



# Full wwPDB X-ray Structure Validation Report ⓘ

Apr 28, 2025 – 08:20 PM JST

PDB ID : 7C90 / pdb\_00007c90  
Title : Crystal structure of Cytochrome CL from the marine methylophilic bacterium *Methylophilaga aminisulfidivorans* MPT (Ma-CytC<sub>L</sub>)  
Authors : Ghosh, S.; Dhanasingh, I.; Lee, S.H.  
Deposited on : 2020-06-04  
Resolution : 2.13 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](mailto: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 Phenix2.0rc1  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 2.0rc1  
EDS : 3.0  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
CCP4 : 9.0.006 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.43.1

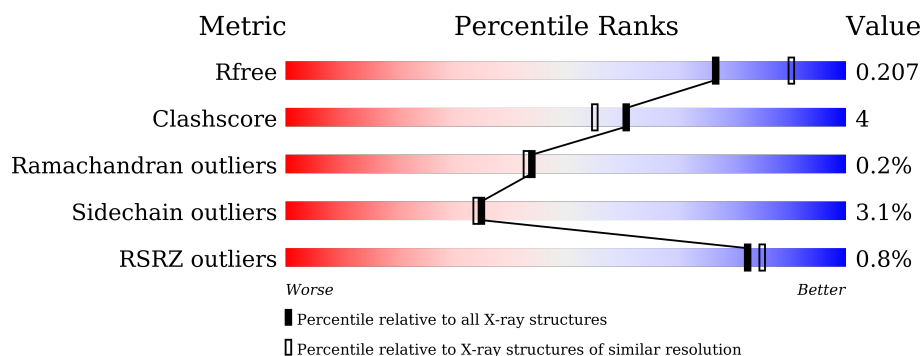
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.13 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))
$R_{free}$	164625	3336 (2.16-2.12)
Clashscore	180529	3585 (2.16-2.12)
Ramachandran outliers	177936	3554 (2.16-2.12)
Sidechain outliers	177891	3553 (2.16-2.12)
RSRZ outliers	164620	3337 (2.16-2.12)

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  $\geq 3$ , 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	194	<div> <div></div> <div>69% 8% 22%</div> </div>
1	B	194	<div> <div></div> <div>72% 6% 22%</div> </div>
1	C	194	<div> <div></div> <div>66% 8% 25%</div> </div>
1	D	194	<div> <div></div> <div>68% 7% 23%</div> </div>

## 2 Entry composition [i](#)

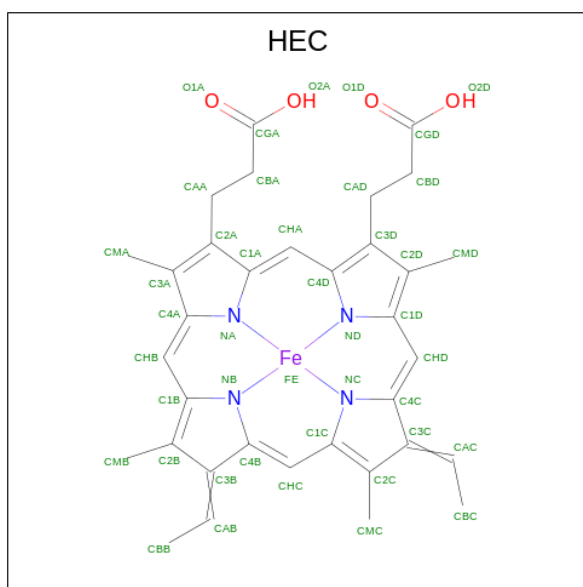
There are 8 unique types of molecules in this entry. The entry contains 5321 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 c, mono-and diheme variant.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	151	Total	C	N	O	S	0	0	0
			1185	748	197	234	6			
1	B	151	Total	C	N	O	S	0	0	0
			1185	748	197	234	6			
1	C	145	Total	C	N	O	S	0	0	0
			1137	717	188	226	6			
1	D	150	Total	C	N	O	S	0	0	0
			1176	742	195	233	6			

- Molecule 2 is HEME C (CCD ID: HEC) (formula:  $C_{34}H_{34}FeN_4O_4$ ) (labeled as "Ligand of Interest" by depositor).



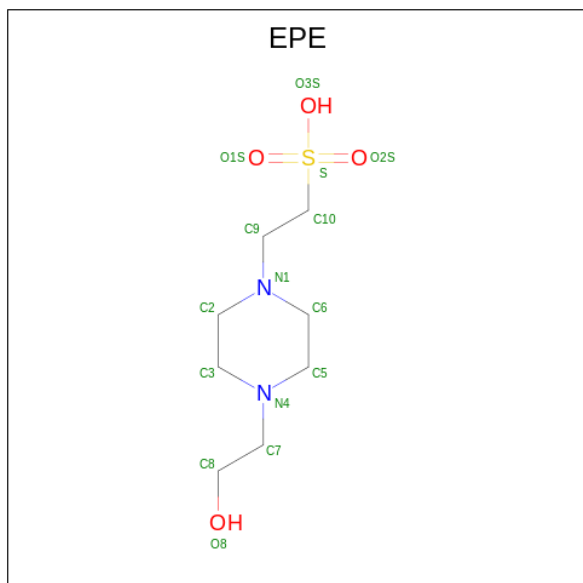
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		
2	B	1	Total	C	Fe	N O	0	0
			43	34	1	4 4		

*Continued on next page...*

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
2	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 3 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula:  $C_8H_{18}N_2O_4S$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
3	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
3	B	1	Total	C	N	O	S	0	0
			15	8	2	4	1		
3	C	1	Total	C	N	O	S	0	0
			15	8	2	4	1		

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



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

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Ca	0	0
			1	1		
5	B	1	Total	Ca	0	0
			1	1		
5	C	1	Total	Ca	0	0
			1	1		
5	D	1	Total	Ca	0	0
			1	1		

- Molecule 6 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	1	Total	Fe	0	0
			1	1		
6	B	1	Total	Fe	0	0
			1	1		

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total 1	Fe 1	0	0
6	D	1	Total 1	Fe 1	0	0

- Molecule 7 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total 2	Na 2	0	0
7	B	2	Total 2	Na 2	0	0
7	C	2	Total 2	Na 2	0	0
7	D	2	Total 2	Na 2	0	0

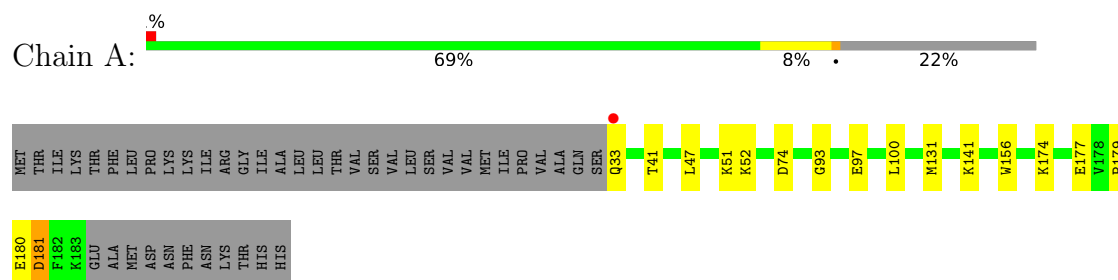
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	108	Total 108	O 108	0	0
8	B	90	Total 90	O 90	0	0
8	C	83	Total 83	O 83	0	0
8	D	91	Total 91	O 91	0	0

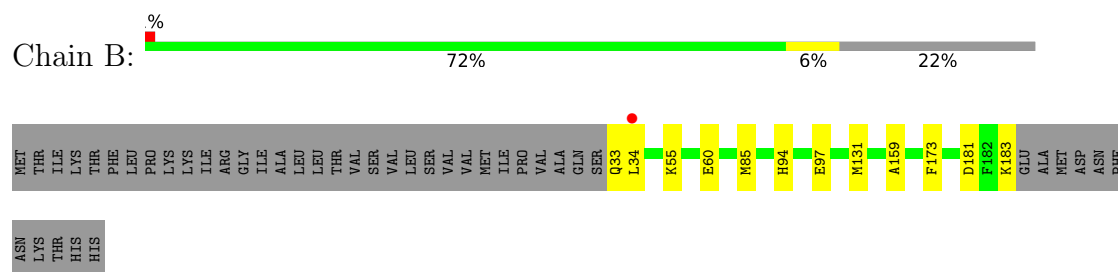
### 3 Residue-property plots [i](#)

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.

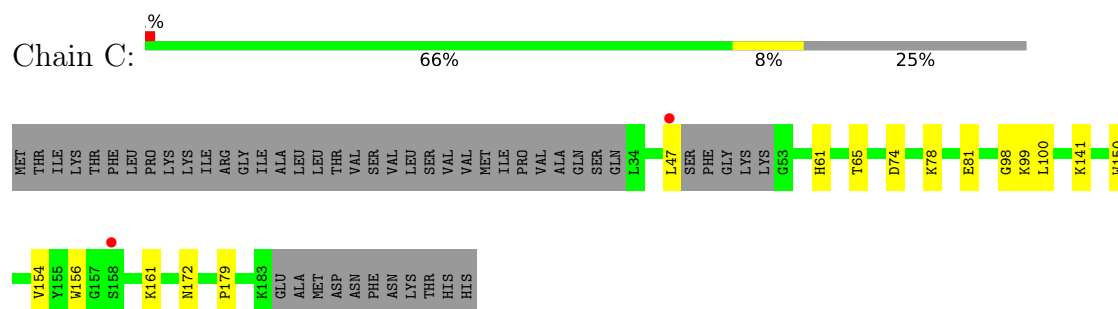
- Molecule 1: Cytochrome c, mono-and diheme variant



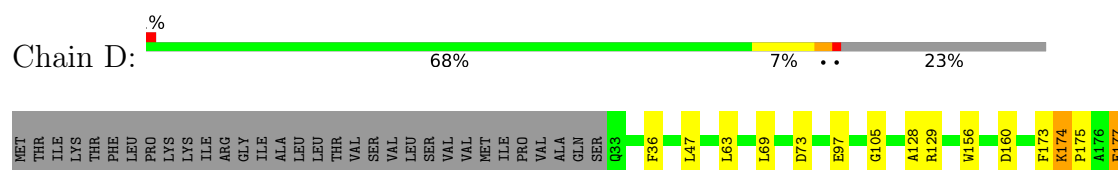
- Molecule 1: Cytochrome c, mono-and diheme variant



- Molecule 1: Cytochrome c, mono-and diheme variant



- Molecule 1: Cytochrome c, mono-and diheme variant



E180	LYS	GLU	ALA	NET	ASP	ASN	PHE	ASN	LYS	THR	HIS	HIS
D181	LYS	GLU	ALA	NET	ASP	ASN	PHE	ASN	LYS	THR	HIS	HIS
F182	LYS	GLU	ALA	NET	ASP	ASN	PHE	ASN	LYS	THR	HIS	HIS



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.76Å 76.06Å 66.44Å 90.00° 106.85° 90.00°	Depositor
Resolution (Å)	30.58 – 2.13 30.58 – 2.13	Depositor EDS
% Data completeness (in resolution range)	97.4 (30.58-2.13) 97.4 (30.58-2.13)	Depositor EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.14 (at 2.14Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
R, $R_{free}$	0.148 , 0.200 0.158 , 0.207	Depositor DCC
$R_{free}$ test set	30898 reflections (6.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	31.7	Xtriage
Anisotropy	0.038	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.40 , 50.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.48$ , $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.97	EDS
Total number of atoms	5321	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.18% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: CA, HEC, NA, FE, EPE, GOL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z  > 5$	RMSZ	$\# Z  > 5$
1	A	1.06	0/1214	1.35	1/1641 (0.1%)
1	B	1.03	0/1214	1.33	0/1641
1	C	1.03	0/1164	1.36	1/1575 (0.1%)
1	D	1.08	2/1205 (0.2%)	1.33	3/1630 (0.2%)
All	All	1.05	2/4797 (0.0%)	1.34	5/6487 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	128	ALA	C-N	-7.72	1.21	1.33
1	D	129	ARG	C-N	-5.08	1.26	1.33

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	129	ARG	CA-C-N	5.68	132.39	121.54
1	D	129	ARG	C-N-CA	5.68	132.39	121.54
1	D	182	PHE	CA-CB-CG	5.65	119.45	113.80
1	C	98	GLY	CA-C-O	-5.16	117.86	121.88
1	A	93	GLY	CA-C-O	-5.03	117.86	122.24

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	1185	0	1113	10	0
1	B	1185	0	1113	9	0
1	C	1137	0	1061	9	0
1	D	1176	0	1100	13	0
2	A	43	0	30	0	0
2	B	43	0	30	0	0
2	C	43	0	30	1	0
2	D	43	0	30	0	0
3	A	15	0	18	0	0
3	B	30	0	35	0	0
3	C	15	0	18	0	0
4	A	12	0	16	0	0
4	B	6	0	8	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
7	A	2	0	0	0	0
7	B	2	0	0	0	0
7	C	2	0	0	0	0
7	D	2	0	0	0	0
8	A	108	0	0	2	0
8	B	90	0	0	2	0
8	C	83	0	0	1	0
8	D	91	0	0	1	0
All	All	5321	0	4602	39	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 4.

All (39) 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:D:177:GLU:H	1:D:177:GLU:CD	1.91	0.77
1:D:97:GLU:HG3	1:D:181:ASP:HB2	1.72	0.72
1:D:97:GLU:HB2	1:D:182:PHE:HB3	1.75	0.67

*Continued on next page...*

*Continued from previous page...*

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:33:GLN:O	1:A:33:GLN:HG3	1.97	0.65
1:A:51:LYS:HE2	8:A:375:HOH:O	2.02	0.59
1:C:74:ASP:O	1:C:78:LYS:HG2	2.04	0.58
1:C:172:ASN:ND2	8:C:302:HOH:O	2.37	0.57
1:B:131:MET:HE2	1:C:100:LEU:HB2	1.87	0.56
1:D:182:PHE:C	1:D:182:PHE:CD1	2.85	0.54
1:C:47:LEU:O	1:C:141:LYS:NZ	2.32	0.51
1:C:81:GLU:HG3	1:C:150:TRP:NE1	2.27	0.49
1:B:97:GLU:HG3	1:B:181:ASP:HB2	1.95	0.48
1:D:173:PHE:CD1	1:D:173:PHE:C	2.91	0.48
1:A:131:MET:HE3	1:A:131:MET:HB3	1.78	0.48
1:D:47:LEU:HD13	1:D:63:LEU:HD22	1.96	0.47
1:D:97:GLU:HB2	1:D:182:PHE:CB	2.42	0.47
1:A:156:TRP:CE3	1:A:179:PRO:HD3	2.51	0.46
1:D:36:PHE:CD1	1:D:47:LEU:HD11	2.51	0.44
1:D:181:ASP:OD1	1:D:181:ASP:N	2.50	0.44
1:B:183:LYS:CG	8:B:383:HOH:O	2.66	0.43
1:B:159:ALA:HB2	1:B:173:PHE:CD2	2.54	0.43
1:D:105:GLY:HA3	1:D:156:TRP:CZ3	2.54	0.43
1:A:97:GLU:HG3	1:A:181:ASP:HB2	2.00	0.42
1:C:61:HIS:CE1	1:C:65:THR:HG21	2.54	0.42
1:B:34:LEU:HD12	1:B:34:LEU:HA	1.93	0.42
1:D:174:LYS:NZ	8:D:303:HOH:O	2.38	0.42
2:C:201:HEC:HBB3	2:C:201:HEC:HMB3	2.02	0.41
1:A:47:LEU:O	1:A:141:LYS:NZ	2.46	0.41
1:A:52:LYS:HG2	8:A:311:HOH:O	2.20	0.41
1:D:173:PHE:HE1	1:D:175:PRO:HG3	1.85	0.41
1:B:183:LYS:CD	8:B:383:HOH:O	2.68	0.41
1:D:73:ASP:OD1	1:D:73:ASP:C	2.64	0.41
1:A:177:GLU:H	1:A:177:GLU:CD	2.28	0.41
1:B:85:MET:HE1	1:B:94:HIS:O	2.21	0.40
1:C:156:TRP:CE3	1:C:179:PRO:HD3	2.57	0.40
1:A:41:THR:HG22	1:B:183:LYS:HD2	2.02	0.40
1:A:180:GLU:CD	1:A:180:GLU:H	2.29	0.40
1:B:131:MET:CE	1:C:100:LEU:HB2	2.51	0.40
1:C:150:TRP:CZ2	1:C:154:VAL:HG11	2.56	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	149/194 (77%)	145 (97%)	4 (3%)	0	100	100
1	B	149/194 (77%)	145 (97%)	4 (3%)	0	100	100
1	C	141/194 (73%)	136 (96%)	4 (3%)	1 (1%)	19	13
1	D	148/194 (76%)	144 (97%)	4 (3%)	0	100	100
All	All	587/776 (76%)	570 (97%)	16 (3%)	1 (0%)	44	43

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	99	LYS

### 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	123/162 (76%)	119 (97%)	4 (3%)	33	31
1	B	123/162 (76%)	120 (98%)	3 (2%)	44	45
1	C	118/162 (73%)	117 (99%)	1 (1%)	79	83
1	D	122/162 (75%)	115 (94%)	7 (6%)	17	13
All	All	486/648 (75%)	471 (97%)	15 (3%)	35	34

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	ASP
1	A	100	LEU
1	A	174	LYS
1	A	181	ASP
1	B	33	GLN
1	B	55	LYS
1	B	60	GLU
1	C	161	LYS
1	D	69	LEU
1	D	160	ASP
1	D	174	LYS
1	D	177	GLU
1	D	180	GLU
1	D	181	ASP
1	D	182	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	114	ASN
1	A	172	ASN
1	B	33	GLN
1	B	61	HIS

### 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 oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 27 ligands modelled in this entry, 16 are monoatomic - leaving 11 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	GOL	A	204	-	5,5,5	0.12	0	5,5,5	0.48	0
3	EPE	A	202	-	15,15,15	1.80	1 (6%)	18,20,20	1.67	1 (5%)
2	HEC	C	201	1	32,50,50	1.86	9 (28%)	24,82,82	2.72	9 (37%)
2	HEC	A	201	1	32,50,50	2.12	10 (31%)	24,82,82	2.81	8 (33%)
3	EPE	B	203	-	15,15,15	1.80	1 (6%)	18,20,20	1.49	2 (11%)
2	HEC	D	201	1	32,50,50	1.94	9 (28%)	24,82,82	2.52	8 (33%)
2	HEC	B	201	1	32,50,50	1.92	9 (28%)	24,82,82	3.11	7 (29%)
4	GOL	B	204	-	5,5,5	0.14	0	5,5,5	0.36	0
4	GOL	A	203	-	5,5,5	0.11	0	5,5,5	0.39	0
3	EPE	B	202	-	15,15,15	1.74	1 (6%)	18,20,20	1.14	3 (16%)
3	EPE	C	202	-	15,15,15	1.98	1 (6%)	18,20,20	2.26	4 (22%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	A	204	-	-	4/4/4/4	-
3	EPE	A	202	-	-	7/9/19/19	0/1/1/1
2	HEC	C	201	1	-	1/10/54/54	-
2	HEC	A	201	1	-	0/10/54/54	-
3	EPE	B	203	-	-	4/9/19/19	0/1/1/1
2	HEC	D	201	1	-	1/10/54/54	-
2	HEC	B	201	1	-	2/10/54/54	-
4	GOL	B	204	-	-	2/4/4/4	-
4	GOL	A	203	-	-	4/4/4/4	-
3	EPE	B	202	-	-	2/9/19/19	0/1/1/1
3	EPE	C	202	-	-	6/9/19/19	0/1/1/1

All (41) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	202	EPE	C10-S	-7.20	1.67	1.77
3	B	202	EPE	C10-S	-6.58	1.68	1.77
3	B	203	EPE	C10-S	-6.52	1.68	1.77
3	A	202	EPE	C10-S	-6.38	1.68	1.77
2	A	201	HEC	C3C-C2C	6.21	1.47	1.40
2	A	201	HEC	C2B-C3B	5.40	1.46	1.40
2	B	201	HEC	C3C-C2C	5.31	1.46	1.40
2	D	201	HEC	C3C-C2C	5.13	1.46	1.40
2	C	201	HEC	C2B-C3B	5.10	1.46	1.40
2	D	201	HEC	C2B-C3B	4.19	1.45	1.40
2	B	201	HEC	C2B-C3B	3.76	1.44	1.40
2	C	201	HEC	C2A-C1A	3.41	1.50	1.42
2	D	201	HEC	C3D-C2D	3.40	1.47	1.37
2	D	201	HEC	C3C-C4C	3.34	1.49	1.43
2	C	201	HEC	C3C-C2C	3.29	1.44	1.40
2	B	201	HEC	CAD-C3D	-3.03	1.47	1.52
2	B	201	HEC	C3C-C4C	2.95	1.48	1.43
2	C	201	HEC	C3A-C4A	2.93	1.49	1.42
2	A	201	HEC	CAD-C3D	-2.93	1.47	1.52
2	C	201	HEC	C3C-C4C	2.93	1.48	1.43
2	D	201	HEC	C2A-C1A	2.87	1.49	1.42
2	C	201	HEC	C2A-C3A	2.75	1.45	1.37
2	A	201	HEC	C2A-C3A	2.71	1.45	1.37
2	B	201	HEC	C2A-C3A	2.65	1.45	1.37
2	A	201	HEC	C3A-C4A	2.63	1.48	1.42
2	A	201	HEC	C1B-NB	-2.54	1.30	1.36
2	C	201	HEC	C4B-C3B	2.53	1.47	1.43
2	D	201	HEC	C2A-C3A	2.51	1.45	1.37
2	A	201	HEC	C2A-C1A	2.49	1.48	1.42
2	D	201	HEC	C3A-C4A	2.47	1.48	1.42
2	B	201	HEC	C3D-C2D	2.46	1.45	1.37
2	B	201	HEC	C4D-ND	-2.43	1.31	1.36
2	B	201	HEC	C2A-C1A	2.41	1.48	1.42
2	A	201	HEC	C3D-C2D	2.41	1.44	1.37
2	B	201	HEC	C4B-C3B	2.30	1.47	1.43
2	C	201	HEC	C1C-NC	-2.29	1.31	1.36
2	D	201	HEC	C4B-C3B	2.22	1.47	1.43
2	C	201	HEC	C3D-C2D	2.17	1.44	1.37
2	D	201	HEC	C4D-CHA	2.15	1.47	1.41
2	A	201	HEC	C1D-CHD	2.03	1.46	1.41
2	A	201	HEC	CAA-C2A	-2.01	1.48	1.52

All (42) bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	201	HEC	C1D-C2D-C3D	-9.91	100.10	107.00
2	C	201	HEC	CMC-C2C-C3C	7.60	134.76	125.82
2	B	201	HEC	CBD-CAD-C3D	-7.34	100.10	112.62
2	C	201	HEC	CBD-CAD-C3D	-7.00	100.68	112.62
2	A	201	HEC	CMC-C2C-C3C	6.91	133.95	125.82
2	D	201	HEC	C1D-C2D-C3D	-6.88	102.21	107.00
3	C	202	EPE	O2S-S-C10	5.73	113.81	106.92
2	A	201	HEC	CBD-CAD-C3D	-5.47	103.28	112.62
2	A	201	HEC	CBA-CAA-C2A	-5.43	103.45	112.60
2	D	201	HEC	CMC-C2C-C3C	5.35	132.11	125.82
3	A	202	EPE	O3S-S-C10	5.07	113.97	105.77
3	C	202	EPE	O3S-S-C10	4.91	113.70	105.77
2	C	201	HEC	C1D-C2D-C3D	-4.70	103.72	107.00
2	A	201	HEC	C1D-C2D-C3D	-4.69	103.73	107.00
3	C	202	EPE	O1S-S-C10	-4.58	101.40	106.92
2	D	201	HEC	CBD-CAD-C3D	-4.41	105.10	112.62
2	B	201	HEC	CMB-C2B-C3B	4.22	130.79	125.82
2	B	201	HEC	CMC-C2C-C3C	4.22	130.78	125.82
3	B	203	EPE	O2S-S-C10	4.03	111.77	106.92
2	A	201	HEC	CMB-C2B-C3B	3.61	130.07	125.82
2	D	201	HEC	CBA-CAA-C2A	-3.59	106.56	112.60
2	B	201	HEC	CBA-CAA-C2A	-3.55	106.63	112.60
2	A	201	HEC	CMA-C3A-C2A	3.40	131.35	124.94
3	B	203	EPE	O1S-S-C10	3.20	110.77	106.92
2	A	201	HEC	C2B-C3B-C4B	-3.12	102.98	106.35
2	C	201	HEC	CBA-CAA-C2A	-3.03	107.50	112.60
2	D	201	HEC	CMB-C2B-C3B	2.69	128.98	125.82
2	B	201	HEC	CMA-C3A-C2A	2.62	129.88	124.94
2	D	201	HEC	CMA-C3A-C2A	2.52	129.70	124.94
2	C	201	HEC	CMA-C3A-C2A	2.52	129.70	124.94
2	A	201	HEC	CMD-C2D-C3D	2.49	129.64	124.94
2	C	201	HEC	CMC-C2C-C1C	-2.37	124.82	128.46
2	C	201	HEC	O1A-CGA-CBA	-2.34	115.56	123.08
3	B	202	EPE	C9-N1-C2	-2.10	105.85	111.23
2	D	201	HEC	CMD-C2D-C3D	2.09	128.88	124.94
2	D	201	HEC	O2A-CGA-CBA	2.08	120.71	114.03
2	C	201	HEC	CMB-C2B-C1B	2.07	131.65	128.46
3	B	202	EPE	O3S-S-C10	2.06	109.10	105.77
3	B	202	EPE	O2S-S-C10	2.03	109.36	106.92
2	B	201	HEC	CMD-C2D-C3D	2.02	128.76	124.94
2	C	201	HEC	CMB-C2B-C3B	2.01	128.18	125.82
3	C	202	EPE	C5-C6-N1	-2.00	106.53	110.64

There are no chirality outliers.

All (33) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	202	EPE	S-C10-C9-N1
3	A	202	EPE	C9-C10-S-O2S
3	B	203	EPE	C8-C7-N4-C5
3	C	202	EPE	C9-C10-S-O1S
4	A	203	GOL	C1-C2-C3-O3
4	B	204	GOL	O1-C1-C2-O2
4	B	204	GOL	O1-C1-C2-C3
3	C	202	EPE	N4-C7-C8-O8
3	A	202	EPE	C9-C10-S-O3S
3	A	202	EPE	N4-C7-C8-O8
4	A	203	GOL	O1-C1-C2-C3
4	A	204	GOL	O1-C1-C2-C3
4	A	204	GOL	C1-C2-C3-O3
4	A	203	GOL	O2-C2-C3-O3
3	C	202	EPE	C9-C10-S-O3S
3	B	203	EPE	N4-C7-C8-O8
4	A	203	GOL	O1-C1-C2-O2
3	C	202	EPE	C10-C9-N1-C2
3	C	202	EPE	C10-C9-N1-C6
4	A	204	GOL	O1-C1-C2-O2
3	A	202	EPE	C9-C10-S-O1S
3	C	202	EPE	C9-C10-S-O2S
3	A	202	EPE	C8-C7-N4-C3
2	C	201	HEC	C3D-CAD-CBD-CGD
4	A	204	GOL	O2-C2-C3-O3
3	B	202	EPE	C10-C9-N1-C6
3	B	203	EPE	C10-C9-N1-C6
3	A	202	EPE	C8-C7-N4-C5
2	D	201	HEC	CAA-CBA-CGA-O1A
2	B	201	HEC	CAD-CBD-CGD-O2D
3	B	202	EPE	C9-C10-S-O1S
2	B	201	HEC	CAA-CBA-CGA-O1A
3	B	203	EPE	C10-C9-N1-C2

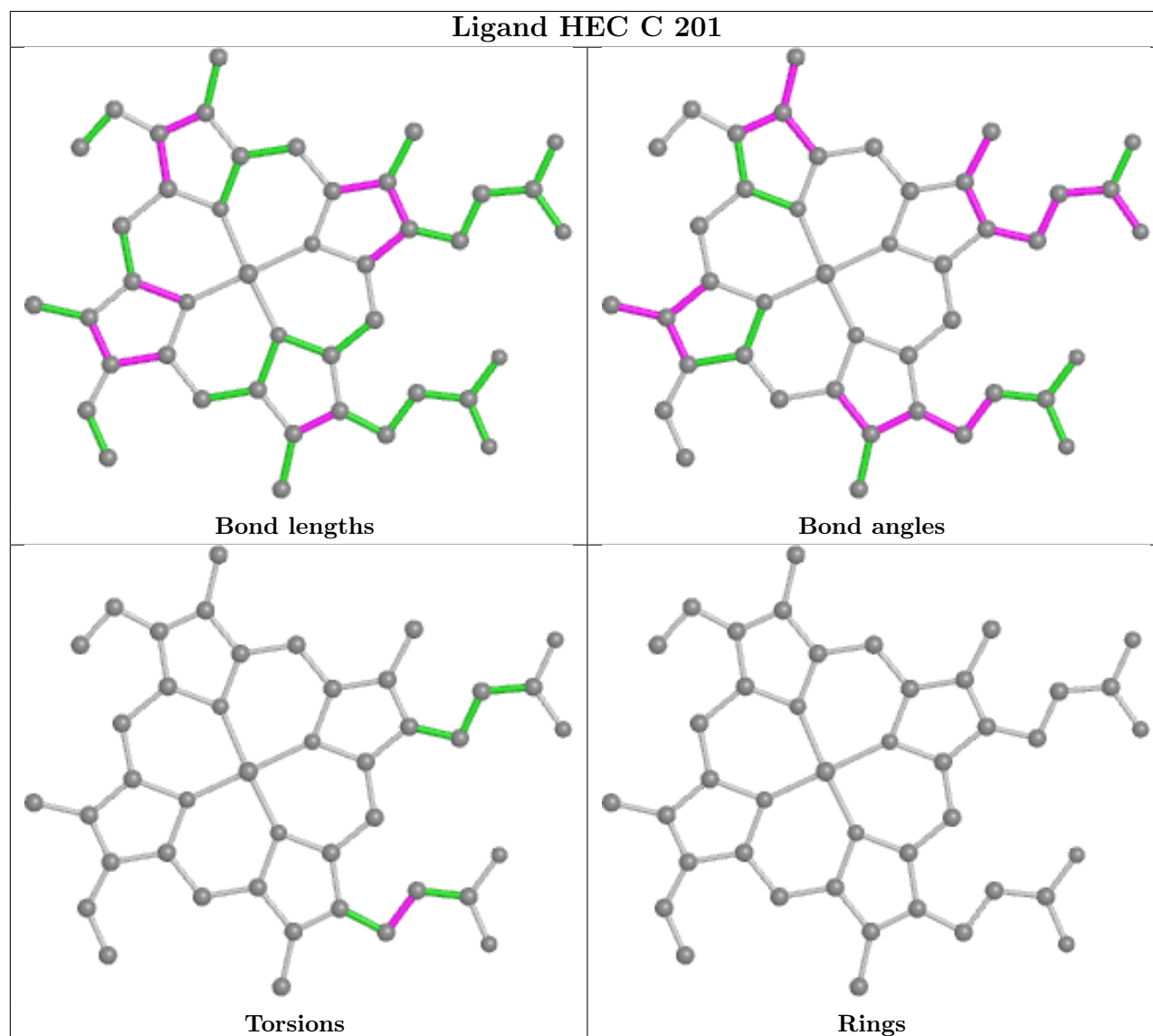
There are no ring outliers.

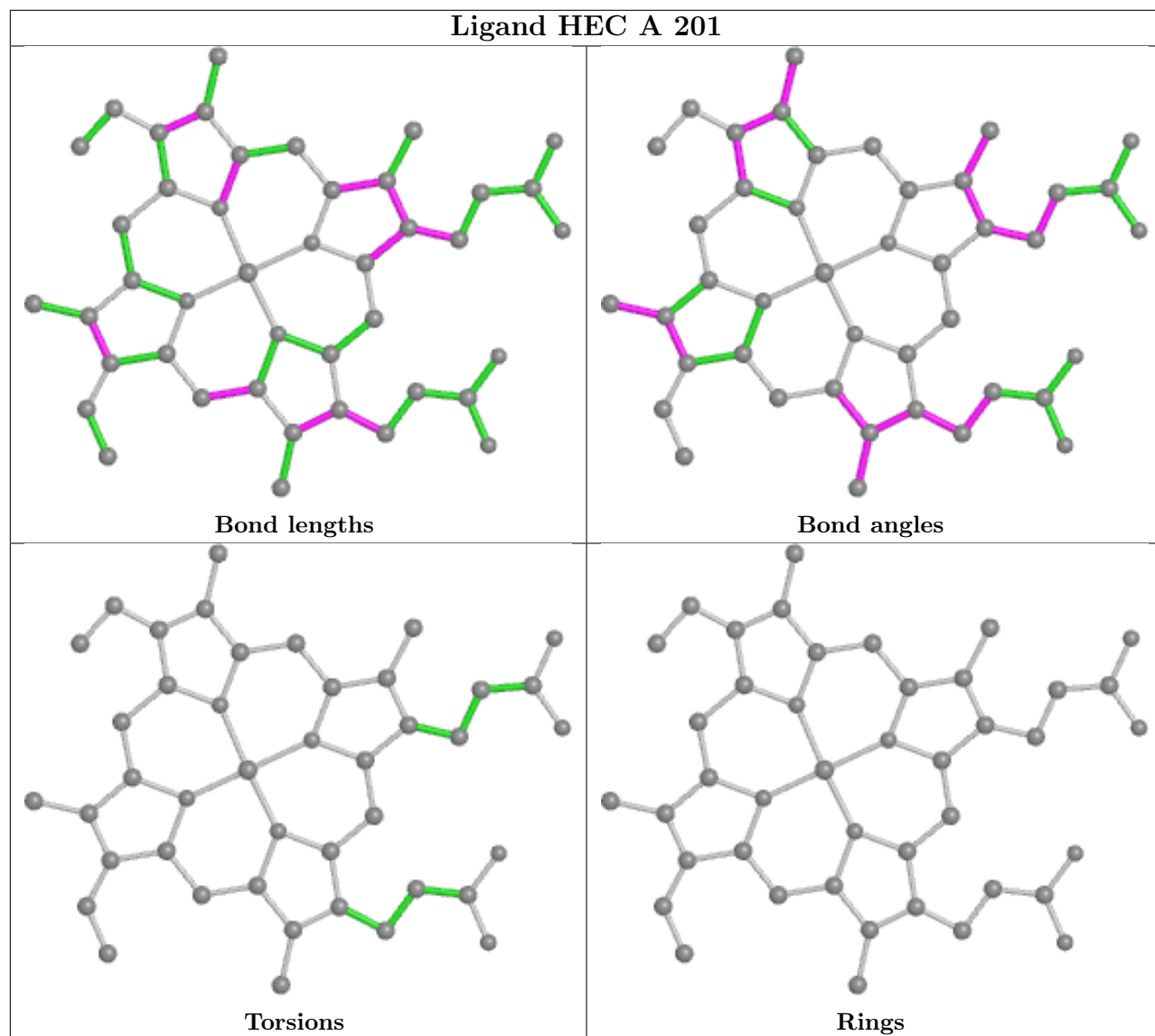
1 monomer is involved in 1 short contact:

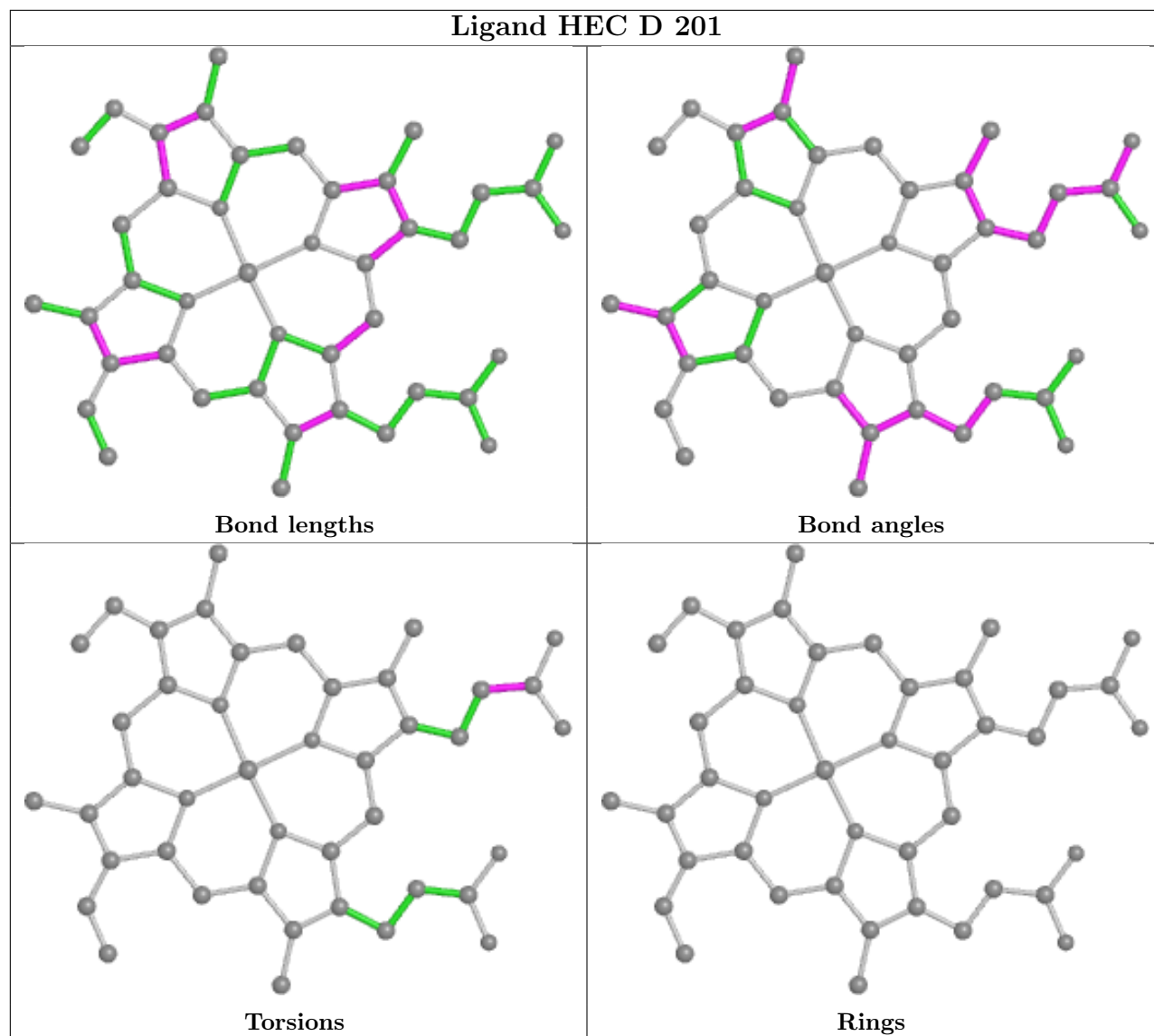
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	201	HEC	1	0

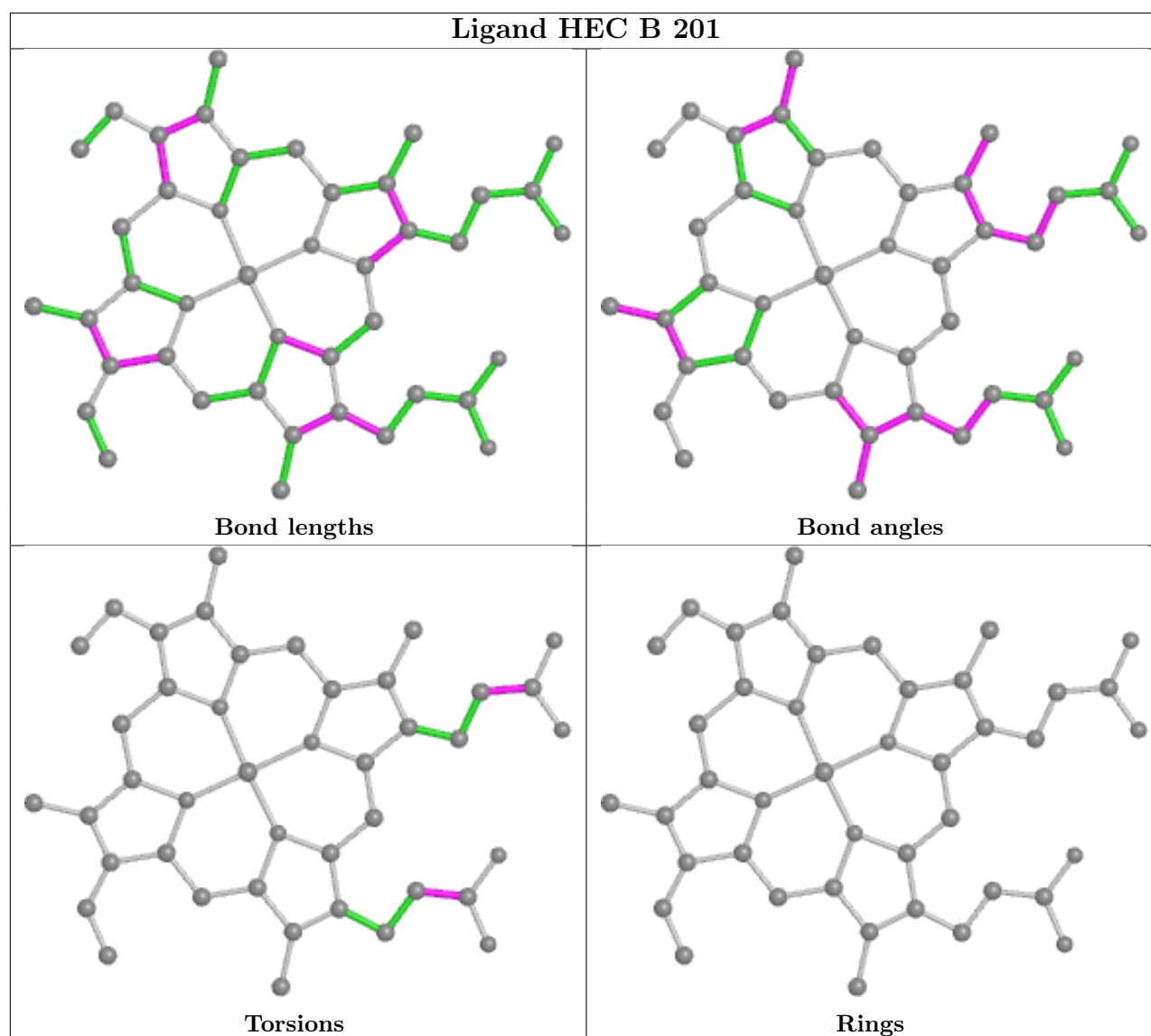
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

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.









## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	151/194 (77%)	-0.29	1 (0%) 84 86	20, 35, 72, 98	0
1	B	151/194 (77%)	-0.29	1 (0%) 84 86	21, 33, 62, 85	0
1	C	145/194 (74%)	-0.13	2 (1%) 73 77	25, 36, 68, 103	0
1	D	150/194 (77%)	-0.27	1 (0%) 84 86	23, 35, 69, 94	0
All	All	597/776 (76%)	-0.24	5 (0%) 82 85	20, 35, 69, 103	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	182	PHE	3.5
1	A	33	GLN	2.2
1	C	47	LEU	2.1
1	B	34	LEU	2.1
1	C	158	SER	2.1

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [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, 95<sup>th</sup> 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.

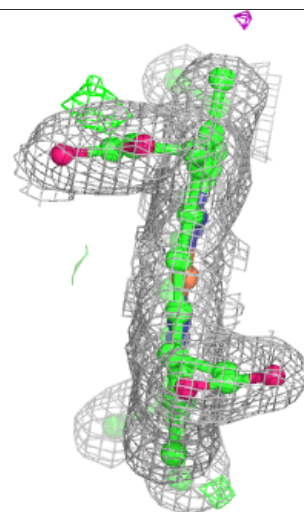
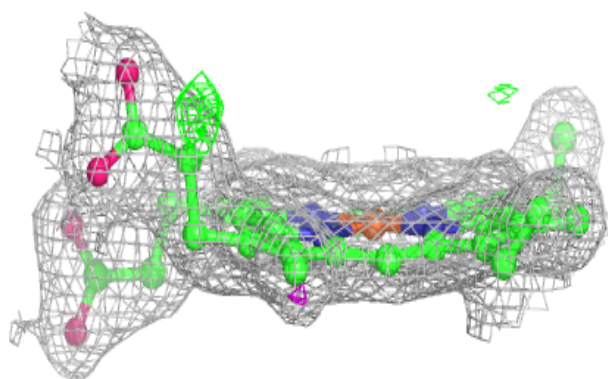
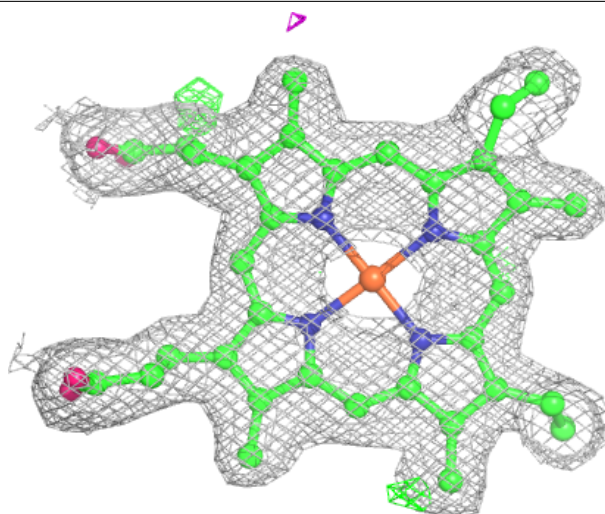
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
3	EPE	B	203	15/15	0.82	0.15	85,97,108,109	0
4	GOL	A	204	6/6	0.83	0.13	64,71,75,85	0
3	EPE	C	202	15/15	0.84	0.16	60,89,116,117	0
4	GOL	B	204	6/6	0.84	0.16	74,80,82,83	0
3	EPE	A	202	15/15	0.85	0.12	64,68,87,88	0
4	GOL	A	203	6/6	0.85	0.12	72,73,77,78	0
6	FE	B	206	1/1	0.92	0.07	90,90,90,90	0
3	EPE	B	202	15/15	0.95	0.11	38,62,85,88	0
6	FE	D	203	1/1	0.95	0.07	92,92,92,92	0
7	NA	A	207	1/1	0.95	0.05	23,23,23,23	0
5	CA	A	205	1/1	0.96	0.06	61,61,61,61	0
7	NA	C	206	1/1	0.96	0.07	31,31,31,31	0
5	CA	D	202	1/1	0.97	0.05	66,66,66,66	0
5	CA	B	205	1/1	0.97	0.05	64,64,64,64	0
7	NA	A	208	1/1	0.97	0.05	31,31,31,31	0
6	FE	C	204	1/1	0.97	0.07	82,82,82,82	0
7	NA	D	205	1/1	0.97	0.05	24,24,24,24	0
6	FE	A	206	1/1	0.98	0.08	80,80,80,80	0
2	HEC	A	201	43/43	0.98	0.05	19,21,28,33	0
7	NA	C	205	1/1	0.98	0.03	28,28,28,28	0
5	CA	C	203	1/1	0.98	0.04	66,66,66,66	0
7	NA	D	204	1/1	0.98	0.06	21,21,21,21	0
2	HEC	C	201	43/43	0.98	0.06	21,23,33,41	0
2	HEC	B	201	43/43	0.99	0.05	18,21,28,30	0
7	NA	B	208	1/1	0.99	0.03	17,17,17,17	0
2	HEC	D	201	43/43	0.99	0.04	20,22,28,31	0
7	NA	B	207	1/1	1.00	0.04	18,18,18,18	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.



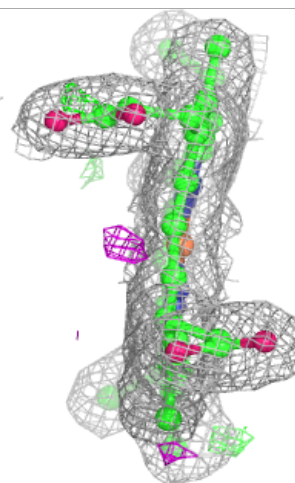
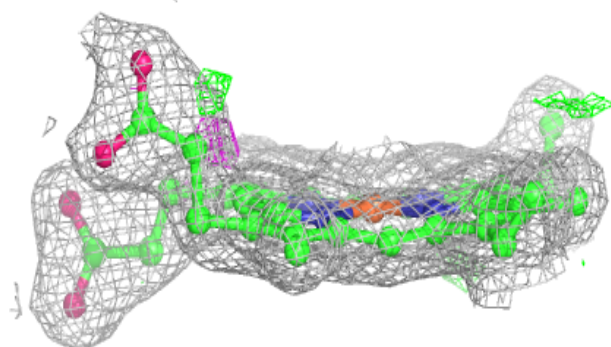
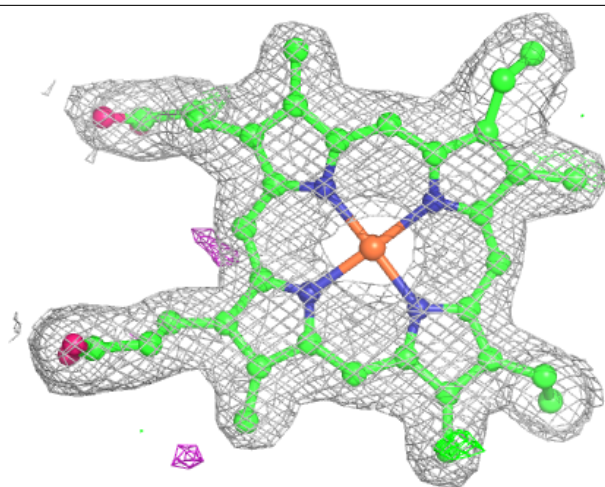
**Electron density around HEC A 201:**

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



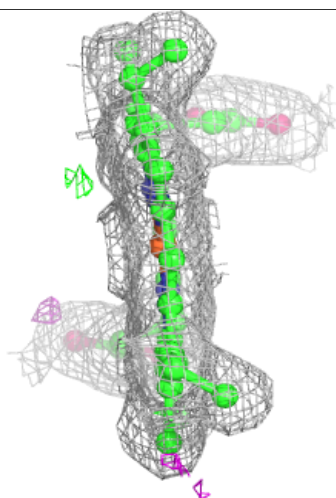
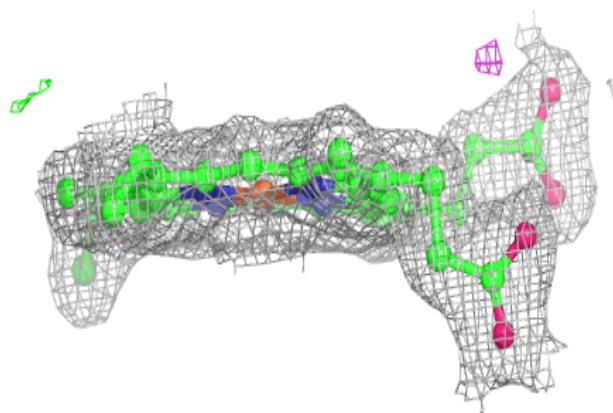
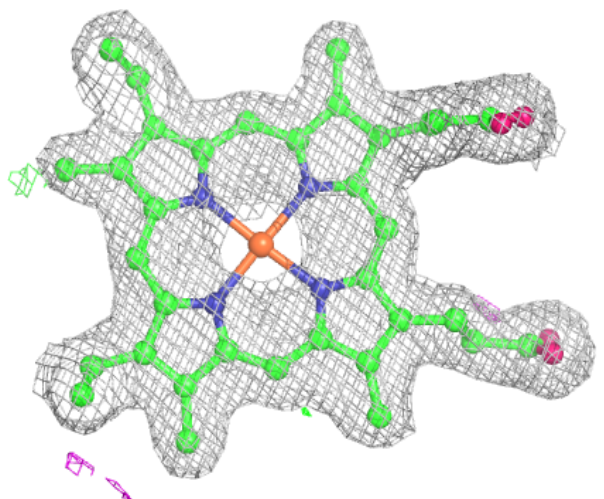
**Electron density around HEC C 201:**

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



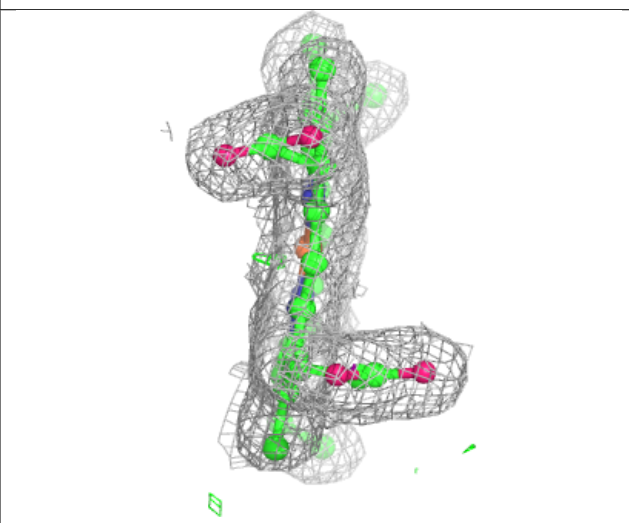
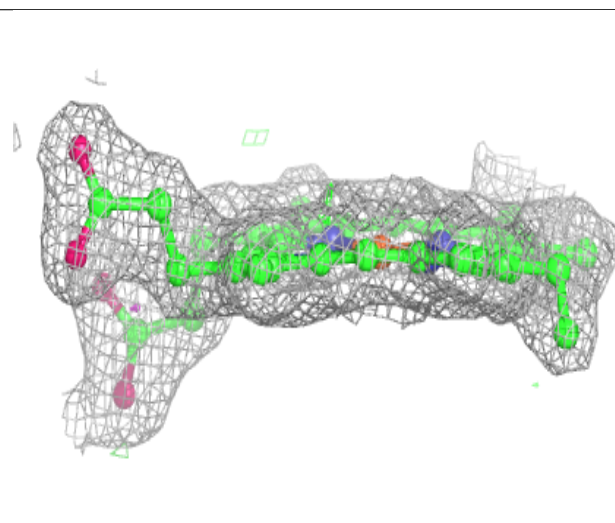
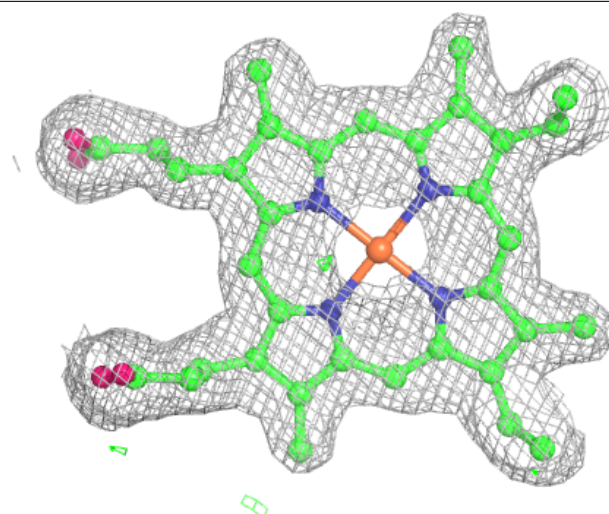
**Electron density around HEC B 201:**

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



**Electron density around HEC D 201:**

$2mF_o-DF_c$  (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.