



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 17, 2026 – 11:55 PM UTC

PDB ID : 4ZGX / pdb_00004zgx
Title : Structure of aldosterone synthase (CYP11B2) in complex with (+)-(R)-N-(4-(4-chloro-3-fluorophenyl)-5,6,7,8-tetrahydroisoquinolin-8-yl)propionamide
Authors : Kuglstatter, A.; Joseph, C.
Deposited on : 2015-04-24
Resolution : 3.20 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

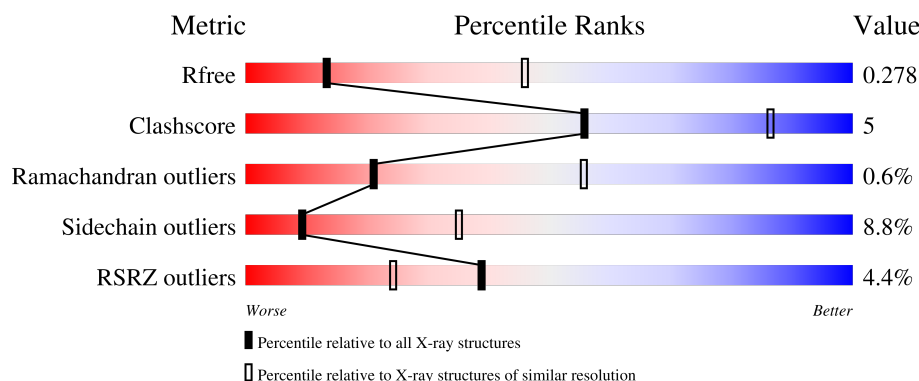
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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.



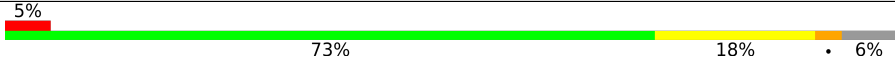
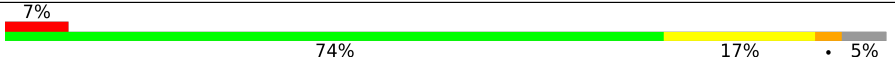
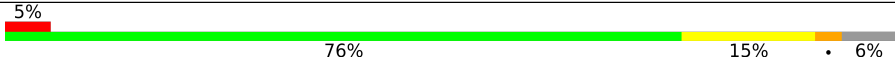
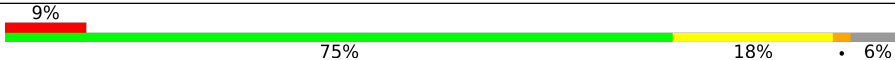
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1466 (3.20-3.20)
Clashscore	190562	1573 (3.20-3.20)
Ramachandran outliers	187476	1548 (3.20-3.20)
Sidechain outliers	187428	1547 (3.20-3.20)
RSRZ outliers	180081	1466 (3.20-3.20)

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	489	
1	B	489	
1	C	489	
1	D	489	
1	E	489	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	F	489	
1	G	489	
1	H	489	
1	I	489	
1	J	489	
1	K	489	
1	L	489	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 46675 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 11B2, mitochondrial.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	465	Total	C	N	O	S	0	0	0
			3774	2439	666	649	20			
1	B	465	Total	C	N	O	S	0	0	0
			3774	2439	666	649	20			
1	C	465	Total	C	N	O	S	0	0	0
			3774	2439	666	649	20			
1	D	465	Total	C	N	O	S	0	0	0
			3774	2439	666	649	20			
1	E	459	Total	C	N	O	S	0	0	0
			3731	2412	658	641	20			
1	F	464	Total	C	N	O	S	0	0	0
			3770	2437	665	648	20			
1	G	465	Total	C	N	O	S	0	0	0
			3774	2439	666	649	20			
1	H	465	Total	C	N	O	S	0	0	0
			3774	2439	666	649	20			
1	I	458	Total	C	N	O	S	0	0	0
			3721	2406	655	640	20			
1	J	463	Total	C	N	O	S	0	0	0
			3762	2431	664	647	20			
1	K	461	Total	C	N	O	S	0	0	0
			3737	2416	657	644	20			
1	L	462	Total	C	N	O	S	0	0	0
			3747	2422	661	644	20			

There are 156 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	24	MET	-	expression tag	UNP P19099
A	25	ALA	-	expression tag	UNP P19099
A	26	THR	-	expression tag	UNP P19099
A	27	LYS	-	expression tag	UNP P19099
A	504	GLY	-	expression tag	UNP P19099

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	505	GLY	-	expression tag	UNP P19099
A	506	ARG	-	expression tag	UNP P19099
A	507	HIS	-	expression tag	UNP P19099
A	508	HIS	-	expression tag	UNP P19099
A	509	HIS	-	expression tag	UNP P19099
A	510	HIS	-	expression tag	UNP P19099
A	511	HIS	-	expression tag	UNP P19099
A	512	HIS	-	expression tag	UNP P19099
B	24	MET	-	expression tag	UNP P19099
B	25	ALA	-	expression tag	UNP P19099
B	26	THR	-	expression tag	UNP P19099
B	27	LYS	-	expression tag	UNP P19099
B	504	GLY	-	expression tag	UNP P19099
B	505	GLY	-	expression tag	UNP P19099
B	506	ARG	-	expression tag	UNP P19099
B	507	HIS	-	expression tag	UNP P19099
B	508	HIS	-	expression tag	UNP P19099
B	509	HIS	-	expression tag	UNP P19099
B	510	HIS	-	expression tag	UNP P19099
B	511	HIS	-	expression tag	UNP P19099
B	512	HIS	-	expression tag	UNP P19099
C	24	MET	-	expression tag	UNP P19099
C	25	ALA	-	expression tag	UNP P19099
C	26	THR	-	expression tag	UNP P19099
C	27	LYS	-	expression tag	UNP P19099
C	504	GLY	-	expression tag	UNP P19099
C	505	GLY	-	expression tag	UNP P19099
C	506	ARG	-	expression tag	UNP P19099
C	507	HIS	-	expression tag	UNP P19099
C	508	HIS	-	expression tag	UNP P19099
C	509	HIS	-	expression tag	UNP P19099
C	510	HIS	-	expression tag	UNP P19099
C	511	HIS	-	expression tag	UNP P19099
C	512	HIS	-	expression tag	UNP P19099
D	24	MET	-	expression tag	UNP P19099
D	25	ALA	-	expression tag	UNP P19099
D	26	THR	-	expression tag	UNP P19099
D	27	LYS	-	expression tag	UNP P19099
D	504	GLY	-	expression tag	UNP P19099
D	505	GLY	-	expression tag	UNP P19099
D	506	ARG	-	expression tag	UNP P19099
D	507	HIS	-	expression tag	UNP P19099

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
D	508	HIS	-	expression tag	UNP P19099
D	509	HIS	-	expression tag	UNP P19099
D	510	HIS	-	expression tag	UNP P19099
D	511	HIS	-	expression tag	UNP P19099
D	512	HIS	-	expression tag	UNP P19099
E	24	MET	-	expression tag	UNP P19099
E	25	ALA	-	expression tag	UNP P19099
E	26	THR	-	expression tag	UNP P19099
E	27	LYS	-	expression tag	UNP P19099
E	504	GLY	-	expression tag	UNP P19099
E	505	GLY	-	expression tag	UNP P19099
E	506	ARG	-	expression tag	UNP P19099
E	507	HIS	-	expression tag	UNP P19099
E	508	HIS	-	expression tag	UNP P19099
E	509	HIS	-	expression tag	UNP P19099
E	510	HIS	-	expression tag	UNP P19099
E	511	HIS	-	expression tag	UNP P19099
E	512	HIS	-	expression tag	UNP P19099
F	24	MET	-	expression tag	UNP P19099
F	25	ALA	-	expression tag	UNP P19099
F	26	THR	-	expression tag	UNP P19099
F	27	LYS	-	expression tag	UNP P19099
F	504	GLY	-	expression tag	UNP P19099
F	505	GLY	-	expression tag	UNP P19099
F	506	ARG	-	expression tag	UNP P19099
F	507	HIS	-	expression tag	UNP P19099
F	508	HIS	-	expression tag	UNP P19099
F	509	HIS	-	expression tag	UNP P19099
F	510	HIS	-	expression tag	UNP P19099
F	511	HIS	-	expression tag	UNP P19099
F	512	HIS	-	expression tag	UNP P19099
G	24	MET	-	expression tag	UNP P19099
G	25	ALA	-	expression tag	UNP P19099
G	26	THR	-	expression tag	UNP P19099
G	27	LYS	-	expression tag	UNP P19099
G	504	GLY	-	expression tag	UNP P19099
G	505	GLY	-	expression tag	UNP P19099
G	506	ARG	-	expression tag	UNP P19099
G	507	HIS	-	expression tag	UNP P19099
G	508	HIS	-	expression tag	UNP P19099
G	509	HIS	-	expression tag	UNP P19099
G	510	HIS	-	expression tag	UNP P19099

Continued on next page...

Continued from previous page...

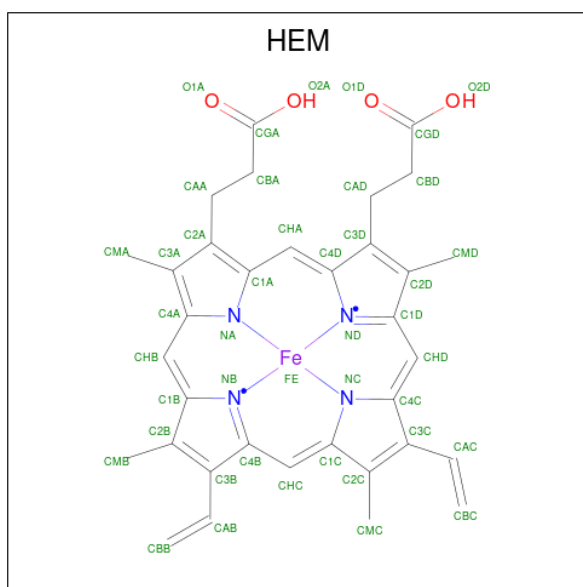
Chain	Residue	Modelled	Actual	Comment	Reference
G	511	HIS	-	expression tag	UNP P19099
G	512	HIS	-	expression tag	UNP P19099
H	24	MET	-	expression tag	UNP P19099
H	25	ALA	-	expression tag	UNP P19099
H	26	THR	-	expression tag	UNP P19099
H	27	LYS	-	expression tag	UNP P19099
H	504	GLY	-	expression tag	UNP P19099
H	505	GLY	-	expression tag	UNP P19099
H	506	ARG	-	expression tag	UNP P19099
H	507	HIS	-	expression tag	UNP P19099
H	508	HIS	-	expression tag	UNP P19099
H	509	HIS	-	expression tag	UNP P19099
H	510	HIS	-	expression tag	UNP P19099
H	511	HIS	-	expression tag	UNP P19099
H	512	HIS	-	expression tag	UNP P19099
I	24	MET	-	expression tag	UNP P19099
I	25	ALA	-	expression tag	UNP P19099
I	26	THR	-	expression tag	UNP P19099
I	27	LYS	-	expression tag	UNP P19099
I	504	GLY	-	expression tag	UNP P19099
I	505	GLY	-	expression tag	UNP P19099
I	506	ARG	-	expression tag	UNP P19099
I	507	HIS	-	expression tag	UNP P19099
I	508	HIS	-	expression tag	UNP P19099
I	509	HIS	-	expression tag	UNP P19099
I	510	HIS	-	expression tag	UNP P19099
I	511	HIS	-	expression tag	UNP P19099
I	512	HIS	-	expression tag	UNP P19099
J	24	MET	-	expression tag	UNP P19099
J	25	ALA	-	expression tag	UNP P19099
J	26	THR	-	expression tag	UNP P19099
J	27	LYS	-	expression tag	UNP P19099
J	504	GLY	-	expression tag	UNP P19099
J	505	GLY	-	expression tag	UNP P19099
J	506	ARG	-	expression tag	UNP P19099
J	507	HIS	-	expression tag	UNP P19099
J	508	HIS	-	expression tag	UNP P19099
J	509	HIS	-	expression tag	UNP P19099
J	510	HIS	-	expression tag	UNP P19099
J	511	HIS	-	expression tag	UNP P19099
J	512	HIS	-	expression tag	UNP P19099
K	24	MET	-	expression tag	UNP P19099

Continued on next page...

Continued from previous page...

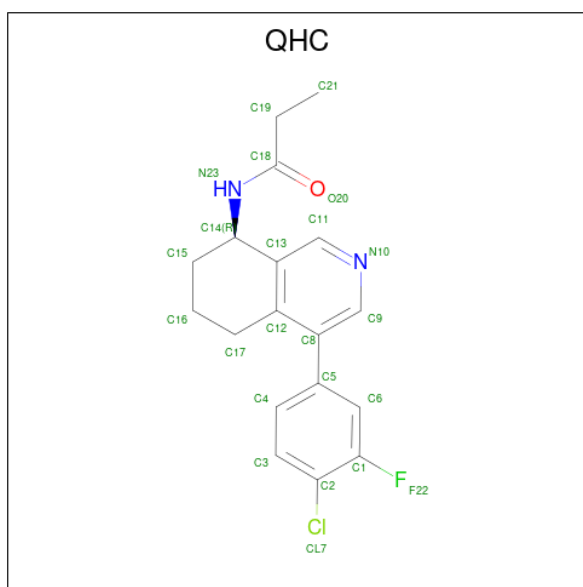
Chain	Residue	Modelled	Actual	Comment	Reference
K	25	ALA	-	expression tag	UNP P19099
K	26	THR	-	expression tag	UNP P19099
K	27	LYS	-	expression tag	UNP P19099
K	504	GLY	-	expression tag	UNP P19099
K	505	GLY	-	expression tag	UNP P19099
K	506	ARG	-	expression tag	UNP P19099
K	507	HIS	-	expression tag	UNP P19099
K	508	HIS	-	expression tag	UNP P19099
K	509	HIS	-	expression tag	UNP P19099
K	510	HIS	-	expression tag	UNP P19099
K	511	HIS	-	expression tag	UNP P19099
K	512	HIS	-	expression tag	UNP P19099
L	24	MET	-	expression tag	UNP P19099
L	25	ALA	-	expression tag	UNP P19099
L	26	THR	-	expression tag	UNP P19099
L	27	LYS	-	expression tag	UNP P19099
L	504	GLY	-	expression tag	UNP P19099
L	505	GLY	-	expression tag	UNP P19099
L	506	ARG	-	expression tag	UNP P19099
L	507	HIS	-	expression tag	UNP P19099
L	508	HIS	-	expression tag	UNP P19099
L	509	HIS	-	expression tag	UNP P19099
L	510	HIS	-	expression tag	UNP P19099
L	511	HIS	-	expression tag	UNP P19099
L	512	HIS	-	expression tag	UNP P19099

- 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
2	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	D	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	F	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	G	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	H	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	I	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	J	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	K	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	L	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is N-[(8R)-4-(4-chloro-3-fluorophenyl)-5,6,7,8-tetrahydroisoquinolin-8-yl]propanamide (CCD ID: QHC) (formula: C₁₈H₁₈ClFN₂O).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	B	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	C	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	D	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	E	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	F	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	I	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	J	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	K	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0
3	L	1	Total 23	C 18	Cl 1	F 1	N 2	O 1	0	0

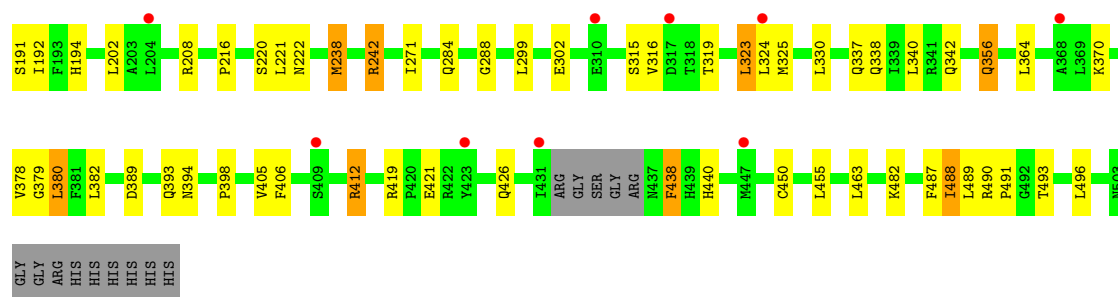
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	108	Total	O	0	0
			108	108		
4	B	81	Total	O	0	0
			81	81		

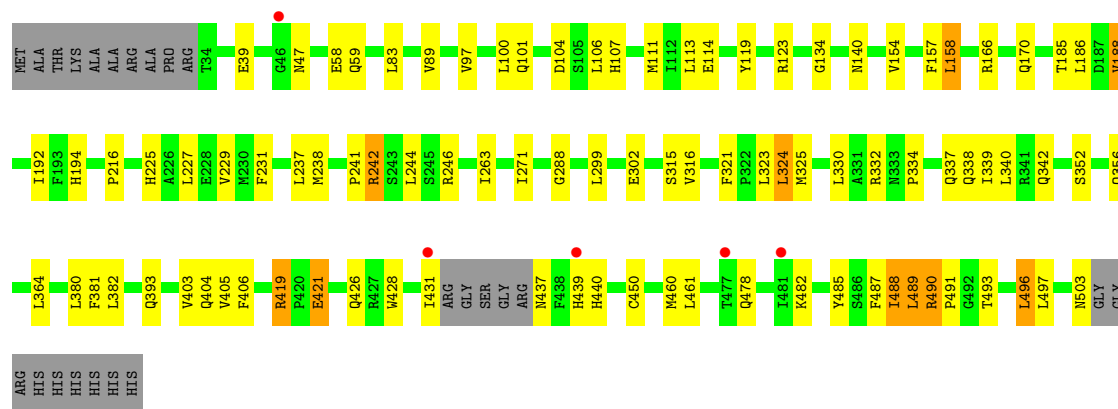
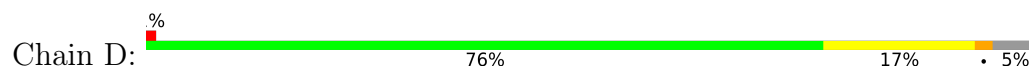
Continued on next page...

Continued from previous page...

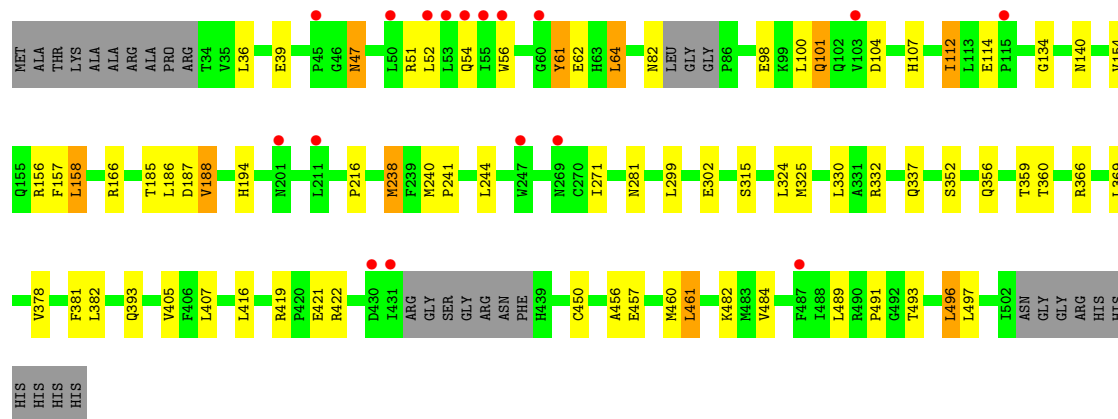
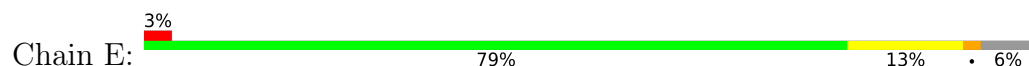
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	C	64	Total 64	O 64	0	0
4	D	100	Total 100	O 100	0	0
4	E	54	Total 54	O 54	0	0
4	F	73	Total 73	O 73	0	0
4	G	45	Total 45	O 45	0	0
4	H	68	Total 68	O 68	0	0
4	I	54	Total 54	O 54	0	0
4	J	50	Total 50	O 50	0	0
4	K	53	Total 53	O 53	0	0
4	L	67	Total 67	O 67	0	0



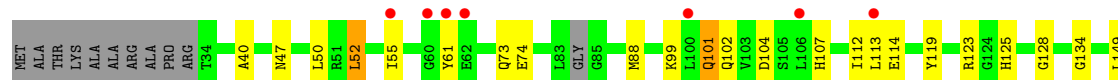
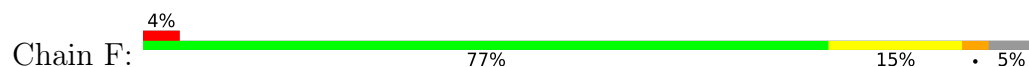
- Molecule 1: Cytochrome P450 11B2, mitochondrial

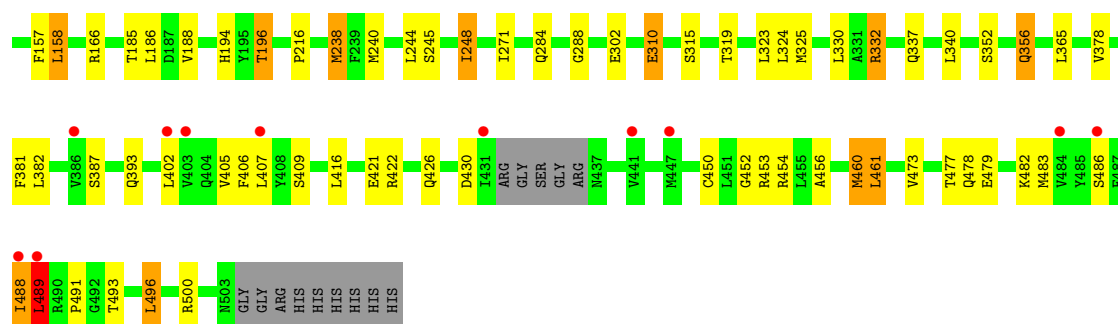


- Molecule 1: Cytochrome P450 11B2, mitochondrial



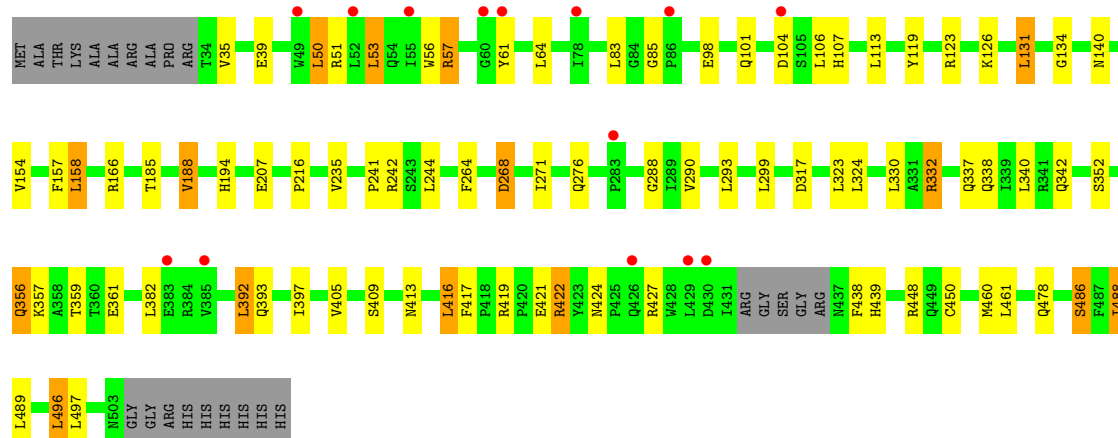
- Molecule 1: Cytochrome P450 11B2, mitochondrial





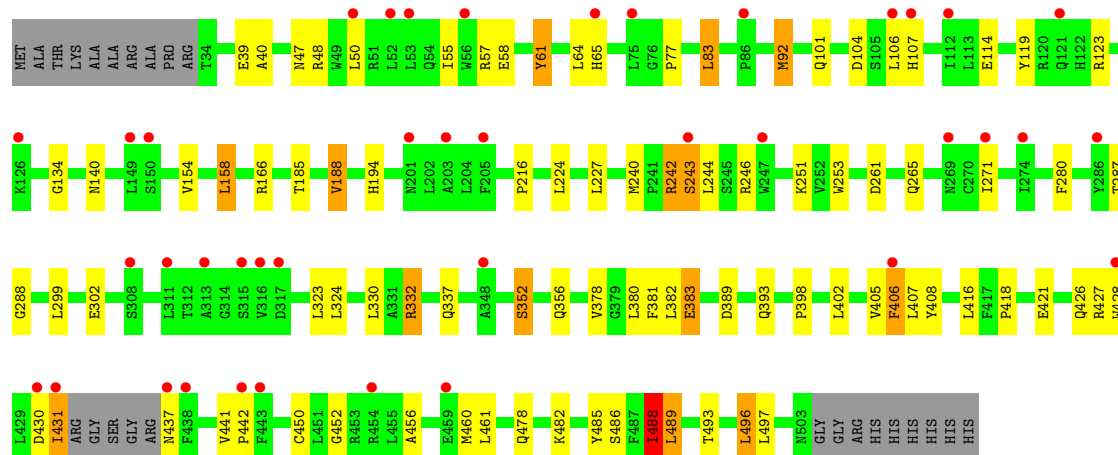
• Molecule 1: Cytochrome P450 11B2, mitochondrial

Chain G: 3% 78% 14% 5%



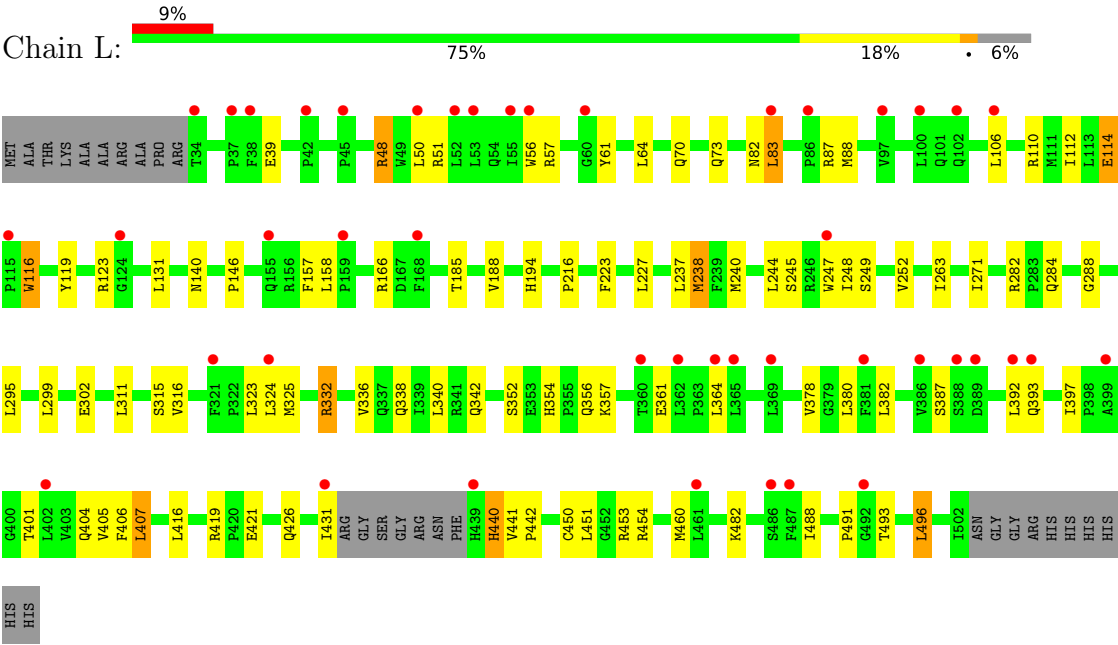
• Molecule 1: Cytochrome P450 11B2, mitochondrial

Chain H: 8% 76% 16% 5%



• Molecule 1: Cytochrome P450 11B2, mitochondrial

Chain I: 5% 73% 18% 6%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	104.70Å 125.96Å 155.23Å 70.25° 85.87° 73.84°	Depositor
Resolution (Å)	88.20 – 3.20 88.20 – 3.20	Depositor EDS
% Data completeness (in resolution range)	97.2 (88.20-3.20) 97.2 (88.20-3.20)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.70 (at 3.19Å)	Xtriage
Refinement program	BUSTER 2.11.4	Depositor
R, R_{free}	0.212 , 0.255 0.234 , 0.278	Depositor DCC
R_{free} test set	5769 reflections (4.88%)	wwPDB-VP
Wilson B-factor (Å ²)	60.1	Xtriage
Anisotropy	0.246	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 79.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	46675	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 21.13 % 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 7.4821e-03. 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 ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, QHC

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	0.80	0/3874	1.35	15/5257 (0.3%)
1	B	0.76	0/3874	1.33	9/5257 (0.2%)
1	C	0.77	0/3874	1.36	18/5257 (0.3%)
1	D	0.79	0/3874	1.33	15/5257 (0.3%)
1	E	0.78	0/3829	1.34	13/5194 (0.3%)
1	F	0.77	0/3869	1.35	13/5249 (0.2%)
1	G	0.77	0/3874	1.33	10/5257 (0.2%)
1	H	0.76	0/3874	1.34	8/5257 (0.2%)
1	I	0.77	0/3818	1.35	16/5179 (0.3%)
1	J	0.76	0/3861	1.34	10/5238 (0.2%)
1	K	0.76	0/3834	1.34	18/5201 (0.3%)
1	L	0.78	0/3846	1.35	18/5219 (0.3%)
All	All	0.77	0/46301	1.34	163/62822 (0.3%)

There are no bond length outliers.

The worst 5 of 163 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	K	381	PHE	CA-CB-CG	8.60	122.40	113.80
1	I	55	ILE	N-CA-C	-7.58	105.04	111.56
1	K	394	ASN	CA-CB-CG	7.06	119.66	112.60
1	F	40	ALA	N-CA-C	-6.99	105.37	114.04
1	E	54	GLN	N-CA-C	-6.91	103.53	112.23

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	3774	0	3808	35	0
1	B	3774	0	3808	24	0
1	C	3774	0	3808	34	0
1	D	3774	0	3808	40	0
1	E	3731	0	3770	30	0
1	F	3770	0	3804	38	0
1	G	3774	0	3809	30	0
1	H	3774	0	3809	42	0
1	I	3721	0	3763	41	0
1	J	3762	0	3793	30	0
1	K	3737	0	3775	39	0
1	L	3747	0	3787	42	0
2	A	43	0	30	6	0
2	B	43	0	30	5	0
2	C	43	0	30	5	0
2	D	43	0	30	2	0
2	E	43	0	30	4	0
2	F	43	0	30	8	0
2	G	43	0	30	1	0
2	H	43	0	30	7	0
2	I	43	0	30	3	0
2	J	43	0	30	1	0
2	K	43	0	30	5	0
2	L	43	0	30	4	0
3	A	23	0	18	7	0
3	B	23	0	18	5	0
3	C	23	0	18	5	0
3	D	23	0	18	0	0
3	E	23	0	18	2	0
3	F	23	0	18	5	0
3	I	23	0	18	4	0
3	J	23	0	18	4	0
3	K	23	0	18	5	0
3	L	23	0	18	3	0
4	A	108	0	0	0	0
4	B	81	0	0	0	0
4	C	64	0	0	1	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	100	0	0	0	0
4	E	54	0	0	1	0
4	F	73	0	0	1	0
4	G	45	0	0	0	0
4	H	68	0	0	0	0
4	I	54	0	0	0	0
4	J	50	0	0	0	0
4	K	53	0	0	1	0
4	L	67	0	0	2	0
All	All	46675	0	46082	432	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 432 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:K:52:LEU:HD11	1:K:240:MET:HB3	1.33	1.08
1:K:323:LEU:HG	1:K:460:MET:HG2	1.46	0.97
1:C:323:LEU:HD21	1:C:463:LEU:HD22	1.54	0.89
2:K:601:HEM:HMA3	3:K:602:QHC:H14	1.60	0.84
1:A:242:ARG:HD3	1:A:246:ARG:HD2	1.60	0.83

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	461/489 (94%)	447 (97%)	13 (3%)	1 (0%)	43 73
1	B	461/489 (94%)	446 (97%)	14 (3%)	1 (0%)	43 73

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	461/489 (94%)	448 (97%)	11 (2%)	2 (0%)	30	62
1	D	461/489 (94%)	441 (96%)	18 (4%)	2 (0%)	30	62
1	E	453/489 (93%)	440 (97%)	12 (3%)	1 (0%)	43	73
1	F	458/489 (94%)	441 (96%)	15 (3%)	2 (0%)	30	62
1	G	461/489 (94%)	440 (95%)	16 (4%)	5 (1%)	11	43
1	H	461/489 (94%)	440 (95%)	17 (4%)	4 (1%)	14	47
1	I	452/489 (92%)	433 (96%)	16 (4%)	3 (1%)	18	52
1	J	457/489 (94%)	433 (95%)	17 (4%)	7 (2%)	8	37
1	K	455/489 (93%)	437 (96%)	15 (3%)	3 (1%)	18	52
1	L	458/489 (94%)	436 (95%)	18 (4%)	4 (1%)	14	47
All	All	5499/5868 (94%)	5282 (96%)	182 (3%)	35 (1%)	21	56

5 of 35 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	59	GLN
1	G	242	ARG
1	J	105	SER
1	L	57	ARG
1	F	489	LEU

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	410/426 (96%)	375 (92%)	35 (8%)	10	37
1	B	410/426 (96%)	379 (92%)	31 (8%)	12	42
1	C	410/426 (96%)	376 (92%)	34 (8%)	10	38
1	D	410/426 (96%)	375 (92%)	35 (8%)	10	37
1	E	406/426 (95%)	377 (93%)	29 (7%)	13	44
1	F	410/426 (96%)	369 (90%)	41 (10%)	7	30

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	G	410/426 (96%)	376 (92%)	34 (8%)	10	38
1	H	410/426 (96%)	372 (91%)	38 (9%)	8	33
1	I	405/426 (95%)	368 (91%)	37 (9%)	9	34
1	J	409/426 (96%)	361 (88%)	48 (12%)	5	23
1	K	407/426 (96%)	368 (90%)	39 (10%)	8	32
1	L	407/426 (96%)	377 (93%)	30 (7%)	13	43
All	All	4904/5112 (96%)	4473 (91%)	431 (9%)	9	35

5 of 431 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	421	GLU
1	I	81	TYR
1	K	493	THR
1	G	496	LEU
1	H	302	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 130 such sidechains are listed below:

Mol	Chain	Res	Type
1	K	354	HIS
1	L	43	GLN
1	E	440	HIS
1	E	393	GLN
1	L	121	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

22 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	D	601	3,1	50,50,50	1.40	8 (16%)	67,82,82	1.46	12 (17%)
2	HEM	I	601	3,1	50,50,50	1.36	6 (12%)	67,82,82	1.59	15 (22%)
3	QHC	K	602	2	25,25,25	0.76	0	31,35,35	1.40	4 (12%)
3	QHC	L	602	2	25,25,25	0.74	0	31,35,35	1.47	7 (22%)
2	HEM	E	601	3,1	50,50,50	1.69	9 (18%)	67,82,82	1.57	14 (20%)
3	QHC	E	602	2	25,25,25	0.80	0	31,35,35	1.81	10 (32%)
2	HEM	J	601	3	50,50,50	1.34	6 (12%)	67,82,82	1.56	15 (22%)
2	HEM	K	601	3,1	50,50,50	1.38	7 (14%)	67,82,82	1.44	10 (14%)
2	HEM	A	601	3,1	50,50,50	1.36	7 (14%)	67,82,82	1.45	9 (13%)
2	HEM	H	601	-	50,50,50	1.29	6 (12%)	67,82,82	1.65	18 (26%)
3	QHC	C	602	2	25,25,25	0.78	0	31,35,35	1.47	5 (16%)
3	QHC	B	602	2	25,25,25	1.01	1 (4%)	31,35,35	1.67	6 (19%)
3	QHC	A	602	2	25,25,25	0.73	0	31,35,35	1.81	11 (35%)
3	QHC	J	602	2	25,25,25	0.84	1 (4%)	31,35,35	1.83	7 (22%)
3	QHC	F	602	2	25,25,25	0.82	0	31,35,35	2.23	8 (25%)
2	HEM	L	601	3,1	50,50,50	1.23	5 (10%)	67,82,82	1.60	14 (20%)
2	HEM	C	601	3,1	50,50,50	1.53	10 (20%)	67,82,82	1.49	10 (14%)
3	QHC	I	602	2	25,25,25	0.93	1 (4%)	31,35,35	1.57	5 (16%)
3	QHC	D	602	2	25,25,25	0.86	1 (4%)	31,35,35	1.85	7 (22%)
2	HEM	G	601	-	50,50,50	1.40	8 (16%)	67,82,82	1.65	15 (22%)
2	HEM	F	601	3,1	50,50,50	1.43	10 (20%)	67,82,82	1.22	7 (10%)
2	HEM	B	601	3,1	50,50,50	1.32	6 (12%)	67,82,82	1.30	9 (13%)

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	D	601	3,1	-	2/14/54/54	-
2	HEM	I	601	3,1	-	5/14/54/54	-
3	QHC	K	602	2	-	2/10/20/20	0/3/3/3
3	QHC	L	602	2	-	2/10/20/20	0/3/3/3
2	HEM	E	601	3,1	-	6/14/54/54	-
3	QHC	E	602	2	-	2/10/20/20	0/3/3/3
2	HEM	J	601	3	-	3/14/54/54	-
2	HEM	K	601	3,1	-	4/14/54/54	-
2	HEM	A	601	3,1	-	6/14/54/54	-
2	HEM	H	601	-	-	4/14/54/54	-
3	QHC	C	602	2	-	1/10/20/20	0/3/3/3
3	QHC	B	602	2	-	3/10/20/20	0/3/3/3
3	QHC	A	602	2	-	2/10/20/20	0/3/3/3
3	QHC	J	602	2	-	1/10/20/20	0/3/3/3
3	QHC	F	602	2	-	1/10/20/20	0/3/3/3
2	HEM	L	601	3,1	-	9/14/54/54	-
2	HEM	C	601	3,1	-	5/14/54/54	-
3	QHC	I	602	2	-	2/10/20/20	0/3/3/3
3	QHC	D	602	2	-	1/10/20/20	0/3/3/3
2	HEM	G	601	-	-	2/14/54/54	-
2	HEM	F	601	3,1	-	5/14/54/54	-
2	HEM	B	601	3,1	-	7/14/54/54	-

The worst 5 of 92 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	601	HEM	C4D-ND	-5.33	1.31	1.40
2	I	601	HEM	C3B-C4B	4.34	1.53	1.44
2	E	601	HEM	C1B-NB	-4.31	1.32	1.40
2	D	601	HEM	C1D-C2D	4.27	1.53	1.44
2	A	601	HEM	C1D-C2D	4.00	1.52	1.44

The worst 5 of 218 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	602	QHC	C15-C14-N23	7.00	117.44	110.64
3	F	602	QHC	C14-N23-C18	6.20	132.69	123.44
3	I	602	QHC	C14-N23-C18	5.70	131.94	123.44
3	B	602	QHC	C14-N23-C18	5.38	131.47	123.44

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	601	HEM	C2B-C1B-NB	5.23	115.85	109.84

There are no chirality outliers.

5 of 75 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	HEM	C2C-C3C-CAC-CBC
2	A	601	HEM	C4C-C3C-CAC-CBC
2	E	601	HEM	C2B-C3B-CAB-CBB
2	H	601	HEM	C2C-C3C-CAC-CBC
2	L	601	HEM	C2C-C3C-CAC-CBC

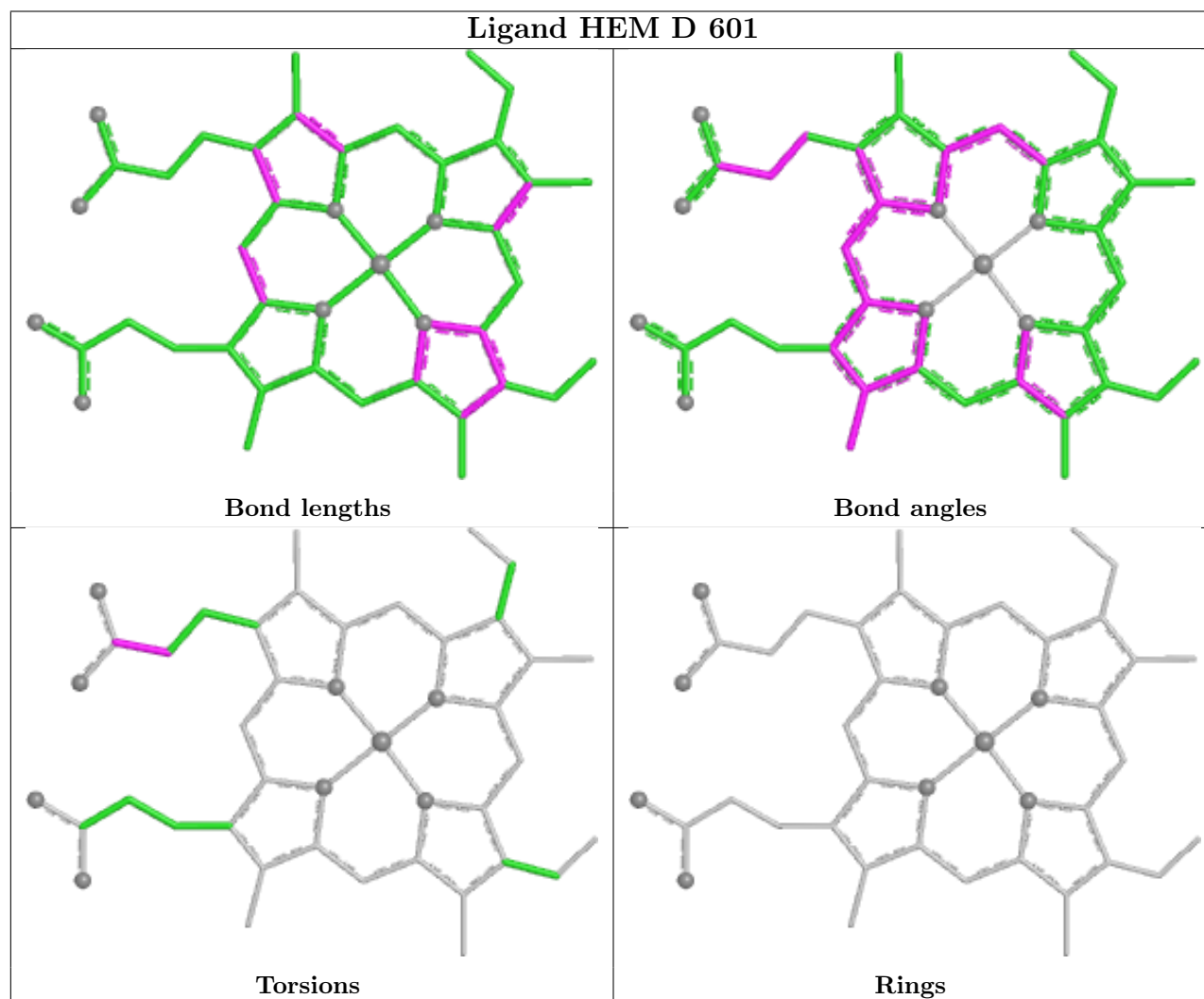
There are no ring outliers.

21 monomers are involved in 76 short contacts:

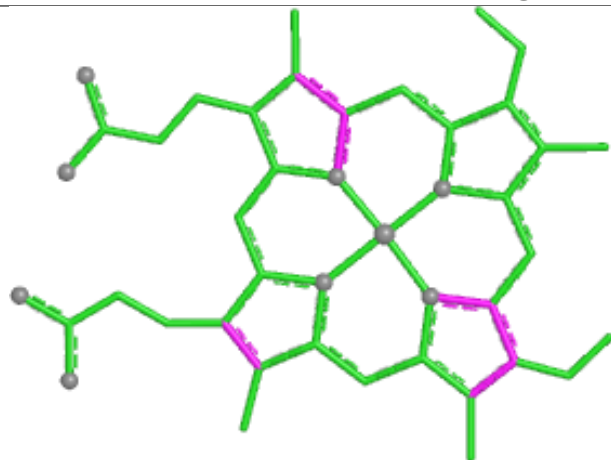
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	601	HEM	2	0
2	I	601	HEM	3	0
3	K	602	QHC	5	0
3	L	602	QHC	3	0
2	E	601	HEM	4	0
3	E	602	QHC	2	0
2	J	601	HEM	1	0
2	K	601	HEM	5	0
2	A	601	HEM	6	0
2	H	601	HEM	7	0
3	C	602	QHC	5	0
3	B	602	QHC	5	0
3	A	602	QHC	7	0
3	J	602	QHC	4	0
3	F	602	QHC	5	0
2	L	601	HEM	4	0
2	C	601	HEM	5	0
3	I	602	QHC	4	0
2	G	601	HEM	1	0
2	F	601	HEM	8	0
2	B	601	HEM	5	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

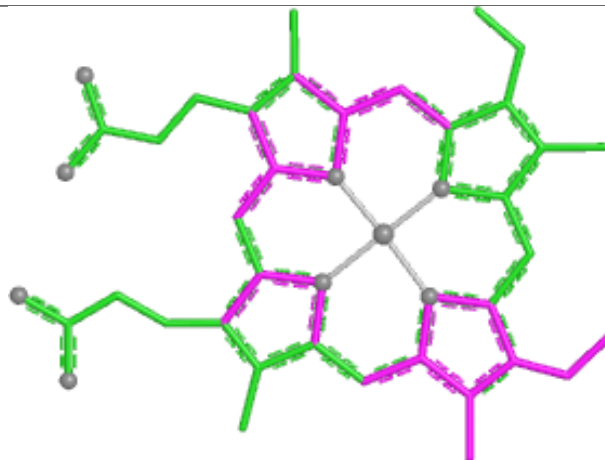
within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question 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.



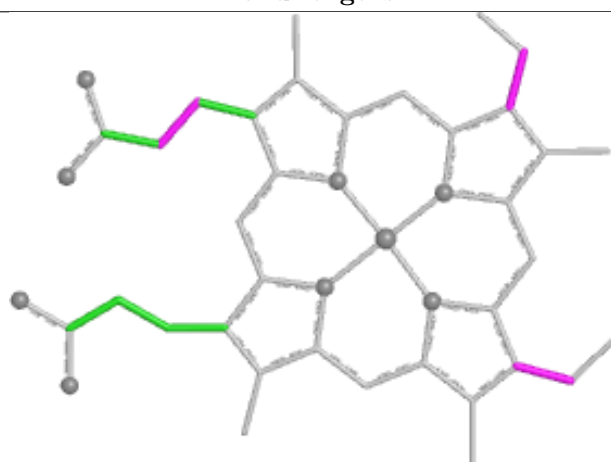
Ligand HEM I 601



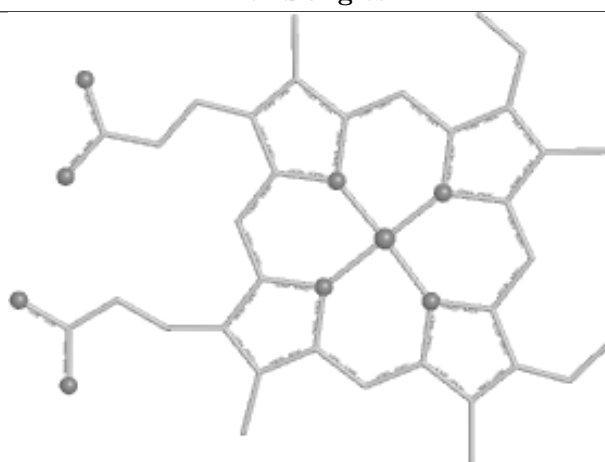
Bond lengths



Bond angles

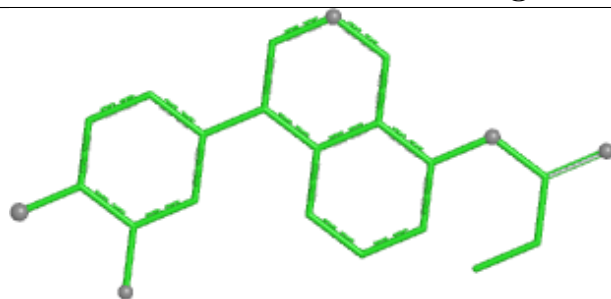


Torsions

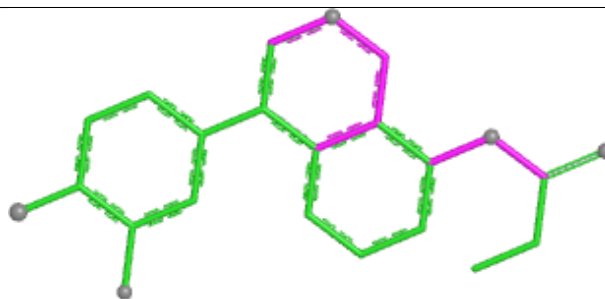


Rings

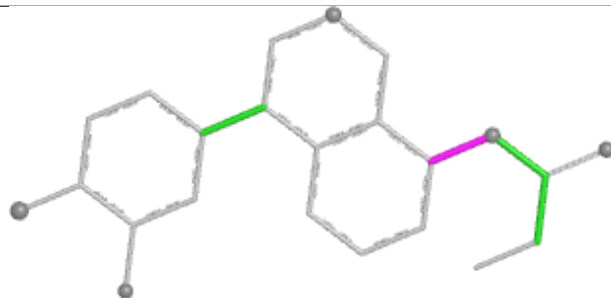
Ligand QHC K 602



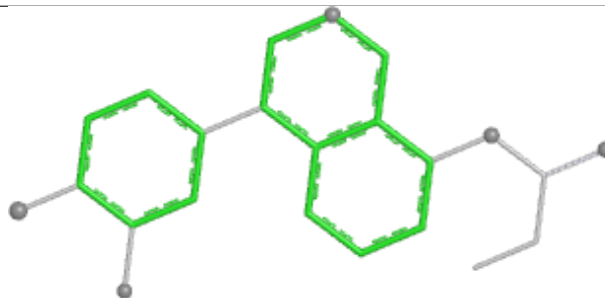
Bond lengths



Bond angles

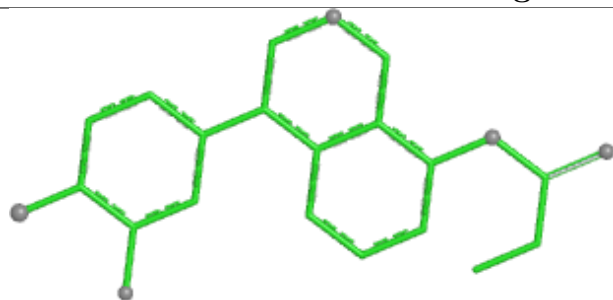


Torsions

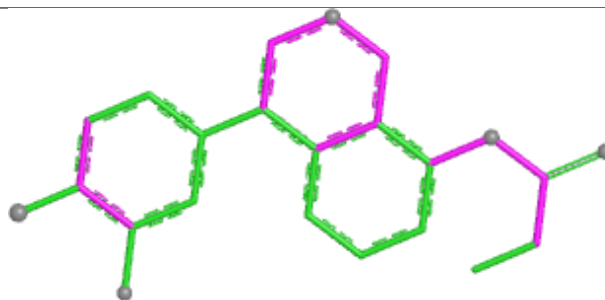


Rings

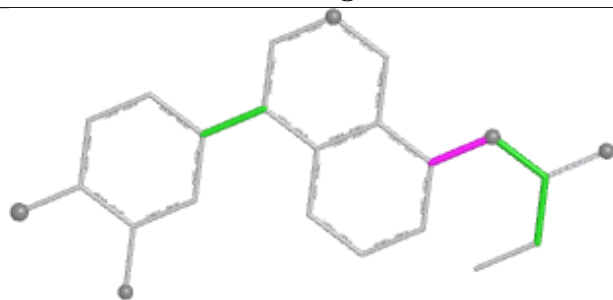
Ligand QHC L 602



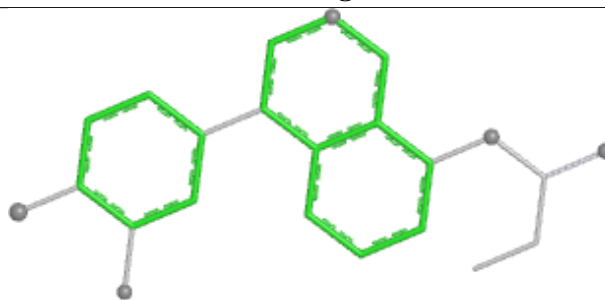
Bond lengths



Bond angles

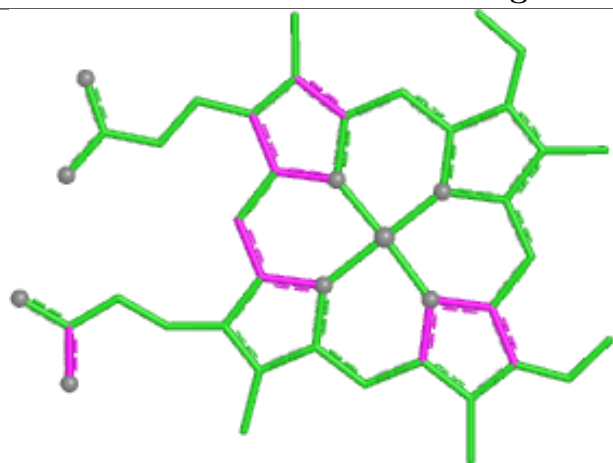


Torsions

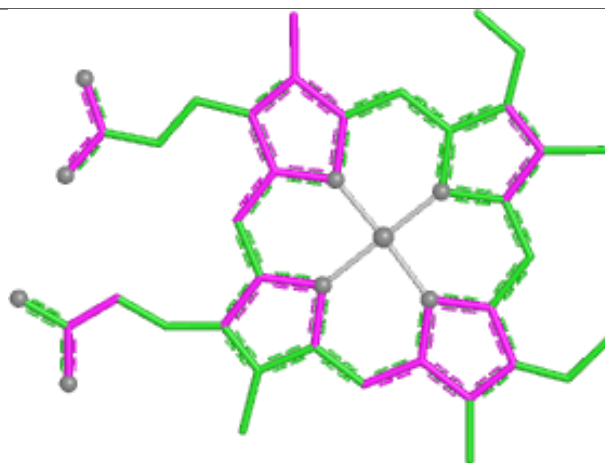


Rings

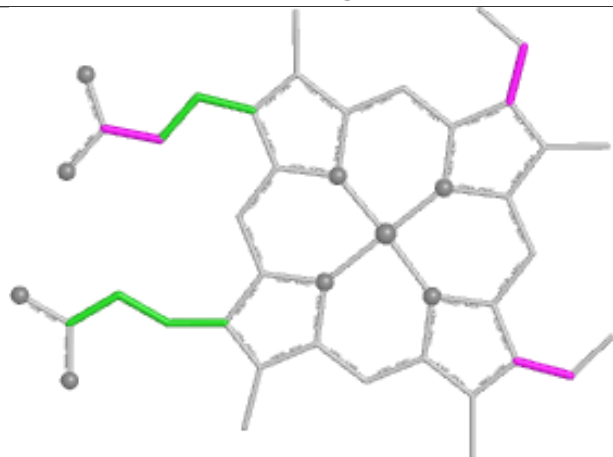
Ligand HEM E 601



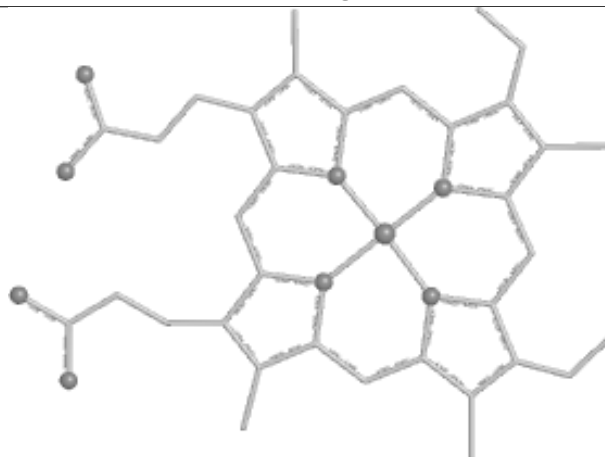
Bond lengths



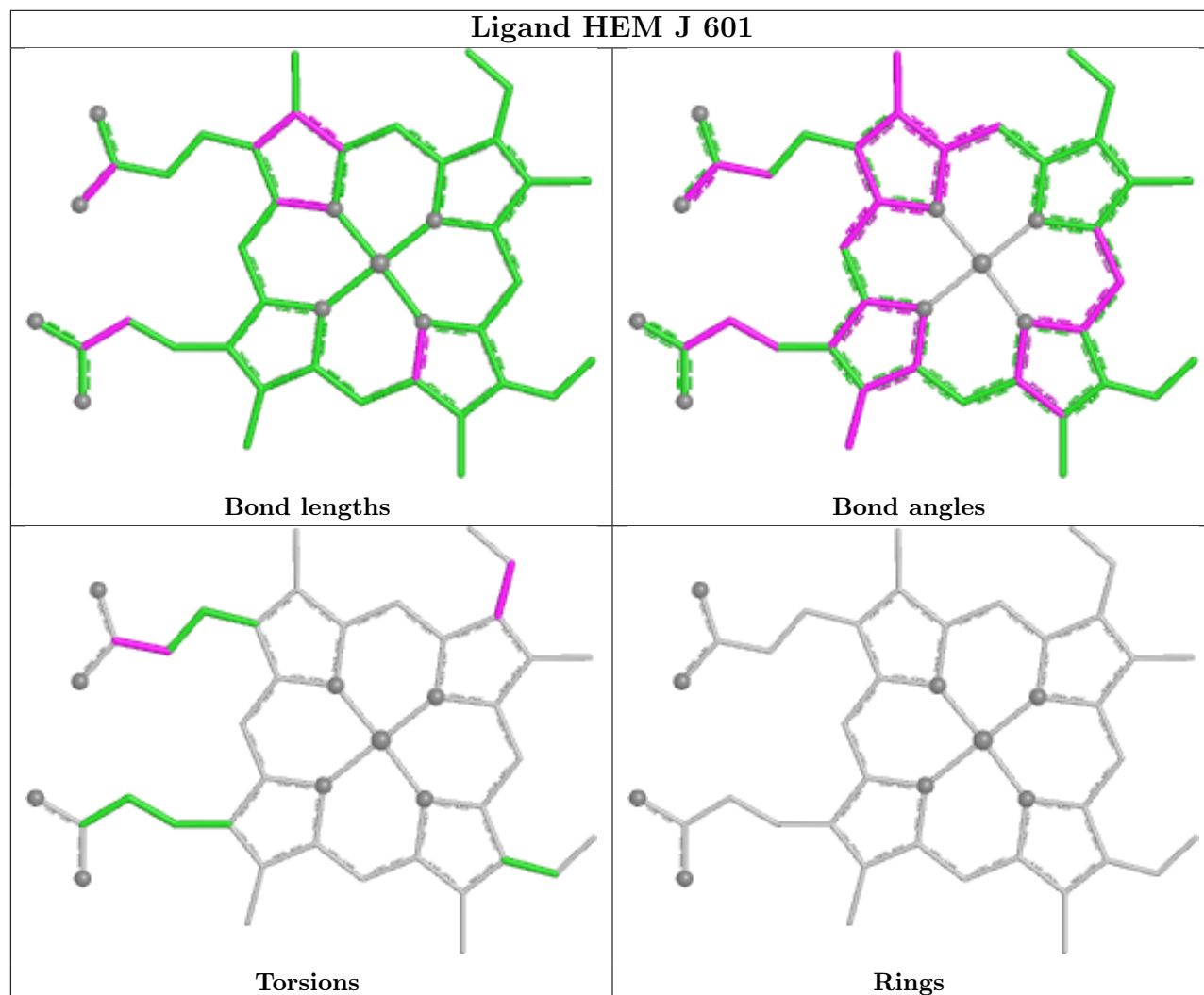
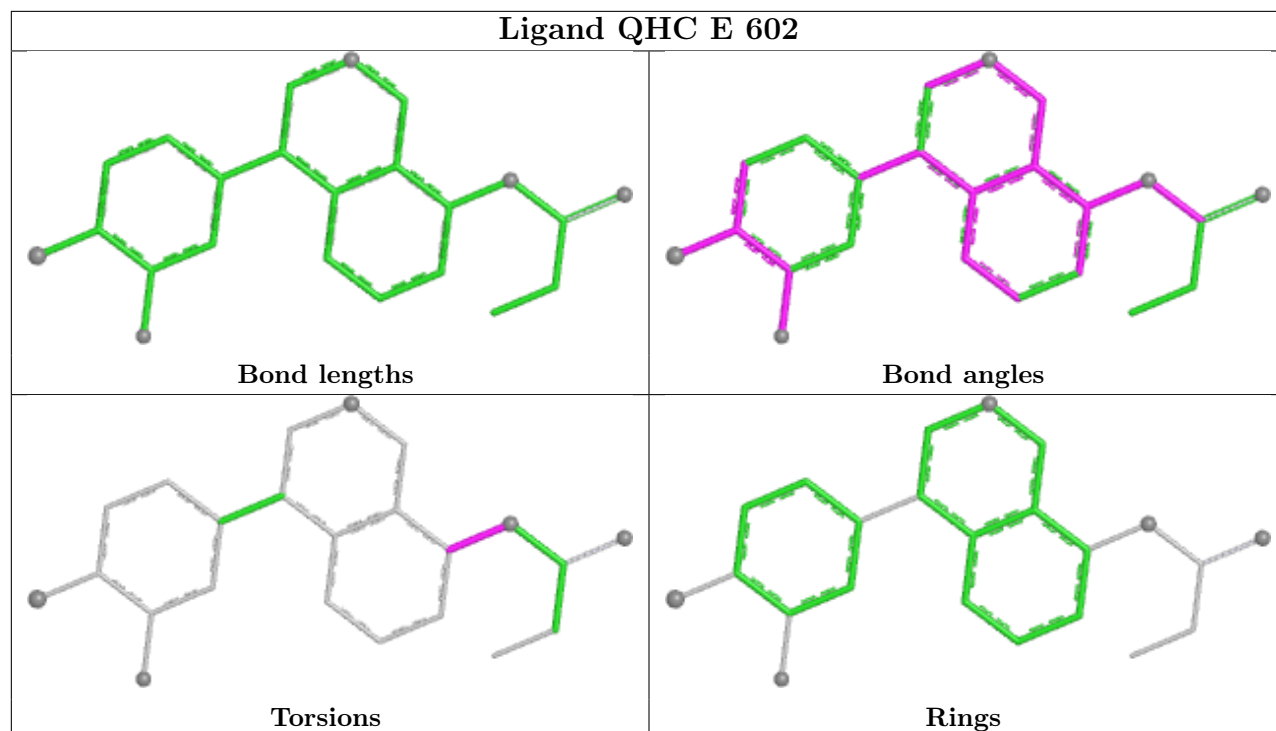
Bond angles

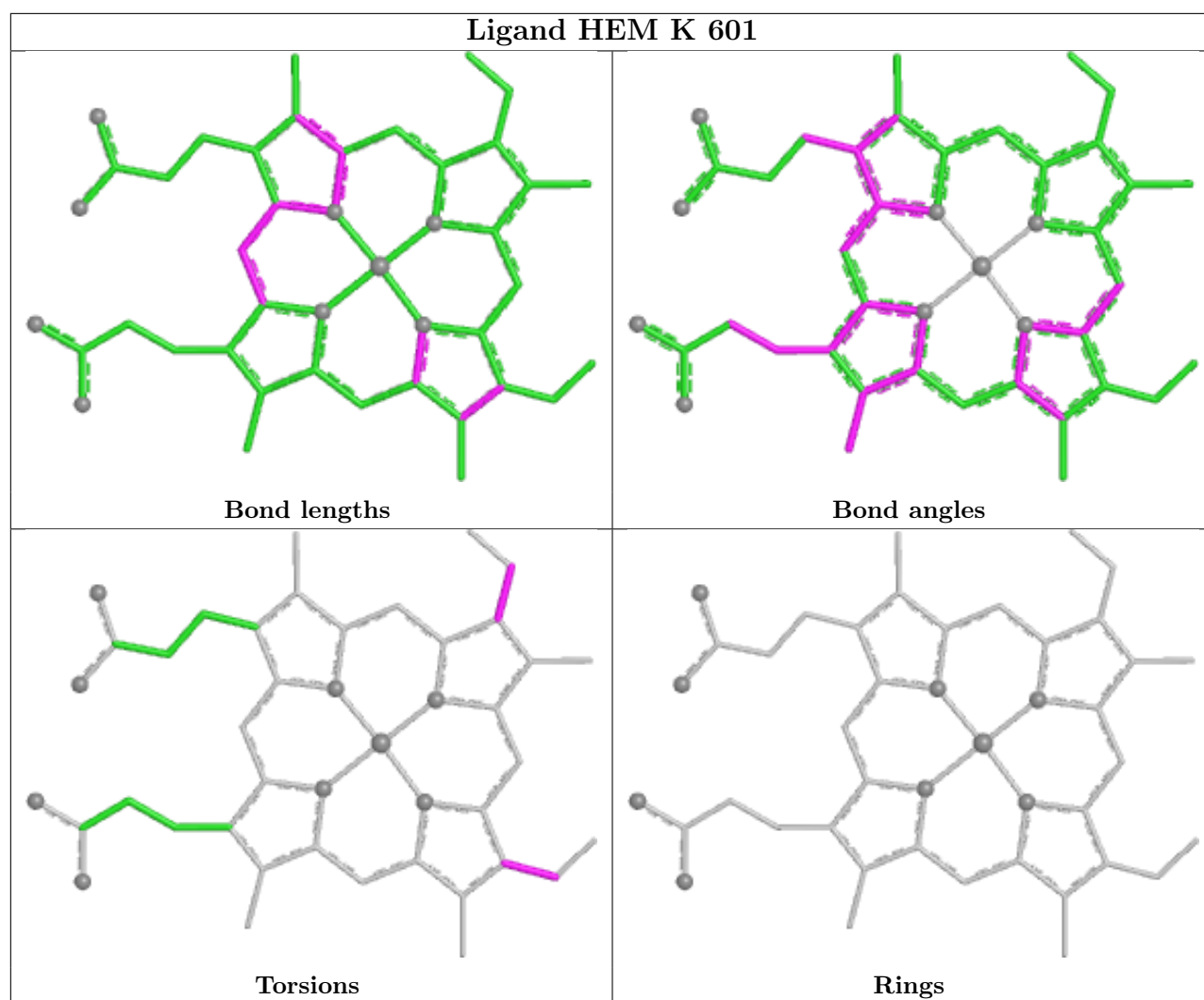


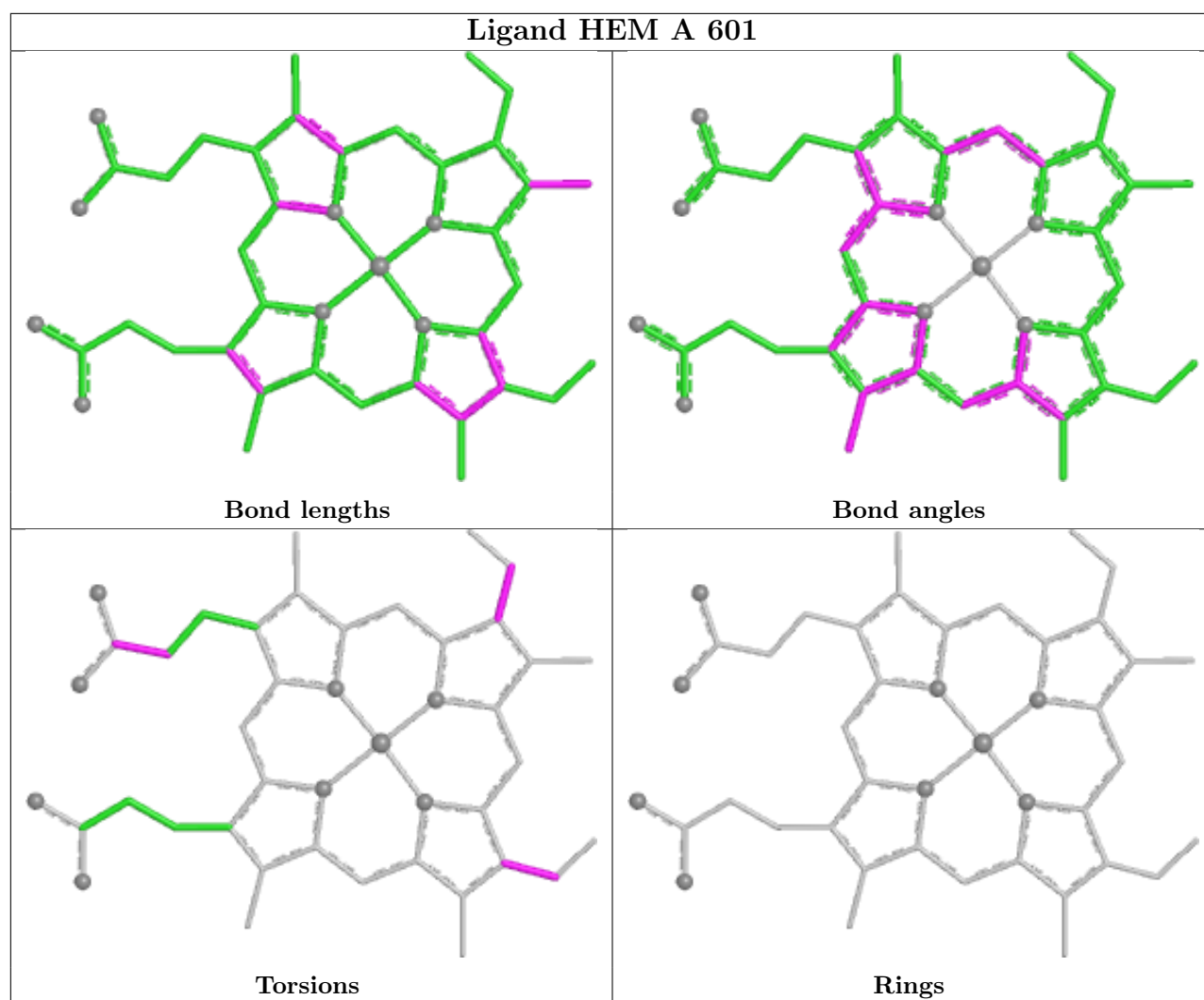
Torsions

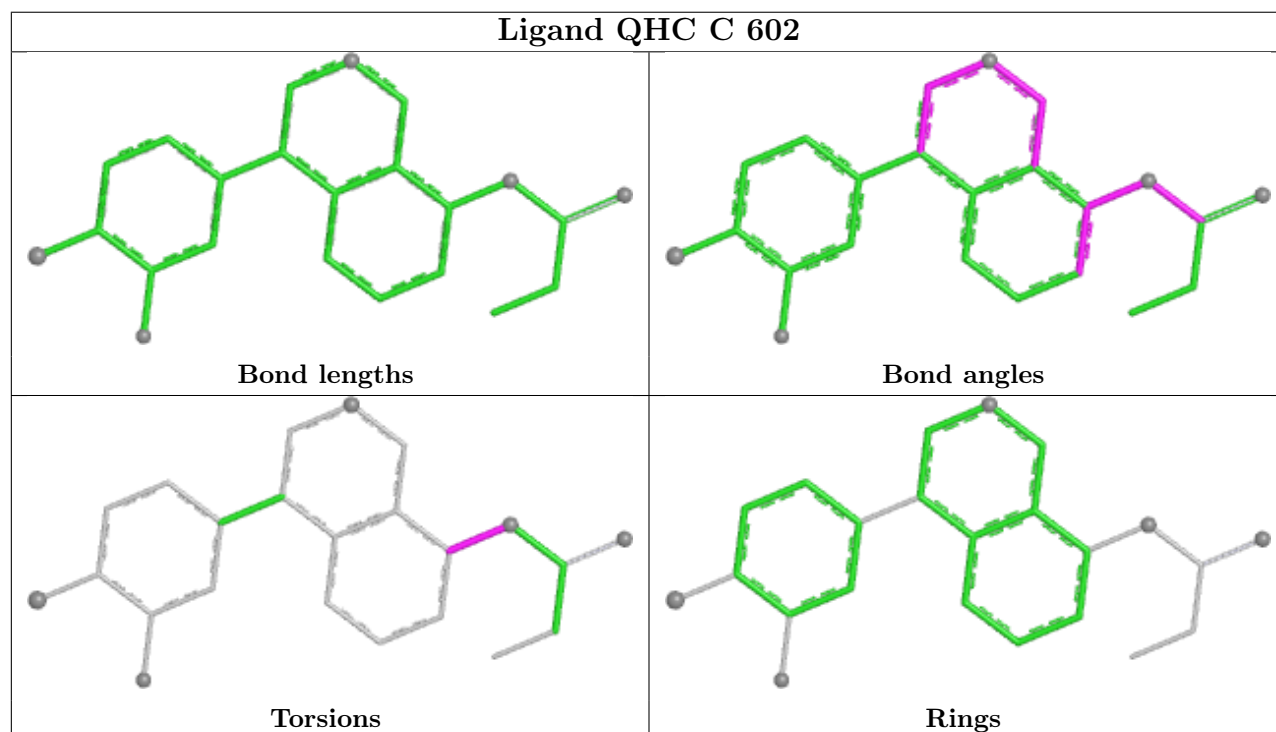
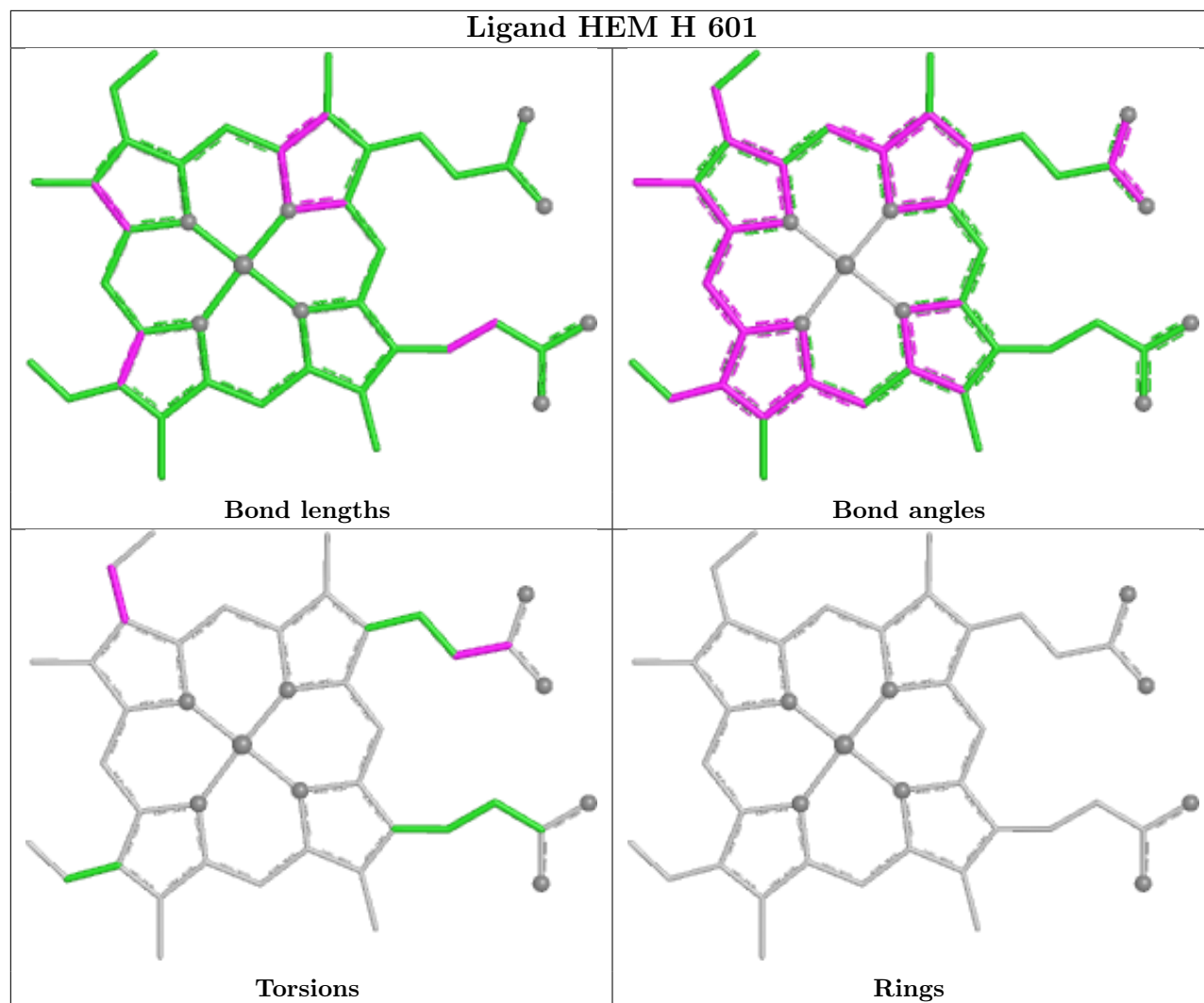


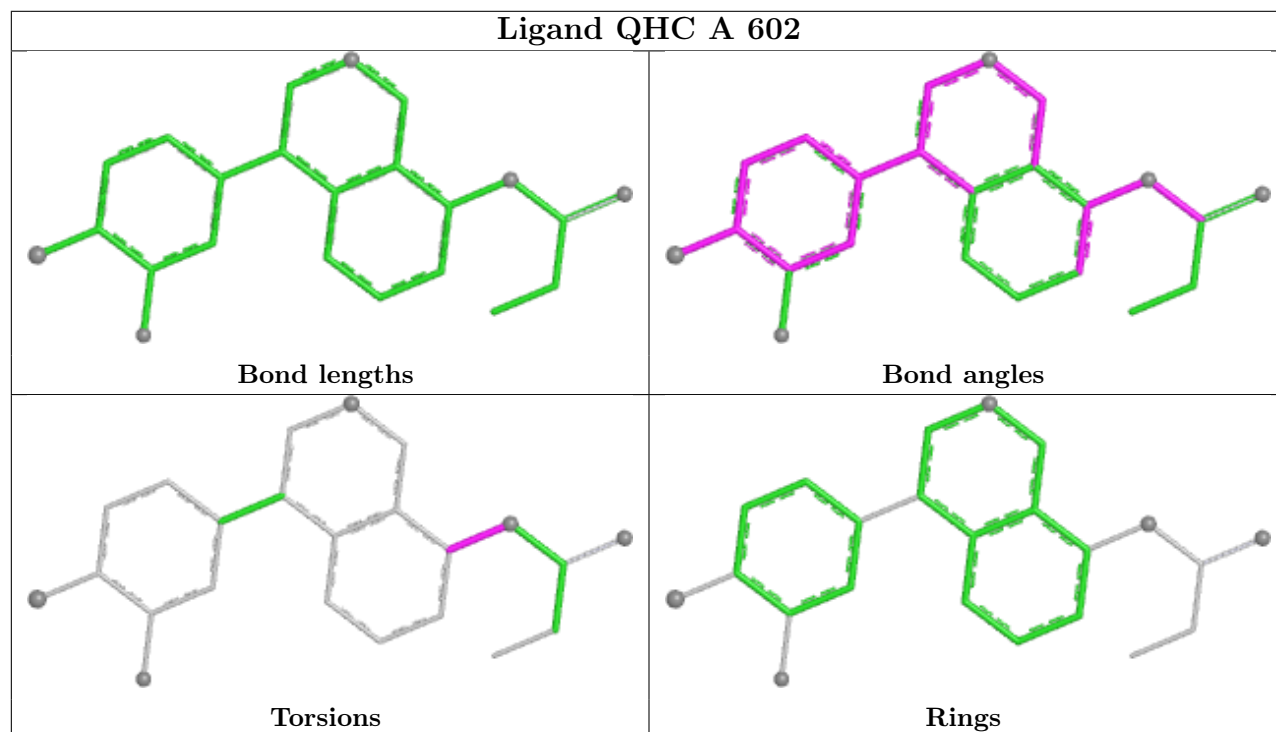
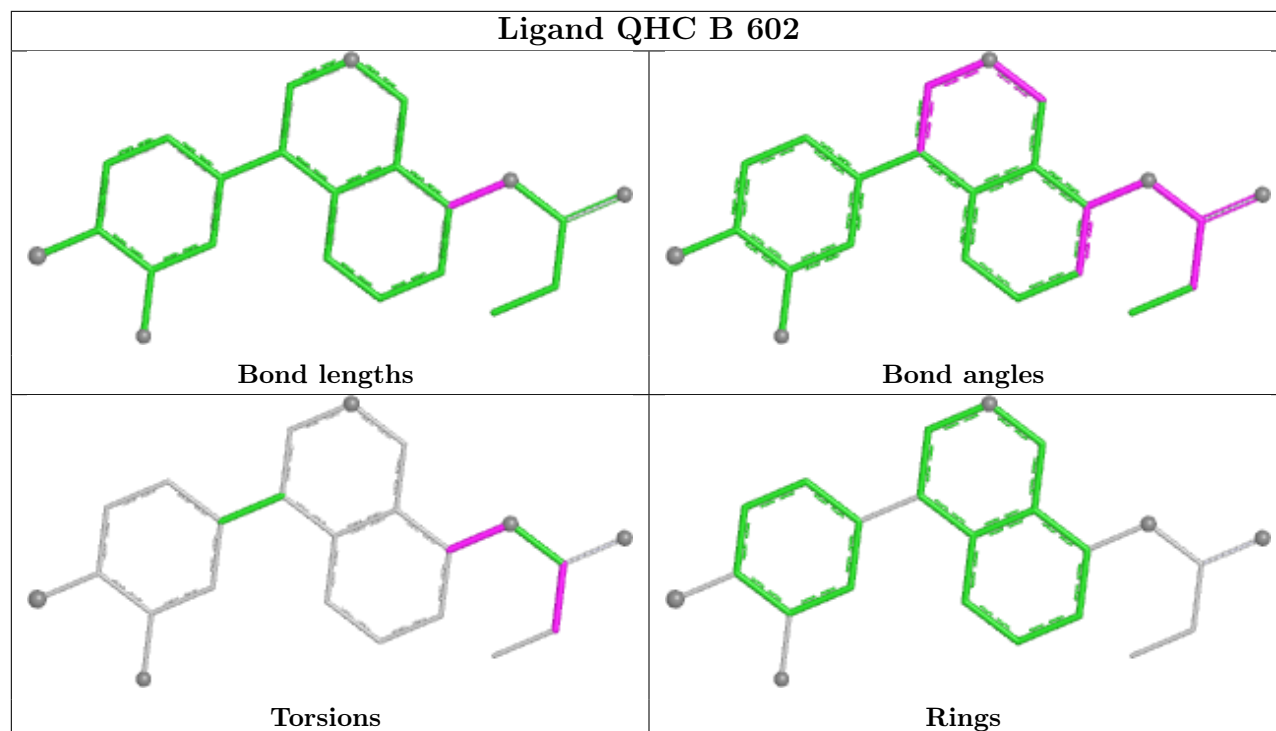
Rings



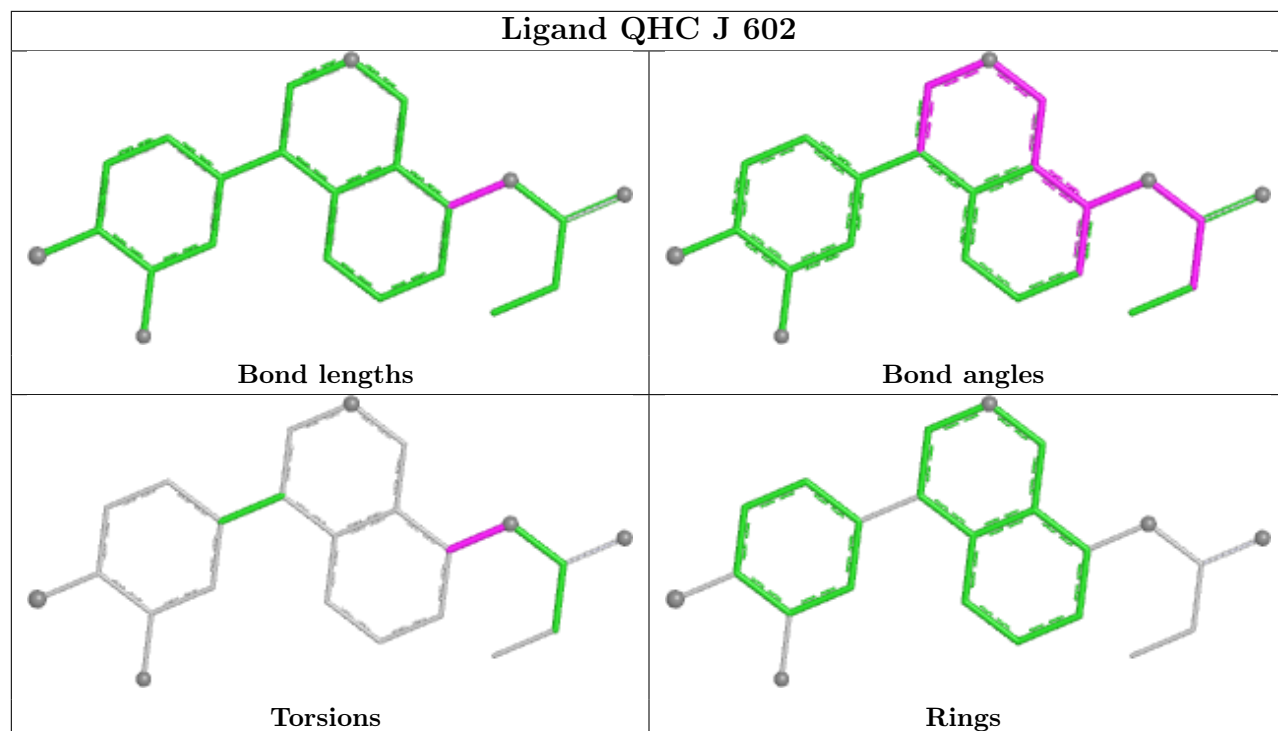




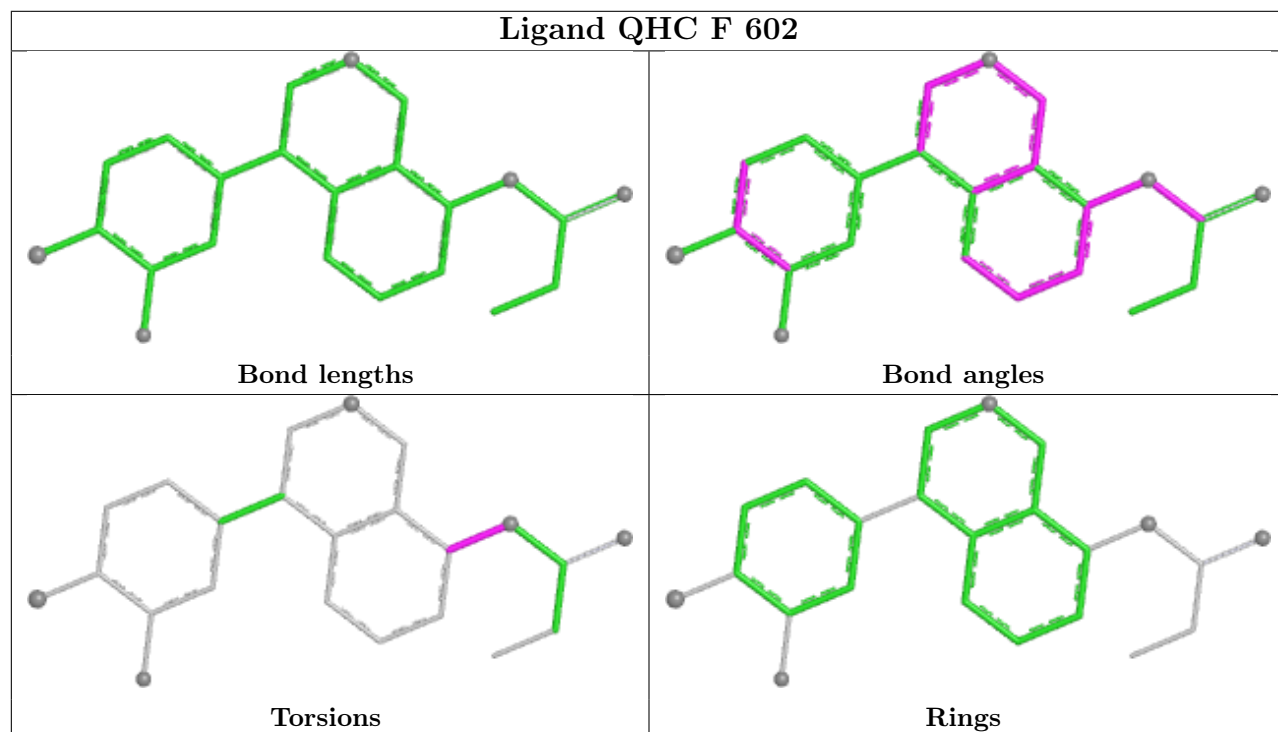


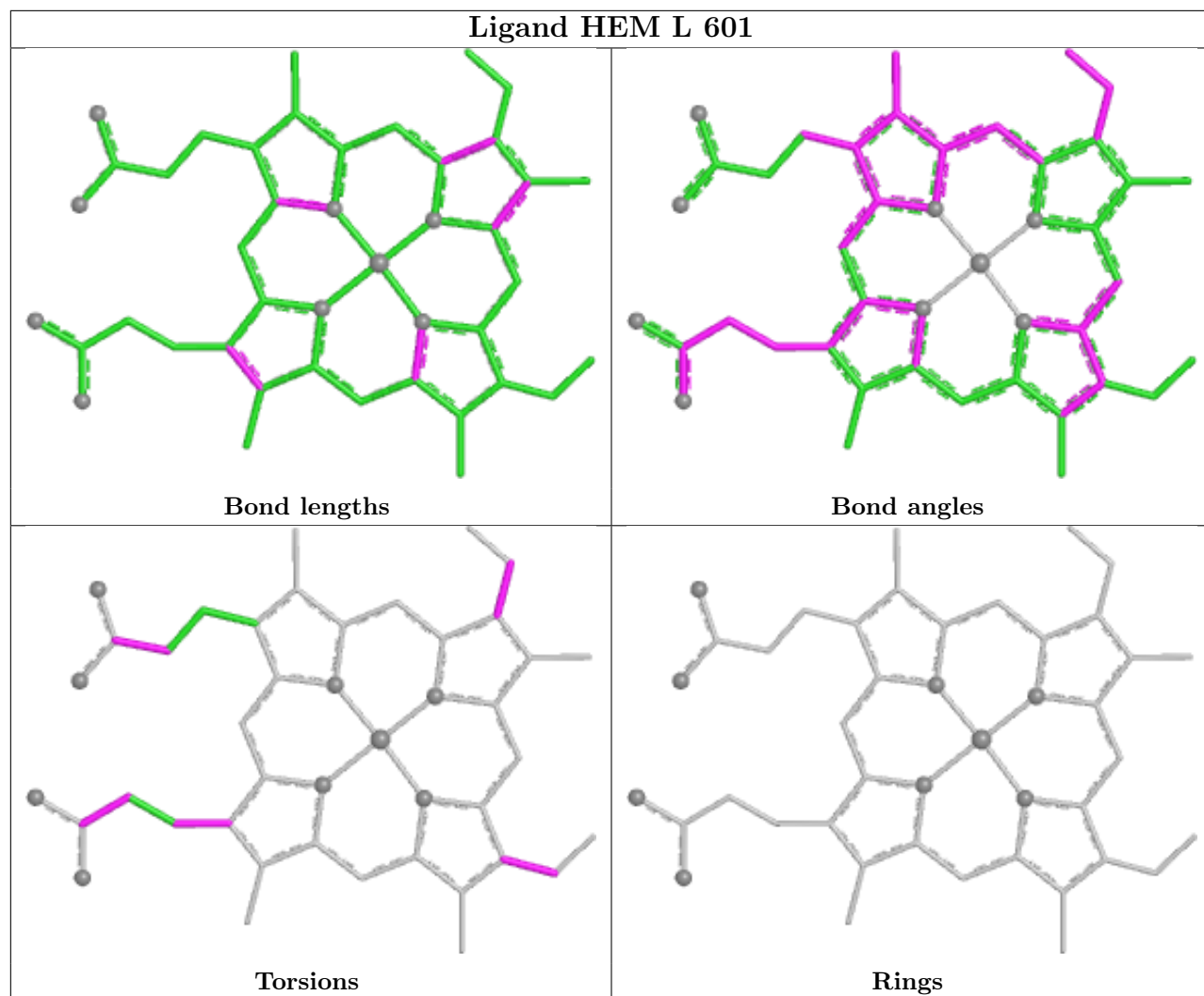


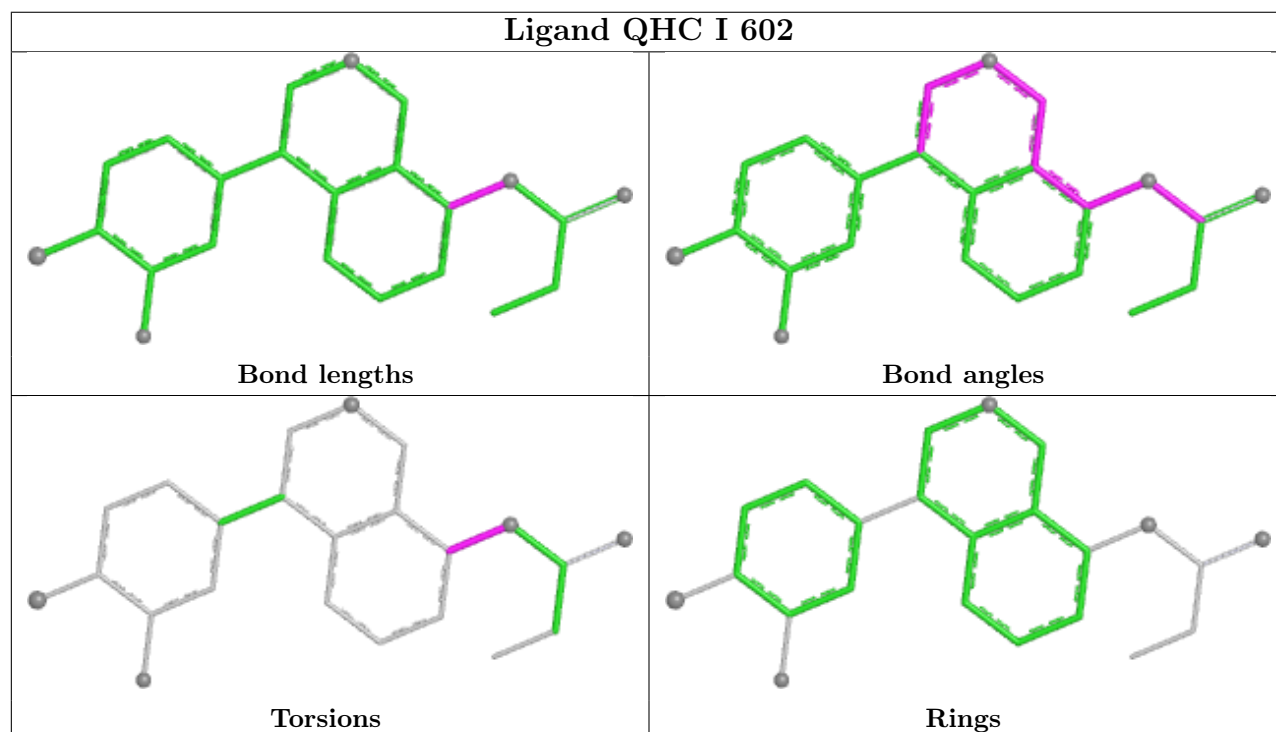
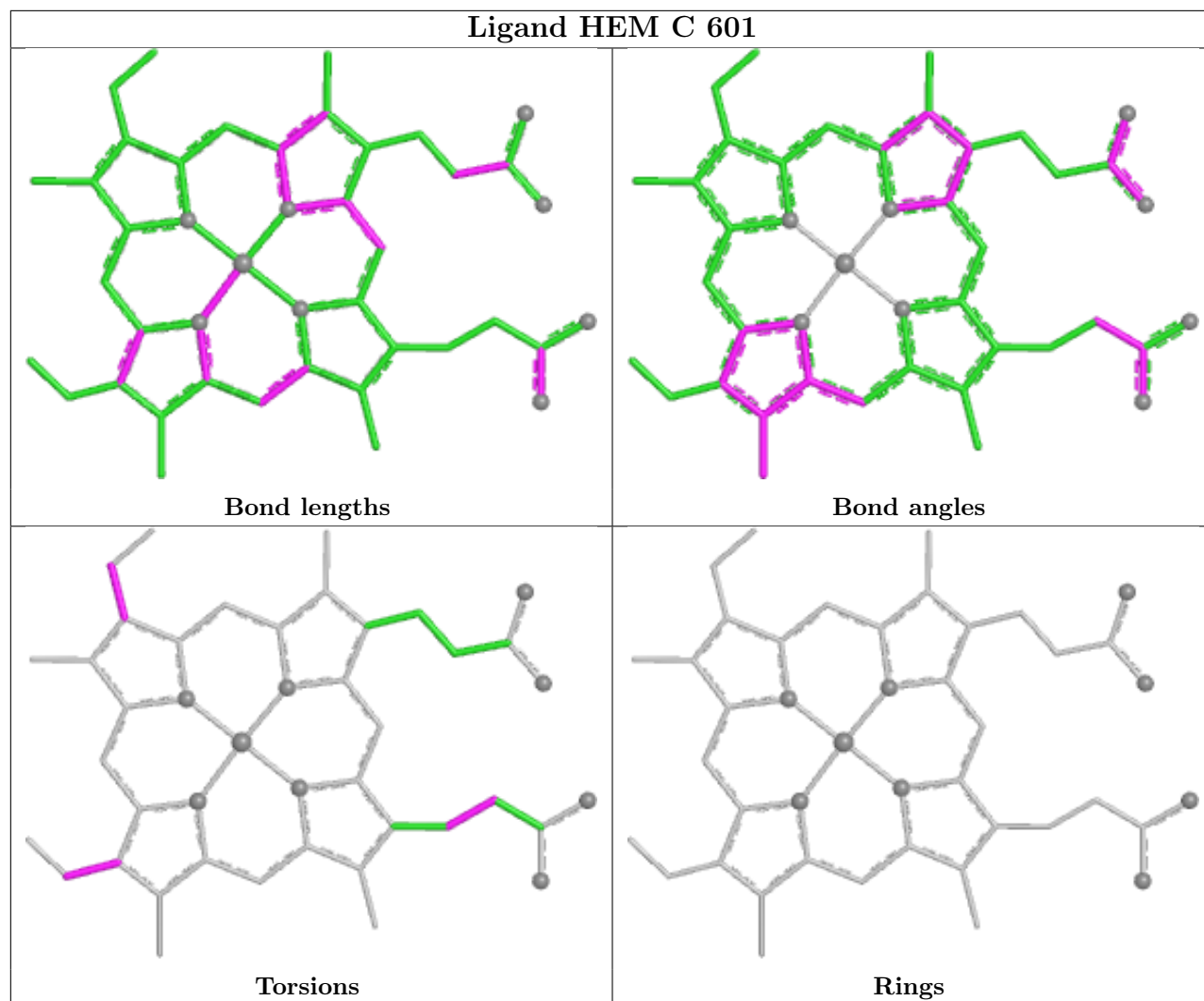
Ligand QHC J 602

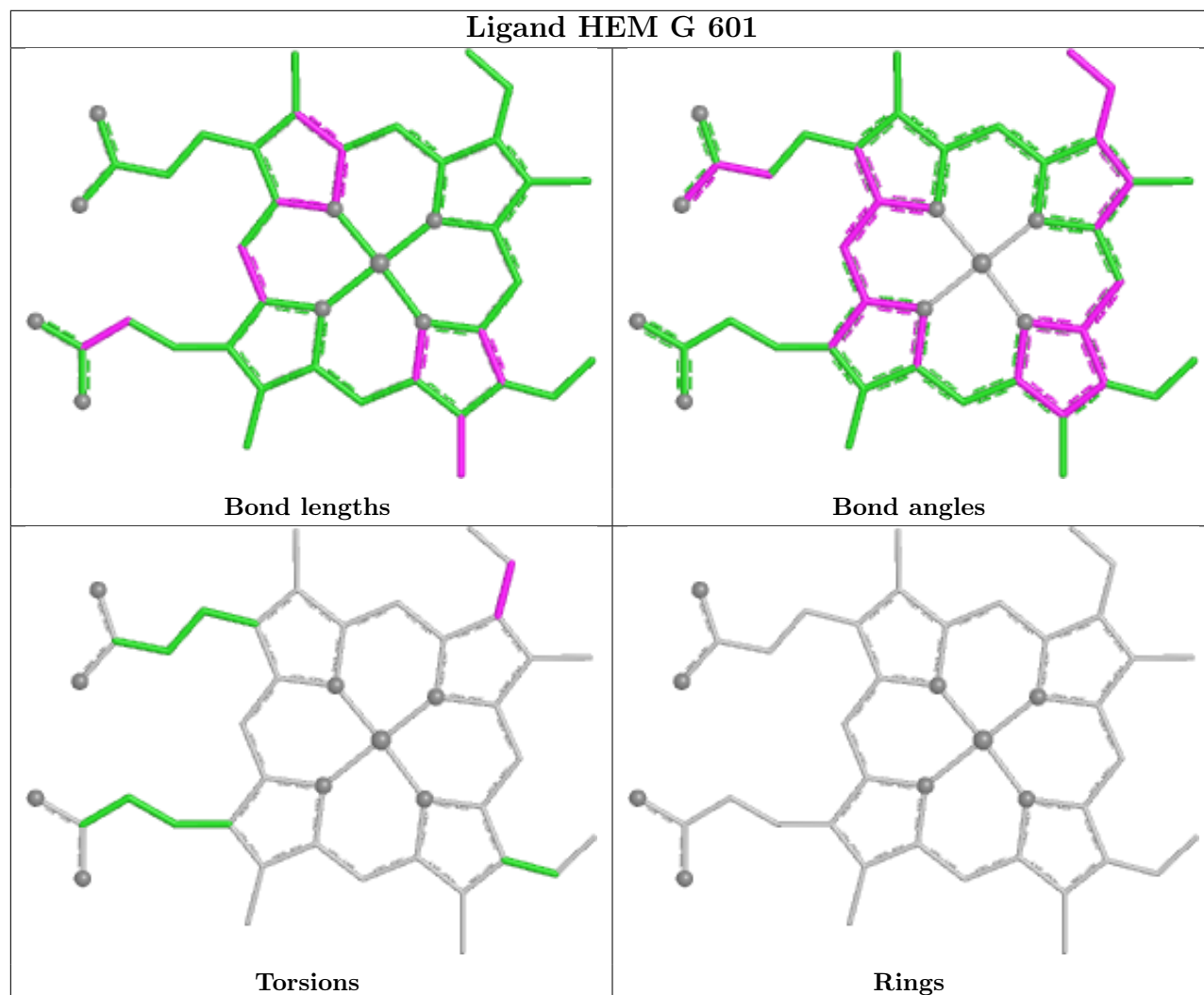
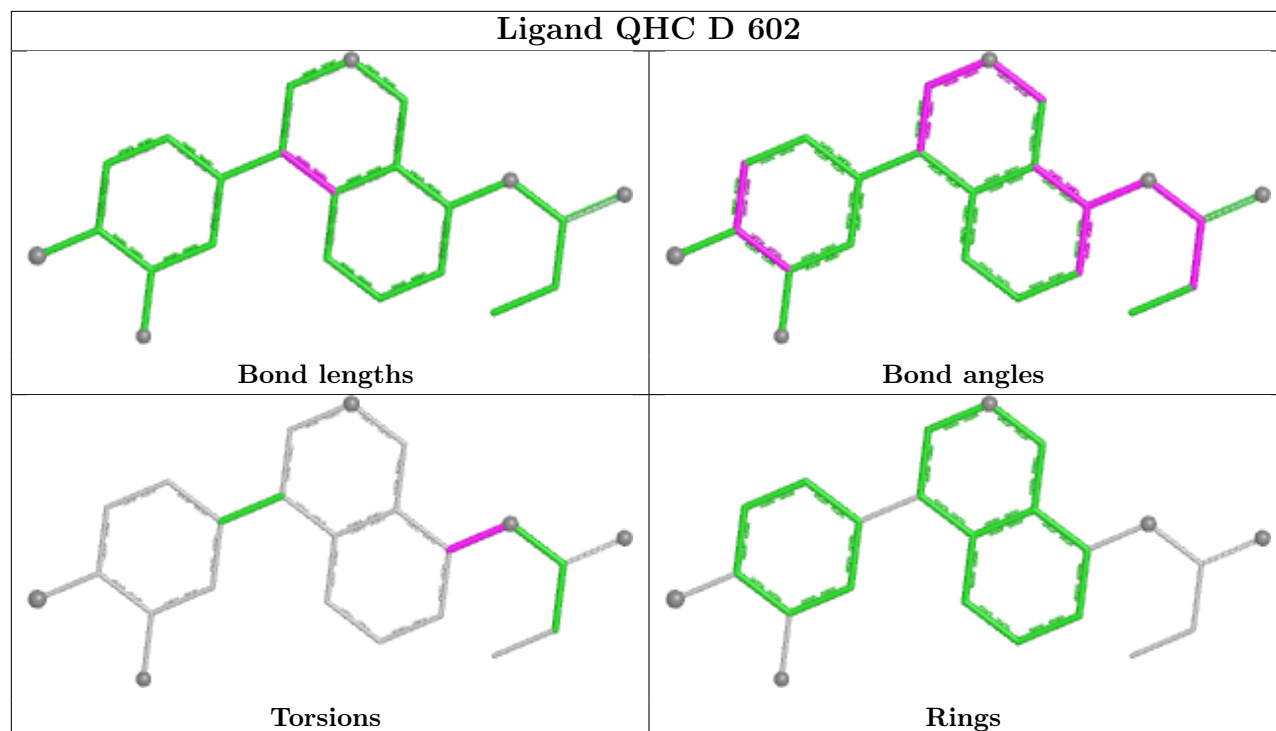


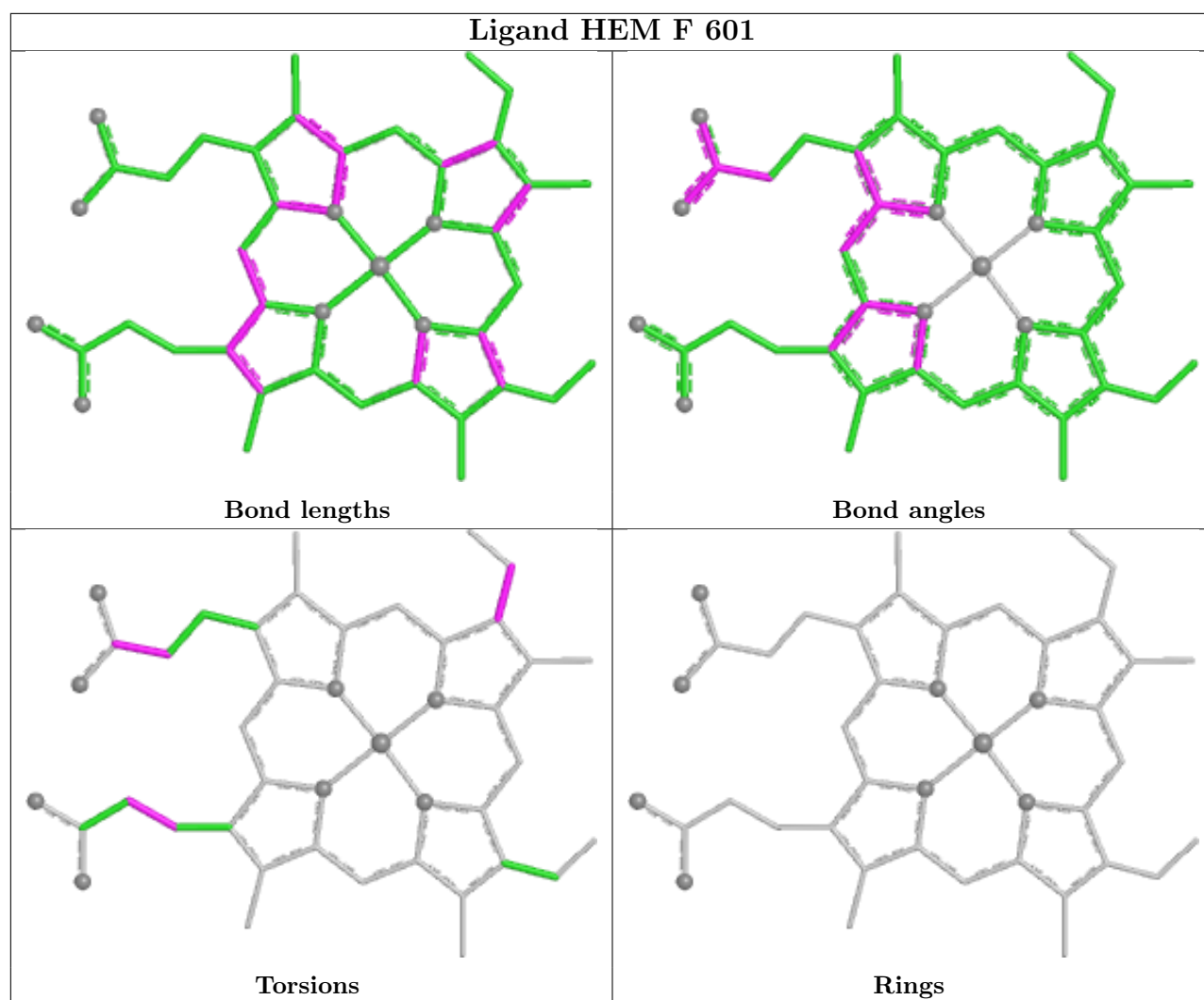
Ligand QHC F 602

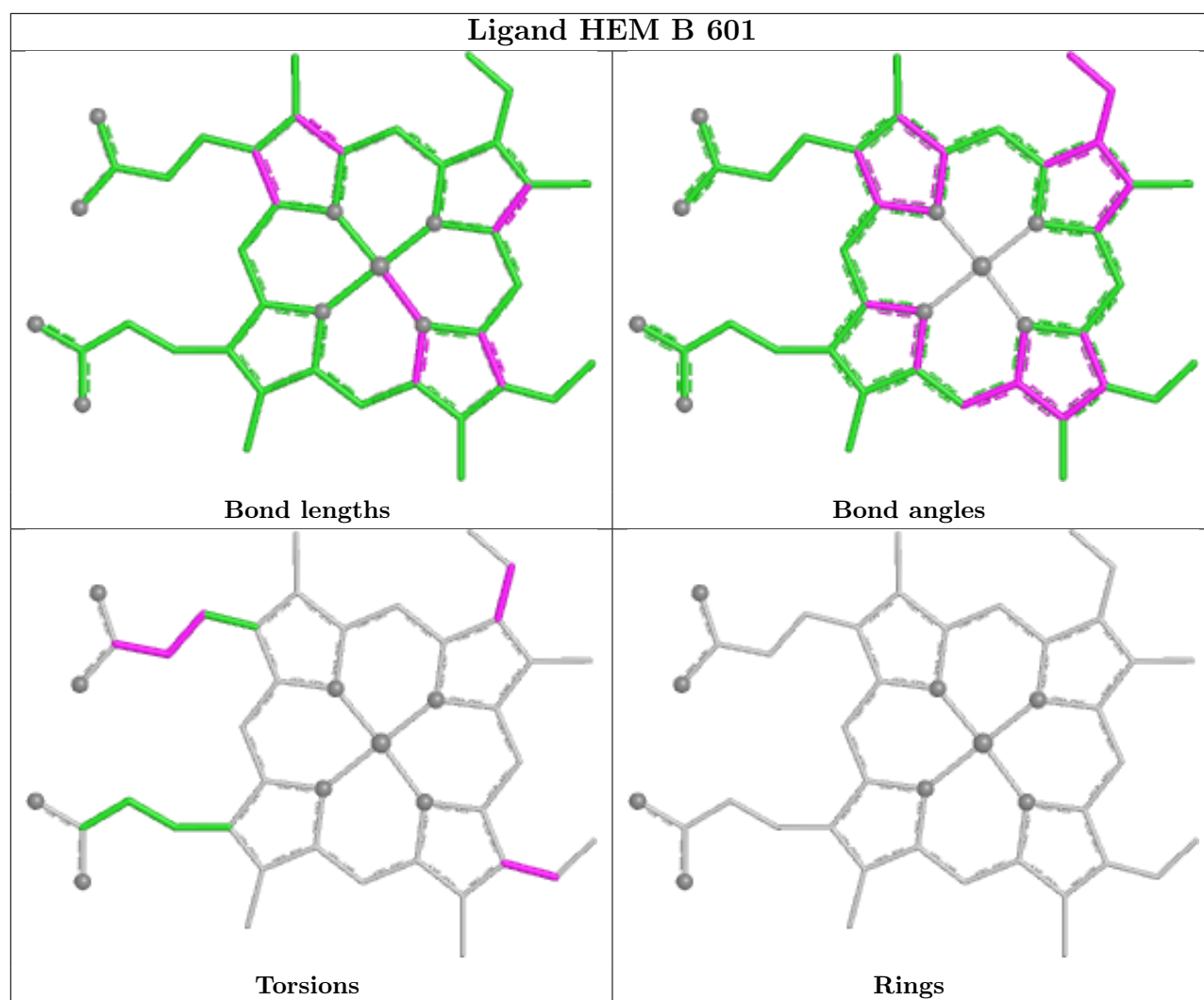












5.7 Other polymers [i](#)

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	465/489 (95%)	0.28	3 (0%) 85 73	25, 50, 78, 112	0
1	B	465/489 (95%)	0.60	12 (2%) 57 37	27, 64, 99, 126	0
1	C	465/489 (95%)	0.69	12 (2%) 57 37	34, 70, 106, 138	0
1	D	465/489 (95%)	0.49	5 (1%) 78 61	27, 58, 99, 136	0
1	E	459/489 (93%)	0.48	17 (3%) 45 28	31, 58, 105, 140	0
1	F	464/489 (94%)	0.68	18 (3%) 43 27	35, 68, 106, 125	0
1	G	465/489 (95%)	0.76	14 (3%) 52 33	44, 76, 124, 145	0
1	H	465/489 (95%)	0.92	40 (8%) 16 11	43, 78, 122, 173	0
1	I	458/489 (93%)	0.73	23 (5%) 34 22	33, 74, 143, 216	0
1	J	463/489 (94%)	0.91	34 (7%) 21 13	44, 82, 145, 230	0
1	K	461/489 (94%)	0.70	25 (5%) 31 20	35, 68, 107, 127	0
1	L	462/489 (94%)	1.03	44 (9%) 14 9	44, 82, 149, 239	0
All	All	5557/5868 (94%)	0.69	247 (4%) 39 24	25, 69, 118, 239	0

The worst 5 of 247 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	484	VAL	4.6
1	K	452	GLY	4.4
1	B	431	ILE	4.4
1	G	429	LEU	4.4
1	L	431	ILE	4.0

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 ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

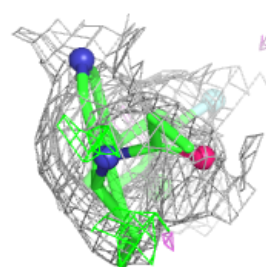
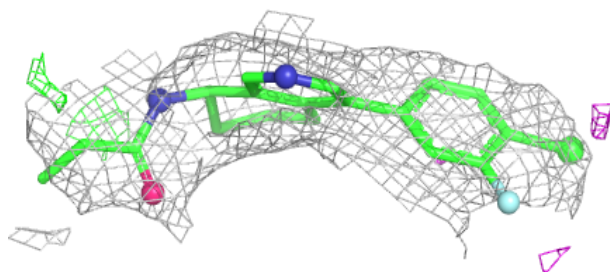
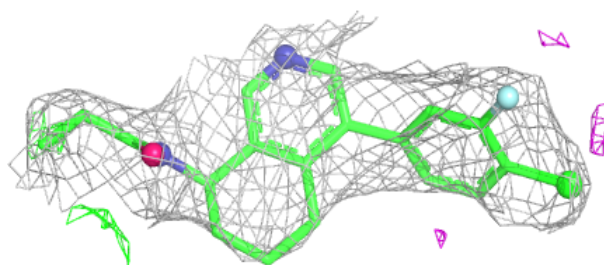
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	QHC	C	602	23/23	0.84	0.20	84,85,89,91	0
3	QHC	D	602	23/23	0.85	0.17	54,60,61,61	0
3	QHC	L	602	23/23	0.85	0.16	65,74,84,86	0
3	QHC	I	602	23/23	0.86	0.16	56,66,73,75	0
3	QHC	J	602	23/23	0.87	0.14	50,54,68,74	0
3	QHC	B	602	23/23	0.87	0.16	56,62,69,72	0
3	QHC	F	602	23/23	0.90	0.13	51,61,68,68	0
3	QHC	E	602	23/23	0.90	0.12	44,49,53,54	0
3	QHC	K	602	23/23	0.91	0.13	51,60,72,72	0
2	HEM	H	601	43/43	0.91	0.15	62,64,67,68	0
2	HEM	L	601	43/43	0.92	0.13	59,60,64,64	0
2	HEM	B	601	43/43	0.92	0.14	54,56,60,61	0
2	HEM	J	601	43/43	0.94	0.12	51,52,56,58	0
2	HEM	K	601	43/43	0.94	0.12	55,56,60,64	0
2	HEM	G	601	43/43	0.94	0.11	51,52,56,58	0
2	HEM	I	601	43/43	0.94	0.12	53,54,58,65	0
2	HEM	F	601	43/43	0.95	0.11	46,48,52,52	0
2	HEM	A	601	43/43	0.95	0.11	39,39,44,51	0
2	HEM	E	601	43/43	0.96	0.10	43,44,48,49	0
2	HEM	C	601	43/43	0.96	0.10	39,41,46,48	0
3	QHC	A	602	23/23	0.96	0.07	3,13,32,37	0
2	HEM	D	601	43/43	0.96	0.10	32,33,38,43	0

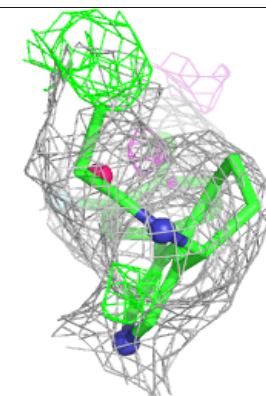
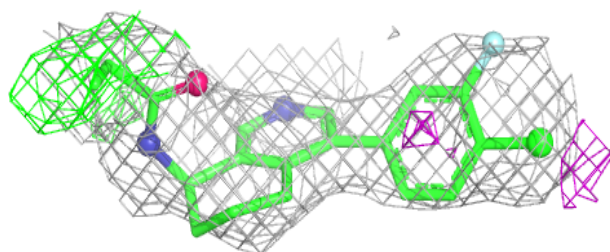
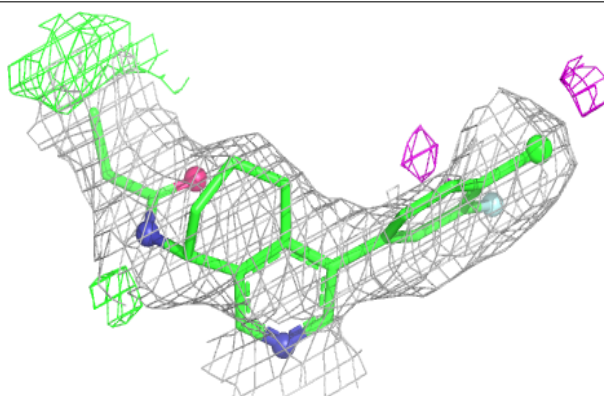
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around QHC 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)

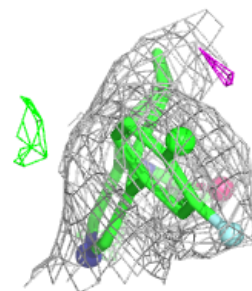
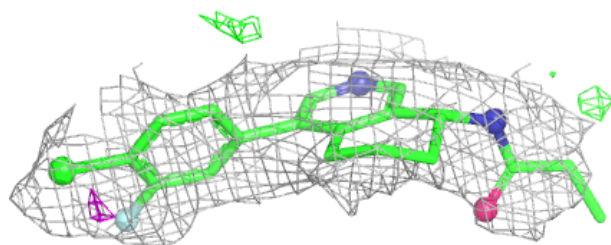
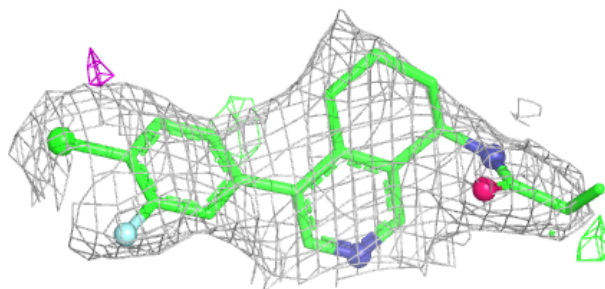
**Electron density around QHC D 602:**

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

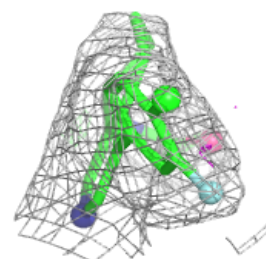
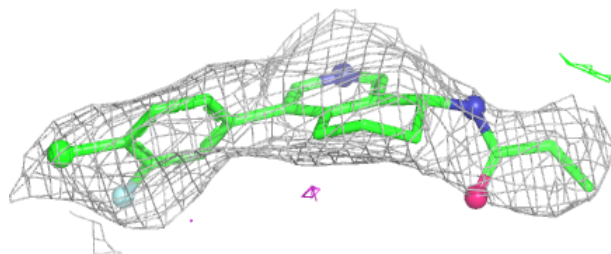
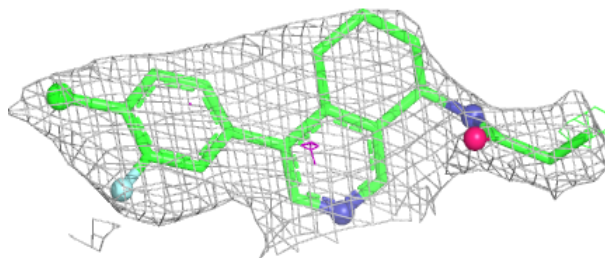


Electron density around QHC L 602:

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

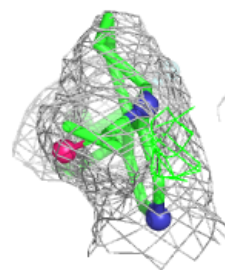
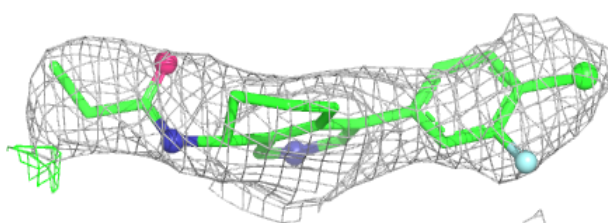
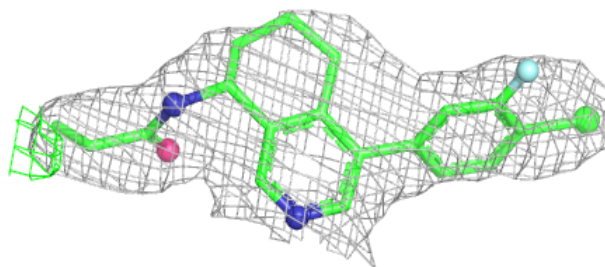
**Electron density around QHC I 602:**

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

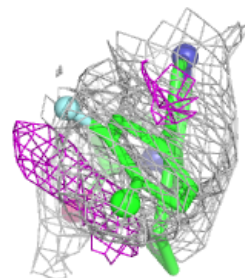
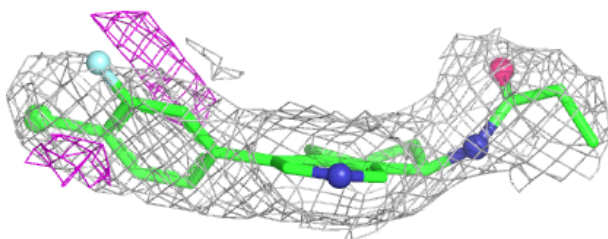
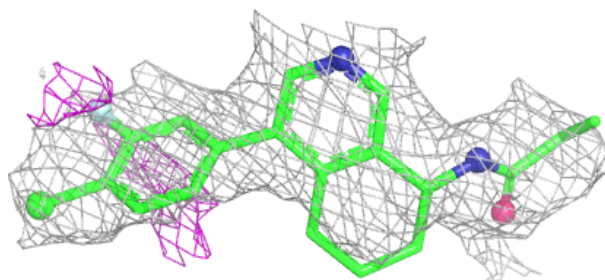


Electron density around QHC J 602:

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

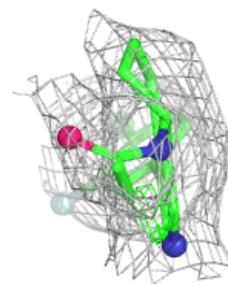
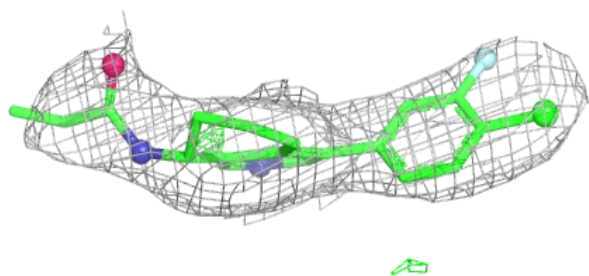
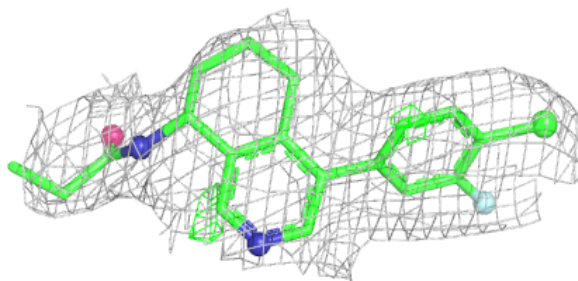
**Electron density around QHC B 602:**

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

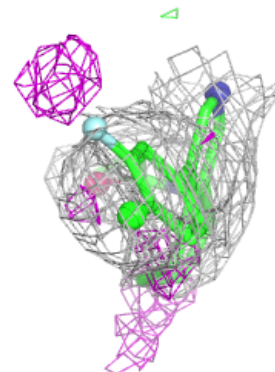
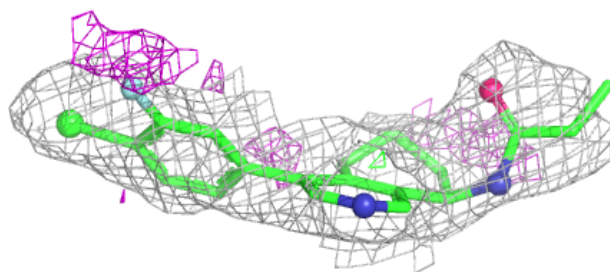
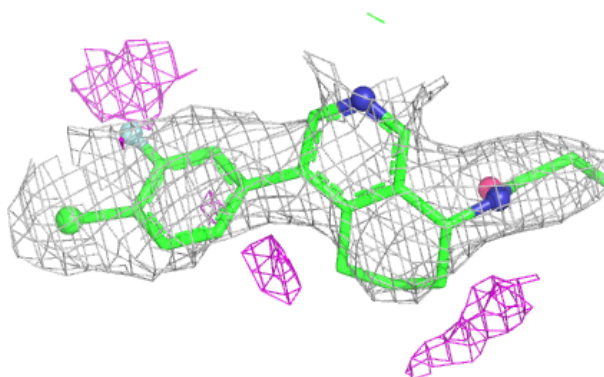


Electron density around QHC F 602:

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

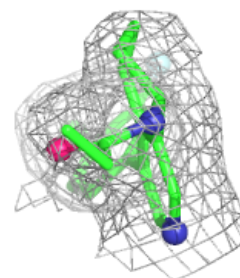
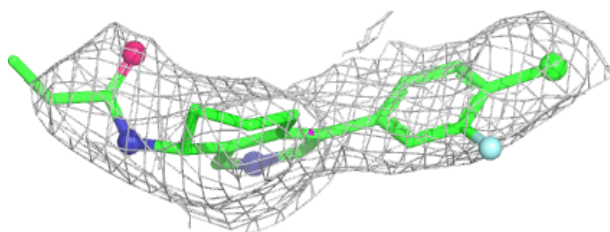
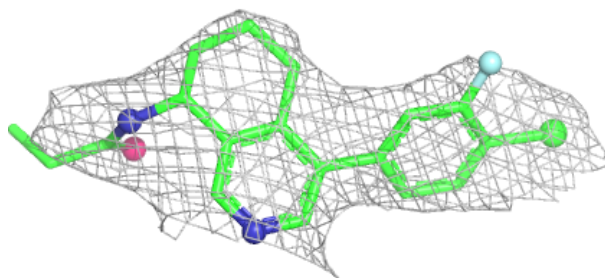
**Electron density around QHC E 602:**

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



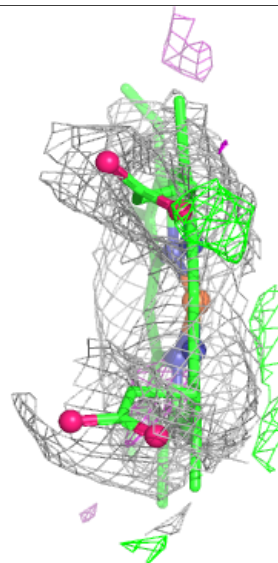
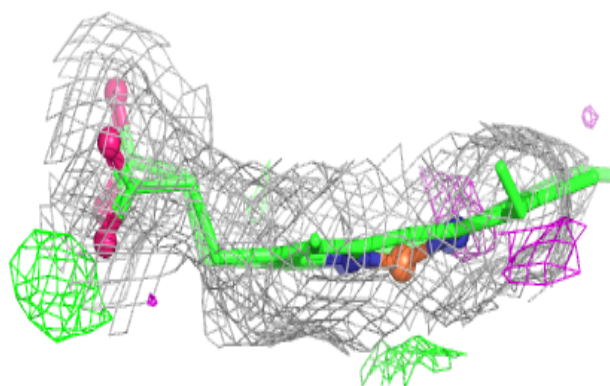
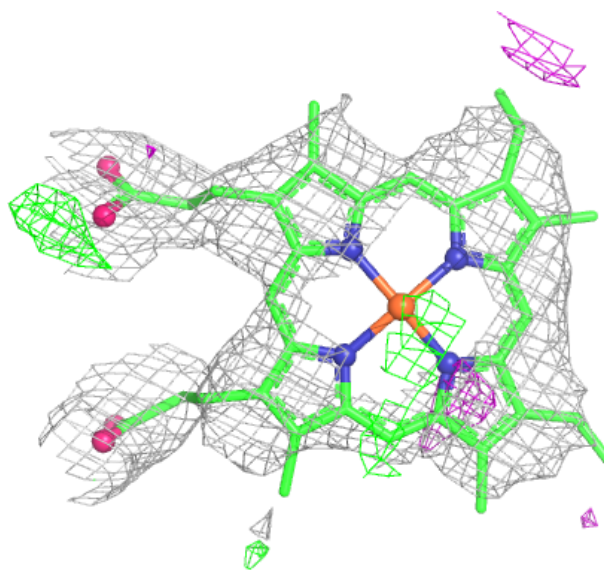
Electron density around QHC K 602:

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



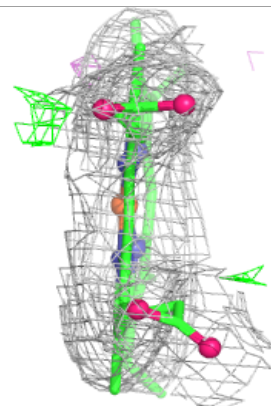
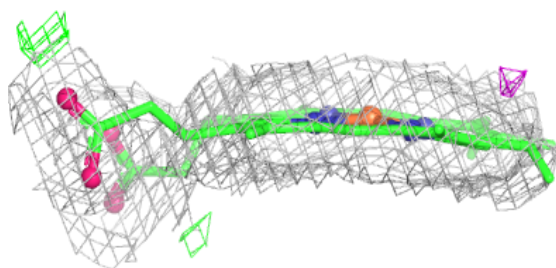
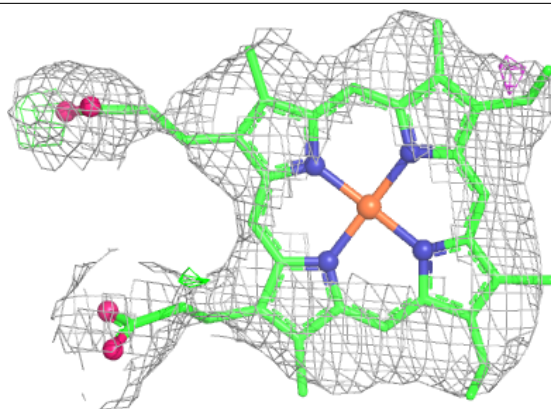
Electron density around HEM H 601:

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



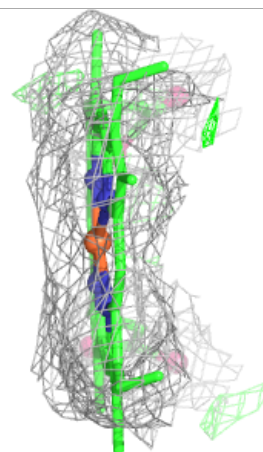
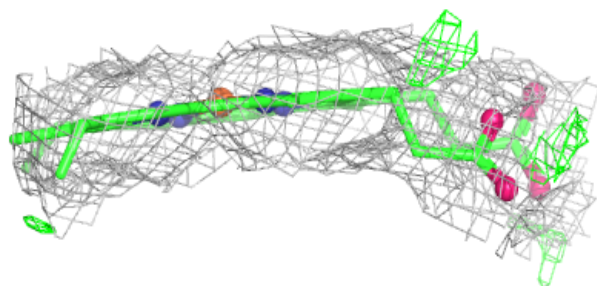
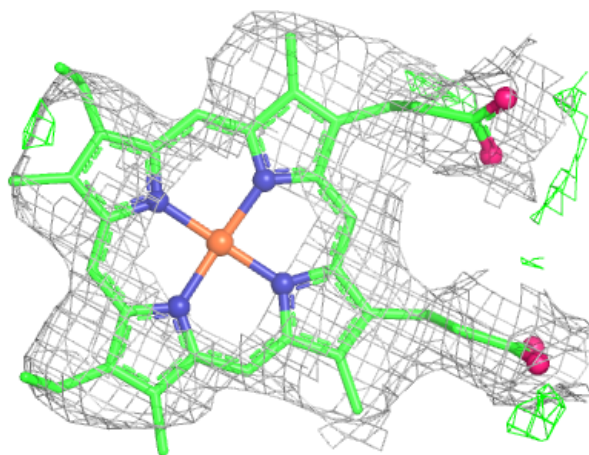
Electron density around HEM L 601:

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



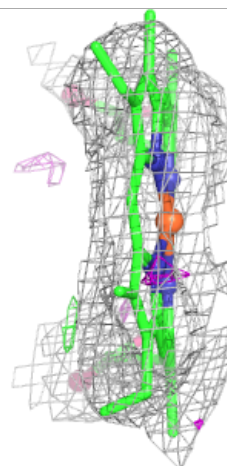
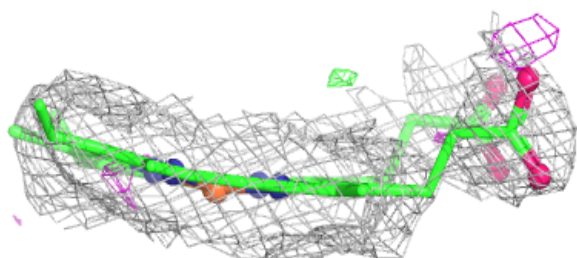
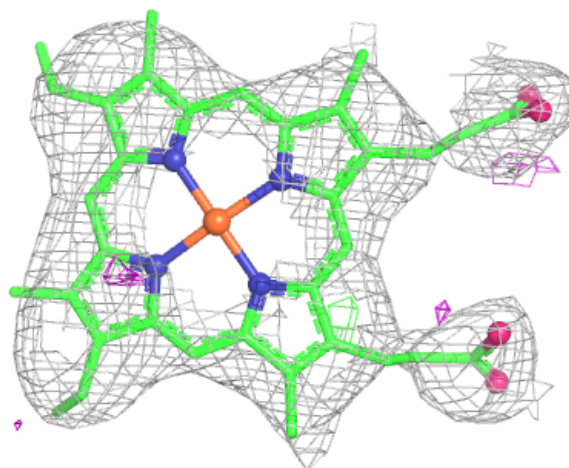
Electron density around HEM B 601:

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



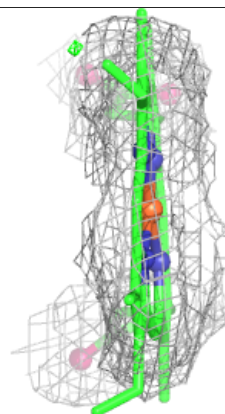
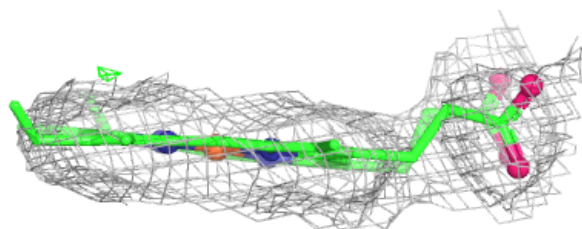
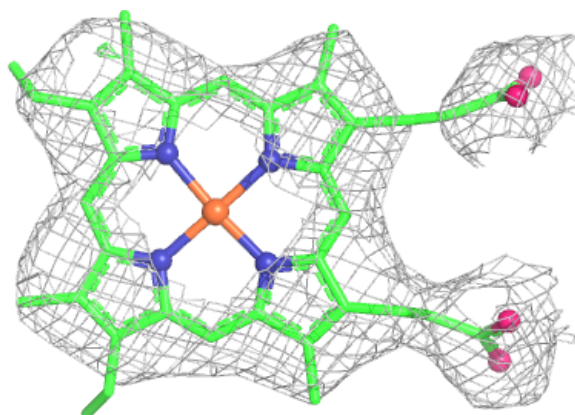
Electron density around HEM J 601:

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



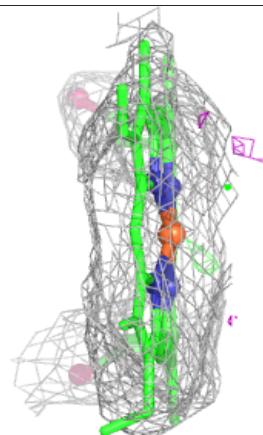
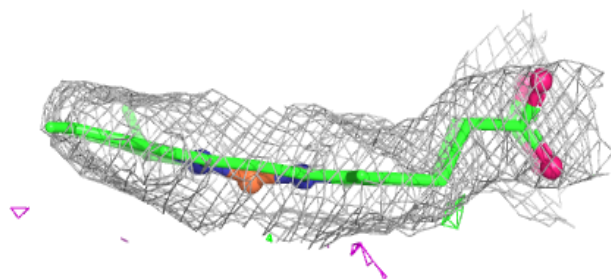
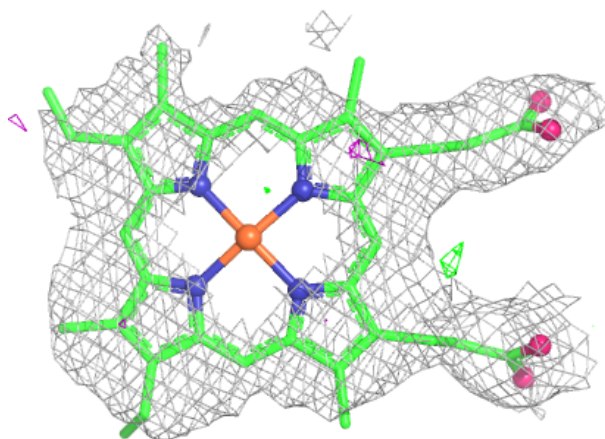
Electron density around HEM K 601:

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



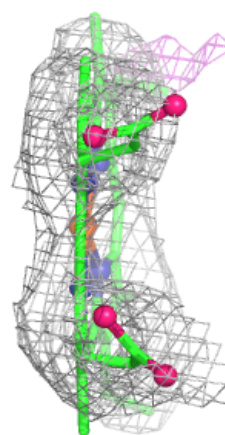
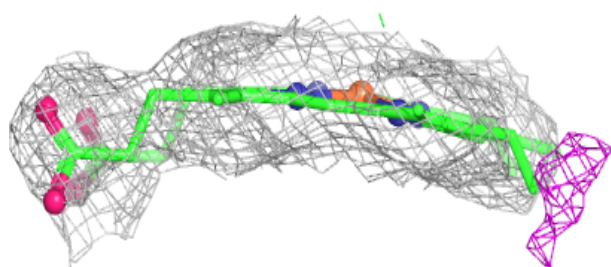
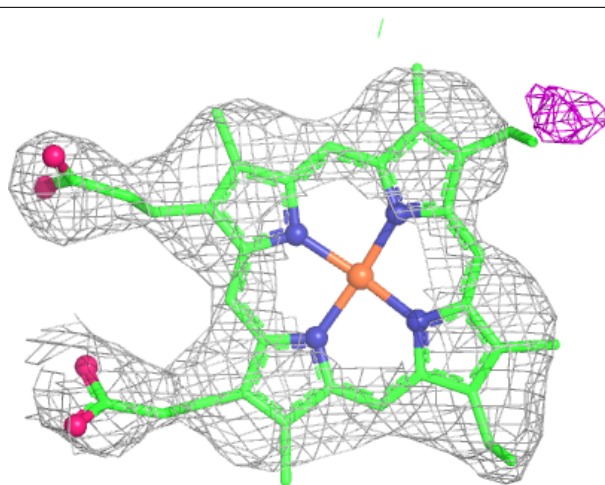
Electron density around HEM G 601:

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



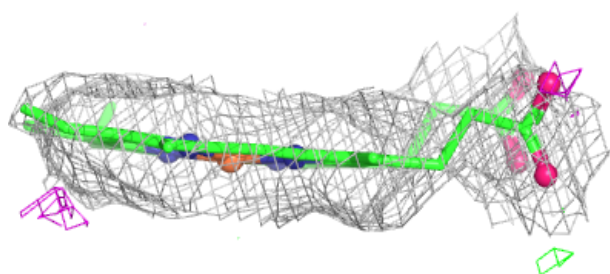
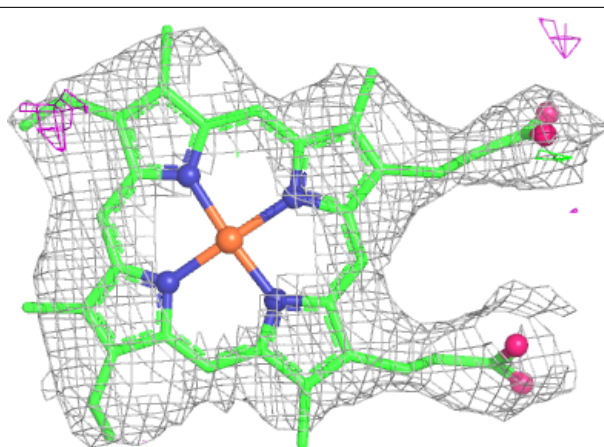
Electron density around HEM I 601:

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



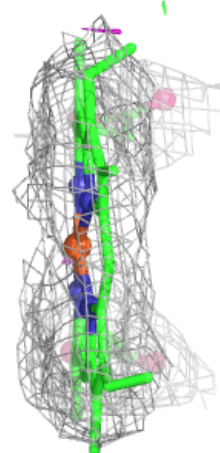
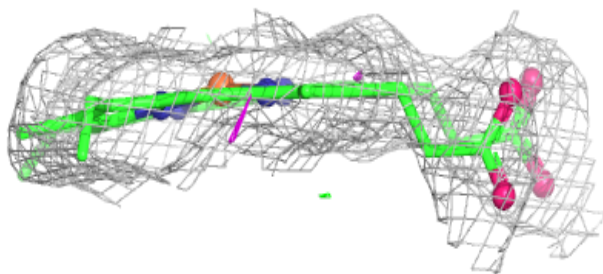
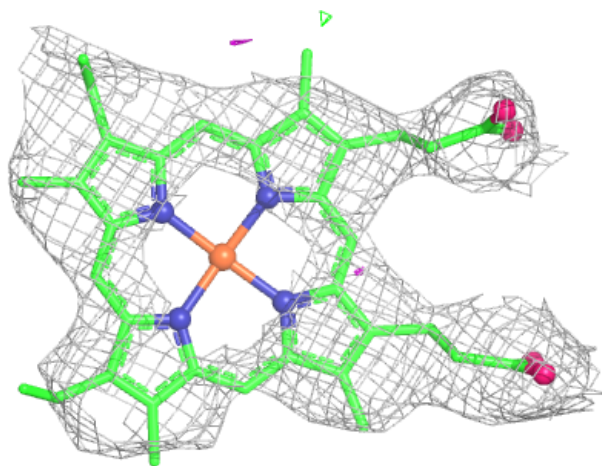
Electron density around HEM F 601:

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



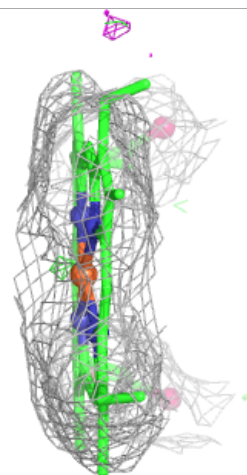
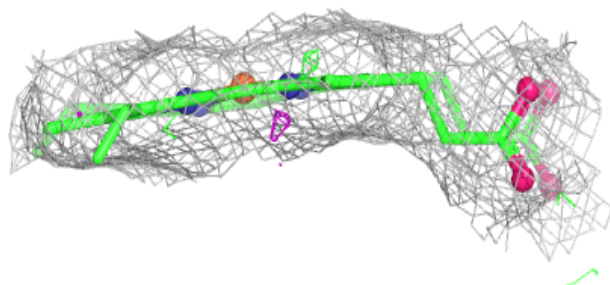
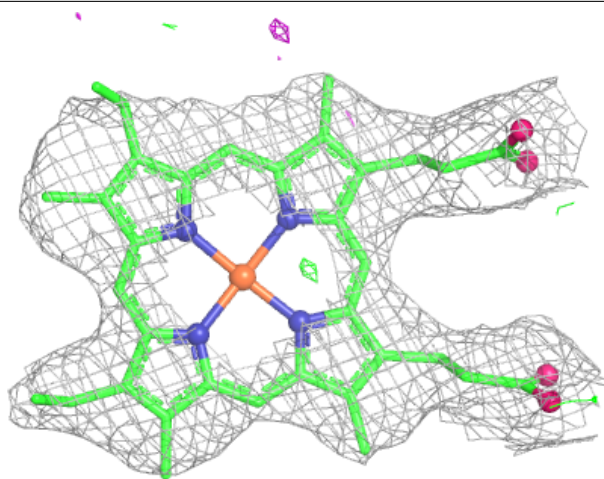
Electron density around HEM A 601:

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



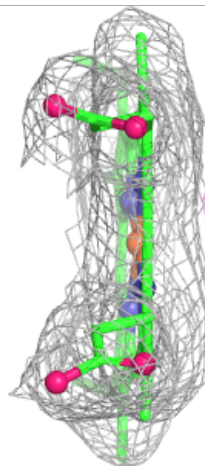
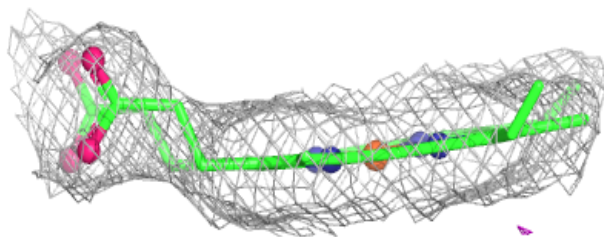
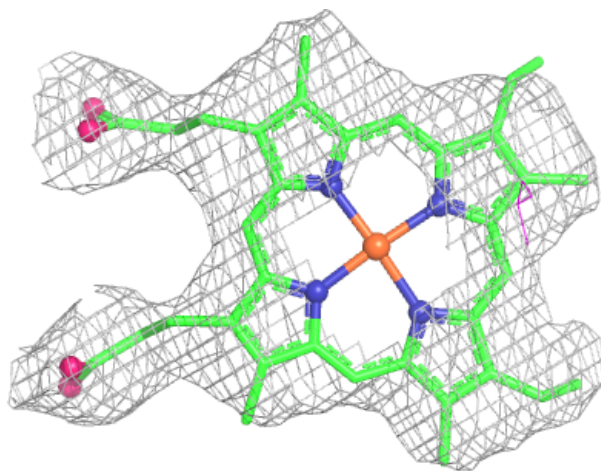
Electron density around HEM E 601:

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



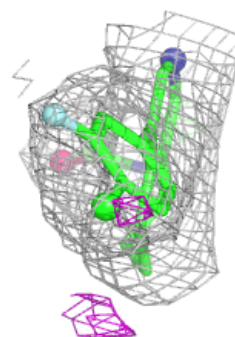
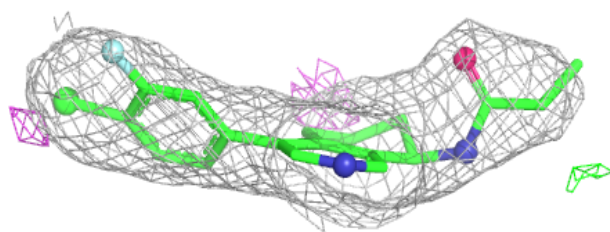
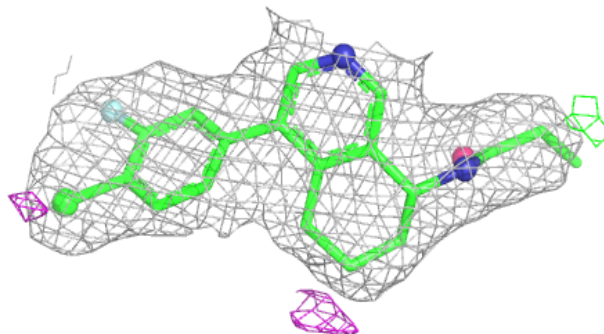
Electron density around HEM C 601:

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

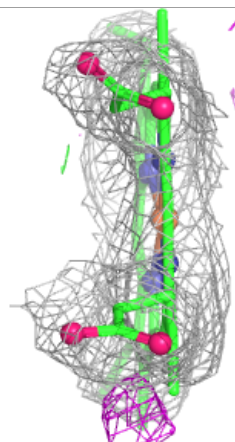
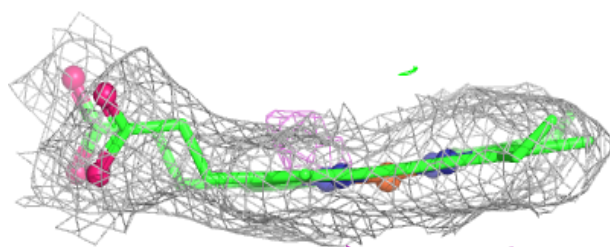
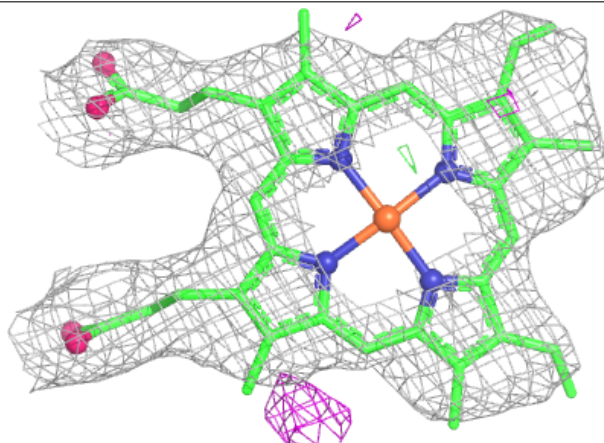


Electron density around QHC A 602:

$2mF_o-DF_c$ (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)



6.5 Other polymers [i](#)

There are no such residues in this entry.