



wwPDB X-ray Structure Validation Summary Report ⓘ

Aug 15, 2026 – 08:05 AM EDT

PDB ID : 9ZS8 / pdb_00009zs8
Title : Structural and electrostatic analysis of a reaction center variant with high-yield charge separation along the alternative electron transfer path
Authors : Keable, S.M.; Kern, J.F.; Laible, P.D.; Hanson, D.K.
Deposited on : 2025-12-22
Resolution : 2.70 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)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

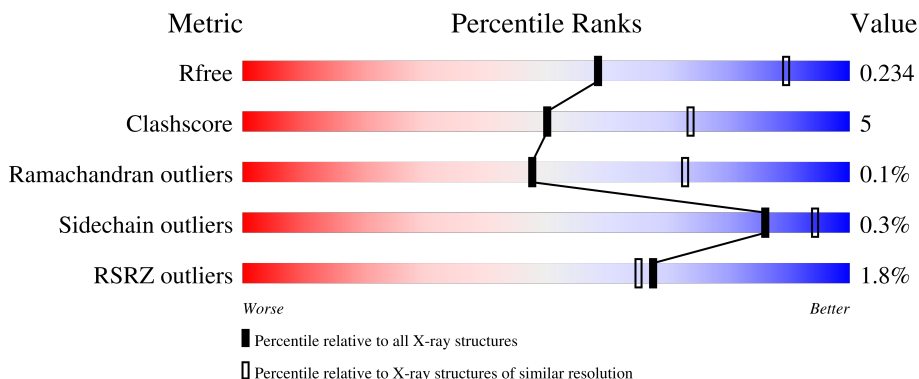
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	H	245	4% 85% 14%
2	L	281	89% 11%
3	M	302	% 89% 11%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 7190 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 Reaction center protein H chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	245	Total	C	N	O	S	0	1	0
			1862	1191	319	342	10			

- Molecule 2 is a protein called Reaction center protein L chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	L	281	Total	C	N	O	S	0	0	0
			2241	1512	357	364	8			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
L	181	TYR	PHE	engineered mutation	UNP Q3J1A5
L	185	TRP	LEU	engineered mutation	UNP Q3J1A5
L	193	THR	LEU	engineered mutation	UNP Q3J1A5
L	216	TRP	PHE	engineered mutation	UNP Q3J1A5

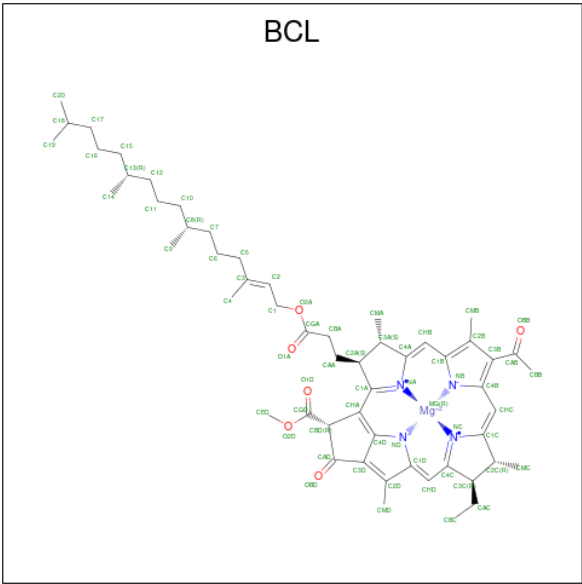
- Molecule 3 is a protein called Reaction center protein M chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	M	302	Total	C	N	O	S	0	1	0
			2417	1613	396	397	11			

There are 5 discrepancies between the modelled and reference sequences:

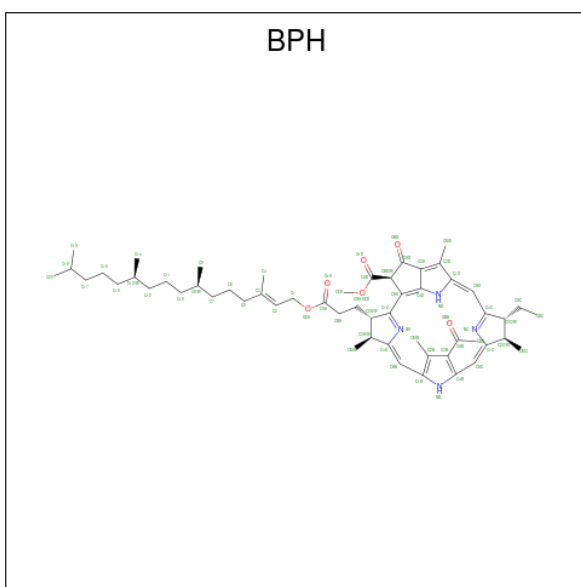
Chain	Residue	Modelled	Actual	Comment	Reference
M	133	GLU	THR	engineered mutation	UNP Q3J1A6
M	210	PHE	TYR	engineered mutation	UNP Q3J1A6
M	214	HIS	LEU	engineered mutation	UNP Q3J1A6
M	252	VAL	TRP	engineered mutation	UNP Q3J1A6
M	260	TRP	ALA	engineered mutation	UNP Q3J1A6

- Molecule 4 is BACTERIOCHLOROPHYLL A (CCD ID: BCL) (formula: $C_{55}H_{74}MgN_4O_6$) (labeled as "Ligand of Interest" by depositor).



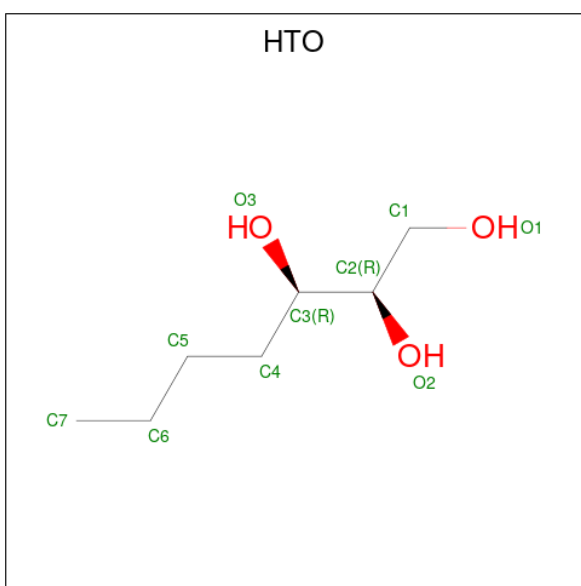
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
4	L	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
4	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
4	M	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

- Molecule 5 is BACTERIOPHEOPHYTIN A (CCD ID: BPH) (formula: $C_{55}H_{76}N_4O_6$) (la-
beled as "Ligand of Interest" by depositor).



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

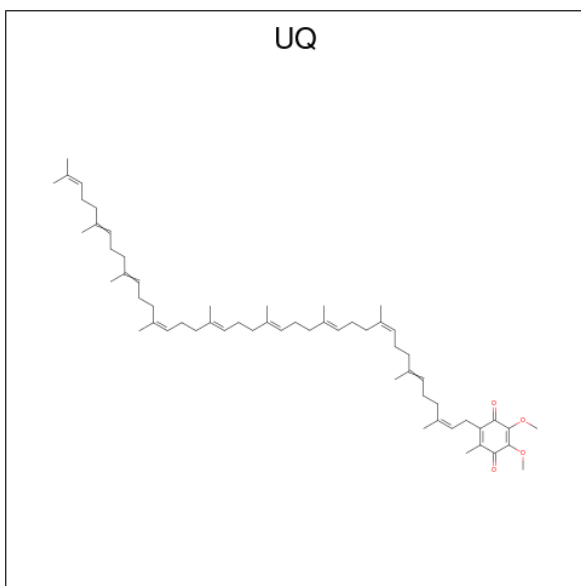
- Molecule 6 is HEPTANE-1,2,3-TRIOL (CCD ID: HTO) (formula: $C_7H_{16}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	L	1	Total	C	O	0	0
			10	7	3		

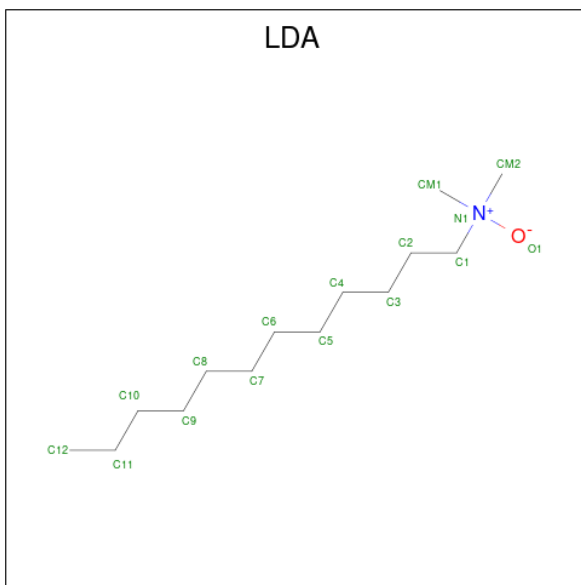
- Molecule 7 is Coenzyme Q10, (2Z,6E,10Z,14E,18E,22E,26Z)-isomer (CCD ID: UQ) (formula:

C₅₉H₉₀O₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	L	1	Total	C	O	0	0
			33	29	4		

- Molecule 8 is LAURYL DIMETHYLAMINE-N-OXIDE (CCD ID: LDA) (formula: C₁₄H₃₁NO).



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

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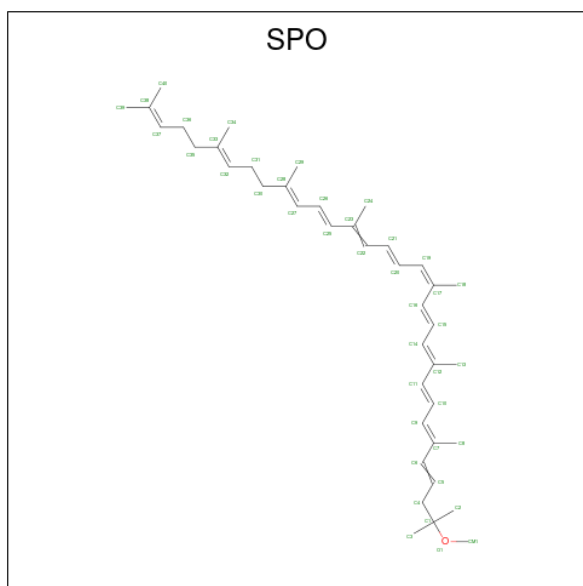
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
8	M	1	Total	C	N	O	0	0
			16	14	1	1		

- Molecule 9 is FE (III) ION (CCD ID: FE) (formula: Fe) (labeled as "Ligand of Interest" by depositor).

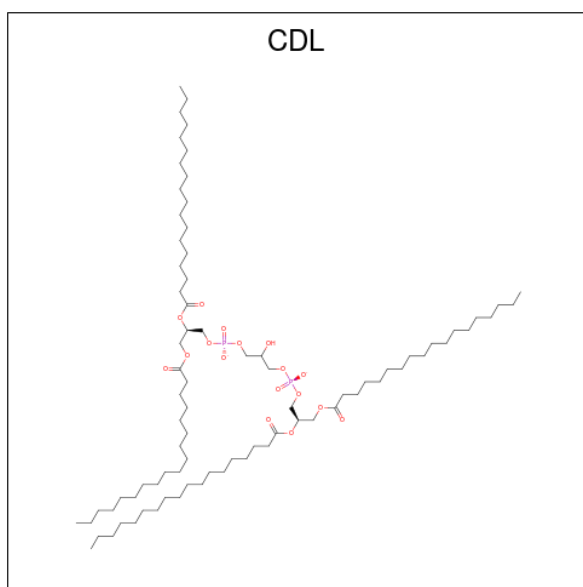
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	M	1	Total	Fe	0	0
			1	1		

- Molecule 10 is SPHEROIDENE (CCD ID: SPO) (formula: C₄₁H₆₀O).



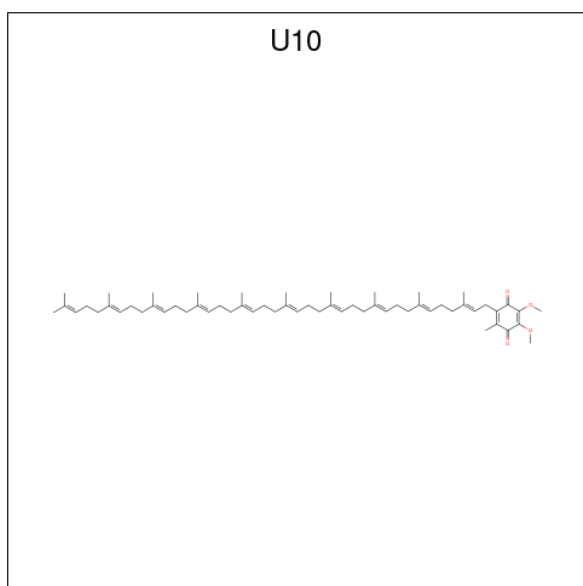
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	M	1	Total	C	O	0	0
			42	41	1		

- Molecule 11 is CARDIOLIPIN (CCD ID: CDL) (formula: C₈₁H₁₅₆O₁₇P₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	M	1	Total	C	O	P	0	0
			82	63	17	2		

- Molecule 12 is UBIQUINONE-10 (CCD ID: U10) (formula: $C_{59}H_{90}O_4$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	M	1	Total	C	O	0	0
			15	11	4		

- Molecule 13 is MAGNESIUM ION (CCD ID: MG) (formula: Mg) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
13	M	1	Total 1	Mg 1	0	0

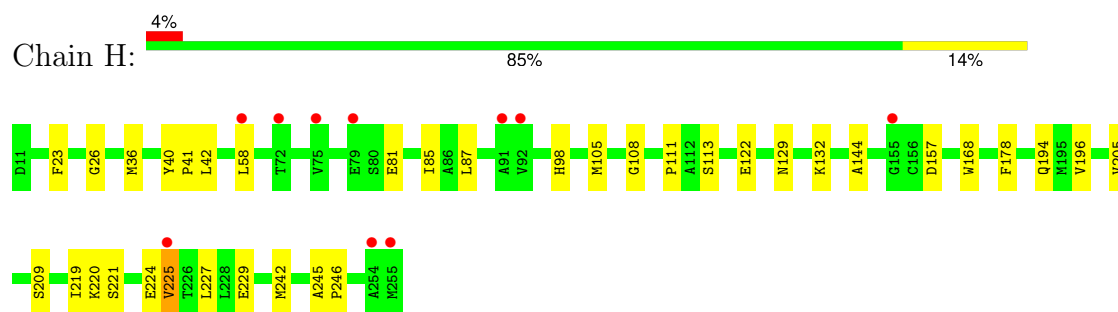
- Molecule 14 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	H	10	Total 10	O 10	0	0
14	L	15	Total 15	O 15	0	0
14	M	19	Total 19	O 19	0	0

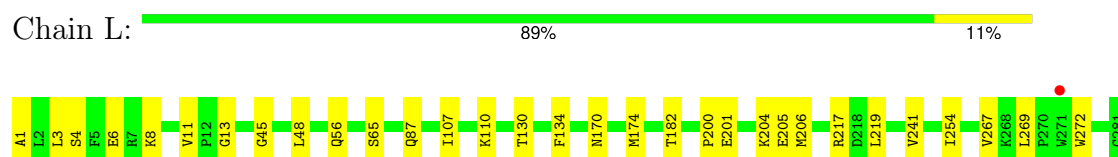
3 Residue-property plots

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.

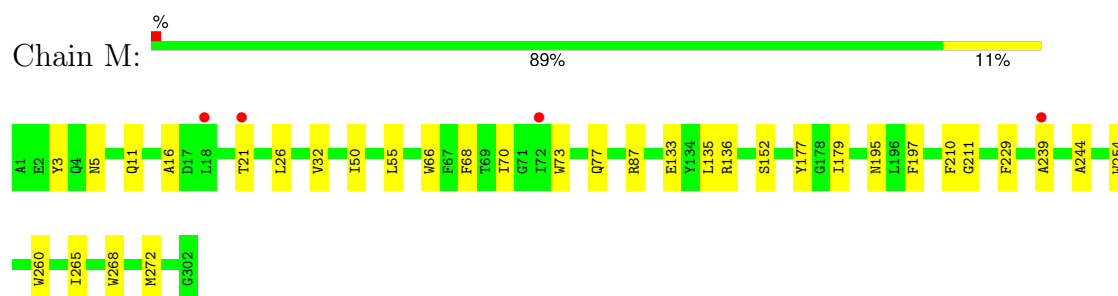
- Molecule 1: Reaction center protein H chain



- Molecule 2: Reaction center protein L chain



- Molecule 3: Reaction center protein M chain



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	142.05Å 139.65Å 77.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.84 – 2.70 44.84 – 2.70	Depositor EDS
% Data completeness (in resolution range)	97.2 (44.84-2.70) 97.1 (44.84-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.00 (at 2.69Å)	Xtriage
Refinement program	PHENIX 2.0_5824	Depositor
R, R_{free}	0.213 , 0.233 0.214 , 0.234	Depositor DCC
R_{free} test set	1996 reflections (4.63%)	wwPDB-VP
Wilson B-factor (Å ²)	79.0	Xtriage
Anisotropy	0.173	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 77.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7190	wwPDB-VP
Average B, all atoms (Å ²)	84.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: SPO, LDA, U10, FE, BPH, UQ, CDL, HTO, BCL, MG

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	H	0.11	0/1917	0.28	0/2608
2	L	0.10	0/2332	0.28	1/3195 (0.0%)
3	M	0.09	0/2513	0.24	0/3428
All	All	0.10	0/6762	0.26	1/9231 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	254	ILE	N-CA-C	-5.09	107.87	112.96

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	H	1862	0	1872	25	0
2	L	2241	0	2183	22	0
3	M	2417	0	2325	27	0
4	L	132	0	148	3	0
4	M	132	0	148	9	0
5	L	65	0	74	5	0
5	M	65	0	76	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	L	10	0	16	1	0
7	L	33	0	39	2	0
8	L	16	0	31	0	0
8	M	32	0	62	0	0
9	M	1	0	0	0	0
10	M	42	0	60	4	0
11	M	82	0	108	4	0
12	M	15	0	11	2	0
13	M	1	0	0	0	0
14	H	10	0	0	0	0
14	L	15	0	0	0	0
14	M	19	0	0	0	0
All	All	7190	0	7153	76	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 5.

The worst 5 of 76 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)
3:M:260:TRP:HE1	3:M:265:ILE:HD12	1.52	0.75
1:H:168:TRP:HB2	1:H:178:PHE:HB2	1.74	0.69
4:M:401:BCL:HBB3	4:M:402:BCL:H41	1.78	0.65
2:L:56:GLN:HE22	2:L:65:SER:H	1.46	0.64
5:L:302:BPH:HBB2	3:M:210:PHE:HB3	1.81	0.61

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	H	244/245 (100%)	236 (97%)	8 (3%)	0	100	100
2	L	279/281 (99%)	269 (96%)	10 (4%)	0	100	100
3	M	301/302 (100%)	292 (97%)	8 (3%)	1 (0%)	36	60
All	All	824/828 (100%)	797 (97%)	26 (3%)	1 (0%)	48	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	M	195	ASN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	199/198 (100%)	198 (100%)	1 (0%)	81	92
2	L	220/220 (100%)	219 (100%)	1 (0%)	81	92
3	M	238/237 (100%)	238 (100%)	0	100	100
All	All	657/655 (100%)	655 (100%)	2 (0%)	86	94

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	225	VAL
2	L	4	SER

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

Mol	Chain	Res	Type
1	H	68	HIS
1	H	126	HIS
2	L	56	GLN
2	L	62	GLN
2	L	159	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 16 ligands modelled in this entry, 2 are monoatomic - leaving 14 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	BCL	L	301	2	69,74,74	1.13	4 (5%)	79,115,115	1.30	7 (8%)
5	BPH	L	302	13	59,70,70	1.00	3 (5%)	59,101,101	1.39	5 (8%)
10	SPO	M	407	-	41,41,41	0.36	0	47,50,50	0.75	3 (6%)
4	BCL	M	402	3	69,74,74	1.15	5 (7%)	79,115,115	1.41	8 (10%)
6	HTO	L	304	-	9,9,9	0.41	0	10,10,10	0.82	0
4	BCL	M	401	3	69,74,74	1.14	4 (5%)	79,115,115	1.42	10 (12%)
8	LDA	L	306	-	13,15,15	0.75	1 (7%)	14,17,17	0.41	0
11	CDL	M	408	-	81,81,99	0.32	0	87,93,111	0.24	0
5	BPH	M	406	-	59,70,70	1.01	3 (5%)	59,101,101	1.28	5 (8%)
4	BCL	L	303	2	69,74,74	1.11	4 (5%)	79,115,115	1.24	8 (10%)
7	UQ	L	305	-	33,33,63	0.46	0	42,43,79	0.67	1 (2%)
12	U10	M	409	-	15,15,63	2.62	7 (46%)	20,21,79	1.51	3 (15%)
8	LDA	M	403	-	13,15,15	0.75	1 (7%)	14,17,17	0.38	0
8	LDA	M	404	-	13,15,15	0.75	1 (7%)	14,17,17	0.36	0

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	BCL	L	301	2	-	5/41/137/137	-
5	BPH	L	302	13	-	9/37/105/105	0/5/6/6
10	SPO	M	407	-	-	8/47/47/47	-
4	BCL	M	402	3	-	3/41/137/137	-
6	HTO	L	304	-	-	4/10/10/10	-
4	BCL	M	401	3	-	6/41/137/137	-
8	LDA	L	306	-	-	4/13/13/13	-
11	CDL	M	408	-	-	41/92/92/110	-
5	BPH	M	406	-	-	10/37/105/105	0/5/6/6
4	BCL	L	303	2	-	0/41/137/137	-
7	UQ	L	305	-	-	9/27/51/87	0/1/1/1
12	U10	M	409	-	-	1/6/30/87	0/1/1/1
8	LDA	M	403	-	-	6/13/13/13	-
8	LDA	M	404	-	-	7/13/13/13	-

The worst 5 of 33 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
12	M	409	U10	O4-C4	-5.57	1.23	1.36
12	M	409	U10	O3-C3	-5.11	1.24	1.36
4	M	402	BCL	MG-NA	4.94	2.18	2.06
5	M	406	BPH	C1B-C2B	4.84	1.44	1.39
4	L	301	BCL	MG-NA	4.77	2.17	2.06

The worst 5 of 50 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L	302	BPH	C4D-CHA-CBD	-7.00	105.10	108.45
5	M	406	BPH	C4D-CHA-CBD	-5.74	105.70	108.45
4	M	402	BCL	C4D-CHA-C1A	5.44	127.73	121.24
4	L	301	BCL	C4D-CHA-C1A	5.21	127.46	121.24
4	M	401	BCL	C4D-CHA-C1A	4.88	127.07	121.24

There are no chirality outliers.

5 of 113 torsion outliers are listed below:

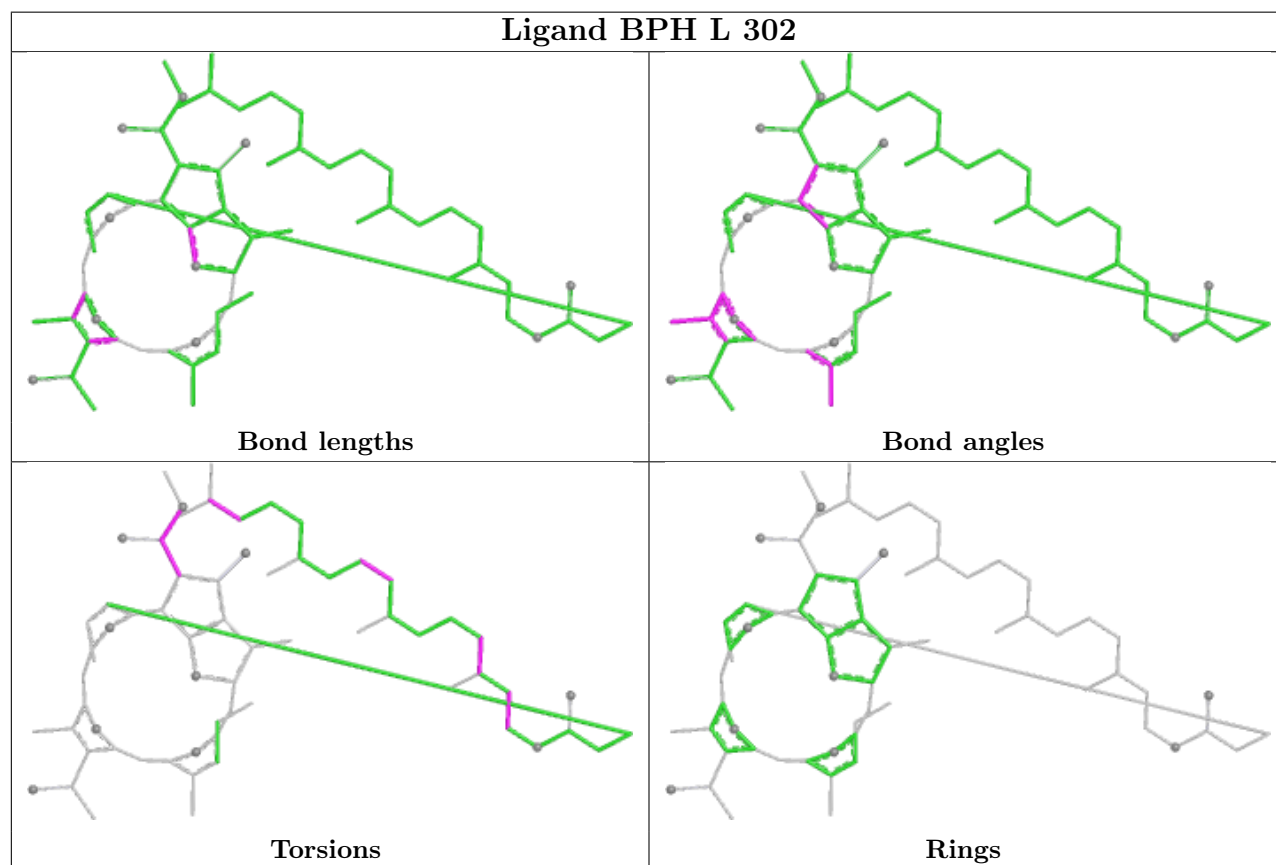
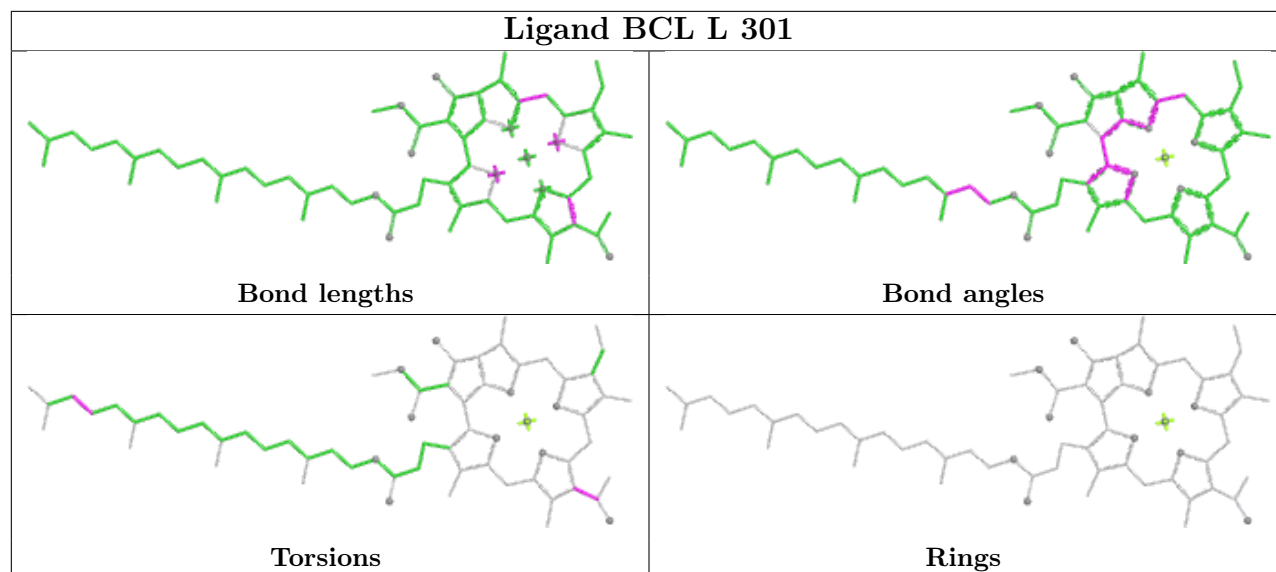
Mol	Chain	Res	Type	Atoms
6	L	304	HTO	C2-C3-C4-C5
7	L	305	UQ	C12-C11-C9-C8
7	L	305	UQ	C12-C11-C9-C10
8	L	306	LDA	N1-C1-C2-C3
10	M	407	SPO	C1-C4-C5-C6

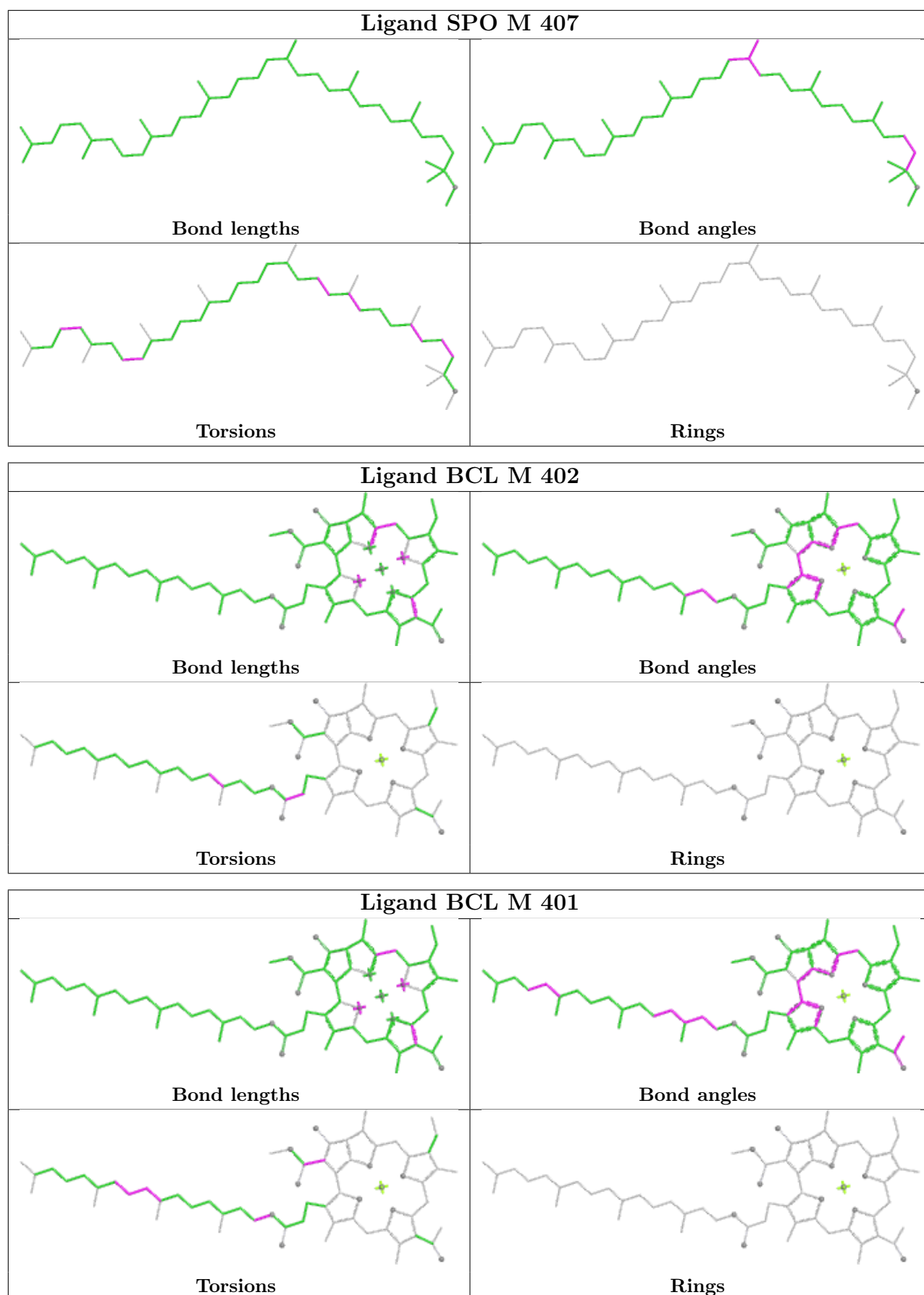
There are no ring outliers.

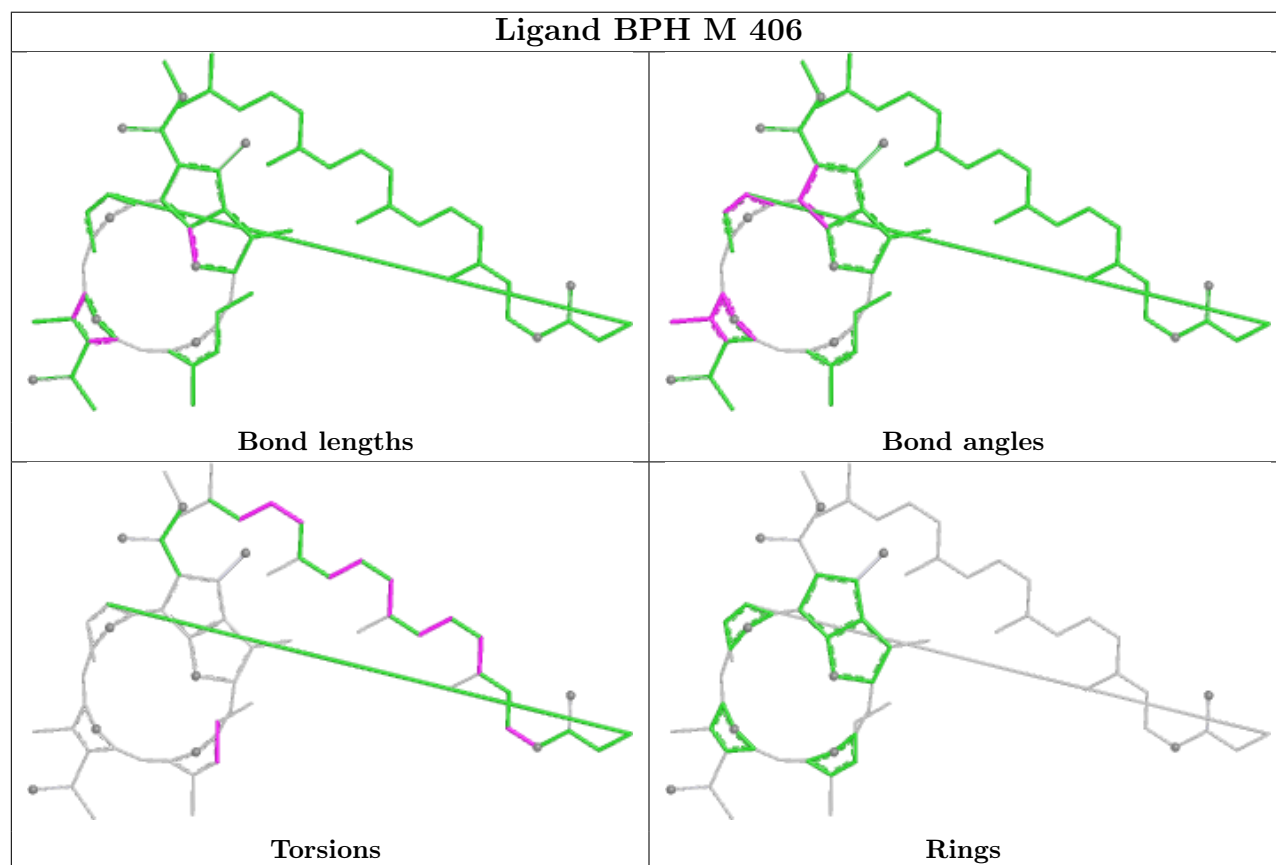
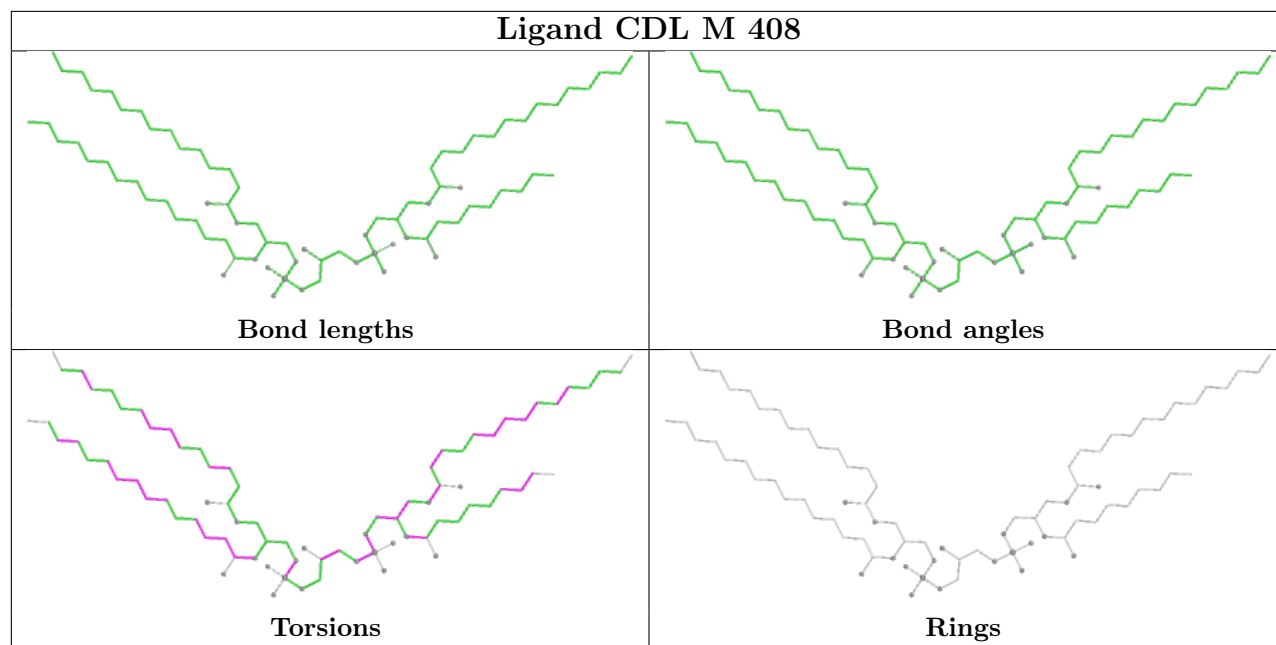
11 monomers are involved in 28 short contacts:

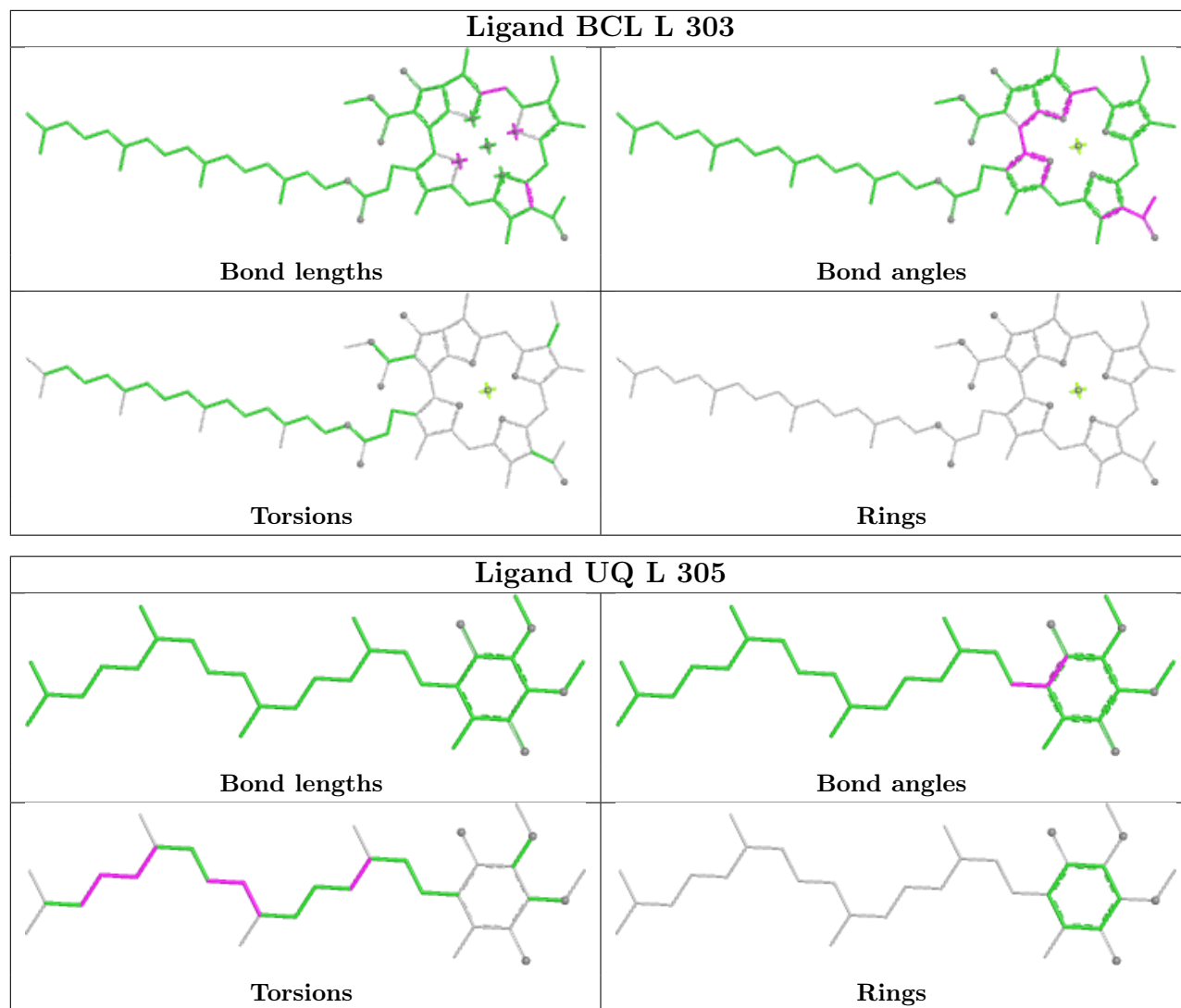
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	301	BCL	1	0
5	L	302	BPH	5	0
10	M	407	SPO	4	0
4	M	402	BCL	5	0
6	L	304	HTO	1	0
4	M	401	BCL	5	0
11	M	408	CDL	4	0
5	M	406	BPH	2	0
4	L	303	BCL	2	0
7	L	305	UQ	2	0
12	M	409	U10	2	0

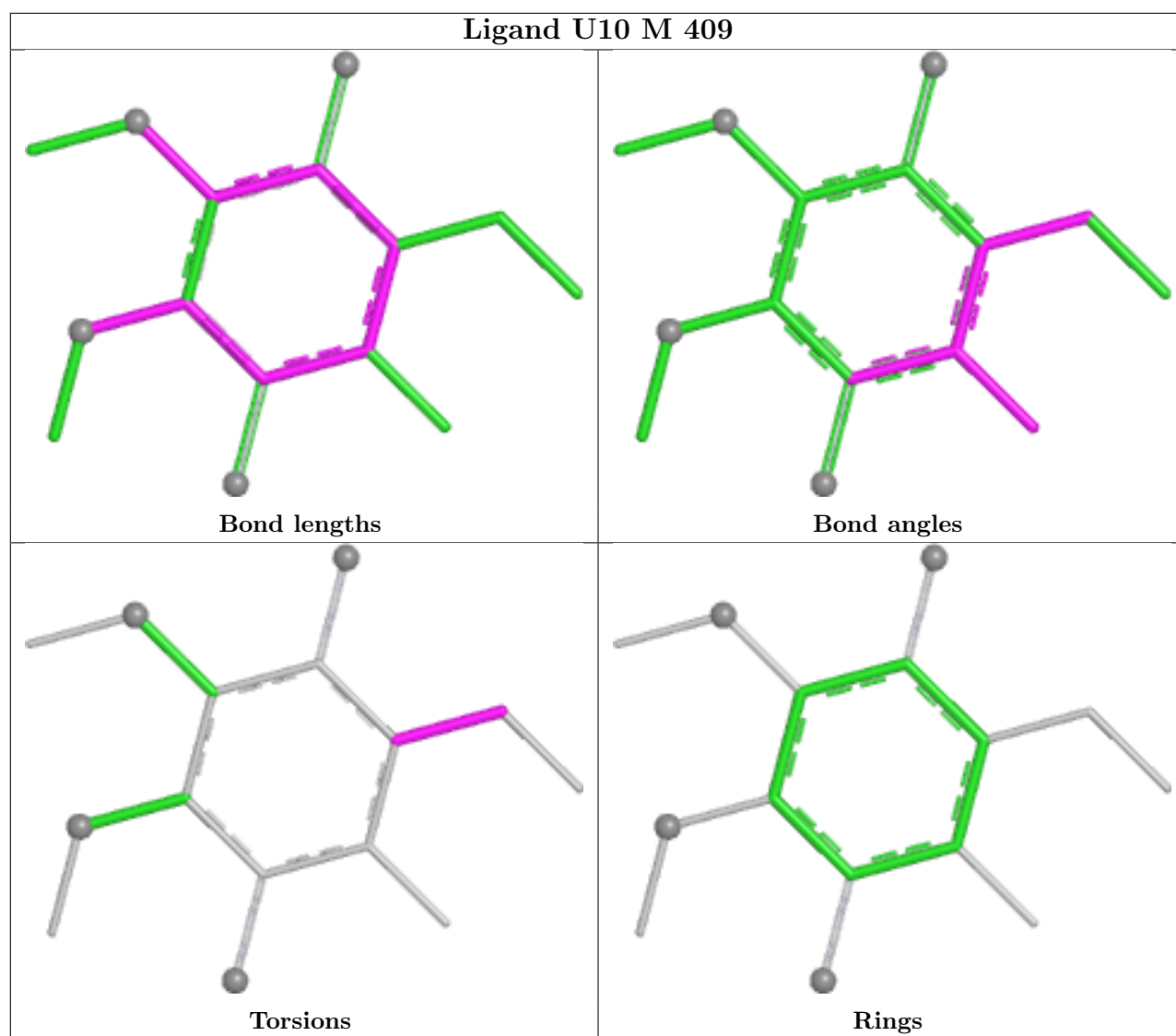
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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	H	245/245 (100%)	0.37	10 (4%) 41 37	66, 89, 125, 147	1 (0%)
2	L	281/281 (100%)	-0.02	1 (0%) 88 87	56, 80, 108, 137	0
3	M	302/302 (100%)	0.05	4 (1%) 75 73	41, 79, 105, 151	1 (0%)
All	All	828/828 (100%)	0.12	15 (1%) 67 65	41, 83, 114, 151	2 (0%)

The worst 5 of 15 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	L	271	TRP	3.5
1	H	254	ALA	3.0
1	H	72	THR	2.9
1	H	75	VAL	2.5
1	H	255	MET	2.5

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

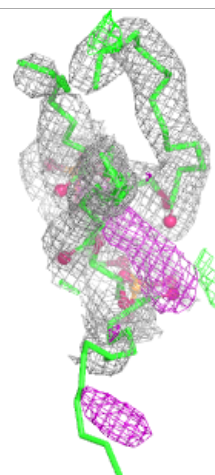
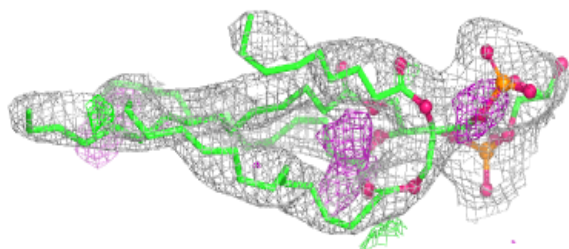
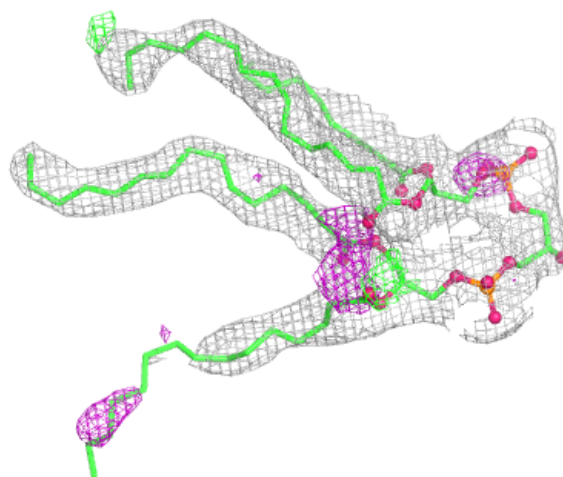
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	HTO	L	304	10/10	0.64	0.20	100,110,113,133	0
8	LDA	L	306	16/16	0.81	0.32	83,104,113,121	0
8	LDA	M	404	16/16	0.82	0.23	80,92,132,133	0
11	CDL	M	408	82/100	0.82	0.19	77,112,132,145	0
10	SPO	M	407	42/42	0.88	0.19	63,85,98,113	0
8	LDA	M	403	16/16	0.89	0.24	87,101,109,110	0
7	UQ	L	305	33/63	0.90	0.20	69,86,113,120	0
12	U10	M	409	15/63	0.91	0.26	93,105,112,115	15
4	BCL	M	401	66/66	0.93	0.14	56,67,104,114	0
5	BPH	M	406	65/65	0.93	0.15	54,73,144,154	0
4	BCL	M	402	66/66	0.95	0.10	44,61,87,123	0
5	BPH	L	302	65/65	0.96	0.10	52,69,81,89	0
4	BCL	L	301	66/66	0.96	0.10	56,68,90,97	0
4	BCL	L	303	66/66	0.97	0.08	51,62,91,95	0
9	FE	M	405	1/1	1.00	0.02	66,66,66,66	0
13	MG	M	410	1/1	1.00	0.04	69,69,69,69	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.

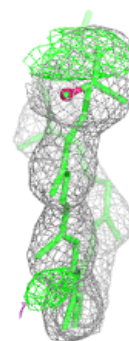
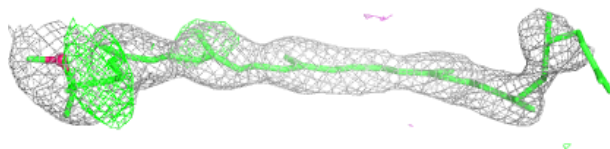
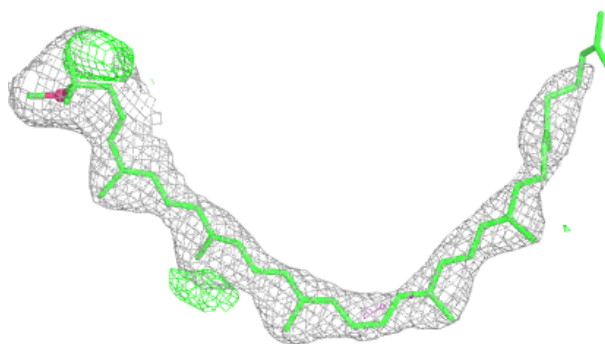
Electron density around CDL M 408:

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

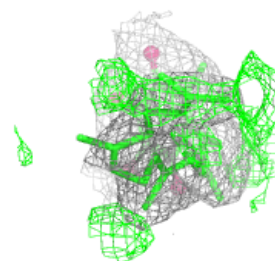
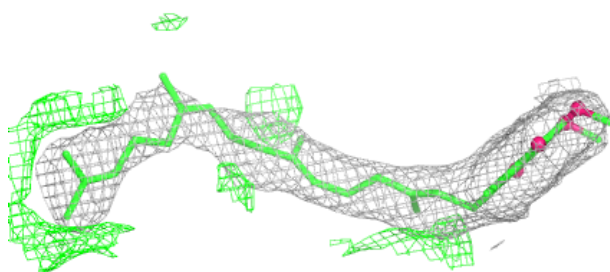
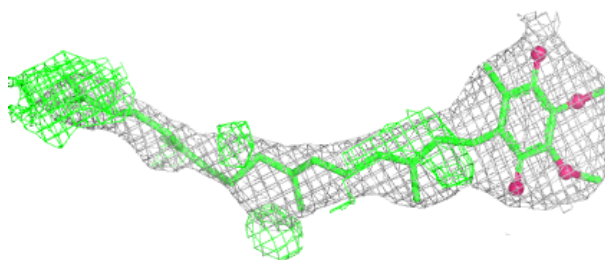


Electron density around SPO M 407:

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

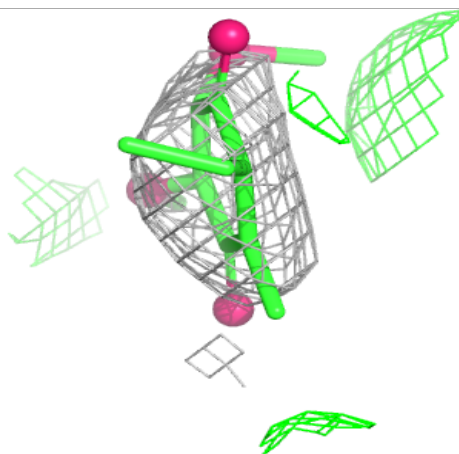
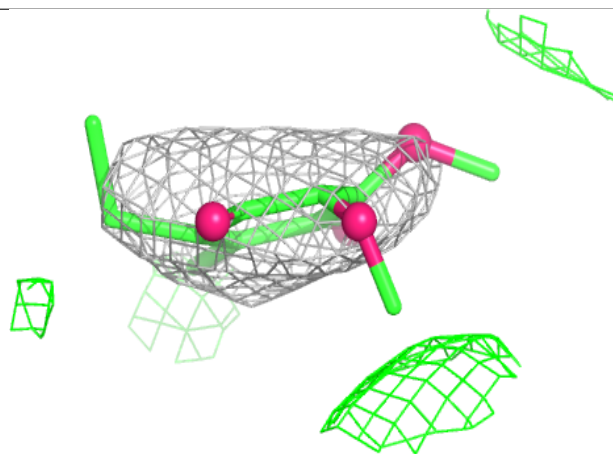
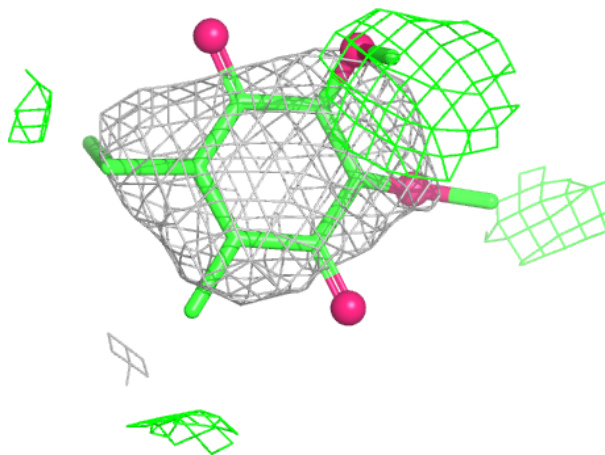
**Electron density around UQ L 305:**

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



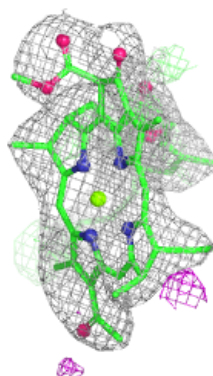
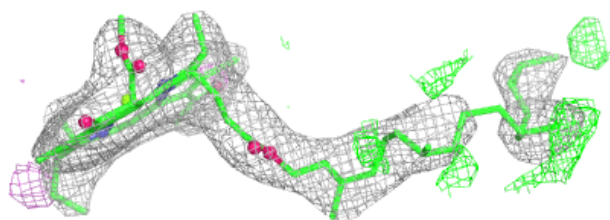
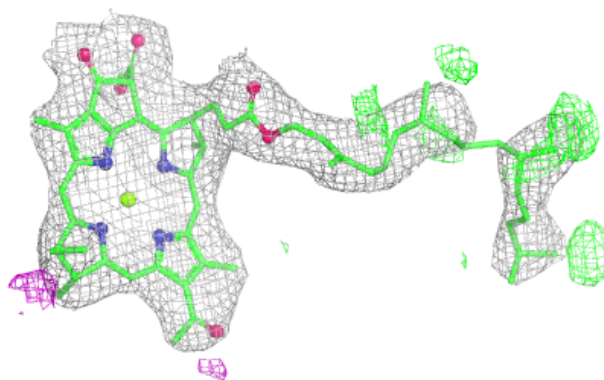
Electron density around U10 M 409:

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

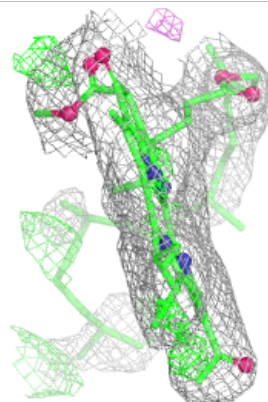
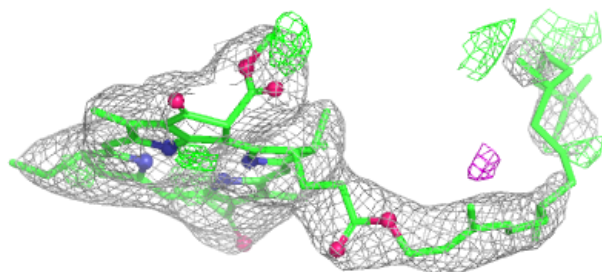
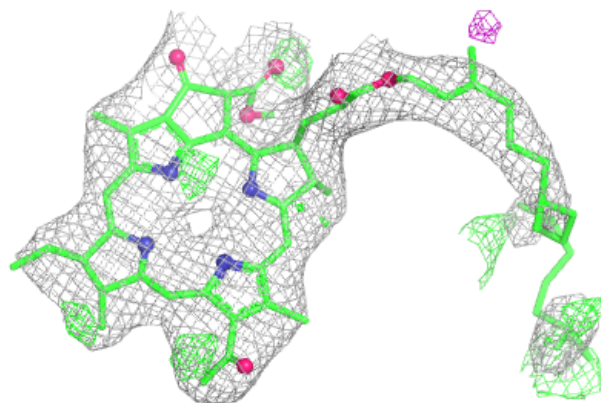


Electron density around BCL M 401:

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

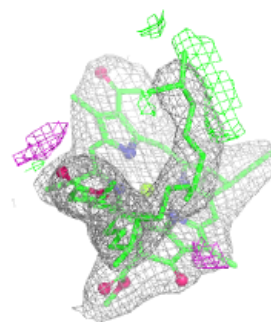
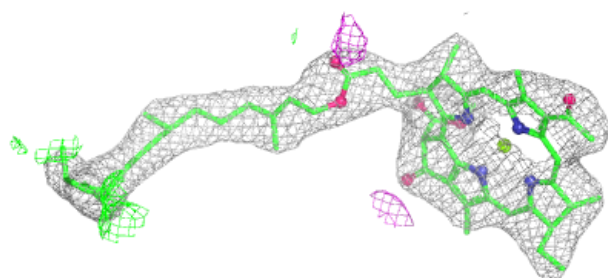
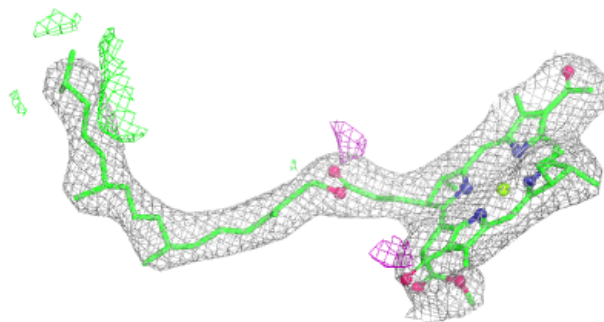
**Electron density around BPH M 406:**

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



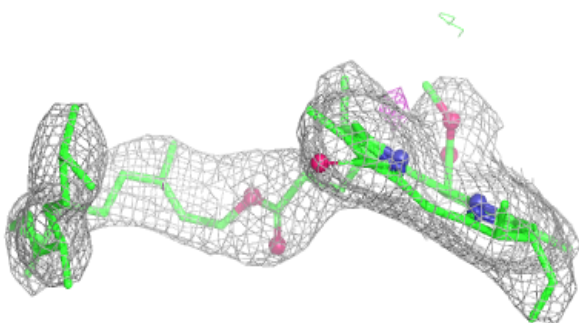
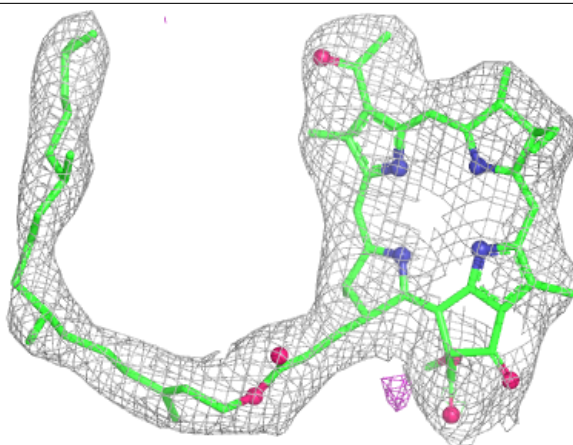
Electron density around BCL M 402:

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



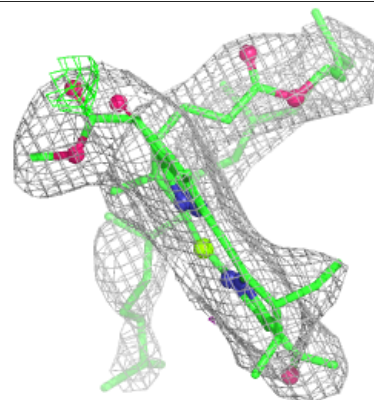
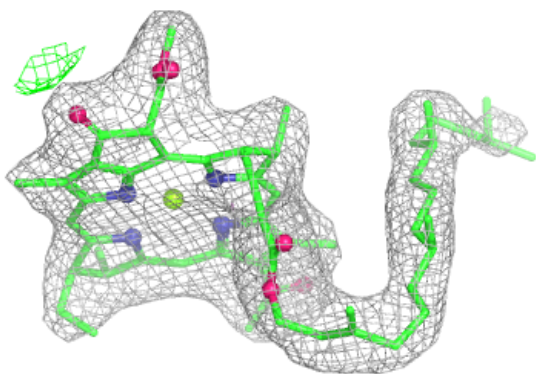
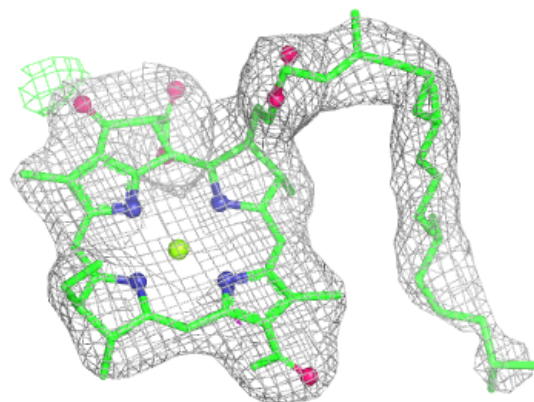
Electron density around BPH L 302:

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

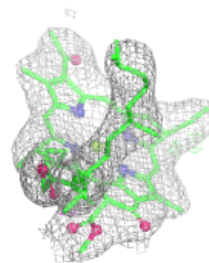
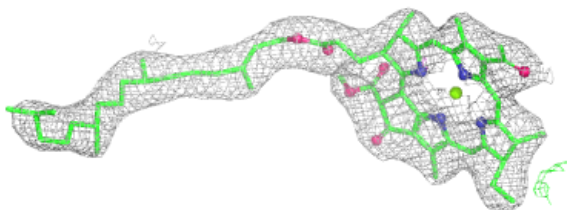
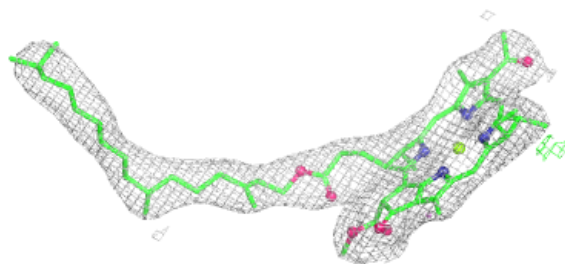


Electron density around BCL L 301:

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

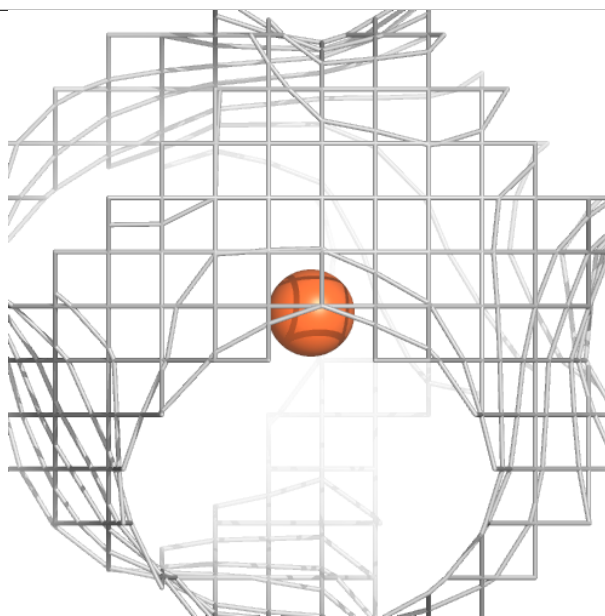
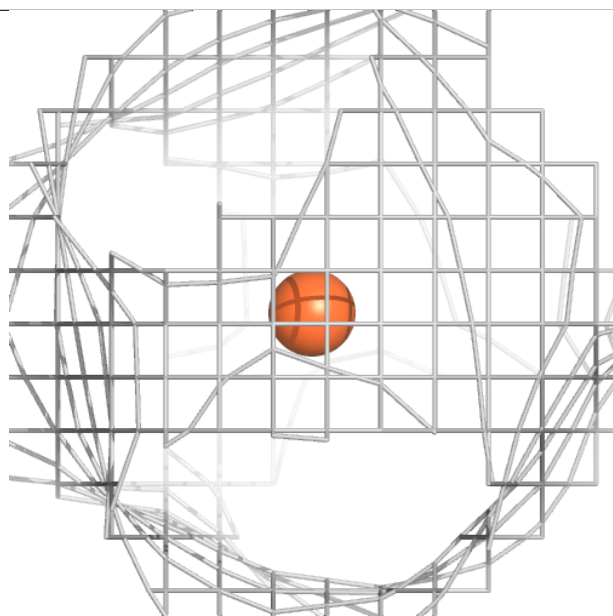
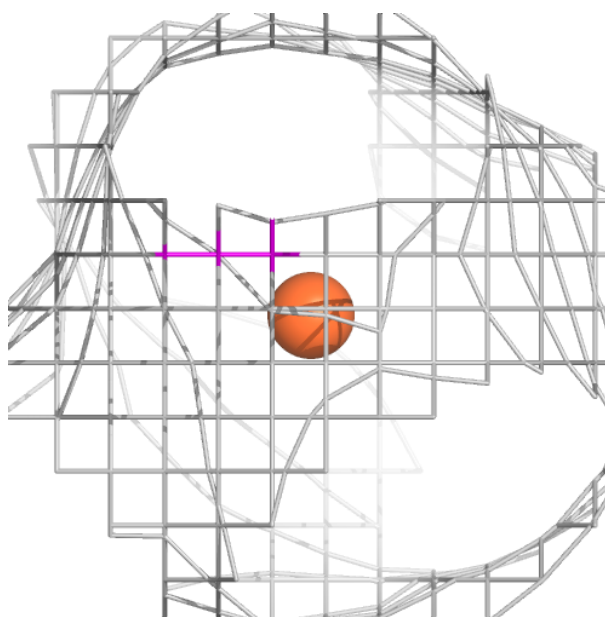
**Electron density around BCL L 303:**

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



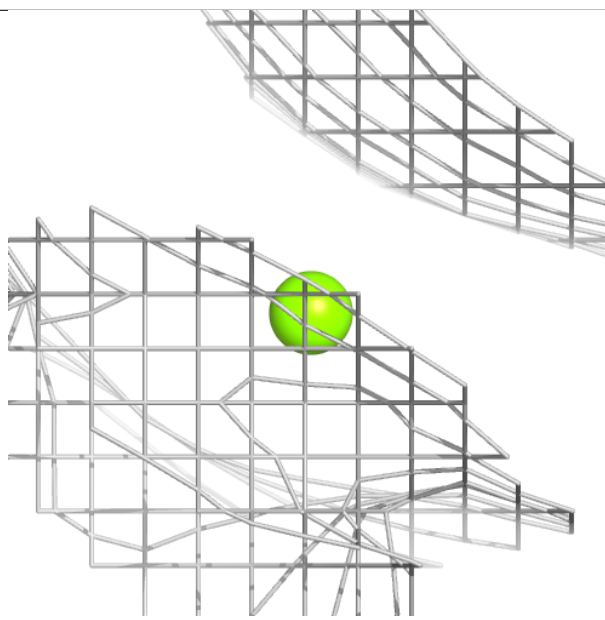
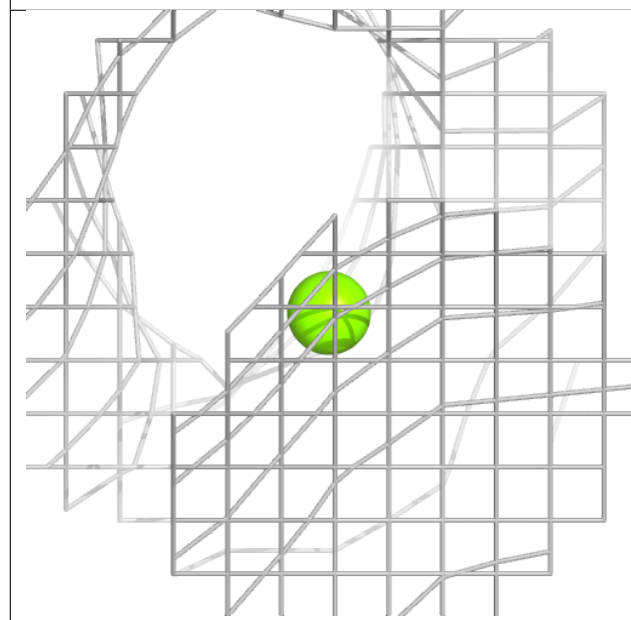
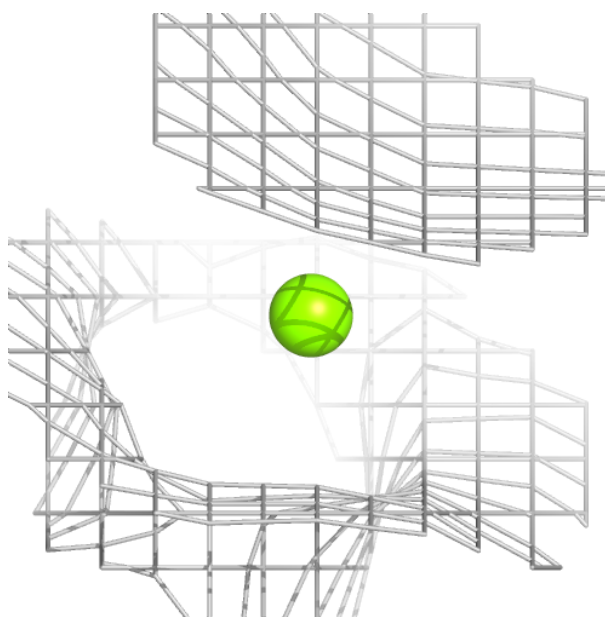
Electron density around FE M 405:

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



Electron density around MG M 410:

$2mF_o-DF_c$ (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.