



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 3, 2026 – 03:56 AM EDT

PDB ID : 5KLI / pdb_00005kli
Title : Rhodobacter sphaeroides bc1 with stigmatellin and antimycin
Authors : Xia, D.; Esser, L.; Zhou, F.; Tang, W.K.; Yu, C.A.
Deposited on : 2016-06-24
Resolution : 3.00 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

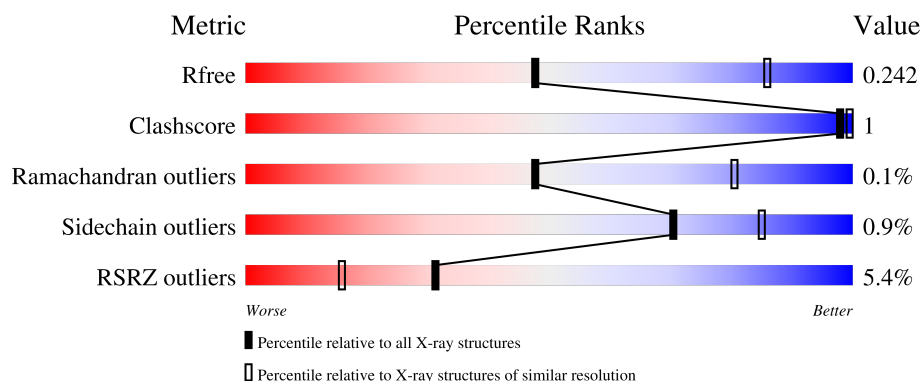
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.00 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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	445	6% 94% • •
1	E	445	4% 94% • •
1	K	445	2% 93% • •
1	O	445	4% 95% • •
2	B	272	7% 92% • 6%

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Mol	Chain	Length	Quality of chain
2	F	272	
2	L	272	
2	P	272	
3	C	187	
3	G	187	
3	M	187	
3	Q	187	

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
9	HEC	B	1001	X	-	-	-
9	HEC	F	1001	X	-	-	-
9	HEC	L	1001	X	-	-	-
9	HEC	P	1001	X	-	-	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 55093 atoms, of which 27155 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 b.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	428	Total	C	H	N	O	S	0	0	0
			6841	2319	3406	545	556	15			
1	E	428	Total	C	H	N	O	S	0	0	0
			6841	2319	3406	545	556	15			
1	K	428	Total	C	H	N	O	S	0	0	0
			6841	2319	3406	545	556	15			
1	O	428	Total	C	H	N	O	S	0	0	0
			6841	2319	3406	545	556	15			

- Molecule 2 is a protein called Cytochrome c1.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
2	B	256	Total	C	H	N	O	S	0	0	0
			3792	1240	1839	326	374	13			
2	F	256	Total	C	H	N	O	S	0	0	0
			3792	1240	1839	326	374	13			
2	L	256	Total	C	H	N	O	S	0	0	0
			3791	1240	1838	326	374	13			
2	P	256	Total	C	H	N	O	S	0	0	0
			3792	1240	1839	326	374	13			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	98	PRO	ALA	variant	UNP Q02760
B	264	GLY	-	expression tag	UNP Q02760
B	265	THR	-	expression tag	UNP Q02760
B	266	GLY	-	expression tag	UNP Q02760
B	267	HIS	-	expression tag	UNP Q02760
B	268	HIS	-	expression tag	UNP Q02760
B	269	HIS	-	expression tag	UNP Q02760
B	270	HIS	-	expression tag	UNP Q02760

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Chain	Residue	Modelled	Actual	Comment	Reference
B	271	HIS	-	expression tag	UNP Q02760
B	272	HIS	-	expression tag	UNP Q02760
F	98	PRO	ALA	variant	UNP Q02760
F	264	GLY	-	expression tag	UNP Q02760
F	265	THR	-	expression tag	UNP Q02760
F	266	GLY	-	expression tag	UNP Q02760
F	267	HIS	-	expression tag	UNP Q02760
F	268	HIS	-	expression tag	UNP Q02760
F	269	HIS	-	expression tag	UNP Q02760
F	270	HIS	-	expression tag	UNP Q02760
F	271	HIS	-	expression tag	UNP Q02760
F	272	HIS	-	expression tag	UNP Q02760
L	98	PRO	ALA	variant	UNP Q02760
L	264	GLY	-	expression tag	UNP Q02760
L	265	THR	-	expression tag	UNP Q02760
L	266	GLY	-	expression tag	UNP Q02760
L	267	HIS	-	expression tag	UNP Q02760
L	268	HIS	-	expression tag	UNP Q02760
L	269	HIS	-	expression tag	UNP Q02760
L	270	HIS	-	expression tag	UNP Q02760
L	271	HIS	-	expression tag	UNP Q02760
L	272	HIS	-	expression tag	UNP Q02760
P	98	PRO	ALA	variant	UNP Q02760
P	264	GLY	-	expression tag	UNP Q02760
P	265	THR	-	expression tag	UNP Q02760
P	266	GLY	-	expression tag	UNP Q02760
P	267	HIS	-	expression tag	UNP Q02760
P	268	HIS	-	expression tag	UNP Q02760
P	269	HIS	-	expression tag	UNP Q02760
P	270	HIS	-	expression tag	UNP Q02760
P	271	HIS	-	expression tag	UNP Q02760
P	272	HIS	-	expression tag	UNP Q02760

- Molecule 3 is a protein called Ubiquinol-cytochrome c reductase iron-sulfur subunit.

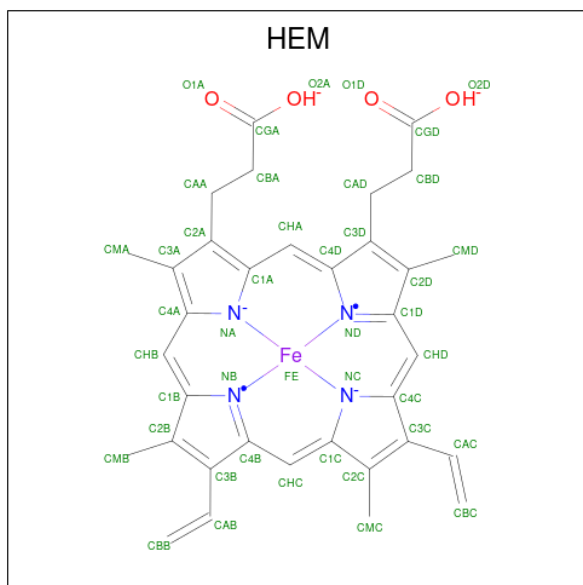
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	C	179	Total	C	H	N	O	S	0	0	0
			2645	845	1304	237	253	6			
3	G	179	Total	C	H	N	O	S	0	0	0
			2645	845	1304	237	253	6			
3	M	179	Total	C	H	N	O	S	0	0	0
			2645	845	1304	237	253	6			

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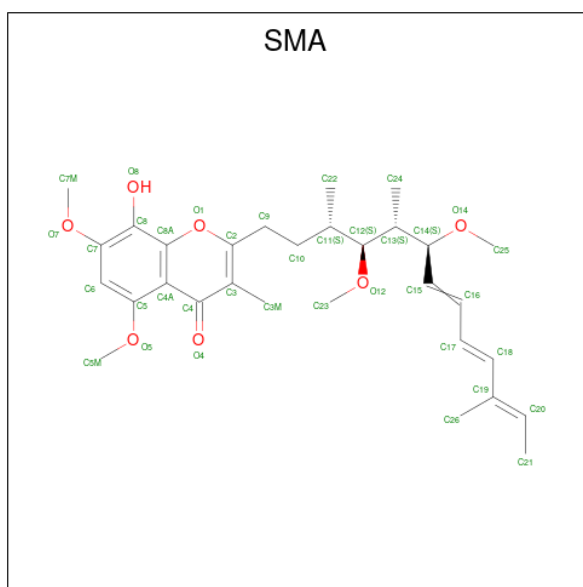
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
3	Q	179	Total	C	H	N	O	S	0	0	0
			2645	845	1304	237	253	6			

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



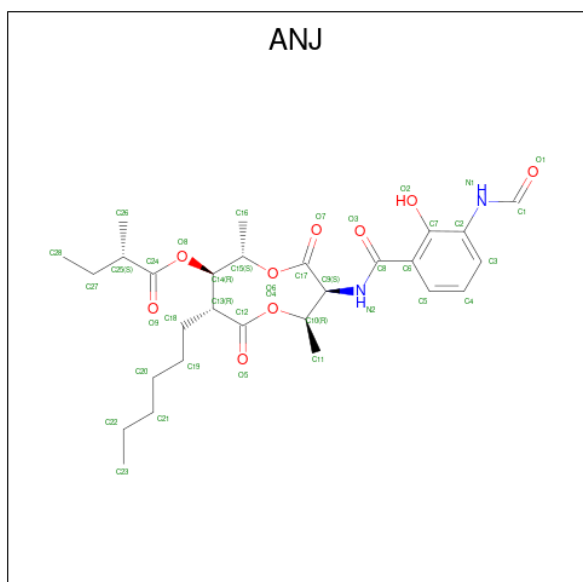
Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
4	A	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
4	A	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
4	E	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
4	E	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
4	K	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
4	K	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
4	O	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		
4	O	1	Total	C	Fe	H	N	O	0	0
			73	34	1	30	4	4		

- Molecule 5 is STIGMATELLIN A (CCD ID: SMA) (formula: $C_{30}H_{42}O_7$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	H	O	0	0
			79	30	42	7		
5	E	1	Total	C	H	O	0	0
			79	30	42	7		
5	K	1	Total	C	H	O	0	0
			79	30	42	7		
5	O	1	Total	C	H	O	0	0
			79	30	42	7		

- Molecule 6 is (2R,3S,6S,7R,8R)-3-[[3-(FORMYLAMINO)-2-HYDROXYBENZOYL]AMINO O}-8-HEXYL-2,6-DIMETHYL-4,9-DIOXO-1,5-DIOXONAN-7-YL (2S)-2-METHYLBUTANOATE (CCD ID: ANJ) (formula: C₂₈H₄₀N₂O₉).

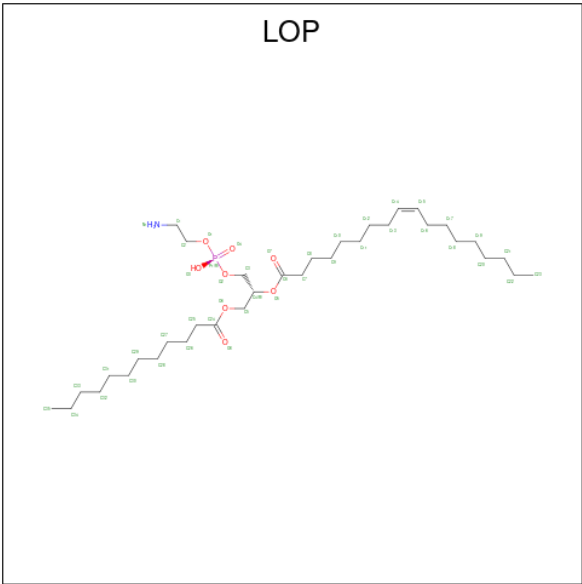


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total	C	H	N	O	0	0
			78	28	39	2	9		
6	E	1	Total	C	H	N	O	0	0
			78	28	39	2	9		
6	K	1	Total	C	H	N	O	0	0
			78	28	39	2	9		
6	O	1	Total	C	H	N	O	0	0
			78	28	39	2	9		

- Molecule 7 is STRONTIUM ION (CCD ID: SR) (formula: Sr).

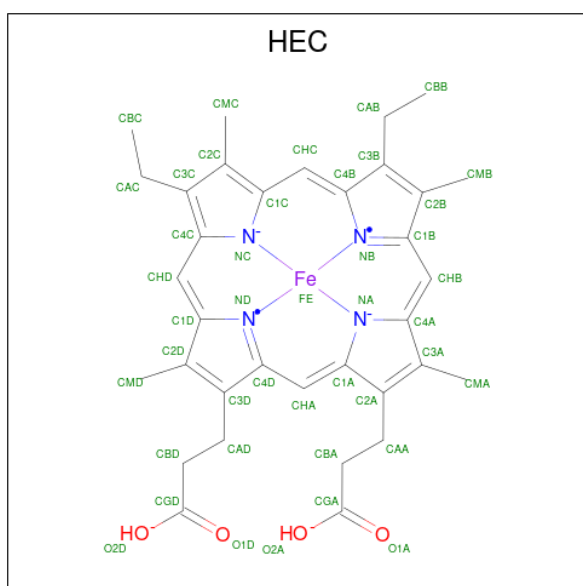
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Sr	0	0
			1	1		
7	B	1	Total	Sr	0	0
			1	1		
7	F	1	Total	Sr	0	0
			1	1		
7	K	1	Total	Sr	0	0
			1	1		
7	L	1	Total	Sr	0	0
			1	1		
7	P	1	Total	Sr	0	0
			1	1		

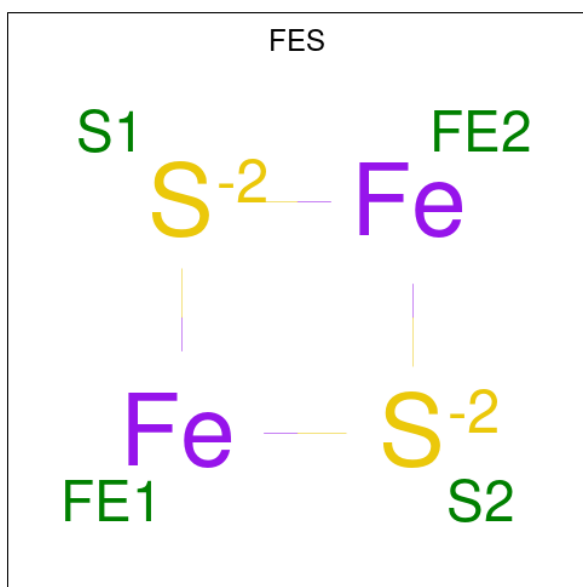
- Molecule 8 is (1R)-2-{[(R)-(2-AMINOETHOXY)(HYDROXY)PHOSPHORYL]OXY}-1-[(DODECANOYLOXY)METHYL]ETHYL (9Z)-OCTADEC-9-ENOATE (CCD ID: LOP) (formula: C₃₅H₆₈NO₈P).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
8	A	1	Total	C	H	N	O	P	
			112	35	67	1	8	1	0
8	E	1	Total	C	H	N	O	P	
			112	35	67	1	8	1	0
8	K	1	Total	C	H	N	O	P	
			112	35	67	1	8	1	0
8	O	1	Total	C	H	N	O	P	
			112	35	67	1	8	1	0

- Molecule 9 is HEME C (CCD ID: HEC) (formula: $C_{34}H_{36}FeN_4O_4$).



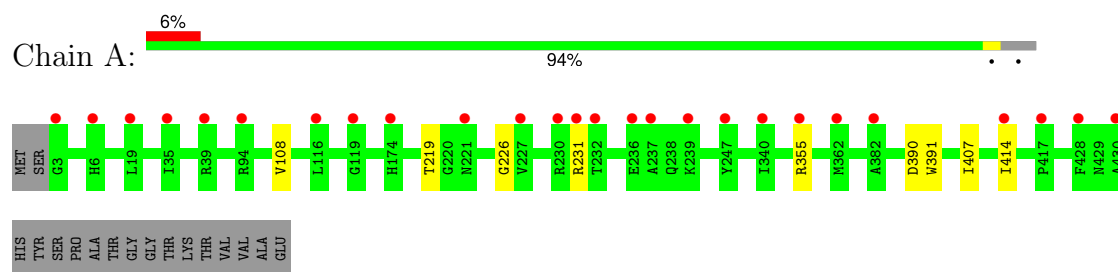


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	C	1	Total	Fe	S	0	0
			4	2	2		
10	G	1	Total	Fe	S	0	0
			4	2	2		
10	M	1	Total	Fe	S	0	0
			4	2	2		
10	Q	1	Total	Fe	S	0	0
			4	2	2		

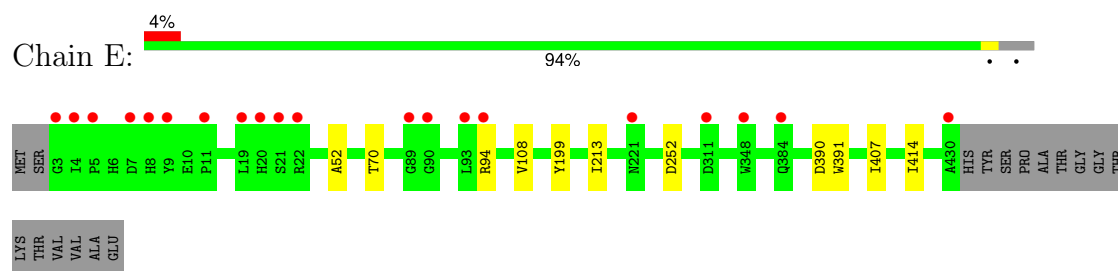
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

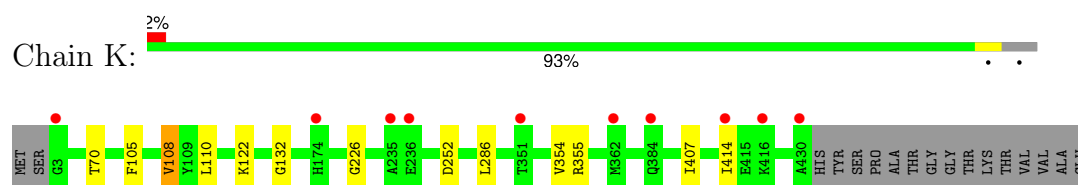
- Molecule 1: Cytochrome b



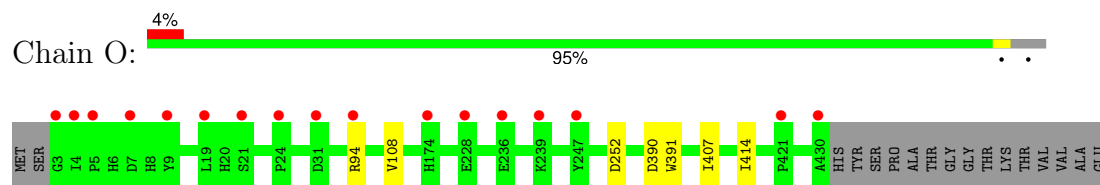
- Molecule 1: Cytochrome b



- Molecule 1: Cytochrome b

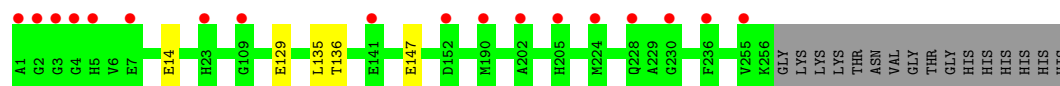


- Molecule 1: Cytochrome b

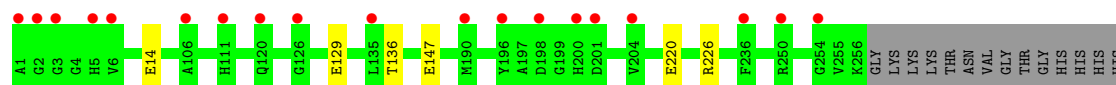


- Molecule 2: Cytochrome c1

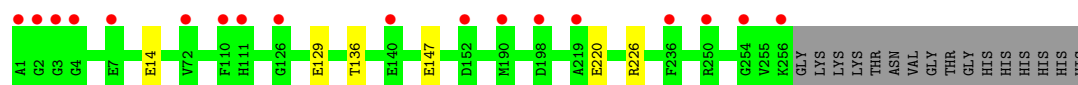
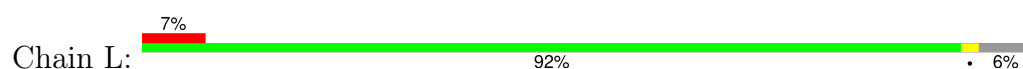




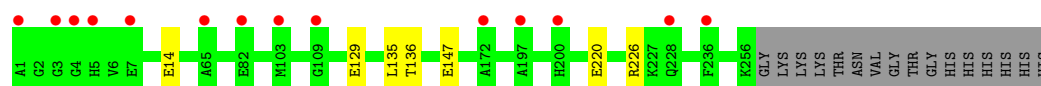
- Molecule 2: Cytochrome c1



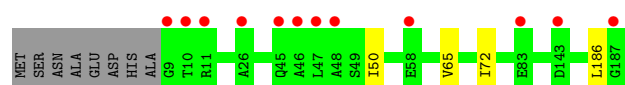
- Molecule 2: Cytochrome c1



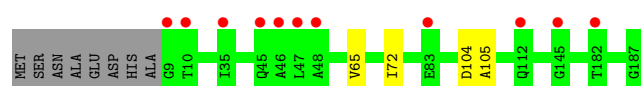
- Molecule 2: Cytochrome c1



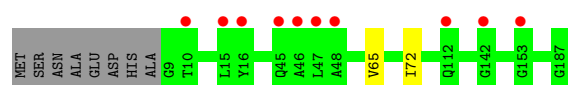
- Molecule 3: Ubiquinol-cytochrome c reductase iron-sulfur subunit



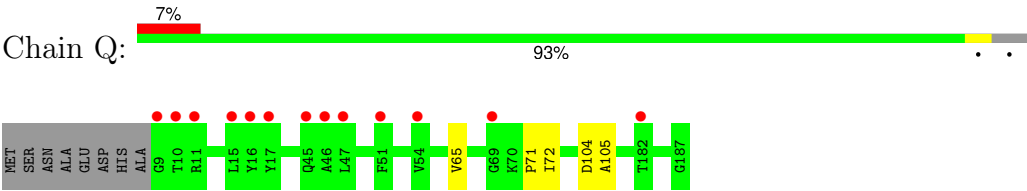
- Molecule 3: Ubiquinol-cytochrome c reductase iron-sulfur subunit



- Molecule 3: Ubiquinol-cytochrome c reductase iron-sulfur subunit



● Molecule 3: Ubiquinol-cytochrome c reductase iron-sulfur subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	118.86Å 126.93Å 127.89Å 64.64° 87.69° 61.80°	Depositor
Resolution (Å)	37.74 – 3.00 37.74 – 3.00	Depositor EDS
% Data completeness (in resolution range)	96.2 (37.74-3.00) 79.4 (37.74-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.08 (at 3.01Å)	Xtriage
Refinement program	PHENIX (1.11rc1_2513: ???)	Depositor
R, R_{free}	0.211 , 0.243 0.214 , 0.242	Depositor DCC
R_{free} test set	2065 reflections (1.80%)	wwPDB-VP
Wilson B-factor (Å ²)	58.7	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 49.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.014 for h,h-k,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	55093	wwPDB-VP
Average B, all atoms (Å ²)	92.0	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 34.19 % 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 7.0909e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ANJ, HEC, SR, SMA, LOP, HEM, FES

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	0.32	0/3565	0.57	0/4891
1	E	0.32	1/3565 (0.0%)	0.56	0/4891
1	K	0.33	1/3565 (0.0%)	0.59	0/4891
1	O	0.30	0/3565	0.55	0/4891
2	B	0.27	0/2010	0.55	0/2733
2	F	0.26	0/2010	0.55	0/2733
2	L	0.27	0/2010	0.58	0/2733
2	P	0.26	0/2010	0.51	0/2733
3	C	0.29	0/1371	0.55	0/1868
3	G	0.28	0/1371	0.52	0/1868
3	M	0.28	0/1371	0.52	0/1868
3	Q	0.28	0/1371	0.51	0/1868
All	All	0.30	2/27784 (0.0%)	0.55	0/37968

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	70	THR	C-N	6.21	1.48	1.33
1	K	70	THR	C-N	5.05	1.45	1.33

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3435	3406	3420	3	0
1	E	3435	3406	3420	6	0
1	K	3435	3406	3420	7	0
1	O	3435	3406	3420	3	0
2	B	1953	1839	1848	2	0
2	F	1953	1839	1848	2	0
2	L	1953	1838	1848	2	0
2	P	1953	1839	1848	3	0
3	C	1341	1304	1307	1	0
3	G	1341	1304	1307	1	0
3	M	1341	1304	1307	0	0
3	Q	1341	1304	1307	2	0
4	A	86	60	60	2	0
4	E	86	60	60	2	0
4	K	86	60	60	3	0
4	O	86	60	60	2	0
5	A	37	42	42	0	0
5	E	37	42	42	0	0
5	K	37	42	42	0	0
5	O	37	42	42	0	0
6	A	39	39	39	0	0
6	E	39	39	39	3	0
6	K	39	39	40	1	0
6	O	39	39	39	1	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	F	1	0	0	0	0
7	K	1	0	0	0	0
7	L	1	0	0	0	0
7	P	1	0	0	0	0
8	A	45	67	67	0	0
8	E	45	67	67	0	0
8	K	45	67	67	1	0
8	O	45	67	67	0	0
9	B	43	32	30	1	0
9	F	43	32	30	1	0
9	L	43	32	30	1	0
9	P	43	32	30	1	0
10	C	4	0	0	0	0
10	G	4	0	0	0	0
10	M	4	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	Q	4	0	0	0	0
All	All	27938	27155	27253	38	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 1.

All (38) 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)
2:B:14:GLU:OE2	2:B:129:GLU:OE2	2.10	0.70
9:F:1001:HEC:HMB3	9:F:1001:HEC:HBB3	1.77	0.67
9:L:1001:HEC:HBB3	9:L:1001:HEC:HMB3	1.77	0.66
2:F:14:GLU:OE2	2:F:129:GLU:OE2	2.18	0.60
2:F:220:GLU:OE1	2:F:226:ARG:NH1	2.35	0.59
2:P:220:GLU:OE1	2:P:226:ARG:NH1	2.38	0.56
1:E:52:ALA:HB2	6:E:1004:ANJ:H163	1.86	0.56
4:O:1001:HEM:HBB2	4:O:1001:HEM:HHC	1.89	0.55
1:A:219:THR:O	1:A:231:ARG:NH2	2.38	0.55
4:K:1001:HEM:HBC2	4:K:1001:HEM:HHD	1.90	0.53
2:L:14:GLU:OE2	2:L:129:GLU:OE2	2.27	0.53
4:K:1001:HEM:HBB2	4:K:1001:HEM:HHC	1.90	0.52
3:C:50:ILE:HD11	3:C:186:LEU:HD12	1.92	0.52
1:A:226:GLY:O	1:A:355:ARG:NH1	2.38	0.51
2:P:14:GLU:OE2	2:P:129:GLU:OE2	2.29	0.51
4:A:1001:HEM:HBB2	4:A:1001:HEM:HHC	1.94	0.49
1:K:110:LEU:HD11	8:K:1006:LOP:H14	1.95	0.49
1:O:94:ARG:NH2	4:O:1001:HEM:O1A	2.45	0.49
2:L:220:GLU:OE1	2:L:226:ARG:NH1	2.46	0.48
1:O:390:ASP:OD1	1:O:391:TRP:N	2.47	0.48
1:E:94:ARG:NH2	4:E:1001:HEM:O1A	2.47	0.48
3:G:104:ASP:OD1	3:G:105:ALA:N	2.47	0.48
1:E:252:ASP:OD2	6:E:1004:ANJ:O2	2.33	0.47
1:A:390:ASP:OD1	1:A:391:TRP:N	2.49	0.46
1:E:213:ILE:HD11	6:E:1004:ANJ:H162	1.97	0.46
1:E:199:TYR:CE2	4:E:1001:HEM:HBC1	2.50	0.46
1:K:122:LYS:NZ	1:K:354:VAL:O	2.47	0.45
1:K:226:GLY:O	1:K:355:ARG:NH1	2.41	0.44
1:K:252:ASP:OD2	6:K:1004:ANJ:O2	2.34	0.43
1:O:252:ASP:OD2	6:O:1004:ANJ:O2	2.36	0.43
1:K:132:GLY:O	4:K:1002:HEM:HMC3	2.18	0.43
3:Q:104:ASP:OD1	3:Q:105:ALA:N	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:1002:HEM:HBA1	4:A:1002:HEM:HHA	2.00	0.43
1:E:390:ASP:OD1	1:E:391:TRP:N	2.53	0.42
2:P:135:LEU:HD21	9:P:1001:HEC:HMB2	2.01	0.41
1:K:105:PHE:HA	1:K:108:VAL:HG22	2.01	0.41
1:K:286:LEU:HA	3:Q:71:PRO:HG3	2.03	0.41
2:B:135:LEU:HD21	9:B:1001:HEC:HMB2	2.03	0.40

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	426/445 (96%)	421 (99%)	5 (1%)	0	100	100
1	E	426/445 (96%)	421 (99%)	5 (1%)	0	100	100
1	K	426/445 (96%)	421 (99%)	5 (1%)	0	100	100
1	O	426/445 (96%)	422 (99%)	4 (1%)	0	100	100
2	B	254/272 (93%)	246 (97%)	7 (3%)	1 (0%)	30	65
2	F	254/272 (93%)	246 (97%)	7 (3%)	1 (0%)	30	65
2	L	254/272 (93%)	246 (97%)	7 (3%)	1 (0%)	30	65
2	P	254/272 (93%)	246 (97%)	7 (3%)	1 (0%)	30	65
3	C	177/187 (95%)	171 (97%)	6 (3%)	0	100	100
3	G	177/187 (95%)	171 (97%)	6 (3%)	0	100	100
3	M	177/187 (95%)	170 (96%)	7 (4%)	0	100	100
3	Q	177/187 (95%)	170 (96%)	7 (4%)	0	100	100
All	All	3428/3616 (95%)	3351 (98%)	73 (2%)	4 (0%)	48	80

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	L	147	GLU
2	B	147	GLU
2	F	147	GLU
2	P	147	GLU

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	353/366 (96%)	350 (99%)	3 (1%)	73	86
1	E	353/366 (96%)	350 (99%)	3 (1%)	73	86
1	K	353/366 (96%)	350 (99%)	3 (1%)	73	86
1	O	353/366 (96%)	350 (99%)	3 (1%)	73	86
2	B	203/216 (94%)	202 (100%)	1 (0%)	81	89
2	F	203/216 (94%)	202 (100%)	1 (0%)	81	89
2	L	203/216 (94%)	202 (100%)	1 (0%)	81	89
2	P	203/216 (94%)	202 (100%)	1 (0%)	81	89
3	C	138/144 (96%)	136 (99%)	2 (1%)	59	80
3	G	138/144 (96%)	136 (99%)	2 (1%)	59	80
3	M	138/144 (96%)	136 (99%)	2 (1%)	59	80
3	Q	138/144 (96%)	136 (99%)	2 (1%)	59	80
All	All	2776/2904 (96%)	2752 (99%)	24 (1%)	70	85

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	VAL
1	A	407	ILE
1	A	414	ILE
2	B	136	THR
3	C	65	VAL
3	C	72	ILE
1	E	108	VAL

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Mol	Chain	Res	Type
1	E	407	ILE
1	E	414	ILE
2	F	136	THR
3	G	65	VAL
3	G	72	ILE
1	K	108	VAL
1	K	407	ILE
1	K	414	ILE
2	L	136	THR
3	M	65	VAL
3	M	72	ILE
1	O	108	VAL
1	O	407	ILE
1	O	414	ILE
2	P	136	THR
3	Q	65	VAL
3	Q	72	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	192	ASN
2	B	91	HIS
2	B	200	HIS
3	C	92	GLN
1	E	20	HIS
1	E	177	GLN
1	E	192	ASN
1	E	291	HIS
2	F	88	HIS
2	F	111	HIS
2	F	200	HIS
3	G	92	GLN
1	K	58	GLN
1	K	192	ASN
1	K	291	HIS
2	L	200	HIS
3	M	92	GLN
1	O	192	ASN
1	O	291	HIS
1	O	319	ASN
2	P	200	HIS

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Mol	Chain	Res	Type
3	Q	92	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 34 ligands modelled in this entry, 6 are monoatomic - leaving 28 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
4	HEM	K	1002	1	50,50,50	1.59	7 (14%)	67,82,82	1.14	5 (7%)
8	LOP	O	1005	-	44,44,44	0.94	1 (2%)	47,49,49	1.18	3 (6%)
8	LOP	A	1006	-	44,44,44	0.93	1 (2%)	47,49,49	1.14	4 (8%)
6	ANJ	E	1004	-	40,40,40	0.92	2 (5%)	34,54,54	1.09	1 (2%)
4	HEM	E	1001	1	50,50,50	1.34	6 (12%)	67,82,82	1.06	4 (5%)
9	HEC	B	1001	2	50,50,50	1.73	6 (12%)	68,82,82	1.31	8 (11%)
4	HEM	A	1002	1	50,50,50	1.39	6 (12%)	67,82,82	1.15	4 (5%)
4	HEM	E	1002	1	50,50,50	1.56	8 (16%)	67,82,82	1.26	8 (11%)
10	FES	Q	1001	3	0,4,4	-	-	-	-	-
5	SMA	E	1003	-	38,38,38	0.77	0	47,52,52	1.58	8 (17%)
10	FES	M	1001	3	0,4,4	-	-	-	-	-

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	HEC	L	1001	2	50,50,50	1.73	7 (14%)	68,82,82	1.33	7 (10%)
10	FES	G	1001	3	0,4,4	-	-	-	-	-
4	HEM	K	1001	1	50,50,50	1.41	7 (14%)	67,82,82	1.24	7 (10%)
4	HEM	O	1002	1	50,50,50	1.45	9 (18%)	67,82,82	1.25	4 (5%)
4	HEM	A	1001	1	50,50,50	1.44	9 (18%)	67,82,82	1.10	5 (7%)
4	HEM	O	1001	1	50,50,50	1.45	8 (16%)	67,82,82	1.08	4 (5%)
6	ANJ	K	1004	-	40,40,40	0.92	1 (2%)	34,54,54	1.20	5 (14%)
9	HEC	F	1001	2	50,50,50	1.72	7 (14%)	68,82,82	1.28	7 (10%)
5	SMA	A	1003	-	38,38,38	0.77	0	47,52,52	1.77	9 (19%)
5	SMA	K	1003	-	38,38,38	0.75	0	47,52,52	1.65	9 (19%)
6	ANJ	A	1004	-	40,40,40	0.86	0	34,54,54	1.42	6 (17%)
10	FES	C	1001	3	0,4,4	-	-	-	-	-
6	ANJ	O	1004	-	40,40,40	0.89	0	34,54,54	1.23	5 (14%)
8	LOP	K	1006	-	44,44,44	0.97	2 (4%)	47,49,49	1.09	4 (8%)
5	SMA	O	1003	-	38,38,38	0.82	0	47,52,52	1.71	7 (14%)
9	HEC	P	1001	2	50,50,50	1.75	8 (16%)	68,82,82	1.33	8 (11%)
8	LOP	E	1005	-	44,44,44	0.95	1 (2%)	47,49,49	1.18	3 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	HEM	K	1002	1	-	2/14/54/54	-
8	LOP	O	1005	-	-	19/48/48/48	-
8	LOP	A	1006	-	-	24/48/48/48	-
6	ANJ	E	1004	-	-	7/40/55/55	0/1/2/2
4	HEM	E	1001	1	-	5/14/54/54	-
9	HEC	B	1001	2	1/1/7/7	6/14/54/54	-
4	HEM	A	1002	1	-	3/14/54/54	-
4	HEM	E	1002	1	-	2/14/54/54	-
10	FES	Q	1001	3	-	-	0/1/1/1
5	SMA	E	1003	-	-	3/34/34/34	0/2/2/2
10	FES	M	1001	3	-	-	0/1/1/1
9	HEC	L	1001	2	1/1/7/7	7/14/54/54	-
10	FES	G	1001	3	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	HEM	K	1001	1	-	7/14/54/54	-
4	HEM	O	1002	1	-	2/14/54/54	-
4	HEM	A	1001	1	-	4/14/54/54	-
9	HEC	F	1001	2	1/1/7/7	6/14/54/54	-
4	HEM	O	1001	1	-	6/14/54/54	-
6	ANJ	K	1004	-	-	12/40/55/55	0/1/2/2
5	SMA	A	1003	-	-	4/34/34/34	0/2/2/2
5	SMA	K	1003	-	-	4/34/34/34	0/2/2/2
6	ANJ	A	1004	-	-	12/40/55/55	0/1/2/2
10	FES	C	1001	3	-	-	0/1/1/1
6	ANJ	O	1004	-	-	12/40/55/55	0/1/2/2
8	LOP	K	1006	-	-	21/48/48/48	-
5	SMA	O	1003	-	-	4/34/34/34	0/2/2/2
9	HEC	P	1001	2	1/1/7/7	6/14/54/54	-
8	LOP	E	1005	-	-	24/48/48/48	-

All (96) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	B	1001	HEC	C3D-C2D	7.32	1.52	1.36
9	F	1001	HEC	C3D-C2D	7.27	1.52	1.36
9	P	1001	HEC	C3D-C2D	7.27	1.52	1.36
9	L	1001	HEC	C3D-C2D	7.09	1.52	1.36
4	K	1002	HEM	FE-NB	6.07	2.13	1.94
4	E	1002	HEM	FE-NB	5.84	2.12	1.94
9	B	1001	HEC	FE-ND	5.43	2.11	1.94
9	P	1001	HEC	FE-ND	5.23	2.11	1.94
4	O	1001	HEM	FE-ND	5.08	2.10	1.94
9	F	1001	HEC	FE-ND	4.66	2.09	1.94
9	L	1001	HEC	FE-ND	4.52	2.08	1.94
9	L	1001	HEC	FE-NC	4.49	2.10	1.95
4	O	1002	HEM	FE-NB	4.30	2.08	1.94
4	E	1001	HEM	FE-ND	3.64	2.06	1.94
4	O	1002	HEM	FE-NA	3.59	2.07	1.95
4	K	1001	HEM	FE-NB	3.56	2.05	1.94
4	A	1002	HEM	CAB-C3B	3.53	1.56	1.47
4	A	1001	HEM	FE-ND	3.52	2.05	1.94
4	K	1002	HEM	CAC-C3C	3.47	1.56	1.47
4	O	1002	HEM	CAC-C3C	3.46	1.56	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	E	1002	HEM	CAC-C3C	3.44	1.56	1.47
9	P	1001	HEC	FE-NC	3.35	2.06	1.95
4	O	1002	HEM	CAB-C3B	3.35	1.56	1.47
4	K	1002	HEM	CAB-C3B	3.33	1.56	1.47
4	A	1002	HEM	CAC-C3C	3.31	1.56	1.47
4	K	1001	HEM	CAC-C3C	3.31	1.56	1.47
4	E	1002	HEM	CAB-C3B	3.30	1.56	1.47
4	A	1002	HEM	FE-NB	3.29	2.05	1.94
4	O	1001	HEM	FE-NA	3.29	2.06	1.95
4	K	1002	HEM	FE-NC	3.25	2.05	1.95
4	A	1001	HEM	FE-NA	3.21	2.05	1.95
4	A	1001	HEM	CAC-C3C	3.14	1.55	1.47
4	K	1002	HEM	FE-NA	3.07	2.05	1.95
9	F	1001	HEC	FE-NC	3.06	2.05	1.95
4	K	1001	HEM	FE-NC	3.06	2.05	1.95
4	E	1002	HEM	FE-NC	3.06	2.05	1.95
9	B	1001	HEC	FE-NA	3.05	2.05	1.95
4	E	1002	HEM	FE-NA	3.01	2.05	1.95
4	A	1001	HEM	CMC-C2C	3.01	1.56	1.50
4	O	1001	HEM	CAC-C3C	2.97	1.55	1.47
4	E	1001	HEM	CAC-C3C	2.95	1.55	1.47
4	E	1001	HEM	CAB-C3B	2.93	1.55	1.47
4	A	1002	HEM	FE-NC	2.88	2.04	1.95
4	O	1001	HEM	CAB-C3B	2.85	1.55	1.47
4	E	1002	HEM	CMB-C2B	2.83	1.56	1.50
4	A	1001	HEM	CAB-C3B	2.83	1.54	1.47
4	K	1001	HEM	CMC-C2C	2.75	1.56	1.50
9	B	1001	HEC	FE-NC	2.73	2.04	1.95
9	F	1001	HEC	FE-NA	2.70	2.04	1.95
4	K	1002	HEM	CMB-C2B	2.68	1.56	1.50
9	F	1001	HEC	FE-NB	2.66	2.03	1.94
4	K	1001	HEM	CAB-C3B	2.65	1.54	1.47
4	E	1001	HEM	CMC-C2C	2.63	1.56	1.50
9	P	1001	HEC	CMC-C2C	2.63	1.56	1.50
4	A	1001	HEM	FE-NB	2.59	2.02	1.94
9	P	1001	HEC	FE-NB	2.55	2.02	1.94
9	L	1001	HEC	FE-NB	2.52	2.02	1.94
4	E	1001	HEM	FE-NC	2.50	2.03	1.95
4	O	1002	HEM	CMC-C2C	2.47	1.55	1.50
4	O	1001	HEM	CMC-C2C	2.43	1.55	1.50
9	L	1001	HEC	FE-NA	2.43	2.03	1.95
4	O	1001	HEM	CMB-C2B	2.43	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	P	1001	HEC	FE-NA	2.36	2.02	1.95
9	B	1001	HEC	CMC-C2C	2.35	1.55	1.50
4	A	1002	HEM	CMB-C2B	2.34	1.55	1.50
4	E	1002	HEM	CMC-C2C	2.34	1.55	1.50
9	F	1001	HEC	CMC-C2C	2.31	1.55	1.50
4	A	1001	HEM	C2A-C3A	-2.31	1.32	1.38
9	F	1001	HEC	CMB-C2B	2.28	1.55	1.50
4	K	1001	HEM	CMB-C2B	2.24	1.55	1.50
4	K	1002	HEM	CMC-C2C	2.24	1.55	1.50
9	B	1001	HEC	CMB-C2B	2.23	1.55	1.50
4	E	1001	HEM	CMB-C2B	2.22	1.55	1.50
4	A	1001	HEM	CMB-C2B	2.22	1.55	1.50
8	K	1006	LOP	C7-C6	2.20	1.57	1.50
4	O	1001	HEM	CMD-C2D	2.20	1.55	1.50
8	E	1005	LOP	C7-C6	2.19	1.57	1.50
4	O	1002	HEM	CMB-C2B	2.18	1.55	1.50
9	L	1001	HEC	CMB-C2B	2.18	1.55	1.50
4	E	1002	HEM	CMA-C3A	2.18	1.55	1.50
9	L	1001	HEC	CMC-C2C	2.15	1.55	1.50
8	A	1006	LOP	C7-C6	2.15	1.56	1.50
9	P	1001	HEC	CMB-C2B	2.15	1.55	1.50
4	O	1002	HEM	CMD-C2D	2.14	1.55	1.50
4	O	1001	HEM	C2A-C3A	-2.13	1.33	1.38
4	O	1002	HEM	C2A-C3A	-2.12	1.33	1.38
4	O	1002	HEM	FE-NC	2.11	2.02	1.95
8	K	1006	LOP	C25-C24	2.09	1.56	1.50
4	A	1002	HEM	CMC-C2C	2.09	1.55	1.50
6	K	1004	ANJ	C1-N1	2.07	1.37	1.34
4	K	1001	HEM	C2A-C3A	-2.07	1.33	1.38
6	E	1004	ANJ	O8-C24	2.05	1.39	1.34
8	O	1005	LOP	C7-C6	2.02	1.56	1.50
4	A	1001	HEM	CMD-C2D	2.01	1.54	1.50
6	E	1004	ANJ	O8-C14	2.01	1.47	1.44
9	P	1001	HEC	CMA-C3A	2.01	1.54	1.50

All (135) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1003	SMA	O5-C5-C4A	7.26	126.12	115.84
5	O	1003	SMA	O5-C5-C4A	6.96	125.70	115.84
5	E	1003	SMA	O5-C5-C4A	6.02	124.36	115.84
5	K	1003	SMA	O5-C5-C4A	5.97	124.30	115.84

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1003	SMA	O5-C5-C6	-5.33	114.90	124.08
5	O	1003	SMA	O5-C5-C6	-5.03	115.41	124.08
9	L	1001	HEC	C1D-ND-C4D	4.77	110.86	105.21
9	P	1001	HEC	C1D-ND-C4D	4.65	110.71	105.21
9	F	1001	HEC	C1D-ND-C4D	4.49	110.53	105.21
9	B	1001	HEC	C1D-ND-C4D	4.37	110.38	105.21
5	K	1003	SMA	O5-C5-C6	-4.22	116.80	124.08
5	E	1003	SMA	O5-C5-C6	-4.19	116.86	124.08
8	O	1005	LOP	O6-C24-C25	3.73	123.19	111.83
4	A	1001	HEM	C3B-C4B-NB	-3.58	106.90	109.47
4	O	1001	HEM	C3B-C4B-NB	-3.54	106.92	109.47
8	O	1005	LOP	O6-C5-C4	3.52	118.55	108.40
4	K	1001	HEM	C3B-C4B-NB	-3.43	107.01	109.47
6	E	1004	ANJ	C14-O8-C24	3.39	123.40	117.76
8	E	1005	LOP	O5-C6-C7	3.39	118.82	111.48
8	E	1005	LOP	O6-C24-C25	3.36	122.08	111.83
9	B	1001	HEC	CAA-CBA-CGA	-3.30	104.92	113.67
4	O	1002	HEM	CMB-C2B-C1B	-3.23	119.99	125.03
6	A	1004	ANJ	O6-C17-C9	3.19	115.29	110.28
8	E	1005	LOP	O6-C5-C4	3.18	117.56	108.40
8	A	1006	LOP	O5-C6-C7	3.17	118.35	111.48
6	K	1004	ANJ	C14-O8-C24	3.12	122.95	117.76
4	E	1002	HEM	C4D-ND-C1D	3.11	108.89	105.21
8	A	1006	LOP	O5-C4-C3	3.11	119.50	108.34
4	E	1002	HEM	CMB-C2B-C1B	-3.07	120.23	125.03
5	K	1003	SMA	C22-C11-C10	2.99	114.96	110.34
9	P	1001	HEC	CAA-CBA-CGA	-2.97	105.79	113.67
4	K	1002	HEM	CMB-C2B-C1B	-2.93	120.46	125.03
9	P	1001	HEC	CMD-C2D-C1D	2.92	129.60	125.03
5	K	1003	SMA	C9-C10-C11	-2.88	109.08	114.46
4	K	1001	HEM	CHD-C4C-C3C	2.87	130.05	125.21
4	E	1001	HEM	CMB-C2B-C1B	-2.87	120.55	125.03
6	A	1004	ANJ	C3-C2-N1	2.85	126.85	122.54
9	F	1001	HEC	CAA-CBA-CGA	-2.85	106.12	113.67
4	O	1002	HEM	C4D-ND-C1D	2.83	108.55	105.21
4	A	1002	HEM	CMB-C2B-C1B	-2.81	120.64	125.03
5	E	1003	SMA	C5M-O5-C5	2.76	121.57	117.51
5	K	1003	SMA	C3M-C3-C2	2.76	127.38	123.28
4	O	1001	HEM	C1B-NB-C4B	2.73	108.44	105.21
4	K	1001	HEM	CHD-C4C-NC	-2.69	121.52	124.45
9	L	1001	HEC	CBD-CAD-C3D	-2.67	105.14	112.53
5	K	1003	SMA	C3M-C3-C4	-2.66	111.77	116.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	O	1004	ANJ	O6-C17-C9	2.66	114.47	110.28
5	E	1003	SMA	C9-C10-C11	-2.65	109.51	114.46
6	A	1004	ANJ	C14-O8-C24	2.64	122.14	117.76
5	O	1003	SMA	C3M-C3-C2	2.63	127.18	123.28
8	K	1006	LOP	O6-C5-C4	2.63	115.97	108.40
9	L	1001	HEC	CMD-C2D-C1D	2.61	129.12	125.03
5	E	1003	SMA	C3M-C3-C4	-2.61	111.87	116.68
9	L	1001	HEC	CHA-C4D-ND	2.59	127.21	124.42
4	O	1002	HEM	CBD-CAD-C3D	-2.58	105.39	112.53
5	O	1003	SMA	C3M-C3-C4	-2.58	111.93	116.68
5	E	1003	SMA	C3M-C3-C2	2.57	127.09	123.28
6	O	1004	ANJ	C14-O8-C24	2.57	122.02	117.76
4	E	1002	HEM	C3D-C4D-ND	-2.55	107.38	110.17
9	F	1001	HEC	C4B-NB-C1B	2.55	108.22	105.21
5	K	1003	SMA	C7M-O7-C7	2.54	121.24	117.51
6	A	1004	ANJ	O6-C17-O7	-2.53	120.96	124.10
8	A	1006	LOP	O5-C6-O7	-2.52	117.80	123.70
6	A	1004	ANJ	O8-C14-C15	-2.52	102.19	107.97
8	K	1006	LOP	C5-C4-C3	2.50	117.62	111.78
8	K	1006	LOP	O5-C6-C7	2.48	116.85	111.48
5	E	1003	SMA	C7M-O7-C7	2.48	121.16	117.51
9	L	1001	HEC	C4B-NB-C1B	2.48	108.15	105.21
6	K	1004	ANJ	O8-C14-C15	-2.44	102.36	107.97
4	K	1001	HEM	CAC-C3C-C4C	2.44	130.64	124.82
4	A	1001	HEM	C1B-NB-C4B	2.43	108.09	105.21
4	A	1001	HEM	CHC-C4B-C3B	2.42	129.91	125.07
5	O	1003	SMA	C22-C11-C10	2.42	114.09	110.34
4	K	1001	HEM	C1B-NB-C4B	2.42	108.07	105.21
4	A	1002	HEM	CMC-C2C-C1C	-2.42	120.47	124.73
4	K	1002	HEM	CBD-CAD-C3D	-2.41	105.87	112.53
4	A	1001	HEM	CAC-C3C-C4C	-2.40	119.08	124.82
4	E	1001	HEM	CAC-C3C-C4C	-2.39	119.11	124.82
9	L	1001	HEC	CAA-CBA-CGA	-2.38	107.34	113.67
6	K	1004	ANJ	O6-C17-O7	-2.37	121.16	124.10
4	O	1002	HEM	C3D-C4D-ND	-2.37	107.57	110.17
6	A	1004	ANJ	C9-N2-C8	2.35	126.27	121.53
9	B	1001	HEC	CMD-C2D-C1D	2.35	128.70	125.03
5	O	1003	SMA	C5M-O5-C5	2.34	120.95	117.51
4	E	1002	HEM	CMC-C2C-C1C	-2.34	120.61	124.73
4	K	1002	HEM	C4D-ND-C1D	2.34	107.97	105.21
9	B	1001	HEC	C1D-C2D-C3D	-2.33	104.53	106.98
5	A	1003	SMA	C22-C11-C10	2.32	113.94	110.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	1003	SMA	C7M-O7-C7	2.32	120.91	117.51
8	O	1005	LOP	O5-C6-C7	2.31	116.47	111.48
4	K	1001	HEM	CHC-C4B-C3B	2.30	129.67	125.07
9	P	1001	HEC	C4B-NB-C1B	2.29	107.92	105.21
4	A	1002	HEM	C3B-C2B-C1B	2.29	108.13	106.41
5	A	1003	SMA	C3M-C3-C4	-2.27	112.50	116.68
8	K	1006	LOP	O6-C24-C25	2.26	118.74	111.83
9	F	1001	HEC	C1D-C2D-C3D	-2.26	104.60	106.98
9	B	1001	HEC	CAC-C3C-C4C	-2.26	121.81	124.91
5	K	1003	SMA	C5M-O5-C5	2.25	120.82	117.51
5	A	1003	SMA	O12-C12-C13	2.24	111.28	107.97
9	F	1001	HEC	CMD-C2D-C1D	2.24	128.53	125.03
5	A	1003	SMA	O14-C14-C13	-2.23	103.36	107.97
5	K	1003	SMA	O14-C14-C13	-2.21	103.41	107.97
4	K	1002	HEM	CHA-C4D-ND	2.17	127.05	124.37
4	K	1002	HEM	CMC-C2C-C1C	-2.16	120.92	124.73
4	O	1001	HEM	CHC-C4B-C3B	2.15	129.36	125.07
4	E	1002	HEM	CAD-CBD-CGD	-2.14	107.98	113.67
6	O	1004	ANJ	O8-C14-C15	-2.14	103.05	107.97
5	A	1003	SMA	C3M-C3-C2	2.14	126.45	123.28
4	E	1002	HEM	CBA-CAA-C2A	-2.14	106.63	112.53
4	E	1002	HEM	C2D-C1D-ND	-2.13	107.44	109.90
9	B	1001	HEC	CAC-C3C-C2C	2.13	130.26	126.89
8	A	1006	LOP	P1-O1-C2	-2.13	111.13	121.26
9	P	1001	HEC	CAD-CBD-CGD	-2.13	108.03	113.67
4	E	1002	HEM	C2A-C1A-NA	-2.12	107.80	110.15
5	E	1003	SMA	C22-C11-C10	2.12	113.61	110.34
5	A	1003	SMA	C9-C10-C11	-2.11	110.51	114.46
9	B	1001	HEC	C3B-C4B-NB	-2.10	107.87	110.17
4	K	1001	HEM	CAC-C3C-C2C	-2.10	121.61	128.43
4	A	1001	HEM	CAB-C3B-C2B	-2.09	121.65	128.43
6	K	1004	ANJ	O1-C1-N1	2.08	128.57	125.72
9	F	1001	HEC	CAC-C3C-C2C	2.08	130.18	126.89
4	E	1001	HEM	C4C-C3C-C2C	2.08	108.62	106.81
6	O	1004	ANJ	O6-C17-O7	-2.07	121.53	124.10
9	P	1001	HEC	C3B-C4B-NB	-2.07	107.90	110.17
9	F	1001	HEC	CAC-C3C-C4C	-2.07	122.07	124.91
9	P	1001	HEC	CAD-C3D-C4D	2.06	128.29	124.70
6	O	1004	ANJ	O8-C24-O9	-2.06	120.23	123.95
6	K	1004	ANJ	C6-C8-N2	2.06	121.13	116.67
5	O	1003	SMA	C7M-O7-C7	2.06	120.53	117.51
4	A	1002	HEM	C1B-NB-C4B	2.03	107.62	105.21

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	O	1001	HEM	CBD-CAD-C3D	-2.03	106.91	112.53
4	E	1001	HEM	CHA-C1A-NA	2.03	127.53	123.86
9	P	1001	HEC	CHA-C1A-NA	2.02	126.66	124.45
9	L	1001	HEC	C2B-C1B-NB	-2.01	107.58	109.90
9	B	1001	HEC	C2A-C1A-NA	-2.00	108.39	110.32

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
9	B	1001	HEC	NA
9	F	1001	HEC	NA
9	L	1001	HEC	NA
9	P	1001	HEC	NA

All (202) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	1004	ANJ	C9-C10-O4-C12
6	A	1004	ANJ	C16-C15-O6-C17
6	A	1004	ANJ	O7-C17-O6-C15
6	A	1004	ANJ	C9-C17-O6-C15
6	A	1004	ANJ	O8-C14-C15-O6
6	A	1004	ANJ	O8-C14-C15-C16
6	A	1004	ANJ	C13-C14-C15-O6
6	E	1004	ANJ	O7-C17-O6-C15
6	E	1004	ANJ	C9-C17-O6-C15
6	K	1004	ANJ	C9-C10-O4-C12
6	K	1004	ANJ	C16-C15-O6-C17
6	K	1004	ANJ	O7-C17-O6-C15
6	K	1004	ANJ	C9-C17-O6-C15
6	K	1004	ANJ	O8-C14-C15-O6
6	K	1004	ANJ	O8-C14-C15-C16
6	K	1004	ANJ	C13-C14-C15-O6
6	O	1004	ANJ	C9-C10-O4-C12
6	O	1004	ANJ	C16-C15-O6-C17
6	O	1004	ANJ	O7-C17-O6-C15
6	O	1004	ANJ	C9-C17-O6-C15
6	O	1004	ANJ	O8-C14-C15-O6
6	O	1004	ANJ	C13-C14-C15-O6
8	A	1006	LOP	C3-O2-P1-O1
8	A	1006	LOP	C3-O2-P1-O4
8	E	1005	LOP	N1-C1-C2-O1

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Mol	Chain	Res	Type	Atoms
8	E	1005	LOP	C2-O1-P1-O2
8	E	1005	LOP	C2-O1-P1-O3
8	E	1005	LOP	C2-O1-P1-O4
8	E	1005	LOP	C3-O2-P1-O1
8	E	1005	LOP	C3-O2-P1-O3
8	E	1005	LOP	C3-O2-P1-O4
8	K	1006	LOP	C2-O1-P1-O2
8	K	1006	LOP	C2-O1-P1-O3
8	K	1006	LOP	C2-O1-P1-O4
8	K	1006	LOP	C3-O2-P1-O3
8	K	1006	LOP	O5-C4-C5-O6
9	B	1001	HEC	C4C-C3C-CAC-CBC
9	F	1001	HEC	C4C-C3C-CAC-CBC
9	L	1001	HEC	C4C-C3C-CAC-CBC
9	P	1001	HEC	C4C-C3C-CAC-CBC
9	B	1001	HEC	C2C-C3C-CAC-CBC
9	F	1001	HEC	C2C-C3C-CAC-CBC
9	L	1001	HEC	C2C-C3C-CAC-CBC
9	P	1001	HEC	C2C-C3C-CAC-CBC
6	A	1004	ANJ	C14-C13-C18-C19
6	K	1004	ANJ	C14-C13-C18-C19
6	E	1004	ANJ	C12-C13-C18-C19
6	K	1004	ANJ	C12-C13-C18-C19
6	O	1004	ANJ	C12-C13-C18-C19
9	L	1001	HEC	C2B-C3B-CAB-CBB
8	A	1006	LOP	C13-C14-C15-C16
5	A	1003	SMA	C6-C5-O5-C5M
9	F	1001	HEC	C2B-C3B-CAB-CBB
5	K	1003	SMA	C6-C5-O5-C5M
5	O	1003	SMA	C6-C5-O5-C5M
5	E	1003	SMA	C6-C5-O5-C5M
8	K	1006	LOP	C13-C14-C15-C16
9	P	1001	HEC	C2B-C3B-CAB-CBB
8	E	1005	LOP	C24-C25-C26-C27
6	A	1004	ANJ	O9-C24-C25-C26
5	A	1003	SMA	C4A-C5-O5-C5M
9	B	1001	HEC	C2B-C3B-CAB-CBB
5	K	1003	SMA	C4A-C5-O5-C5M
5	O	1003	SMA	C4A-C5-O5-C5M
8	A	1006	LOP	O7-C6-O5-C4
5	E	1003	SMA	C4A-C5-O5-C5M
8	O	1005	LOP	C24-C25-C26-C27

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Mol	Chain	Res	Type	Atoms
8	O	1005	LOP	O7-C6-O5-C4
6	E	1004	ANJ	C14-C13-C18-C19
8	E	1005	LOP	C17-C18-C19-C20
8	O	1005	LOP	C17-C18-C19-C20
9	L	1001	HEC	C4B-C3B-CAB-CBB
8	E	1005	LOP	C9-C10-C11-C12
9	F	1001	HEC	C4B-C3B-CAB-CBB
8	E	1005	LOP	C16-C17-C18-C19
8	A	1006	LOP	C16-C17-C18-C19
8	K	1006	LOP	C15-C16-C17-C18
8	K	1006	LOP	C9-C10-C11-C12
8	O	1005	LOP	C16-C17-C18-C19
8	O	1005	LOP	C9-C10-C11-C12
8	O	1005	LOP	C30-C31-C32-C33
4	K	1001	HEM	C4C-C3C-CAC-CBC
8	K	1006	LOP	O8-C24-O6-C5
8	A	1006	LOP	C17-C18-C19-C20
8	A	1006	LOP	C24-C25-C26-C27
8	K	1006	LOP	C11-C12-C13-C14
6	E	1004	ANJ	C13-C18-C19-C20
8	K	1006	LOP	C3-C4-C5-O6
8	E	1005	LOP	C15-C16-C17-C18
8	O	1005	LOP	C11-C12-C13-C14
8	K	1006	LOP	C27-C28-C29-C30
4	K	1002	HEM	C1A-C2A-CAA-CBA
8	E	1005	LOP	C13-C14-C15-C16
8	O	1005	LOP	C13-C14-C15-C16
8	A	1006	LOP	C20-C21-C22-C23
8	A	1006	LOP	C32-C33-C34-C35
6	K	1004	ANJ	C19-C20-C21-C22
6	A	1004	ANJ	O8-C24-C25-C26
8	A	1006	LOP	C28-C29-C30-C31
8	E	1005	LOP	C30-C31-C32-C33
8	E	1005	LOP	C18-C19-C20-C21
9	P	1001	HEC	C4B-C3B-CAB-CBB
4	K	1001	HEM	C3D-CAD-CBD-CGD
8	O	1005	LOP	C18-C19-C20-C21
8	E	1005	LOP	C6-C7-C8-C9
9	B	1001	HEC	C4B-C3B-CAB-CBB
8	A	1006	LOP	O2-C3-C4-O5
8	A	1006	LOP	O5-C4-C5-O6
8	K	1006	LOP	C28-C29-C30-C31

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Mol	Chain	Res	Type	Atoms
5	K	1003	SMA	C9-C10-C11-C22
6	O	1004	ANJ	O9-C24-C25-C26
6	E	1004	ANJ	O8-C24-C25-C26
8	K	1006	LOP	O2-C3-C4-O5
4	K	1001	HEM	C4B-C3B-CAB-CBB
4	O	1001	HEM	C4B-C3B-CAB-CBB
8	A	1006	LOP	C9-C10-C11-C12
8	A	1006	LOP	C11-C12-C13-C14
8	A	1006	LOP	C27-C28-C29-C30
5	O	1003	SMA	C9-C10-C11-C22
6	K	1004	ANJ	O9-C24-C25-C26
6	O	1004	ANJ	O5-C12-O4-C10
8	A	1006	LOP	C18-C19-C20-C21
8	A	1006	LOP	C2-O1-P1-O3
8	K	1006	LOP	C3-O2-P1-O1
8	O	1005	LOP	C2-O1-P1-O4
8	E	1005	LOP	O2-C3-C4-C5
8	K	1006	LOP	C12-C13-C14-C15
8	A	1006	LOP	C29-C30-C31-C32
8	A	1006	LOP	C12-C13-C14-C15
5	A	1003	SMA	C26-C19-C20-C21
4	O	1001	HEM	C3D-CAD-CBD-CGD
8	O	1005	LOP	C28-C29-C30-C31
8	E	1005	LOP	C28-C29-C30-C31
6	K	1004	ANJ	C26-C25-C27-C28
8	K	1006	LOP	C30-C31-C32-C33
4	A	1001	HEM	CAD-CBD-CGD-O1D
4	O	1001	HEM	CAA-CBA-CGA-O1A
8	K	1006	LOP	O7-C6-O5-C4
4	K	1001	HEM	CAA-CBA-CGA-O1A
8	O	1005	LOP	C7-C8-C9-C10
4	K	1002	HEM	C3A-C2A-CAA-CBA
4	O	1001	HEM	CAA-CBA-CGA-O2A
4	A	1001	HEM	CAD-CBD-CGD-O2D
8	E	1005	LOP	O6-C24-C25-C26
8	E	1005	LOP	C3-C4-C5-O6
4	A	1001	HEM	CAA-CBA-CGA-O2A
4	K	1001	HEM	CAA-CBA-CGA-O2A
8	O	1005	LOP	O6-C24-C25-C26
8	O	1005	LOP	O2-C3-C4-O5
4	A	1002	HEM	CAD-CBD-CGD-O2D
4	K	1001	HEM	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
8	E	1005	LOP	C25-C26-C27-C28
8	K	1006	LOP	O2-C3-C4-C5
4	E	1001	HEM	CAD-CBD-CGD-O2D
9	B	1001	HEC	CAA-CBA-CGA-O2A
8	O	1005	LOP	C25-C26-C27-C28
8	E	1005	LOP	C14-C15-C16-C17
8	O	1005	LOP	C12-C13-C14-C15
8	K	1006	LOP	C25-C24-O6-C5
4	A	1001	HEM	CAA-CBA-CGA-O1A
9	F	1001	HEC	CAA-CBA-CGA-O2A
8	O	1005	LOP	C29-C30-C31-C32
5	K	1003	SMA	C15-C14-O14-C25
8	E	1005	LOP	C12-C13-C14-C15
4	A	1002	HEM	CAD-CBD-CGD-O1D
4	K	1001	HEM	CAD-CBD-CGD-O1D
4	O	1001	HEM	CAD-CBD-CGD-O2D
8	E	1005	LOP	C26-C27-C28-C29
8	A	1006	LOP	C26-C27-C28-C29
6	E	1004	ANJ	C9-C10-O4-C12
4	E	1001	HEM	CAD-CBD-CGD-O1D
8	A	1006	LOP	C7-C8-C9-C10
6	A	1004	ANJ	C19-C20-C21-C22
9	L	1001	HEC	CAA-CBA-CGA-O2A
9	F	1001	HEC	CAA-CBA-CGA-O1A
4	O	1001	HEM	CAD-CBD-CGD-O1D
9	B	1001	HEC	CAA-CBA-CGA-O1A
8	K	1006	LOP	O5-C6-C7-C8
9	L	1001	HEC	CAA-CBA-CGA-O1A
4	O	1002	HEM	CAA-CBA-CGA-O1A
8	A	1006	LOP	C25-C26-C27-C28
5	E	1003	SMA	C13-C14-C15-C16
4	E	1001	HEM	C3D-CAD-CBD-CGD
8	A	1006	LOP	C7-C6-O5-C4
9	P	1001	HEC	CAA-CBA-CGA-O2A
8	O	1005	LOP	O5-C6-C7-C8
4	E	1001	HEM	CAA-CBA-CGA-O1A
4	E	1001	HEM	CAA-CBA-CGA-O2A
4	E	1002	HEM	CAA-CBA-CGA-O1A
6	O	1004	ANJ	C18-C19-C20-C21
5	O	1003	SMA	C11-C10-C9-C2
6	O	1004	ANJ	O8-C14-C15-C16
8	A	1006	LOP	O5-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
4	E	1002	HEM	CAA-CBA-CGA-O2A
4	A	1002	HEM	C1A-C2A-CAA-CBA
9	P	1001	HEC	CAA-CBA-CGA-O1A
4	O	1002	HEM	CAA-CBA-CGA-O2A
9	L	1001	HEC	CAD-CBD-CGD-O2D
6	O	1004	ANJ	O8-C24-C25-C26
5	A	1003	SMA	C11-C10-C9-C2
8	O	1005	LOP	C26-C27-C28-C29
6	A	1004	ANJ	C12-C13-C18-C19

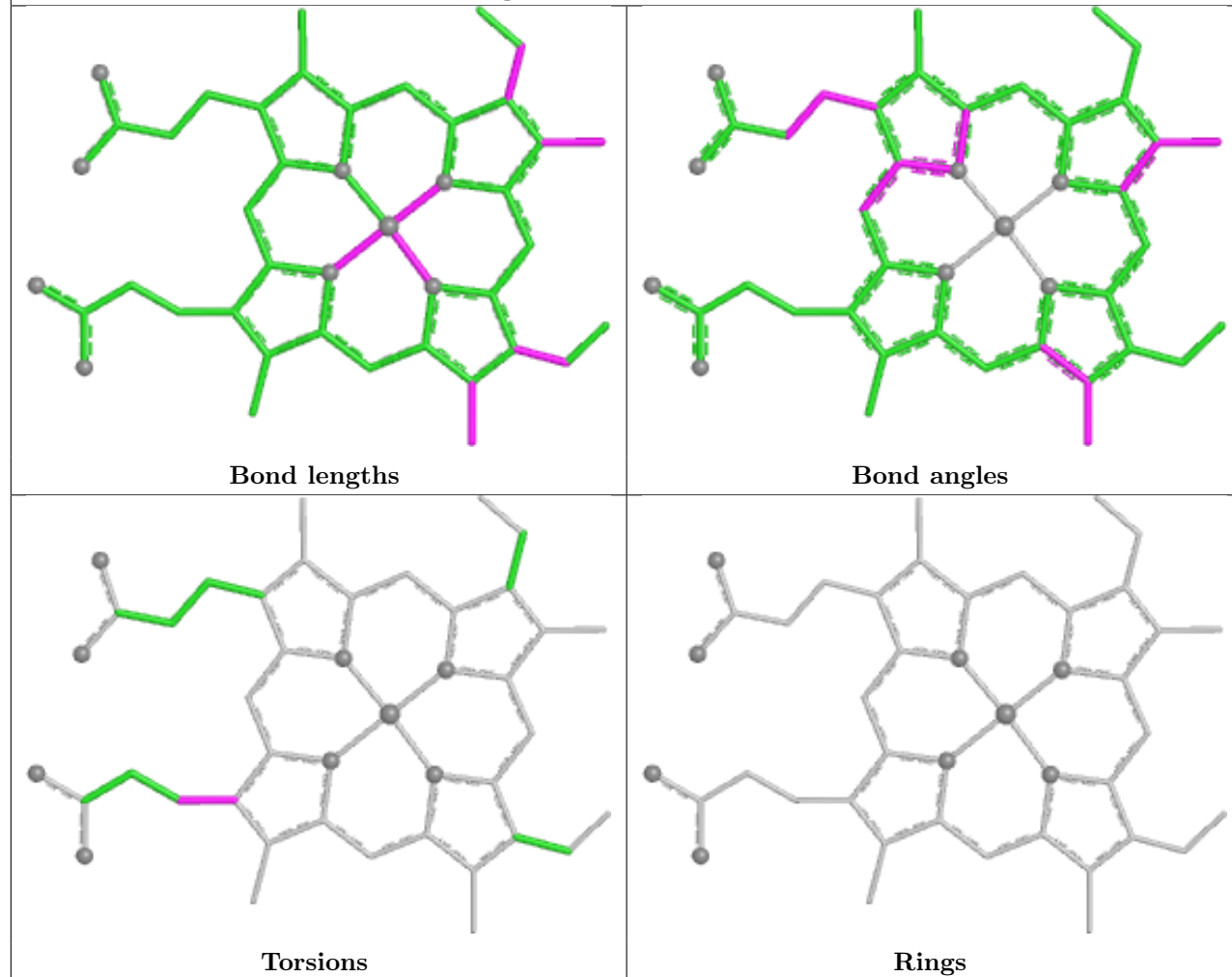
There are no ring outliers.

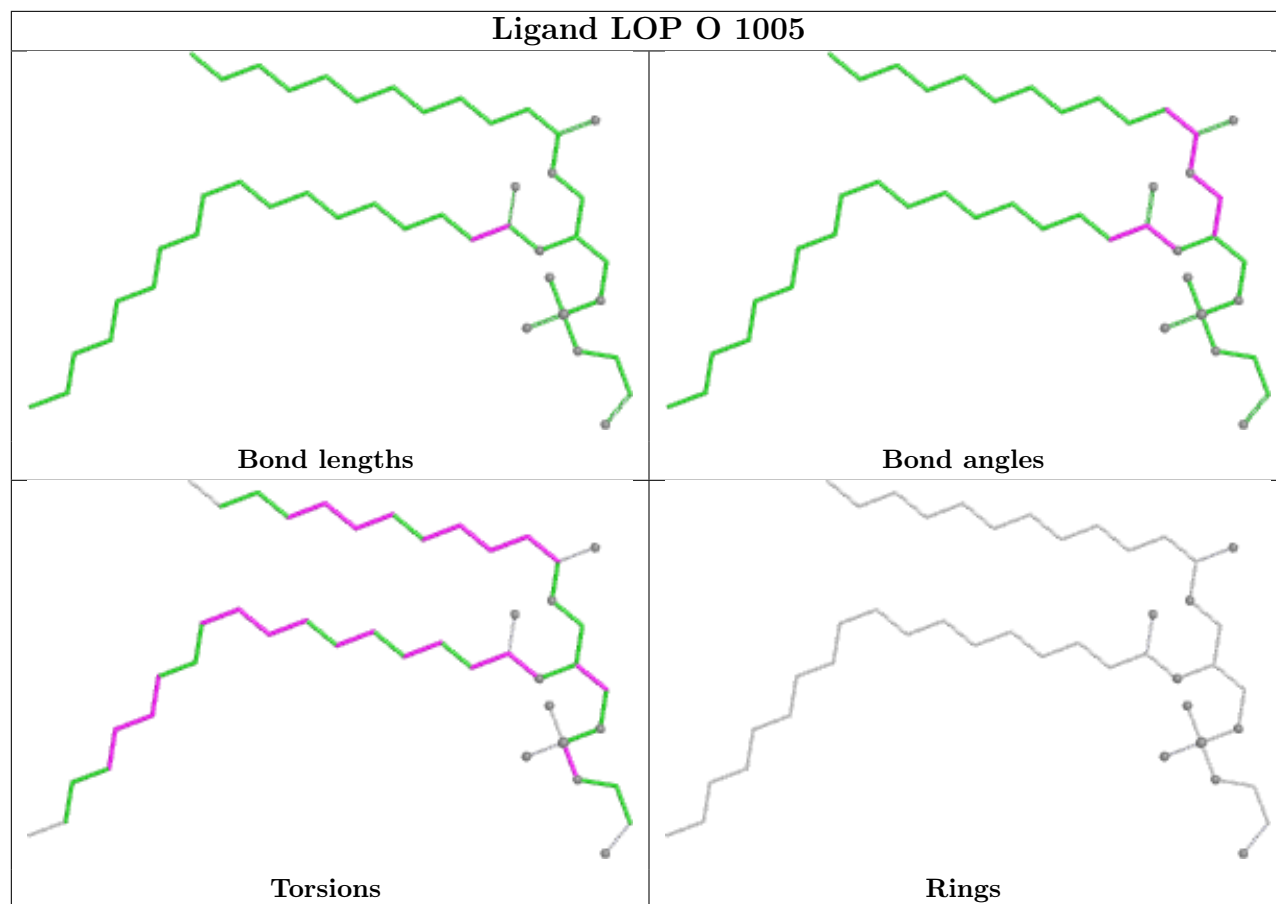
14 monomers are involved in 19 short contacts:

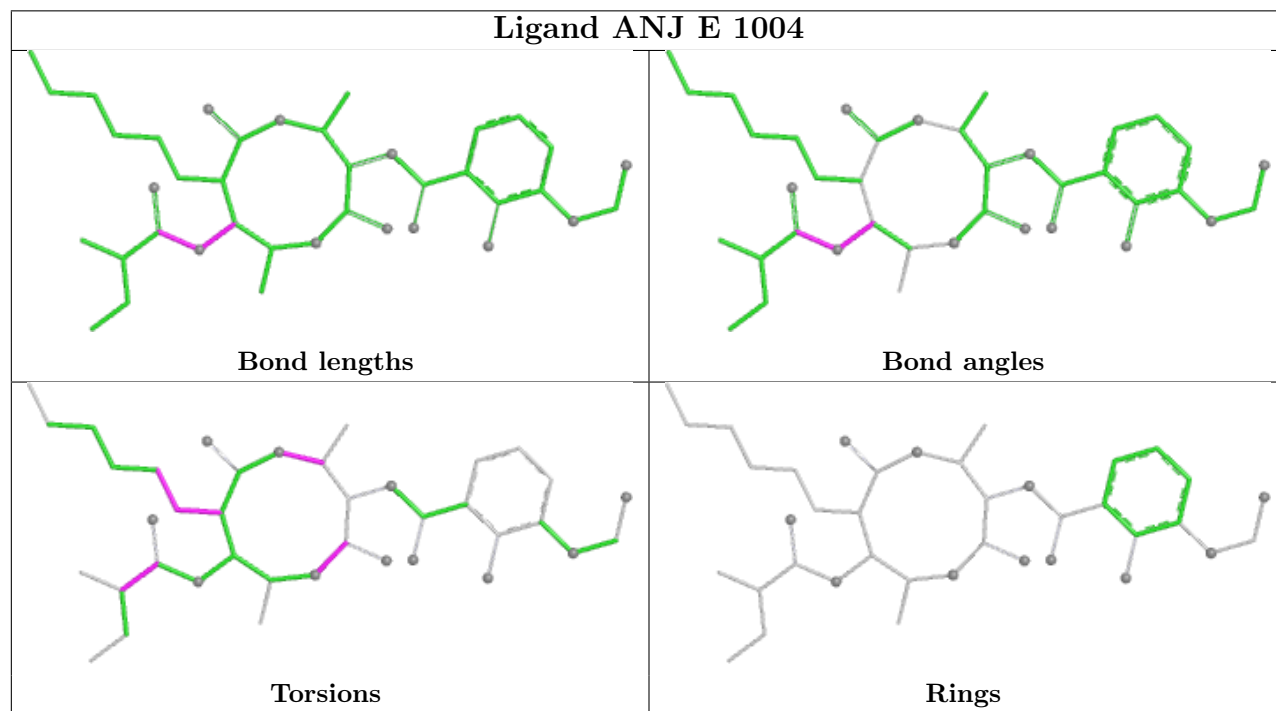
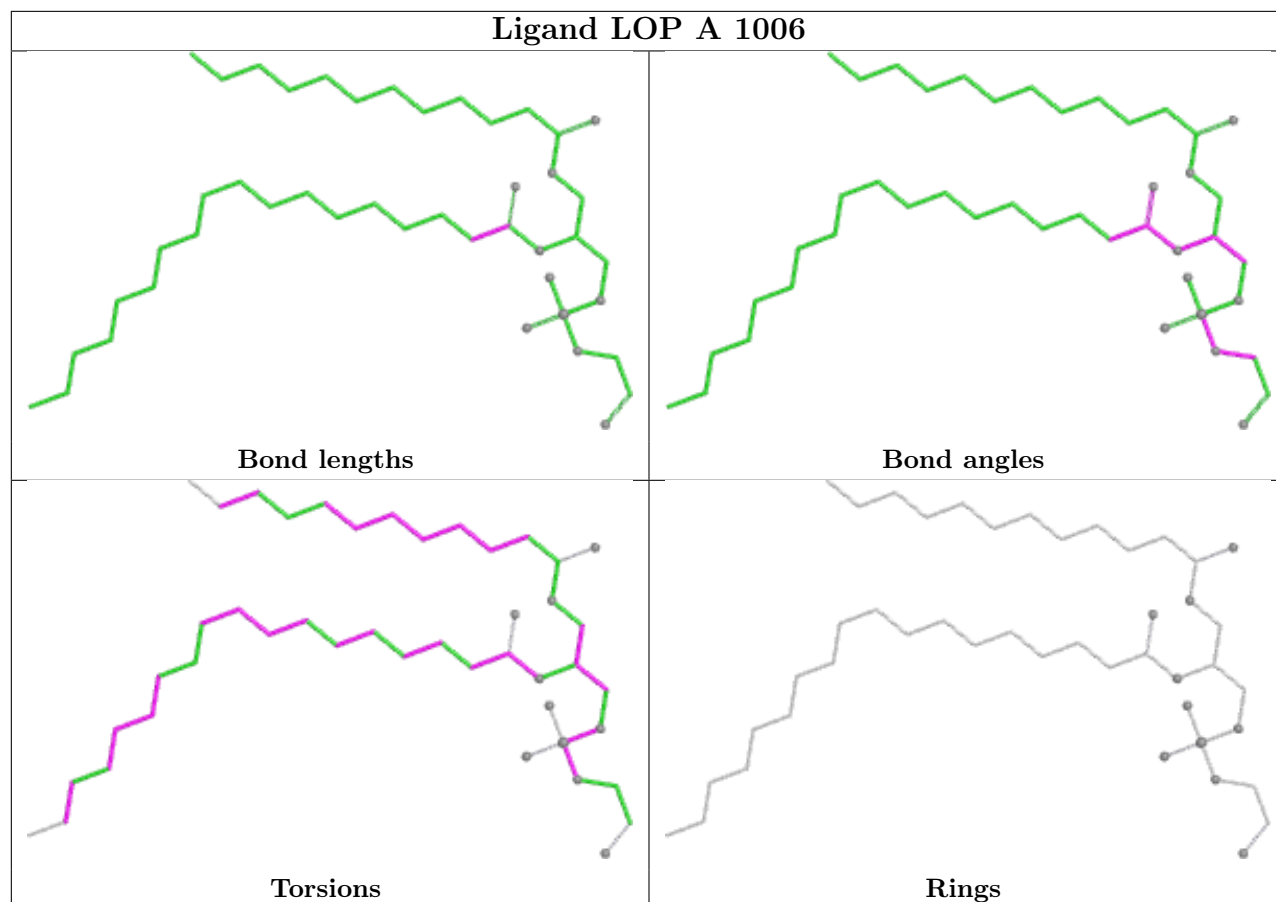
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	K	1002	HEM	1	0
6	E	1004	ANJ	3	0
4	E	1001	HEM	2	0
9	B	1001	HEC	1	0
4	A	1002	HEM	1	0
9	L	1001	HEC	1	0
4	K	1001	HEM	2	0
4	A	1001	HEM	1	0
4	O	1001	HEM	2	0
6	K	1004	ANJ	1	0
9	F	1001	HEC	1	0
6	O	1004	ANJ	1	0
8	K	1006	LOP	1	0
9	P	1001	HEC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

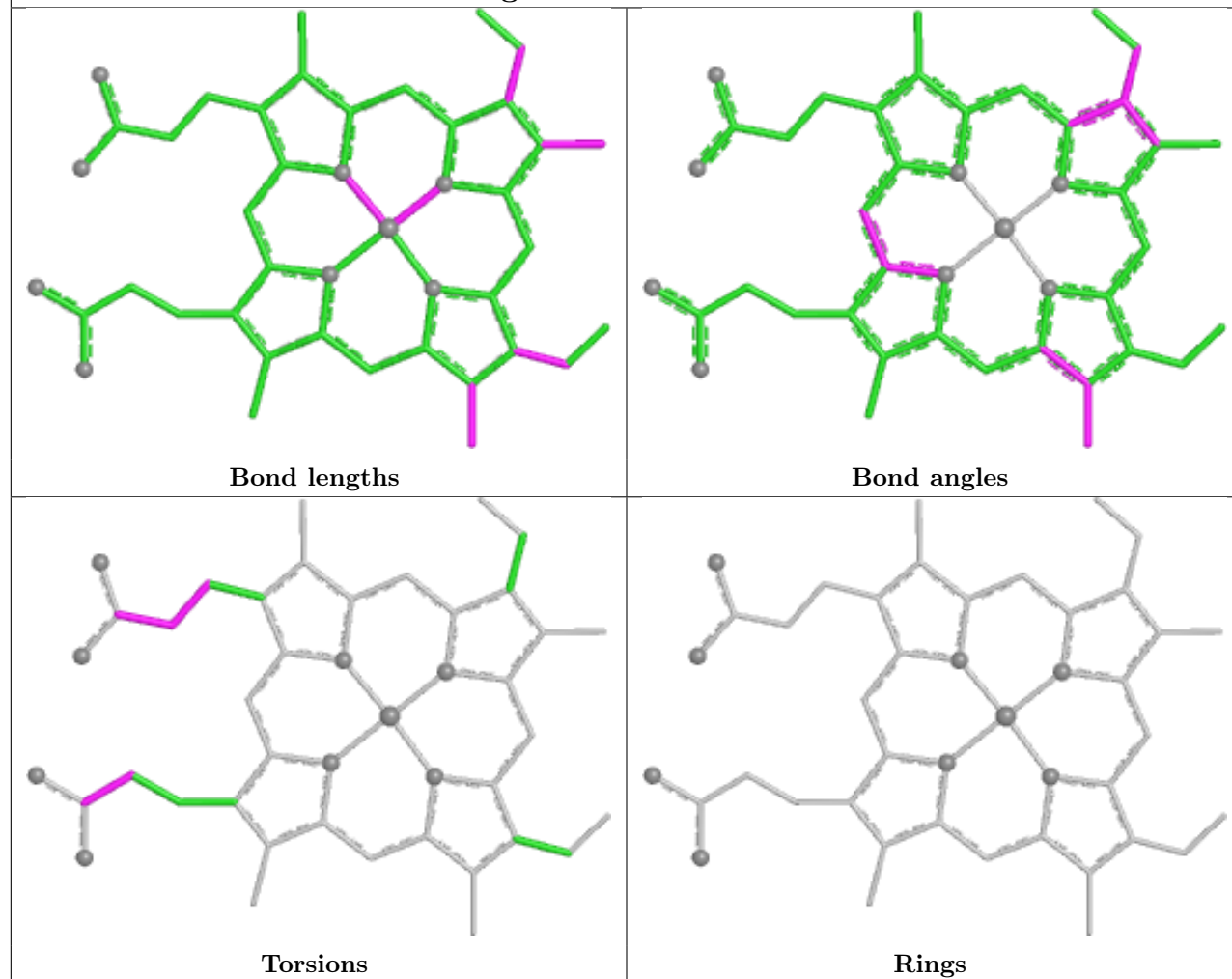
Ligand HEM K 1002



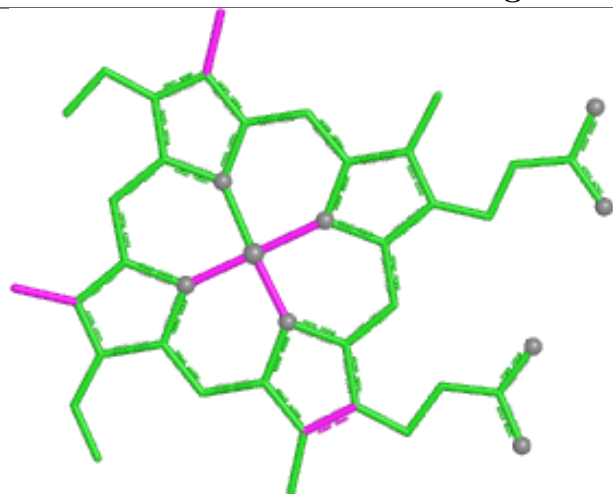




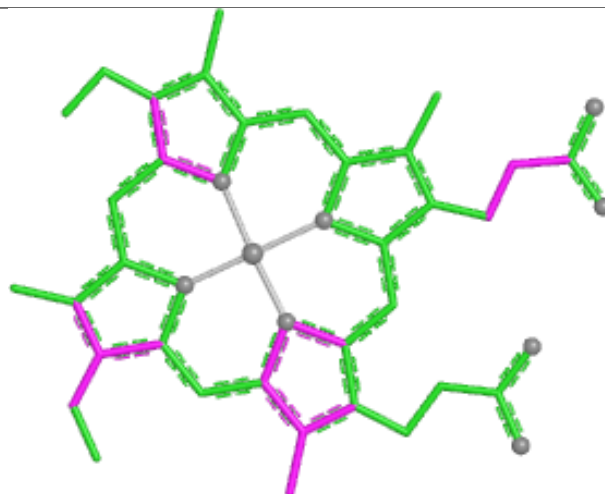
Ligand HEM E 1001



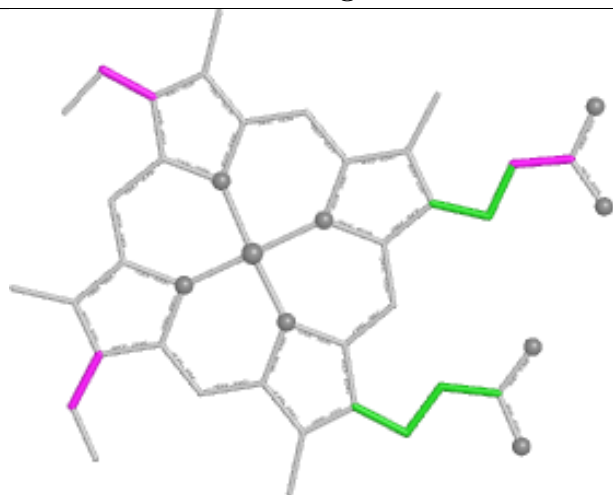
Ligand HEC B 1001



Bond lengths



Bond angles

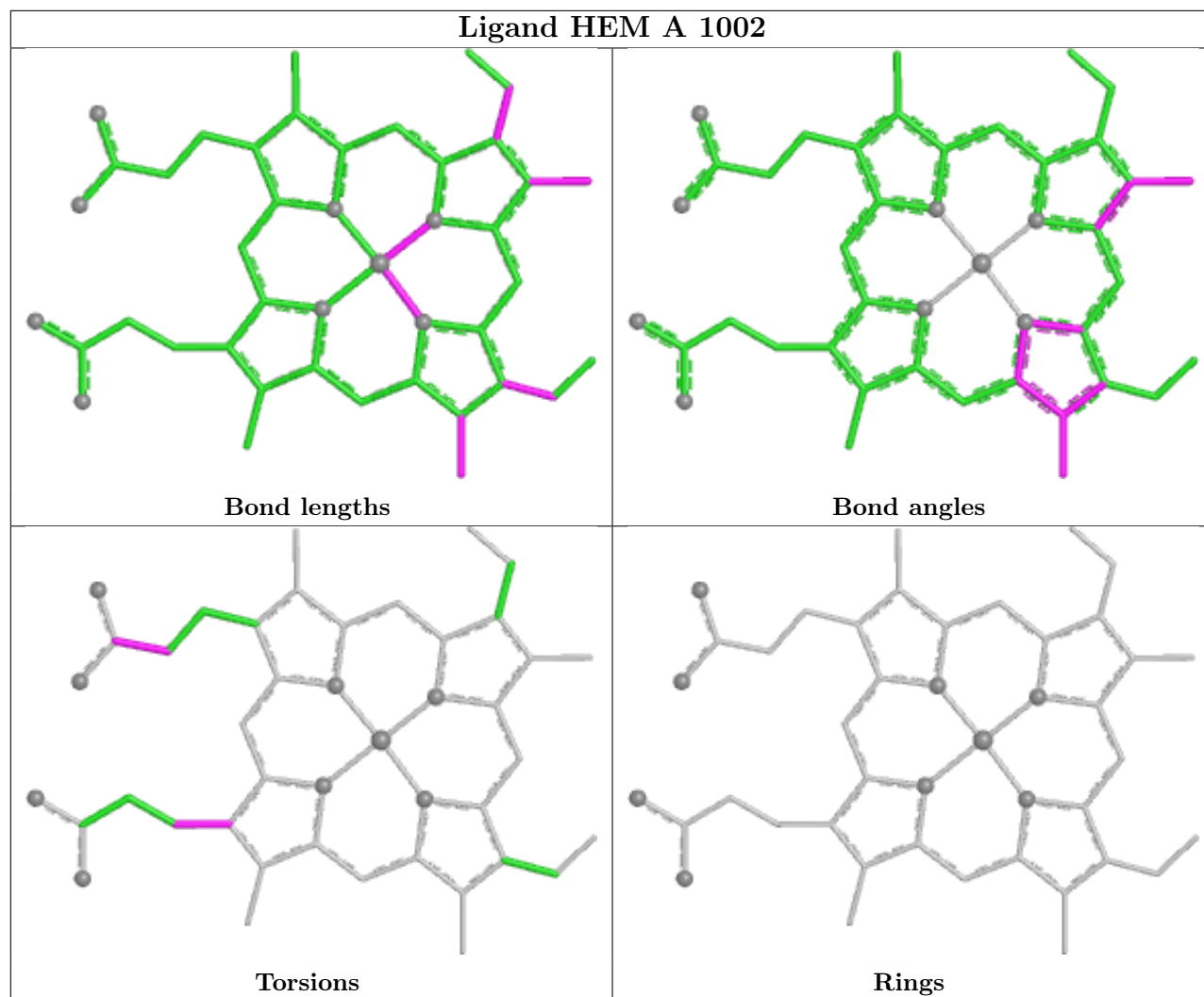


Torsions

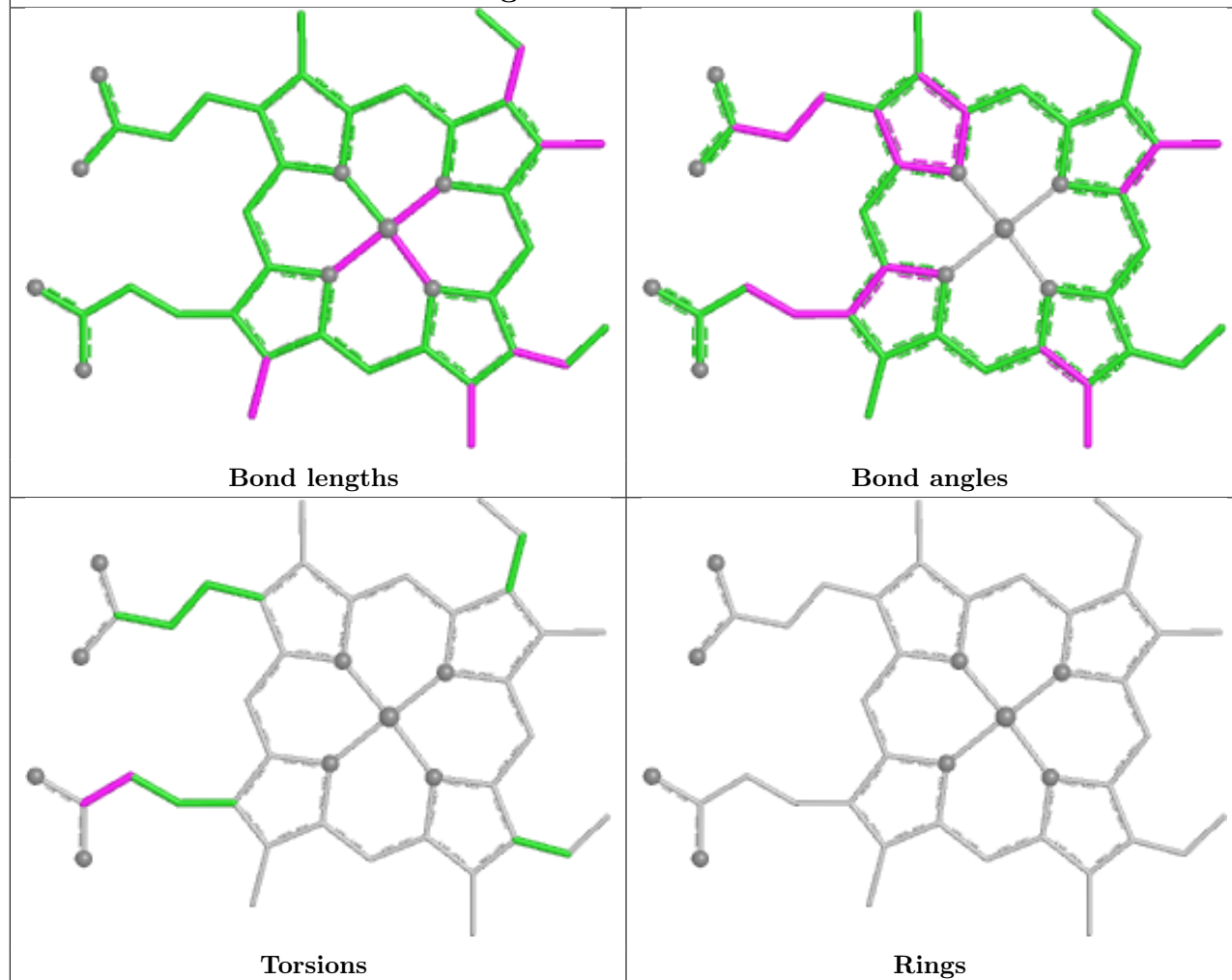


Rings

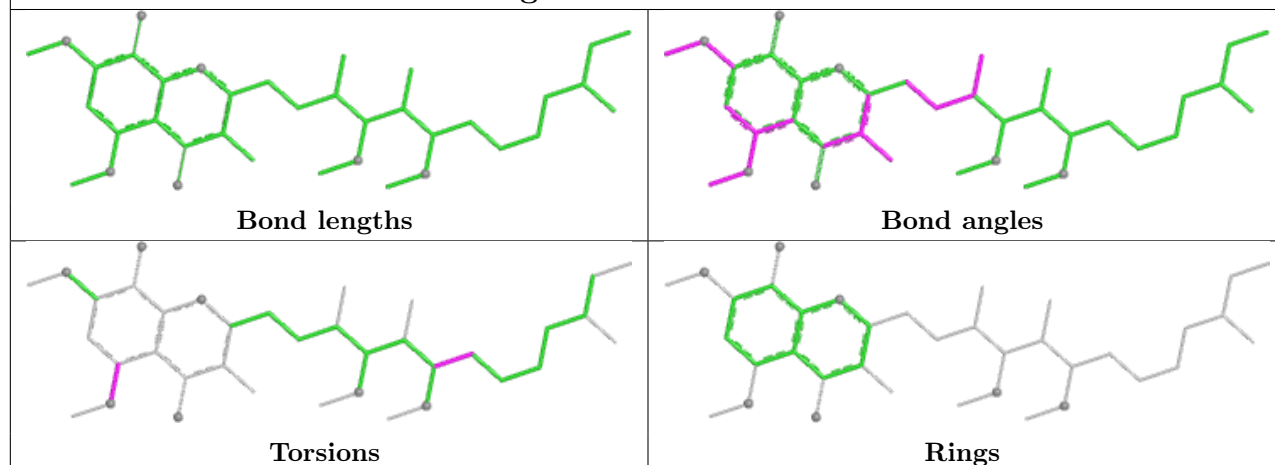
Ligand HEM A 1002



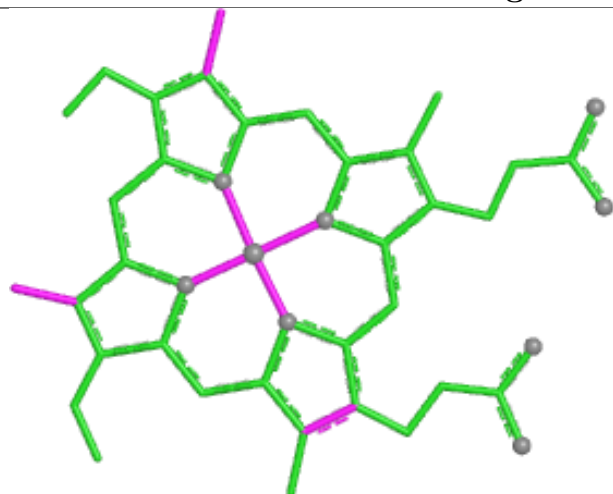
Ligand HEM E 1002



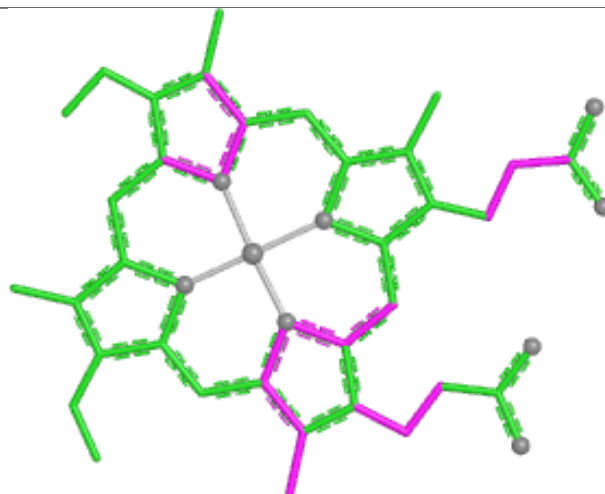
Ligand SMA E 1003



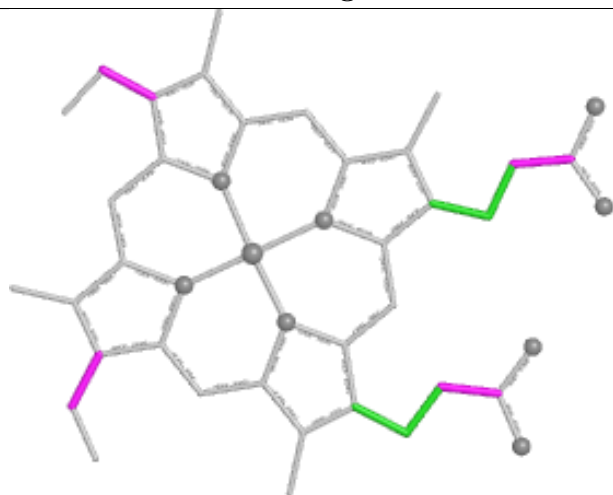
Ligand HEC L 1001



Bond lengths



Bond angles

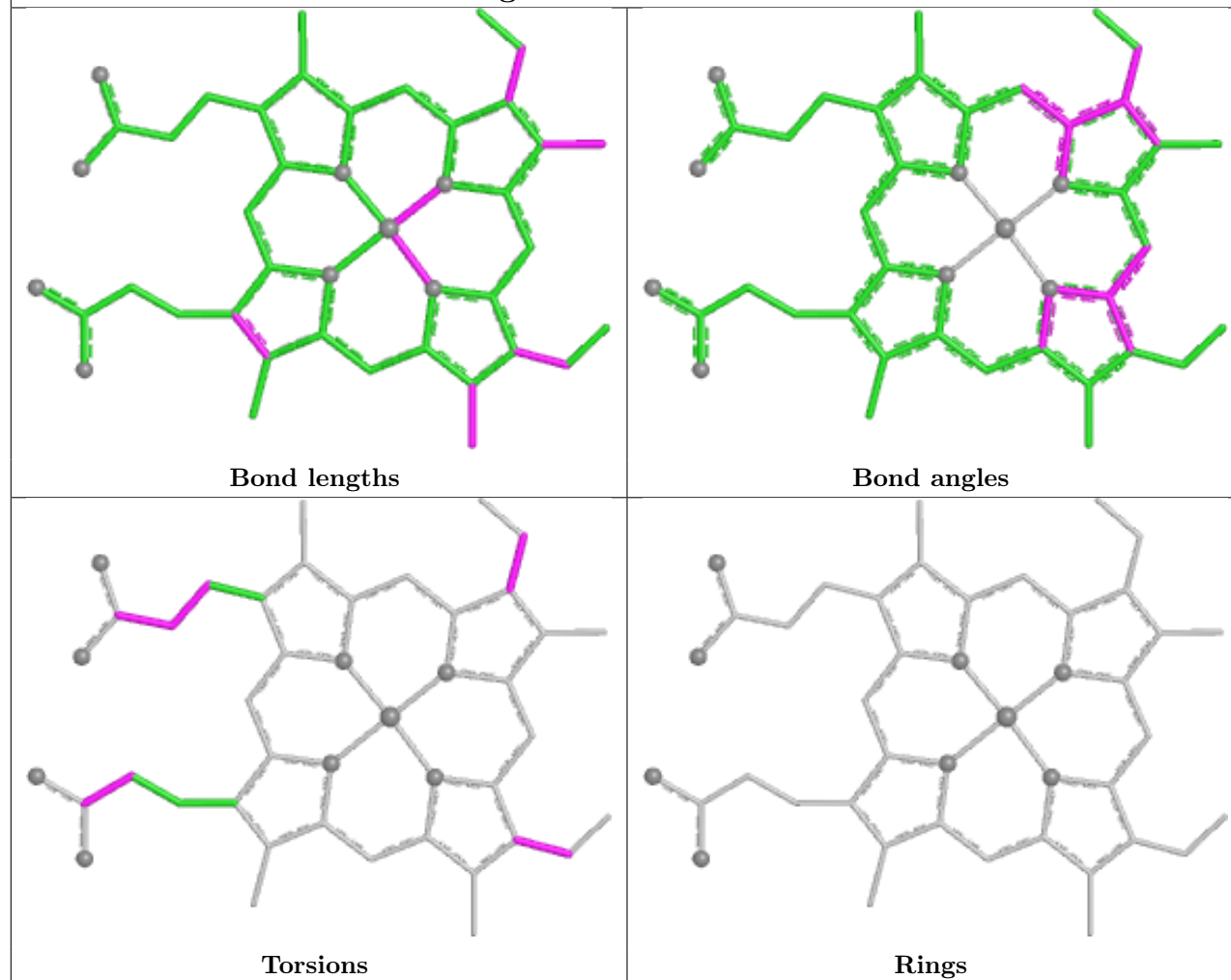


Torsions

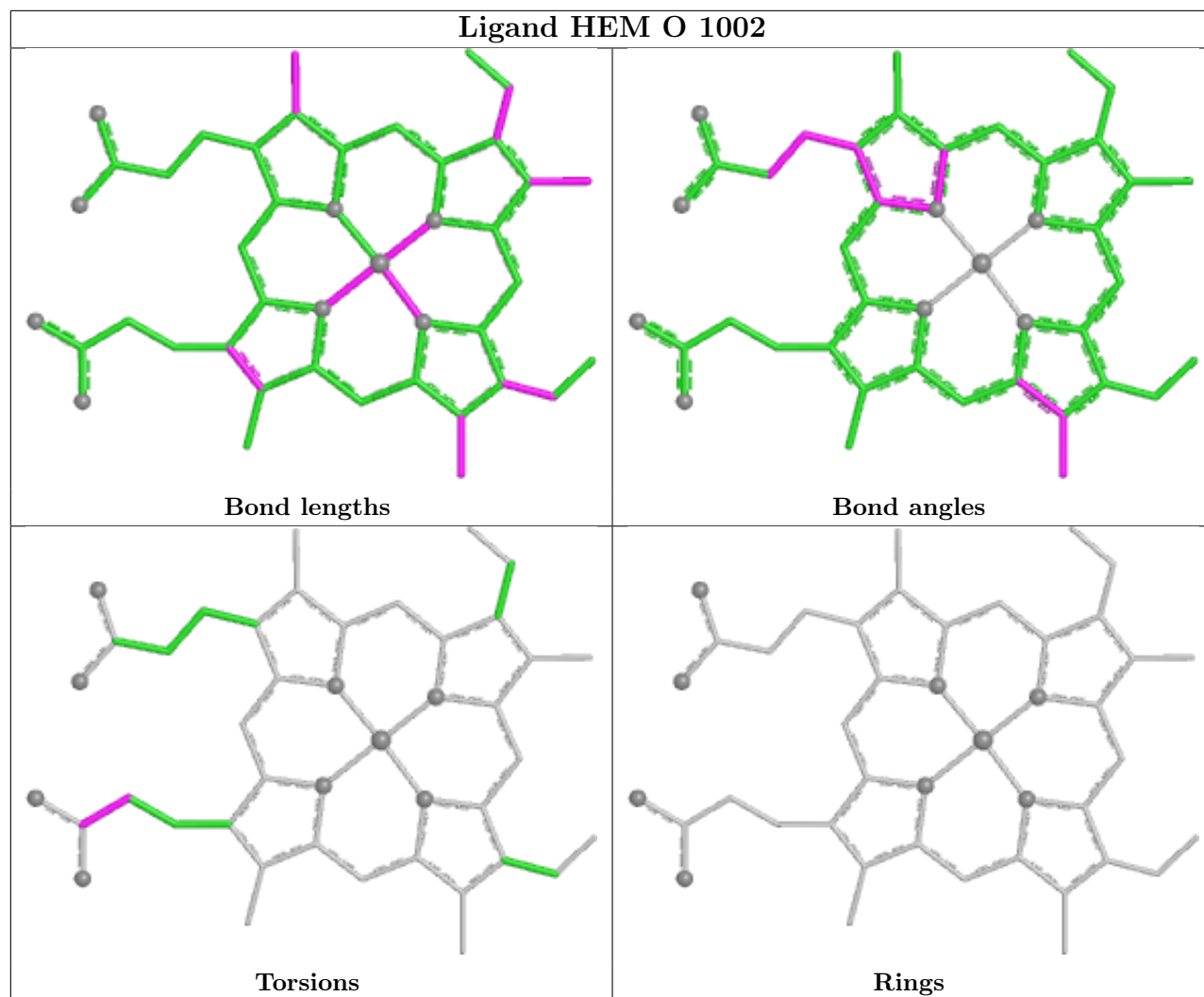


Rings

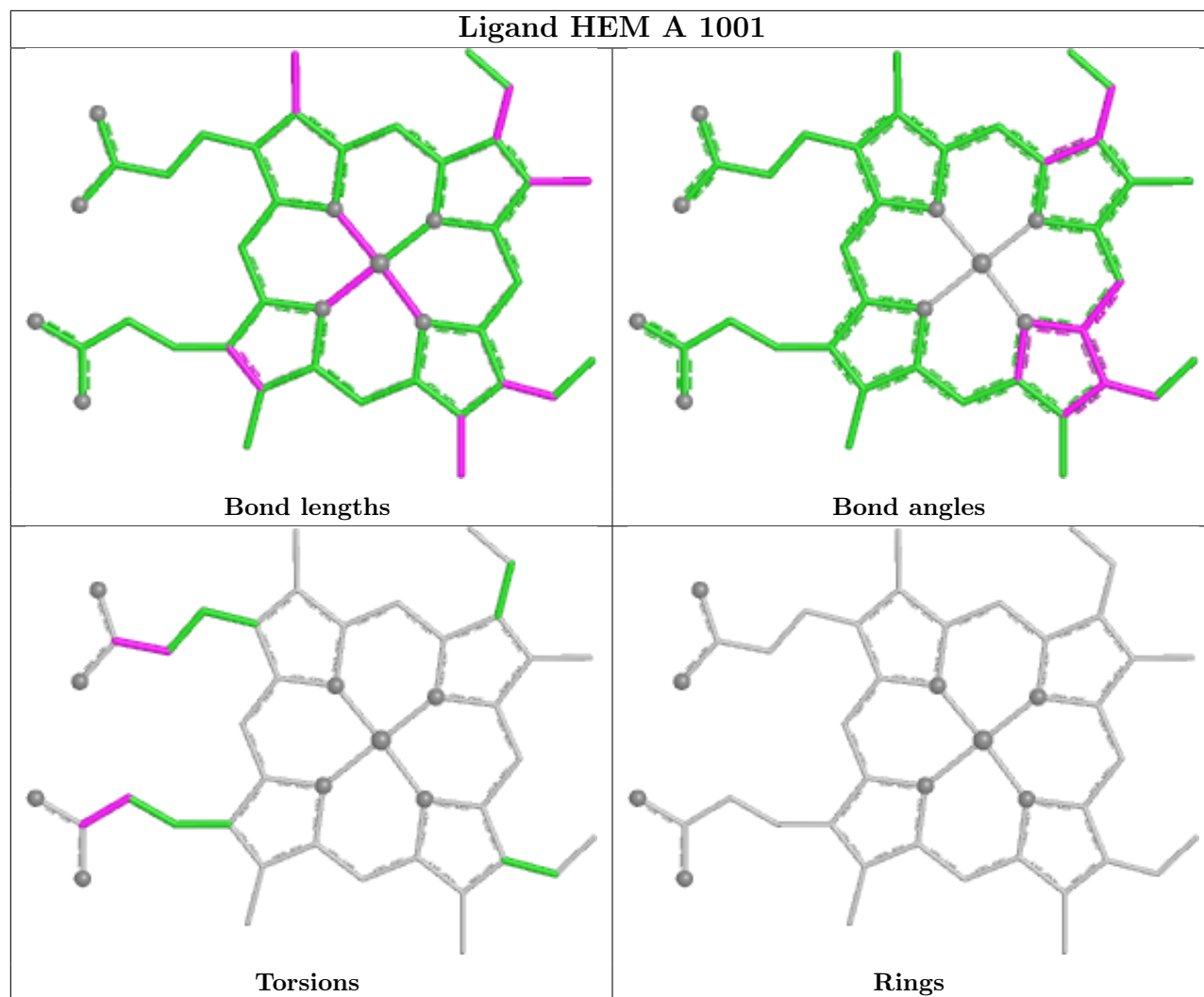
Ligand HEM K 1001



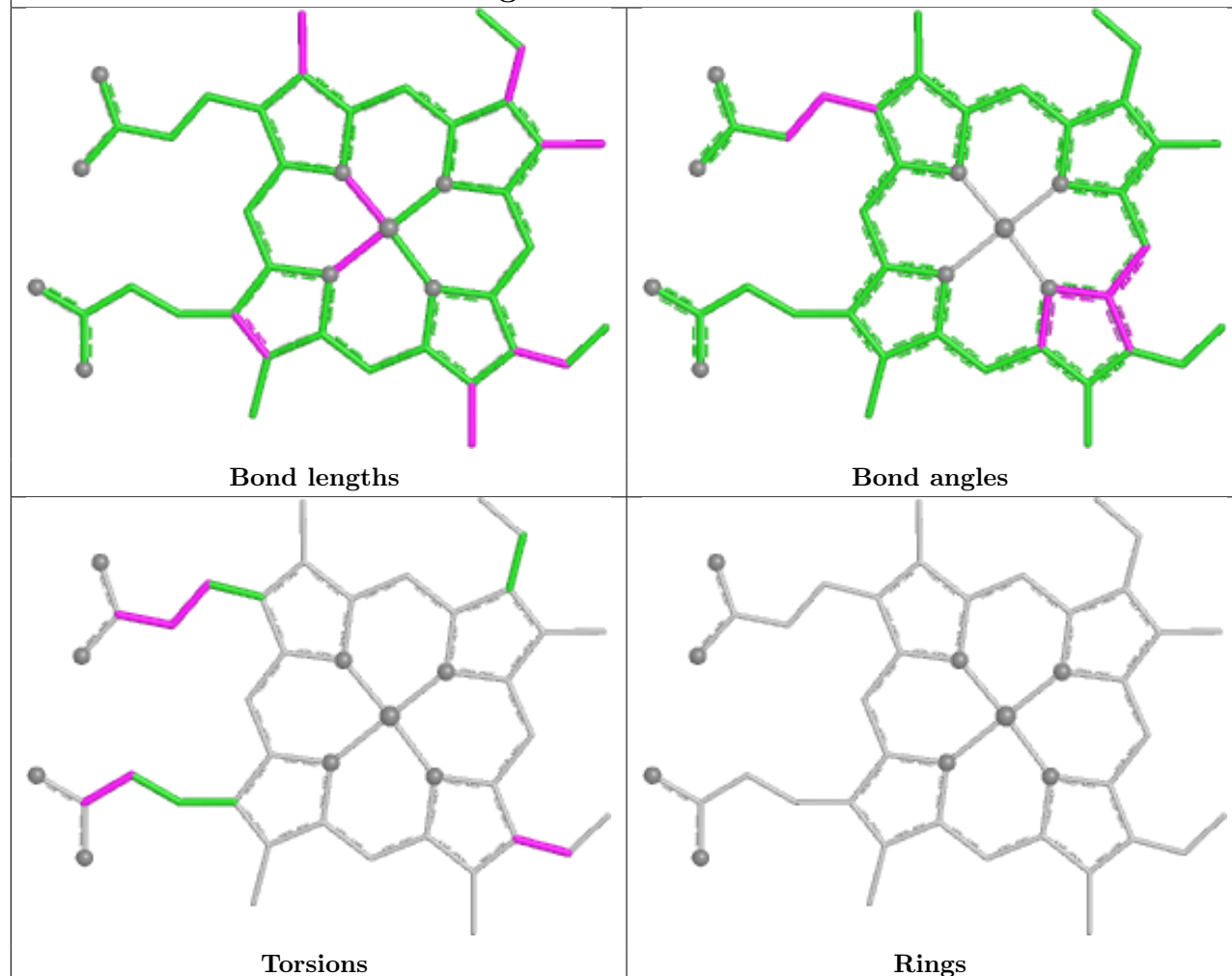
Ligand HEM O 1002



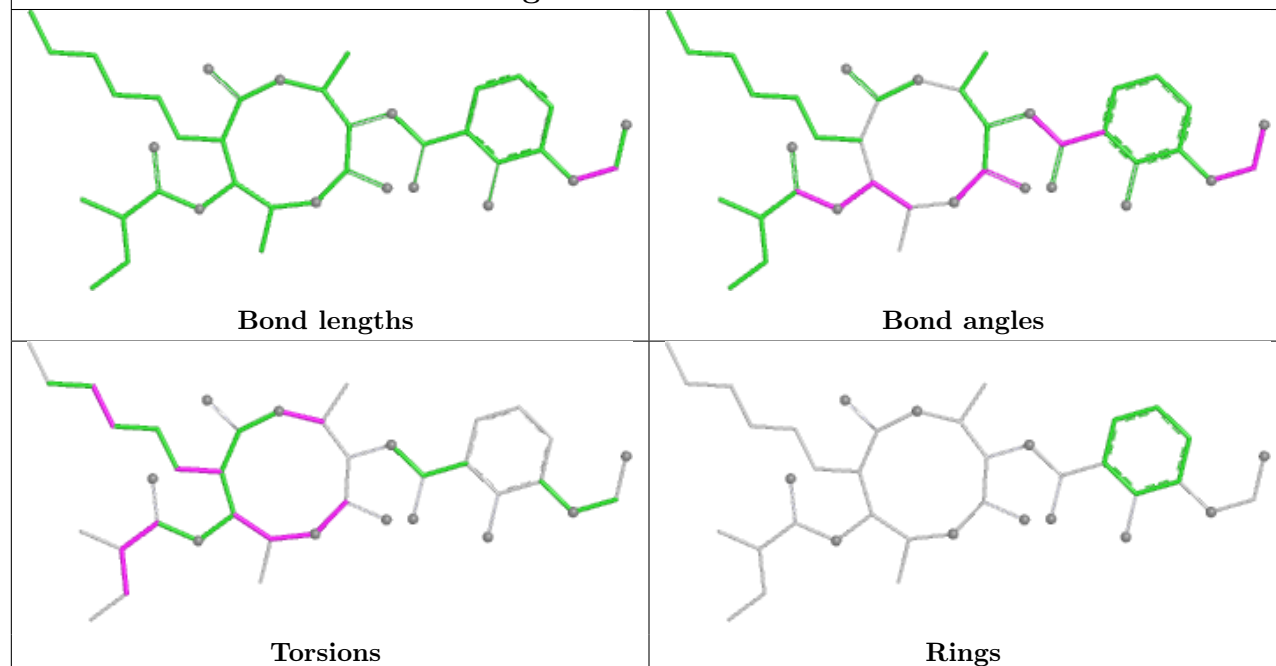
Ligand HEM A 1001



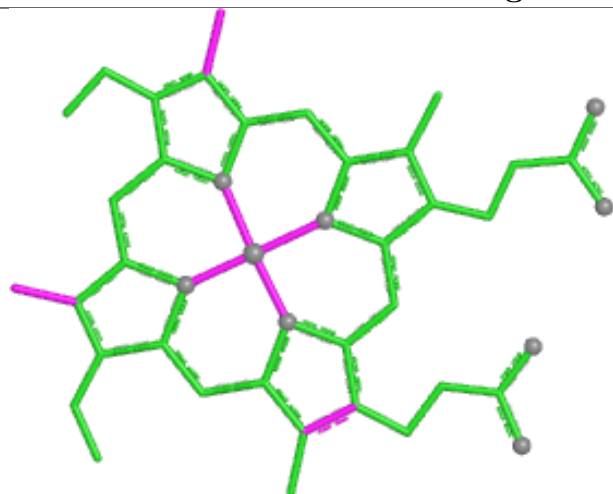
Ligand HEM O 1001



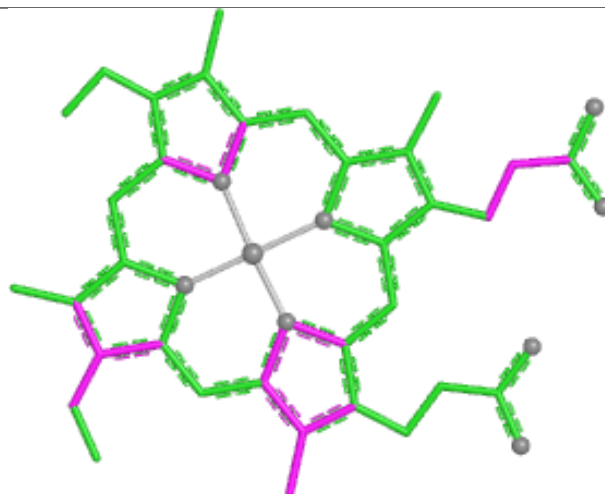
Ligand ANJ K 1004



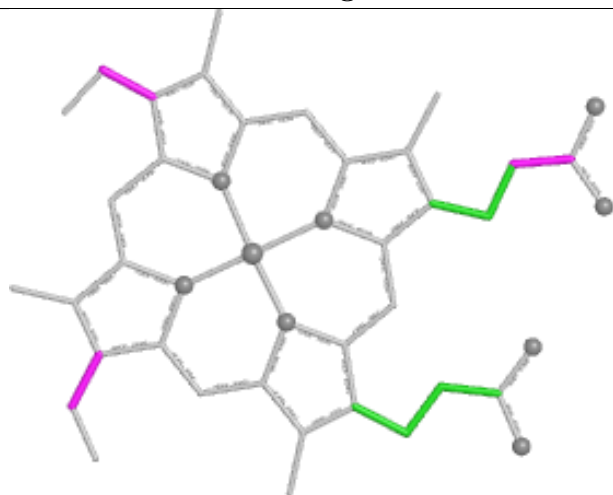
Ligand HEC F 1001



Bond lengths



Bond angles

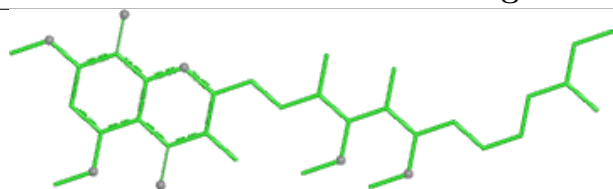


Torsions

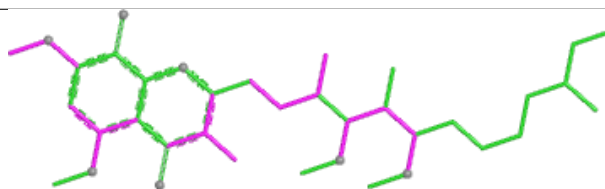


Rings

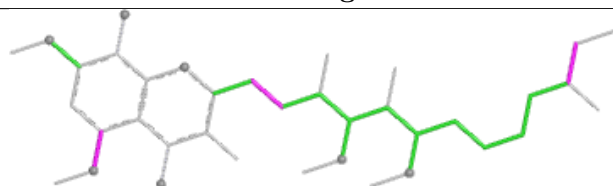
Ligand SMA A 1003



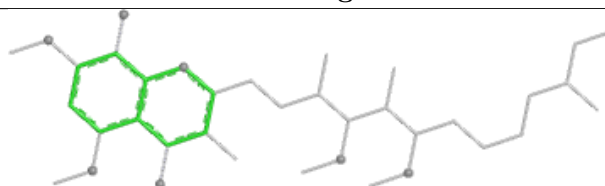
Bond lengths



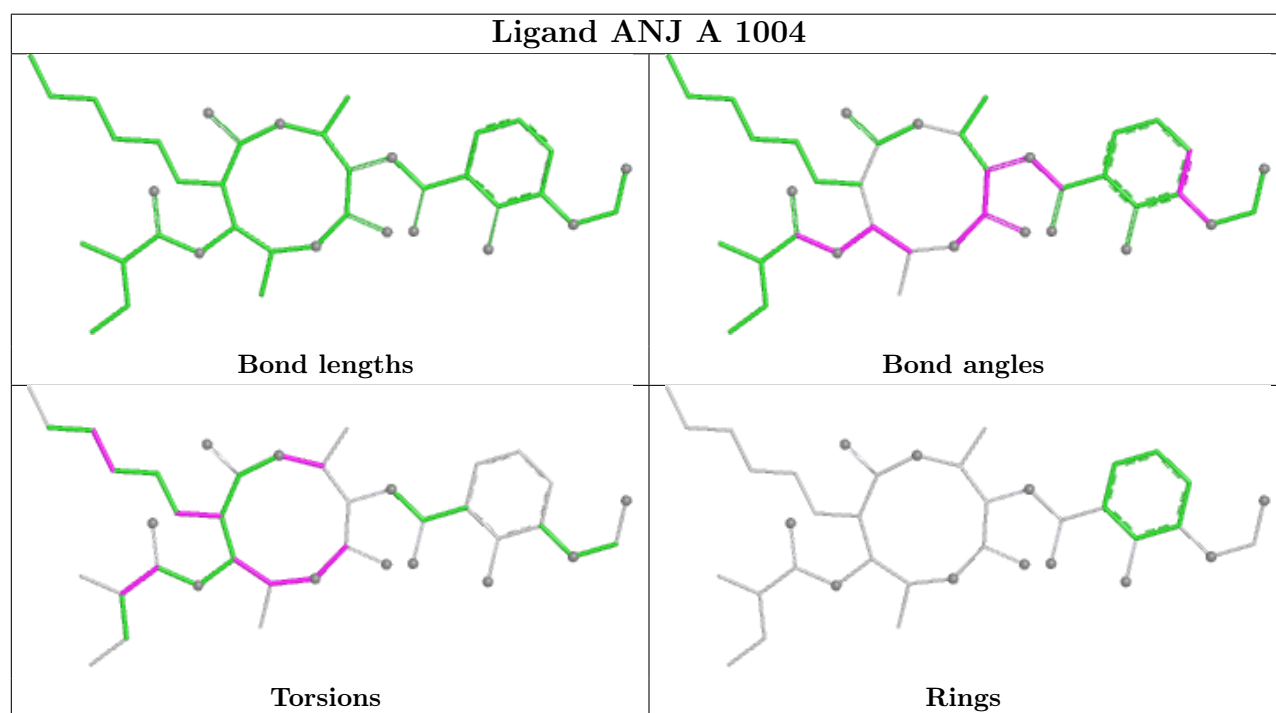
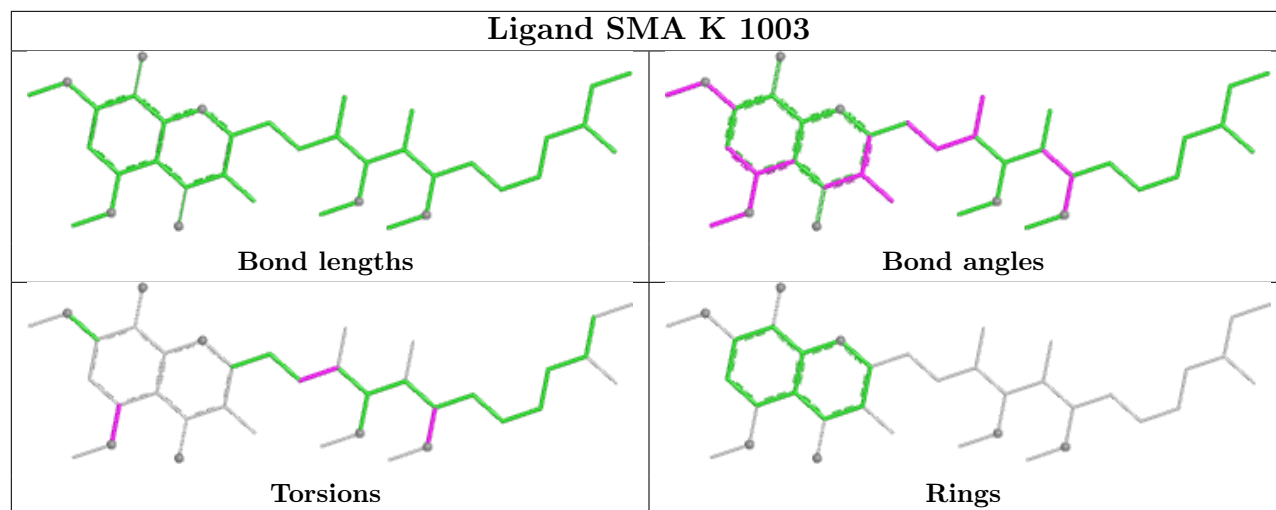
Bond angles

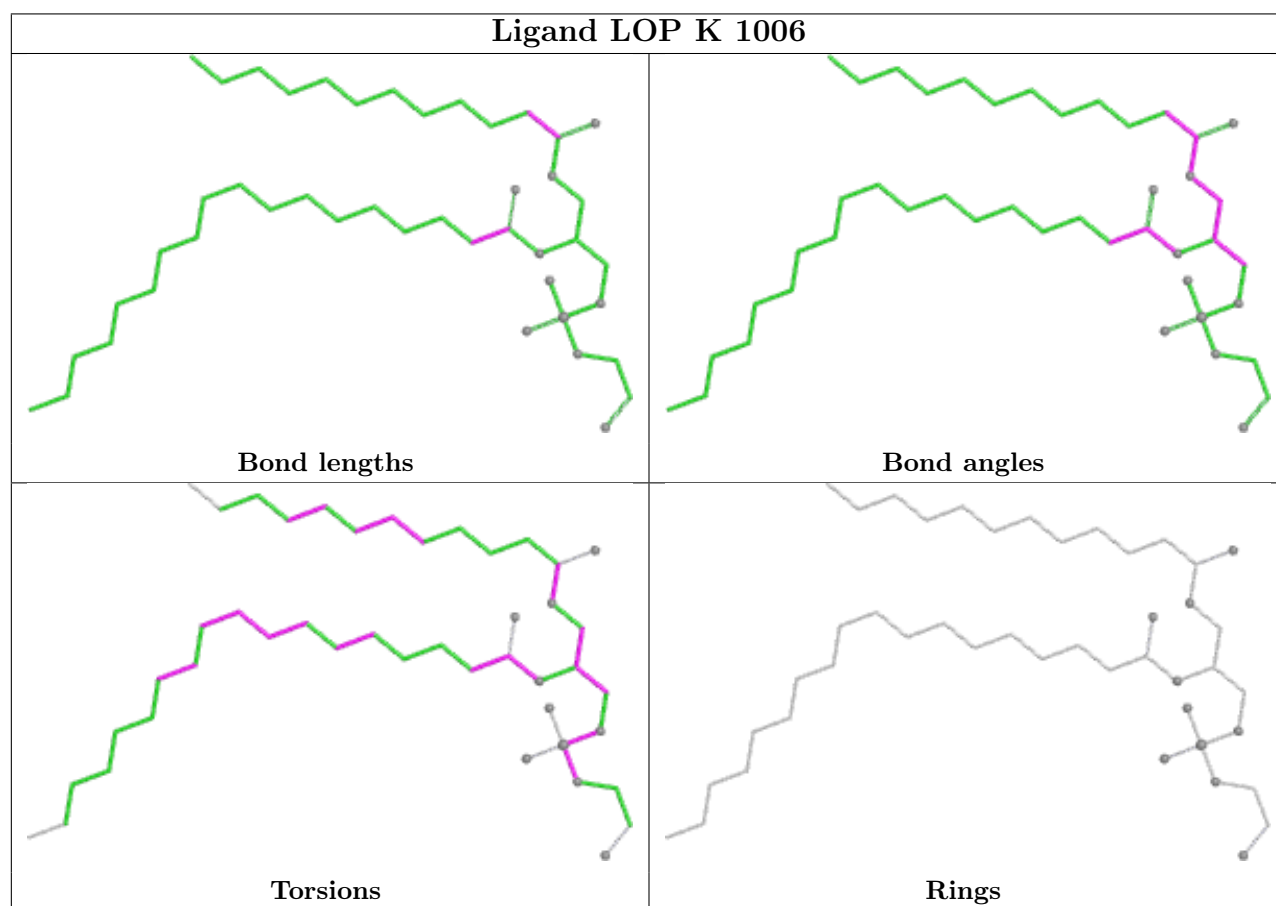
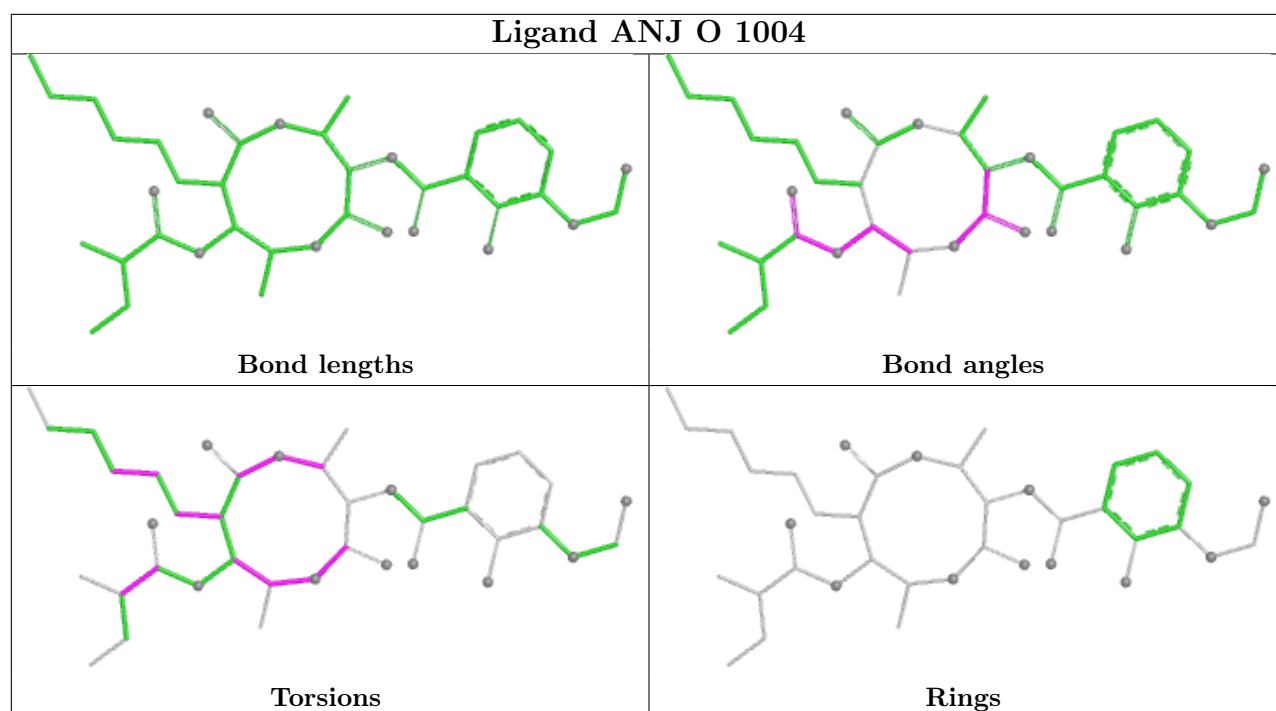


Torsions

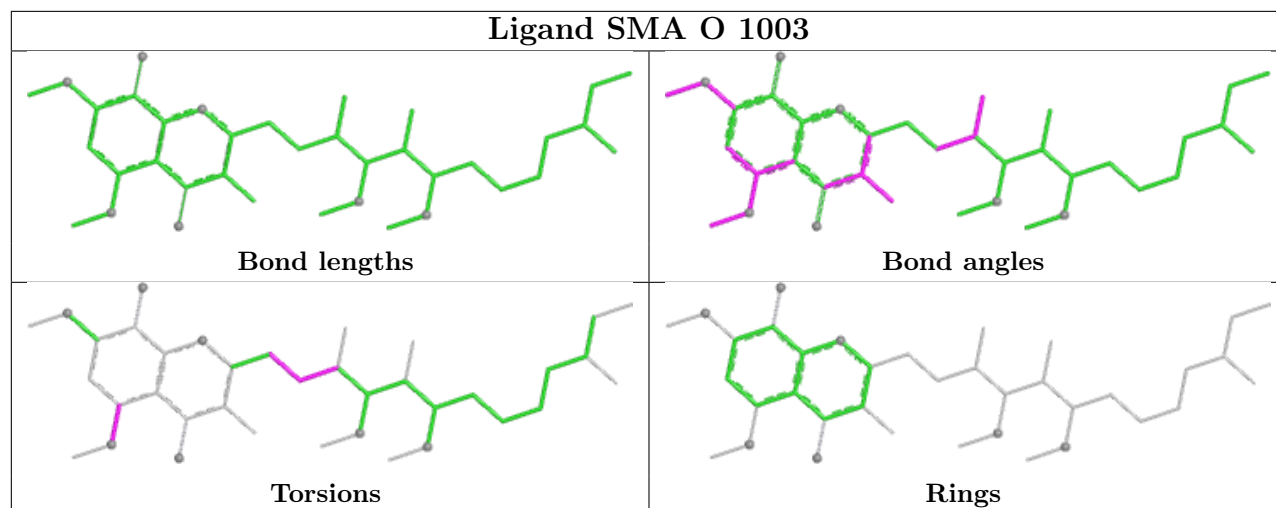


Rings

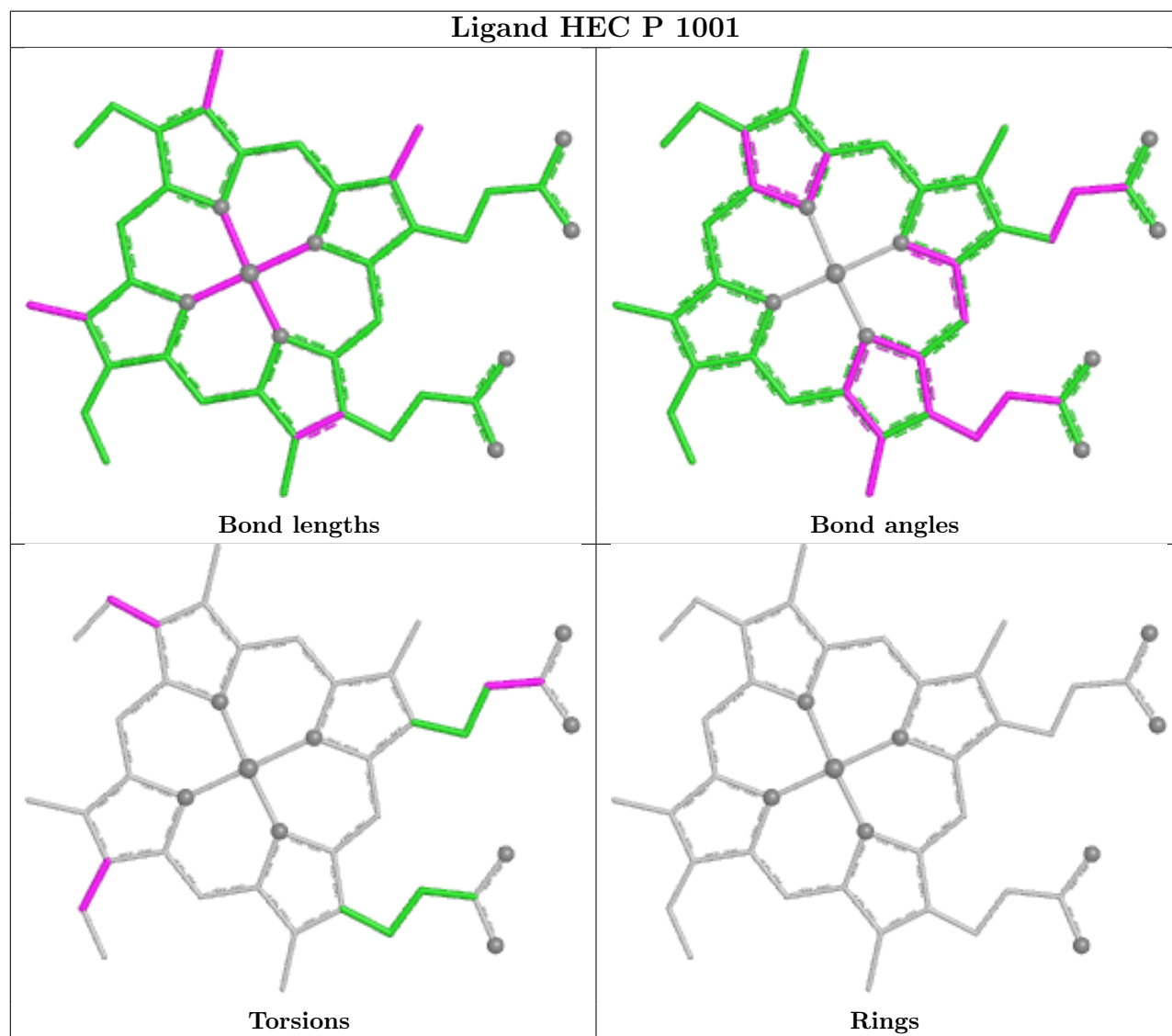


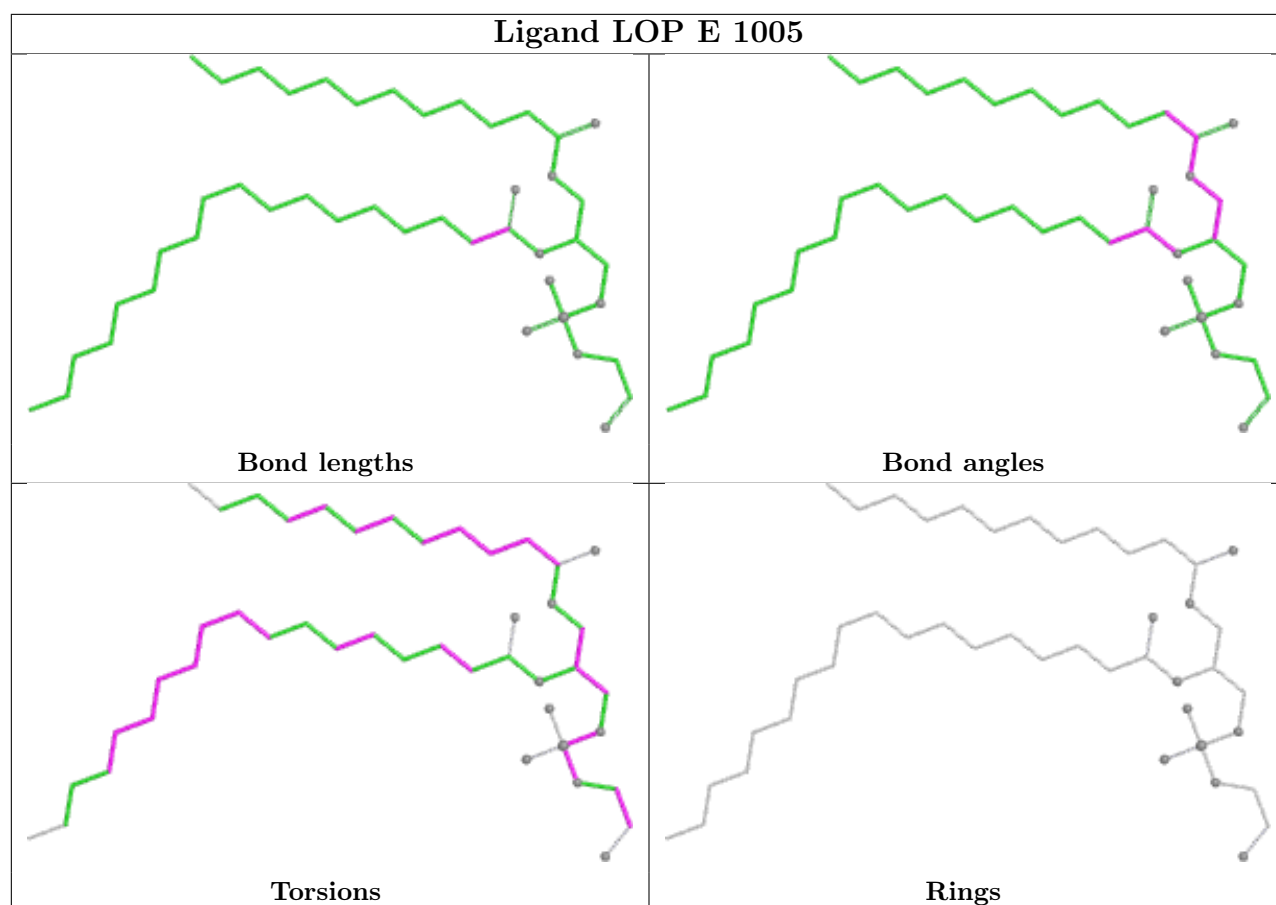


Ligand SMA O 1003



Ligand HEC P 1001





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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	428/445 (96%)	0.47	26 (6%)	27	14	58, 96, 145, 190	0
1	E	428/445 (96%)	0.37	20 (4%)	36	19	53, 81, 131, 179	0
1	K	428/445 (96%)	0.19	10 (2%)	61	38	43, 69, 108, 164	0
1	O	428/445 (96%)	0.37	17 (3%)	42	23	55, 84, 128, 166	0
2	B	256/272 (94%)	0.62	18 (7%)	22	12	62, 93, 147, 215	0
2	F	256/272 (94%)	0.57	19 (7%)	20	10	62, 98, 143, 178	0
2	L	256/272 (94%)	0.51	18 (7%)	22	12	47, 83, 134, 185	0
2	P	256/272 (94%)	0.66	14 (5%)	30	15	69, 105, 141, 183	0
3	C	179/187 (95%)	0.63	12 (6%)	24	12	58, 94, 152, 207	0
3	G	179/187 (95%)	0.58	11 (6%)	27	14	72, 97, 146, 202	0
3	M	179/187 (95%)	0.55	10 (5%)	30	15	68, 93, 142, 194	0
3	Q	179/187 (95%)	0.69	13 (7%)	21	11	53, 94, 152, 210	0
All	All	3452/3616 (95%)	0.48	188 (5%)	31	16	43, 89, 140, 215	0

All (188) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	Q	46	ALA	6.1
3	C	46	ALA	5.7
3	G	46	ALA	5.5
2	L	1	ALA	5.3
1	O	430	ALA	5.1
1	E	221	ASN	4.5
2	B	1	ALA	4.5
1	E	430	ALA	4.4
3	M	46	ALA	4.3
1	A	362	MET	4.3
2	B	2	GLY	4.1

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Mol	Chain	Res	Type	RSRZ
1	A	174	HIS	4.1
1	O	3	GLY	4.1
2	L	250	ARG	3.9
2	P	1	ALA	3.9
1	A	236	GLU	3.9
2	L	152	ASP	3.7
2	P	228	GLN	3.7
1	E	90	GLY	3.7
2	F	3	GLY	3.6
1	O	4	ILE	3.6
3	Q	47	LEU	3.6
1	A	239	LYS	3.5
2	L	2	GLY	3.5
1	A	428	PHE	3.4
1	O	19	LEU	3.4
3	C	47	LEU	3.4
3	G	47	LEU	3.4
1	E	3	GLY	3.4
3	C	10	THR	3.4
2	L	3	GLY	3.3
1	A	382	ALA	3.3
1	E	19	LEU	3.3
2	F	2	GLY	3.3
2	F	135	LEU	3.3
1	E	9	TYR	3.2
3	G	9	GLY	3.2
3	C	9	GLY	3.2
3	C	187	GLY	3.2
1	K	236	GLU	3.2
2	P	3	GLY	3.2
3	M	10	THR	3.2
3	M	47	LEU	3.1
3	Q	11	ARG	3.1
2	F	126	GLY	3.1
2	B	202	ALA	3.1
2	B	3	GLY	3.0
2	F	1	ALA	3.0
1	K	3	GLY	3.0
2	L	190	MET	3.0
2	B	152	ASP	2.9
3	Q	10	THR	2.9
2	F	250	ARG	2.9

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Mol	Chain	Res	Type	RSRZ
3	C	11	ARG	2.9
3	M	45	GLN	2.9
2	P	236	PHE	2.9
1	A	35	ILE	2.9
3	M	142	GLY	2.9
1	E	20	HIS	2.9
2	B	205	HIS	2.9
1	A	232	THR	2.8
1	K	384	GLN	2.8
1	A	39	ARG	2.8
3	Q	182	THR	2.8
1	A	430	ALA	2.8
3	Q	9	GLY	2.8
1	E	5	PRO	2.8
2	B	228	GLN	2.8
1	A	340	ILE	2.8
2	F	200	HIS	2.8
1	O	7	ASP	2.7
1	O	94	ARG	2.7
2	L	72	VAL	2.7
1	O	236	GLU	2.7
3	C	83	GLU	2.7
3	Q	16	TYR	2.7
2	P	200	HIS	2.6
3	G	145	GLY	2.6
2	B	7	GLU	2.6
1	A	3	GLY	2.6
2	B	255	VAL	2.6
3	Q	54	VAL	2.6
2	B	141	GLU	2.6
1	E	93	LEU	2.5
2	F	6	VAL	2.5
1	K	235	ALA	2.5
2	F	236	PHE	2.5
2	F	201	ASP	2.5
3	C	143	ASP	2.5
2	P	82	GLU	2.5
3	G	112	GLN	2.5
3	C	48	ALA	2.5
3	G	10	THR	2.5
1	O	31	ASP	2.5
2	F	198	ASP	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	190	MET	2.5
2	B	224	MET	2.5
2	L	4	GLY	2.5
2	L	7	GLU	2.5
3	Q	45	GLN	2.5
1	K	362	MET	2.4
1	O	24	PRO	2.4
1	A	355	ARG	2.4
1	E	8	HIS	2.4
2	B	23	HIS	2.4
1	A	221	ASN	2.4
1	A	414	ILE	2.4
1	O	228	GLU	2.4
1	E	384	GLN	2.4
2	F	204	VAL	2.4
1	E	21	SER	2.4
1	A	119	GLY	2.4
3	G	83	GLU	2.4
3	C	26	ALA	2.4
1	K	174	HIS	2.4
1	O	5	PRO	2.4
3	Q	69	GLY	2.3
2	L	198	ASP	2.3
2	L	140	GLU	2.3
3	G	48	ALA	2.3
1	O	174	HIS	2.3
2	L	236	PHE	2.3
3	Q	51	PHE	2.3
1	A	237	ALA	2.3
1	E	348	TRP	2.3
2	L	110	PHE	2.3
3	Q	15	LEU	2.3
2	F	5	HIS	2.3
2	P	5	HIS	2.3
1	A	231	ARG	2.3
3	M	153	GLY	2.3
3	G	35	ILE	2.3
1	A	230	ARG	2.2
1	E	94	ARG	2.2
1	E	89	GLY	2.2
2	L	256	LYS	2.2
1	K	414	ILE	2.2

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Mol	Chain	Res	Type	RSRZ
3	M	112	GLN	2.2
3	M	16	TYR	2.2
2	L	111	HIS	2.2
1	A	417	PRO	2.2
2	F	196	TYR	2.2
2	L	126	GLY	2.2
2	P	109	GLY	2.2
1	A	116	LEU	2.2
1	K	430	ALA	2.2
3	C	45	GLN	2.2
1	A	227	VAL	2.2
2	P	103	MET	2.2
3	G	182	THR	2.2
1	A	19	LEU	2.2
1	A	94	ARG	2.1
2	B	5	HIS	2.1
2	F	190	MET	2.1
1	K	351	THR	2.1
2	B	109	GLY	2.1
2	P	4	GLY	2.1
1	O	9	TYR	2.1
2	P	172	ALA	2.1
1	E	11	PRO	2.1
2	L	254	GLY	2.1
1	O	247	TYR	2.1
2	F	106	ALA	2.1
2	F	120	GLN	2.1
1	O	421	PRO	2.1
2	F	254	GLY	2.1
2	L	219	ALA	2.1
2	P	65	ALA	2.1
1	A	247	TYR	2.1
2	P	7	GLU	2.1
1	A	6	HIS	2.1
1	O	239	LYS	2.1
3	M	15	LEU	2.1
2	B	230	GLY	2.1
1	E	7	ASP	2.1
1	E	22	ARG	2.1
1	E	4	ILE	2.1
2	P	197	ALA	2.1
1	K	416	LYS	2.0

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Mol	Chain	Res	Type	RSRZ
2	F	111	HIS	2.0
3	G	45	GLN	2.0
2	B	236	PHE	2.0
1	E	311	ASP	2.0
1	O	21	SER	2.0
2	B	4	GLY	2.0
3	M	48	ALA	2.0
3	C	58	GLU	2.0
3	Q	17	TYR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	LOP	A	1006	45/45	0.80	0.21	80,104,114,116	0
6	ANJ	A	1004	39/39	0.81	0.16	77,89,107,108	0
8	LOP	E	1005	45/45	0.81	0.19	75,95,109,113	0
8	LOP	O	1005	45/45	0.81	0.19	78,100,116,117	0
8	LOP	K	1006	45/45	0.85	0.19	64,85,101,106	0
6	ANJ	E	1004	39/39	0.88	0.14	58,74,101,101	0
6	ANJ	O	1004	39/39	0.88	0.14	67,84,107,108	0
6	ANJ	K	1004	39/39	0.89	0.12	49,71,100,101	0
7	SR	K	1005	1/1	0.91	0.09	121,121,121,121	0
7	SR	A	1005	1/1	0.93	0.07	133,133,133,133	0
5	SMA	A	1003	37/37	0.93	0.12	56,69,83,89	0
5	SMA	O	1003	37/37	0.93	0.13	53,68,90,93	0
5	SMA	E	1003	37/37	0.94	0.12	47,67,88,88	0
4	HEM	A	1002	43/43	0.94	0.14	80,96,114,126	0

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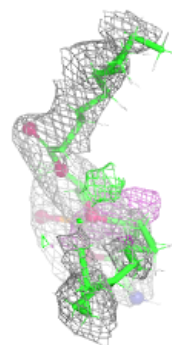
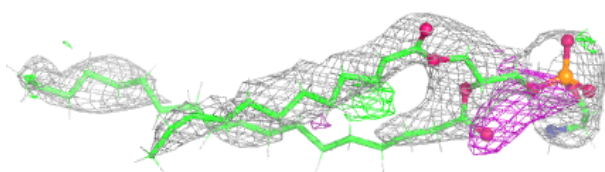
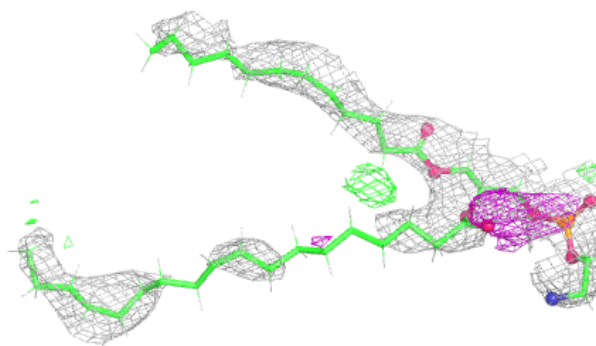
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	SMA	K	1003	37/37	0.95	0.12	47,64,83,86	0
4	HEM	K	1002	43/43	0.95	0.12	64,72,89,96	0
4	HEM	O	1002	43/43	0.95	0.12	75,80,97,98	0
4	HEM	A	1001	43/43	0.96	0.14	61,72,87,95	0
4	HEM	O	1001	43/43	0.96	0.12	48,64,79,81	0
4	HEM	E	1002	43/43	0.96	0.11	65,78,94,98	0
7	SR	B	1002	1/1	0.96	0.04	117,117,117,117	0
7	SR	F	1002	1/1	0.96	0.05	132,132,132,132	0
9	HEC	F	1001	43/43	0.96	0.10	47,61,82,85	0
9	HEC	P	1001	43/43	0.96	0.09	47,62,75,83	0
4	HEM	E	1001	43/43	0.97	0.12	56,69,85,89	0
9	HEC	B	1001	43/43	0.97	0.09	41,56,69,76	0
7	SR	L	1002	1/1	0.97	0.04	113,113,113,113	0
9	HEC	L	1001	43/43	0.97	0.08	40,55,69,69	0
7	SR	P	1002	1/1	0.97	0.04	144,144,144,144	0
4	HEM	K	1001	43/43	0.98	0.10	51,62,79,83	0
10	FES	Q	1001	4/4	0.98	0.08	58,63,64,66	0
10	FES	G	1001	4/4	0.99	0.07	71,74,74,76	0
10	FES	M	1001	4/4	0.99	0.05	72,72,75,77	0
10	FES	C	1001	4/4	0.99	0.05	59,62,64,64	0

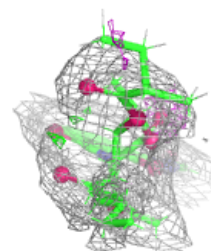
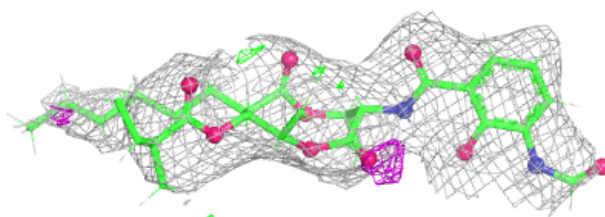
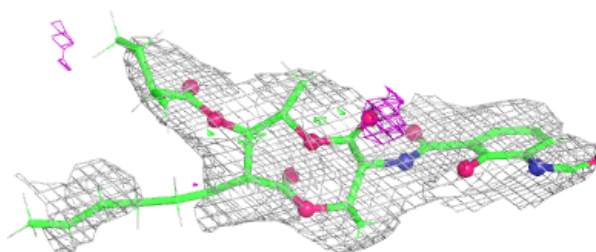
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around LOP A 1006:

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

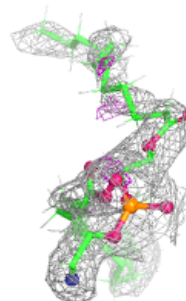
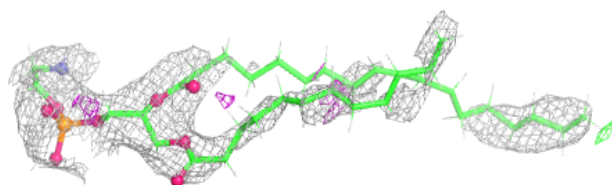
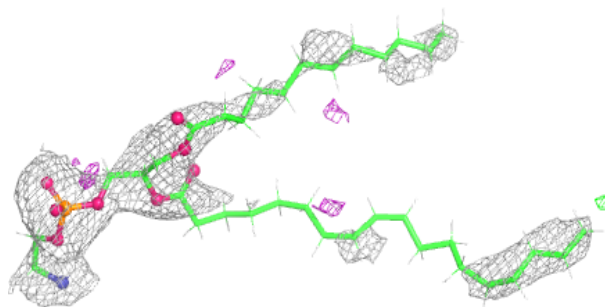
**Electron density around ANJ A 1004:**

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

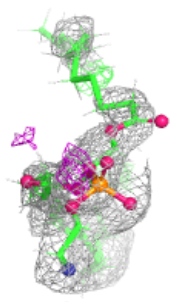
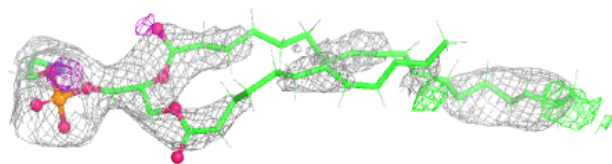
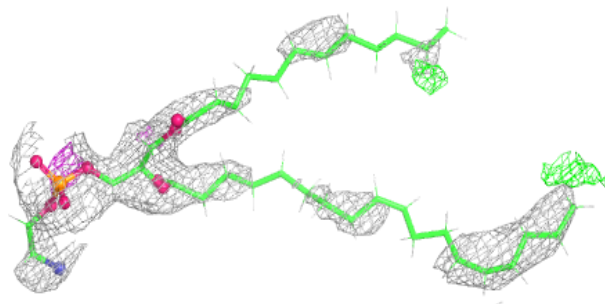


Electron density around LOP E 1005:

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

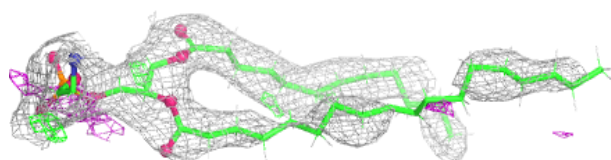
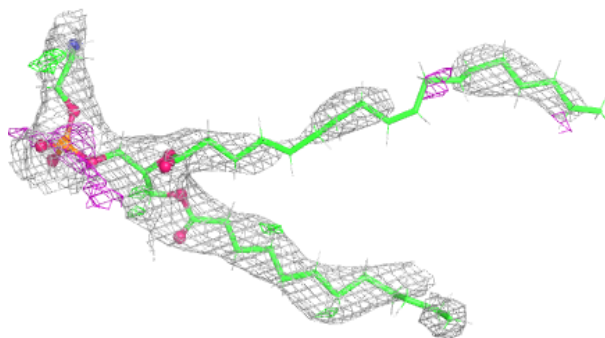
**Electron density around LOP O 1005:**

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

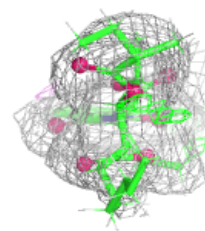
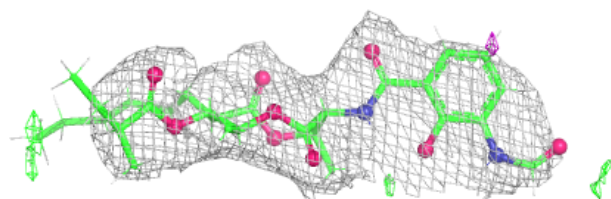
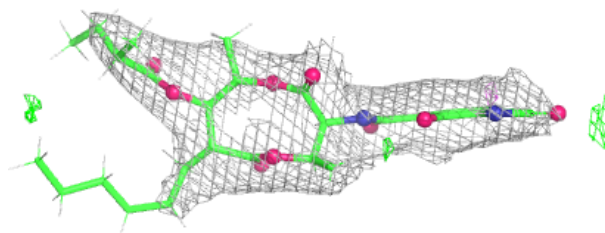


Electron density around LOP K 1006:

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and green (positive)

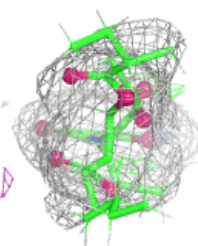
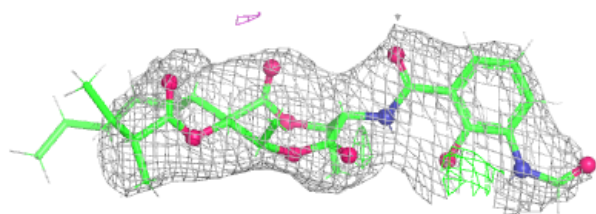
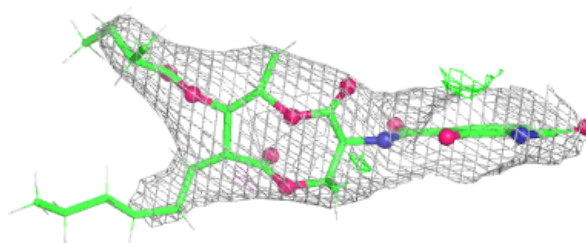
**Electron density around ANJ E 1004:**

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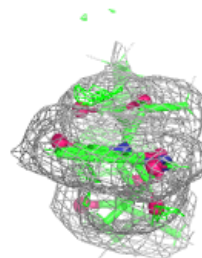
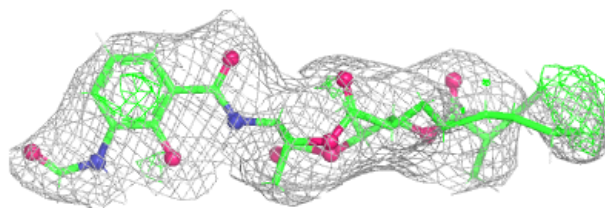
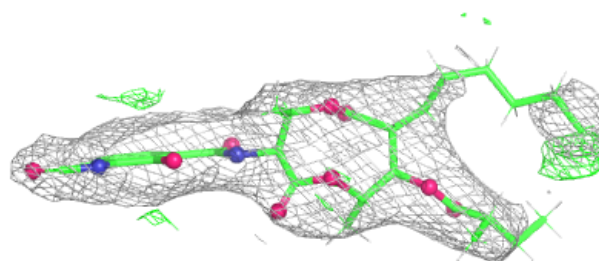


Electron density around ANJ O 1004:

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

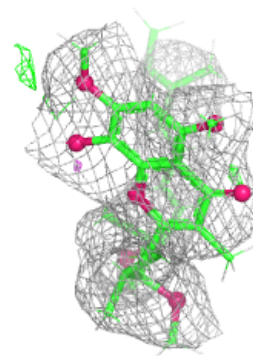
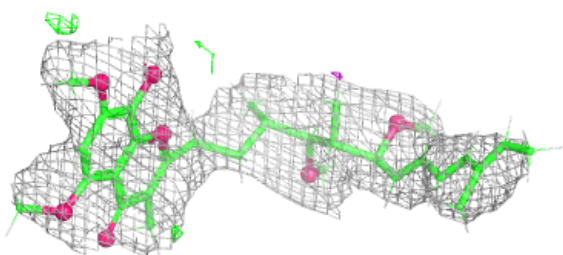
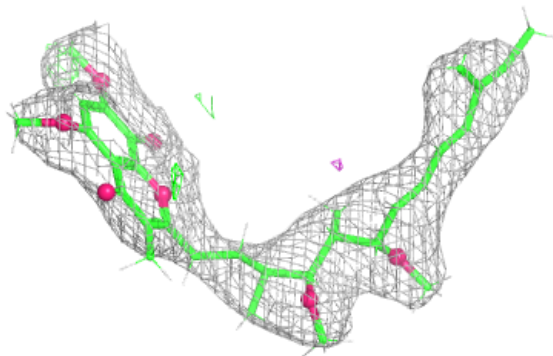
**Electron density around ANJ K 1004:**

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and green (positive)

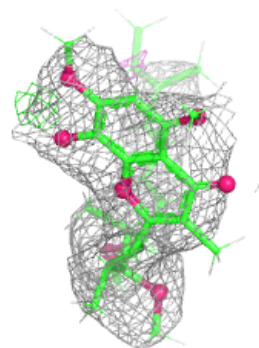
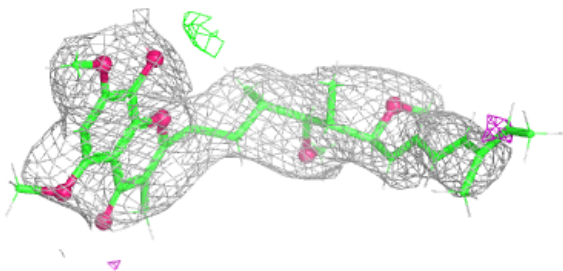
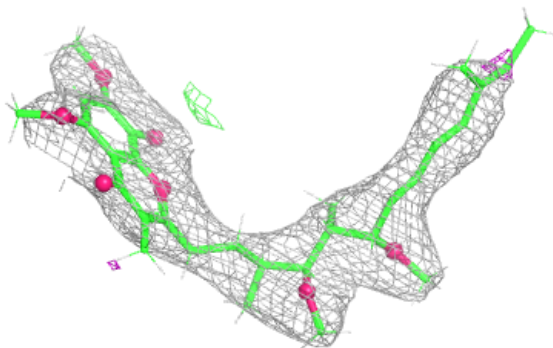


Electron density around SMA A 1003:

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

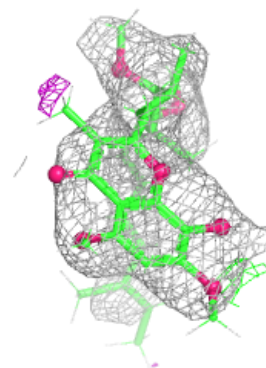
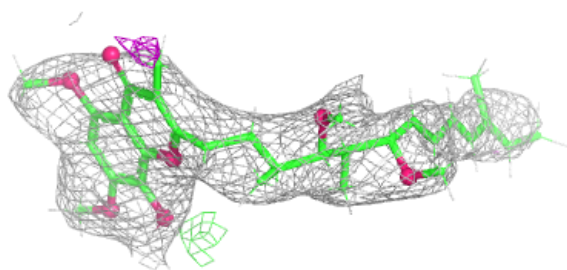
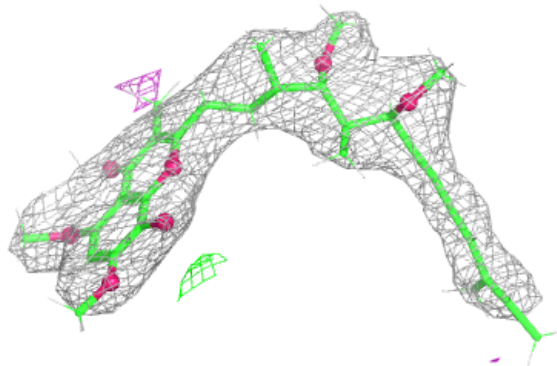
**Electron density around SMA O 1003:**

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



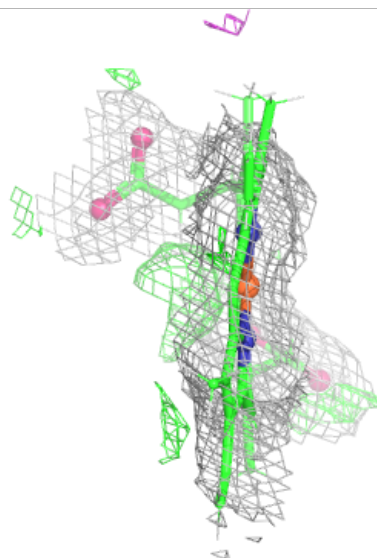
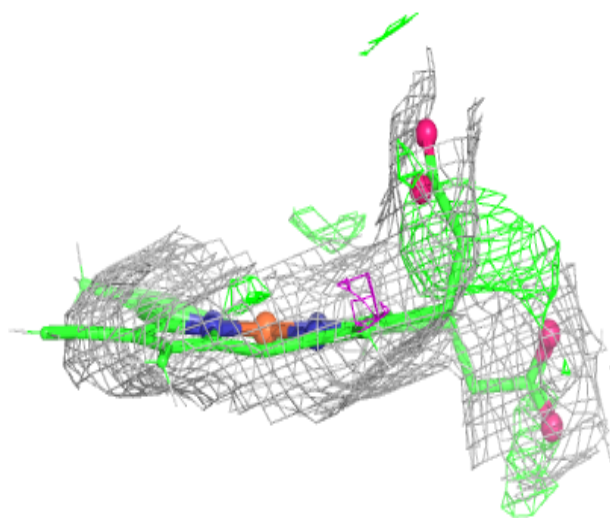
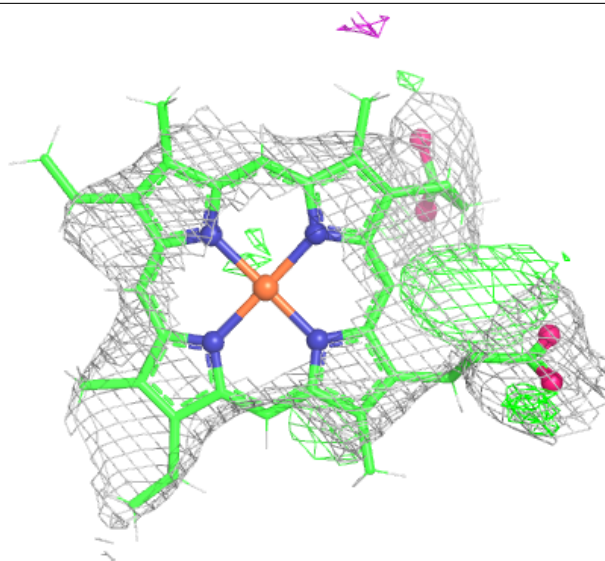
Electron density around SMA E 1003:

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and green (positive)



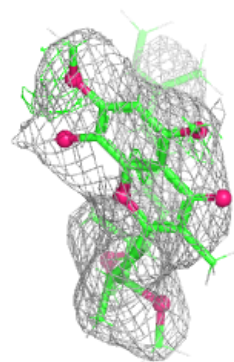
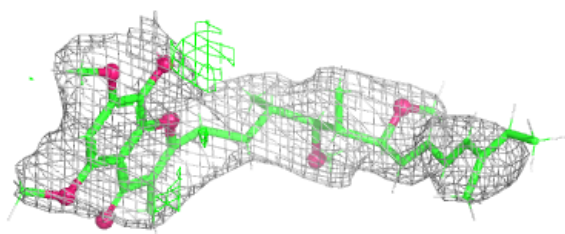
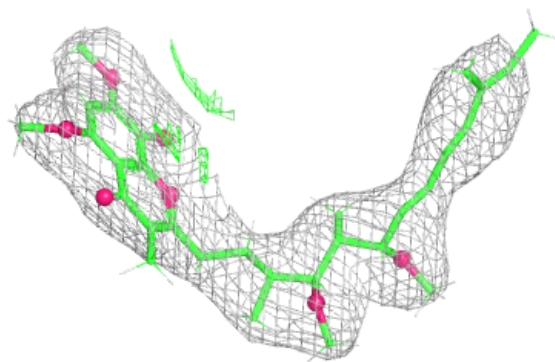
Electron density around HEM A 1002:

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and green (positive)



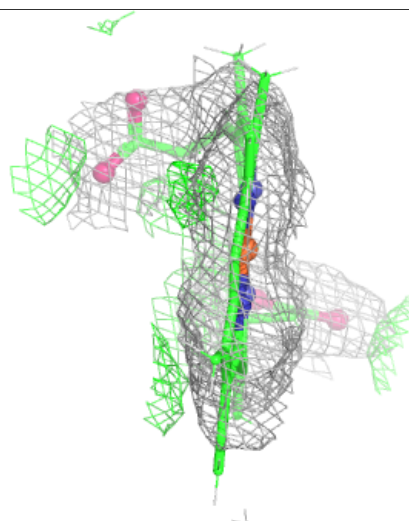
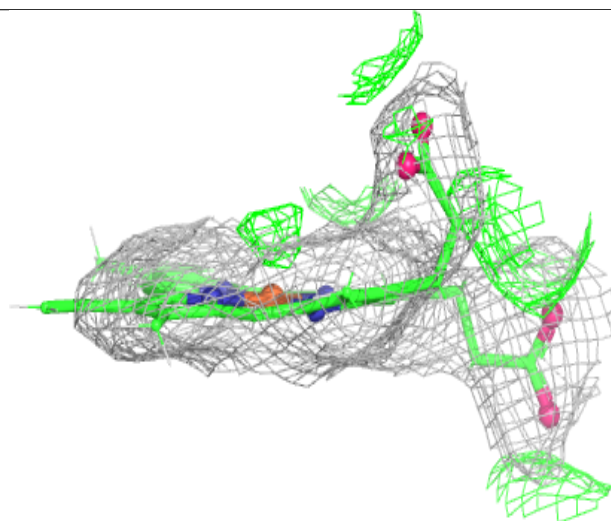
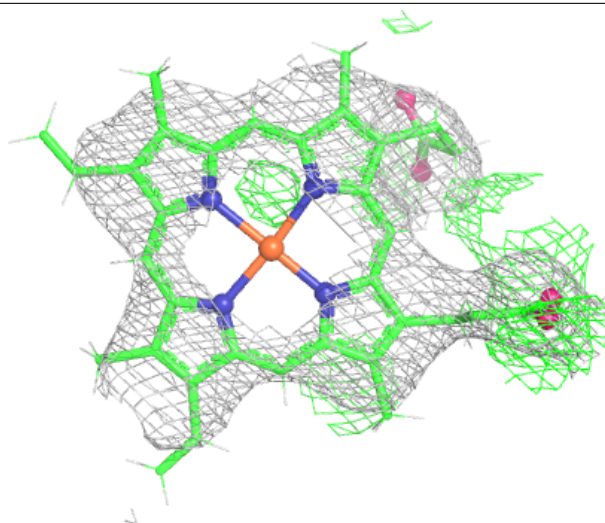
Electron density around SMA K 1003:

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and green (positive)



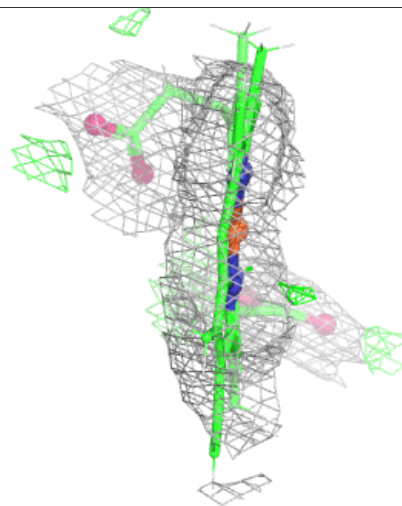
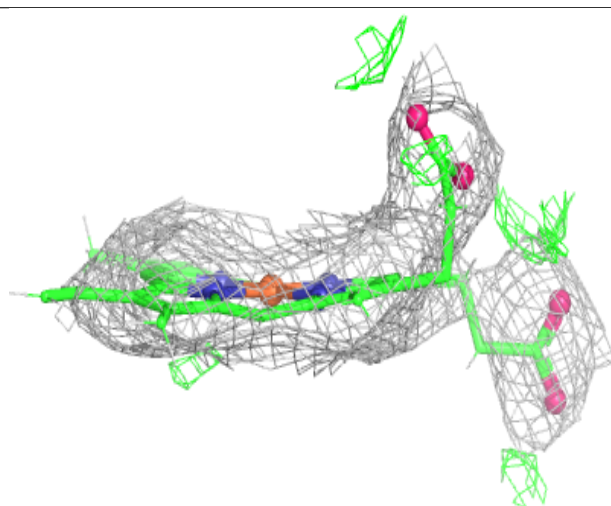
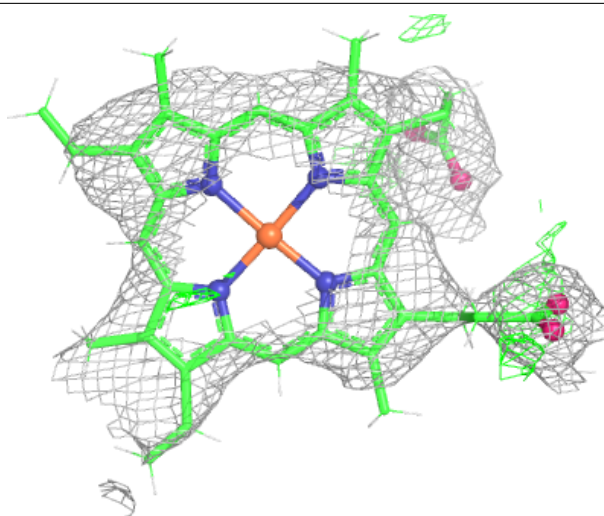
Electron density around HEM K 1002:

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and green (positive)



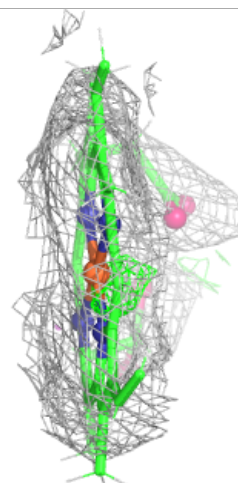
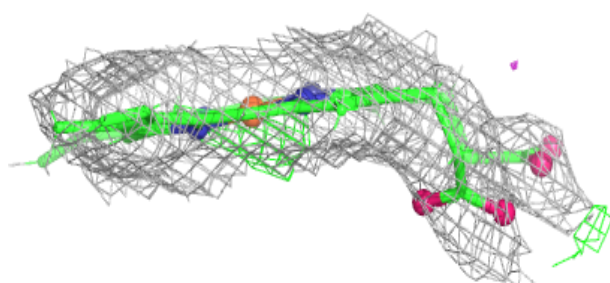
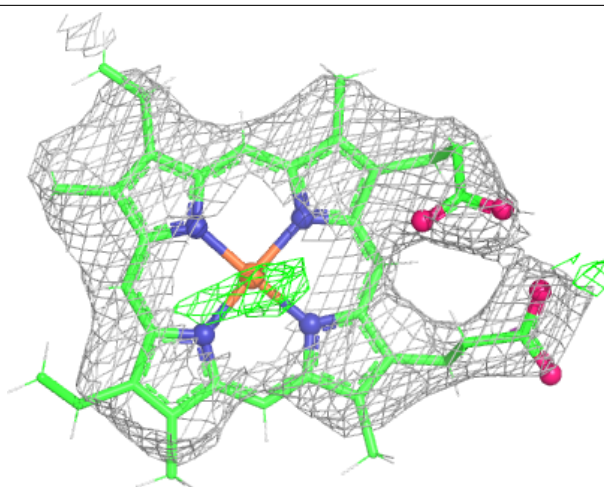
Electron density around HEM O 1002:

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and green (positive)



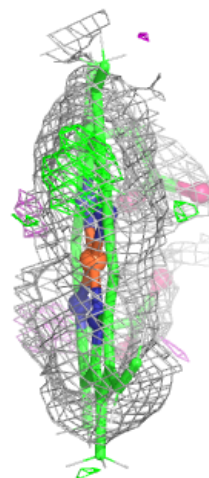
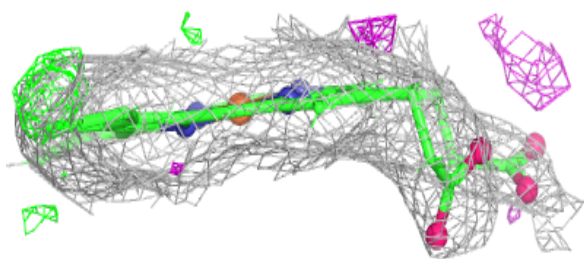
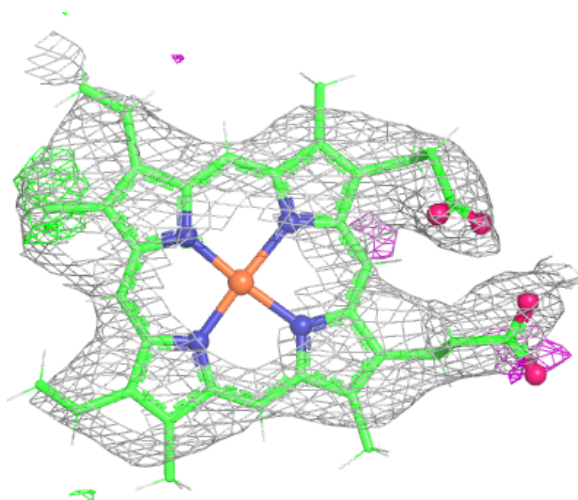
Electron density around HEM A 1001:

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and green (positive)



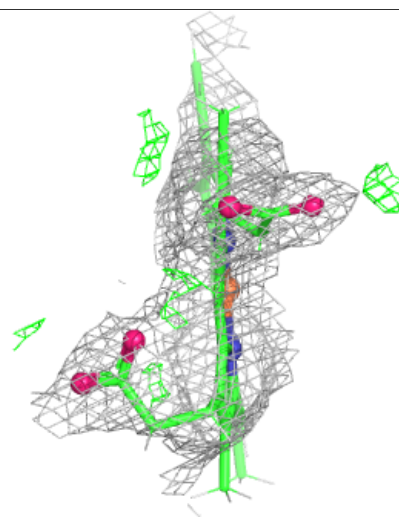
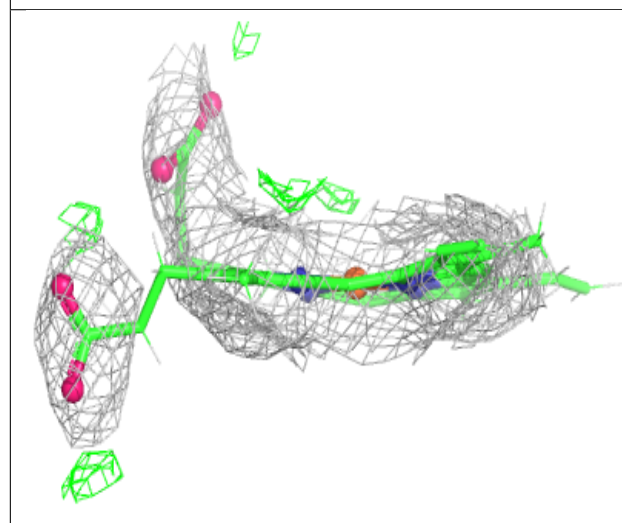
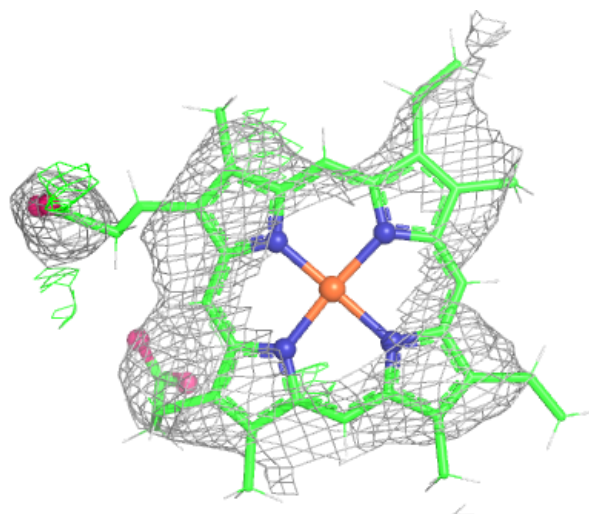
Electron density around HEM O 1001:

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and green (positive)



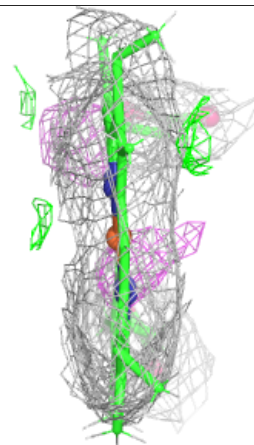
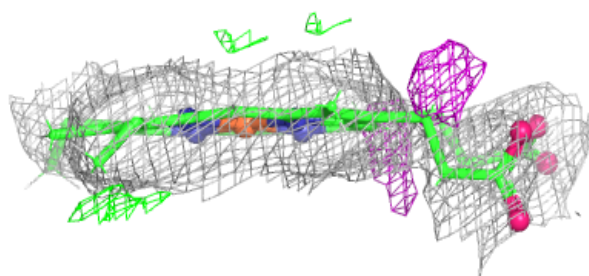
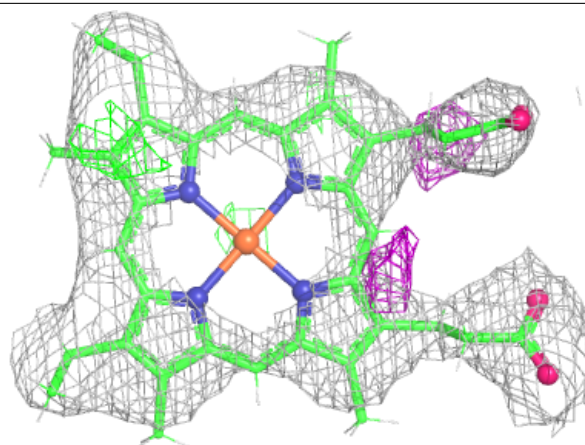
Electron density around HEM E 1002:

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and green (positive)



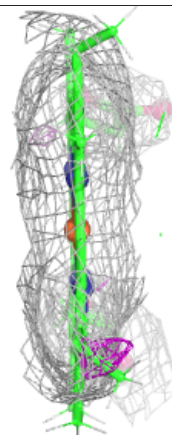
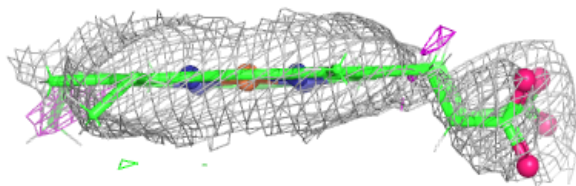
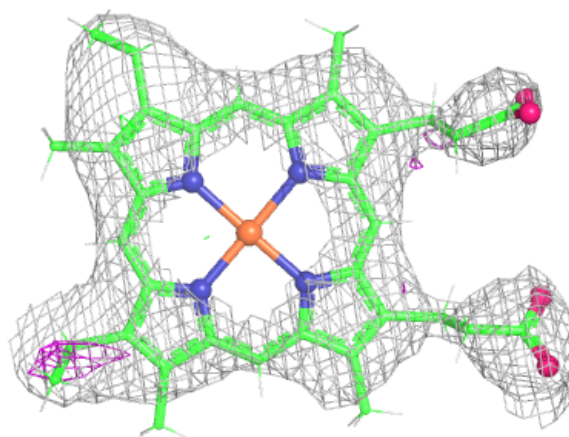
Electron density around HEC F 1001:

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and green (positive)



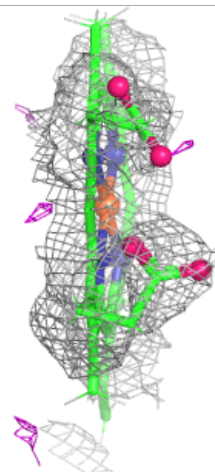
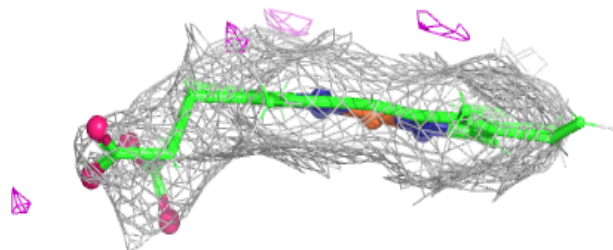
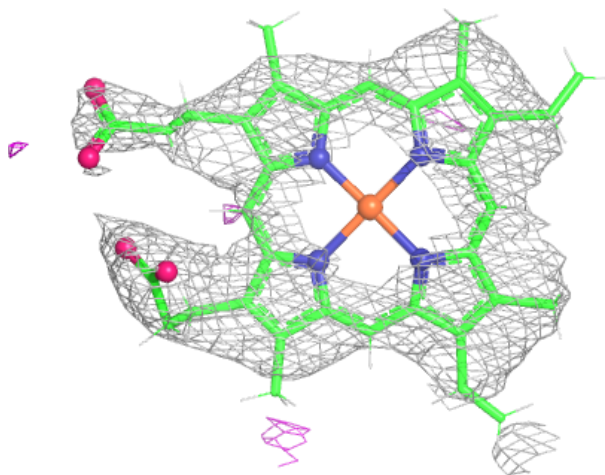
Electron density around HEC P 1001:

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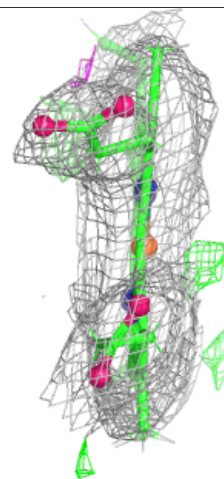
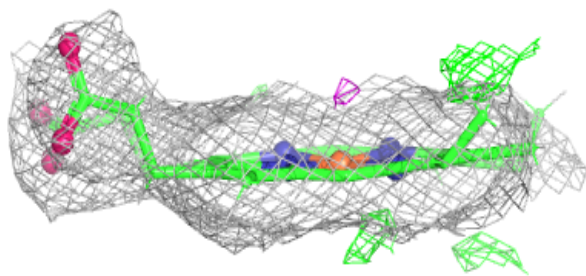
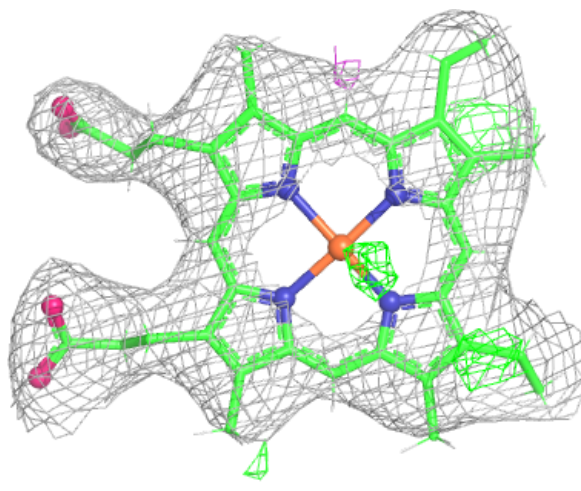
Electron density around HEM E 1001:

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



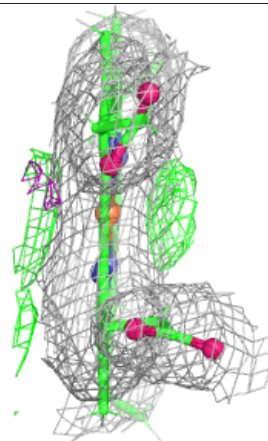
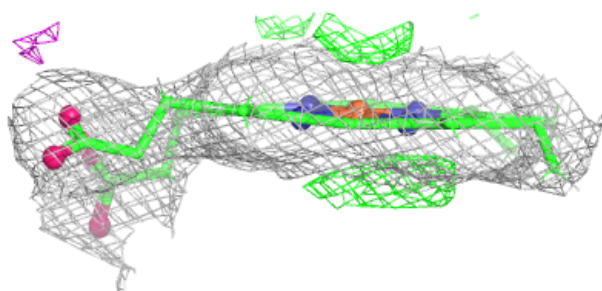
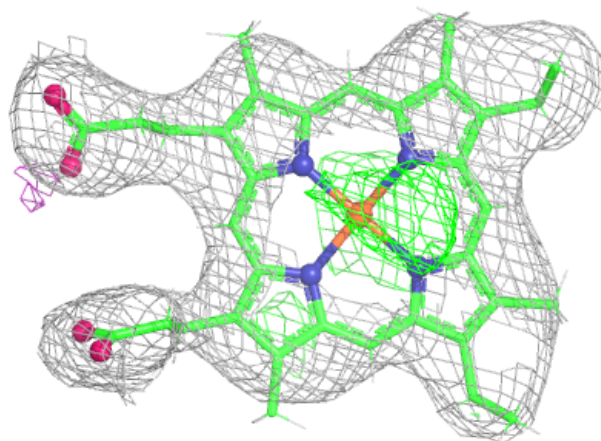
Electron density around HEC B 1001:

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



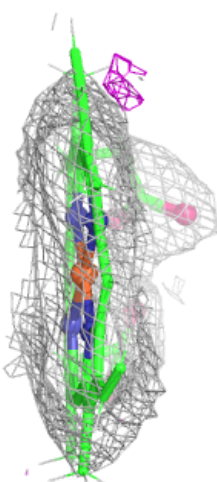
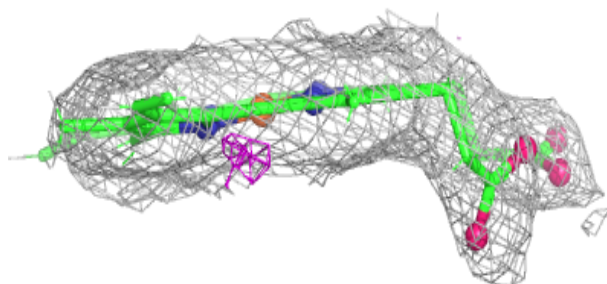
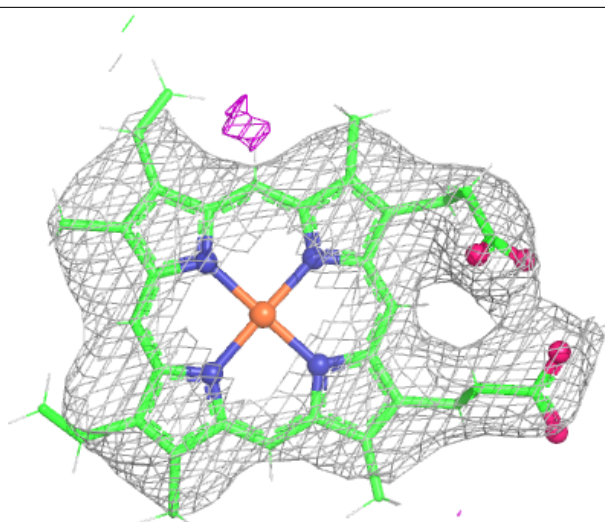
Electron density around HEC L 1001:

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



Electron density around HEM K 1001:

$2mF_o-DF_c$ (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.