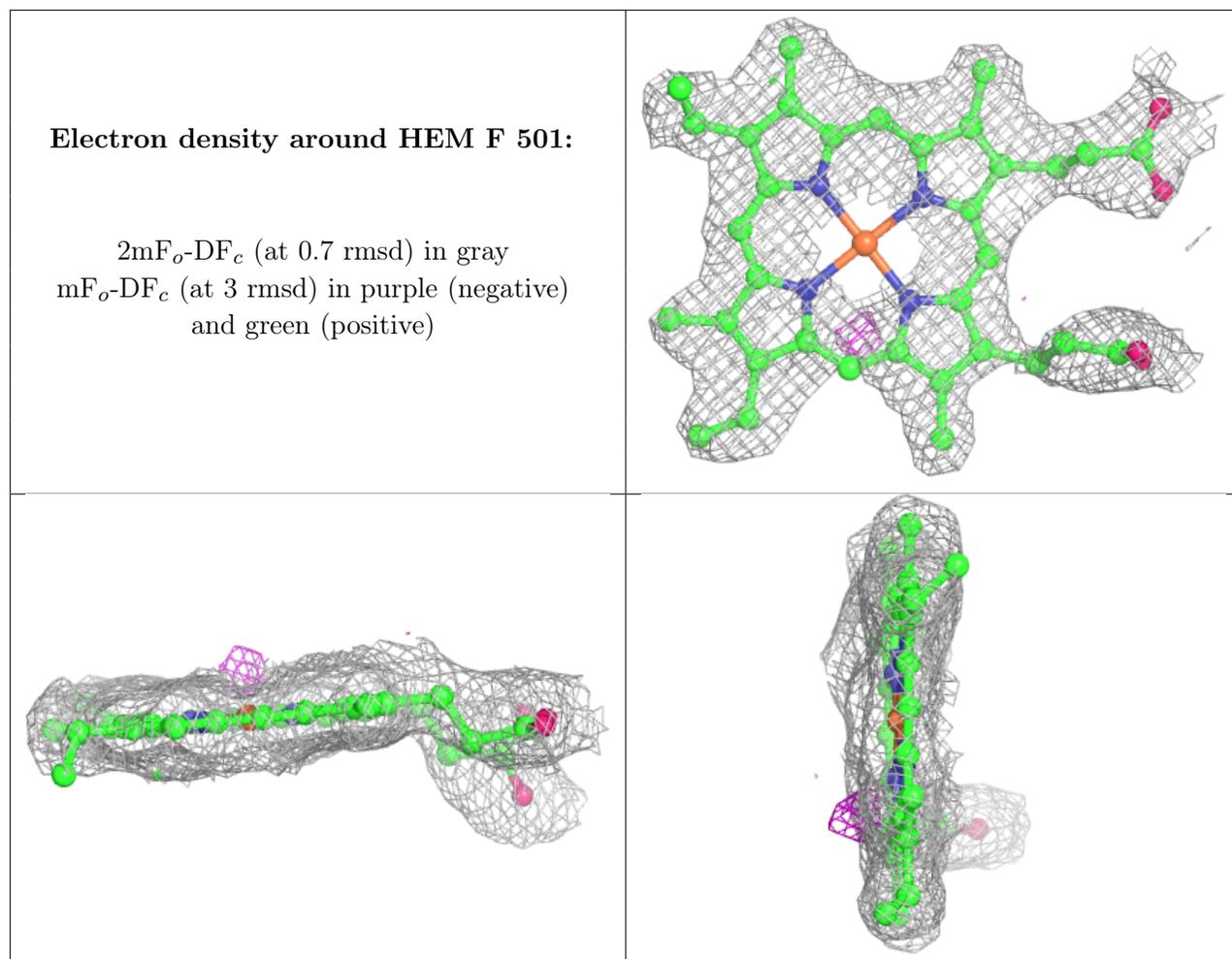
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

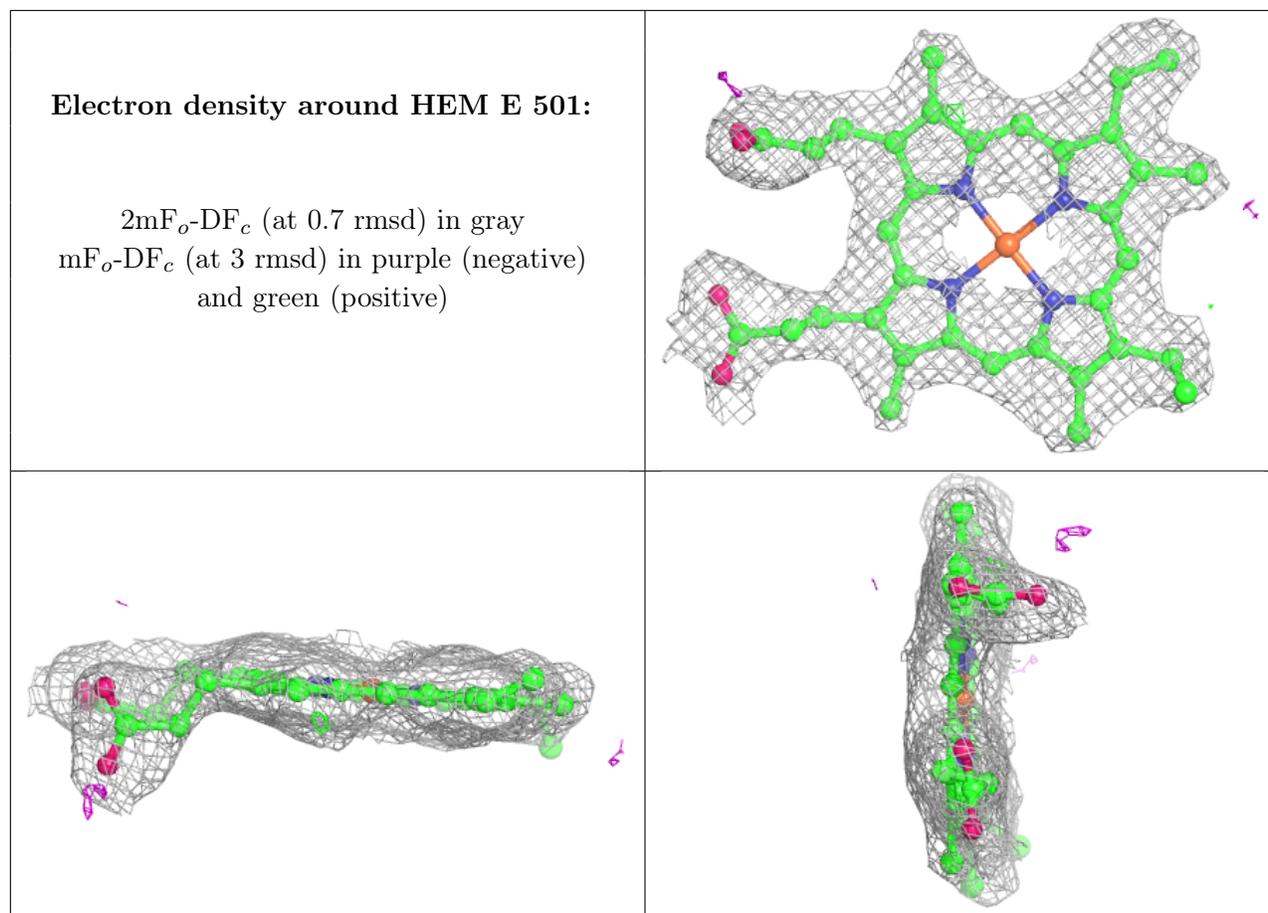


Electron density around HEM D 501:

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







6.5 Other polymers [i](#)

There are no such residues in this entry.