



wwPDB X-ray Structure Validation Summary Report ⓘ

May 15, 2020 – 10:16 am BST

PDB ID : 2ZZS
Title : Crystal structure of cytochrome c554 from *Vibrio parahaemolyticus* strain RIMD2210633
Authors : Akazaki, H.; Kawai, F.; Kumaki, Y.; Sekine, K.; Hakamata, W.; Nishio, T.; Park, S.-Y.; Oku, T.
Deposited on : 2009-02-24
Resolution : 1.80 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 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)
Xtriage (Phenix) : 1.13
EDS : 2.11
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) : 2.11

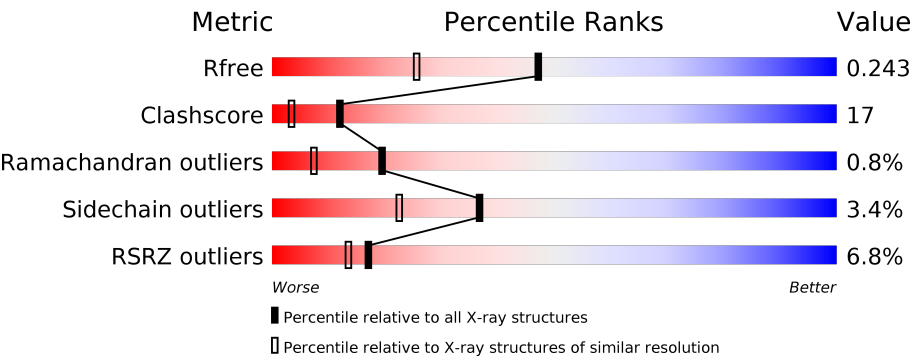
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 on 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	1	103	<div><div>7%</div><div><div></div><div></div><div></div><div></div></div><div>63%15%•21%</div></div>
1	2	103	<div><div>16%</div><div><div></div><div></div><div></div><div></div></div><div>58%16%••21%</div></div>
1	3	103	<div><div>11%</div><div><div></div><div></div><div></div><div></div></div><div>59%16%••21%</div></div>
1	4	103	<div><div>8%</div><div><div></div><div></div><div></div><div></div></div><div>54%23%•21%</div></div>
1	5	103	<div><div>30%</div><div><div></div><div></div><div></div><div></div></div><div>41%29%7%•21%</div></div>
1	6	103	<div><div>55%</div><div><div></div><div></div><div></div><div></div></div><div>29%32%11%5%23%</div></div>

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Mol	Chain	Length	Quality of chain
1	A	103	
1	B	103	
1	C	103	
1	D	103	
1	E	103	
1	F	103	
1	G	103	
1	H	103	
1	I	103	
1	J	103	
1	K	103	
1	L	103	
1	M	103	
1	N	103	
1	O	103	
1	P	103	
1	Q	103	
1	R	103	
1	S	103	
1	T	103	
1	U	103	
1	V	103	
1	W	103	
1	X	103	
1	Y	103	

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Mol	Chain	Length	Quality of chain
1	Z	103	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HEC	6	220	-	-	X	-
3	GOL	R	104	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 21402 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 c554.

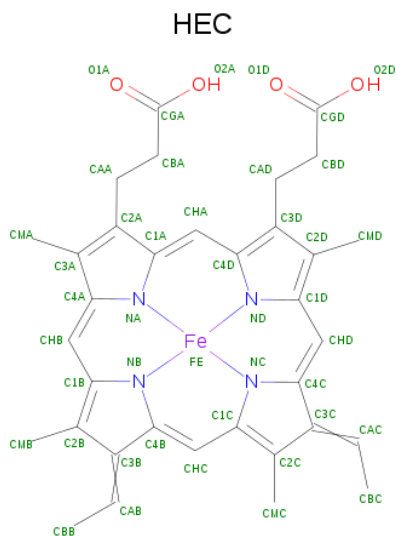
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	B	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	C	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	D	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	E	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	F	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	G	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	H	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	I	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	J	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	K	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	L	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	M	80	Total	C	N	O	S	0	0	0
			567	347	100	117	3			
1	N	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	O	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	P	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	Q	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	R	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	S	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	T	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	U	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	V	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	W	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	X	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	Y	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	Z	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	1	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	2	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	3	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	4	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	5	81	Total	C	N	O	S	0	0	0
			575	353	101	118	3			
1	6	79	Total	C	N	O	S	0	0	0
			561	344	99	115	3			

- Molecule 2 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).



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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	O	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	P	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	Q	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	R	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	S	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	T	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	U	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	V	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	W	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	X	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	Y	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	Z	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	1	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	2	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	3	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	4	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	5	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	6	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	C	1	Total	C	O	0	0
			6	3	3		
3	E	1	Total	C	O	0	0
			6	3	3		
3	P	1	Total	C	O	0	0
			6	3	3		
3	R	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	91	Total	O	0	0
			91	91		
4	B	101	Total	O	0	0
			101	101		
4	C	73	Total	O	0	0
			73	73		
4	D	62	Total	O	0	0
			62	62		
4	E	76	Total	O	0	0
			76	76		
4	F	64	Total	O	0	0
			64	64		
4	G	62	Total	O	0	0
			62	62		
4	H	70	Total	O	0	0
			70	70		

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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	I	62	Total 62	O 62	0	0
4	J	24	Total 24	O 24	0	0
4	K	35	Total 35	O 35	0	0
4	L	53	Total 53	O 53	0	0
4	M	51	Total 51	O 51	0	0
4	N	37	Total 37	O 37	0	0
4	O	71	Total 71	O 71	0	0
4	P	35	Total 35	O 35	0	0
4	Q	63	Total 63	O 63	0	0
4	R	46	Total 46	O 46	0	0
4	S	42	Total 42	O 42	0	0
4	T	69	Total 69	O 69	0	0
4	U	67	Total 67	O 67	0	0
4	V	59	Total 59	O 59	0	0
4	W	53	Total 53	O 53	0	0
4	X	21	Total 21	O 21	0	0
4	Y	55	Total 55	O 55	0	0
4	Z	48	Total 48	O 48	0	0
4	1	34	Total 34	O 34	0	0
4	2	15	Total 15	O 15	0	0
4	3	21	Total 21	O 21	0	0

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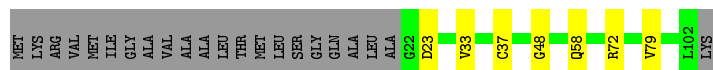
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	4	28	Total 28	O 28	0	0
4	5	27	Total 27	O 27	0	0
4	6	9	Total 9	O 9	0	0

3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

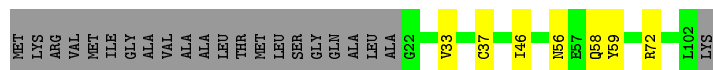
- Molecule 1: Cytochrome c554

Chain A: 



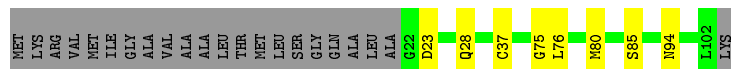
- Molecule 1: Cytochrome c554

Chain B: 



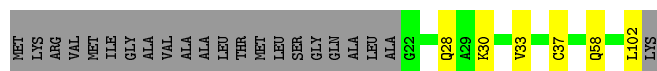
- Molecule 1: Cytochrome c554

Chain C: 



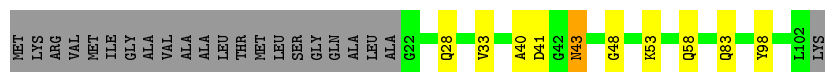
- Molecule 1: Cytochrome c554

Chain D: 



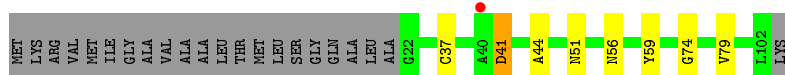
- Molecule 1: Cytochrome c554

Chain E: 

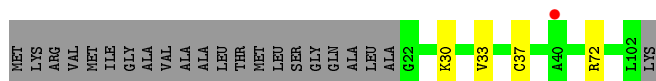
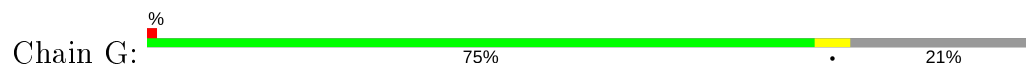


- Molecule 1: Cytochrome c554

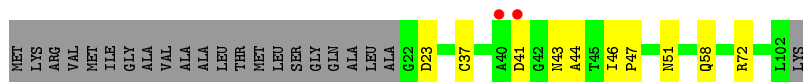
Chain F: 



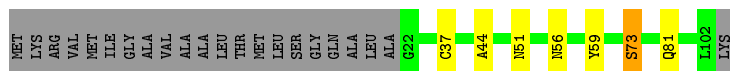
- Molecule 1: Cytochrome c554



- Molecule 1: Cytochrome c554



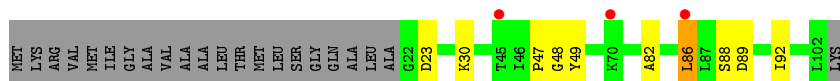
- Molecule 1: Cytochrome c554



- Molecule 1: Cytochrome c554



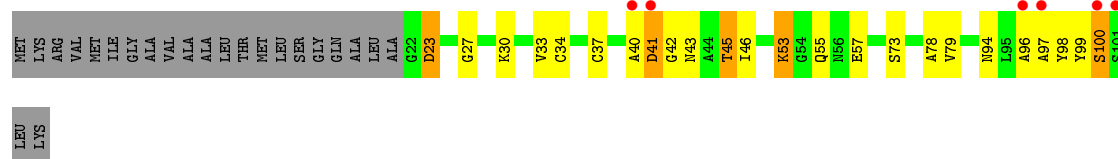
- Molecule 1: Cytochrome c554



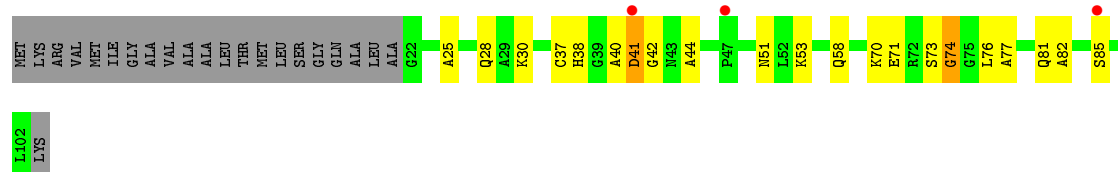
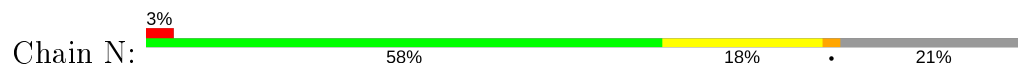
- Molecule 1: Cytochrome c554



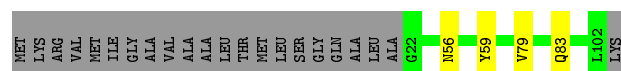
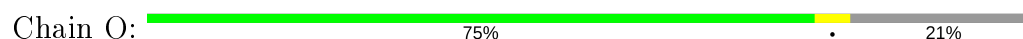
- Molecule 1: Cytochrome c554



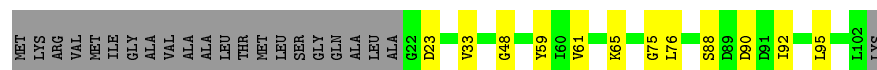
- Molecule 1: Cytochrome c554



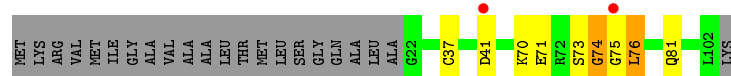
- Molecule 1: Cytochrome c554



- Molecule 1: Cytochrome c554



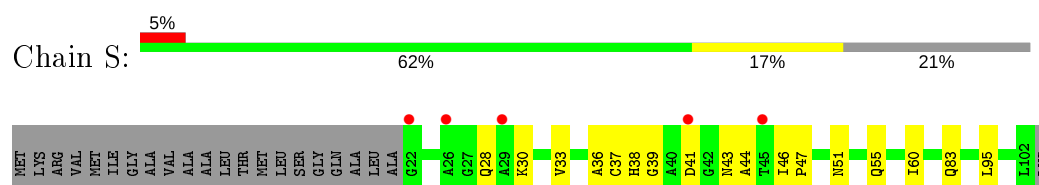
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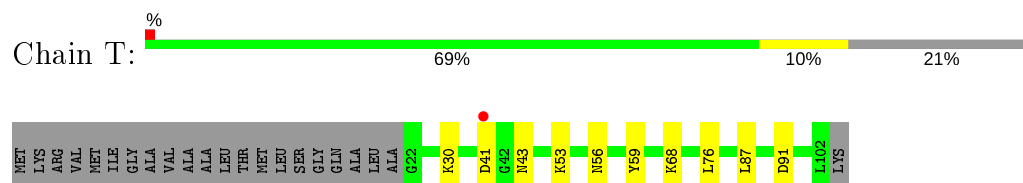
- Molecule 1: Cytochrome c554



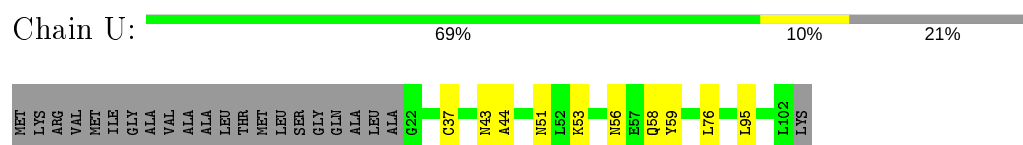
- Molecule 1: Cytochrome c554



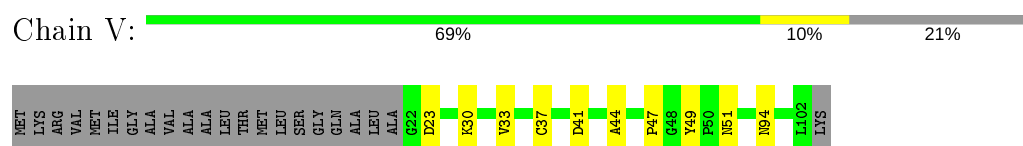
- Molecule 1: Cytochrome c554



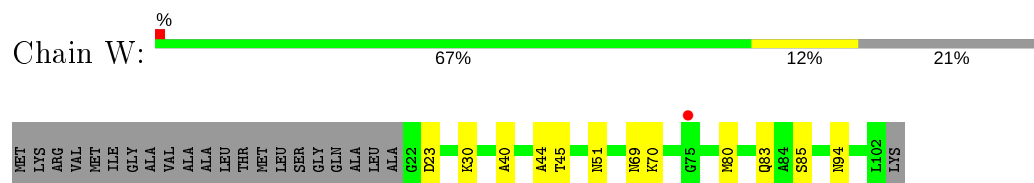
- Molecule 1: Cytochrome c554



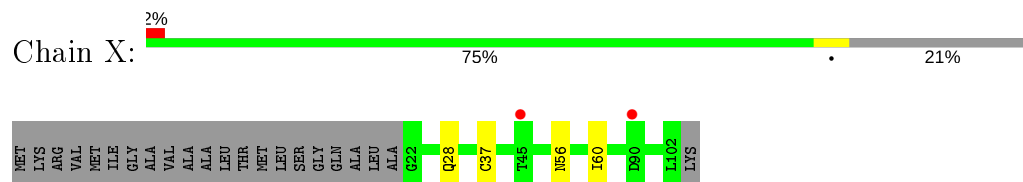
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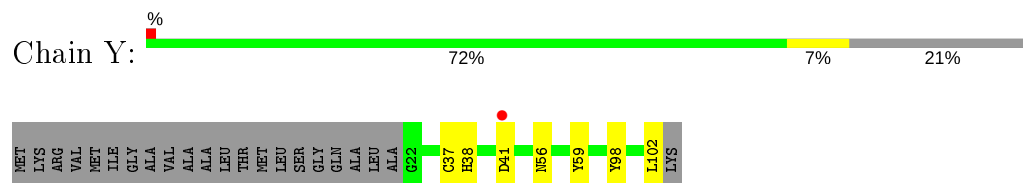
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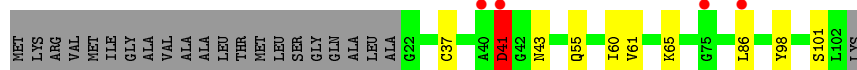
- Molecule 1: Cytochrome c554



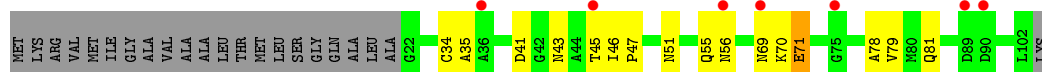
- Molecule 1: Cytochrome c554



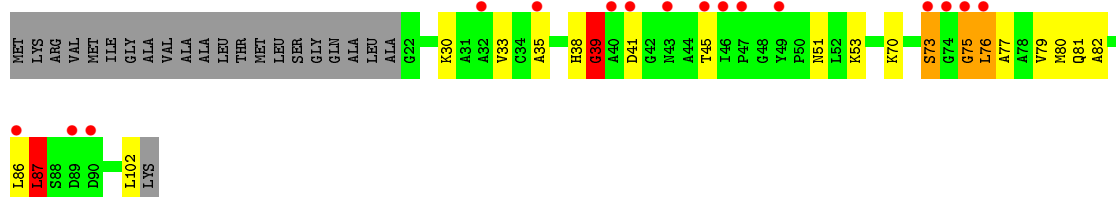
- Molecule 1: Cytochrome c554



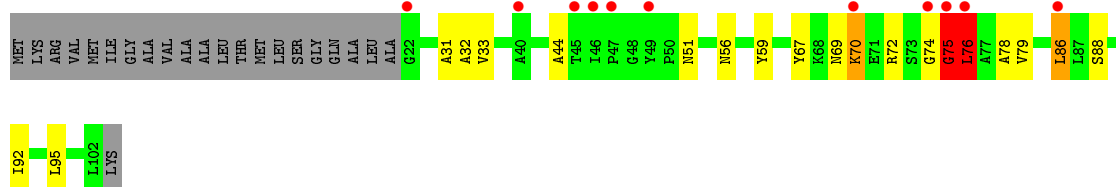
- Molecule 1: Cytochrome c554



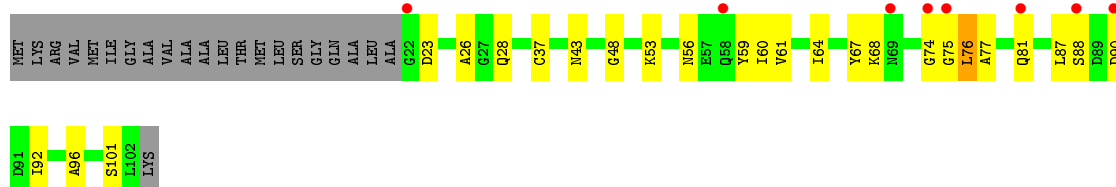
- Molecule 1: Cytochrome c554



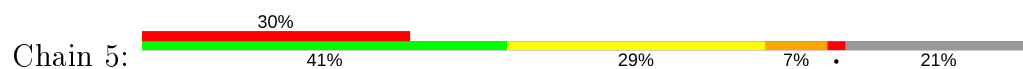
- Molecule 1: Cytochrome c554

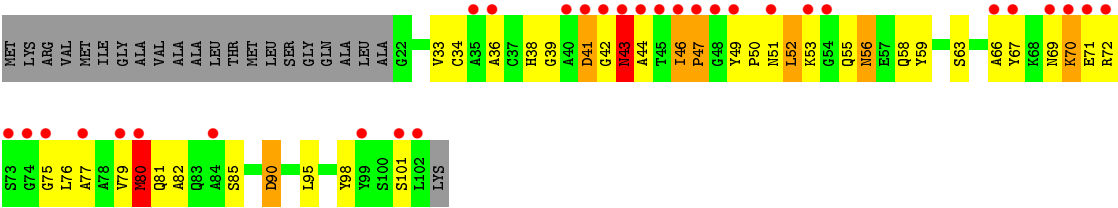


- Molecule 1: Cytochrome c554

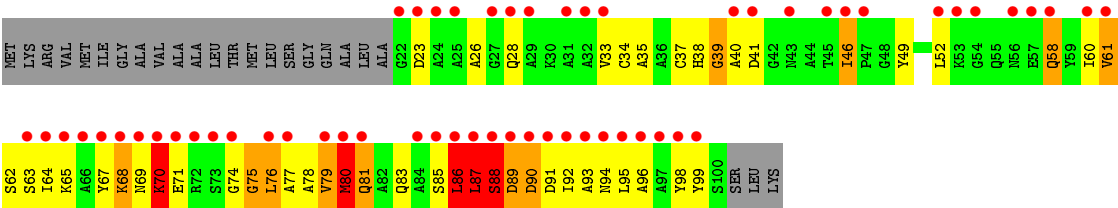
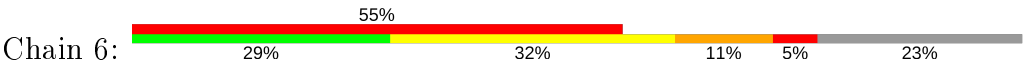


- Molecule 1: Cytochrome c554





● Molecule 1: Cytochrome c554



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	84.95Å 87.61Å 103.85Å 71.50° 72.98° 83.68°	Depositor
Resolution (Å)	20.00 – 1.80 20.03 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-1.80) 94.2 (20.03-1.80)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.4.0067	Depositor
R, R_{free}	0.193 , 0.242 0.194 , 0.243	Depositor DCC
R_{free} test set	11820 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	18.2	Xtriage
Anisotropy	0.301	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 55.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.026 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	21402	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 8.22% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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	1	0.78	1/582 (0.2%)	0.79	0/787
1	2	0.62	0/582	0.75	1/787 (0.1%)
1	3	0.65	0/582	0.77	2/787 (0.3%)
1	4	0.62	0/582	0.65	0/787
1	5	0.71	0/582	0.89	1/787 (0.1%)
1	6	0.64	0/568	1.06	4/768 (0.5%)
1	A	1.04	1/582 (0.2%)	0.89	1/787 (0.1%)
1	B	0.97	0/582	0.86	0/787
1	C	0.86	0/582	0.78	0/787
1	D	0.79	0/582	0.75	0/787
1	E	0.88	0/582	0.75	0/787
1	F	0.89	0/582	0.86	0/787
1	G	0.81	0/582	0.71	0/787
1	H	0.80	0/582	0.80	1/787 (0.1%)
1	I	0.74	0/582	0.76	0/787
1	J	0.65	0/582	0.71	0/787
1	K	0.74	0/582	0.73	0/787
1	L	0.80	0/582	0.82	3/787 (0.4%)
1	M	0.92	0/574	0.97	1/776 (0.1%)
1	N	0.81	1/582 (0.2%)	0.75	0/787
1	O	0.85	0/582	0.81	0/787
1	P	0.80	1/582 (0.2%)	0.78	0/787
1	Q	0.73	0/582	0.93	2/787 (0.3%)
1	R	0.80	0/582	0.72	0/787
1	S	0.80	0/582	0.80	0/787
1	T	0.91	0/582	0.84	0/787
1	U	0.85	0/582	0.75	0/787
1	V	0.88	1/582 (0.2%)	0.79	0/787
1	W	0.75	0/582	0.69	0/787
1	X	0.68	0/582	0.70	0/787
1	Y	0.79	0/582	0.73	0/787
1	Z	0.83	0/582	0.79	0/787

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
All	All	0.80	5/18602 (0.0%)	0.80	16/25154 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	2	0	3
1	3	0	2
1	6	0	5
1	L	0	2
1	Q	0	2
1	Z	0	1
All	All	0	15

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	N	37	CYS	CB-SG	-5.86	1.72	1.81
1	A	79	VAL	CB-CG2	5.42	1.64	1.52
1	P	59	TYR	CD1-CE1	5.15	1.47	1.39
1	V	49	TYR	CD2-CE2	5.12	1.47	1.39
1	1	34	CYS	CB-SG	-5.10	1.73	1.81

The worst 5 of 16 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Q	74	GLY	N-CA-C	11.32	141.40	113.10
1	6	87	LEU	N-CA-C	6.35	128.15	111.00
1	5	80	MET	CB-CG-SD	-6.07	94.18	112.40
1	L	76	LEU	N-CA-C	-5.98	94.86	111.00
1	2	39	GLY	N-CA-C	5.98	128.04	113.10

There are no chirality outliers.

5 of 15 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	2	38	HIS	Peptide
1	L	75	GLY	Mainchain,Peptide
1	Q	73	SER	Peptide

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Mol	Chain	Res	Type	Group
1	Q	75	GLY	Peptide
1	Z	41	ASP	Peptide

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	1	575	0	555	16	0
1	2	575	0	555	36	0
1	3	575	0	555	26	0
1	4	575	0	556	21	0
1	5	575	0	555	72	0
1	6	561	0	539	128	0
1	A	575	0	556	10	0
1	B	575	0	555	10	0
1	C	575	0	556	9	0
1	D	575	0	556	9	0
1	E	575	0	555	13	0
1	F	575	0	556	12	0
1	G	575	0	556	6	0
1	H	575	0	556	15	0
1	I	575	0	556	9	0
1	J	575	0	556	31	0
1	K	575	0	555	12	0
1	L	575	0	555	17	0
1	M	567	0	545	31	0
1	N	575	0	555	24	0
1	O	575	0	555	9	0
1	P	575	0	555	9	0
1	Q	575	0	556	6	0
1	R	575	0	555	15	0
1	S	575	0	556	13	0
1	T	575	0	555	6	0
1	U	575	0	555	17	0
1	V	575	0	556	12	0
1	W	575	0	555	10	0
1	X	575	0	556	4	0
1	Y	575	0	556	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Z	575	0	555	13	0
2	1	43	0	30	2	0
2	2	43	0	30	8	0
2	3	43	0	30	6	0
2	4	43	0	31	7	0
2	5	43	0	30	10	0
2	6	43	0	30	28	0
2	A	43	0	31	8	0
2	B	43	0	30	3	0
2	C	43	0	31	7	0
2	D	43	0	31	8	0
2	E	43	0	30	2	0
2	F	43	0	31	6	0
2	G	43	0	31	9	0
2	H	43	0	31	7	0
2	I	43	0	31	6	0
2	J	43	0	31	7	0
2	K	43	0	30	2	0
2	L	43	0	30	6	0
2	M	43	0	31	8	0
2	N	43	0	30	5	0
2	O	43	0	30	3	0
2	P	43	0	30	2	0
2	Q	43	0	31	6	0
2	R	43	0	30	1	0
2	S	43	0	31	10	0
2	T	43	0	30	2	0
2	U	43	0	31	6	0
2	V	43	0	31	5	0
2	W	43	0	30	2	0
2	X	43	0	31	7	0
2	Y	43	0	31	8	0
2	Z	43	0	30	4	0
3	C	6	0	8	0	0
3	E	6	0	8	0	0
3	P	6	0	8	0	0
3	R	6	0	8	7	0
4	1	34	0	0	2	0
4	2	15	0	0	0	0
4	3	21	0	0	0	0
4	4	28	0	0	1	0
4	5	27	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	6	9	0	0	0	0
4	A	91	0	0	1	0
4	B	101	0	0	1	0
4	C	73	0	0	1	0
4	D	62	0	0	2	0
4	E	76	0	0	2	0
4	F	64	0	0	2	0
4	G	62	0	0	0	0
4	H	70	0	0	2	0
4	I	62	0	0	1	0
4	J	24	0	0	0	0
4	K	35	0	0	0	0
4	L	53	0	0	0	0
4	M	51	0	0	3	0
4	N	37	0	0	0	0
4	O	71	0	0	0	0
4	P	35	0	0	3	0
4	Q	63	0	0	0	0
4	R	46	0	0	1	0
4	S	42	0	0	1	0
4	T	69	0	0	0	0
4	U	67	0	0	0	0
4	V	59	0	0	0	0
4	W	53	0	0	0	0
4	X	21	0	0	0	0
4	Y	55	0	0	0	0
4	Z	48	0	0	0	0
All	All	21402	0	18756	643	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 17.

The worst 5 of 643 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)
1:6:60:ILE:HD11	1:6:96:ALA:CB	1.43	1.45
1:5:44:ALA:CB	1:5:49:TYR:HB3	1.47	1.41
1:5:44:ALA:HB1	1:5:49:TYR:CB	1.53	1.36
1:6:88:SER:OG	1:6:92:ILE:HG13	1.25	1.34
1:5:39:GLY:HA3	1:5:43:ASN:ND2	1.44	1.31

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	1	79/103 (77%)	77 (98%)	2 (2%)	0	100	100
1	2	79/103 (77%)	72 (91%)	4 (5%)	3 (4%)	3	0
1	3	79/103 (77%)	76 (96%)	2 (2%)	1 (1%)	12	3
1	4	79/103 (77%)	75 (95%)	2 (2%)	2 (2%)	5	1
1	5	79/103 (77%)	70 (89%)	5 (6%)	4 (5%)	2	0
1	6	77/103 (75%)	63 (82%)	8 (10%)	6 (8%)	1	0
1	A	79/103 (77%)	79 (100%)	0	0	100	100
1	B	79/103 (77%)	79 (100%)	0	0	100	100
1	C	79/103 (77%)	76 (96%)	2 (2%)	1 (1%)	12	3
1	D	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
1	E	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
1	F	79/103 (77%)	76 (96%)	3 (4%)	0	100	100
1	G	79/103 (77%)	77 (98%)	2 (2%)	0	100	100
1	H	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
1	I	79/103 (77%)	77 (98%)	2 (2%)	0	100	100
1	J	79/103 (77%)	76 (96%)	3 (4%)	0	100	100
1	K	79/103 (77%)	79 (100%)	0	0	100	100
1	L	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
1	M	78/103 (76%)	75 (96%)	3 (4%)	0	100	100
1	N	79/103 (77%)	76 (96%)	1 (1%)	2 (2%)	5	1
1	O	79/103 (77%)	79 (100%)	0	0	100	100
1	P	79/103 (77%)	78 (99%)	1 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	Q	79/103 (77%)	75 (95%)	2 (2%)	2 (2%)	5	1
1	R	79/103 (77%)	76 (96%)	3 (4%)	0	100	100
1	S	79/103 (77%)	79 (100%)	0	0	100	100
1	T	79/103 (77%)	79 (100%)	0	0	100	100
1	U	79/103 (77%)	79 (100%)	0	0	100	100
1	V	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
1	W	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
1	X	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
1	Y	79/103 (77%)	79 (100%)	0	0	100	100
1	Z	79/103 (77%)	78 (99%)	1 (1%)	0	100	100
All	All	2525/3296 (77%)	2451 (97%)	53 (2%)	21 (1%)	19	7

5 of 21 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	N	41	ASP
1	Q	74	GLY
1	Q	76	LEU
1	2	39	GLY
1	5	43	ASN

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	1	55/70 (79%)	52 (94%)	3 (6%)	21	8
1	2	55/70 (79%)	51 (93%)	4 (7%)	14	4
1	3	55/70 (79%)	51 (93%)	4 (7%)	14	4
1	4	55/70 (79%)	51 (93%)	4 (7%)	14	4
1	5	55/70 (79%)	47 (86%)	8 (14%)	3	0
1	6	53/70 (76%)	43 (81%)	10 (19%)	1	0

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	55/70 (79%)	55 (100%)	0	100	100
1	B	55/70 (79%)	55 (100%)	0	100	100
1	C	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	D	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	E	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	F	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	G	55/70 (79%)	55 (100%)	0	100	100
1	H	55/70 (79%)	55 (100%)	0	100	100
1	I	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	J	55/70 (79%)	53 (96%)	2 (4%)	35	20
1	K	55/70 (79%)	51 (93%)	4 (7%)	14	4
1	L	55/70 (79%)	53 (96%)	2 (4%)	35	20
1	M	54/70 (77%)	50 (93%)	4 (7%)	13	4
1	N	55/70 (79%)	55 (100%)	0	100	100
1	O	55/70 (79%)	55 (100%)	0	100	100
1	P	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	Q	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	R	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	S	55/70 (79%)	53 (96%)	2 (4%)	35	20
1	T	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	U	55/70 (79%)	55 (100%)	0	100	100
1	V	55/70 (79%)	55 (100%)	0	100	100
1	W	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	X	55/70 (79%)	54 (98%)	1 (2%)	59	48
1	Y	55/70 (79%)	55 (100%)	0	100	100
1	Z	55/70 (79%)	53 (96%)	2 (4%)	35	20
All	All	1757/2240 (78%)	1697 (97%)	60 (3%)	37	22

5 of 60 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	1	41	ASP
1	2	87	LEU

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Mol	Chain	Res	Type
1	6	79	VAL
1	2	30	LYS
1	3	70	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 74 such sidechains are listed below:

Mol	Chain	Res	Type
1	Q	81	GLN
1	U	51	ASN
1	5	56	ASN
1	R	43	ASN
1	T	43	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

36 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	HEC	R	220	1	26,50,50	2.13	6 (23%)	18,82,82	2.76	8 (44%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HEC	I	220	1	26,50,50	2.56	8 (30%)	18,82,82	1.99	5 (27%)
2	HEC	Q	220	1	26,50,50	2.12	8 (30%)	18,82,82	2.54	7 (38%)
2	HEC	C	220	1	26,50,50	2.55	8 (30%)	18,82,82	1.94	6 (33%)
2	HEC	B	220	1	26,50,50	2.55	9 (34%)	18,82,82	2.40	6 (33%)
2	HEC	D	220	1	26,50,50	2.04	6 (23%)	18,82,82	2.27	6 (33%)
2	HEC	L	220	1	26,50,50	1.96	5 (19%)	18,82,82	1.80	5 (27%)
2	HEC	4	220	1	26,50,50	2.40	6 (23%)	18,82,82	2.16	6 (33%)
2	HEC	W	220	1	26,50,50	2.11	7 (26%)	18,82,82	2.34	8 (44%)
2	HEC	Z	220	1	26,50,50	2.23	5 (19%)	18,82,82	2.52	6 (33%)
2	HEC	F	220	1	26,50,50	2.38	5 (19%)	18,82,82	2.23	7 (38%)
2	HEC	Y	220	1	26,50,50	2.07	5 (19%)	18,82,82	2.29	6 (33%)
2	HEC	K	220	1	26,50,50	2.03	6 (23%)	18,82,82	2.14	6 (33%)
2	HEC	M	220	1	26,50,50	2.34	6 (23%)	18,82,82	2.21	7 (38%)
2	HEC	5	220	1	26,50,50	2.14	7 (26%)	18,82,82	2.75	7 (38%)
2	HEC	T	220	1	26,50,50	2.42	7 (26%)	18,82,82	2.08	5 (27%)
2	HEC	6	220	1	26,50,50	2.18	8 (30%)	18,82,82	2.36	7 (38%)
3	GOL	C	104	-	5,5,5	0.25	0	5,5,5	0.93	0
3	GOL	R	104	-	5,5,5	0.64	0	5,5,5	0.91	0
3	GOL	E	104	-	5,5,5	0.25	0	5,5,5	0.80	0
3	GOL	P	104	-	5,5,5	0.29	0	5,5,5	0.48	0
2	HEC	N	220	1	26,50,50	1.99	6 (23%)	18,82,82	2.32	5 (27%)
2	HEC	A	220	1	26,50,50	2.46	12 (46%)	18,82,82	2.57	7 (38%)
2	HEC	3	220	1	26,50,50	2.35	7 (26%)	18,82,82	1.81	7 (38%)
2	HEC	2	220	1	26,50,50	2.18	8 (30%)	18,82,82	1.98	4 (22%)
2	HEC	U	220	1	26,50,50	2.28	8 (30%)	18,82,82	2.33	6 (33%)
2	HEC	V	220	1	26,50,50	2.43	6 (23%)	18,82,82	2.21	8 (44%)
2	HEC	X	220	1	26,50,50	2.45	9 (34%)	18,82,82	1.81	6 (33%)
2	HEC	1	220	1	26,50,50	2.57	8 (30%)	18,82,82	2.41	8 (44%)
2	HEC	E	220	1	26,50,50	2.17	5 (19%)	18,82,82	2.51	7 (38%)
2	HEC	G	220	1	26,50,50	2.03	9 (34%)	18,82,82	2.59	10 (55%)
2	HEC	H	220	1	26,50,50	2.21	10 (38%)	18,82,82	2.22	5 (27%)
2	HEC	J	220	1	26,50,50	2.35	11 (42%)	18,82,82	1.81	4 (22%)
2	HEC	O	220	1	26,50,50	2.28	7 (26%)	18,82,82	1.81	6 (33%)
2	HEC	P	220	1	26,50,50	2.12	8 (30%)	18,82,82	2.82	9 (50%)
2	HEC	S	220	1	26,50,50	2.77	10 (38%)	18,82,82	1.96	6 (33%)

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	HEC	R	220	1	-	0/6/54/54	-
2	HEC	I	220	1	-	0/6/54/54	-
2	HEC	Q	220	1	-	0/6/54/54	-
2	HEC	C	220	1	-	0/6/54/54	-
2	HEC	B	220	1	-	0/6/54/54	-
2	HEC	D	220	1	-	0/6/54/54	-
2	HEC	L	220	1	-	0/6/54/54	-
2	HEC	4	220	1	-	0/6/54/54	-
2	HEC	W	220	1	-	0/6/54/54	-
2	HEC	Z	220	1	-	0/6/54/54	-
2	HEC	F	220	1	-	0/6/54/54	-
2	HEC	Y	220	1	-	0/6/54/54	-
2	HEC	K	220	1	-	0/6/54/54	-
2	HEC	M	220	1	-	0/6/54/54	-
2	HEC	5	220	1	-	0/6/54/54	-
2	HEC	T	220	1	-	0/6/54/54	-
2	HEC	6	220	1	-	0/6/54/54	-
3	GOL	C	104	-	-	2/4/4/4	-
3	GOL	R	104	-	-	4/4/4/4	-
3	GOL	E	104	-	-	1/4/4/4	-
3	GOL	P	104	-	-	2/4/4/4	-
2	HEC	N	220	1	-	0/6/54/54	-
2	HEC	A	220	1	-	0/6/54/54	-
2	HEC	3	220	1	-	0/6/54/54	-
2	HEC	2	220	1	-	0/6/54/54	-
2	HEC	U	220	1	-	0/6/54/54	-
2	HEC	V	220	1	-	0/6/54/54	-
2	HEC	X	220	1	-	0/6/54/54	-
2	HEC	1	220	1	-	0/6/54/54	-
2	HEC	E	220	1	-	0/6/54/54	-
2	HEC	G	220	1	-	0/6/54/54	-
2	HEC	H	220	1	-	0/6/54/54	-
2	HEC	J	220	1	-	0/6/54/54	-
2	HEC	O	220	1	-	0/6/54/54	-
2	HEC	P	220	1	-	0/6/54/54	-
2	HEC	S	220	1	-	0/6/54/54	-

The worst 5 of 236 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	220	HEC	C3B-C2B	-8.21	1.32	1.40
2	I	220	HEC	C3B-C2B	-7.91	1.32	1.40
2	C	220	HEC	C3C-C2C	-7.44	1.33	1.40
2	S	220	HEC	C3B-C2B	-7.44	1.33	1.40
2	V	220	HEC	C3B-C2B	-7.25	1.33	1.40

The worst 5 of 206 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	5	220	HEC	CBD-CAD-C3D	-7.44	98.76	112.49
2	R	220	HEC	CMC-C2C-C1C	-6.12	119.06	128.46
2	A	220	HEC	CMC-C2C-C1C	-5.90	119.40	128.46
2	Z	220	HEC	CMB-C2B-C1B	-5.70	119.70	128.46
2	E	220	HEC	CMC-C2C-C3C	5.63	132.45	125.82

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	104	GOL	O1-C1-C2-C3
3	R	104	GOL	O1-C1-C2-C3
3	R	104	GOL	C1-C2-C3-O3
3	P	104	GOL	C1-C2-C3-O3
3	C	104	GOL	O1-C1-C2-O2

There are no ring outliers.

33 monomers are involved in 191 short contacts:

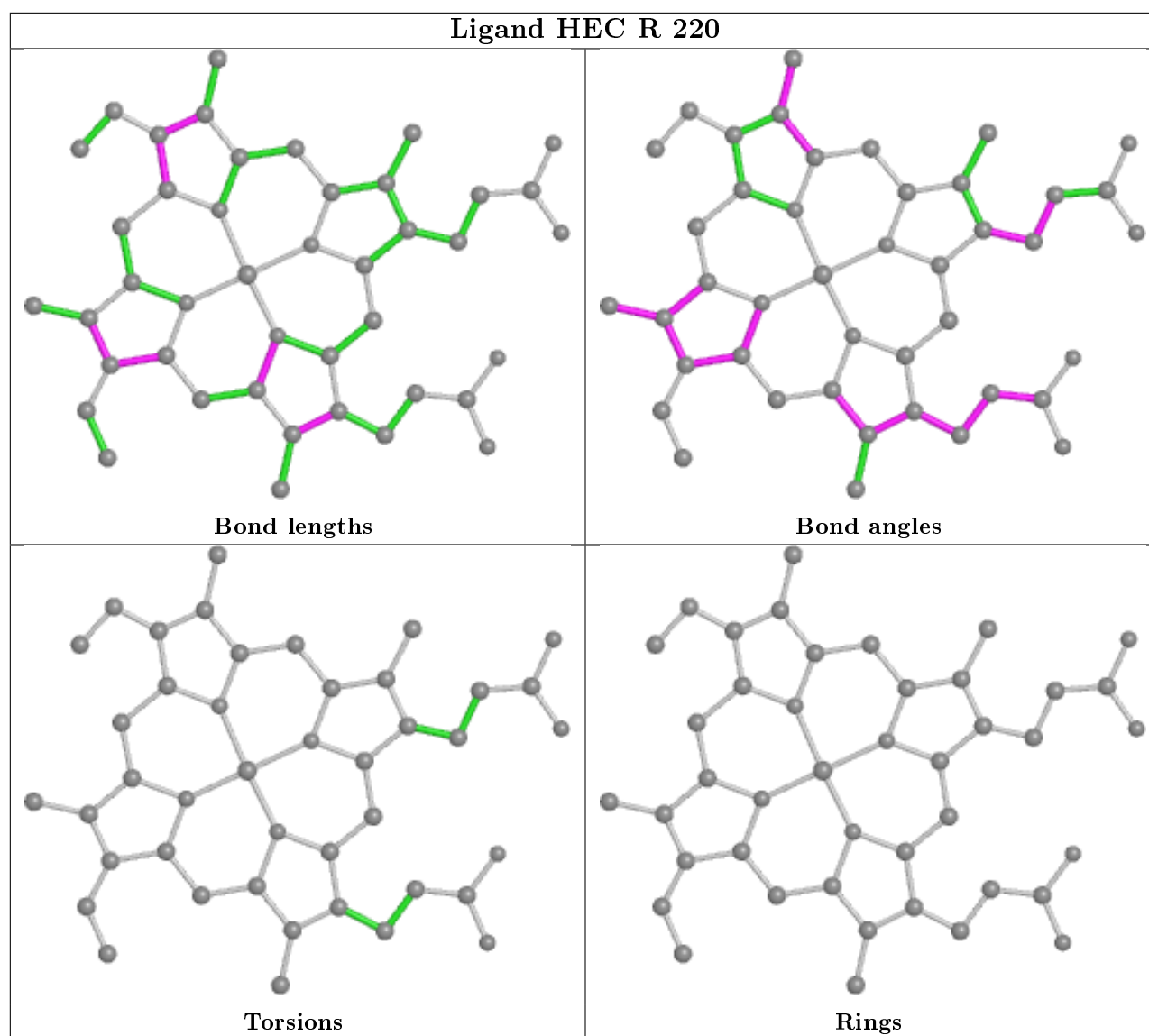
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	R	220	HEC	1	0
2	I	220	HEC	6	0
2	Q	220	HEC	6	0
2	C	220	HEC	7	0
2	B	220	HEC	3	0
2	D	220	HEC	8	0
2	L	220	HEC	6	0
2	4	220	HEC	7	0
2	W	220	HEC	2	0
2	Z	220	HEC	4	0
2	F	220	HEC	6	0
2	Y	220	HEC	8	0
2	K	220	HEC	2	0

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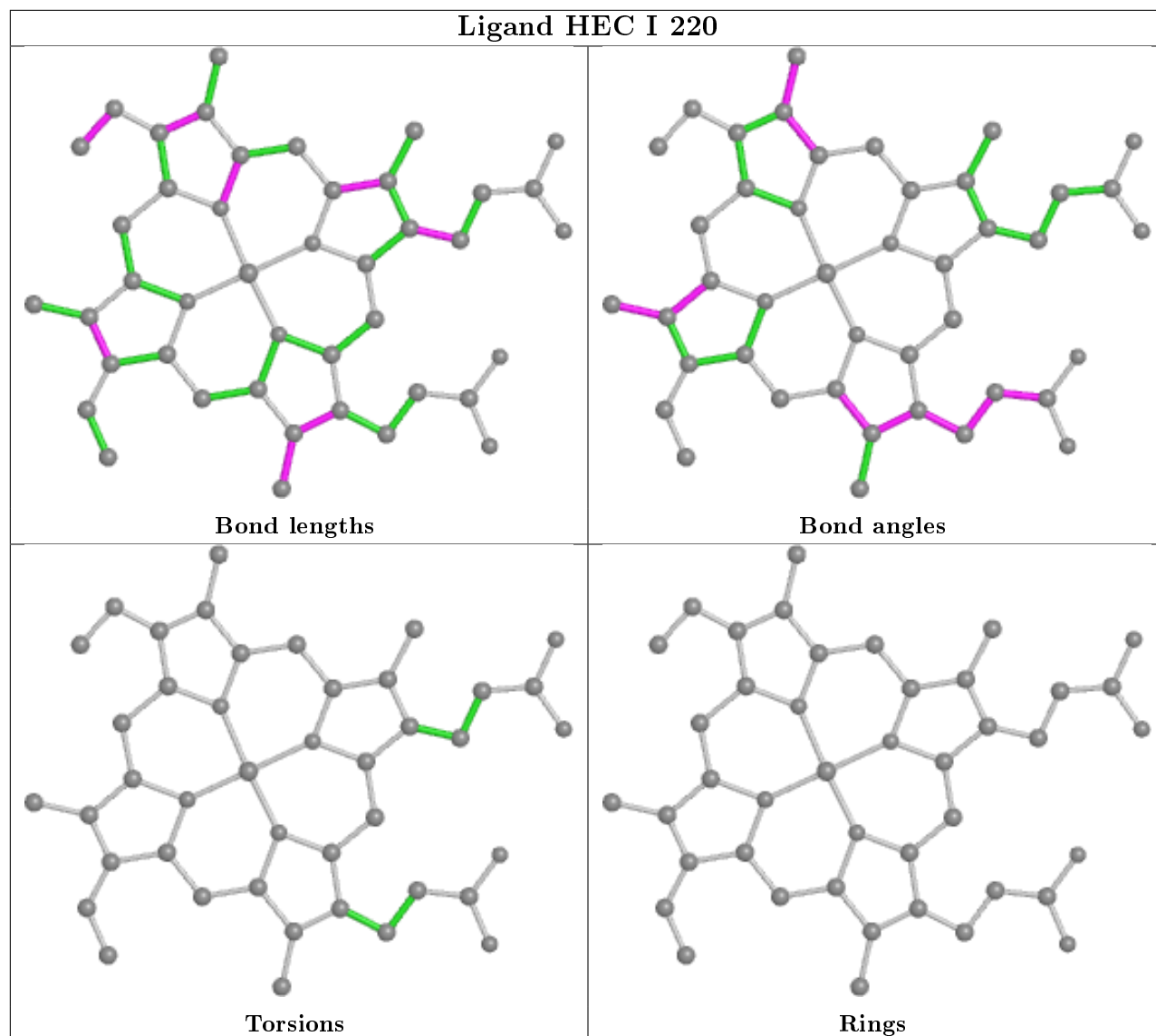
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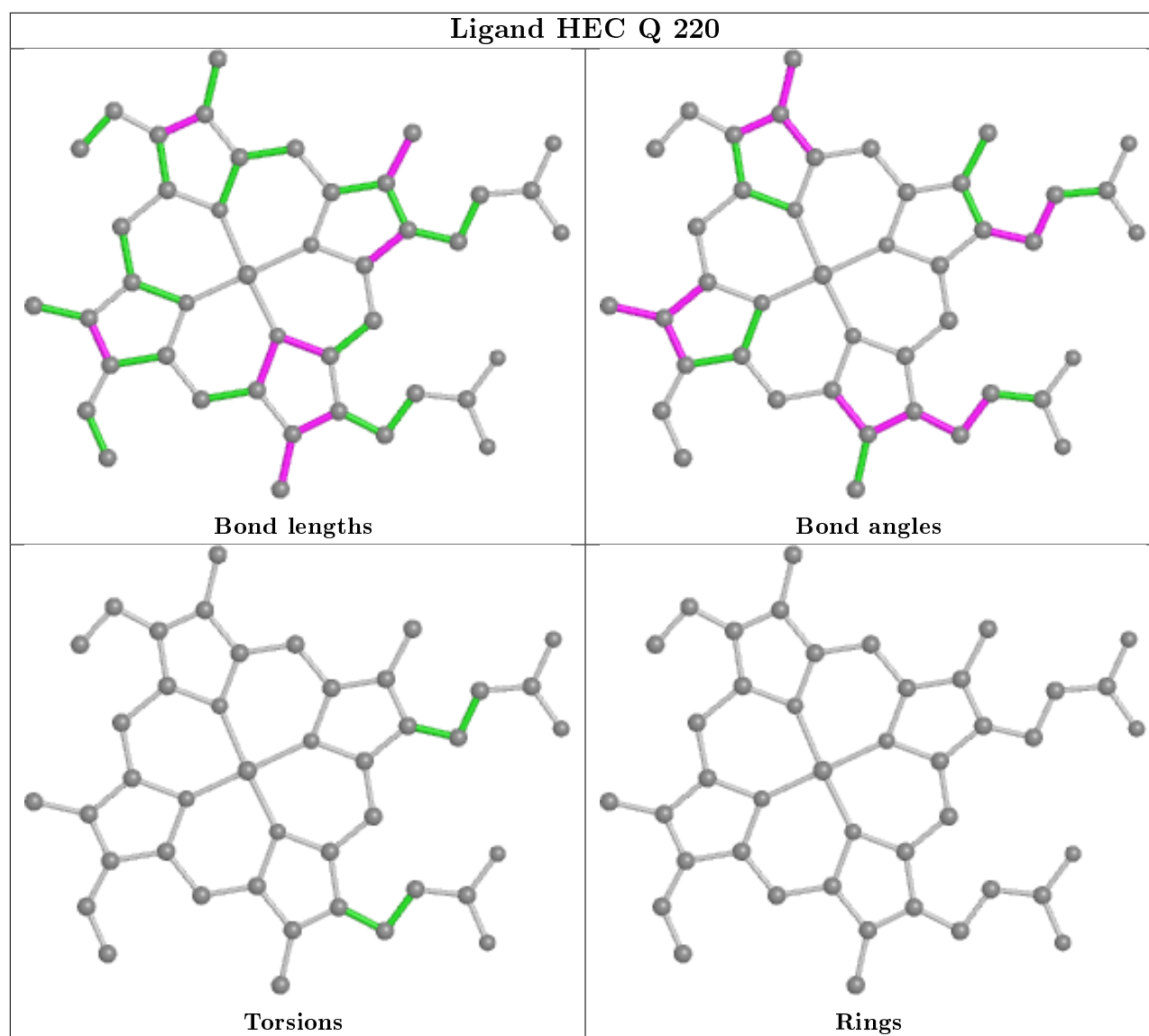
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	M	220	HEC	8	0
2	5	220	HEC	10	0
2	T	220	HEC	2	0
2	6	220	HEC	28	0
3	R	104	GOL	7	0
2	N	220	HEC	5	0
2	A	220	HEC	8	0
2	3	220	HEC	6	0
2	2	220	HEC	8	0
2	U	220	HEC	6	0
2	V	220	HEC	5	0
2	X	220	HEC	7	0
2	1	220	HEC	2	0
2	E	220	HEC	2	0
2	G	220	HEC	9	0
2	H	220	HEC	7	0
2	J	220	HEC	7	0
2	O	220	HEC	3	0
2	P	220	HEC	2	0
2	S	220	HEC	10	0

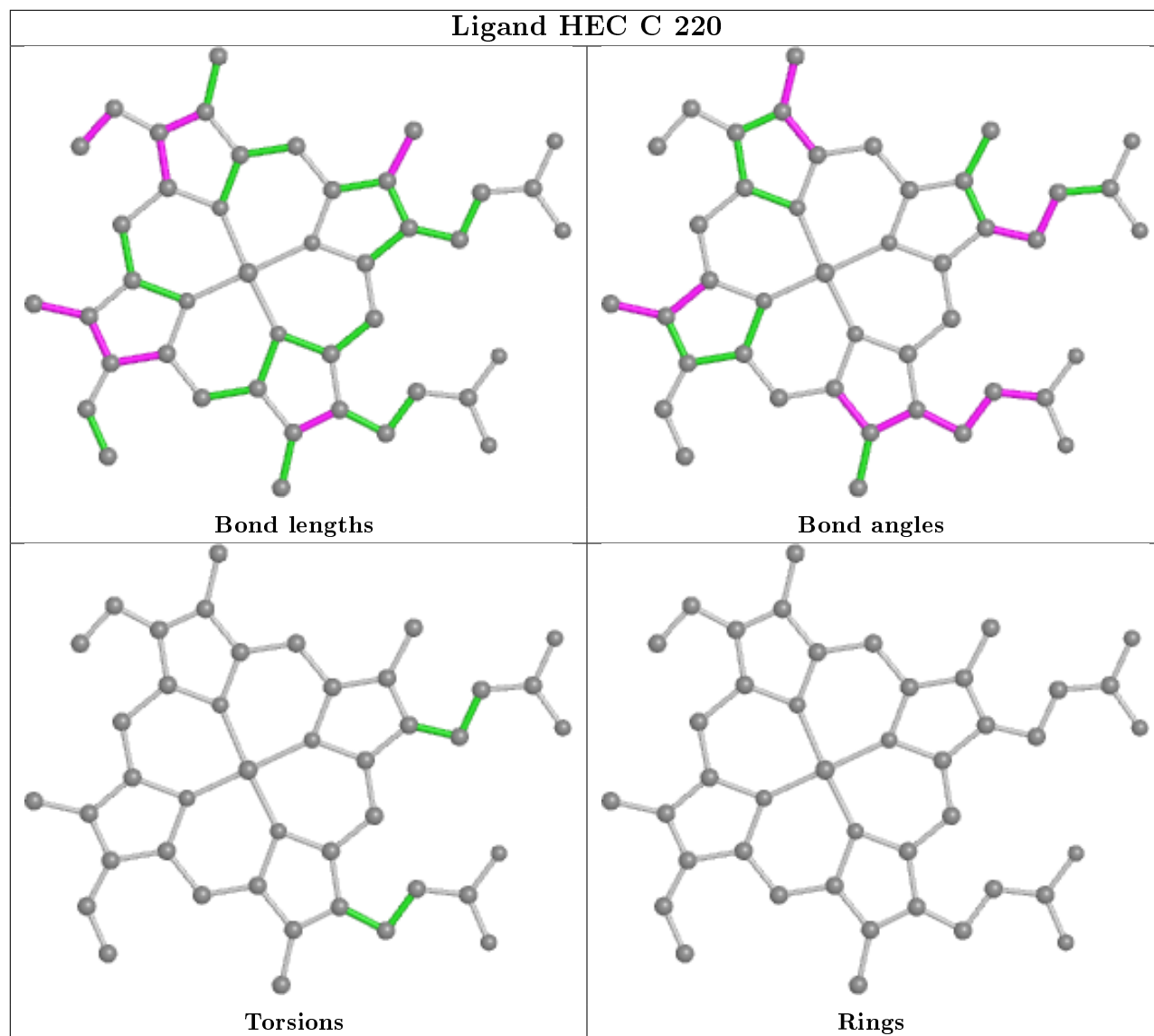
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

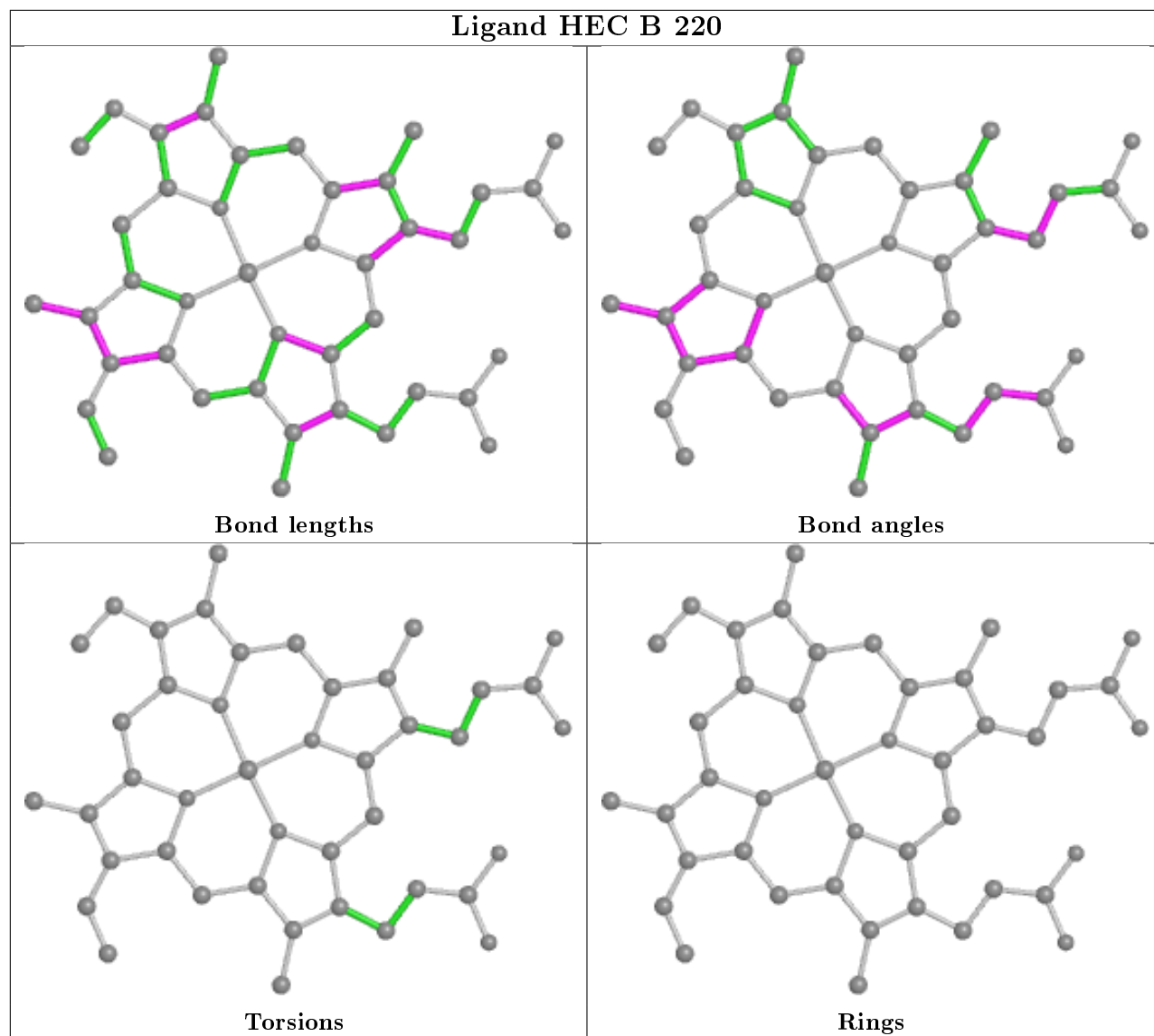


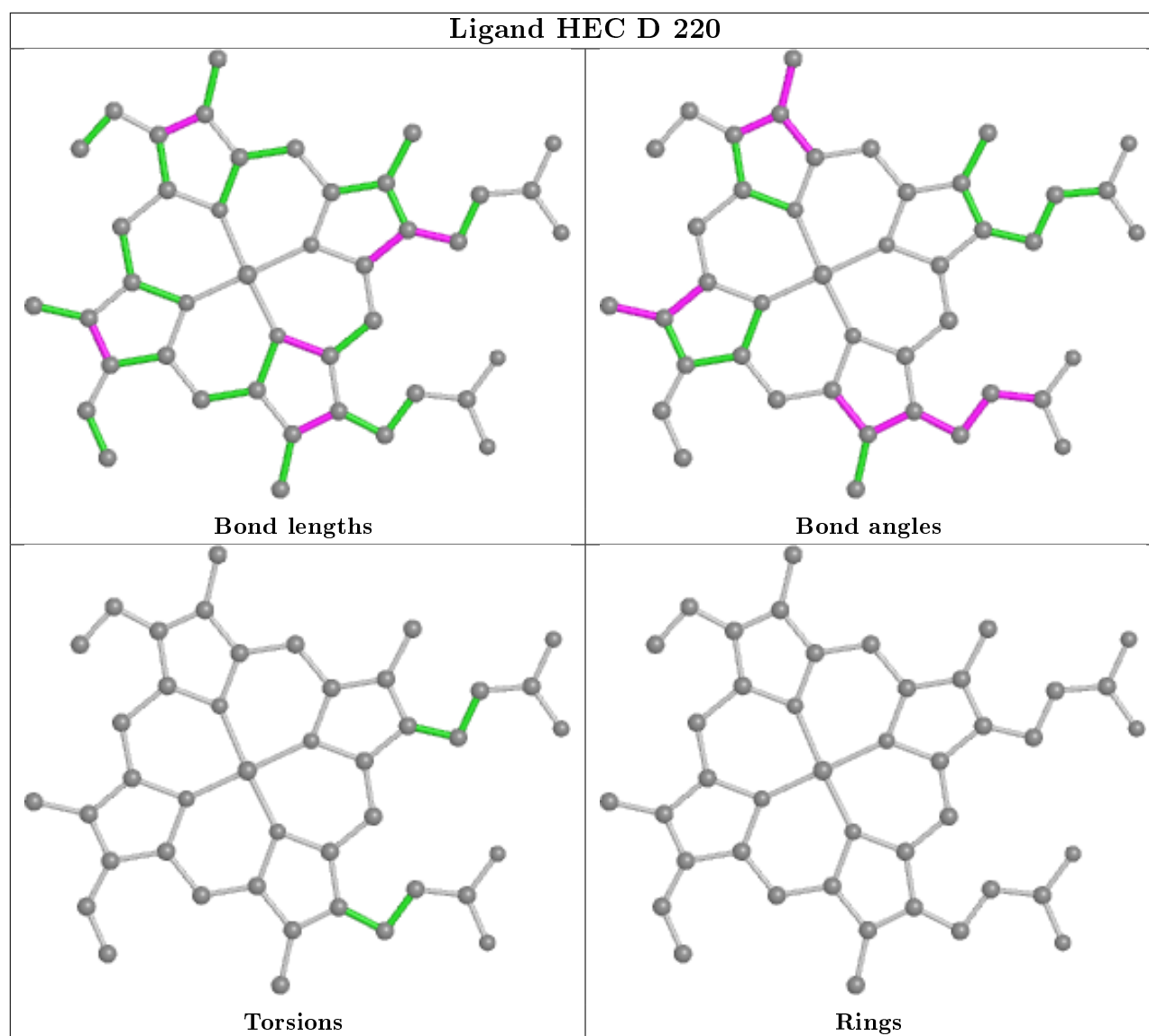
Ligand HEC I 220



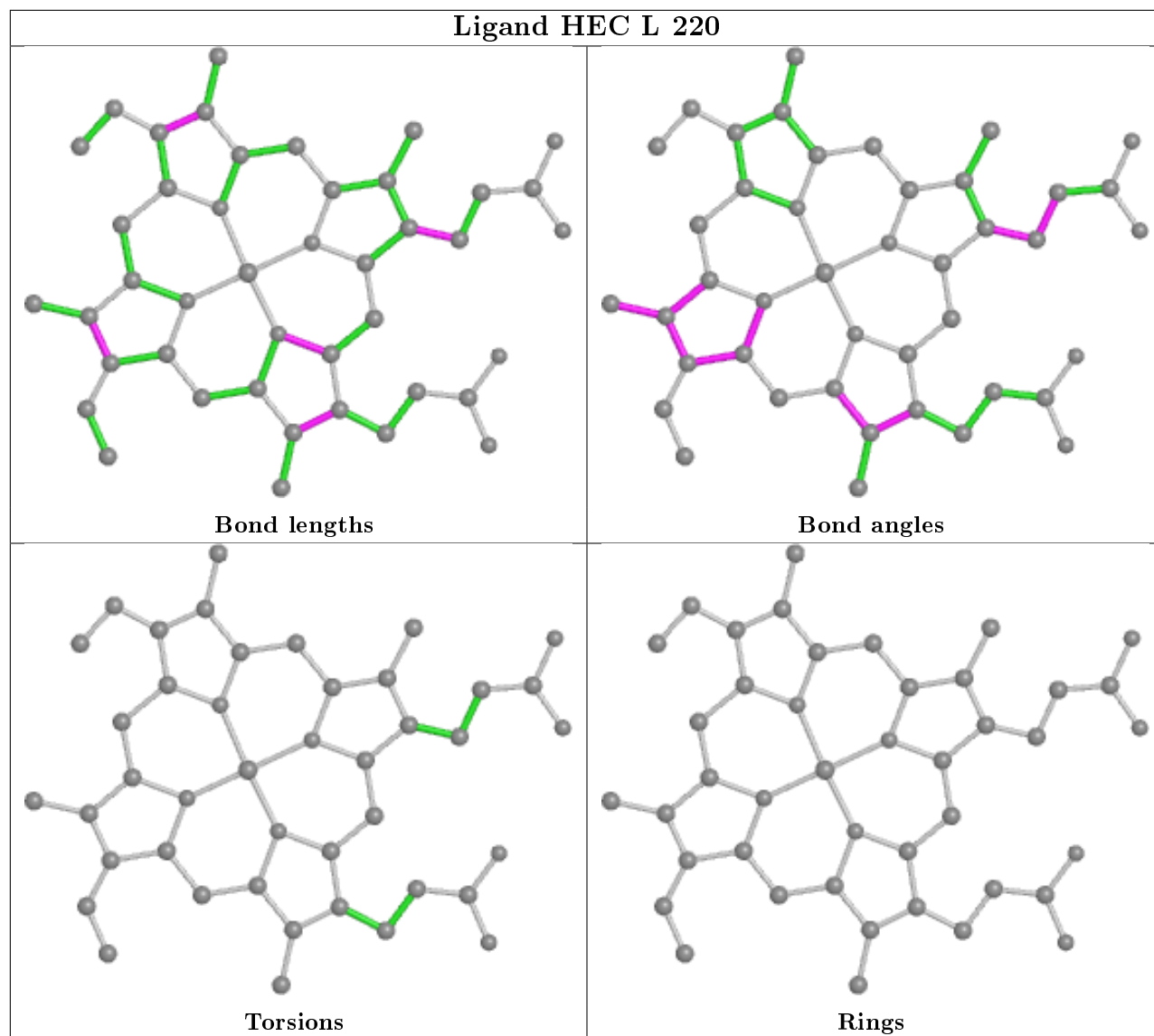




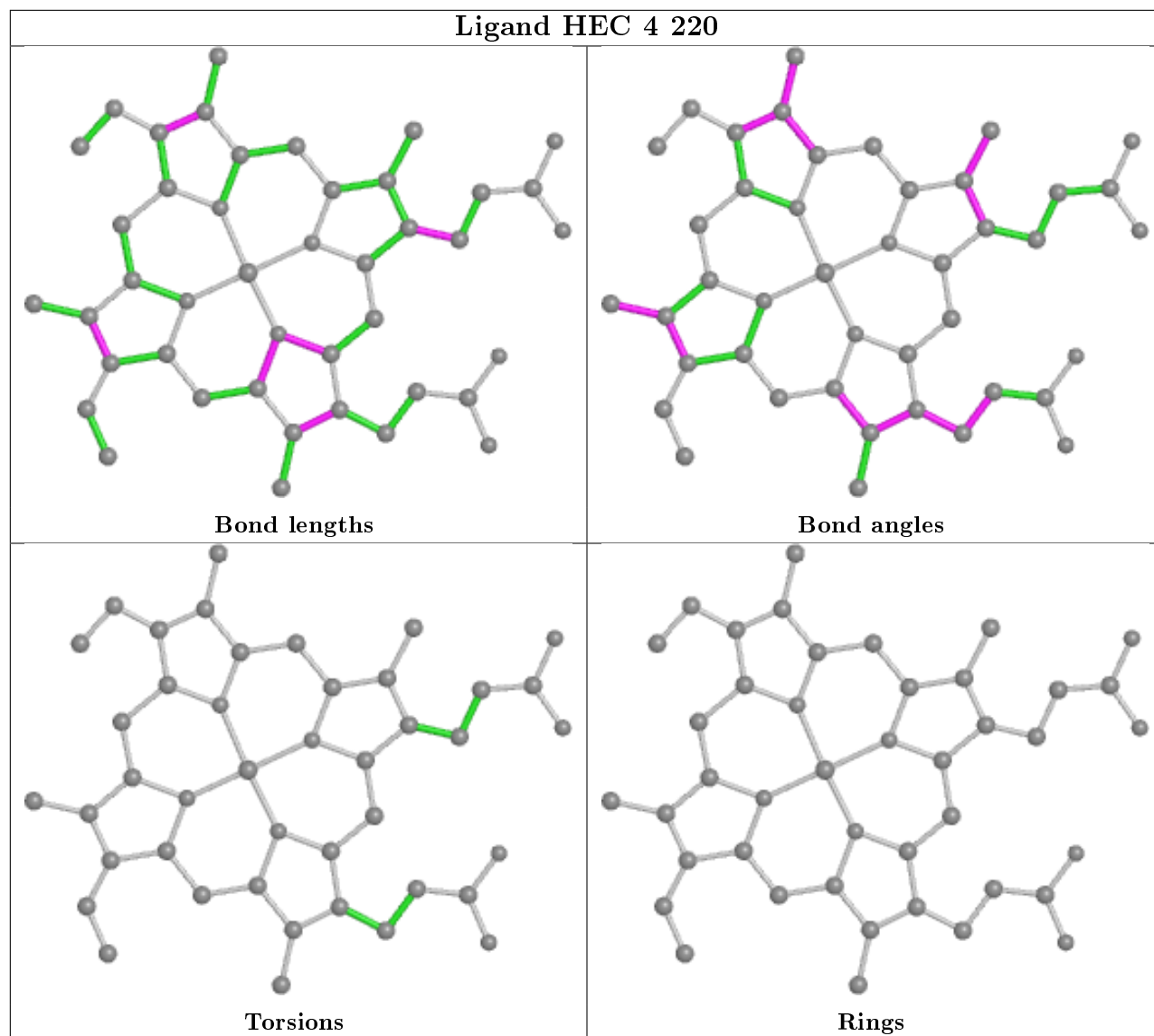


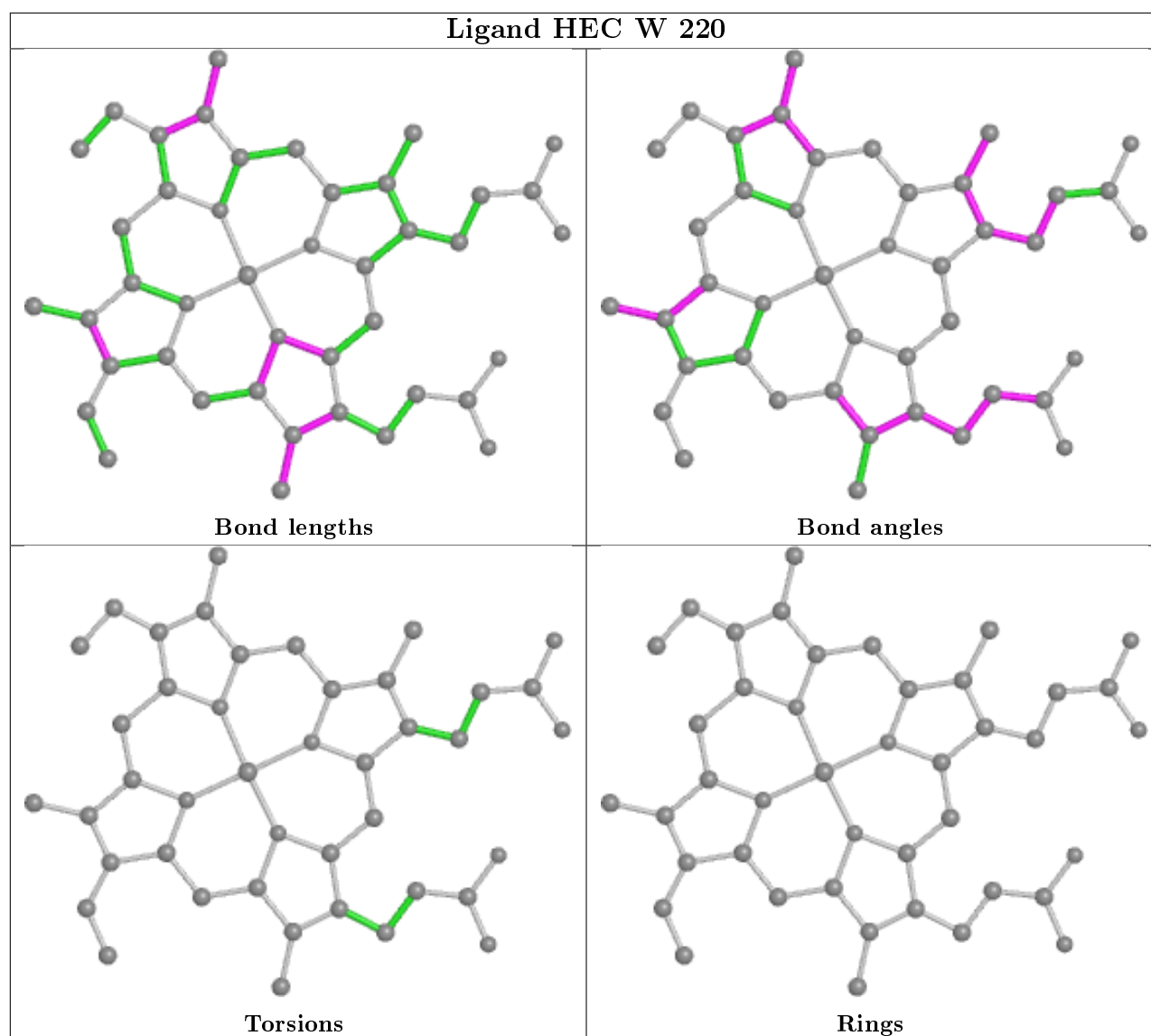


Ligand HEC L 220

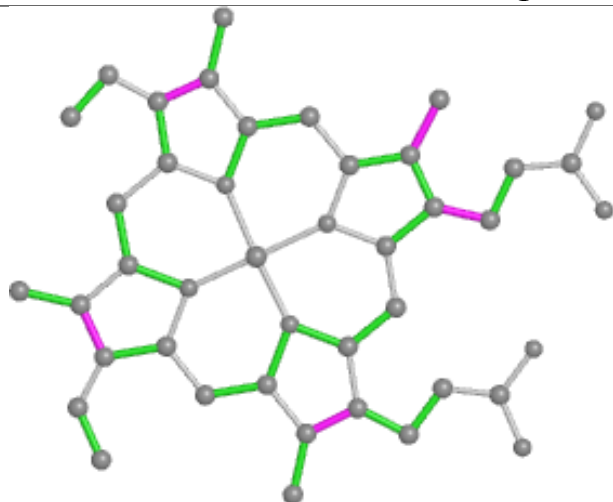


Ligand HEC 4 220

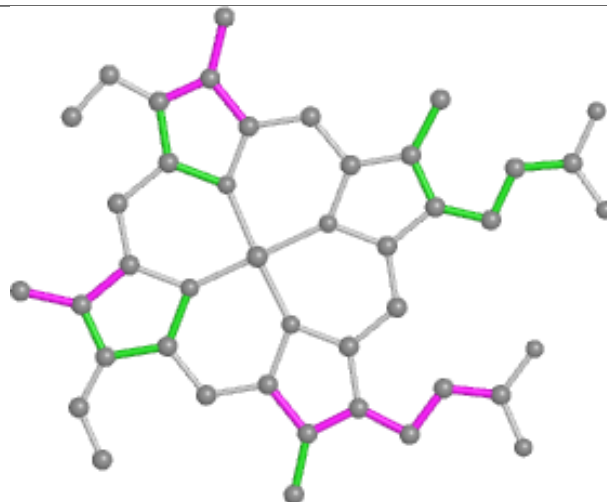




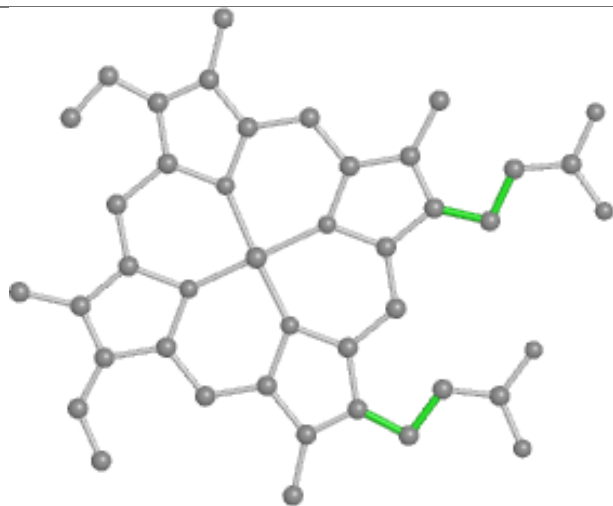
Ligand HEC Z 220



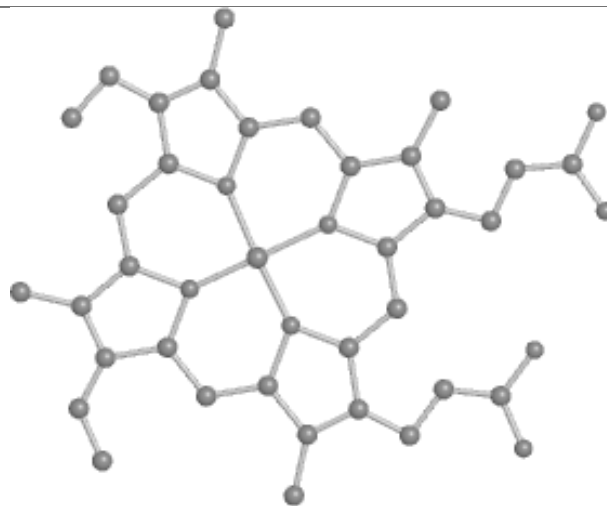
Bond lengths



Bond angles

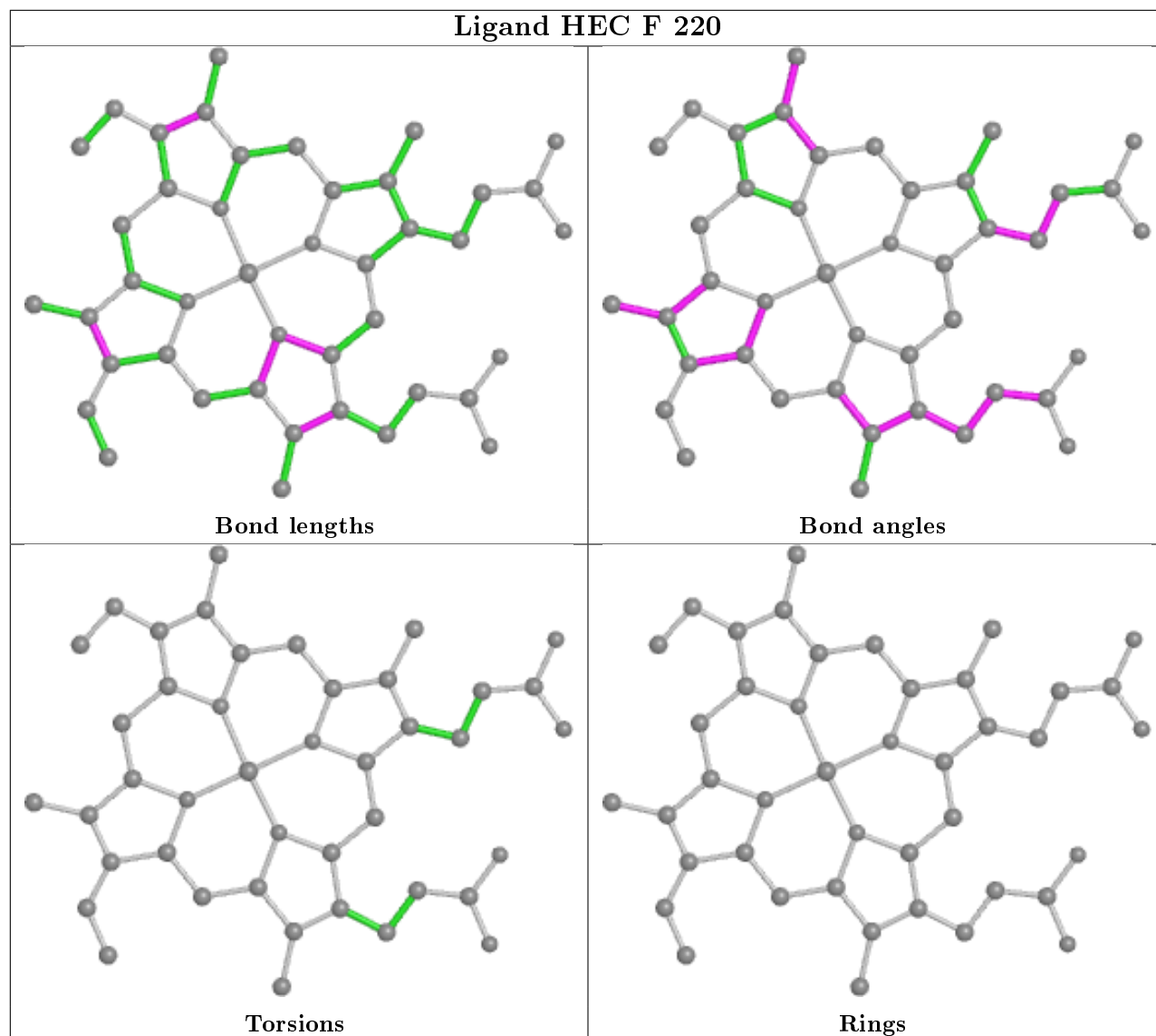


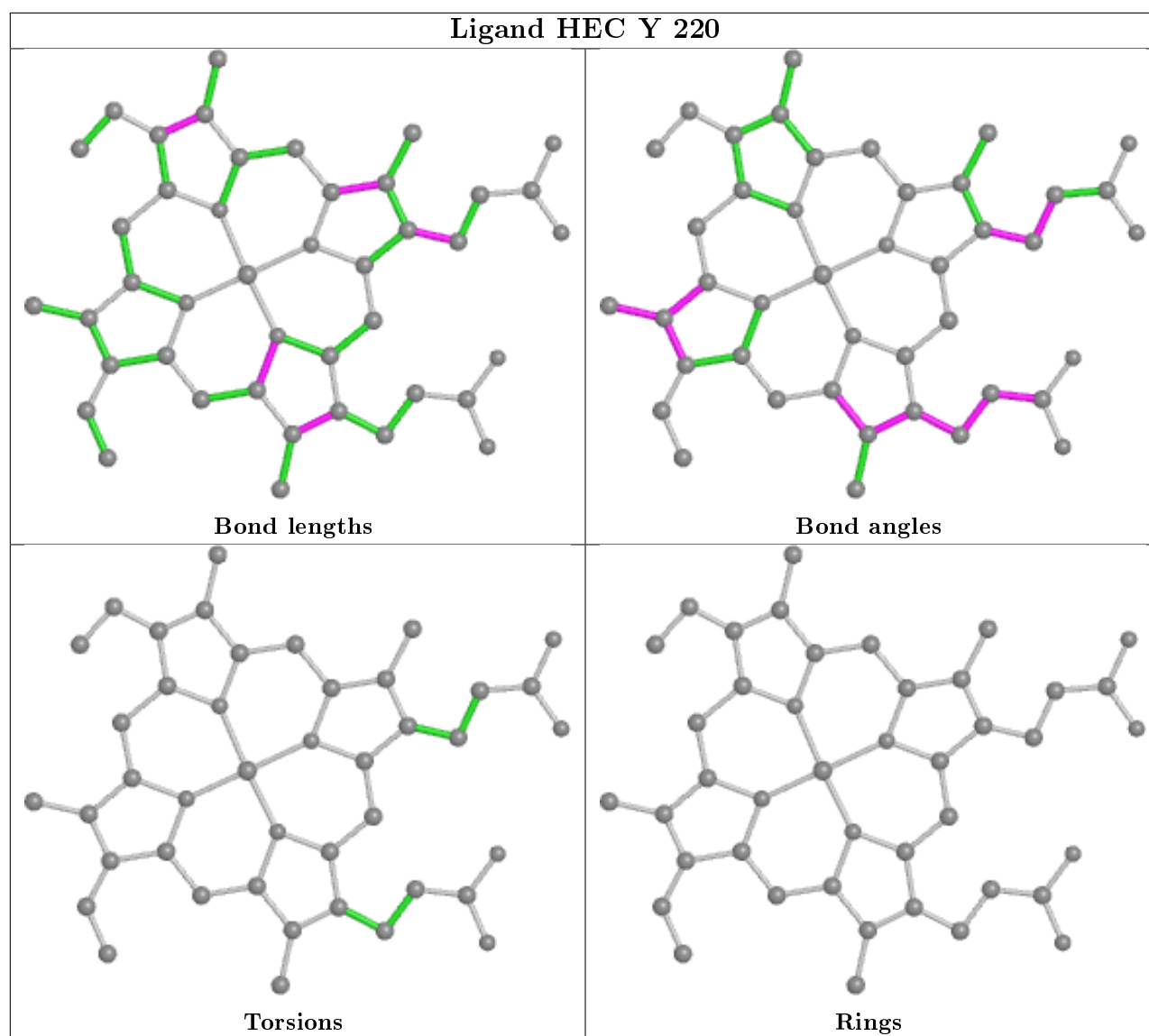
Torsions

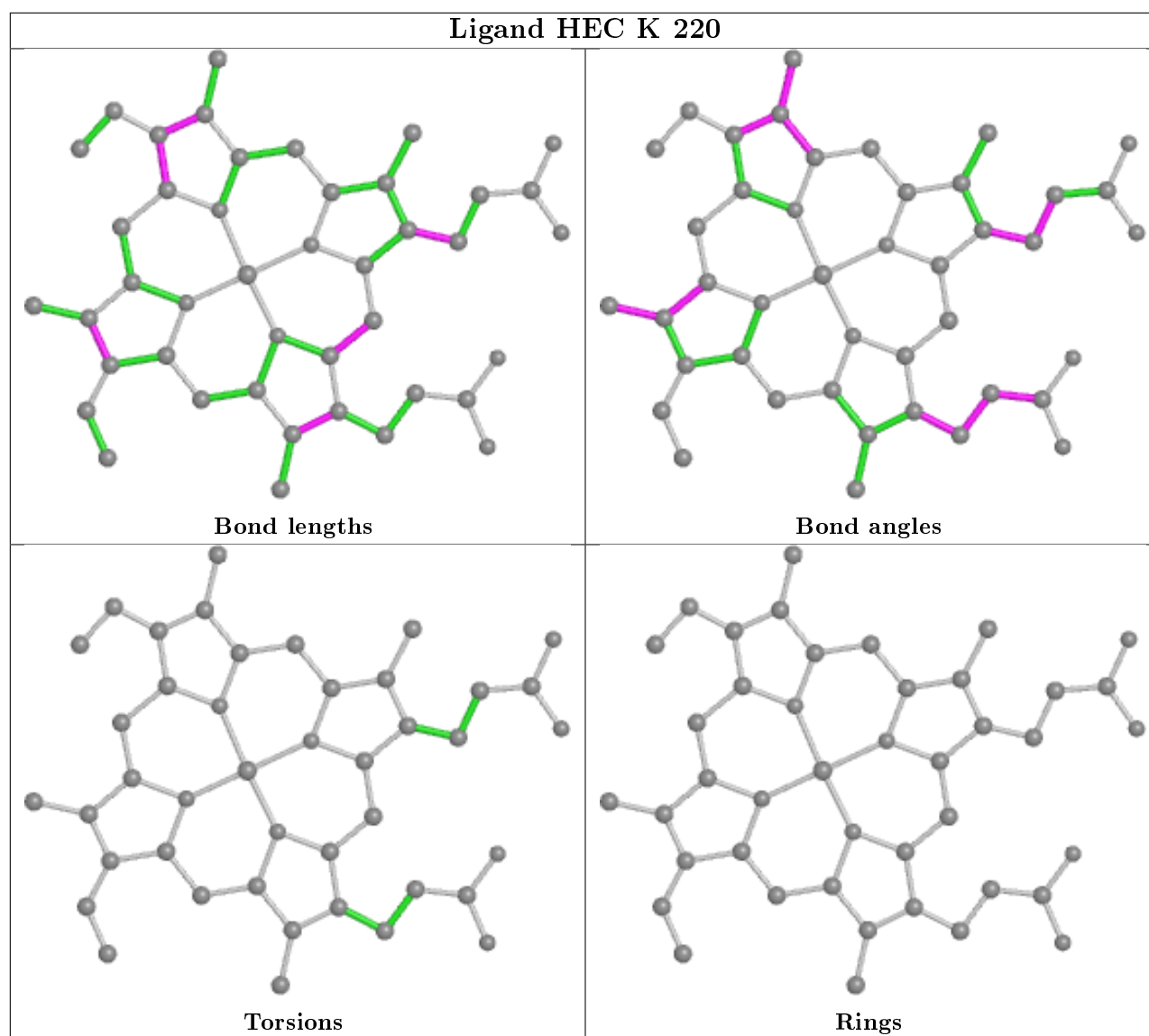


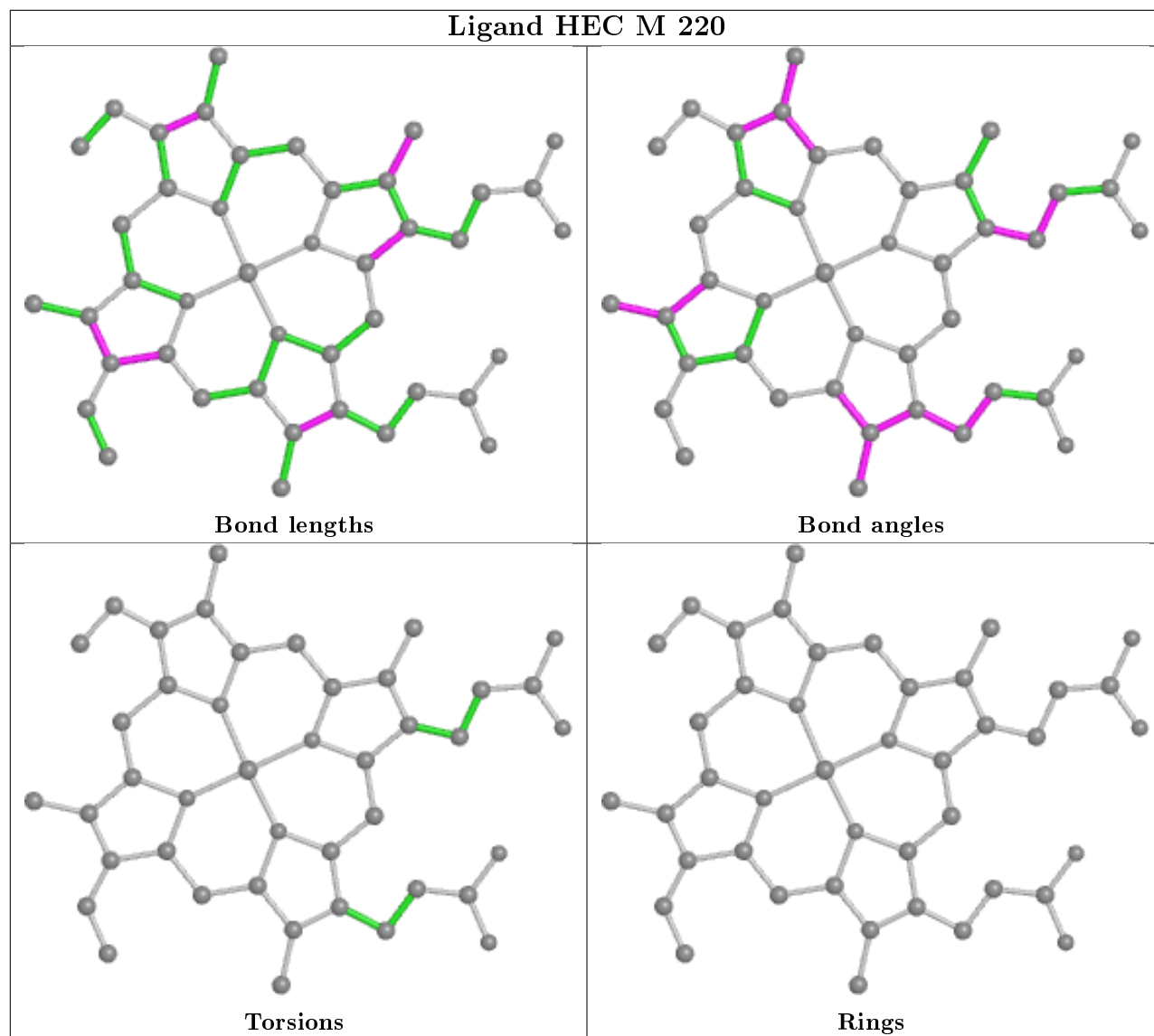
Rings

Ligand HEC F 220

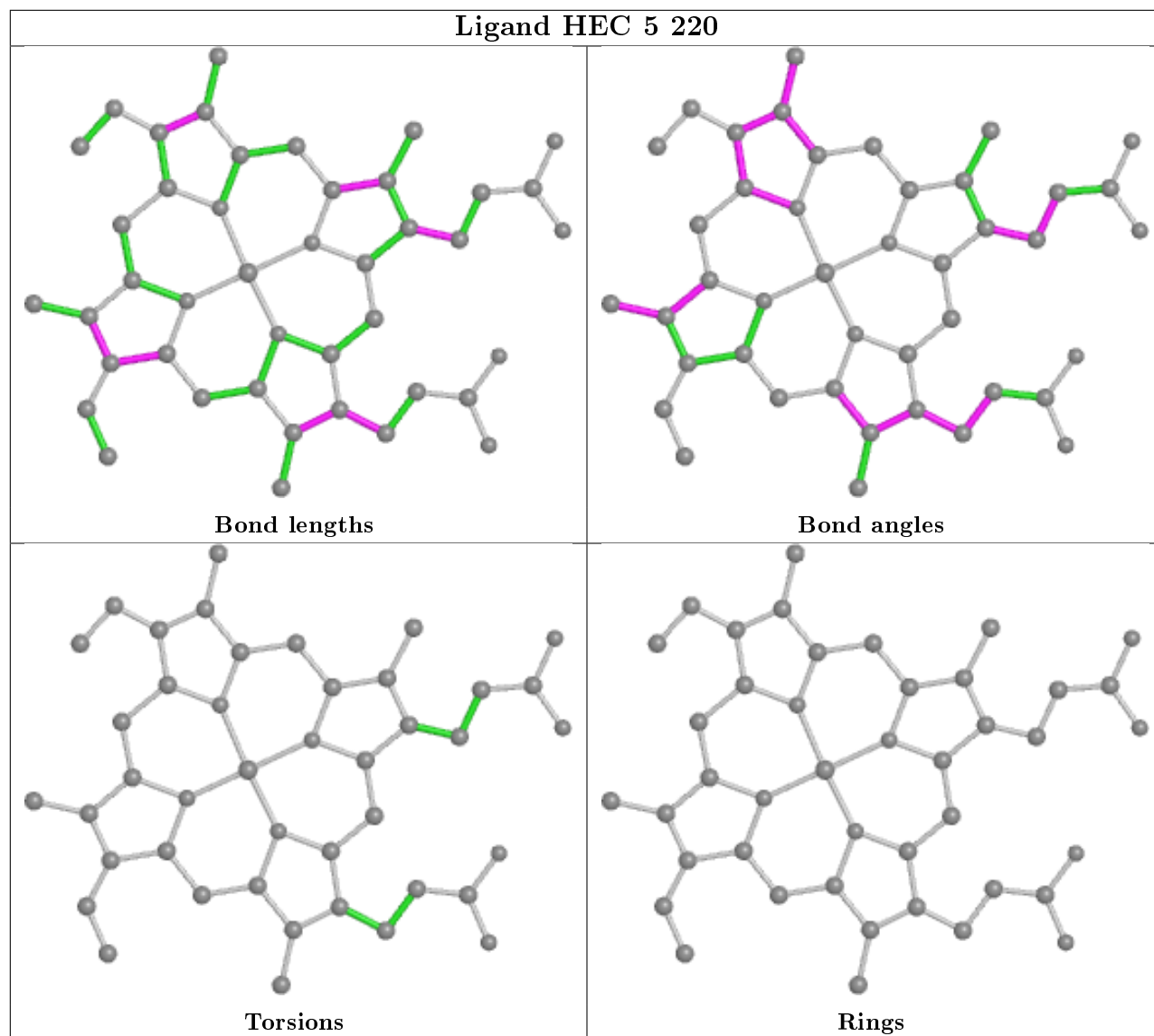




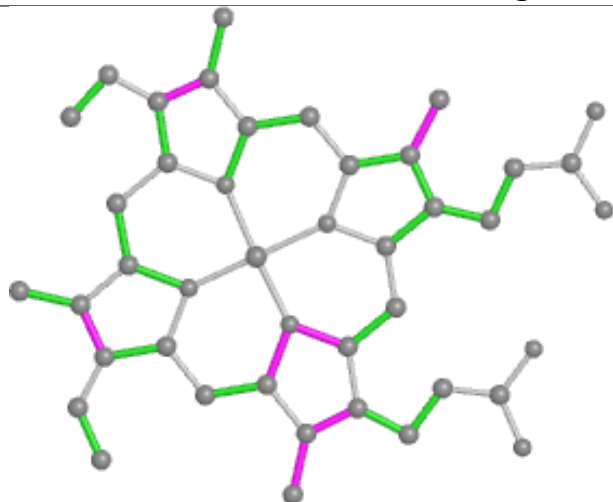




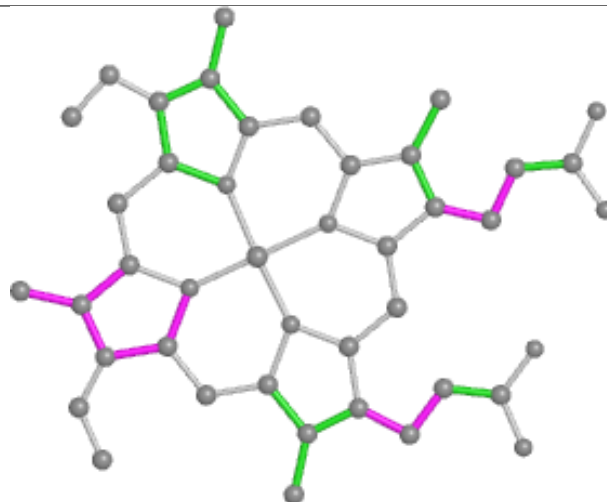
Ligand HEC 5 220



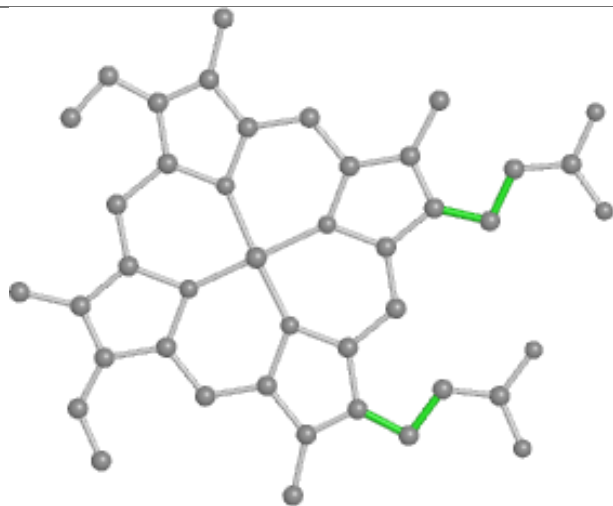
Ligand HEC T 220



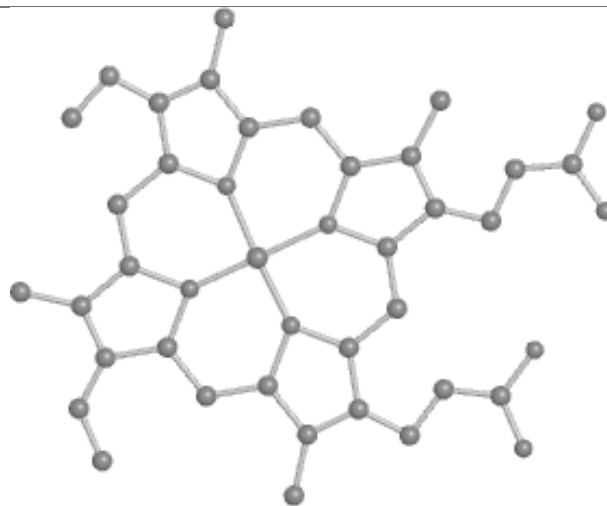
Bond lengths



Bond angles

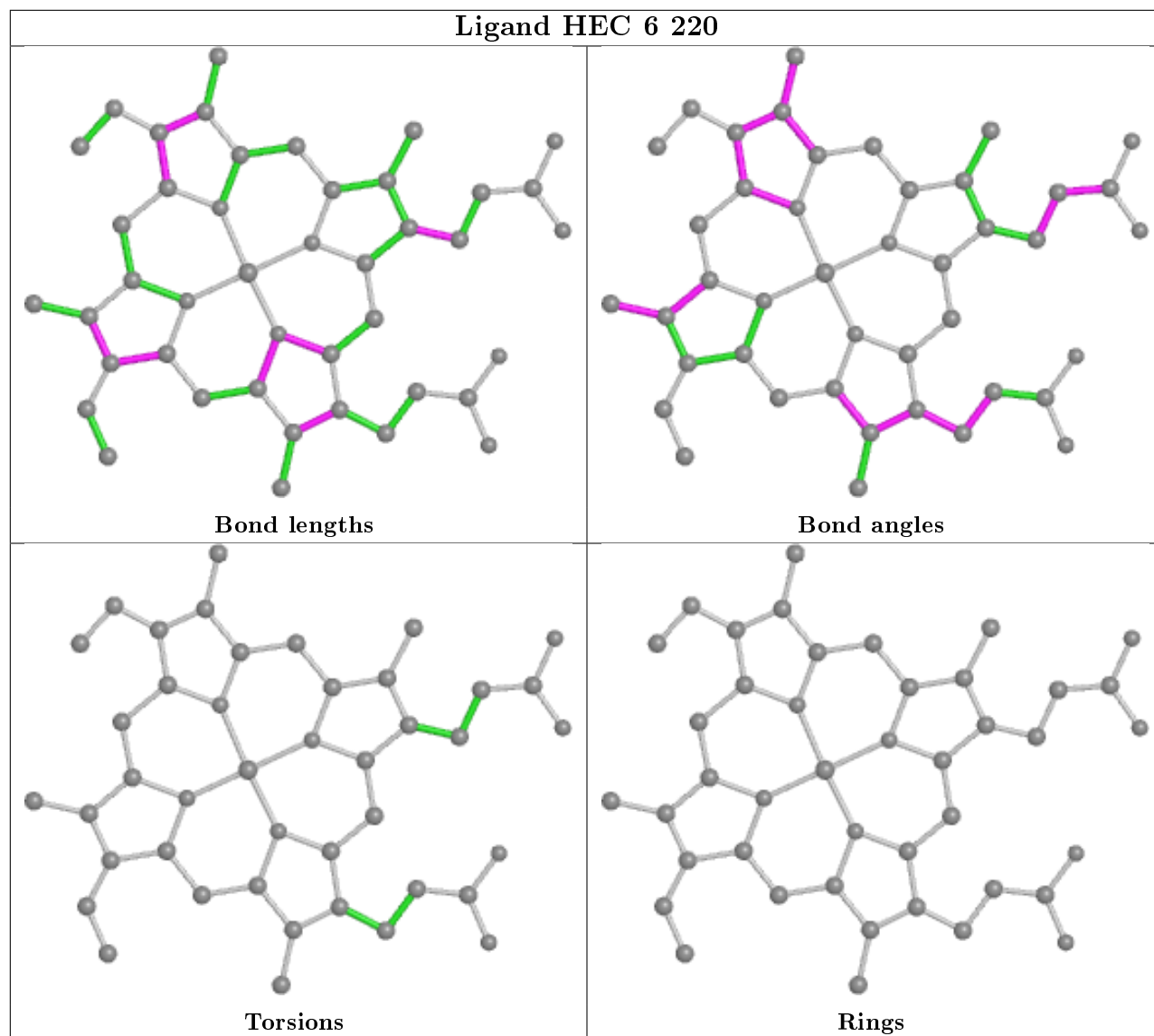


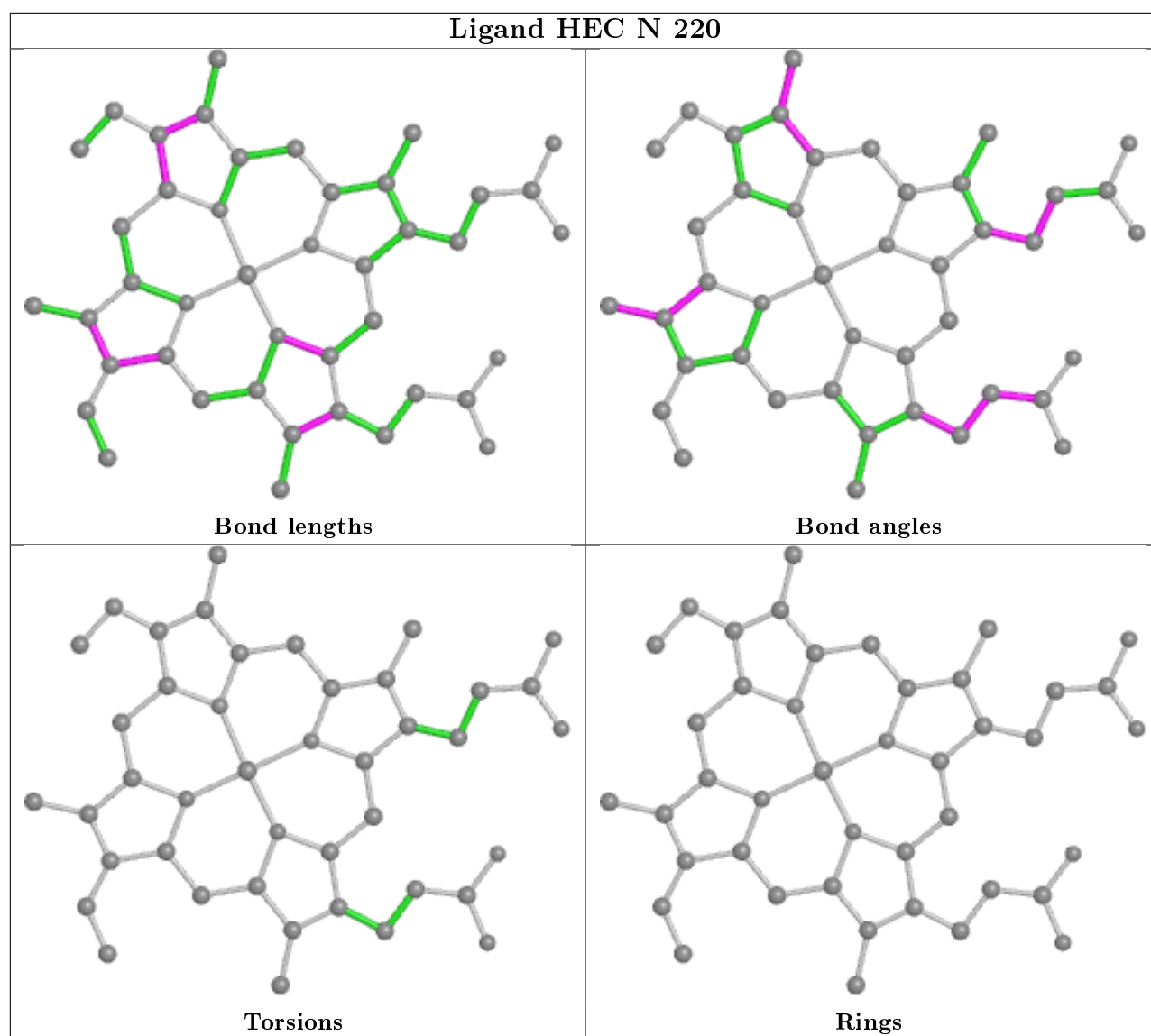
Torsions

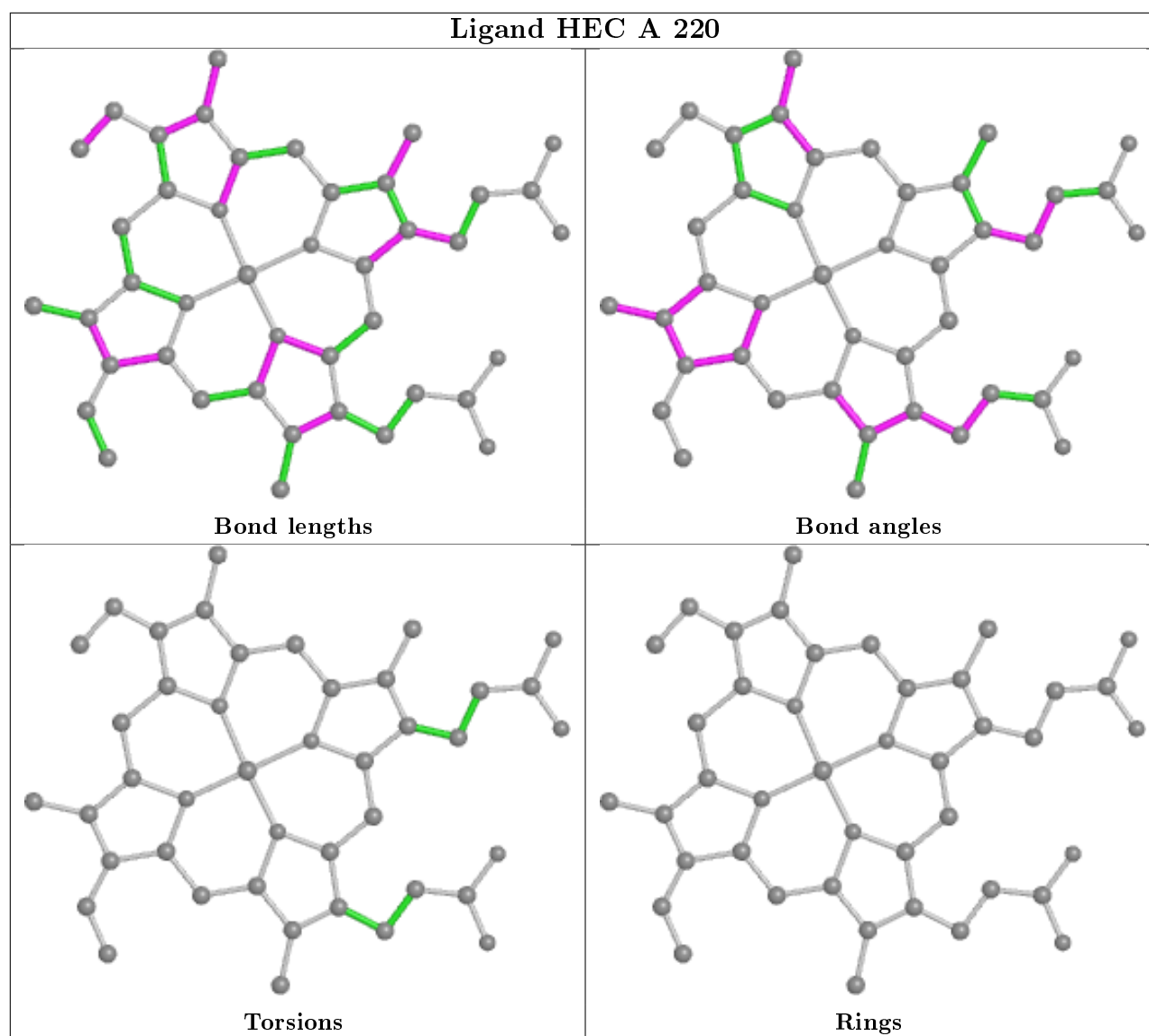


Rings

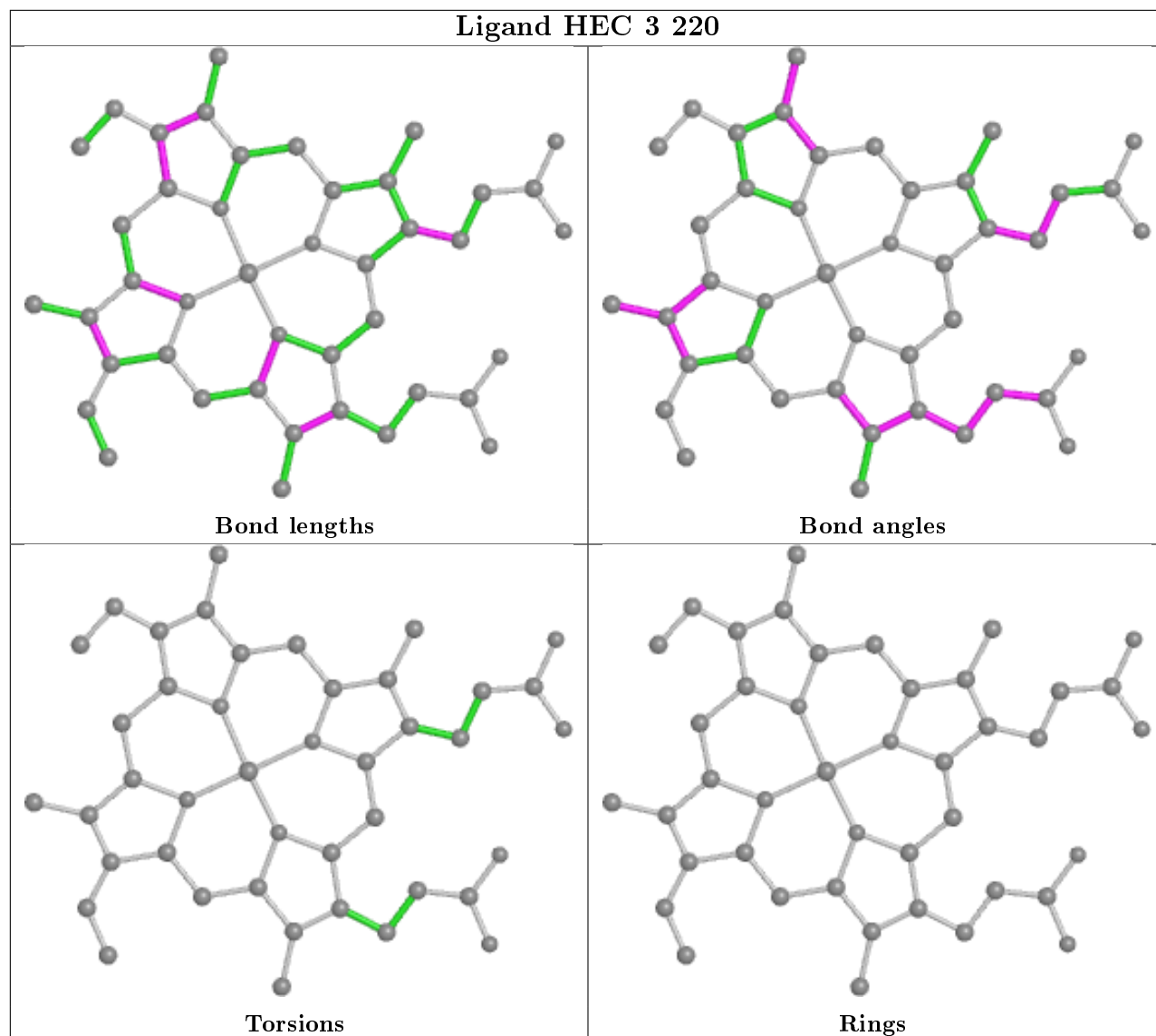
Ligand HEC 6 220



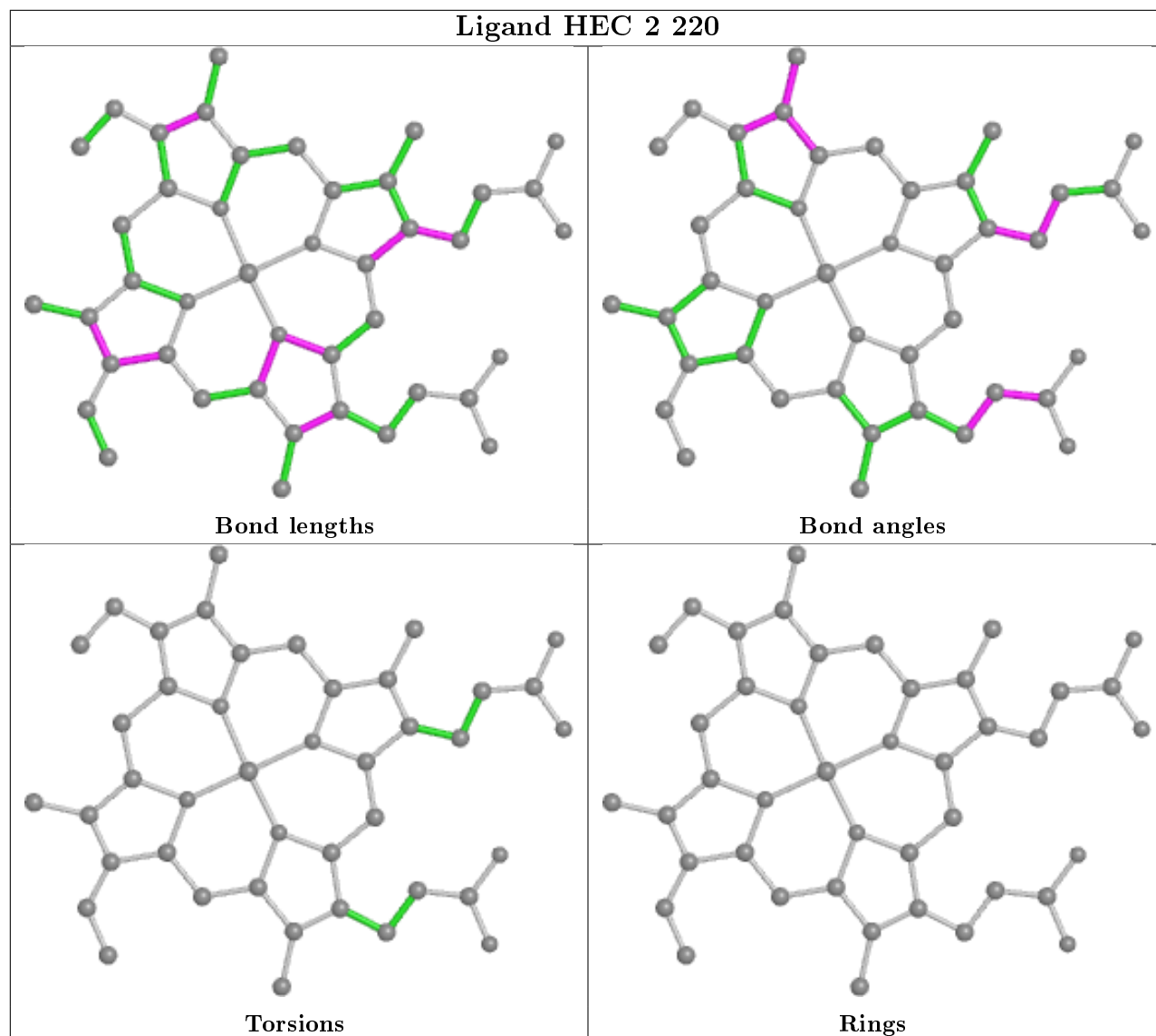


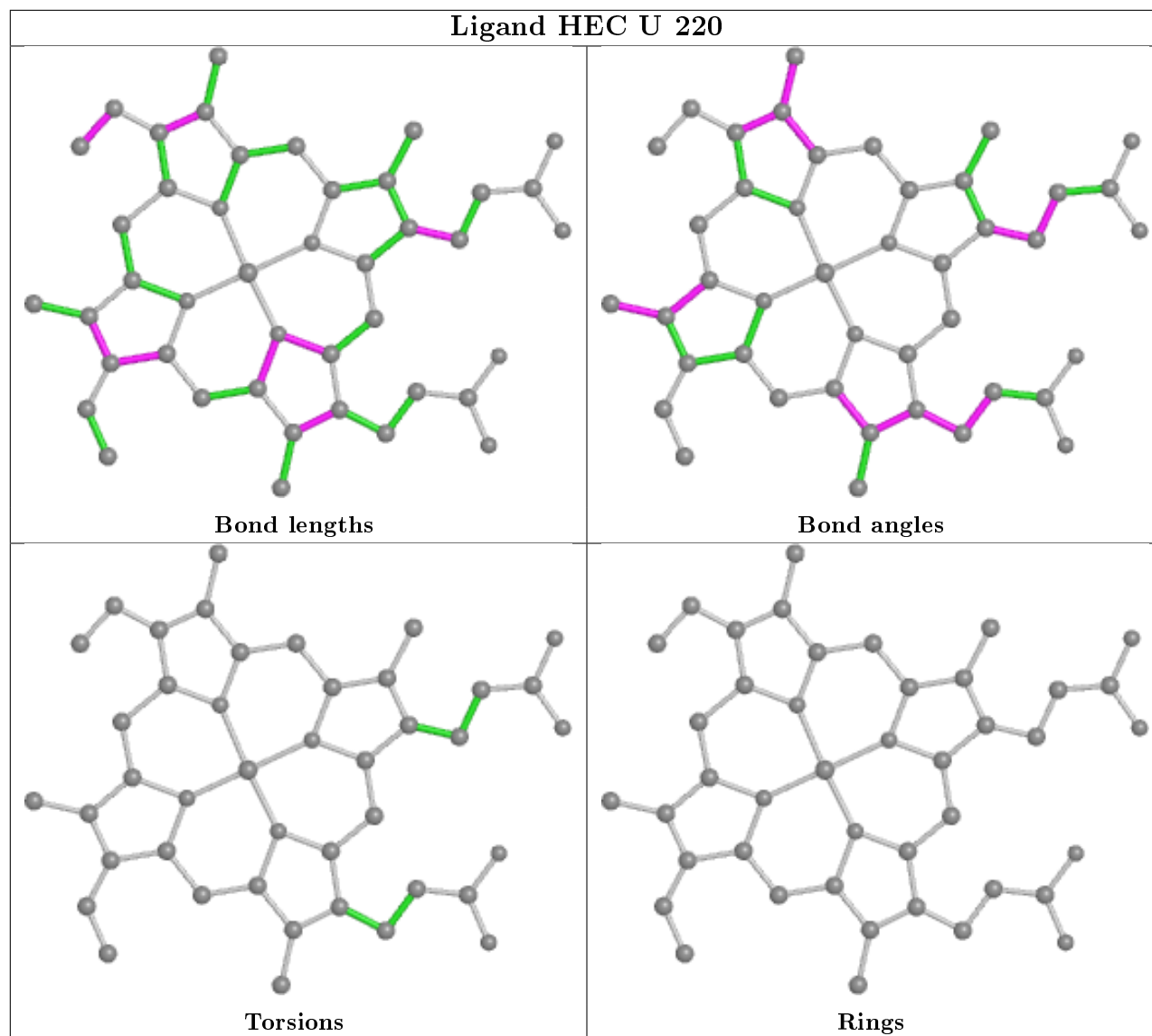


Ligand HEC 3 220

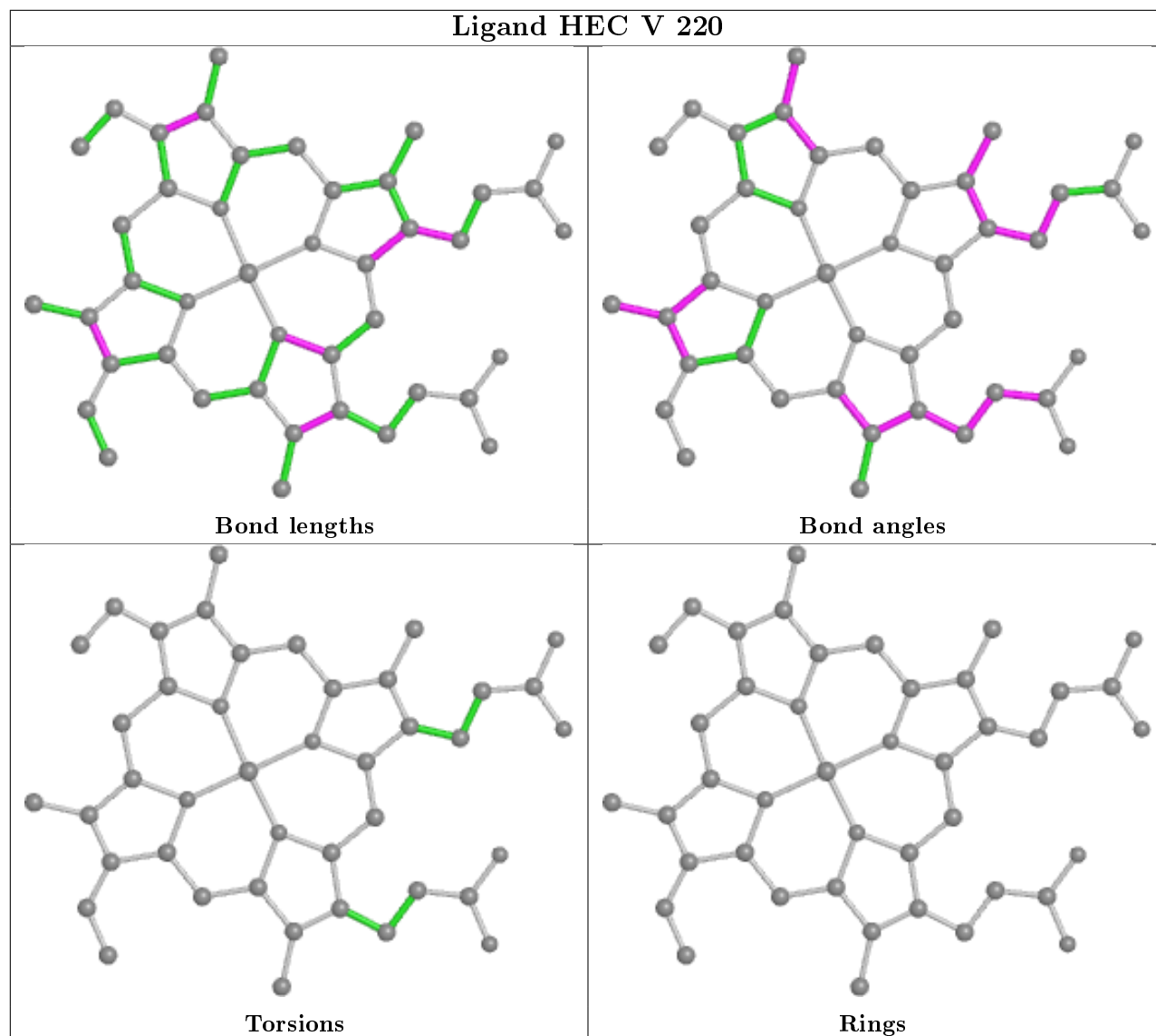


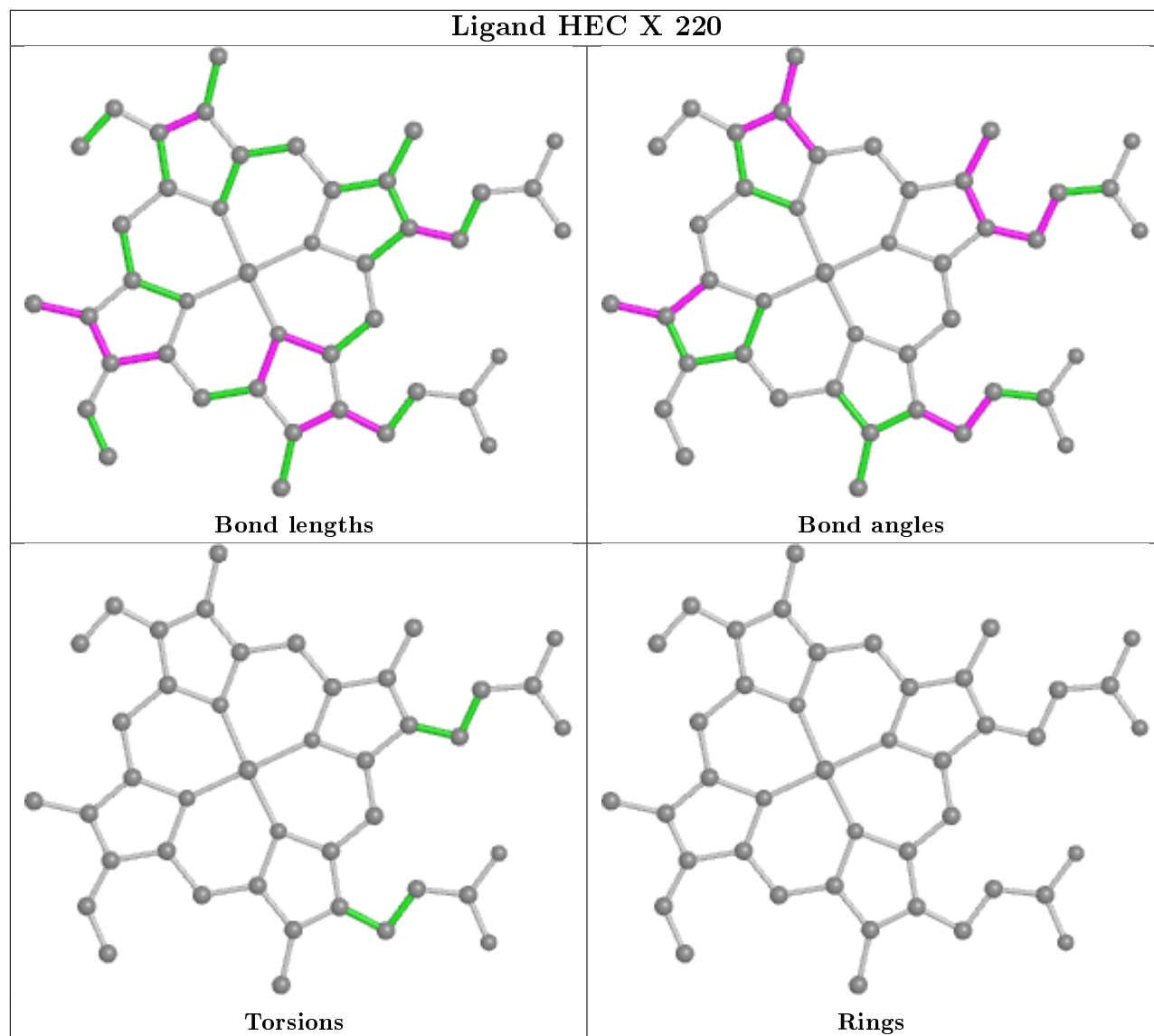
Ligand HEC 2 220



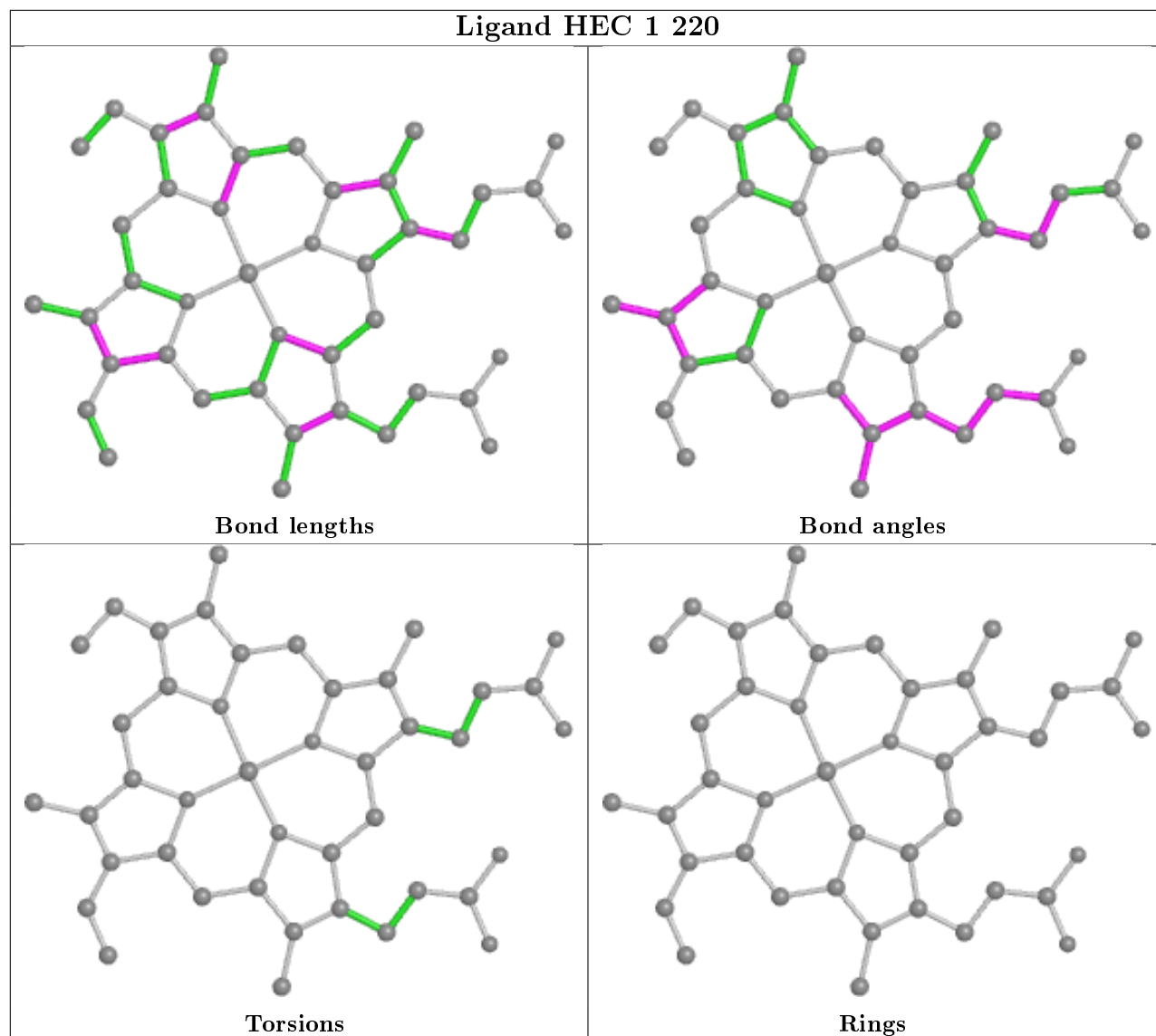


Ligand HEC V 220

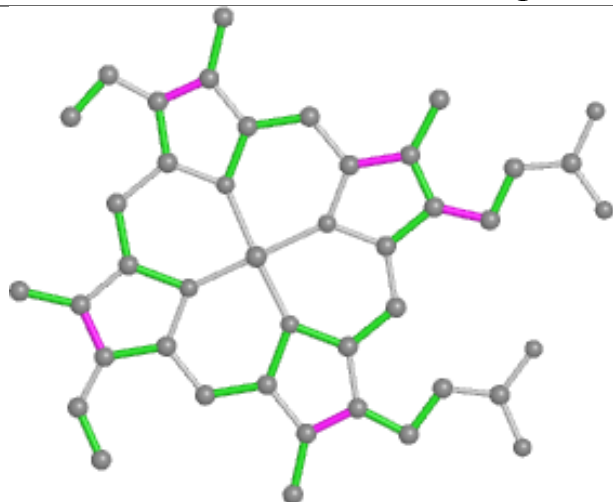




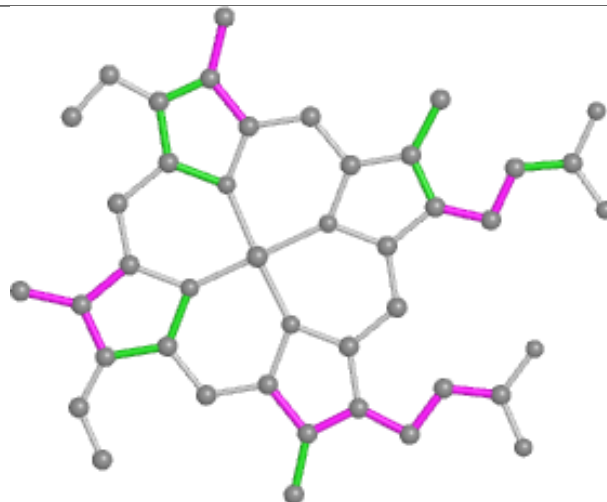
Ligand HEC 1 220



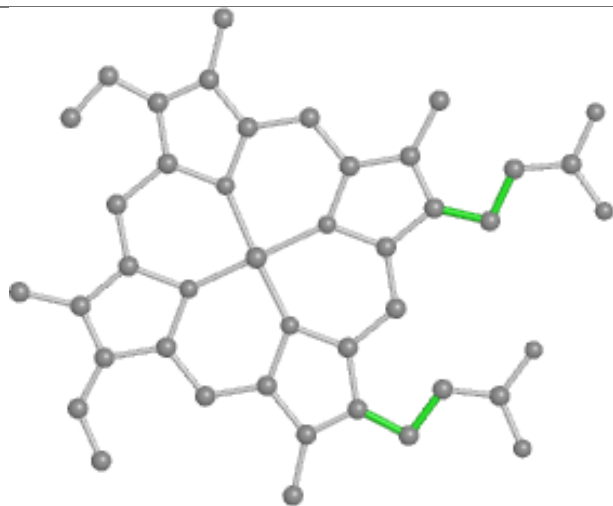
Ligand HEC E 220



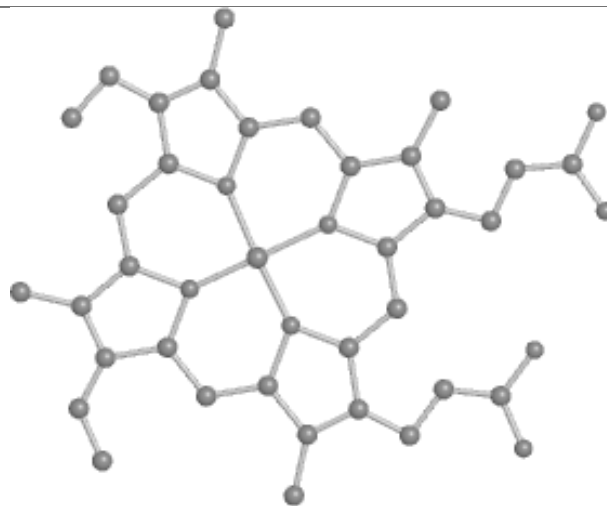
Bond lengths



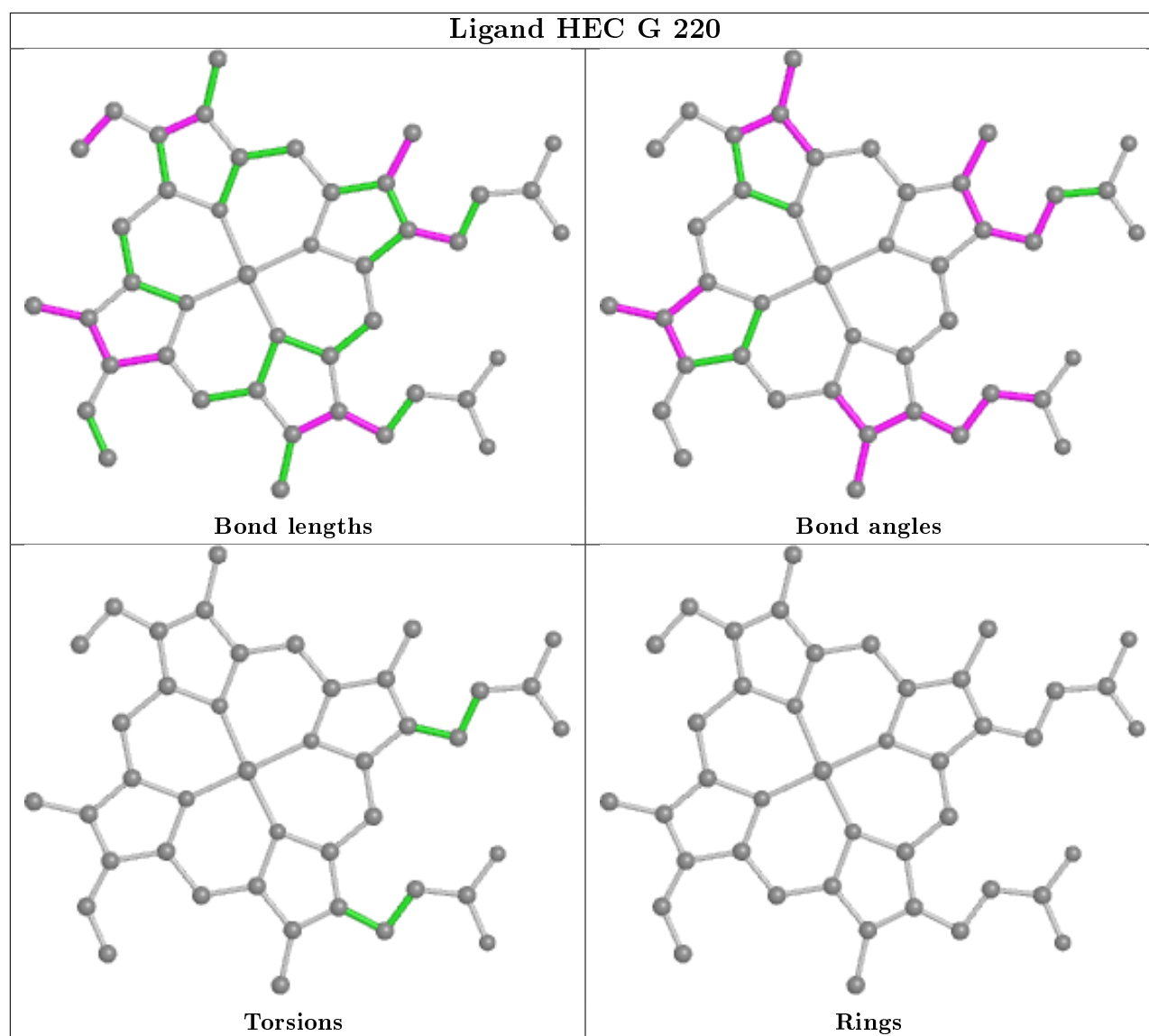
Bond angles

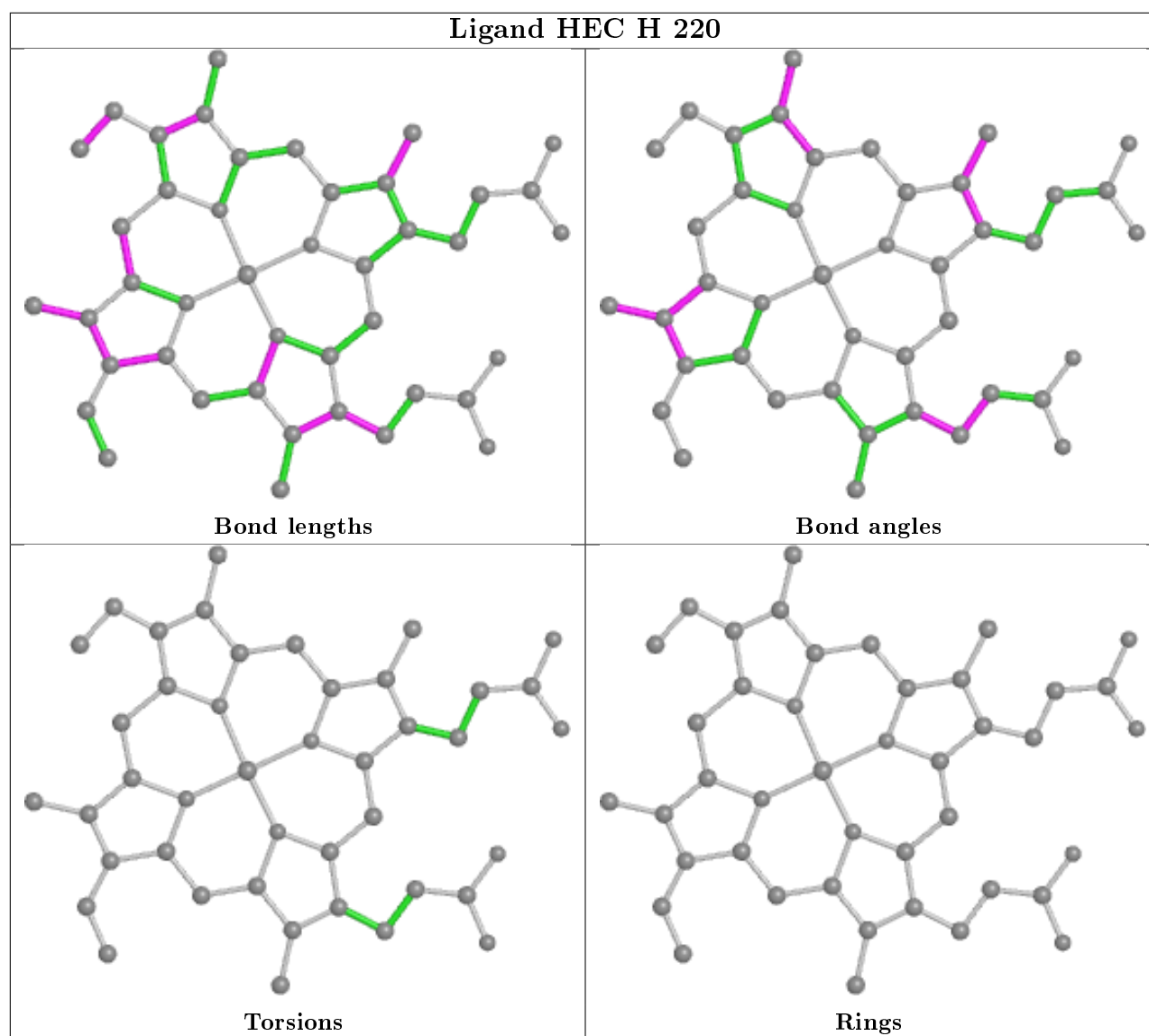


Torsions

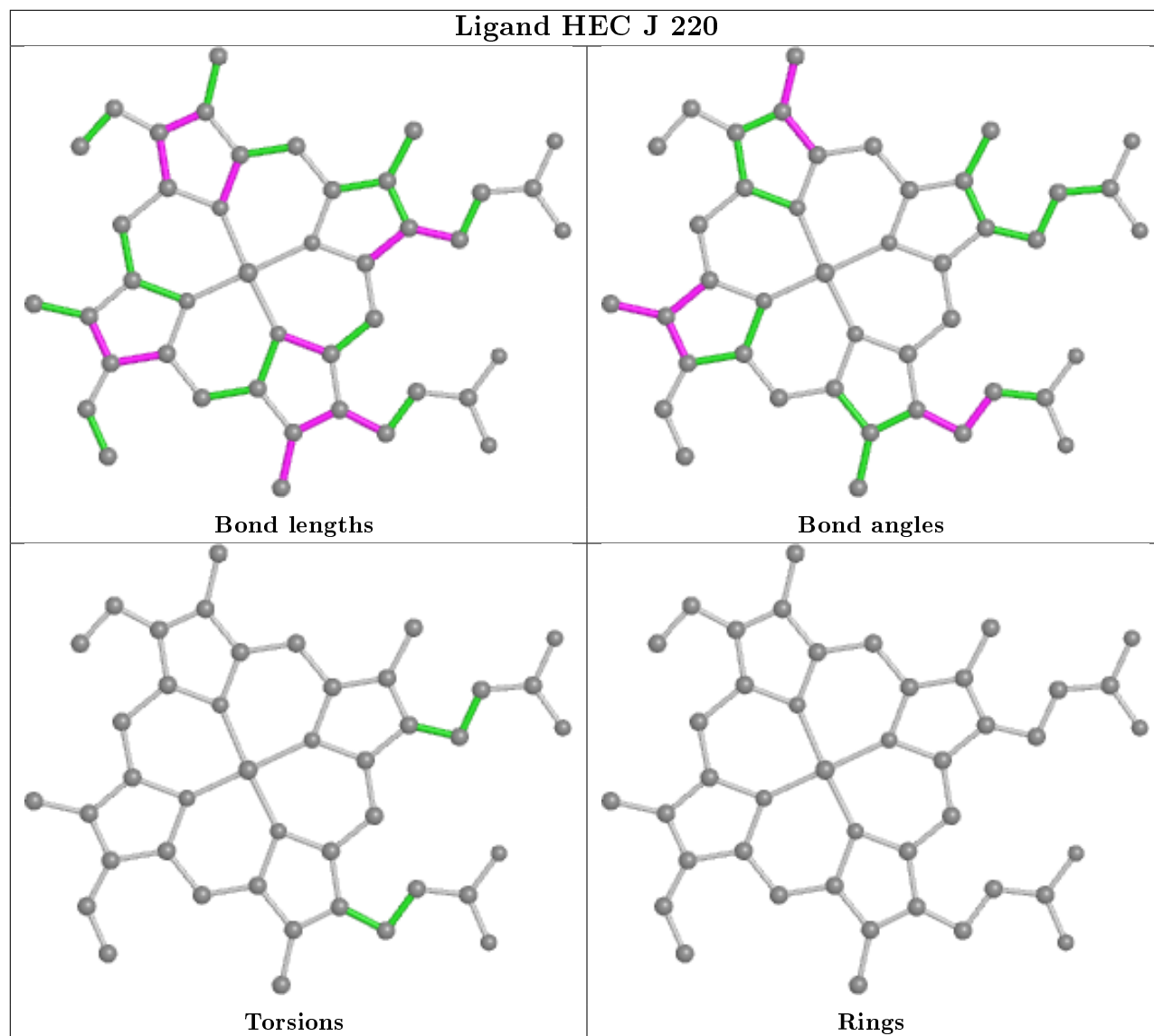


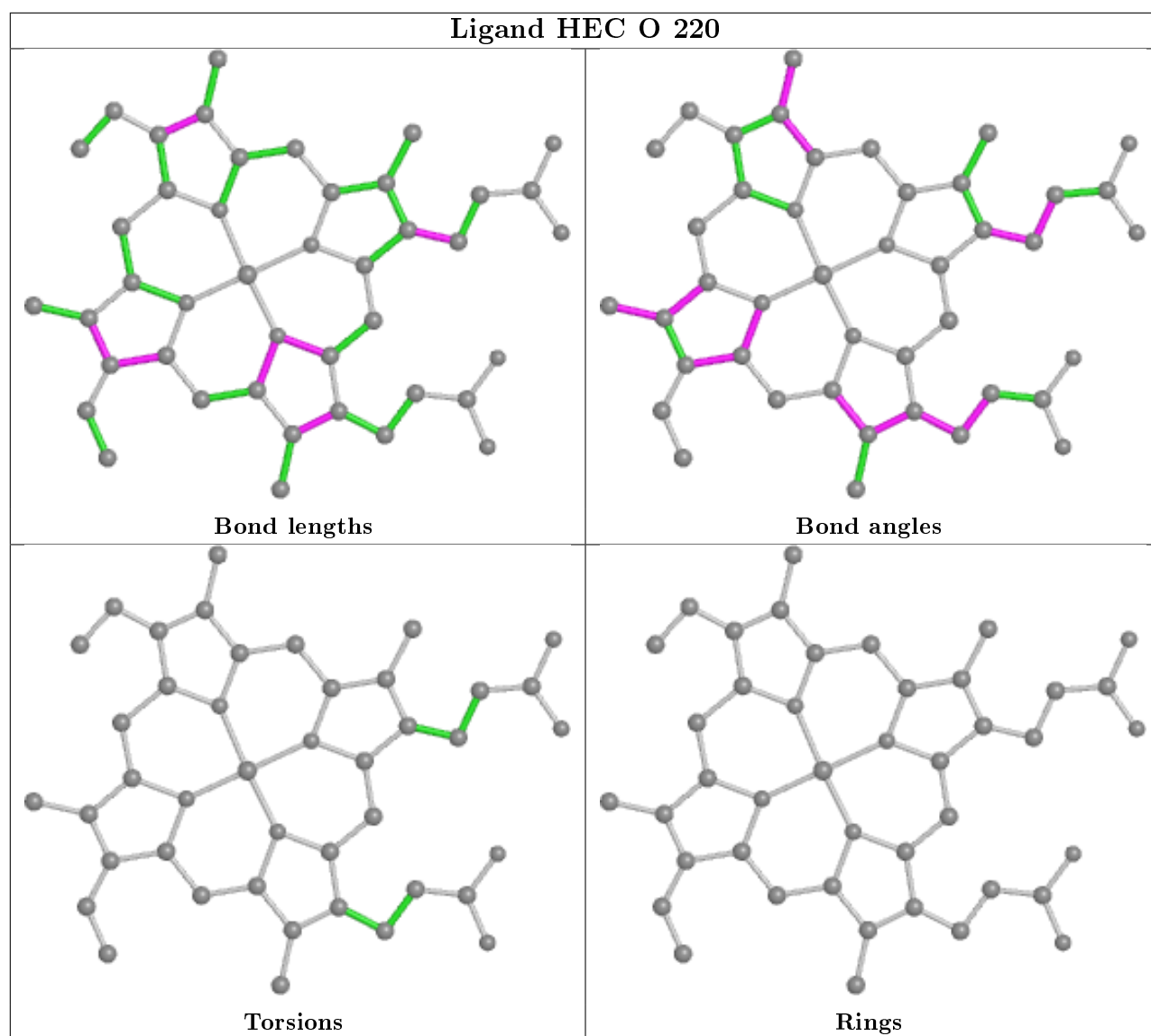
Rings

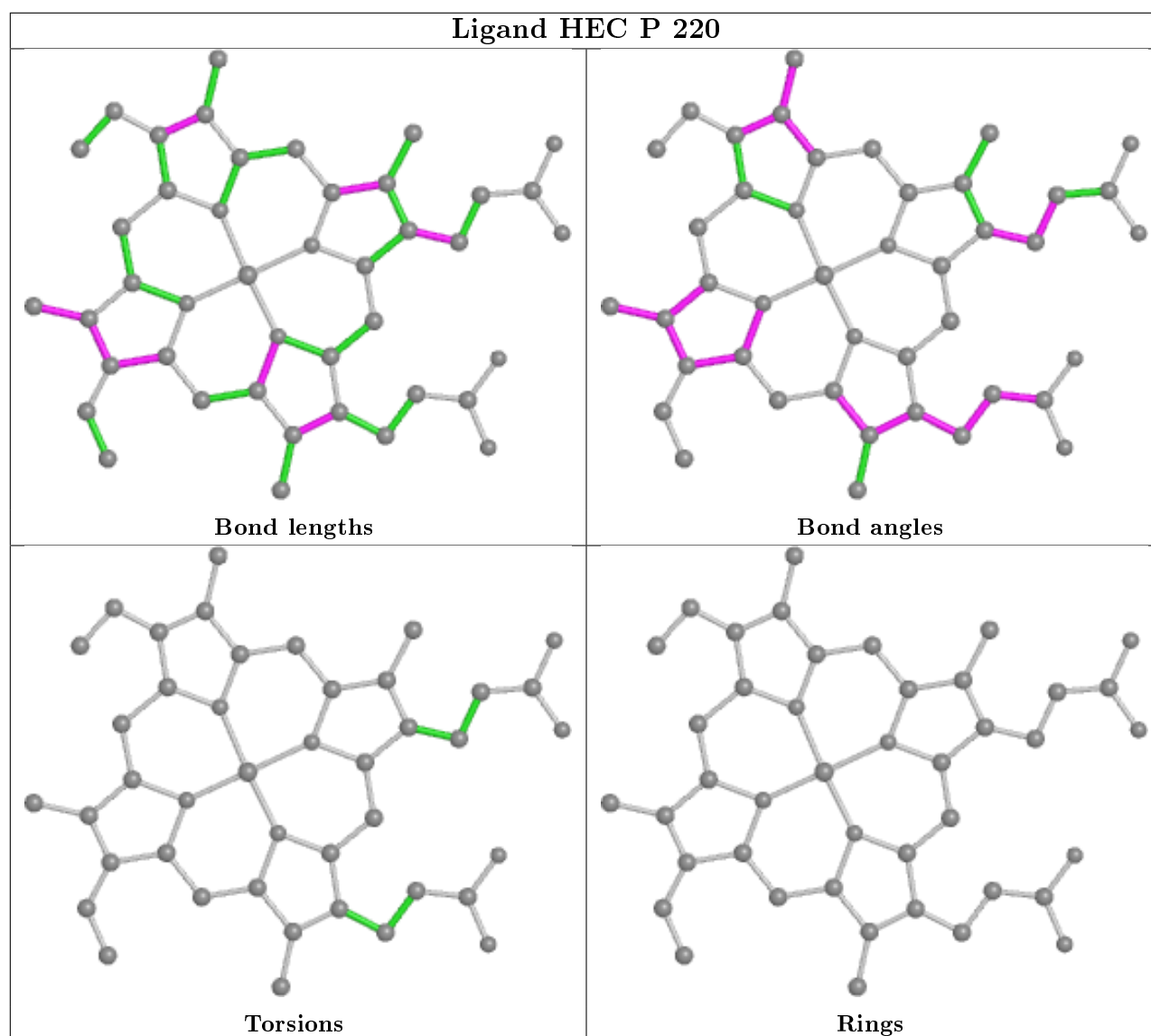


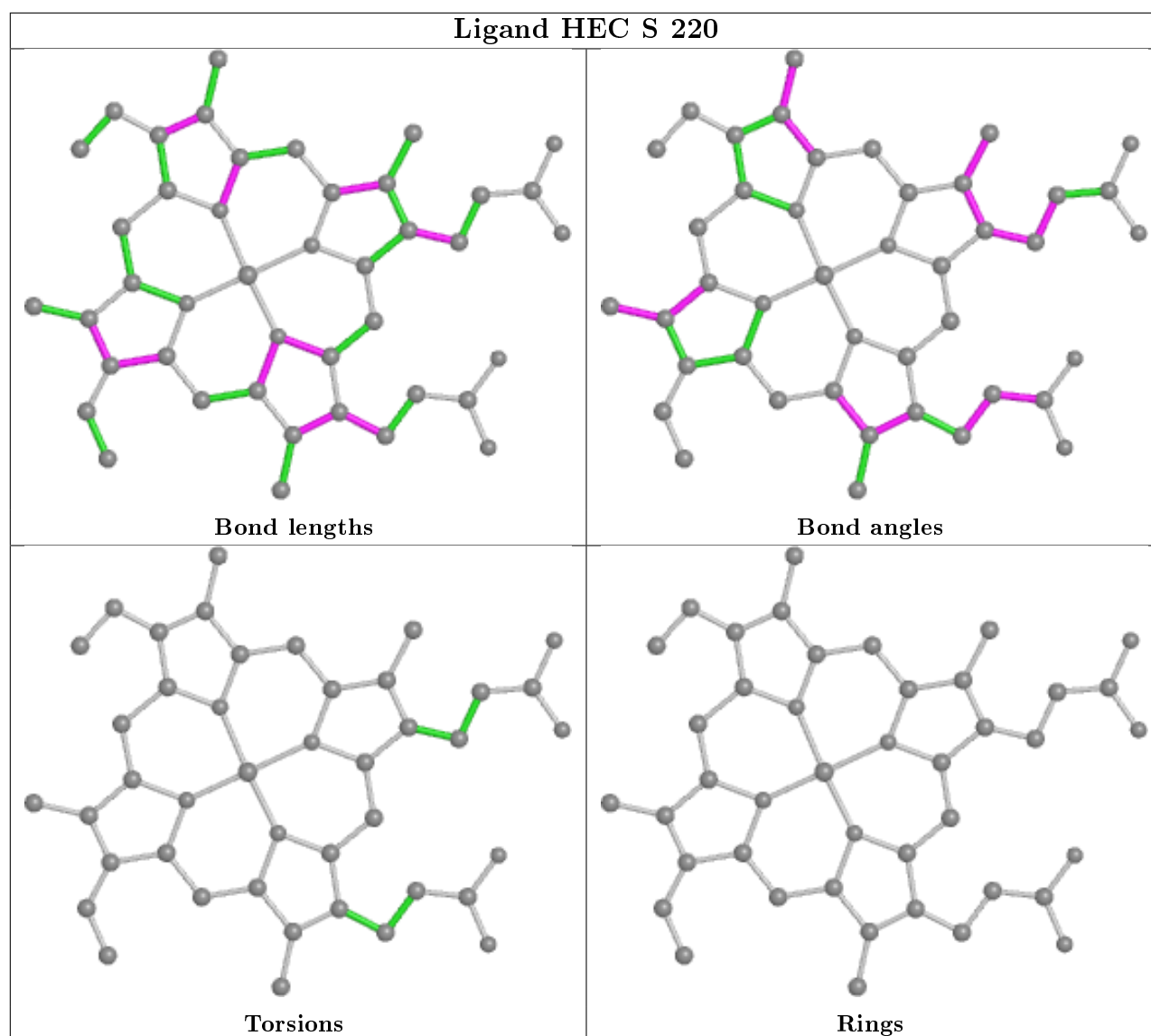


Ligand HEC J 220









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	1	81/103 (78%)	0.52	7 (8%) 10 8	17, 30, 42, 48	0
1	2	81/103 (78%)	1.03	16 (19%) 1 0	21, 38, 54, 57	0
1	3	81/103 (78%)	0.78	11 (13%) 3 2	22, 33, 46, 48	0
1	4	81/103 (78%)	0.67	8 (9%) 7 5	24, 33, 45, 49	0
1	5	81/103 (78%)	1.82	31 (38%) 0 0	25, 37, 58, 63	0
1	6	79/103 (76%)	3.12	57 (72%) 0 0	34, 54, 62, 64	0
1	A	81/103 (78%)	-0.40	0 100 100	10, 17, 24, 30	0
1	B	81/103 (78%)	-0.33	0 100 100	11, 17, 25, 25	0
1	C	81/103 (78%)	-0.26	0 100 100	14, 21, 28, 33	0
1	D	81/103 (78%)	-0.22	0 100 100	14, 22, 31, 34	0
1	E	81/103 (78%)	-0.36	0 100 100	11, 18, 25, 34	0
1	F	81/103 (78%)	-0.20	1 (1%) 79 76	11, 21, 29, 35	0
1	G	81/103 (78%)	-0.24	1 (1%) 79 76	15, 21, 27, 34	0
1	H	81/103 (78%)	-0.27	2 (2%) 57 52	14, 20, 32, 40	0
1	I	81/103 (78%)	-0.20	0 100 100	16, 23, 29, 35	0
1	J	81/103 (78%)	0.52	7 (8%) 10 8	23, 31, 41, 44	0
1	K	81/103 (78%)	0.37	3 (3%) 41 36	18, 28, 42, 47	0
1	L	81/103 (78%)	-0.03	2 (2%) 57 52	16, 22, 35, 40	0
1	M	80/103 (77%)	0.57	6 (7%) 14 11	14, 26, 37, 44	0
1	N	81/103 (78%)	0.36	3 (3%) 41 36	15, 27, 41, 50	0
1	O	81/103 (78%)	-0.24	0 100 100	12, 18, 29, 33	0
1	P	81/103 (78%)	0.06	0 100 100	16, 24, 32, 34	0
1	Q	81/103 (78%)	-0.17	2 (2%) 57 52	15, 20, 33, 37	0
1	R	81/103 (78%)	0.13	4 (4%) 29 24	14, 23, 34, 38	0

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Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	S	81/103 (78%)	0.25	5 (6%) 20 16	15, 26, 36, 39	0
1	T	81/103 (78%)	-0.13	1 (1%) 79 76	13, 21, 33, 36	0
1	U	81/103 (78%)	-0.18	0 100 100	14, 22, 30, 33	0
1	V	81/103 (78%)	-0.23	0 100 100	14, 21, 28, 32	0
1	W	81/103 (78%)	-0.03	1 (1%) 79 76	18, 24, 32, 35	0
1	X	81/103 (78%)	0.12	2 (2%) 57 52	21, 28, 36, 40	0
1	Y	81/103 (78%)	-0.02	1 (1%) 79 76	16, 23, 32, 37	0
1	Z	81/103 (78%)	0.05	4 (4%) 29 24	17, 23, 33, 38	0
All	All	2589/3296 (78%)	0.21	175 (6%) 17 13	10, 24, 45, 64	0

The worst 5 of 175 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	6	87	LEU	8.9
1	6	45	THR	7.4
1	6	77	ALA	7.3
1	5	47	PRO	7.0
1	6	74	GLY	6.8

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates 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
2	HEC	5	220	43/43	0.85	0.24	26,35,43,48	0
2	HEC	6	220	43/43	0.85	0.24	19,38,44,48	0

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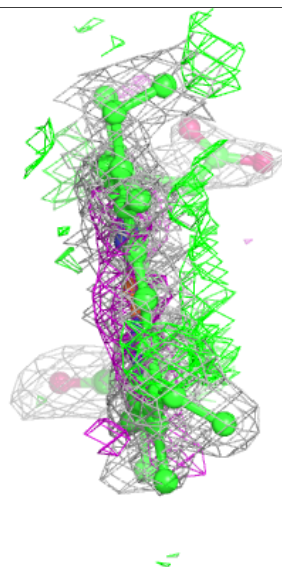
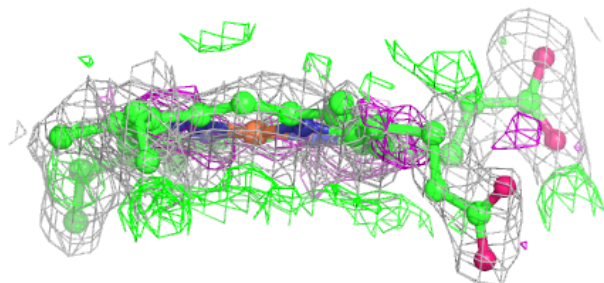
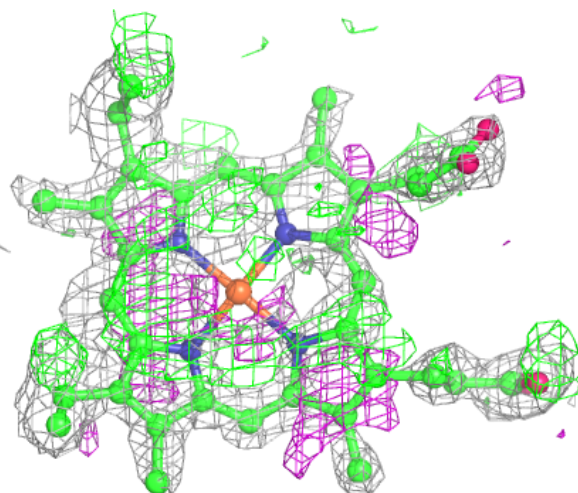
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	P	104	6/6	0.90	0.17	24,32,34,37	0
3	GOL	E	104	6/6	0.92	0.15	18,26,29,34	0
3	GOL	C	104	6/6	0.92	0.10	26,31,33,34	0
3	GOL	R	104	6/6	0.93	0.17	32,36,37,41	0
2	HEC	2	220	43/43	0.93	0.13	17,31,35,36	0
2	HEC	N	220	43/43	0.95	0.12	14,23,28,33	0
2	HEC	3	220	43/43	0.95	0.11	21,29,32,34	0
2	HEC	4	220	43/43	0.95	0.12	19,25,30,31	0
2	HEC	J	220	43/43	0.96	0.13	11,21,24,32	0
2	HEC	Z	220	43/43	0.97	0.11	10,15,20,25	0
2	HEC	X	220	43/43	0.97	0.11	16,21,25,27	0
2	HEC	Y	220	43/43	0.97	0.10	12,16,20,22	0
2	HEC	S	220	43/43	0.97	0.11	12,20,23,26	0
2	HEC	L	220	43/43	0.98	0.11	11,16,19,23	0
2	HEC	R	220	43/43	0.98	0.12	9,14,18,20	0
2	HEC	W	220	43/43	0.98	0.12	14,19,23,24	0
2	HEC	I	220	43/43	0.98	0.11	13,16,22,24	0
2	HEC	Q	220	43/43	0.98	0.11	10,14,18,22	0
2	HEC	A	220	43/43	0.98	0.12	4,11,16,17	0
2	HEC	K	220	43/43	0.98	0.13	12,16,19,21	0
2	HEC	M	220	43/43	0.98	0.11	11,15,18,24	0
2	HEC	U	220	43/43	0.98	0.13	10,14,19,22	0
2	HEC	V	220	43/43	0.98	0.12	10,14,18,19	0
2	HEC	C	220	43/43	0.98	0.12	9,13,18,20	0
2	HEC	1	220	43/43	0.98	0.12	12,19,24,28	0
2	HEC	G	220	43/43	0.98	0.12	12,16,19,22	0
2	HEC	H	220	43/43	0.98	0.11	11,14,20,21	0
2	HEC	T	220	43/43	0.98	0.13	8,14,18,22	0
2	HEC	O	220	43/43	0.98	0.11	7,12,19,20	0
2	HEC	P	220	43/43	0.98	0.11	10,15,19,21	0
2	HEC	D	220	43/43	0.98	0.12	11,15,18,19	0
2	HEC	B	220	43/43	0.99	0.12	8,11,15,17	0
2	HEC	F	220	43/43	0.99	0.13	7,12,17,19	0
2	HEC	E	220	43/43	0.99	0.11	7,11,16,19	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.

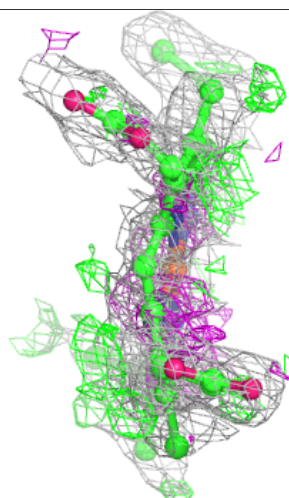
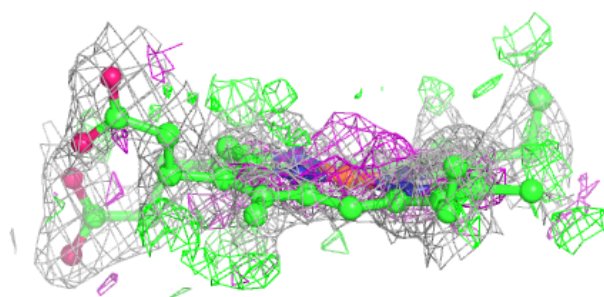
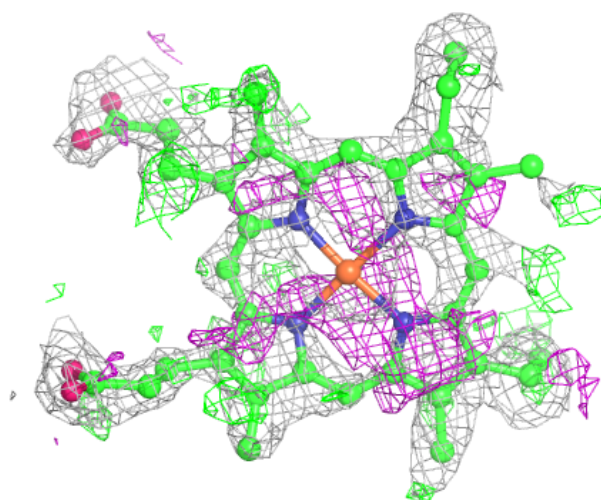
Electron density around HEC 5 220:

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



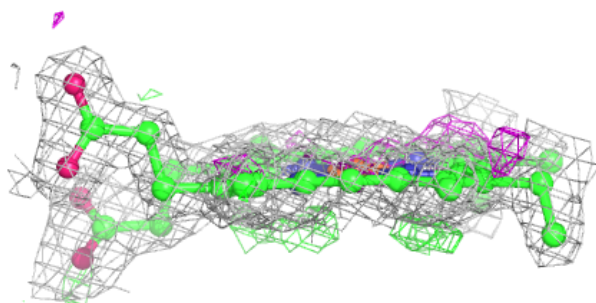
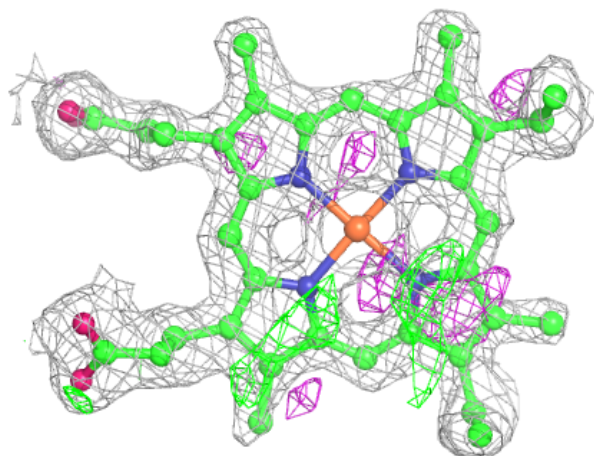
Electron density around HEC 6 220:

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



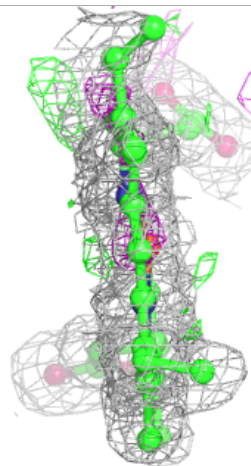
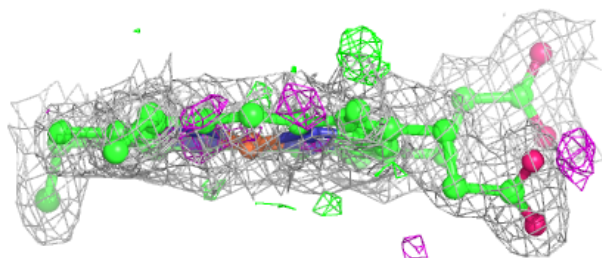
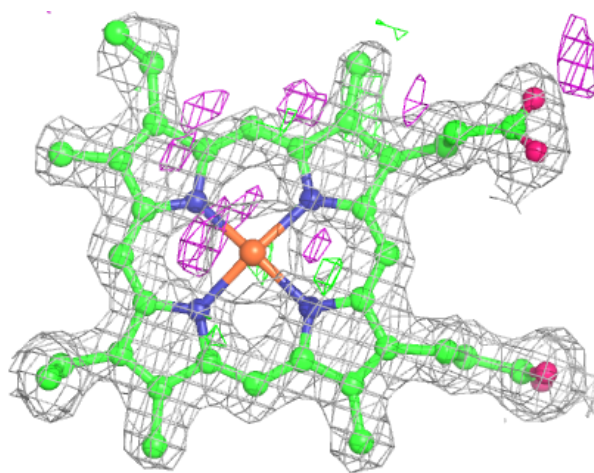
Electron density around HEC 2 220:

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and green (positive)



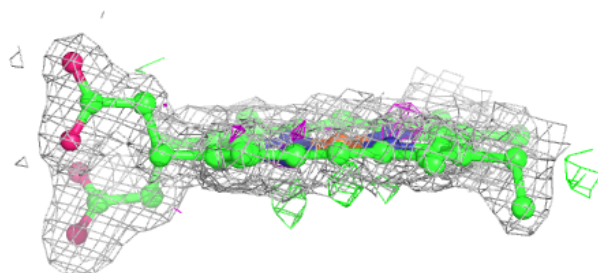
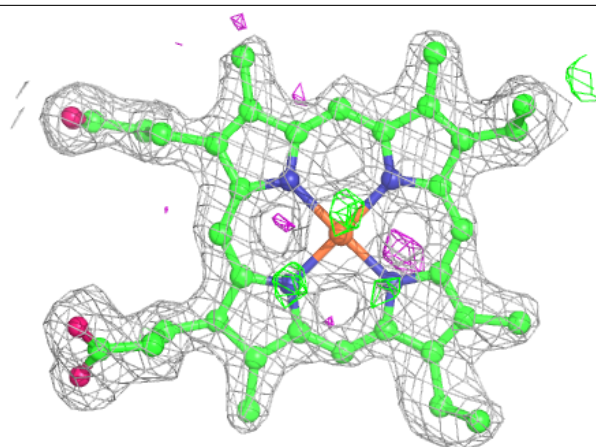
Electron density around HEC N 220:

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



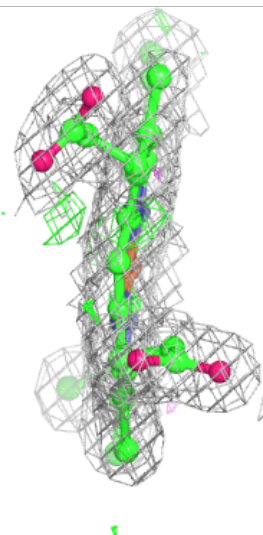
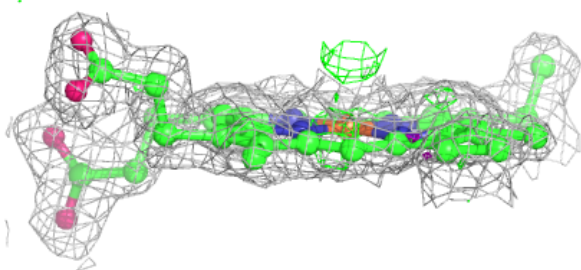
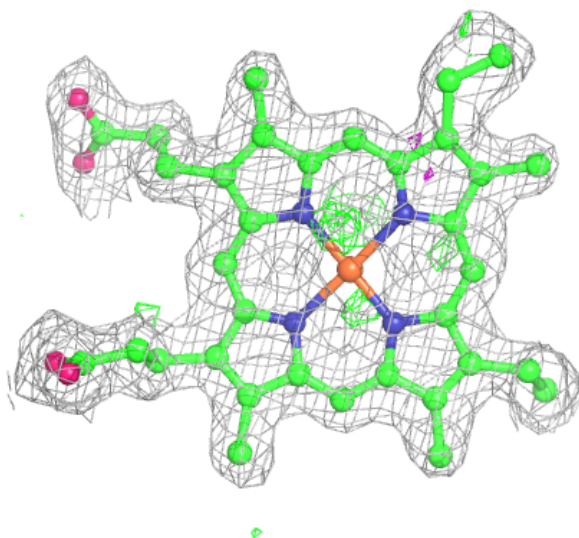
Electron density around HEC 3 220:

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



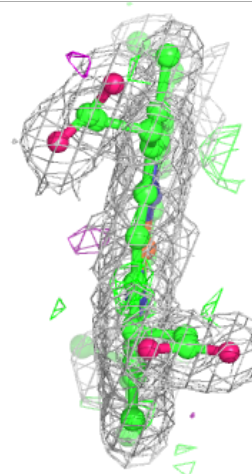
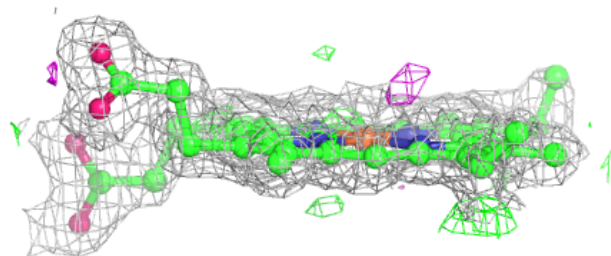
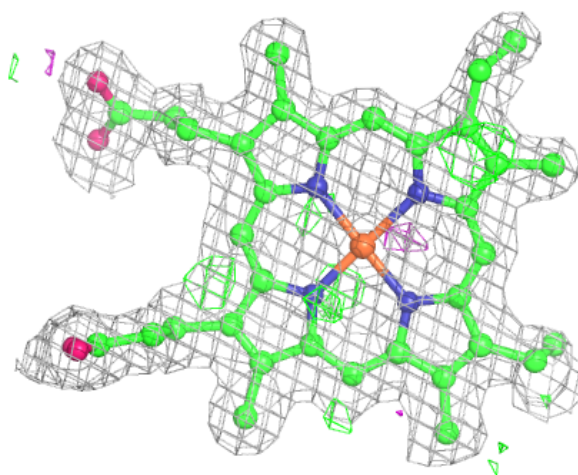
Electron density around HEC 4 220:

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



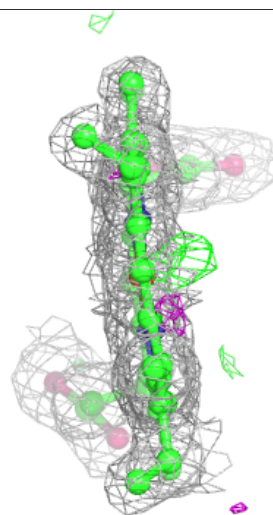
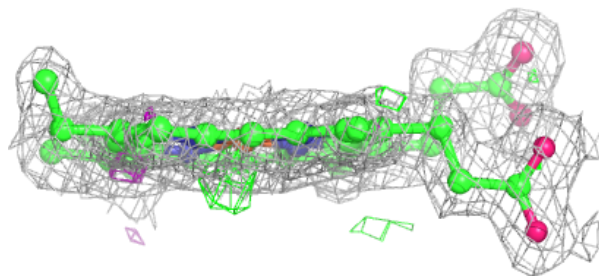
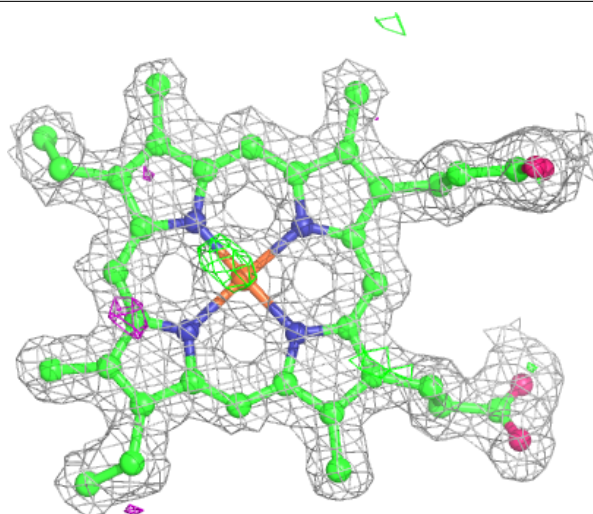
Electron density around HEC J 220:

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



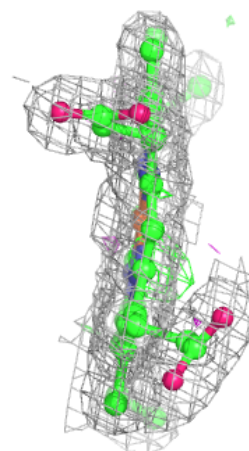
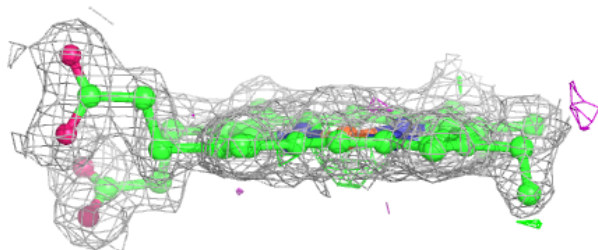
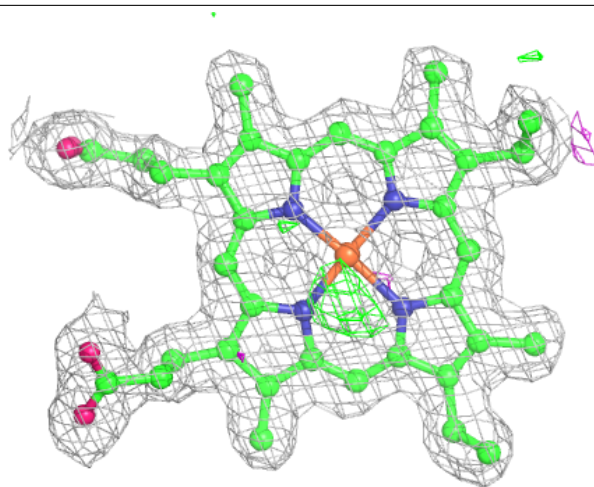
Electron density around HEC Z 220:

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



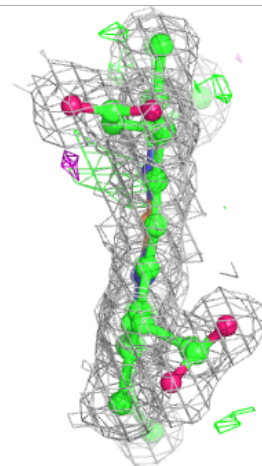
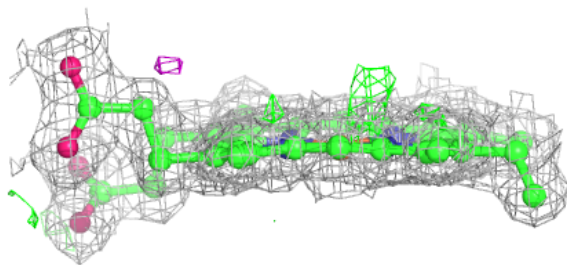
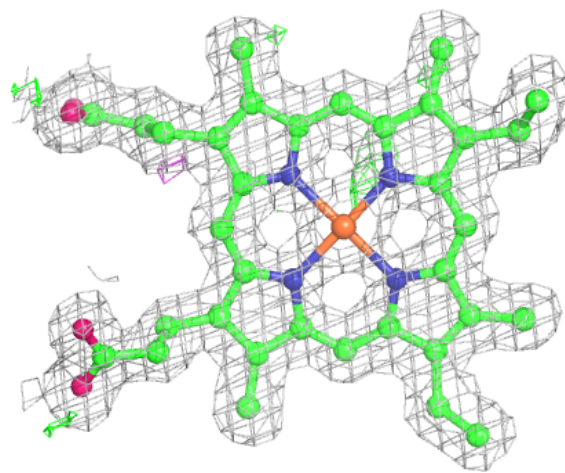
Electron density around HEC X 220:

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



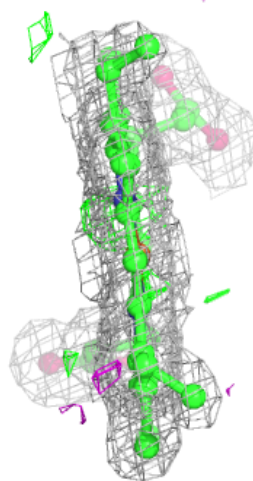
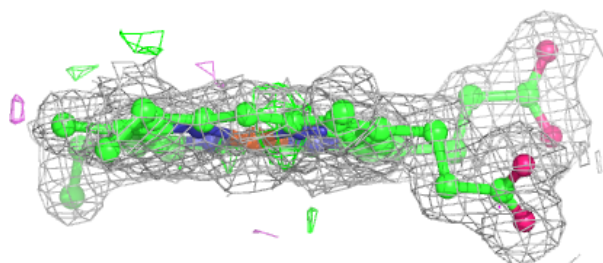
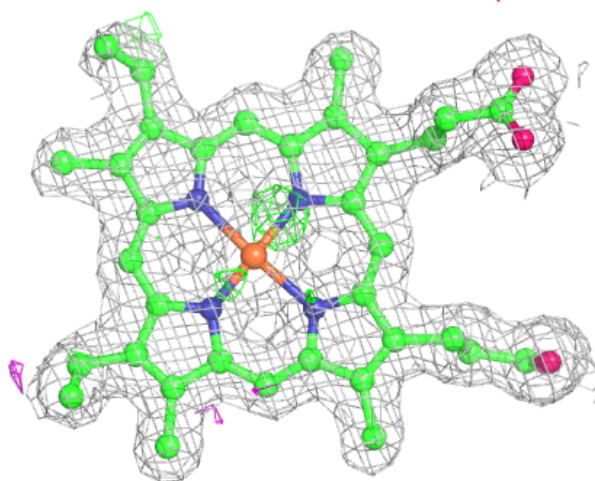
Electron density around HEC Y 220:

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



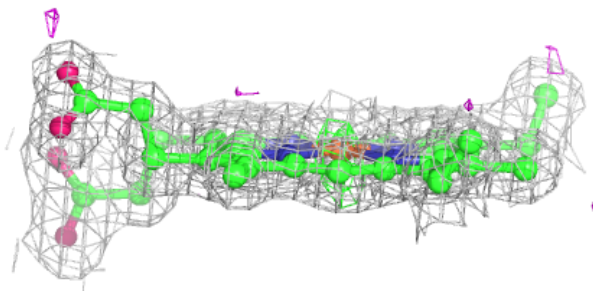
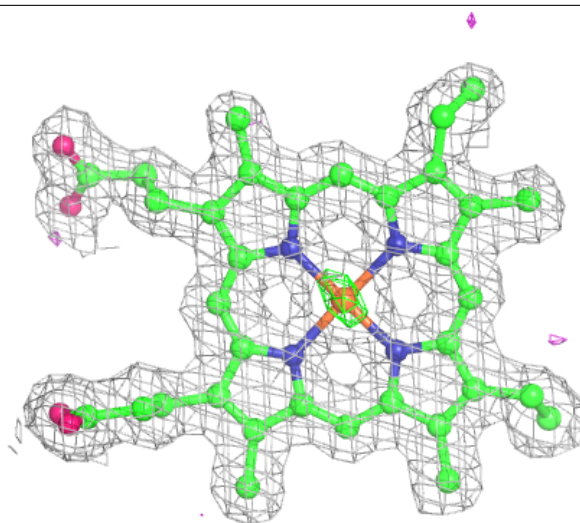
Electron density around HEC S 220:

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



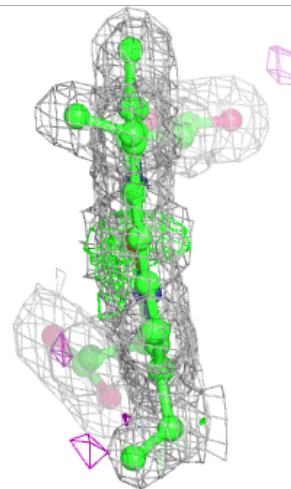
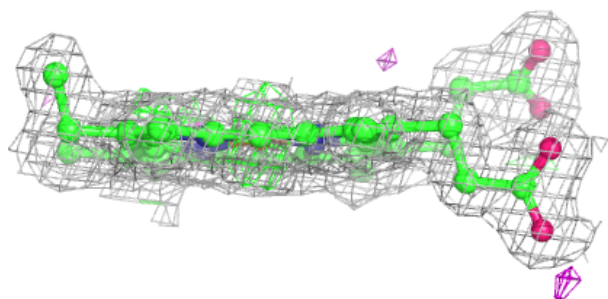
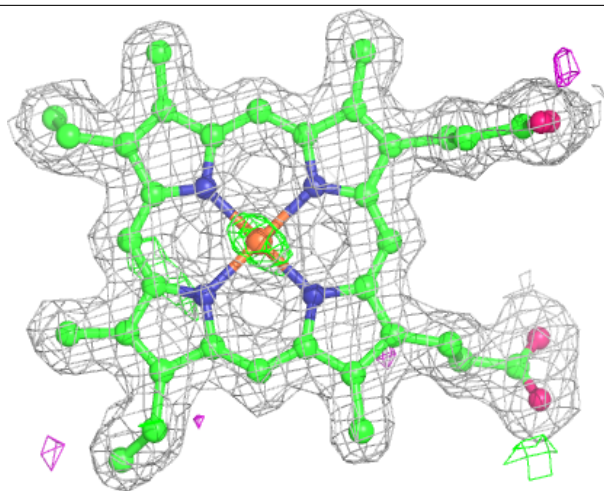
Electron density around HEC L 220:

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



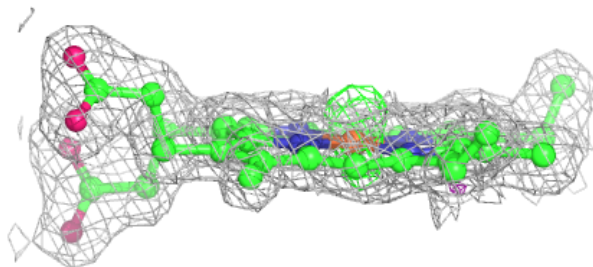
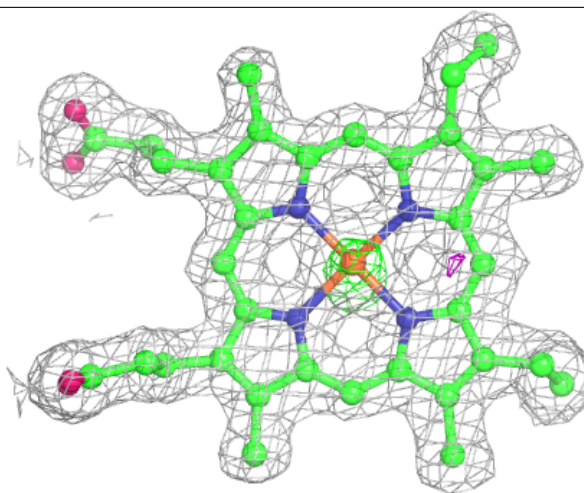
Electron density around HEC R 220:

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



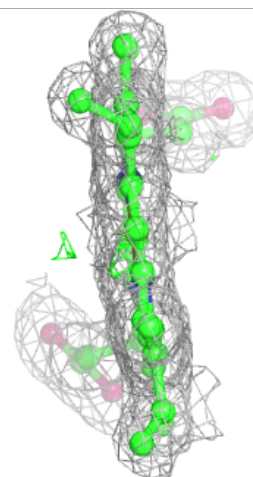
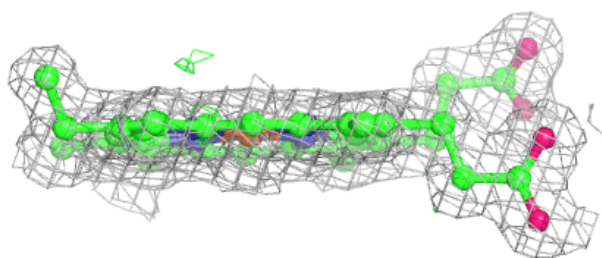
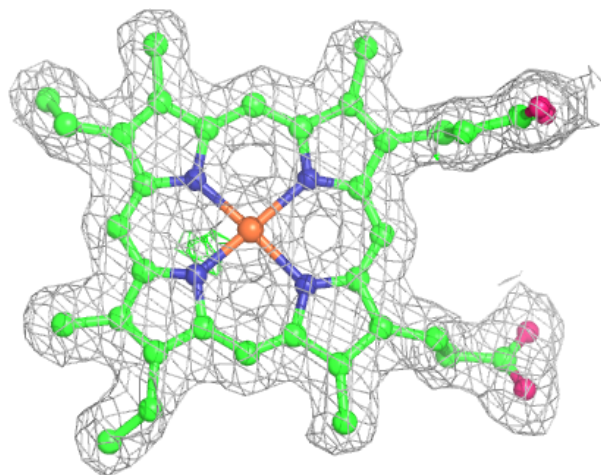
Electron density around HEC W 220:

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



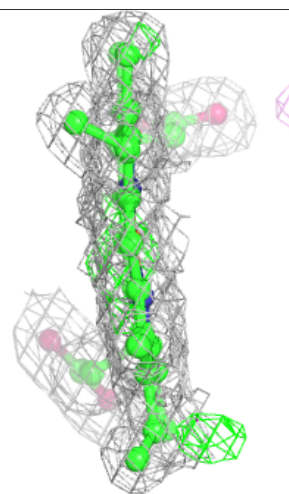
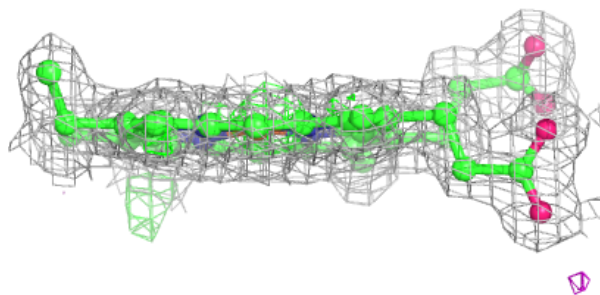
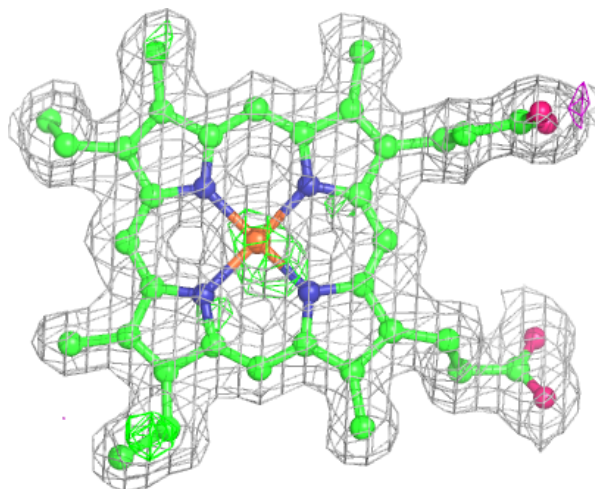
Electron density around HEC I 220:

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



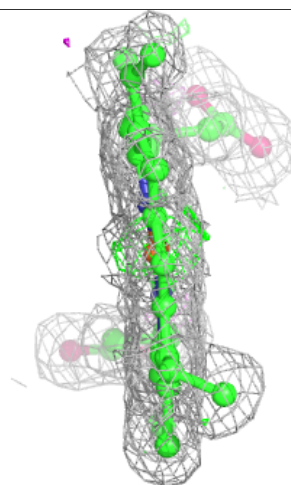
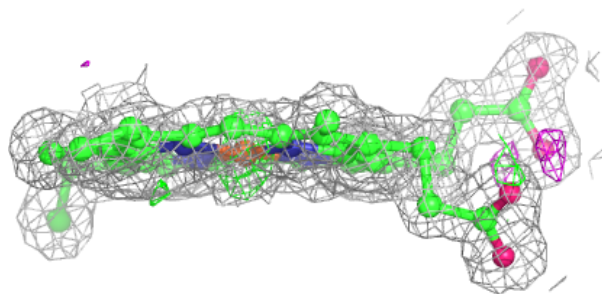
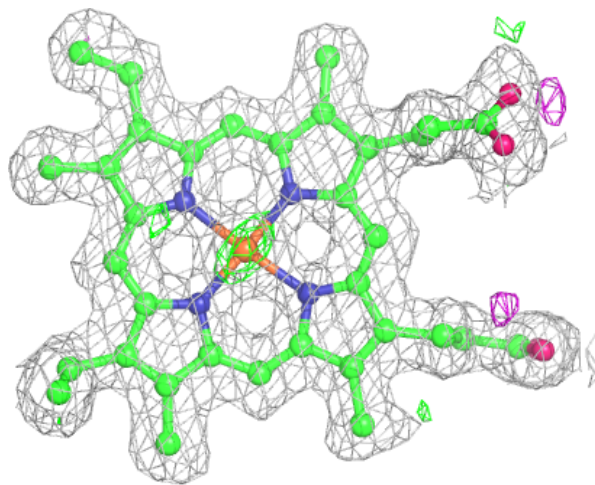
Electron density around HEC Q 220:

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



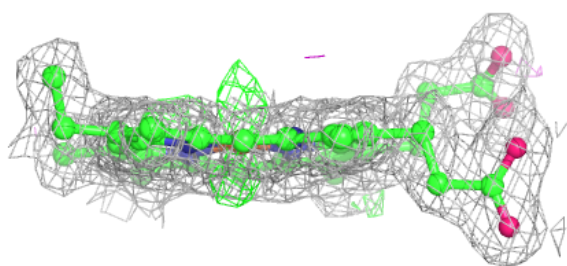
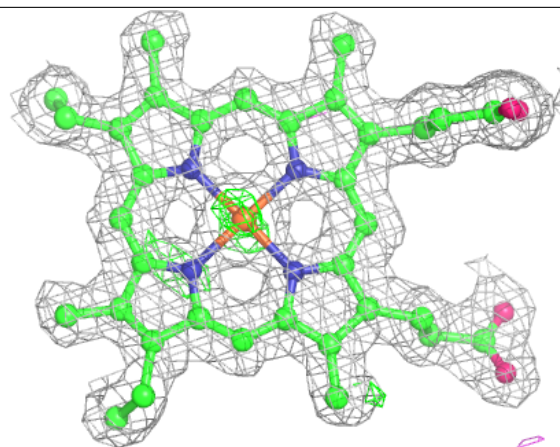
Electron density around HEC A 220:

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



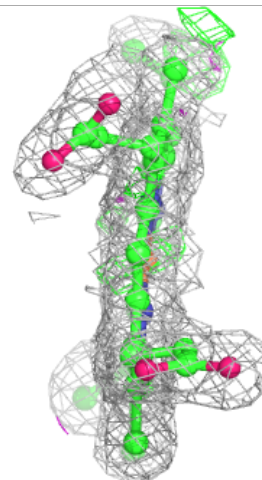
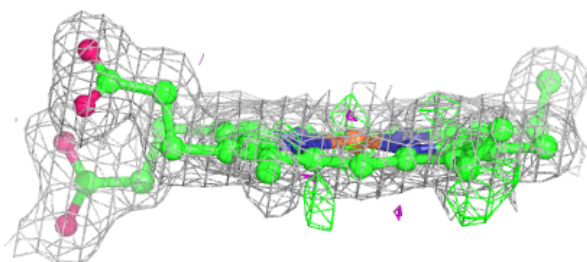
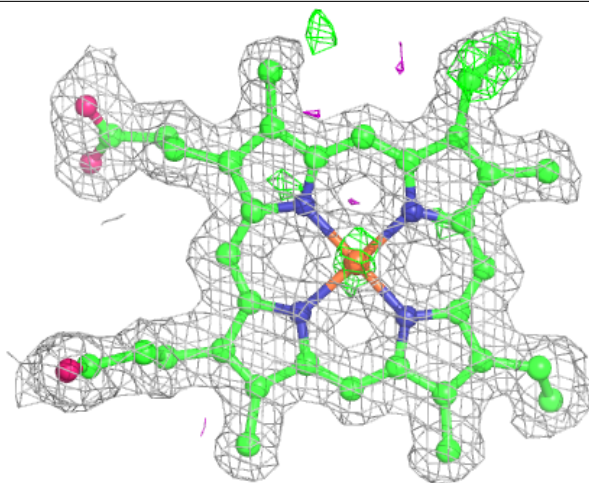
Electron density around HEC K 220:

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and green (positive)



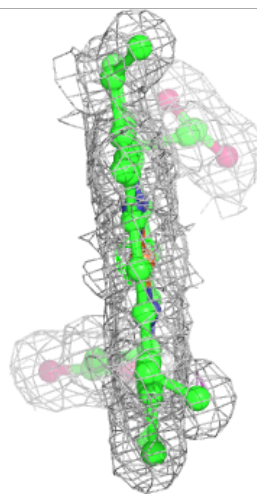
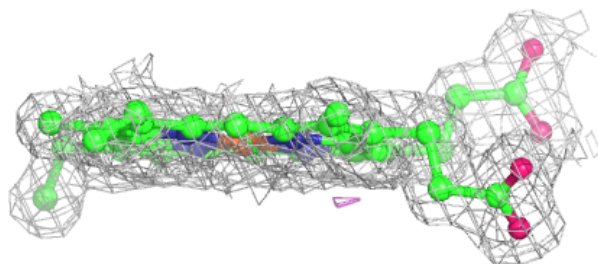
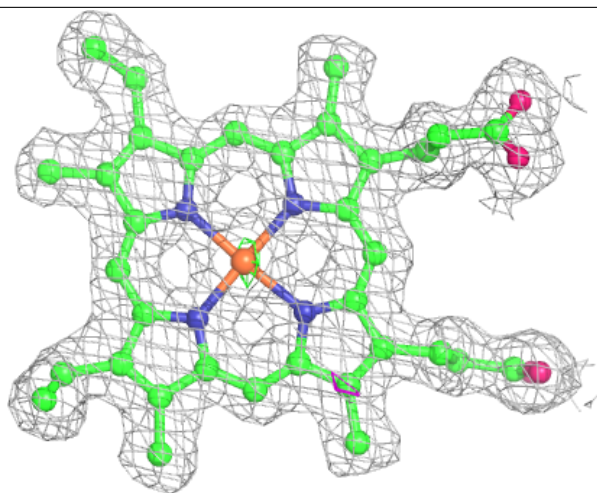
Electron density around HEC M 220:

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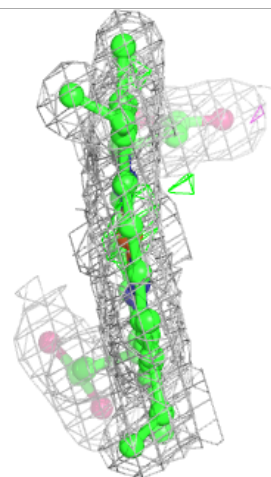
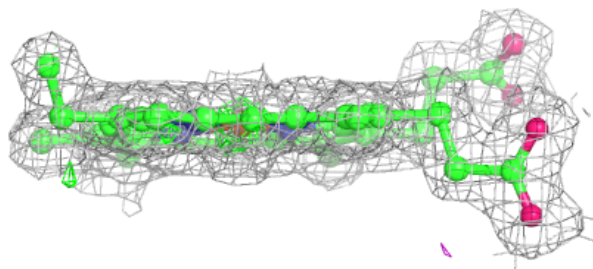
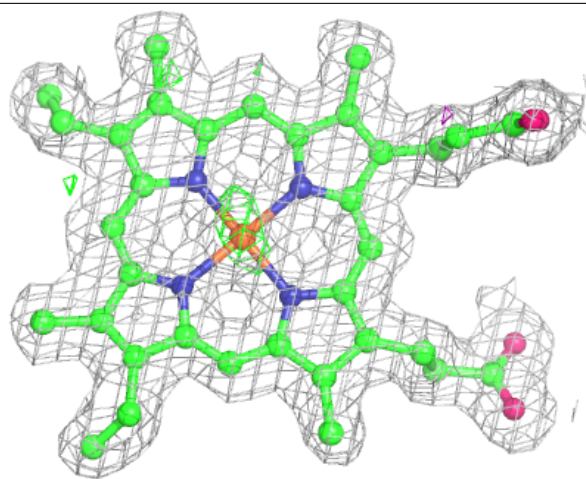
Electron density around HEC U 220:

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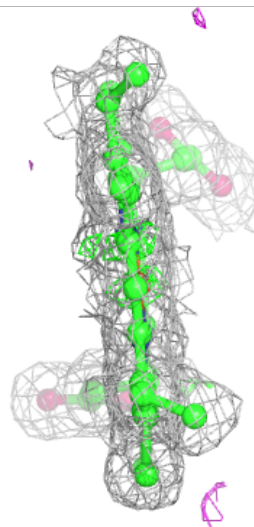
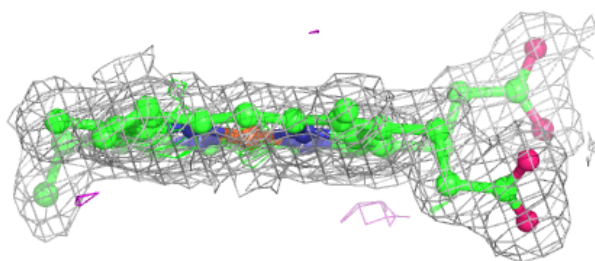
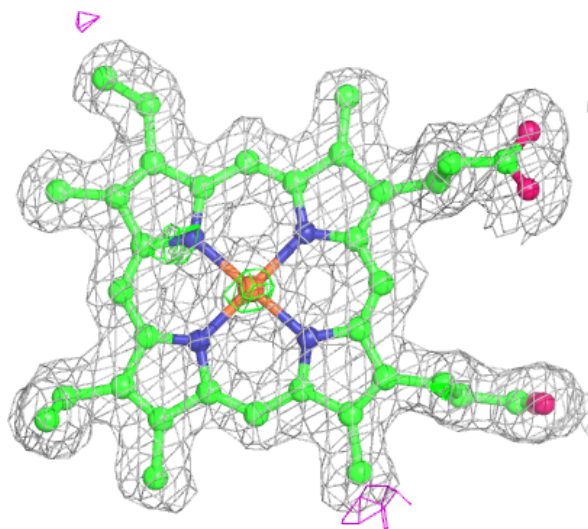
Electron density around HEC V 220:

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and green (positive)



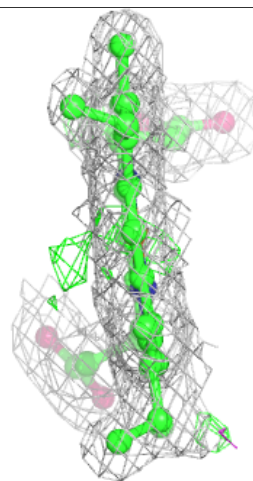
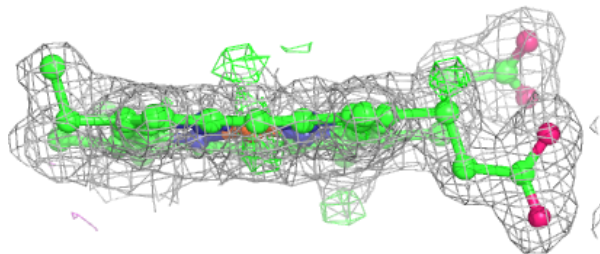
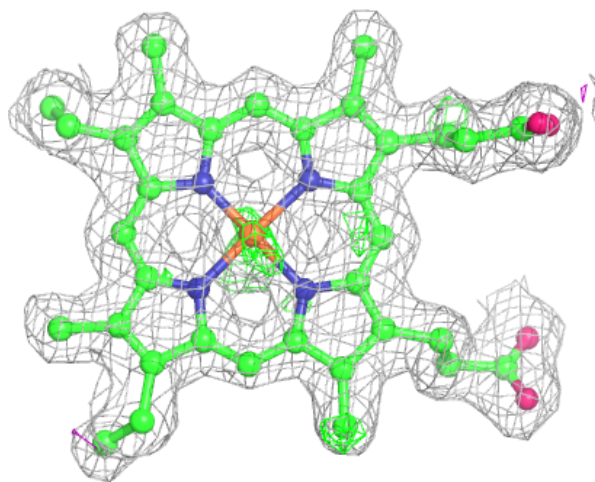
Electron density around HEC C 220:

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



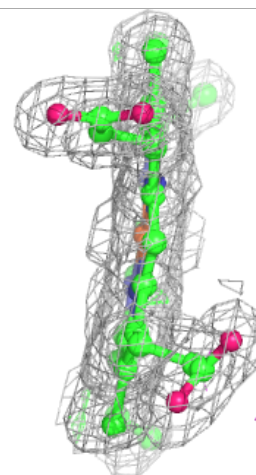
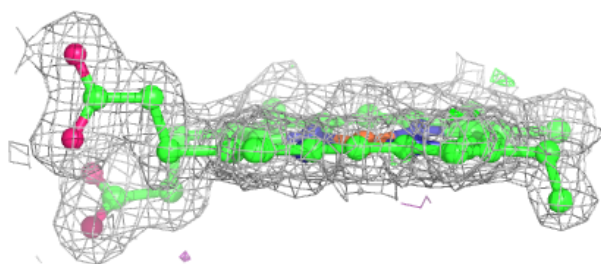
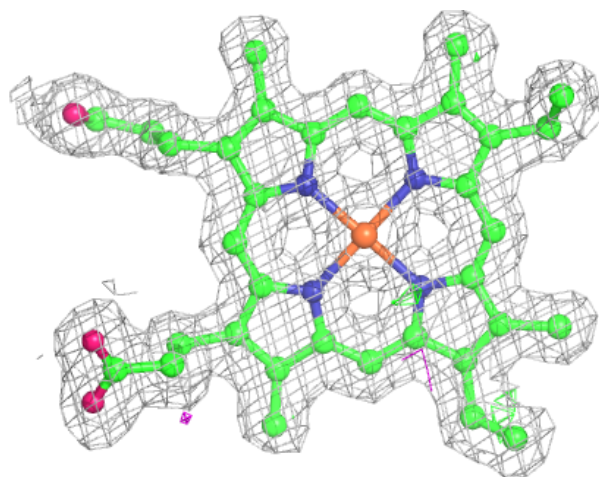
Electron density around HEC 1 220:

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



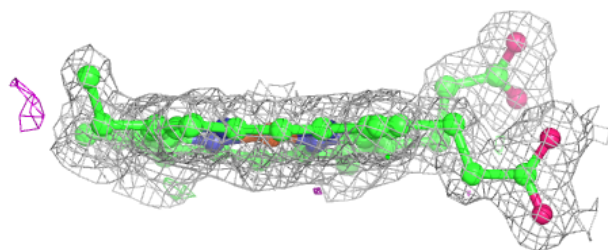
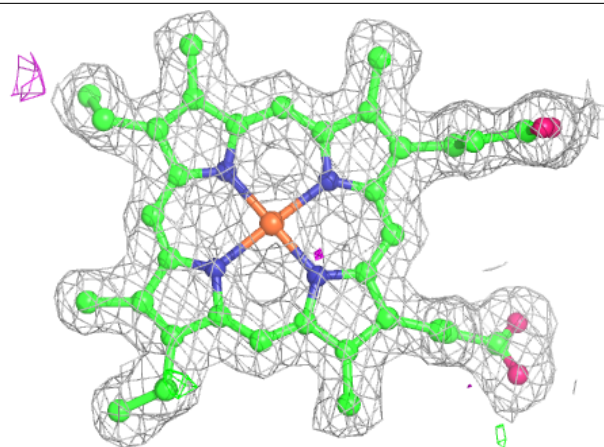
Electron density around HEC G 220:

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



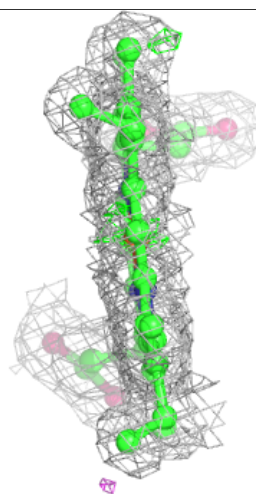
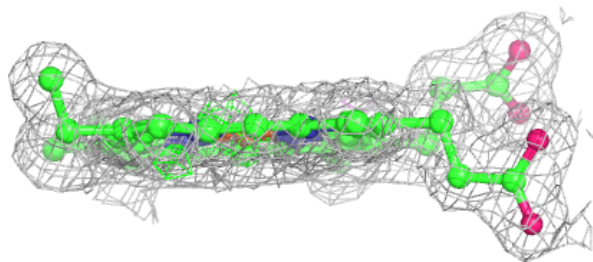
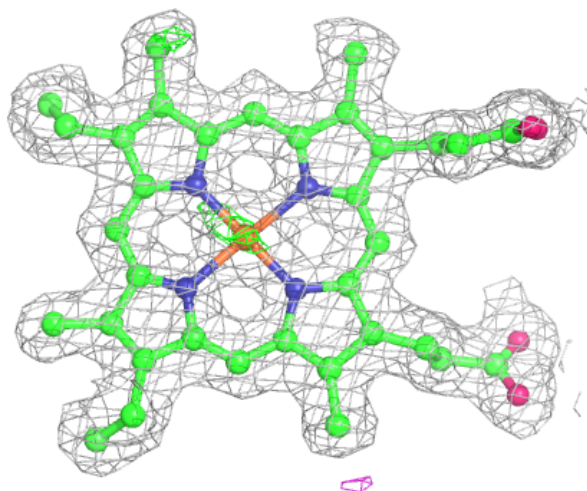
Electron density around HEC H 220:

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and green (positive)



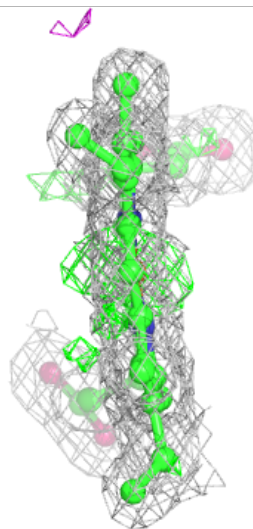
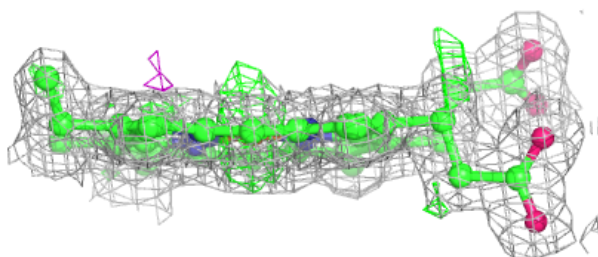
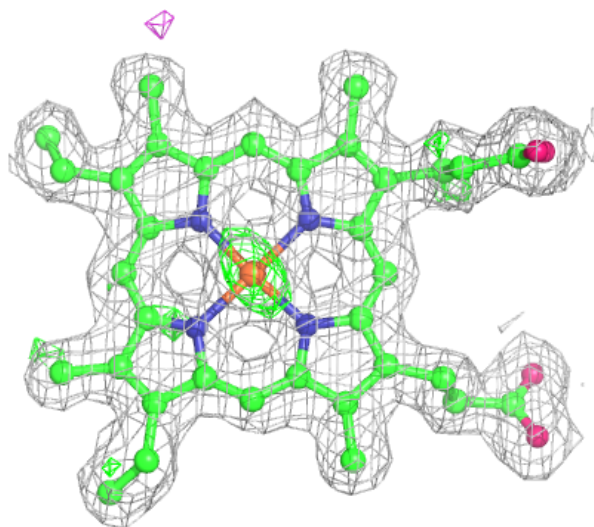
Electron density around HEC T 220:

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and green (positive)



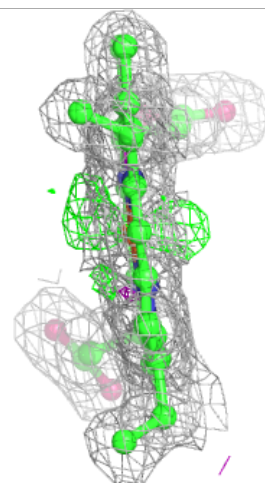
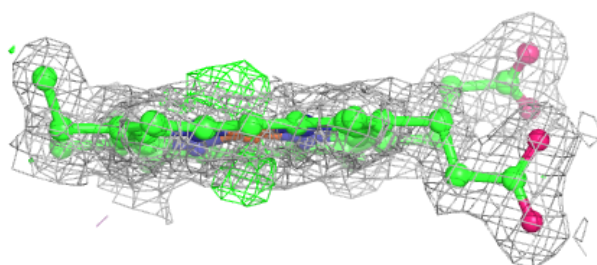
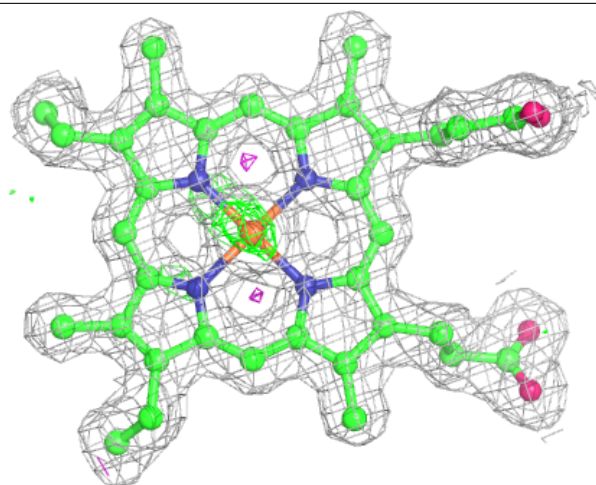
Electron density around HEC O 220:

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



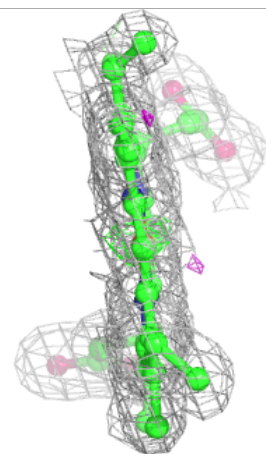
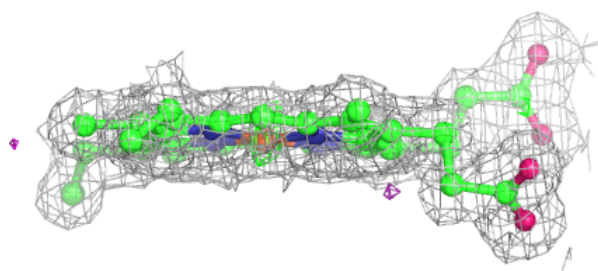
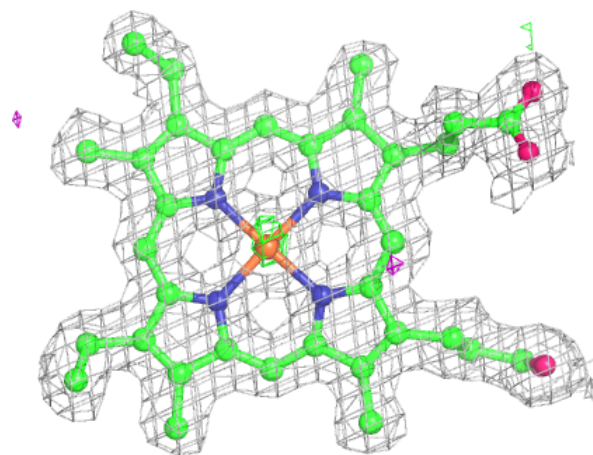
Electron density around HEC P 220:

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



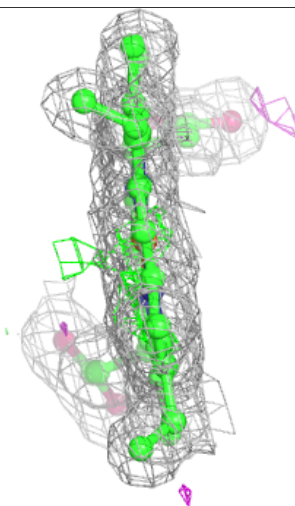
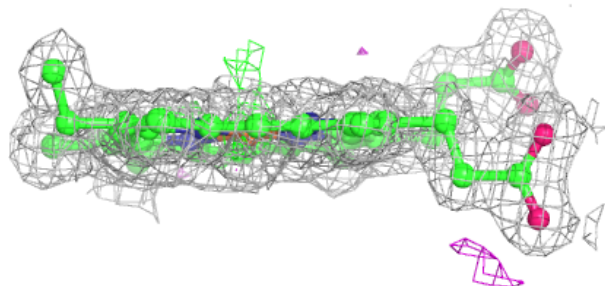
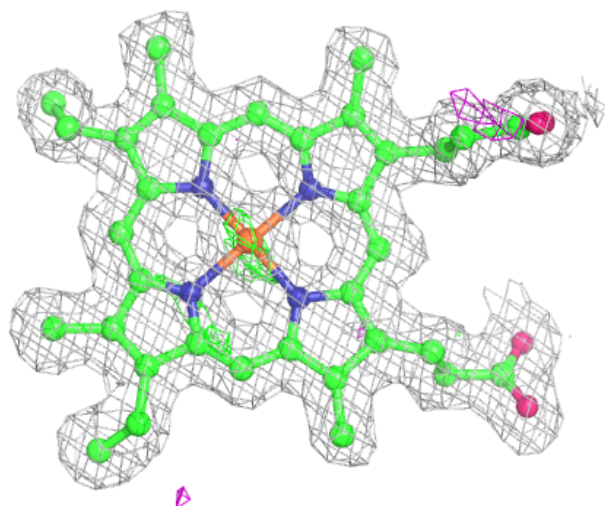
Electron density around HEC D 220:

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



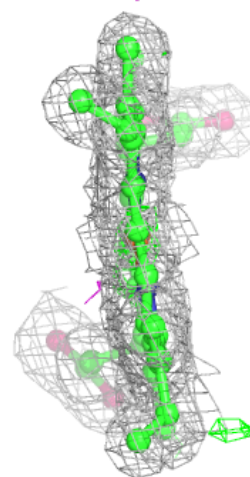
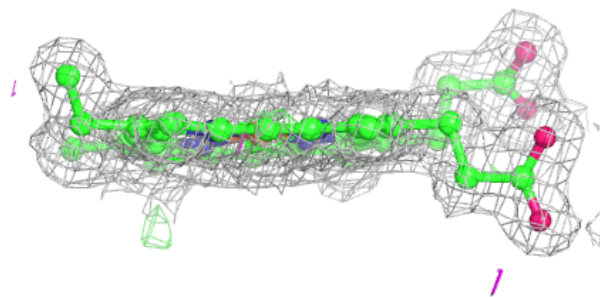
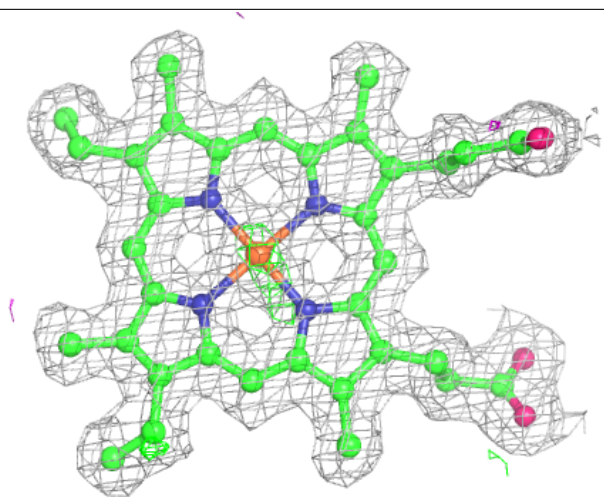
Electron density around HEC B 220:

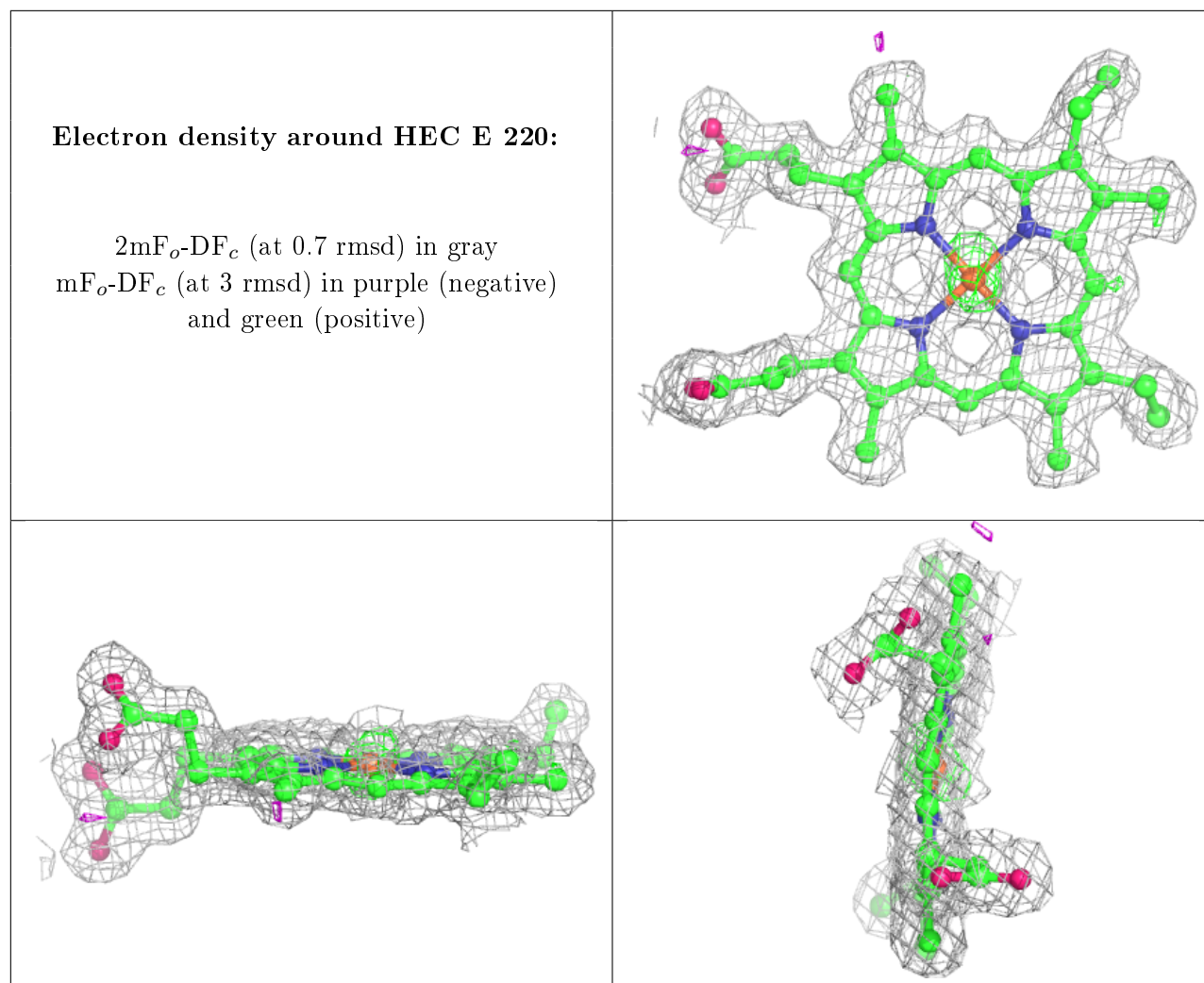
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around HEC F 220:

$2mF_o-DF_c$ (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.