



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 6, 2026 – 08:59 AM UTC

PDB ID : 3PRC / pdb_00003prc
Title : PHOTOSYNTHETIC REACTION CENTER FROM RHODOPSEU-
DOMONAS VIRIDIS (QB-DEPLETED)
Authors : Lancaster, C.R.D.; Michel, H.
Deposited on : 1997-07-29
Resolution : 2.40 Å (reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtrriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

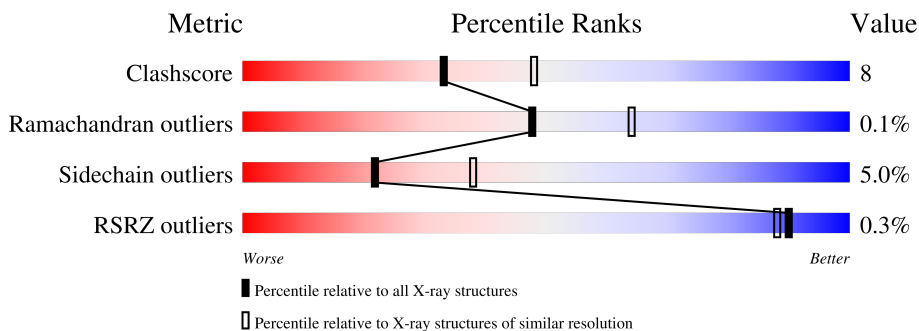
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 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	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	C	336	
2	L	273	
3	M	323	
4	H	258	

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
6	BCB	L	302	X	-	-	-
6	BCB	L	304	X	-	-	-
6	BCB	M	805	X	-	-	-
6	BCB	M	806	X	-	-	-

2 Entry composition

There are 13 unique types of molecules in this entry. The entry contains 10606 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
			Total	C	N	O	S			
1	C	332	2607	1642	467	480	18	38	1	0

- Molecule 2 is a protein called PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	273	2193	1471	358	357	7	14	2	0

- Molecule 3 is a protein called PHOTOSYNTHETIC REACTION CENTER.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	M	323	2566	1711	420	424	11	14	1	0

- Molecule 4 is a protein called PHOTOSYNTHETIC REACTION CENTER.

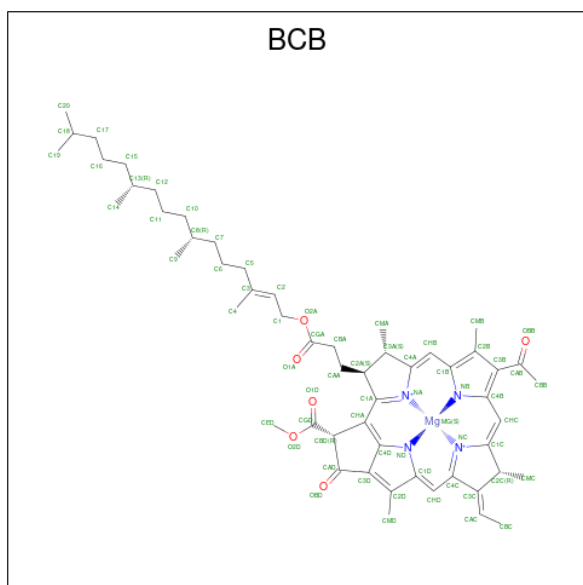
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	H	258	2028	1298	345	382	3	129	1	0

- Molecule 5 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: C₃₄H₃₂FeN₄O₄).



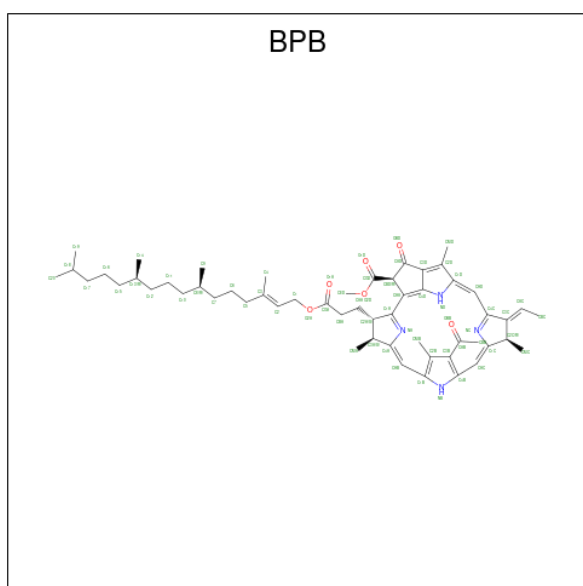
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
5	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
5	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
5	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
5	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

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



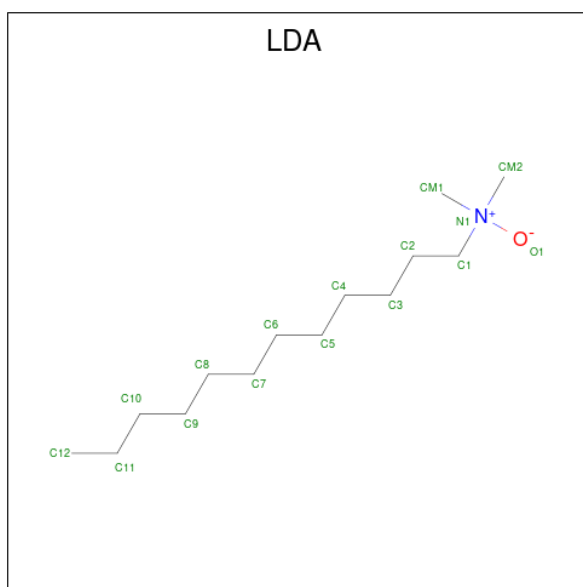
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
6	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

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



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

- Molecule 8 is LAURYL DIMETHYLAMINE-N-OXIDE (CCD ID: LDA) (formula: $C_{14}H_{31}NO$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
8	L	1	16	14	1	1	0	0
8	L	1	16	14	1	1	0	0
8	M	1	16	14	1	1	3	0
8	M	1	16	14	1	1	5	0
8	M	1	16	14	1	1	4	0
8	H	1	16	14	1	1	0	0
8	H	1	16	14	1	1	1	0

- Molecule 9 is FE (II) ION (CCD ID: FE2) (formula: Fe).

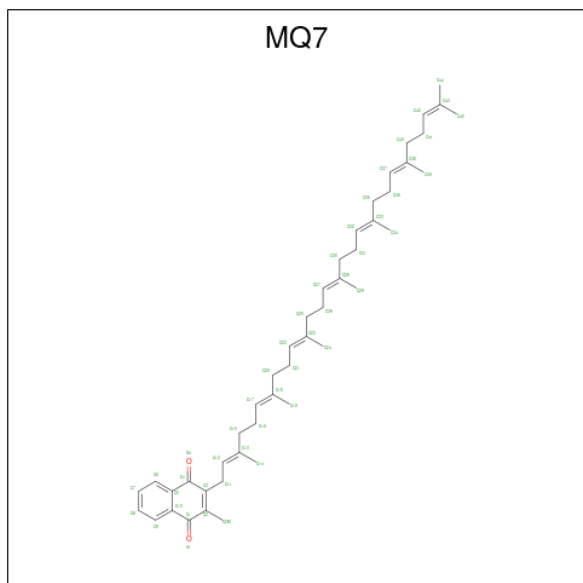
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Fe		
9	M	1	1	1	0	0

- Molecule 10 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



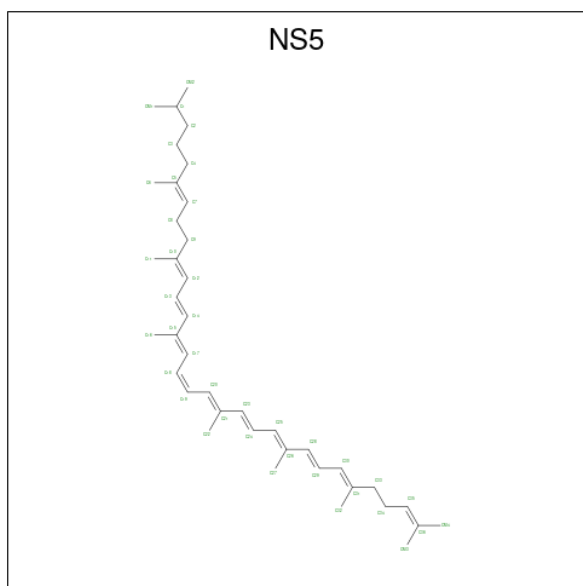
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
10	M	1	Total O S 5 4 1	0	0
10	M	1	Total O S 5 4 1	0	0
10	M	1	Total O S 5 4 1	0	0
10	H	1	Total O S 5 4 1	0	0

- Molecule 11 is MENAQUINONE-7 (CCD ID: MQ7) (formula: $C_{46}H_{64}O_2$).



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

- Molecule 12 is 15-cis-1,2-dihydroneurosporene (CCD ID: NS5) (formula: C₄₀H₆₀).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	M	1	Total	C	9	0
			40	40		

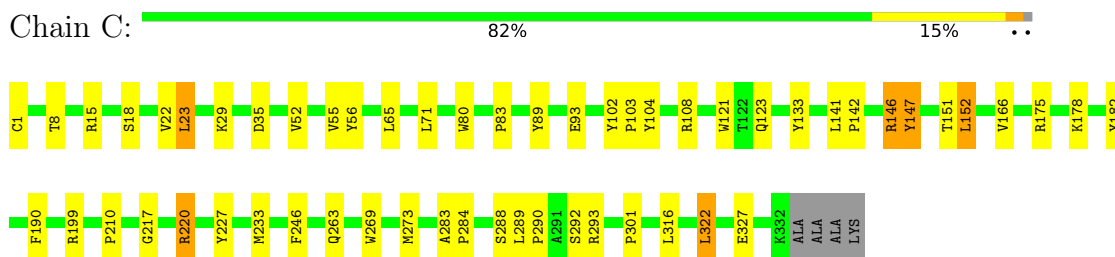
- Molecule 13 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	C	163	Total	O	0	0
			163	163		
13	L	65	Total	O	0	0
			65	65		
13	M	97	Total	O	0	0
			97	97		
13	H	100	Total	O	0	0
			100	100		

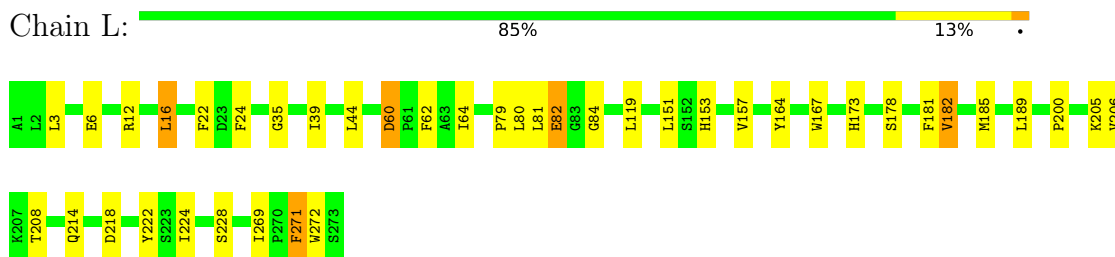
3 Residue-property plots [i](#)

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.

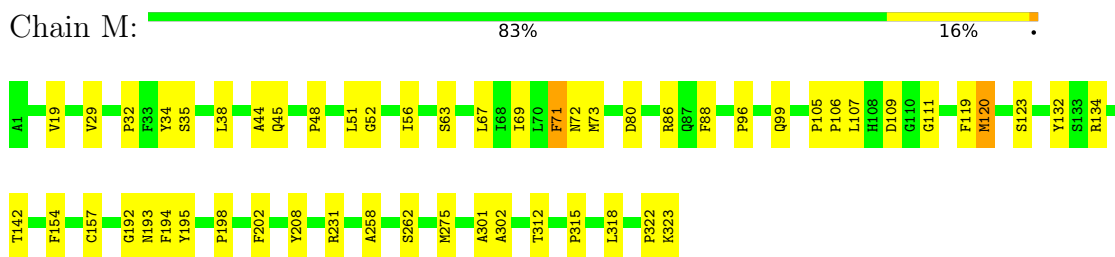
- Molecule 1: PHOTOSYNTHETIC REACTION CENTER



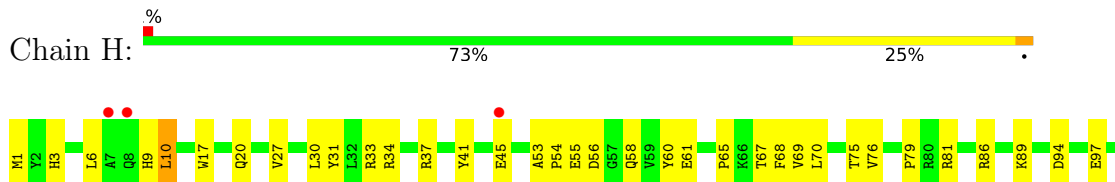
- Molecule 2: PHOTOSYNTHETIC REACTION CENTER

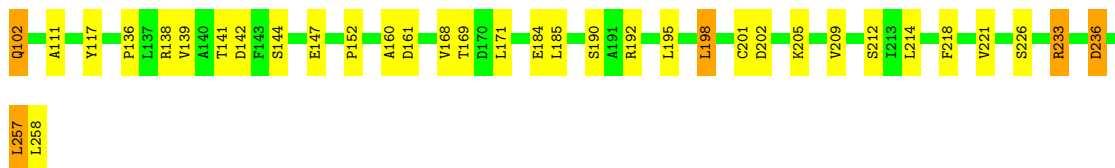


- Molecule 3: PHOTOSYNTHETIC REACTION CENTER



- Molecule 4: PHOTOSYNTHETIC REACTION CENTER





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	223.50Å 223.50Å 113.60Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.40 10.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	79.5 (10.00-2.40) 78.2 (10.00-2.40)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.69 (at 2.39Å)	Xtrriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.178 , 0.215 0.170 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	30.3	Xtrriage
Anisotropy	0.030	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 87.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	10606	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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: FME, NS5, FE2, HEM, LDA, BCB, MQ7, SO4, BPB

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	0.64	0/2674	0.84	1/3645 (0.0%)
2	L	0.64	0/2281	0.85	4/3112 (0.1%)
3	M	0.63	0/2671	0.87	5/3653 (0.1%)
4	H	0.67	3/2055 (0.1%)	0.88	4/2807 (0.1%)
All	All	0.64	3/9681 (0.0%)	0.86	14/13217 (0.1%)

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

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

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	54	PRO	N-CA	-6.81	1.39	1.47
4	H	53	ALA	CA-C	-6.46	1.44	1.52
4	H	53	ALA	C-N	-5.92	1.27	1.33

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	53	ALA	N-CA-C	9.72	131.30	109.81
4	H	53	ALA	CB-CA-C	-9.04	92.36	110.17
4	H	45	GLU	CB-CA-C	-7.85	94.70	110.17

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	M	80	ASP	CA-C-N	7.15	126.64	118.85
3	M	80	ASP	C-N-CA	7.15	126.64	118.85

There are no chirality outliers.

5 of 7 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	147	TYR	Sidechain
1	C	182	TYR	Sidechain
1	C	190	PHE	Sidechain
1	C	227	TYR	Sidechain
1	C	89	TYR	Sidechain

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	C	2607	0	2575	33	0
2	L	2193	0	2122	31	0
3	M	2566	0	2460	36	0
4	H	2028	0	2029	39	0
5	C	172	0	120	1	0
6	L	132	0	144	5	0
6	M	132	0	144	12	0
7	L	65	0	74	4	0
7	M	65	0	74	11	0
8	H	32	0	62	0	0
8	L	32	0	62	10	0
8	M	48	0	93	2	0
9	M	1	0	0	0	0
10	H	5	0	0	0	0
10	M	15	0	0	0	0
11	M	48	0	64	0	0
12	M	40	0	60	0	0
13	C	163	0	0	5	0
13	H	100	0	0	2	0
13	L	65	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	M	97	0	0	0	0
All	All	10606	0	10083	153	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 8.

The worst 5 of 153 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)
7:L:402:BPB:HBBB	7:L:402:BPB:HHC	1.53	0.89
7:M:401:BPB:HBBB	7:M:401:BPB:HHC	1.50	0.89
6:M:805:BCB:HHC	6:M:805:BCB:HBB2	1.58	0.82
1:C:121:TRP:HA	1:C:123[A]:GLN:HE21	1.43	0.81
7:L:402:BPB:HHC	7:L:402: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%)	320 (97%)	11 (3%)	0	100	100
2	L	273/273 (100%)	264 (97%)	9 (3%)	0	100	100
3	M	322/323 (100%)	308 (96%)	13 (4%)	1 (0%)	36	50
4	H	256/258 (99%)	247 (96%)	9 (4%)	0	100	100
All	All	1182/1190 (99%)	1139 (96%)	42 (4%)	1 (0%)	48	64

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	M	322	PRO

5.3.2 Protein sidechains [i](#)

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%)	266 (95%)	15 (5%)	20	36
2	L	220/218 (101%)	214 (97%)	6 (3%)	39	62
3	M	250/249 (100%)	246 (98%)	4 (2%)	55	76
4	H	212/212 (100%)	189 (89%)	23 (11%)	6	9
All	All	963/961 (100%)	915 (95%)	48 (5%)	22	38

5 of 48 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	H	69	VAL
4	H	144	SER
4	H	81	ARG
4	H	97	GLU
4	H	169	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such sidechains are listed below:

Mol	Chain	Res	Type
2	L	239	ASN
3	M	16	HIS
4	H	102	GLN
1	C	297	GLN
1	C	37	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	FME	H	1[B]	4	8,9,10	0.61	0	8,9,11	2.76	2 (25%)
4	FME	H	1[A]	4	8,9,10	0.67	0	8,9,11	2.09	2 (25%)

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[B]	4	-	4/7/9/11	-
4	FME	H	1[A]	4	-	3/7/9/11	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	1[B]	FME	CA-N-CN	-5.33	114.62	122.82
4	H	1[B]	FME	O1-CN-N	-5.12	112.09	125.32
4	H	1[A]	FME	O1-CN-N	-3.90	115.25	125.32
4	H	1[A]	FME	CA-N-CN	-3.81	116.96	122.82

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

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

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides 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$
8	LDA	M	706	-	13,15,15	2.54	2 (15%)	14,17,17	0.50	0
5	HEM	C	339	1	50,50,50	1.57	7 (14%)	67,82,82	1.01	4 (5%)
8	LDA	H	703	-	13,15,15	2.42	2 (15%)	14,17,17	0.59	0
10	SO4	H	801	-	4,4,4	0.65	0	6,6,6	0.37	0
5	HEM	C	338	1	50,50,50	1.48	5 (10%)	67,82,82	1.05	5 (7%)
10	SO4	M	803	-	4,4,4	0.55	0	6,6,6	0.51	0
6	BCB	L	302	2	60,74,74	1.67	9 (15%)	59,115,115	2.52	15 (25%)
8	LDA	L	702	-	13,15,15	2.09	2 (15%)	14,17,17	0.69	0
7	BPB	M	401	-	57,70,70	1.63	11 (19%)	55,101,101	2.37	12 (21%)
6	BCB	M	806	3	60,74,74	1.50	8 (13%)	59,115,115	2.88	17 (28%)
5	HEM	C	337	1	50,50,50	1.50	5 (10%)	67,82,82	1.33	10 (14%)
5	HEM	C	340	1	50,50,50	1.38	6 (12%)	67,82,82	1.00	4 (5%)
6	BCB	L	304	2	60,74,74	1.68	11 (18%)	59,115,115	2.85	15 (25%)
8	LDA	L	707	-	13,15,15	1.84	1 (7%)	14,17,17	0.59	0
7	BPB	L	402	-	57,70,70	1.73	10 (17%)	55,101,101	2.52	15 (27%)
10	SO4	M	804	-	4,4,4	0.88	0	6,6,6	0.69	0
12	NS5	M	600	-	39,39,39	0.68	0	46,46,46	1.12	5 (10%)
6	BCB	M	805	3	60,74,74	1.68	9 (15%)	59,115,115	2.82	18 (30%)
8	LDA	H	701	-	13,15,15	2.42	2 (15%)	14,17,17	0.76	0
10	SO4	M	802	-	4,4,4	0.91	0	6,6,6	0.85	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	LDA	M	705	-	13,15,15	2.44	2 (15%)	14,17,17	0.52	0
8	LDA	M	704	-	13,15,15	2.43	2 (15%)	14,17,17	0.54	0
11	MQ7	M	501	-	49,49,49	1.57	9 (18%)	61,63,63	1.31	8 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	LDA	M	706	-	-	5/13/13/13	-
5	HEM	C	339	1	-	4/14/54/54	-
8	LDA	H	703	-	-	5/13/13/13	-
5	HEM	C	338	1	-	7/14/54/54	-
6	BCB	L	302	2	3/3/21/26	4/37/137/137	-
8	LDA	L	702	-	-	3/13/13/13	-
7	BPB	M	401	-	-	6/37/105/105	0/5/6/6
6	BCB	M	806	3	3/3/21/26	3/37/137/137	-
5	HEM	C	337	1	-	6/14/54/54	-
5	HEM	C	340	1	-	5/14/54/54	-
6	BCB	L	304	2	2/2/21/26	6/37/137/137	-
8	LDA	L	707	-	-	5/13/13/13	-
7	BPB	L	402	-	-	4/37/105/105	0/5/6/6
12	NS5	M	600	-	-	11/43/43/43	-
6	BCB	M	805	3	3/3/21/26	7/37/137/137	-
8	LDA	H	701	-	-	4/13/13/13	-
8	LDA	M	705	-	-	2/13/13/13	-
8	LDA	M	704	-	-	4/13/13/13	-
11	MQ7	M	501	-	-	1/41/61/61	0/2/2/2

The worst 5 of 103 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	H	701	LDA	O1-N1	-8.26	1.21	1.42
8	H	703	LDA	O1-N1	-7.53	1.23	1.42
8	M	704	LDA	O1-N1	-7.35	1.24	1.42
8	M	706	LDA	O1-N1	-7.33	1.24	1.42
8	L	702	LDA	O1-N1	-7.04	1.24	1.42

The worst 5 of 128 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	M	401	BPB	O2D-CGD-CBD	12.03	124.16	110.95
6	L	304	BCB	O2D-CGD-CBD	11.56	123.64	110.95
6	M	805	BCB	O2D-CGD-CBD	11.20	123.26	110.95
7	L	402	BPB	O2D-CGD-CBD	10.58	122.57	110.95
6	M	806	BCB	O2D-CGD-CBD	10.19	122.14	110.95

5 of 11 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	L	302	BCB	NA
6	L	302	BCB	ND
6	L	302	BCB	NC
6	L	304	BCB	NA
6	L	304	BCB	NC

5 of 92 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	M	806	BCB	CAD-CBD-CGD-O1D
6	M	806	BCB	CAD-CBD-CGD-O2D
12	M	600	NS5	C9-C10-C12-C13
12	M	600	NS5	C11-C10-C12-C13
12	M	600	NS5	C10-C12-C13-C14

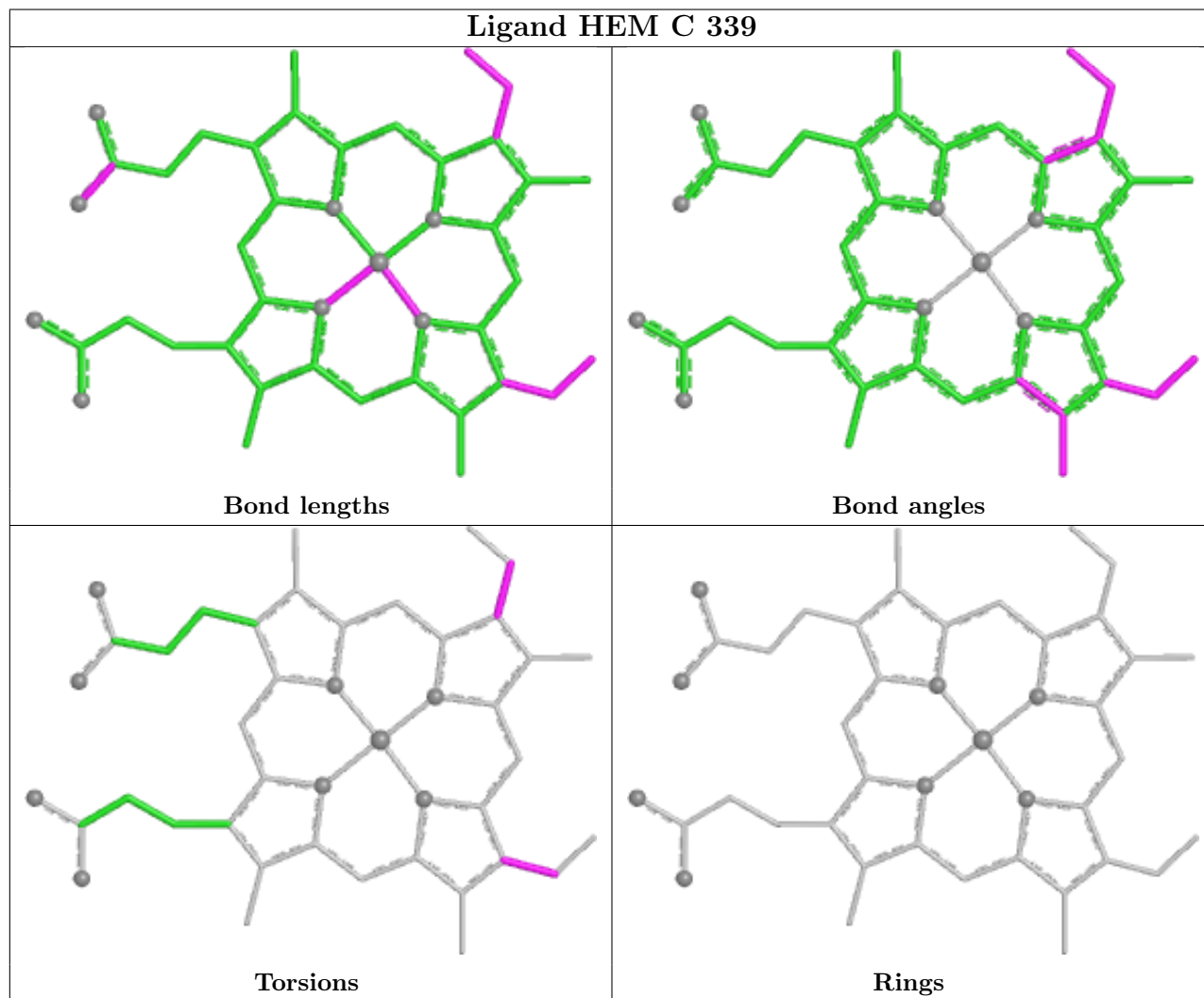
There are no ring outliers.

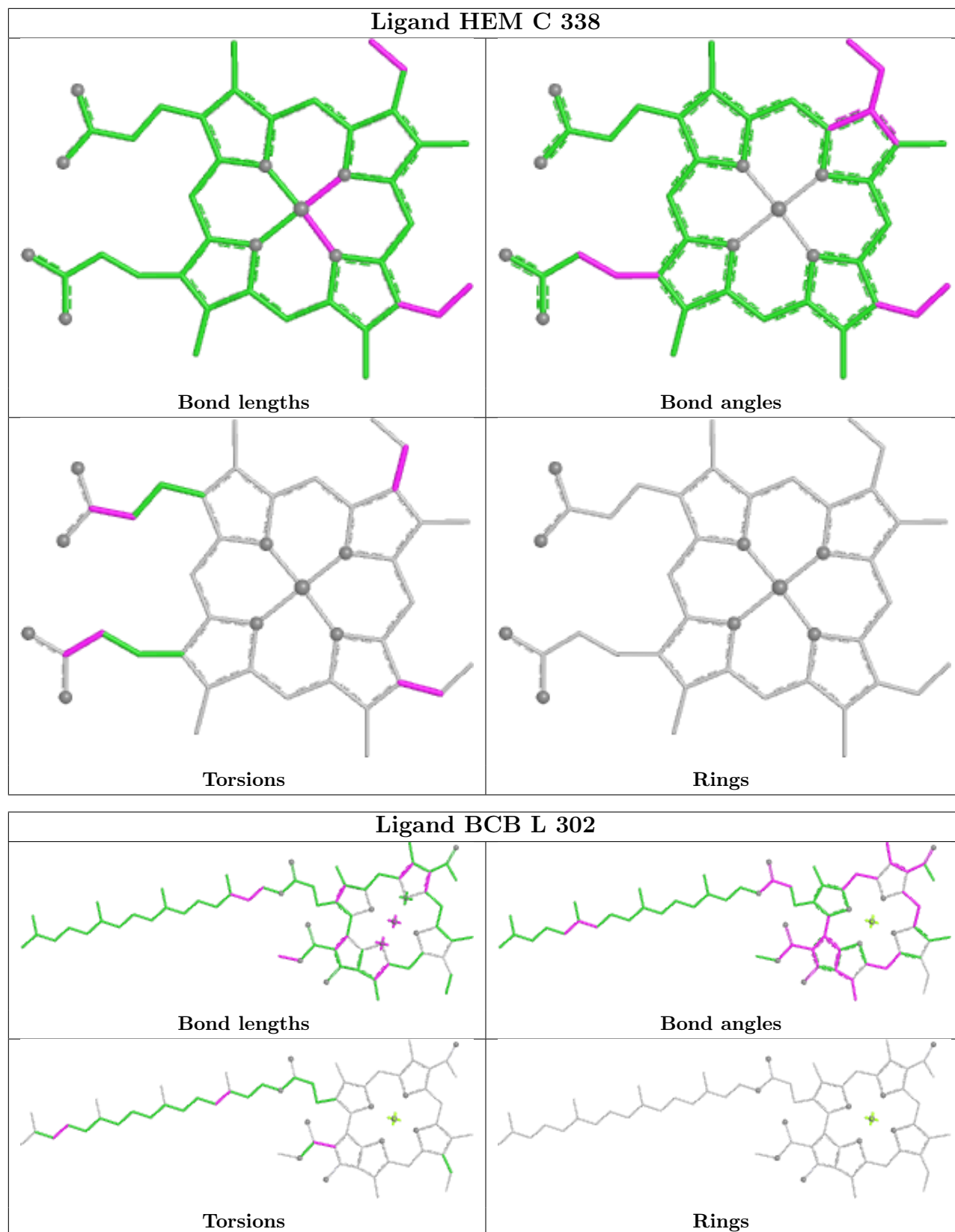
10 monomers are involved in 43 short contacts:

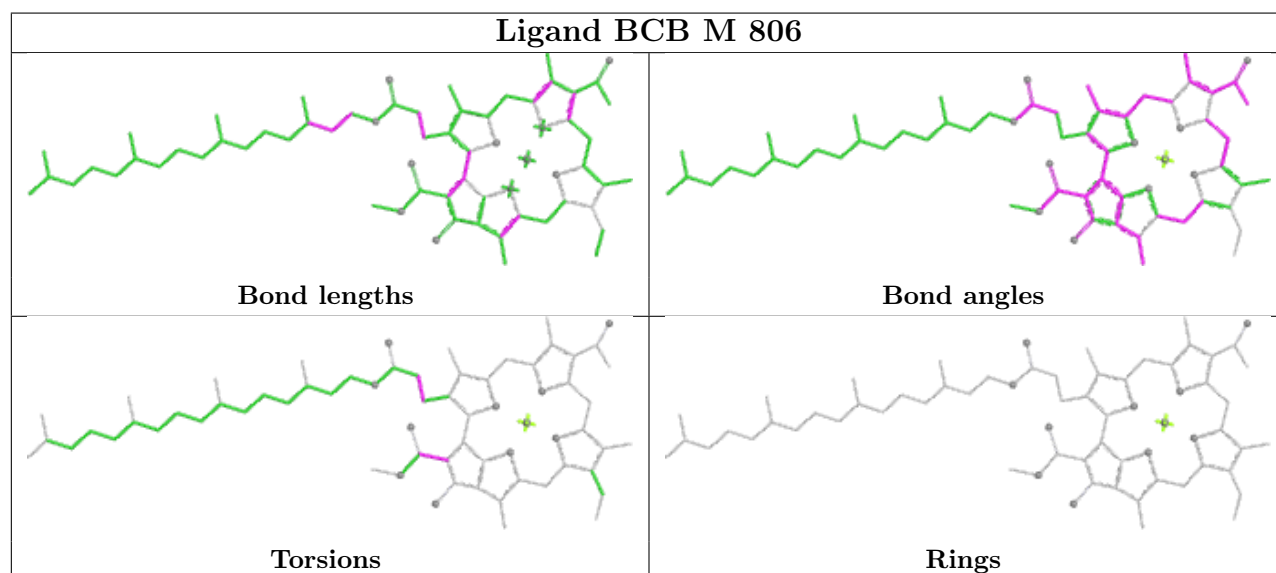
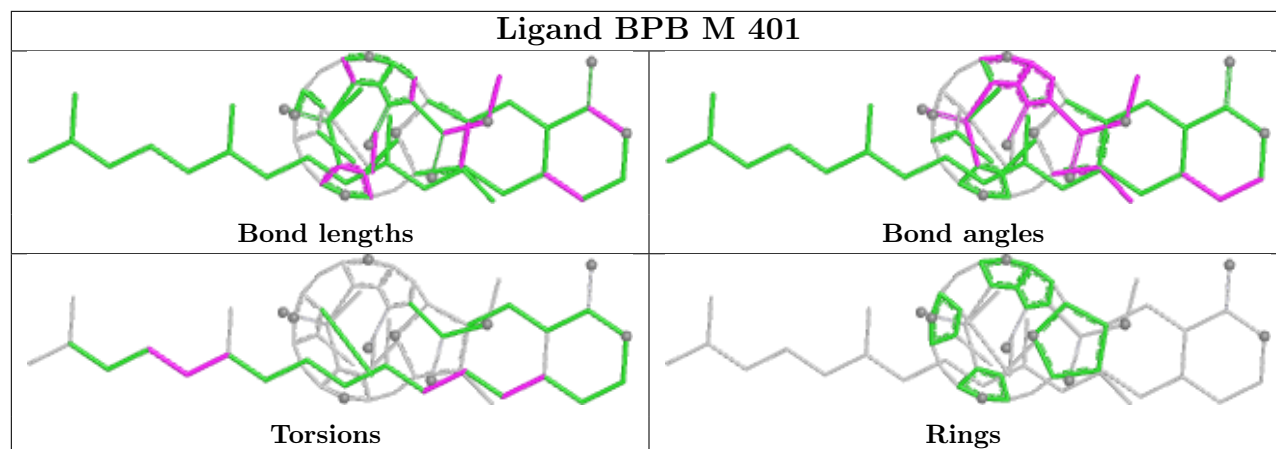
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	M	706	LDA	2	0
5	C	339	HEM	1	0
6	L	302	BCB	3	0
8	L	702	LDA	8	0
7	M	401	BPB	11	0
6	M	806	BCB	6	0
6	L	304	BCB	2	0
8	L	707	LDA	2	0
7	L	402	BPB	4	0
6	M	805	BCB	6	0

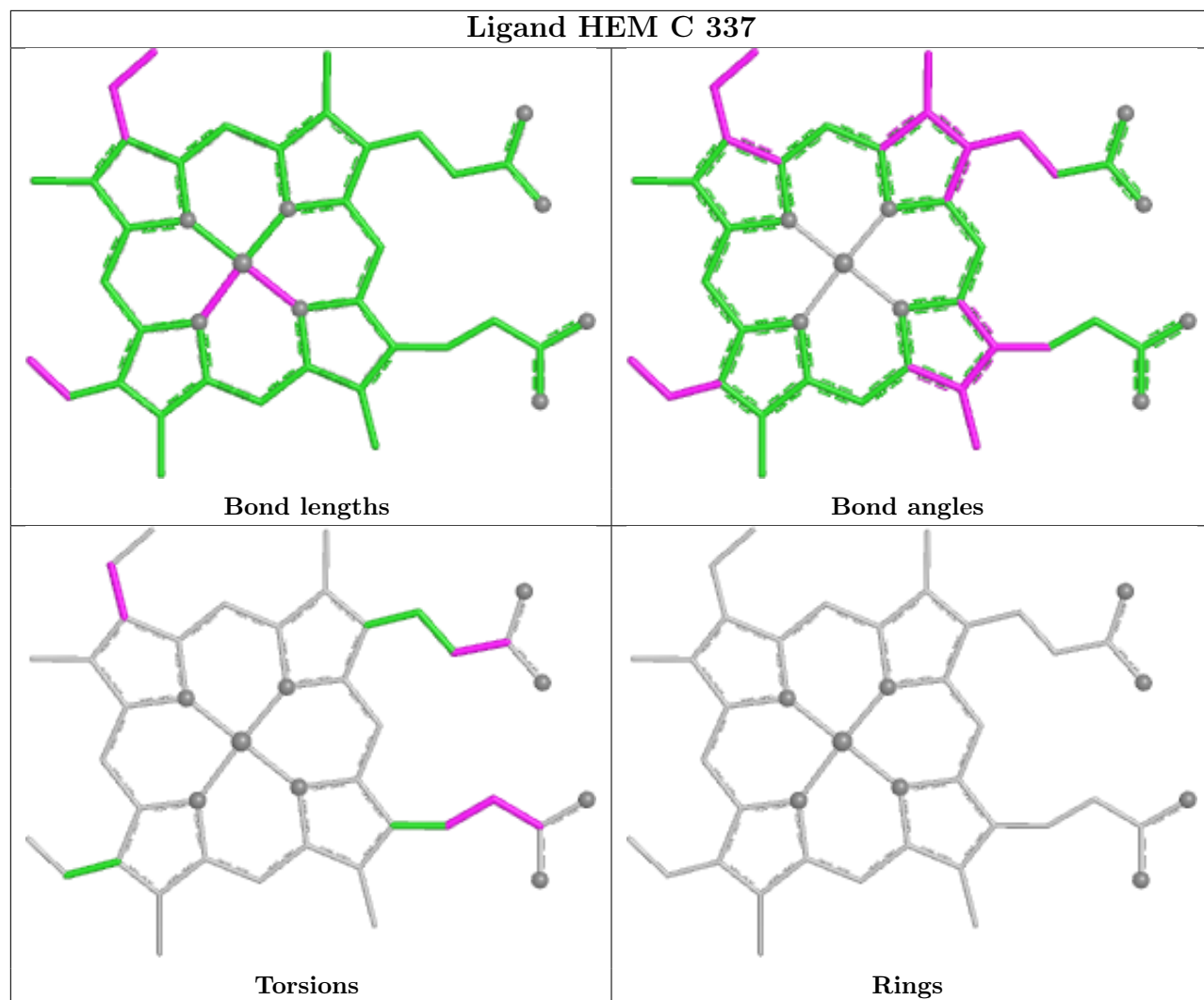
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

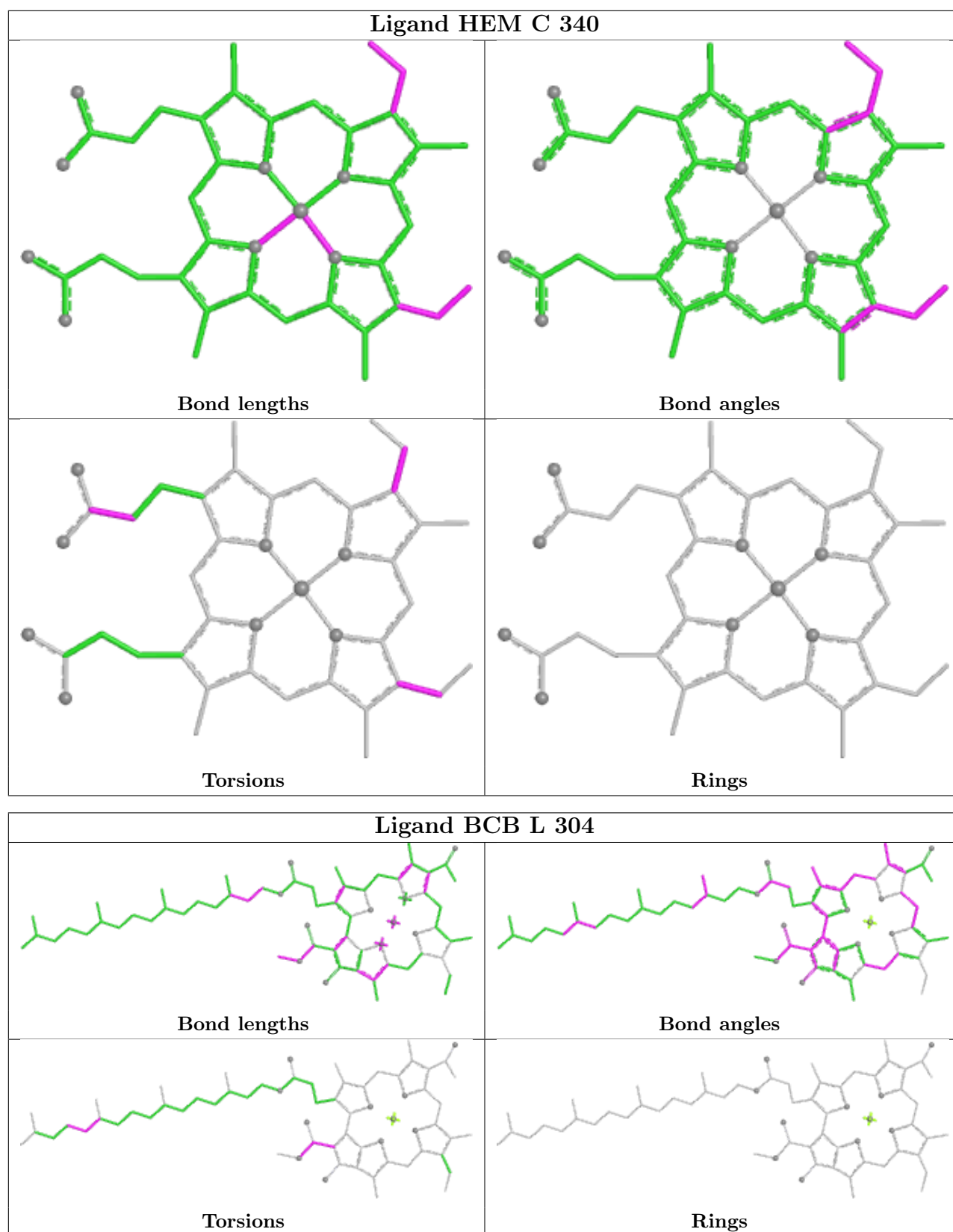
addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. 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.

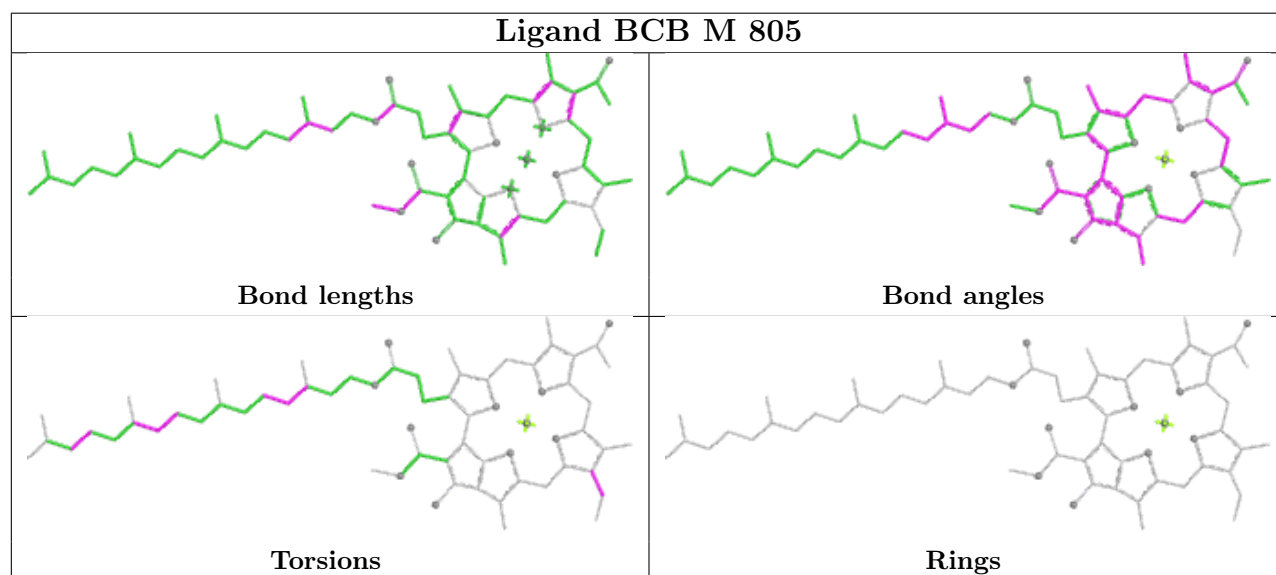
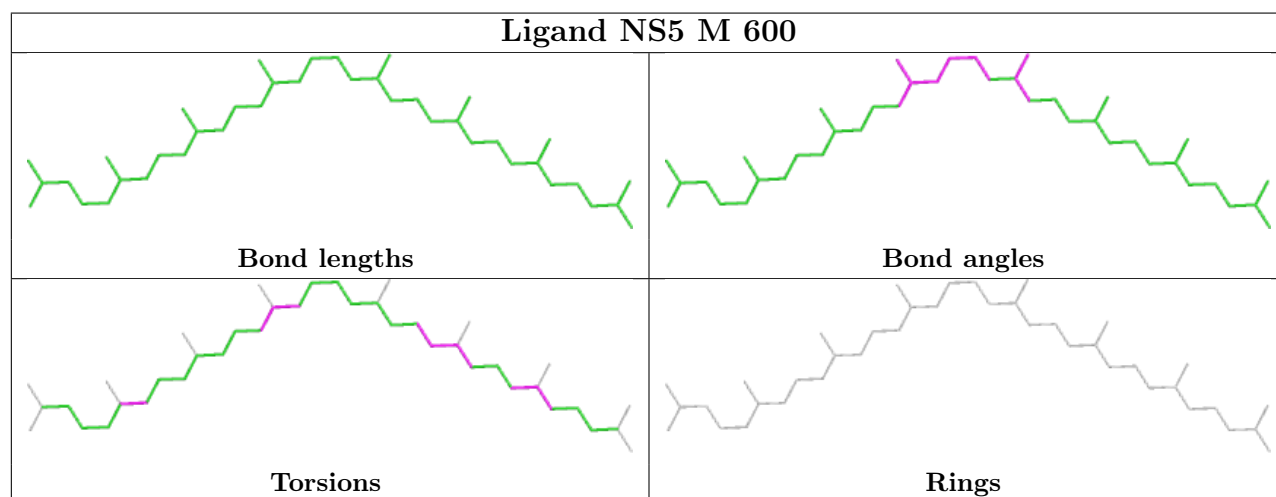
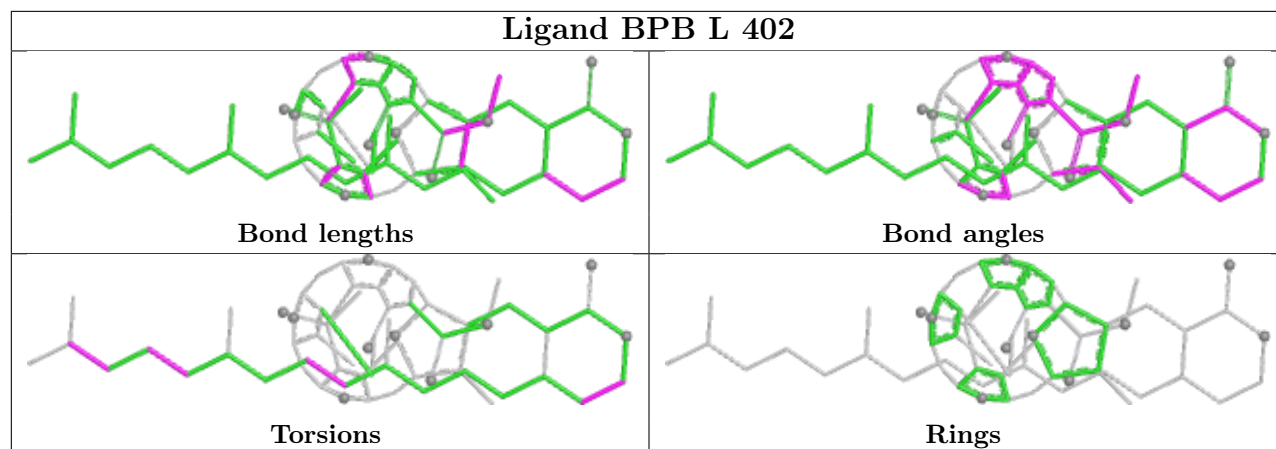


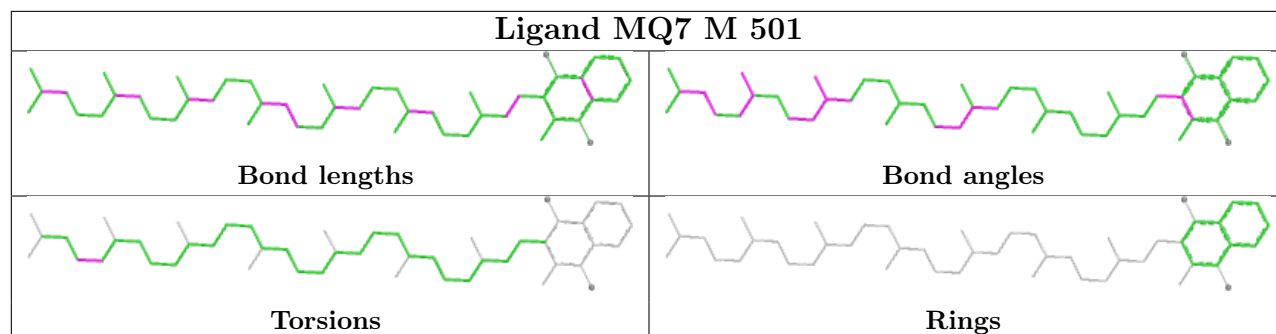












5.7 Other polymers [i](#)

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, 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	C	332/336 (98%)	-0.90	0	100	100	11, 25, 46, 67	16 (4%)
2	L	273/273 (100%)	-0.99	0	100	100	9, 19, 39, 50	9 (3%)
3	M	323/323 (100%)	-0.91	0	100	100	9, 22, 50, 61	9 (2%)
4	H	249/258 (96%)	-0.65	3 (1%)	76	73	12, 28, 46, 61	22 (8%)
All	All	1177/1190 (98%)	-0.87	3 (0%)	90	88	9, 23, 46, 67	56 (4%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
4	H	45	GLU	3.0
4	H	8	GLN	2.8
4	H	7	ALA	2.3

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[A]	10/11	0.94	0.08	22,25,27,27	10
4	FME	H	1[B]	10/11	0.94	0.08	27,30,30,30	10

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

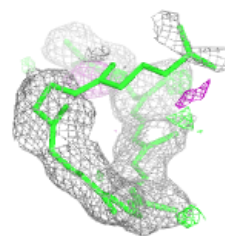
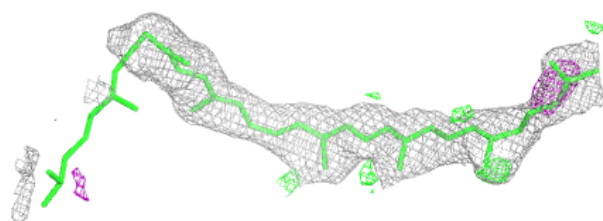
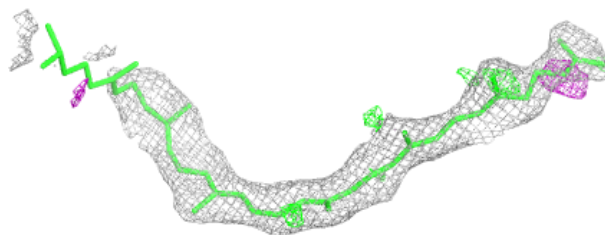
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	LDA	L	707	16/16	0.72	0.13	54,64,65,66	0
8	LDA	M	705	16/16	0.74	0.11	51,55,59,59	5
8	LDA	M	706	16/16	0.77	0.15	69,70,72,72	4
8	LDA	M	704	16/16	0.82	0.11	53,59,61,62	3
12	NS5	M	600	40/40	0.87	0.09	31,49,68,68	9
8	LDA	H	703	16/16	0.91	0.08	30,41,52,52	1
10	SO4	M	803	5/5	0.92	0.10	85,85,86,86	0
6	BCB	M	805	66/66	0.94	0.08	8,18,37,38	20
8	LDA	H	701	16/16	0.95	0.05	28,30,31,32	0
8	LDA	L	702	16/16	0.95	0.09	19,44,66,66	0
10	SO4	H	801	5/5	0.96	0.07	64,65,65,67	0
7	BPB	M	401	65/65	0.97	0.05	10,24,61,62	7
7	BPB	L	402	65/65	0.98	0.04	7,12,18,20	0
6	BCB	L	304	66/66	0.98	0.04	6,13,26,36	0
10	SO4	M	804	5/5	0.98	0.07	40,41,42,43	0
6	BCB	L	302	66/66	0.98	0.03	7,11,18,19	0
11	MQ7	M	501	48/48	0.98	0.04	11,17,43,49	0
6	BCB	M	806	66/66	0.98	0.04	4,12,30,31	0
5	HEM	C	338	43/43	0.99	0.04	16,25,32,38	0
5	HEM	C	339	43/43	0.99	0.04	8,14,18,25	0
5	HEM	C	340	43/43	0.99	0.05	11,19,34,42	0
5	HEM	C	337	43/43	0.99	0.05	21,27,38,43	0
10	SO4	M	802	5/5	0.99	0.06	30,30,32,39	0
9	FE2	M	500	1/1	1.00	0.01	14,14,14,14	0

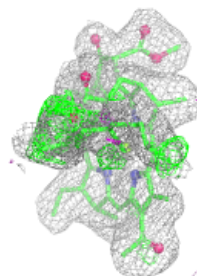
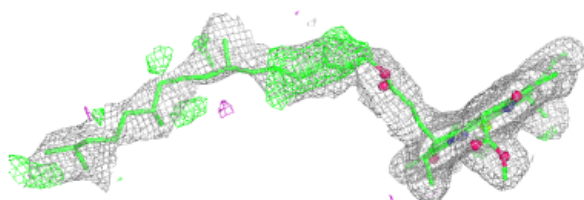
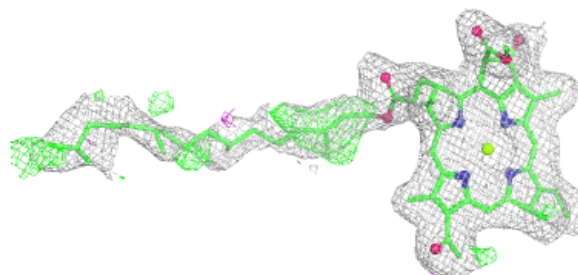
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

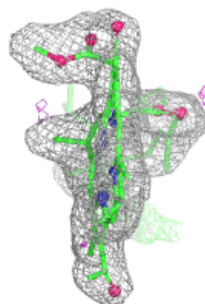
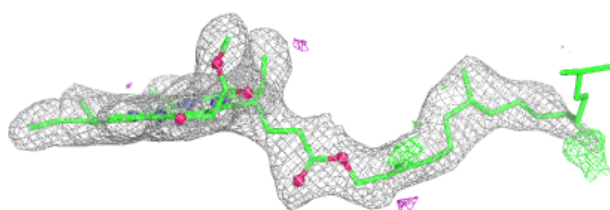
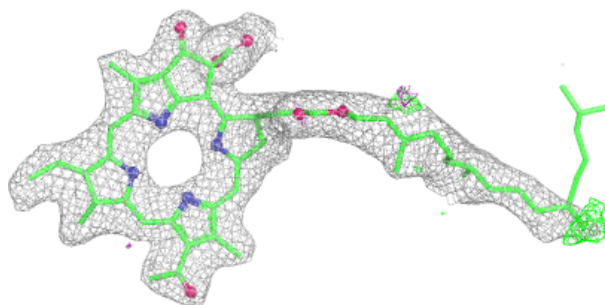
**Electron density around BCB M 805:**

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

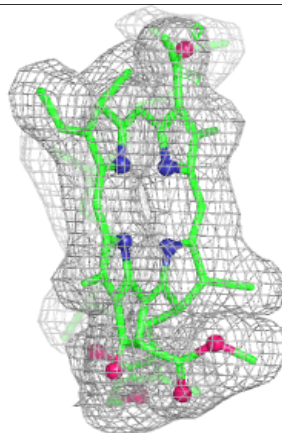
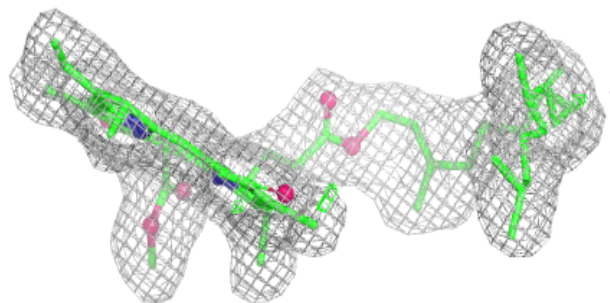
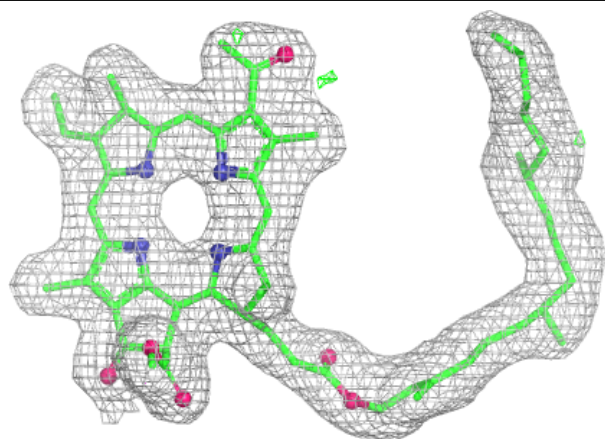


Electron density around BPB M 401:

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

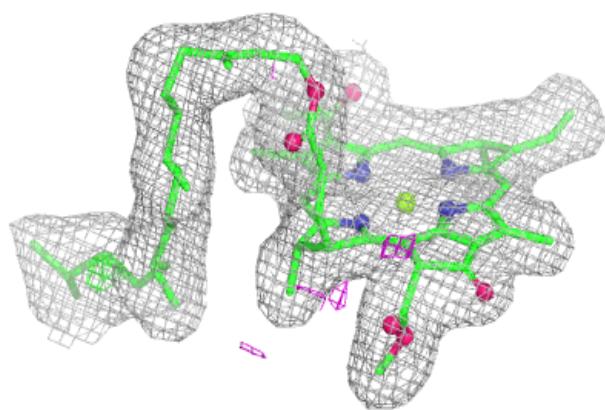
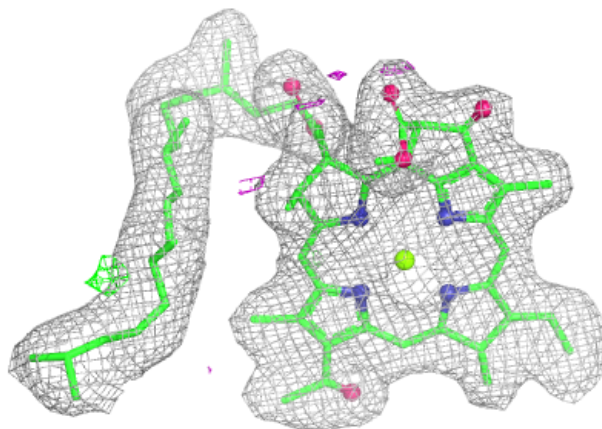
**Electron density around BPB L 402:**

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



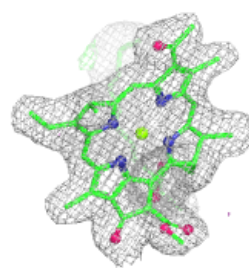
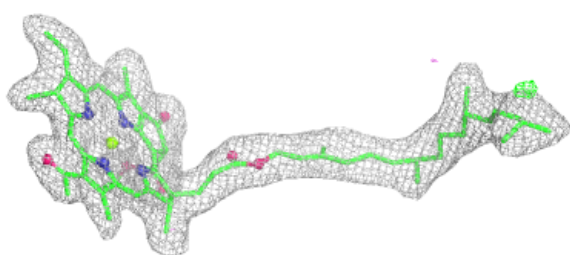
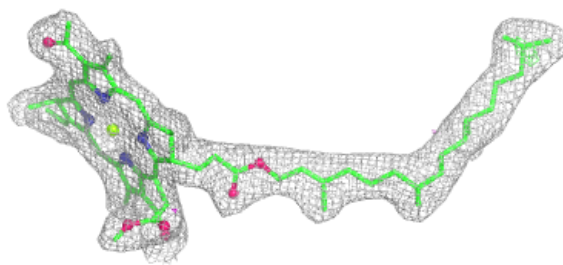
Electron density around BCB L 304:

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and green (positive)

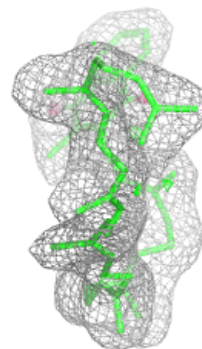
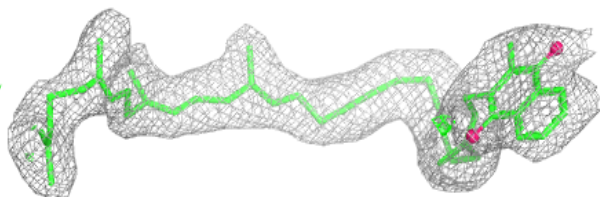
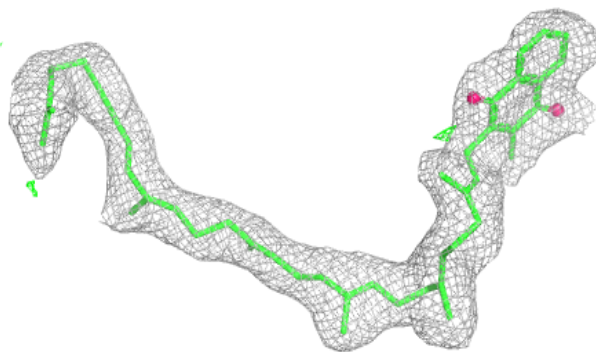


Electron density around BCB L 302:

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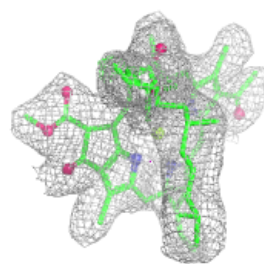
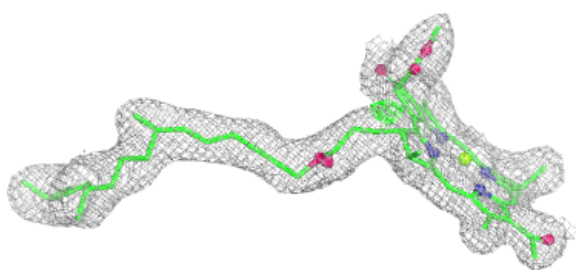
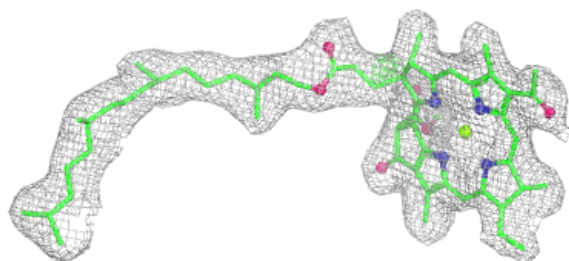
**Electron density around MQ7 M 501:**

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and green (positive)



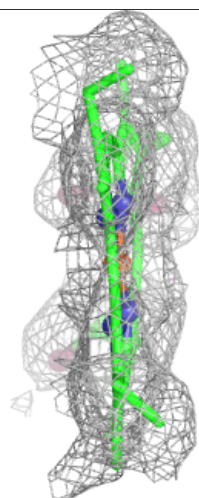
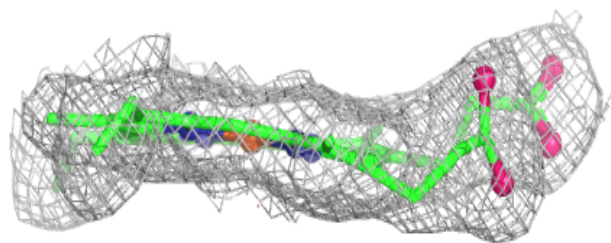
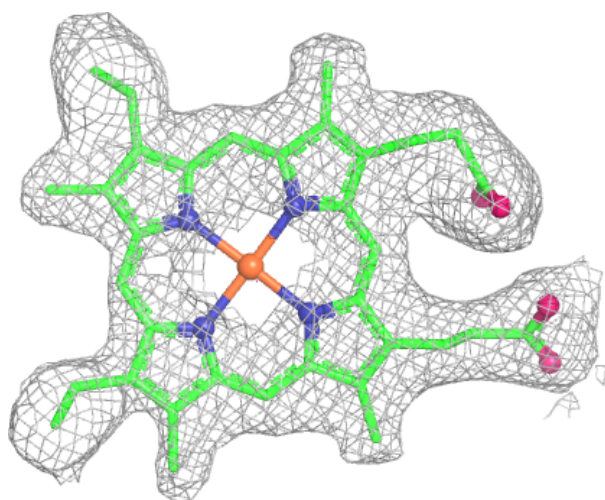
Electron density around BCB M 806:

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and green (positive)



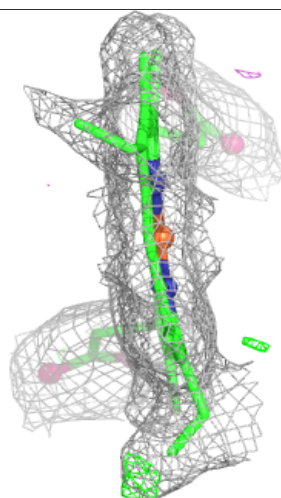
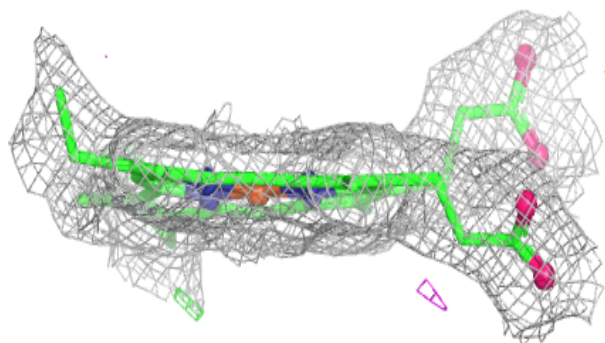
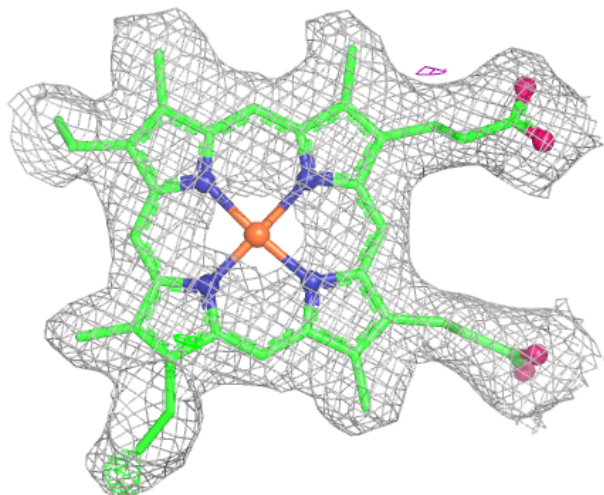
Electron density around HEM C 338:

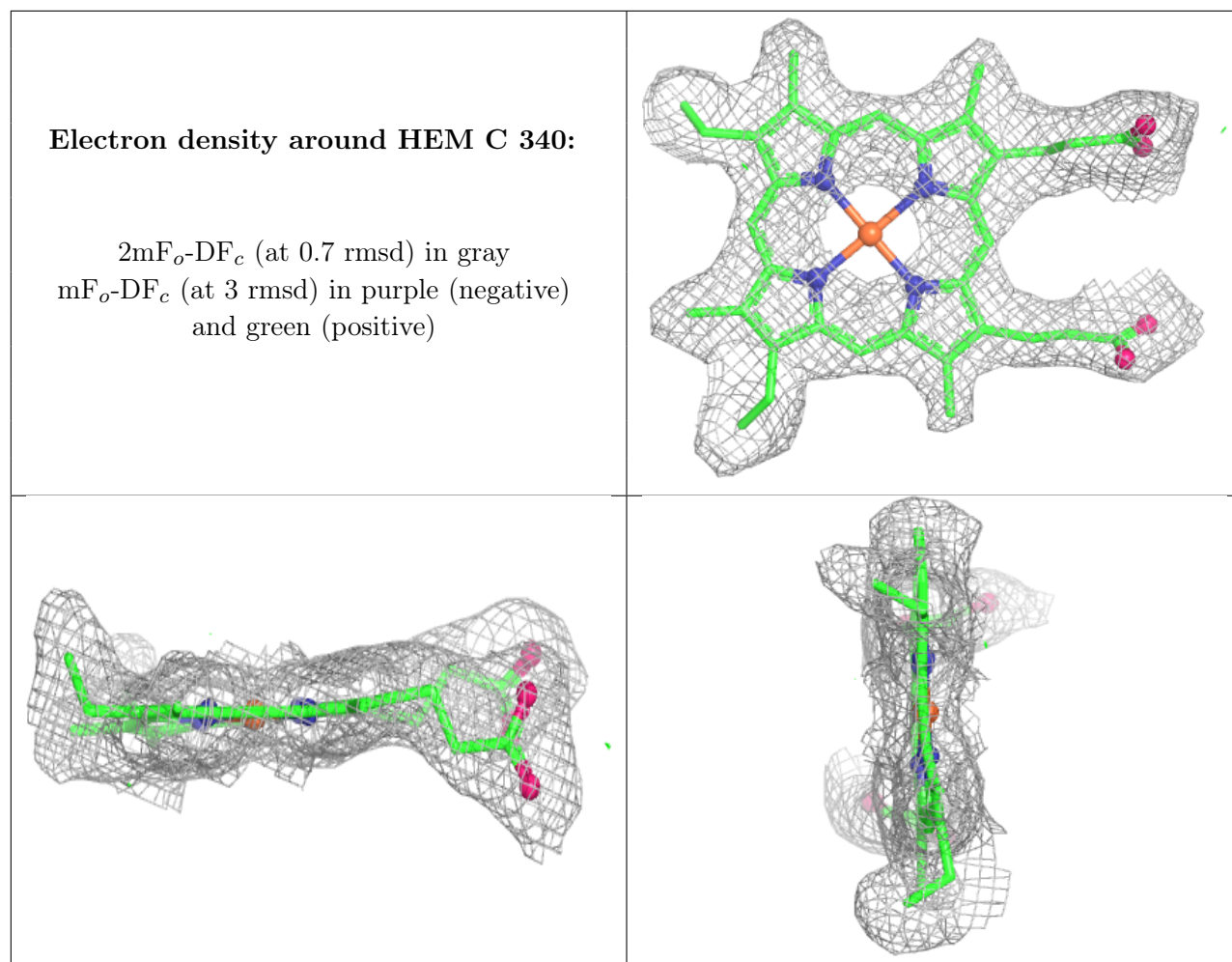
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

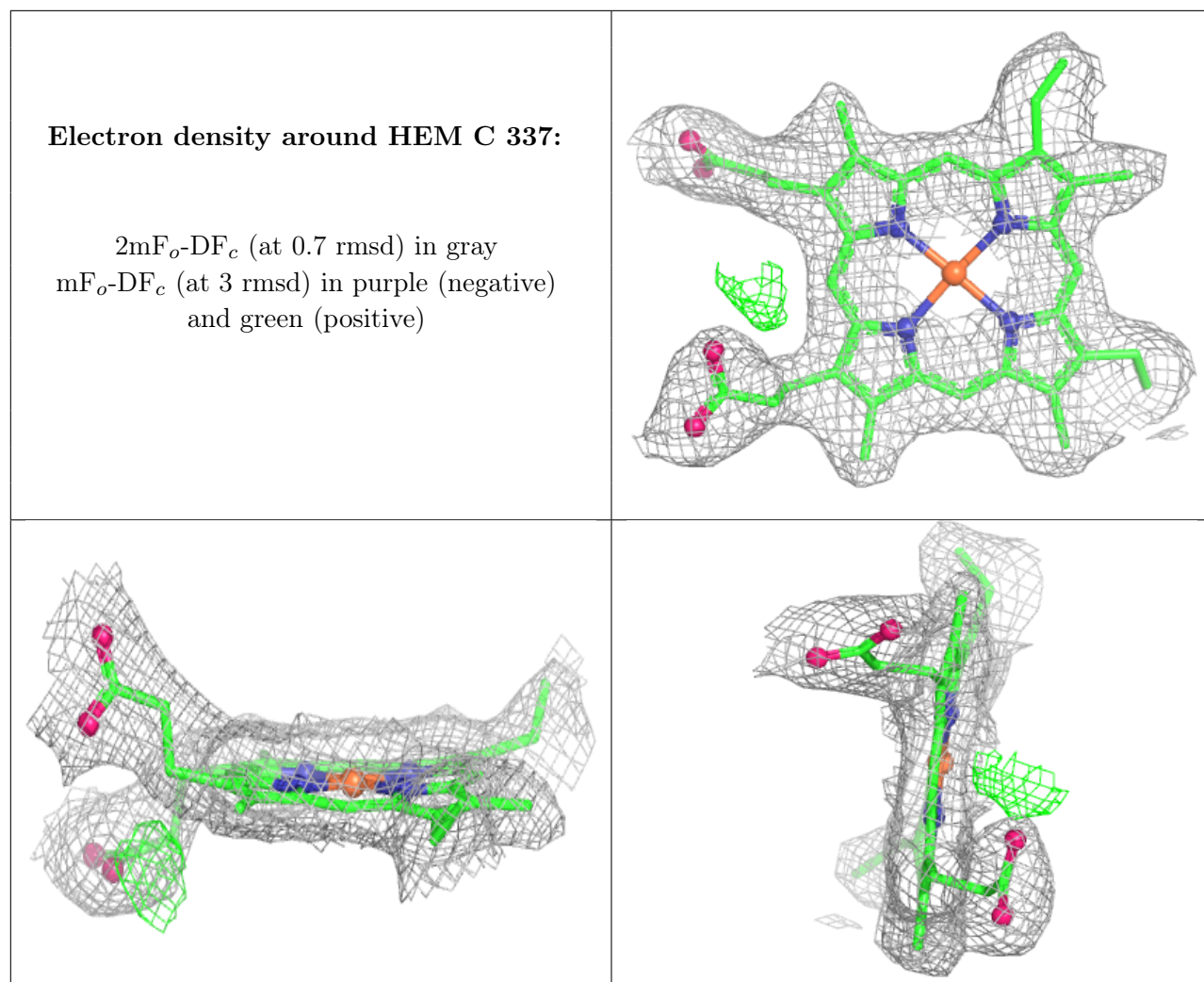


Electron density around HEM C 339:

$2mF_o-DF_c$ (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.