



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 22, 2020 – 08:17 PM BST

PDB ID : 7PRC
Title : PHOTOSYNTHETIC REACTION CENTER FROM RHODOPSEUDOMONAS VIRIDIS (DG-420315 (TRIAZINE) COMPLEX)
Authors : Lancaster, C.R.D.; Michel, H.
Deposited on : 1997-08-01
Resolution : 2.65 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

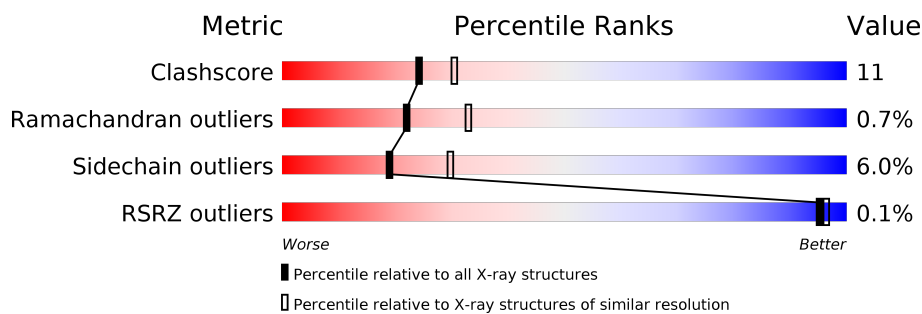
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION





The reported resolution of this entry is 2.65 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))
Clashscore	141614	1374 (2.68-2.64)
Ramachandran outliers	138981	1349 (2.68-2.64)
Sidechain outliers	138945	1349 (2.68-2.64)
RSRZ outliers	127900	1318 (2.68-2.64)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	C	336	 76% 20% . .
2	L	273	 78% 21% .
3	M	323	 76% 22% .
4	H	258	 70% 26% .

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
7	BPB	L	402	X	-	-	-

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 10476 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 PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	C	332	Total	C	N	O	S	51	1	0
			2607	1642	467	480	18			

- Molecule 2 is a protein called PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	273	Total	C	N	O	S	14	0	0
			2171	1459	350	355	7			

- Molecule 3 is a protein called PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	323	Total	C	N	O	S	19	2	0
			2577	1720	421	425	11			

- Molecule 4 is a protein called PHOTOSYNTHETIC REACTION CENTER.

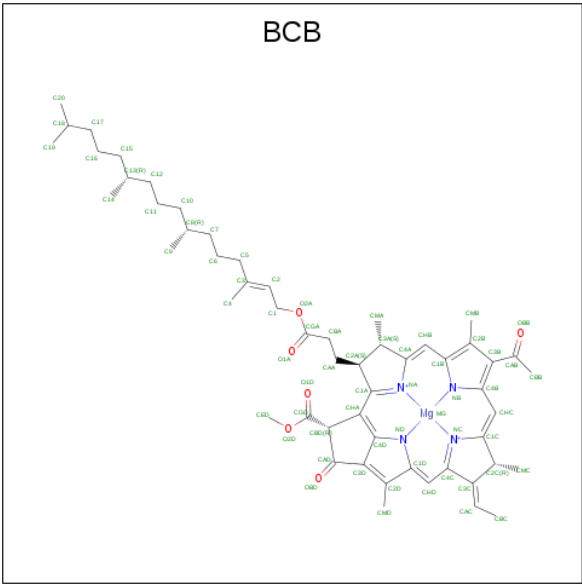
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	258	Total	C	N	O	S	122	0	0
			2018	1292	344	380	2			

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



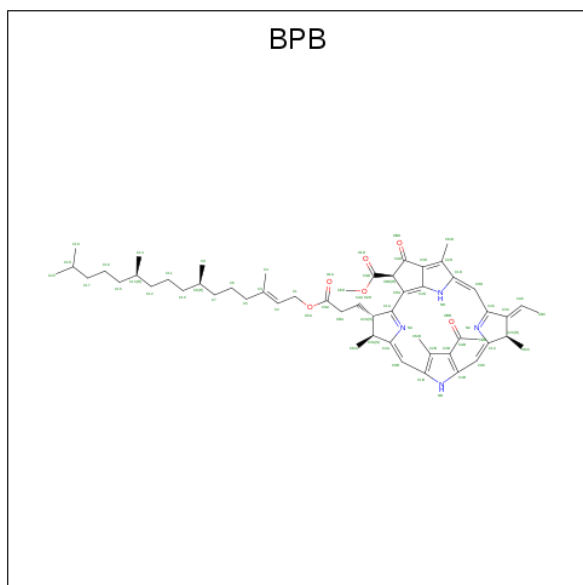
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
5	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 6 is BACTERIOCHLOROPHYLL B (three-letter code: BCB) (formula: $C_{55}H_{72}MgN_4O_6$).



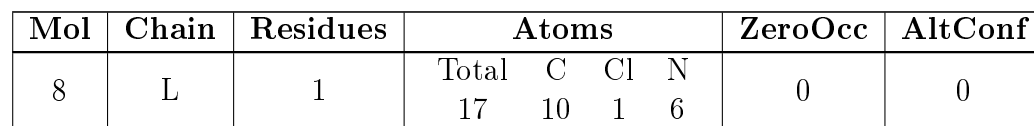
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	L	1	Total 66	C 55	Mg 1	N 4	O 6	0	0
6	L	1	Total 66	C 55	Mg 1	N 4	O 6	0	0
6	M	1	Total 66	C 55	Mg 1	N 4	O 6	0	0
6	M	1	Total 66	C 55	Mg 1	N 4	O 6	0	0

- Molecule 7 is BACTERIOPHEOPHYTIN B (three-letter code: BPB) (formula: $C_{55}H_{74}N_4O_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	L	1	Total	C	N	O	7	0
			65	55	4	6		
7	L	1	Total	C	N	O	0	0
			65	55	4	6		

- Molecule 8 is 2-CHLORO-4-ETHYLAMINO-6-(R(+)-2'-CYANO-4-BUTYLAMINO)-1,3,5-TRIAZINE (three-letter code: CET) (formula: $C_{10}H_{15}ClN_6$).



- # LDA

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	L	1	Total 16	C 14	N 1	O 1	0	0
9	M	1	Total 16	C 14	N 1	O 1	0	0



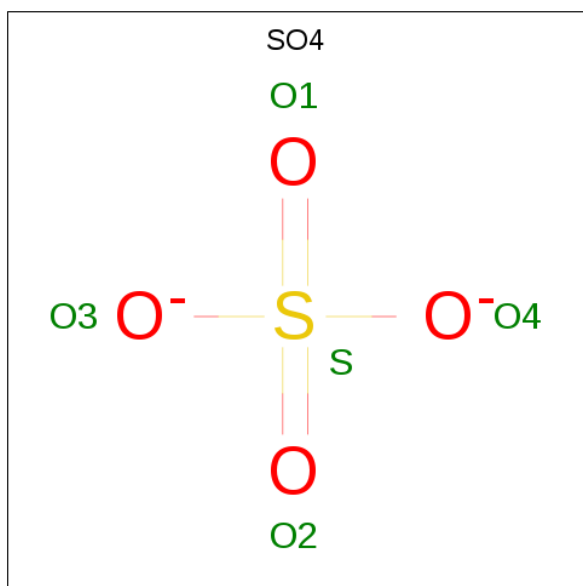
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	M	1	Total	C	N	O	0	0
			16	14	1	1		
9	M	1	Total	C	N	O	5	0
			16	14	1	1		
9	M	1	Total	C	N	O	4	0
			16	14	1	1		
9	H	1	Total	C	N	O	0	0
			16	14	1	1		

- Molecule 10 is FE (II) ION (three-letter code: FE2) (formula: Fe).

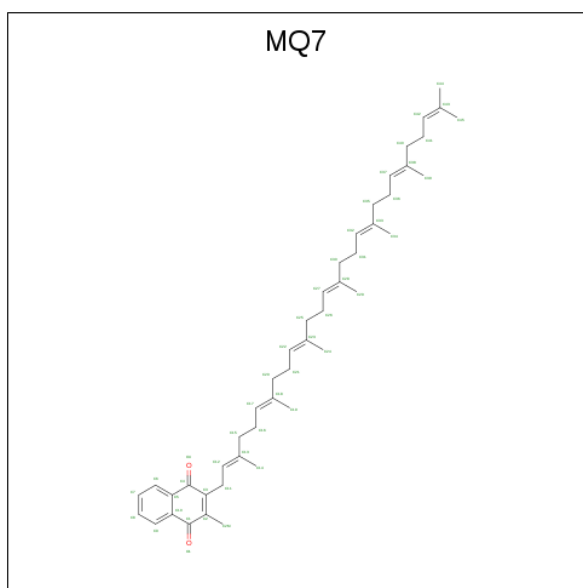
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	M	1	Total	Fe	0	0
			1	1		

- Molecule 11 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



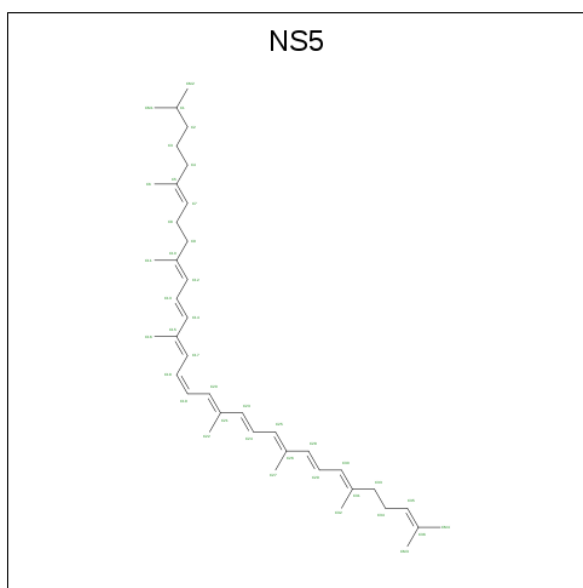
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
11	M	1	Total	O	S	0	0
			5	4	1		
11	M	1	Total	O	S	0	0
			5	4	1		
11	M	1	Total	O	S	0	0
			5	4	1		
11	H	1	Total	O	S	0	0
			5	4	1		

- Molecule 12 is MENAQUINONE-7 (three-letter code: MQ7) (formula: $C_{46}H_{64}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	M	1	Total	C	O	0	0
			48	46	2		

- Molecule 13 is 15-cis-1,2-dihydroneurosporene (three-letter code: NS5) (formula: $C_{40}H_{60}$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	M	1	Total	C	4	0
			40	40		

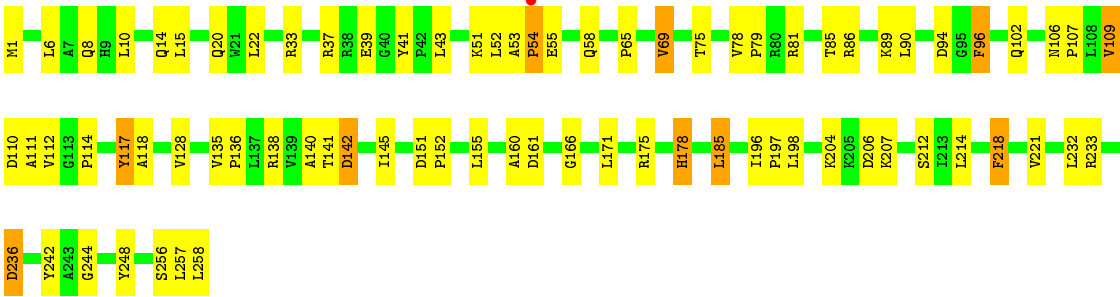
- Molecule 14 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	C	118	Total 118	O 118	0	0
14	L	55	Total 55	O 55	0	0
14	M	78	Total 78	O 78	0	0
14	H	64	Total 64	O 64	0	0

Chain H:

70%

26%



4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	223.50 Å 223.50 Å 113.60 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.65 29.74 – 2.60	Depositor EDS
% Data completeness (in resolution range)	88.3 (10.00-2.65) 87.0 (29.74-2.60)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.59 (at 2.61 Å)	Xtriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.184 , 0.231 0.180 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	23.1	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 90.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10476	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.19% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: LDA, BPB, BCB, CET, FE2, SO4, MQ7, HEM, FME, NS5

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	C	0.55	0/2674	0.67	2/3645 (0.1%)
2	L	0.58	0/2259	0.62	0/3084
3	M	0.58	0/2683	0.62	0/3669
4	H	0.67	2/2055 (0.1%)	0.86	4/2807 (0.1%)
All	All	0.59	2/9671 (0.0%)	0.69	6/13205 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	4
2	L	0	7
3	M	0	5
4	H	0	2
All	All	0	18

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	54	PRO	C-N	14.77	1.68	1.34
4	H	52	LEU	C-N	-6.29	1.19	1.34

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	52	LEU	O-C-N	-18.31	93.41	122.70
4	H	52	LEU	CA-C-N	13.07	145.95	117.20
4	H	52	LEU	C-N-CA	8.96	144.09	121.70

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	53	ALA	CB-CA-C	-6.78	99.92	110.10
1	C	50	PRO	CB-CA-C	-5.63	97.93	112.00

There are no chirality outliers.

5 of 18 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	15	ARG	Sidechain
1	C	187	TYR	Sidechain
1	C	202	ARG	Sidechain
1	C	227	TYR	Sidechain
2	L	85	PHE	Sidechain

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2607	0	2575	54	0
2	L	2171	0	2098	46	0
3	M	2577	0	2468	60	0
4	H	2018	0	2019	51	0
5	C	172	0	120	6	0
6	L	132	0	144	10	0
6	M	132	0	144	14	0
7	L	130	0	148	18	0
8	L	17	0	15	1	0
9	H	16	0	31	1	0
9	L	16	0	31	0	0
9	M	64	0	124	5	0
10	M	1	0	0	0	0
11	H	5	0	0	0	0
11	M	15	0	0	2	0
12	M	48	0	64	1	0
13	M	40	0	60	7	0
14	C	118	0	0	3	0
14	H	64	0	0	0	0
14	L	55	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
14	M	78	0	0	2	0
All	All	10476	0	10041	216	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 11.

The worst 5 of 216 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)
4:H:54:PRO:C	4:H:55:GLU:N	1.68	1.44
7:L:402:BPB:HBBB	7:L:402:BPB:HHC	1.54	0.90
6:M:805:BCB:HBB2	6:M:805:BCB:HHC	1.57	0.85
7:L:402:BPB:HHC	7:L:402:BPB:CBB	2.09	0.83
2:L:181:PHE:HB3	7:L:401:BPB:CBB	2.10	0.81

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	C	331/336 (98%)	314 (95%)	14 (4%)	3 (1%)	17	26
2	L	271/273 (99%)	255 (94%)	14 (5%)	2 (1%)	22	33
3	M	323/323 (100%)	308 (95%)	12 (4%)	3 (1%)	17	26
4	H	256/258 (99%)	241 (94%)	15 (6%)	0	100	100
All	All	1181/1190 (99%)	1118 (95%)	55 (5%)	8 (1%)	22	33

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	M	51	LEU

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Mol	Chain	Res	Type
3	M	193	ASN
2	L	57	PRO
1	C	67	GLU
2	L	23	ASP

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	C	281/282 (100%)	263 (94%)	18 (6%)	17	27
2	L	218/218 (100%)	212 (97%)	6 (3%)	43	61
3	M	251/249 (101%)	238 (95%)	13 (5%)	23	36
4	H	212/212 (100%)	190 (90%)	22 (10%)	7	10
All	All	962/961 (100%)	903 (94%)	59 (6%)	19	29

5 of 59 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	M	40	LYS
3	M	181	ILE
4	H	212	SER
3	M	71	PHE
3	M	114	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	L	183	ASN
2	L	214	GLN
2	L	239	ASN
3	M	147	ASN
4	H	102	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	FME	H	1	4	8,9,10	0.61	0	7,9,11	2.81	2 (28%)

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	FME	H	1	4	-	3/7/9/11	-

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1	FME	O1-CN-N	-5.63	110.45	125.27
4	H	1	FME	CA-N-CN	-4.04	116.60	122.82

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	1	FME	O1-CN-N-CA
4	H	1	FME	CB-CG-SD-CE
4	H	1	FME	CA-CB-CG-SD

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 24 ligands modelled in this entry, 1 is monoatomic - leaving 23 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	$\# Z > 2$
9	LDA	M	706	-	12,15,15	2.19	1 (8%)	14,17,17	0.50	0
8	CET	L	502	-	16,17,17	1.34	2 (12%)	21,23,23	0.89	1 (4%)
5	HEM	C	340	1	27,50,50	1.94	8 (29%)	17,82,82	1.48	3 (17%)
9	LDA	M	701	-	12,15,15	2.49	1 (8%)	14,17,17	0.71	0
9	LDA	H	703	-	12,15,15	2.27	1 (8%)	14,17,17	0.68	0
5	HEM	C	337	1	27,50,50	1.79	5 (18%)	17,82,82	0.94	1 (5%)
12	MQ7	M	501	-	49,49,49	1.49	12 (24%)	60,63,63	1.74	12 (20%)
5	HEM	C	338	1	27,50,50	1.54	4 (14%)	17,82,82	1.27	2 (11%)
9	LDA	M	704	-	12,15,15	1.93	1 (8%)	14,17,17	0.54	0
9	LDA	M	705	-	12,15,15	2.19	1 (8%)	14,17,17	0.50	0
11	SO4	M	802	-	4,4,4	0.57	0	6,6,6	0.83	0
9	LDA	L	702	-	12,15,15	2.27	1 (8%)	14,17,17	0.54	0
7	BPB	L	402	-	64,70,70	1.39	9 (14%)	64,101,101	1.63	9 (14%)
6	BCB	M	805	3	60,74,74	3.77	25 (41%)	48,115,115	2.87	18 (37%)
11	SO4	H	801	-	4,4,4	0.55	0	6,6,6	0.33	0
11	SO4	M	803	-	4,4,4	0.67	0	6,6,6	0.78	0
13	NS5	M	600	-	39,39,39	0.81	1 (2%)	44,46,46	1.20	6 (13%)
11	SO4	M	804	-	4,4,4	0.90	0	6,6,6	0.70	0
6	BCB	L	304	2	60,74,74	3.86	24 (40%)	48,115,115	2.56	17 (35%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	BPB	L	401	-	64,70,70	1.21	8 (12%)	64,101,101	2.04	14 (21%)
6	BCB	L	302	2	60,74,74	3.68	21 (35%)	48,115,115	2.38	14 (29%)
6	BCB	M	806	3	60,74,74	3.58	22 (36%)	48,115,115	2.92	16 (33%)
5	HEM	C	339	1	27,50,50	1.76	6 (22%)	17,82,82	1.66	5 (29%)

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
9	LDA	M	706	-	-	4/13/13/13	-
8	CET	L	502	-	-	0/11/14/14	0/1/1/1
12	MQ7	M	501	-	-	2/41/61/61	0/2/2/2
5	HEM	C	338	1	-	0/6/54/54	-
5	HEM	C	339	1	-	0/6/54/54	-
7	BPB	L	401	-	-	8/47/105/105	0/5/6/6
5	HEM	C	337	1	-	0/6/54/54	-
9	LDA	M	704	-	-	6/13/13/13	-
6	BCB	M	805	3	-	12/41/177/177	-
5	HEM	C	340	1	-	0/6/54/54	-
9	LDA	M	705	-	-	1/13/13/13	-
9	LDA	M	701	-	-	5/13/13/13	-
13	NS5	M	600	-	-	14/43/43/43	-
6	BCB	L	304	2	-	11/41/177/177	-
9	LDA	L	702	-	-	4/13/13/13	-
9	LDA	H	703	-	-	3/13/13/13	-
6	BCB	L	302	2	-	9/41/177/177	-
6	BCB	M	806	3	-	14/41/177/177	-
7	BPB	L	402	-	1/1/18/23	5/47/105/105	0/5/6/6

The worst 5 of 153 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	304	BCB	C1A-CHA	-10.41	1.37	1.54
6	L	302	BCB	C1A-CHA	-9.79	1.38	1.54
6	M	805	BCB	CHB-C1B	-9.29	1.38	1.53
6	L	304	BCB	CHD-C4C	-9.27	1.37	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	M	805	BCB	CHD-C1D	-9.09	1.39	1.53

The worst 5 of 118 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	M	806	BCB	CMB-C2B-C3B	8.96	136.54	114.29
6	M	805	BCB	CBB-CAB-C3B	8.60	125.58	116.80
7	L	401	BPB	O2D-CGD-CBD	7.52	124.63	111.27
6	M	806	BCB	OBB-CAB-C3B	7.42	129.34	121.52
6	M	805	BCB	C1D-CHD-C4C	7.41	128.16	112.37

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
7	L	402	BPB	C13

5 of 98 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
9	M	706	LDA	N1-C1-C2-C3
9	M	704	LDA	C2-C1-N1-CM1
9	M	704	LDA	N1-C1-C2-C3
6	M	805	BCB	C3A-C2A-CAA-CBA
6	M	805	BCB	C2-C3-C5-C6

There are no ring outliers.

19 monomers are involved in 61 short contacts:

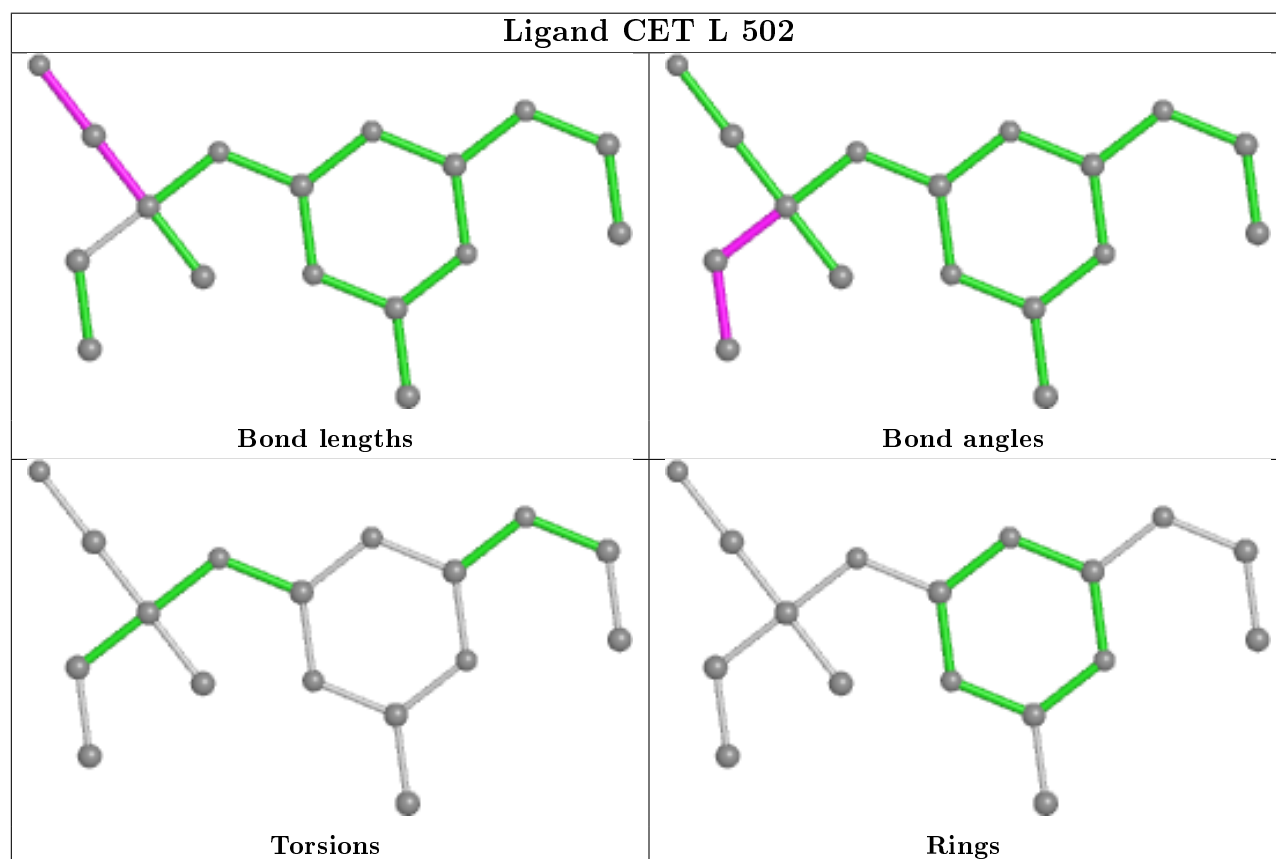
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	M	706	LDA	1	0
8	L	502	CET	1	0
9	M	701	LDA	2	0
9	H	703	LDA	1	0
5	C	337	HEM	3	0
12	M	501	MQ7	1	0
5	C	338	HEM	2	0
9	M	704	LDA	2	0
9	M	705	LDA	1	0
7	L	402	BPB	5	0
6	M	805	BCB	8	0
11	M	803	SO4	1	0

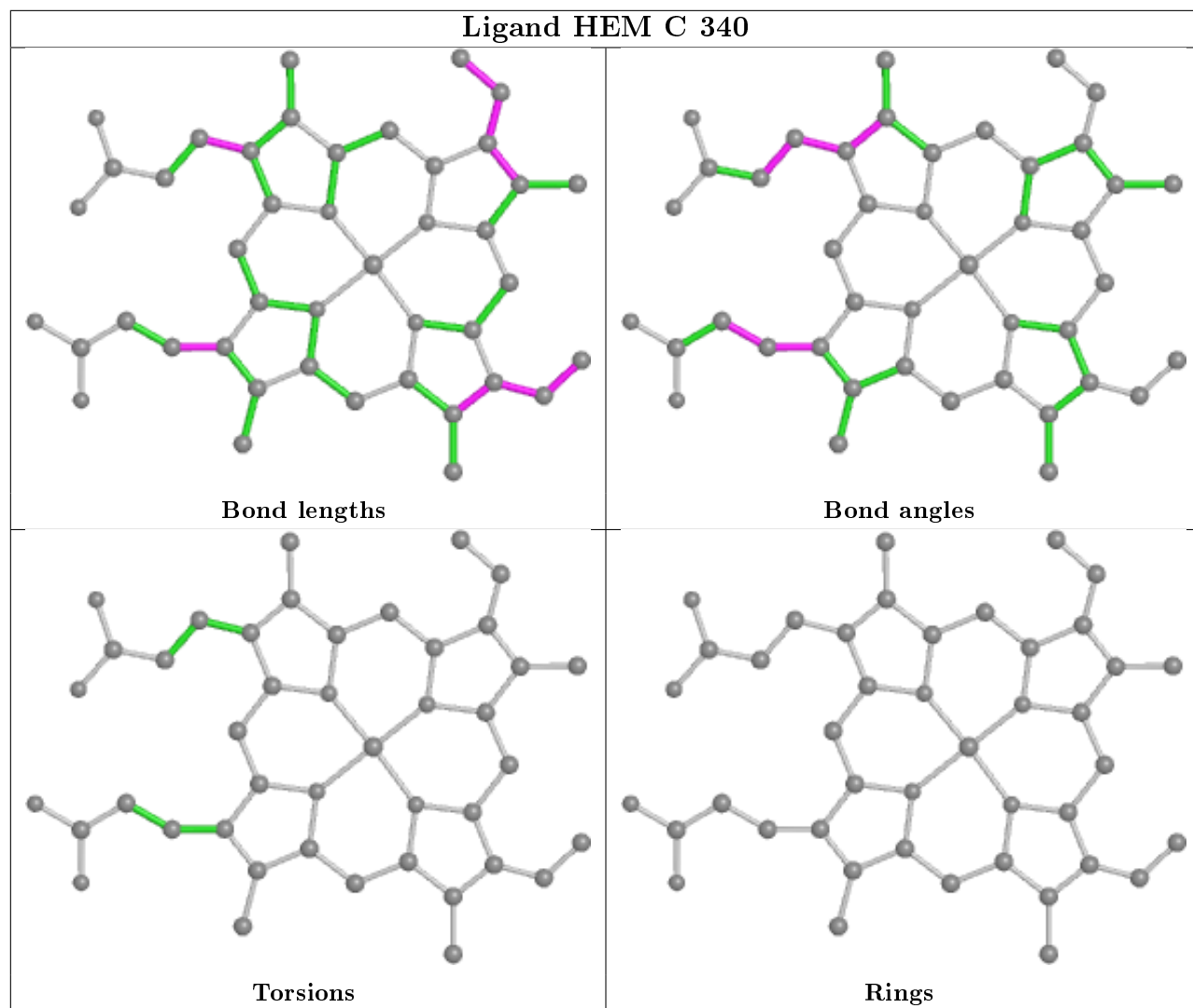
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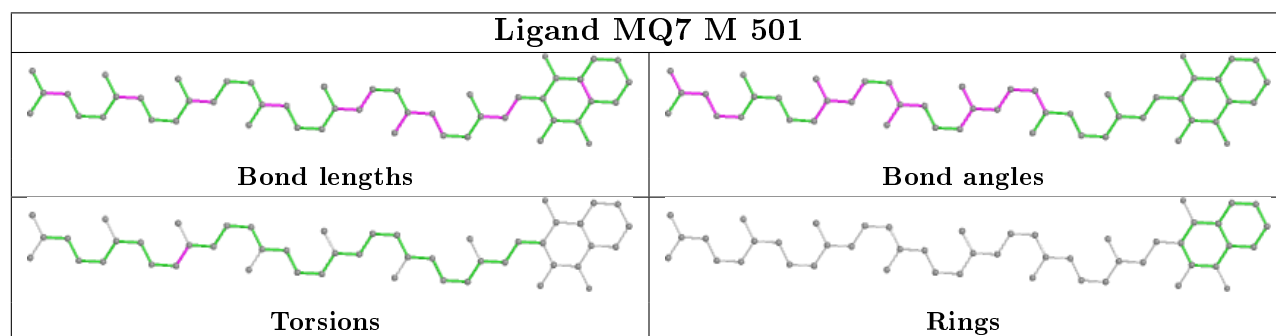
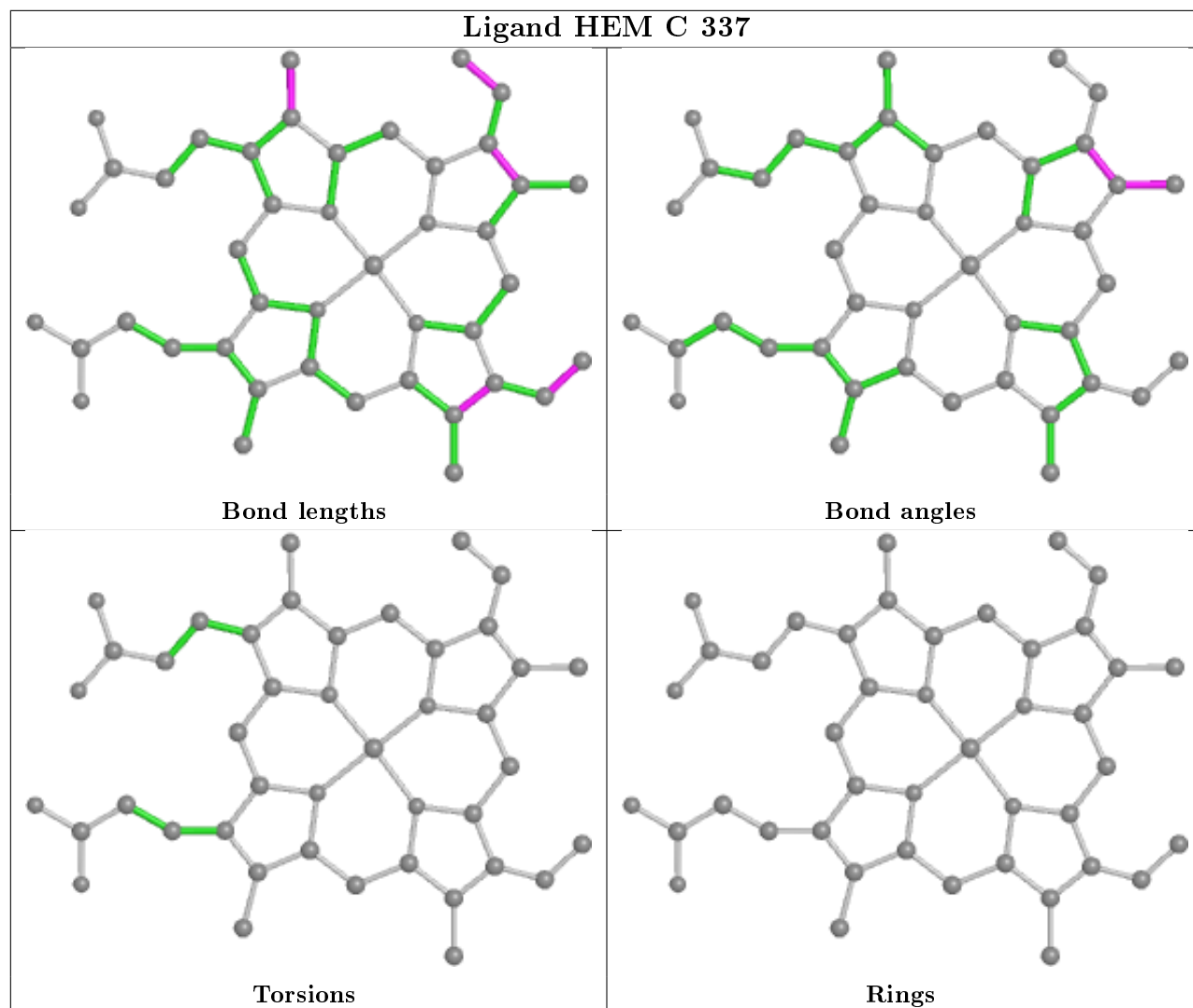
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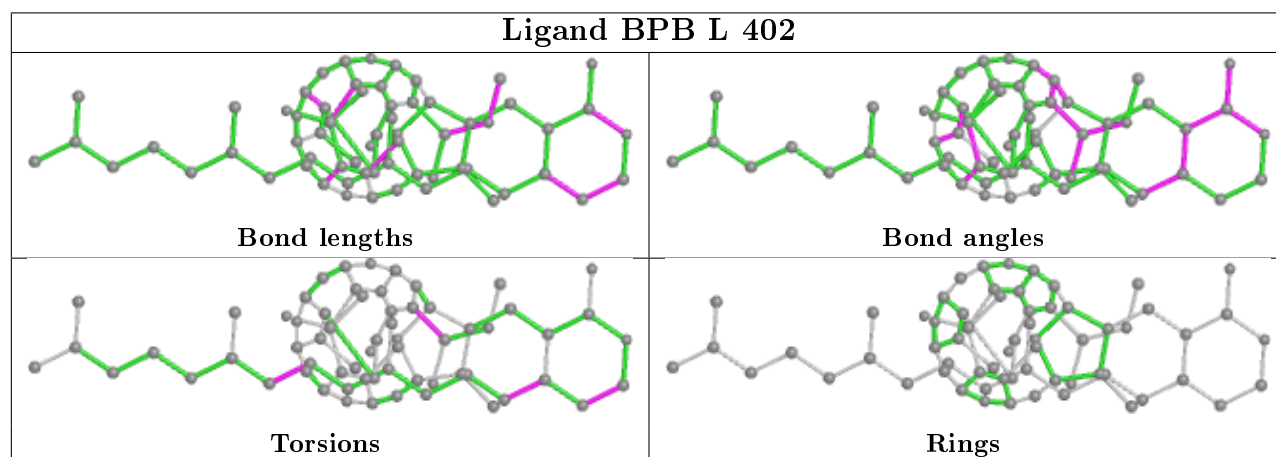
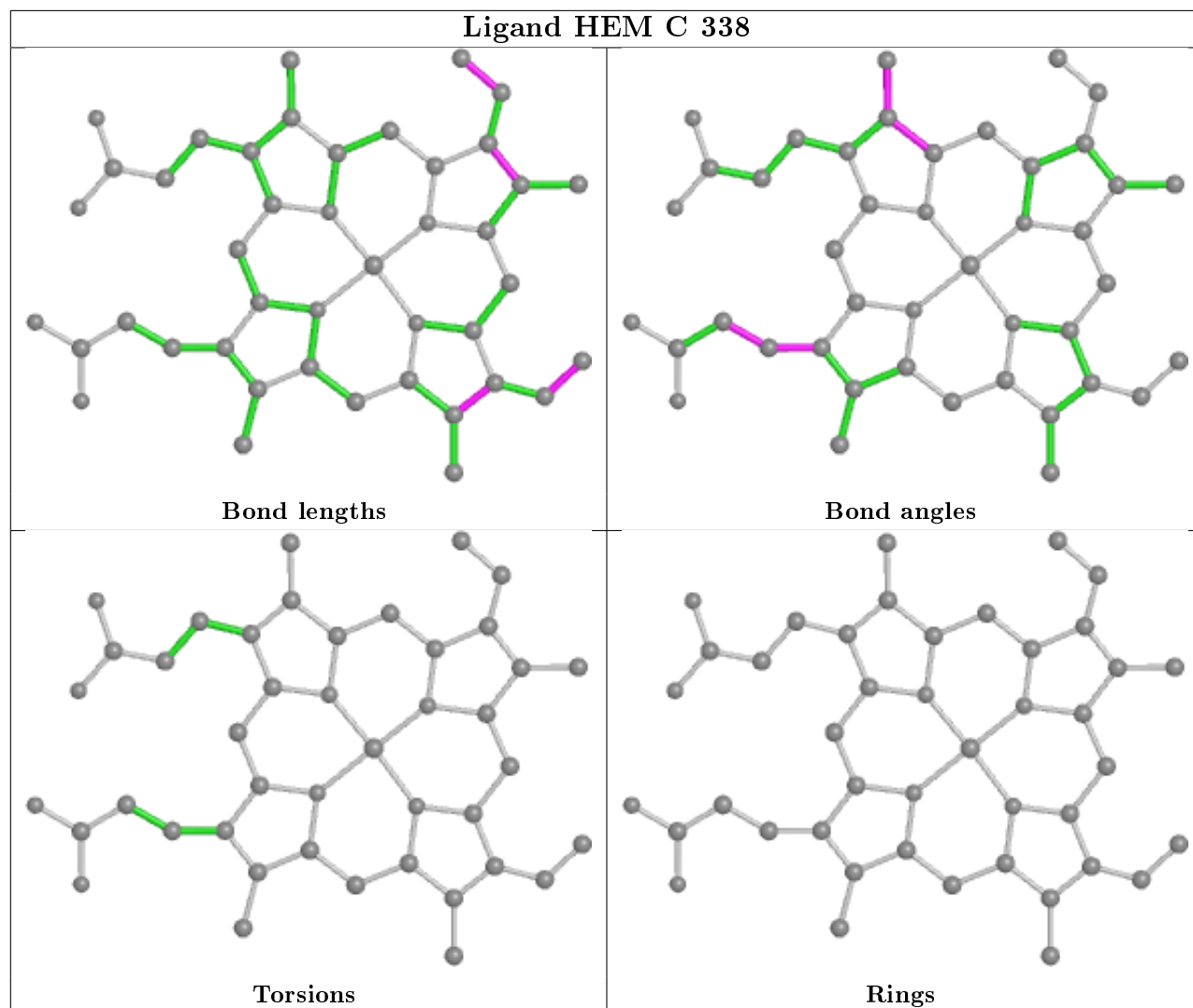
Mol	Chain	Res	Type	Clashes	Symm-Clashes
13	M	600	NS5	7	0
11	M	804	SO4	1	0
6	L	304	BCB	4	0
7	L	401	BPB	13	0
6	L	302	BCB	6	0
6	M	806	BCB	7	0
5	C	339	HEM	1	0

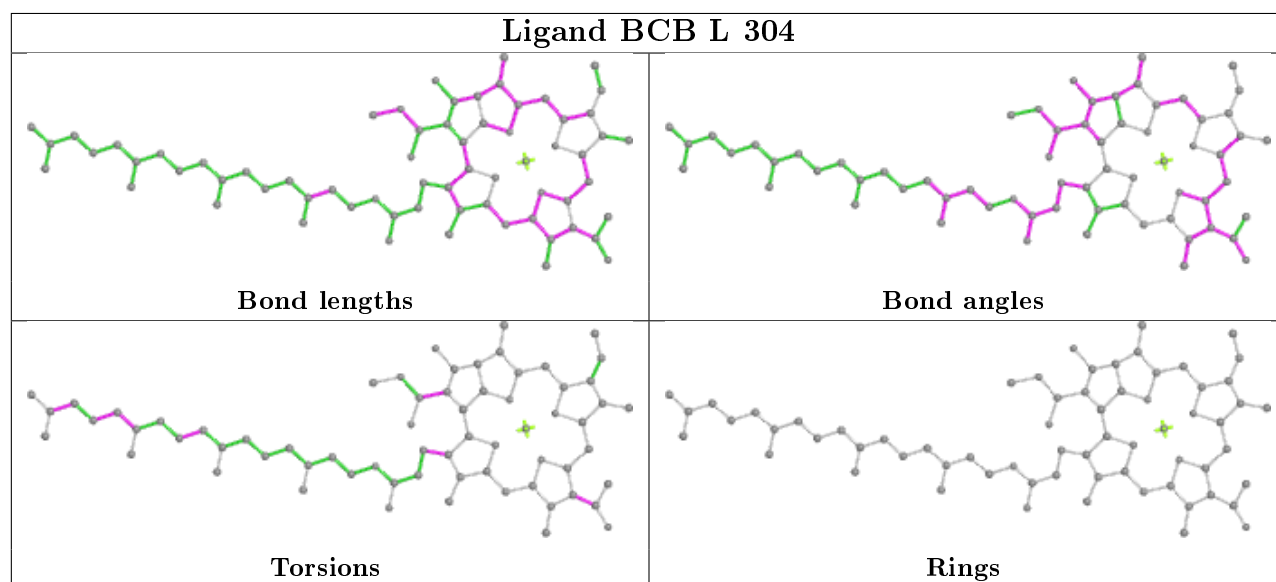
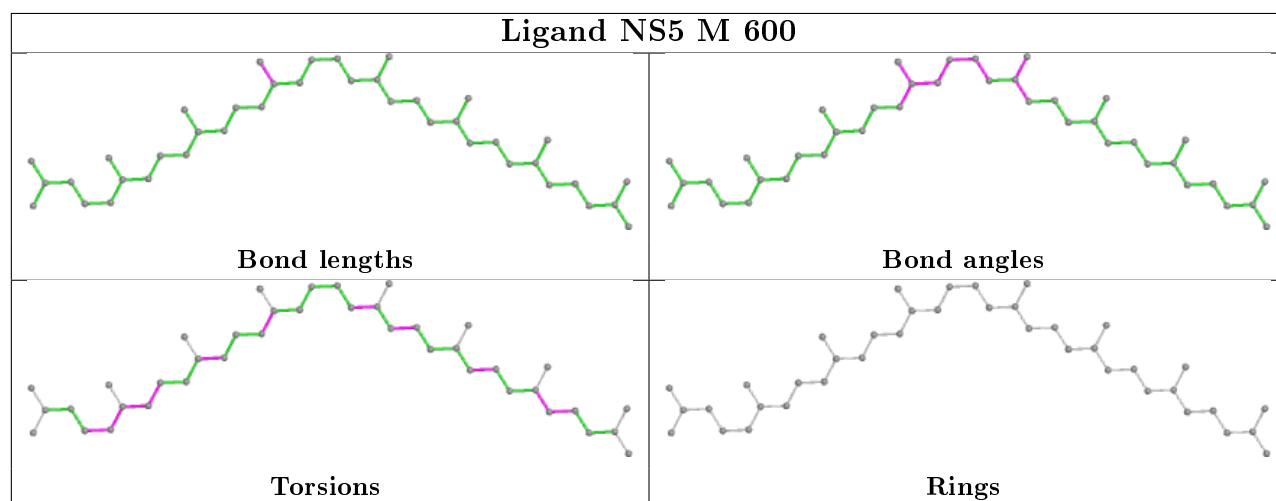
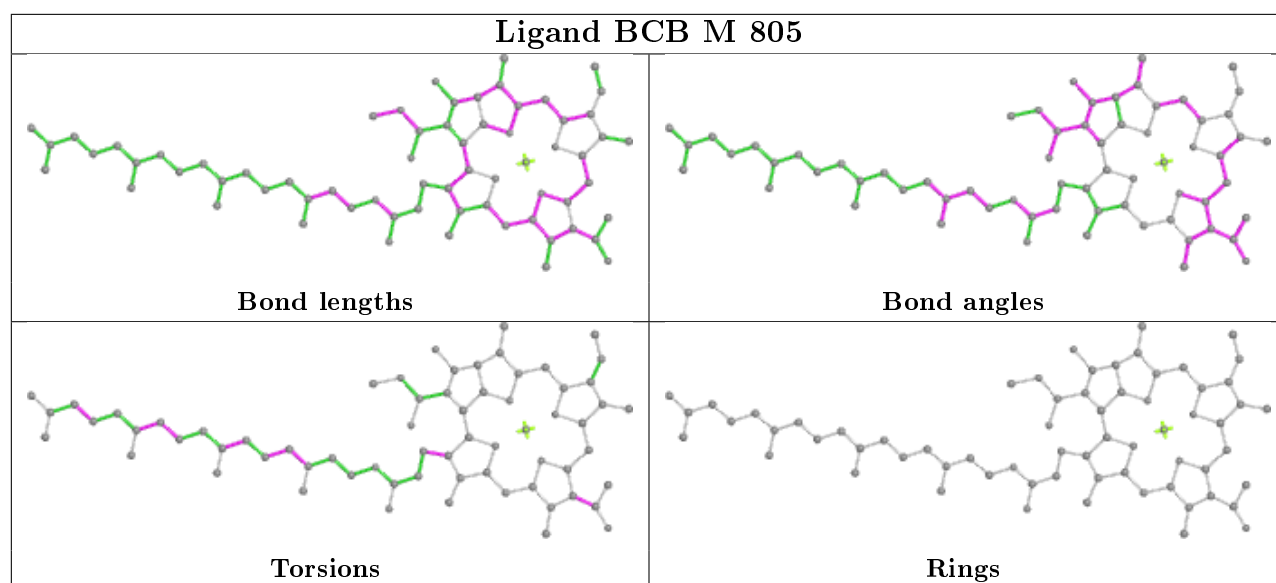
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

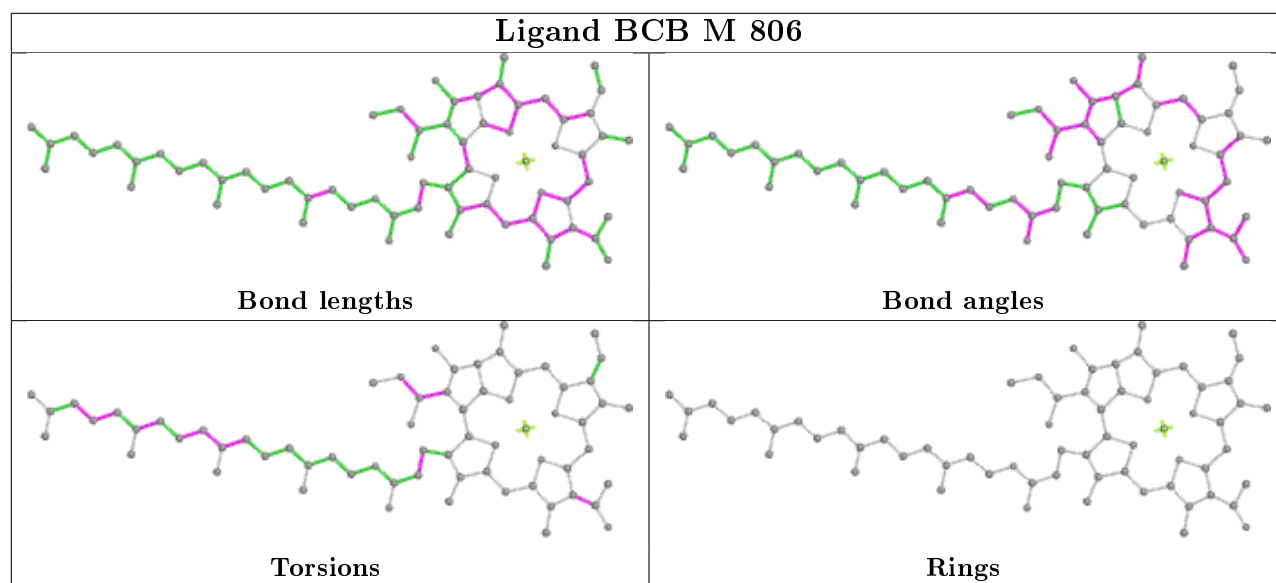
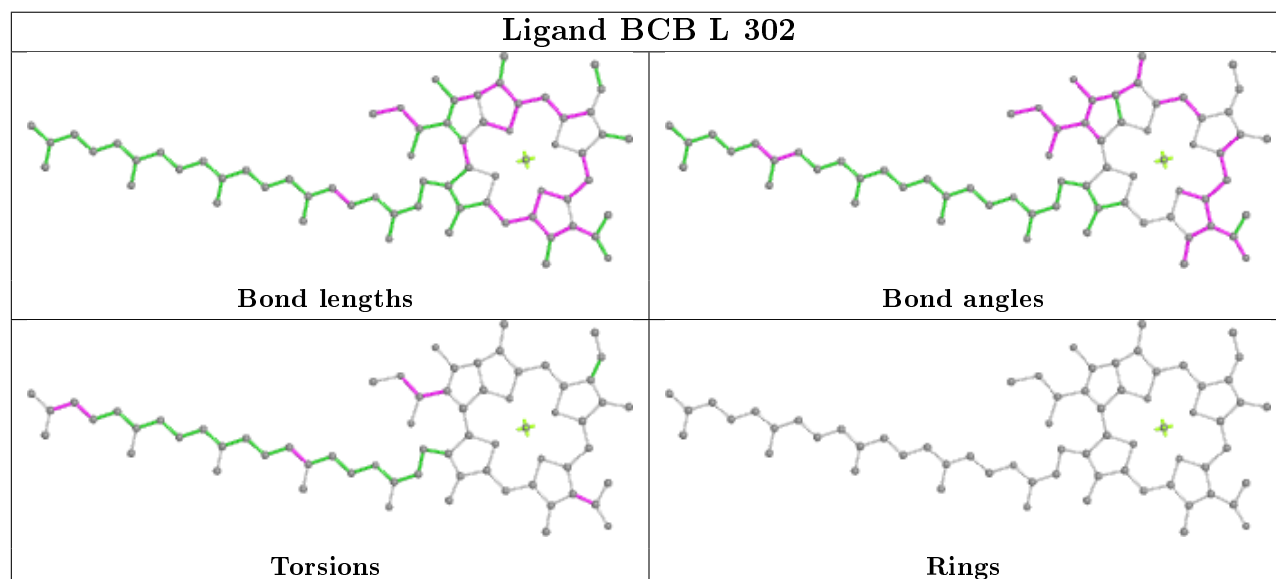
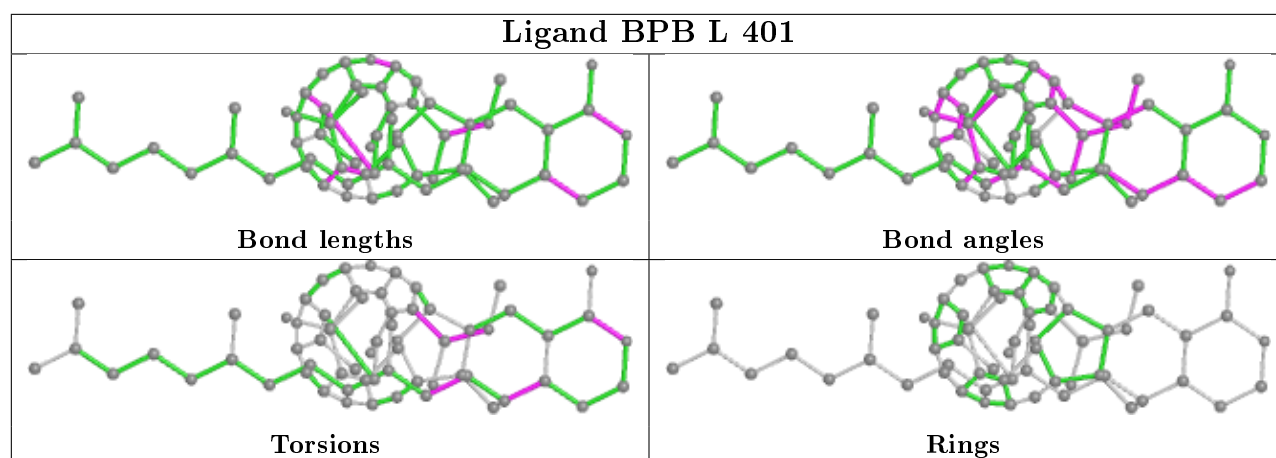


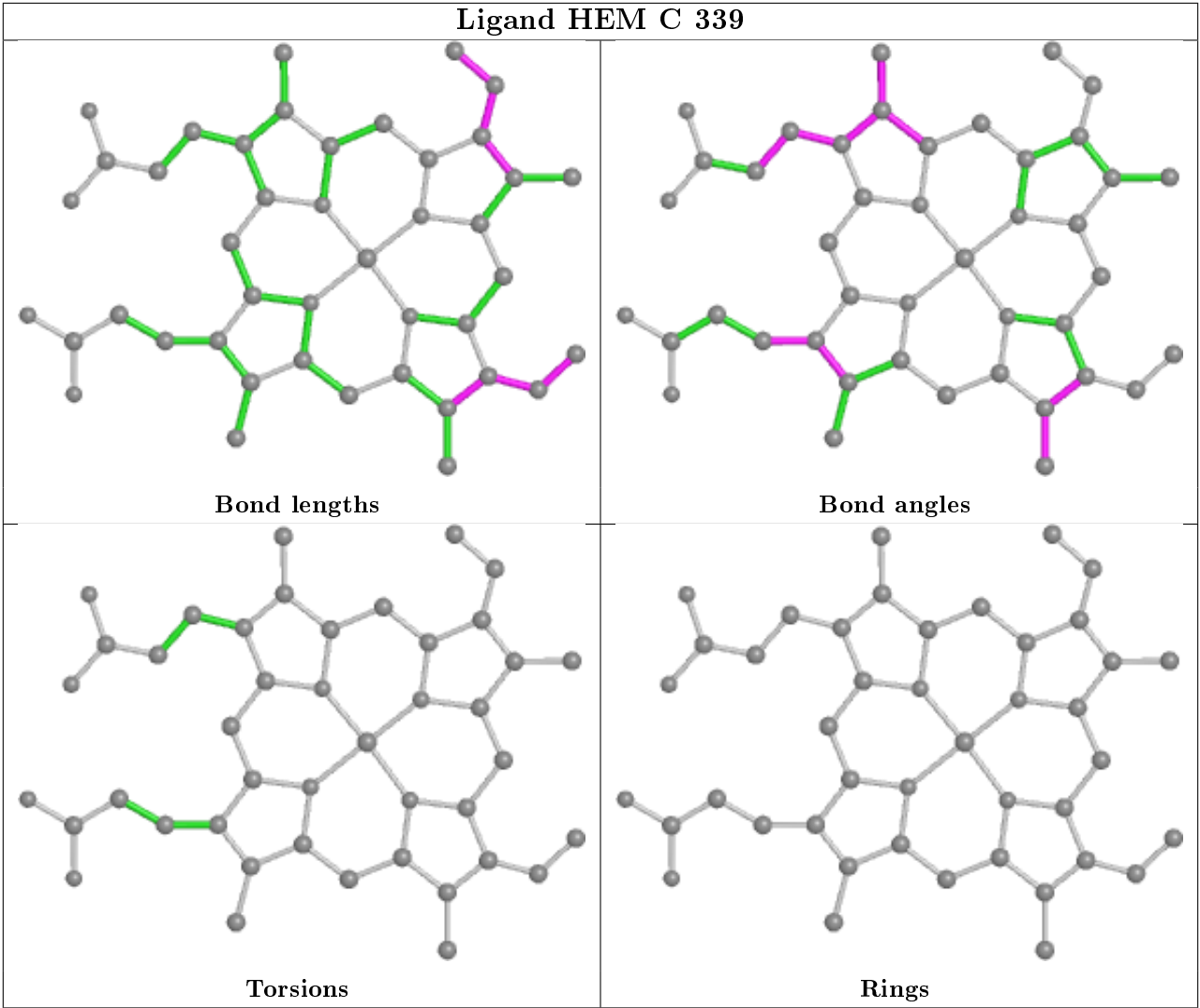












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
4	H	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	54:PRO	C	55:GLU	N	1.68

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Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	H	52:LEU	C	53:ALA	N	1.19

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å²)	Q<0.9	
1	C	332/336 (98%)	-0.77	0	100	100	4, 20, 44, 66	18 (5%)
2	L	273/273 (100%)	-0.91	0	100	100	4, 16, 41, 61	6 (2%)
3	M	323/323 (100%)	-0.78	0	100	100	4, 18, 48, 60	10 (3%)
4	H	249/258 (96%)	-0.65	1 (0%)	92	93	6, 27, 58, 76	19 (7%)
All	All	1177/1190 (98%)	-0.78	1 (0%)	95	96	4, 20, 48, 76	53 (4%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	H	54	PRO	3.3

6.2 Non-standard residues in protein, DNA, RNA chains [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
4	FME	H	1	10/11	0.96	0.10	23,30,40,42	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands ⓘ

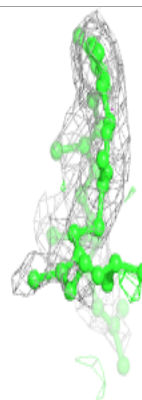
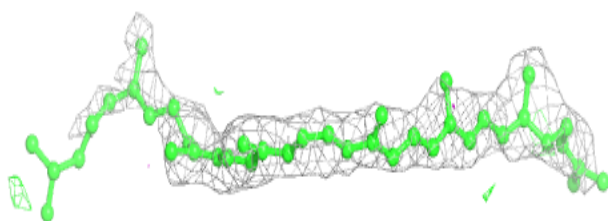
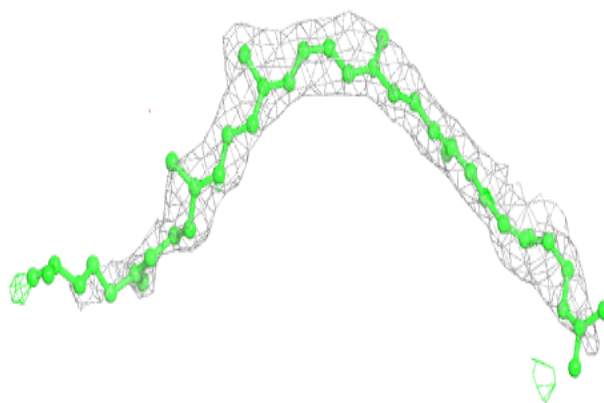
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
9	LDA	L	702	16/16	0.73	0.38	17,45,71,72	0
9	LDA	M	706	16/16	0.77	0.36	68,71,72,73	4
9	LDA	M	705	16/16	0.80	0.23	53,58,69,69	5
9	LDA	M	704	16/16	0.80	0.34	55,59,79,81	0
13	NS5	M	600	40/40	0.84	0.23	23,40,76,76	4
9	LDA	H	703	16/16	0.93	0.22	27,38,57,58	0
12	MQ7	M	501	48/48	0.94	0.13	3,15,34,38	0
11	SO4	M	803	5/5	0.95	0.18	73,74,76,77	0
6	BCB	M	805	66/66	0.95	0.12	2,15,44,46	0
7	BPB	L	401	65/65	0.95	0.13	2,21,70,70	7
9	LDA	M	701	16/16	0.96	0.11	11,21,24,27	0
5	HEM	C	337	43/43	0.97	0.10	8,21,30,39	0
5	HEM	C	340	43/43	0.97	0.11	5,13,30,44	0
5	HEM	C	338	43/43	0.97	0.14	11,22,28,33	0
7	BPB	L	402	65/65	0.97	0.10	2,6,13,17	0
6	BCB	L	302	66/66	0.97	0.11	2,8,15,19	0
6	BCB	M	806	66/66	0.97	0.12	2,10,25,29	0
6	BCB	L	304	66/66	0.98	0.12	2,6,21,27	0
8	CET	L	502	17/17	0.98	0.11	10,18,21,25	0
11	SO4	H	801	5/5	0.98	0.09	56,56,57,57	0
11	SO4	M	804	5/5	0.98	0.12	39,39,44,44	0
5	HEM	C	339	43/43	0.98	0.10	2,10,15,23	0
10	FE2	M	500	1/1	0.99	0.04	16,16,16,16	0
11	SO4	M	802	5/5	0.99	0.09	21,24,27,28	0

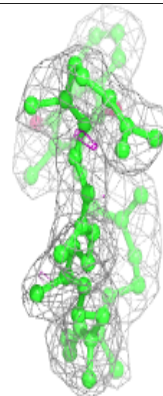
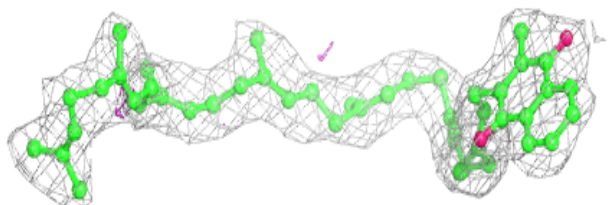
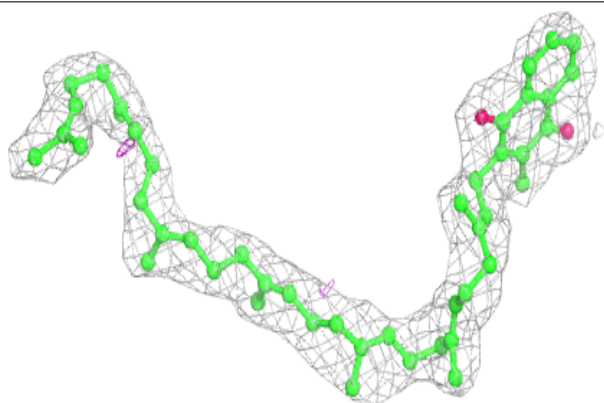
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around NS5 M 600:

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

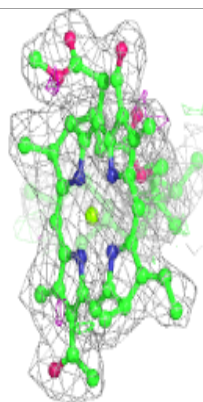
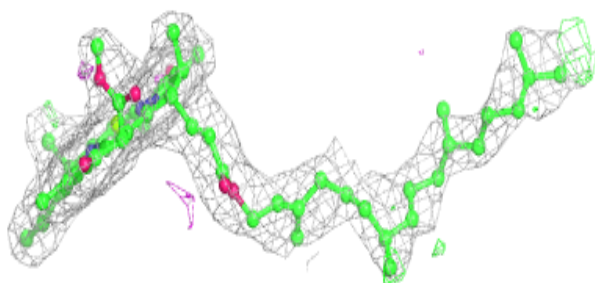
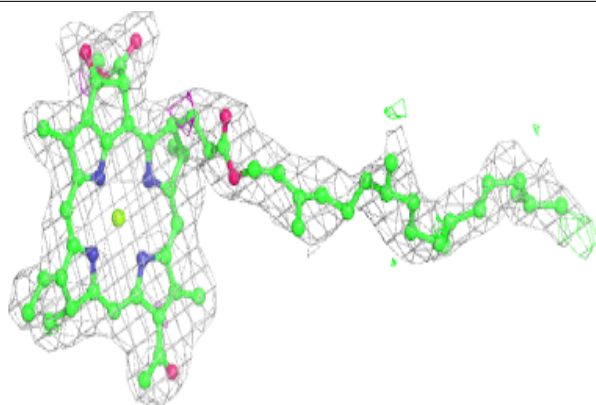
**Electron density around MQ7 M 501:**

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

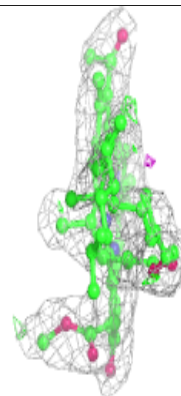
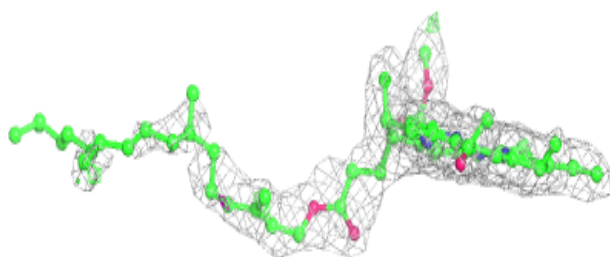
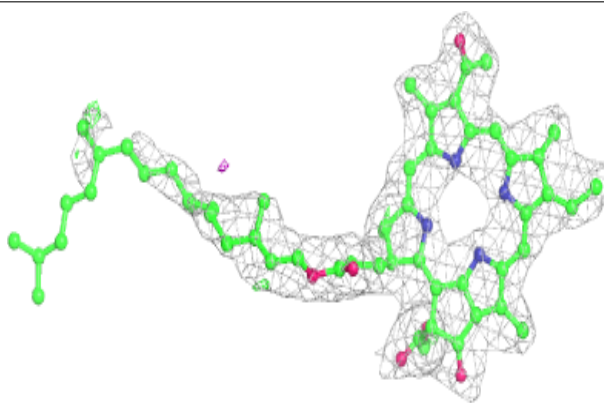


Electron density around BCB M 805:

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and green (positive)

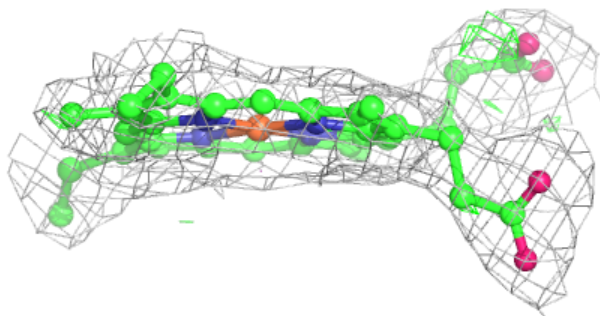
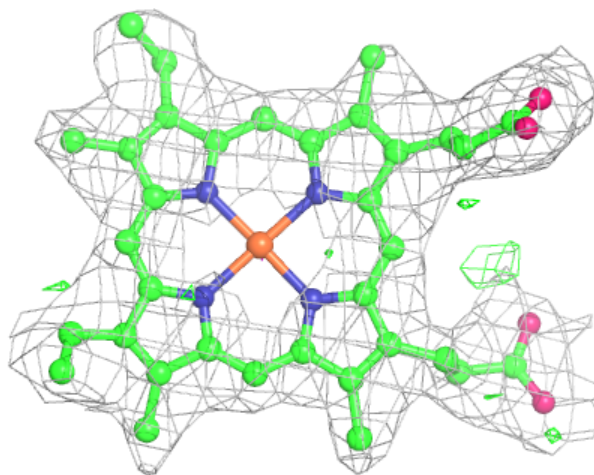
**Electron density around BPB L 401:**

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



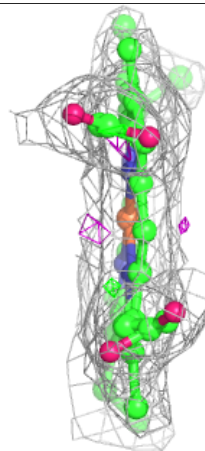
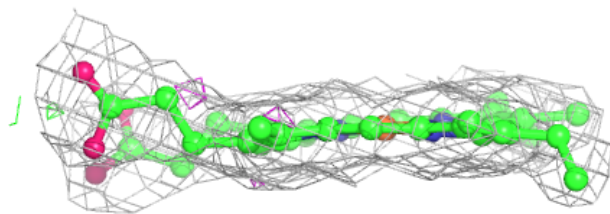
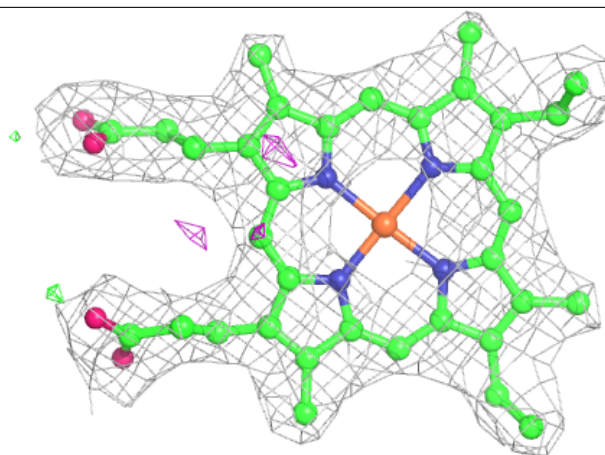
Electron density around HEM C 337:

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and green (positive)



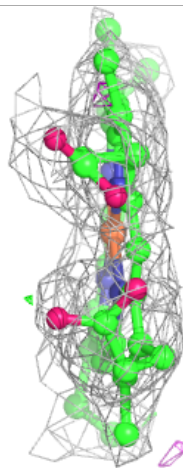
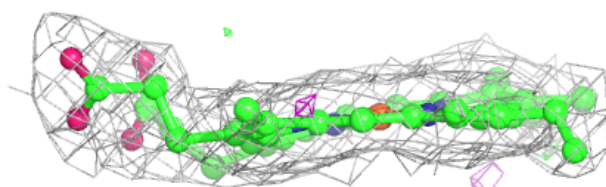
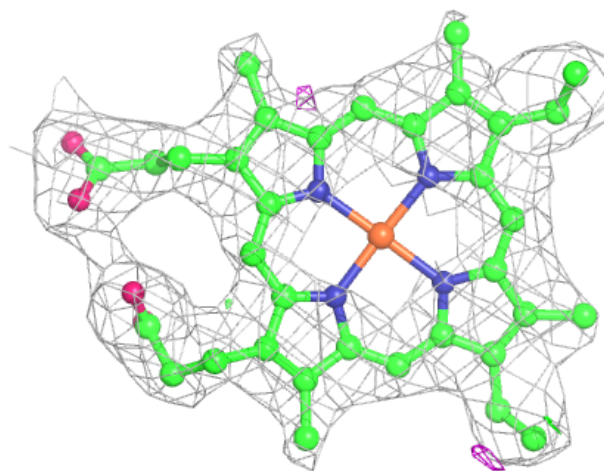
Electron density around HEM C 340:

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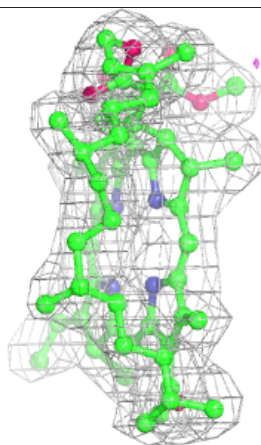
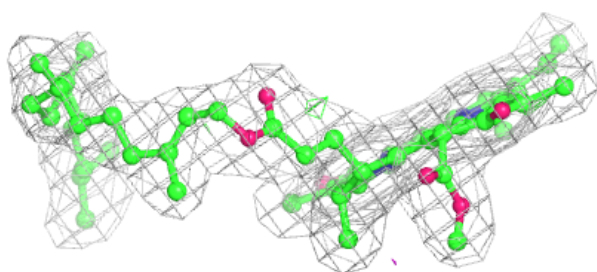
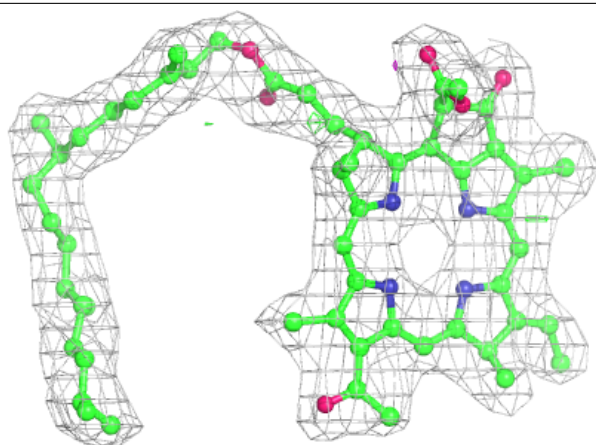
Electron density around HEM C 338:

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and green (positive)

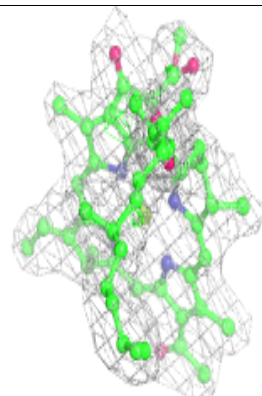
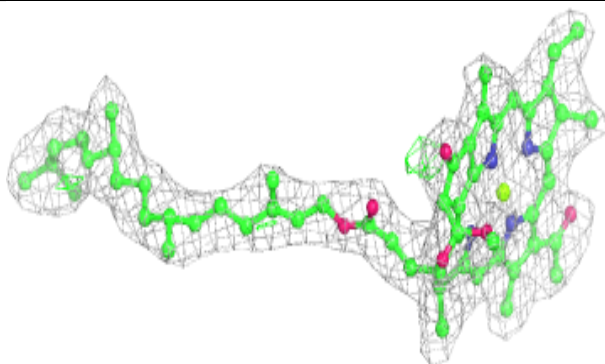
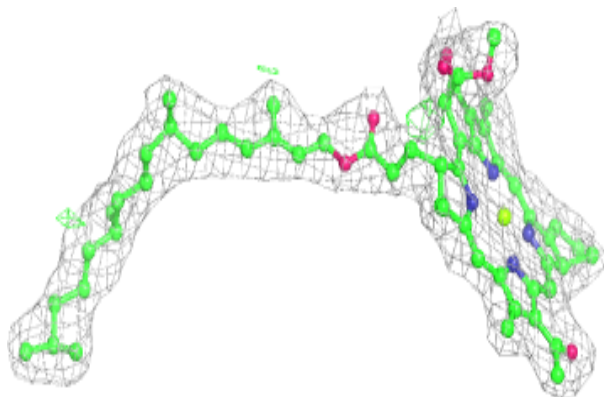


Electron density around BPB L 402:

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and green (positive)

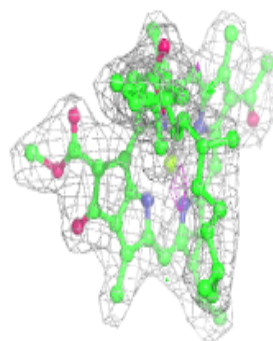
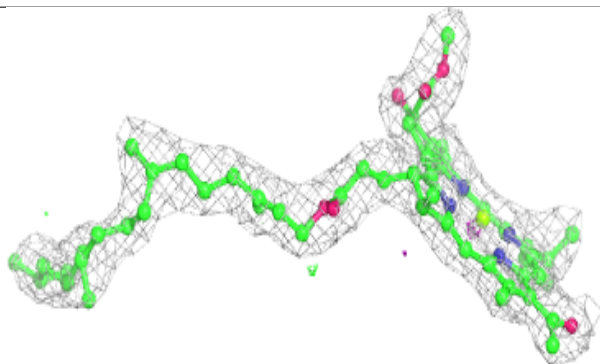
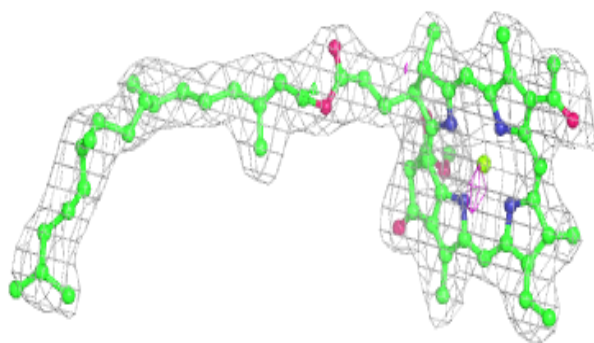
**Electron density around BCB L 302:**

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and green (positive)

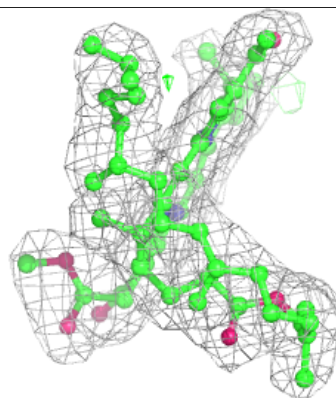
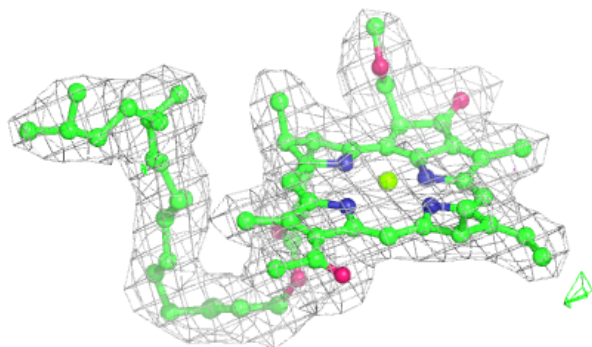
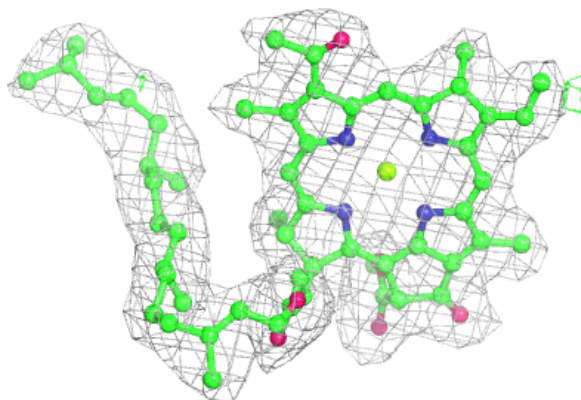


Electron density around BCB M 806:

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and green (positive)

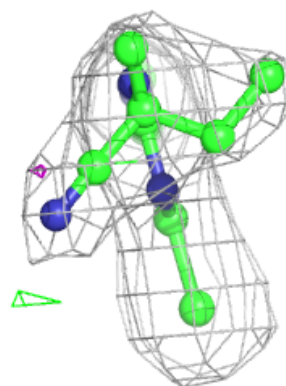
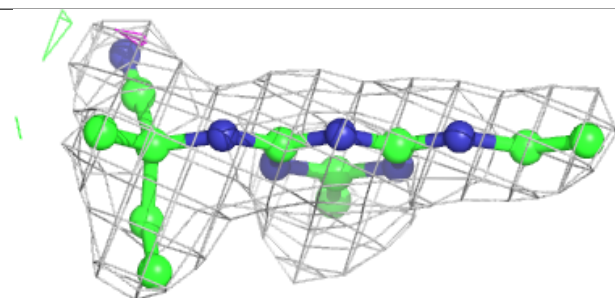
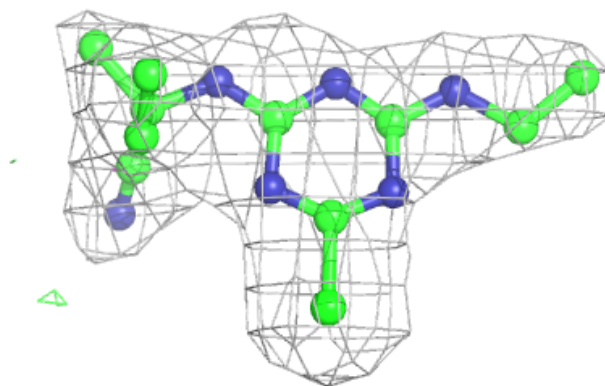
**Electron density around BCB L 304:**

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



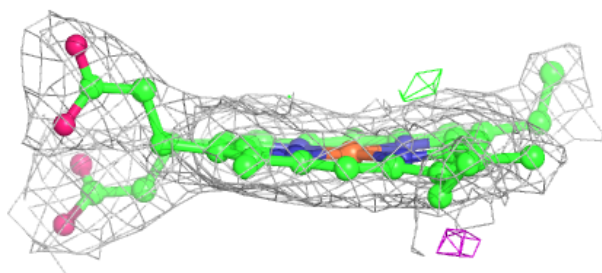
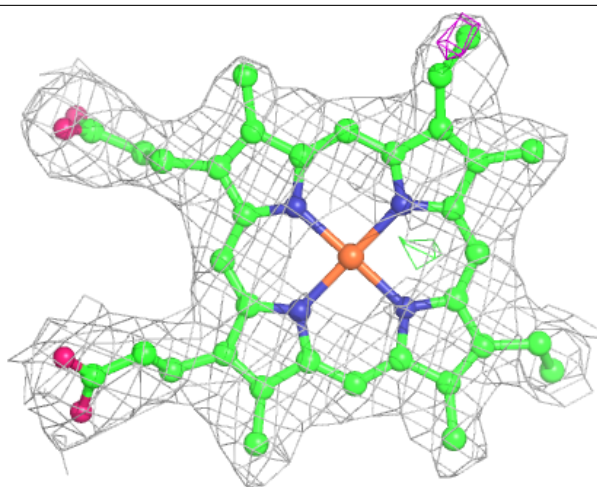
Electron density around CET L 502:

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



Electron density around HEM C 339:

$2mF_o-DF_c$ (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.