



# wwPDB X-ray Structure Validation Summary Report ⓘ

Dec 5, 2023 – 10:26 am GMT

PDB ID : 2UXJ  
Title : X-ray high resolution structure of the photosynthetic reaction center from Rb. sphaeroides at pH 10 in the neutral state  
Authors : Koepke, J.; Diehm, R.; Fritzsich, G.  
Deposited on : 2007-03-28  
Resolution : 2.25 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](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-467  
Mogul : 1.8.4, CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.36  
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) : 2.36

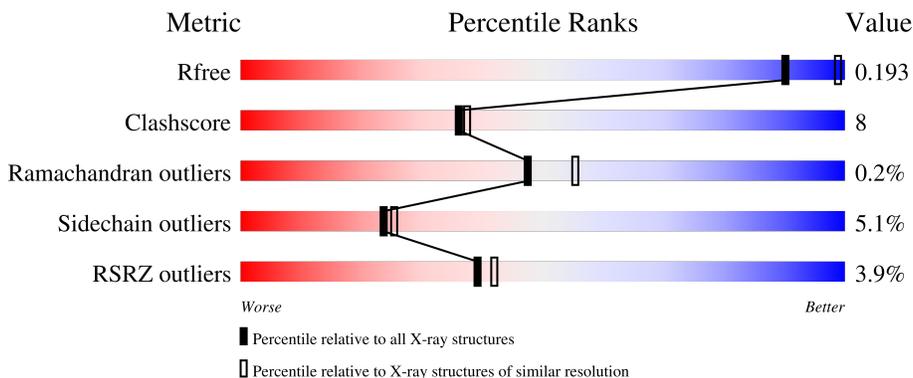
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 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))
$R_{free}$	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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

Mol	Chain	Length	Quality of chain
1	H	260	
2	L	281	
3	M	307	

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
10	HTO	L	1904	-	-	-	X
14	CDL	M	1316	-	-	-	X
5	BCL	L	1282	X	-	-	-
5	BCL	L	1287	X	-	-	-
5	BCL	M	1303	X	-	-	-
5	BCL	M	1304	X	-	-	-
6	LDA	L	1284	-	-	-	X
8	UQ2	L	1286[B]	-	-	X	-

## 2 Entry composition [i](#)

There are 15 unique types of molecules in this entry. The entry contains 7707 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
			Total	C	N	O	S			
1	H	241	1846	1181	319	337	9	0	3	1

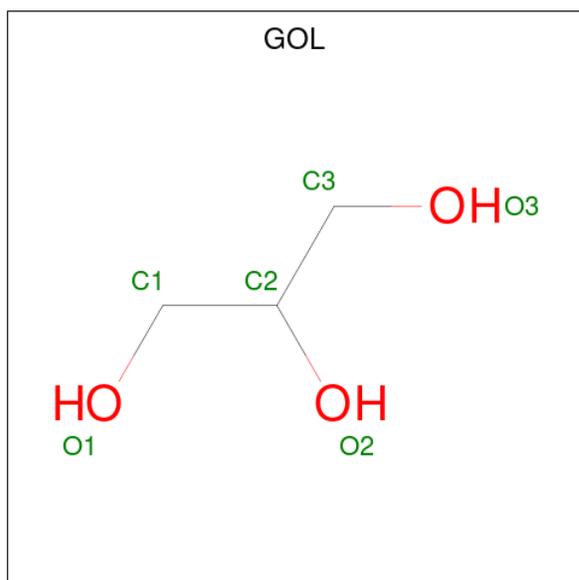
- Molecule 2 is a protein called REACTION CENTER PROTEIN L CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	L	281	2232	1507	355	362	8	0	0	0

- Molecule 3 is a protein called REACTION CENTER PROTEIN M CHAIN.

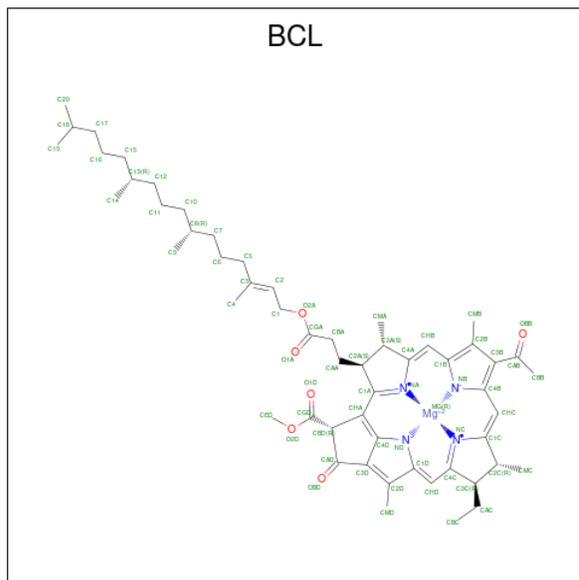
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	M	303	2409	1607	395	397	10	0	0	1

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: C<sub>3</sub>H<sub>8</sub>O<sub>3</sub>).



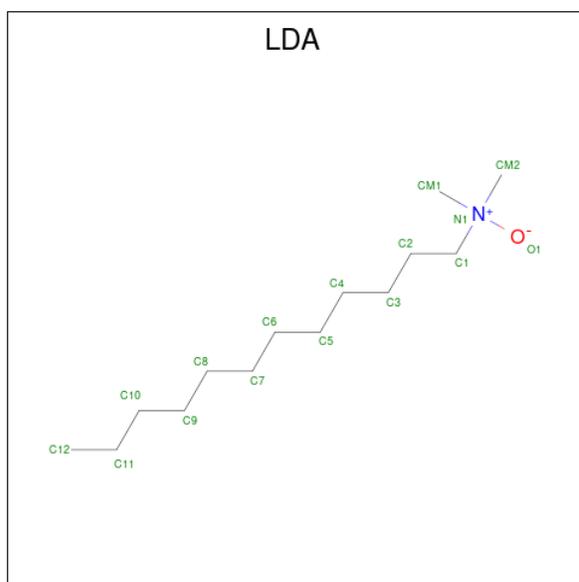
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	H	1	Total C O 6 3 3	0	0
4	H	1	Total C O 6 3 3	0	0
4	H	1	Total C O 6 3 3	0	0
4	L	1	Total C O 6 3 3	0	0
4	L	1	Total C O 6 3 3	0	0
4	L	1	Total C O 6 3 3	0	0

- Molecule 5 is BACTERIOCHLOROPHYLL A (three-letter code: BCL) (formula:  $C_{55}H_{74}MgN_4O_6$ ).



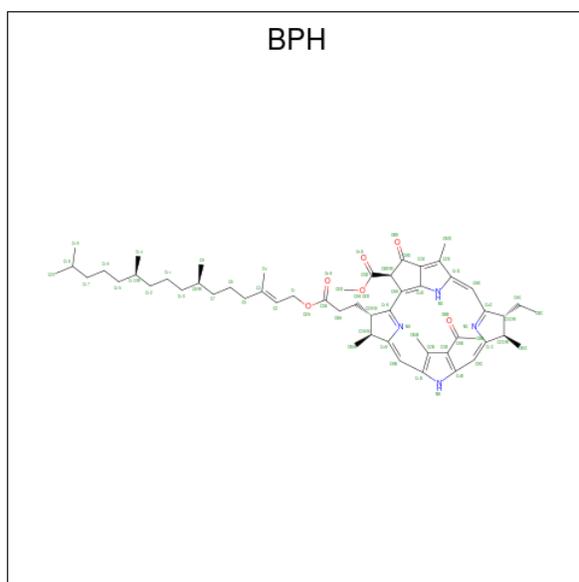
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	L	1	Total C Mg N O 66 55 1 4 6	0	0
5	L	1	Total C Mg N O 66 55 1 4 6	0	0
5	M	1	Total C Mg N O 66 55 1 4 6	0	0
5	M	1	Total C Mg N O 66 55 1 4 6	0	0

- Molecule 6 is LAURYL DIMETHYLAMINE-N-OXIDE (three-letter code: LDA) (formula:  $C_{14}H_{31}NO$ ).



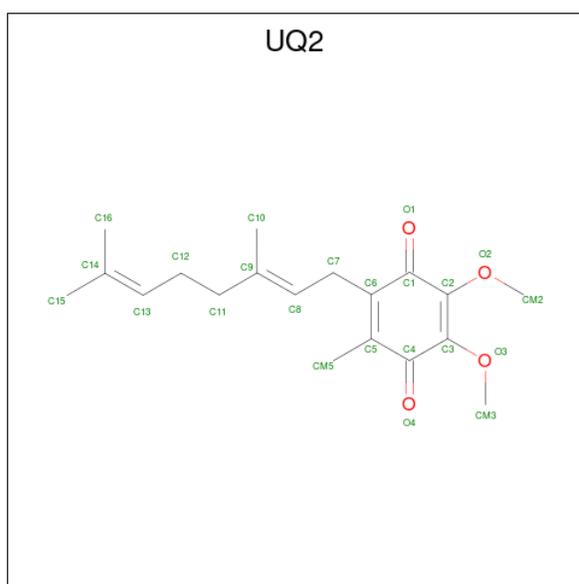
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
6	L	1	Total 16	C 14	N 1	O 1	0	0
6	L	1	Total 16	C 14	N 1	O 1	0	0
6	M	1	Total 16	C 14	N 1	O 1	0	0
6	M	1	Total 16	C 14	N 1	O 1	0	0
6	M	1	Total 16	C 14	N 1	O 1	0	0
6	M	1	Total 16	C 14	N 1	O 1	0	0
6	M	1	Total 16	C 14	N 1	O 1	0	0
6	M	1	Total 16	C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

- Molecule 7 is BACTERIOPHEOPHYTIN A (three-letter code: BPH) (formula: C<sub>55</sub>H<sub>76</sub>N<sub>4</sub>O<sub>6</sub>).



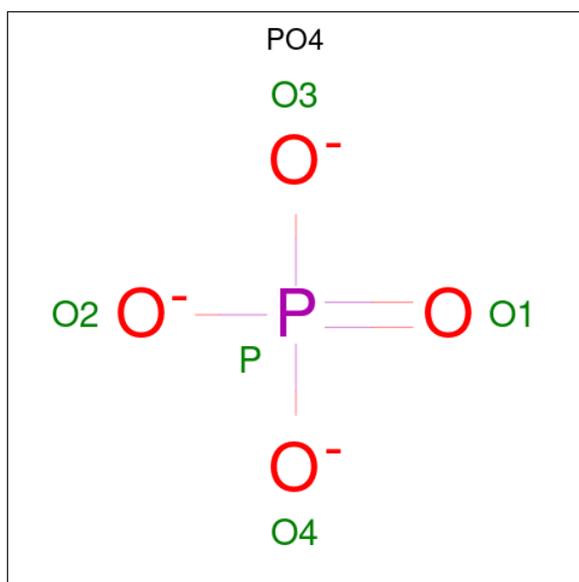
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	L	1	Total	C	N	O	0	0
			65	55	4	6		
7	M	1	Total	C	N	O	0	0
			65	55	4	6		

- Molecule 8 is UBIQUINONE-2 (three-letter code: UQ2) (formula:  $C_{19}H_{26}O_4$ ).



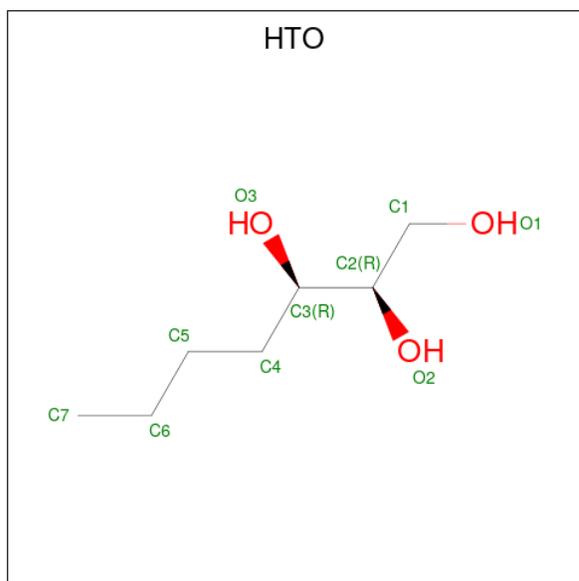
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	L	1	Total	C	O	0	1
			46	38	8		

- Molecule 9 is PHOSPHATE ION (three-letter code: PO4) (formula:  $O_4P$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	L	1	Total	O	P	0	0
			5	4	1		

- Molecule 10 is HEPTANE-1,2,3-TRIOL (three-letter code: HTO) (formula:  $C_7H_{16}O_3$ ).

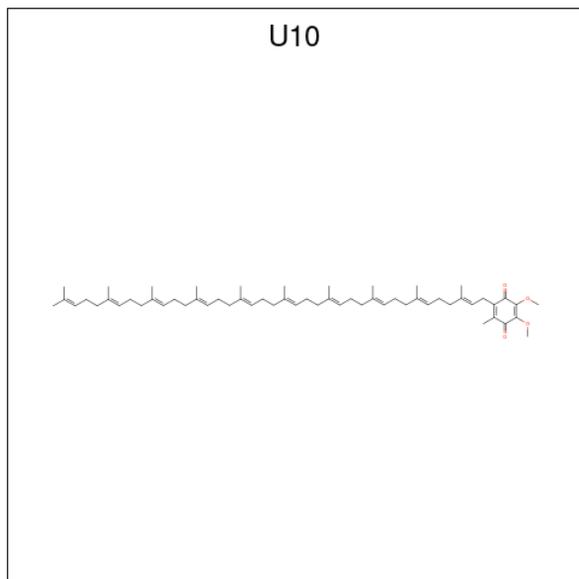


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	L	1	Total	C	O	0	0
			10	7	3		
10	L	1	Total	C	O	0	0
			10	7	3		

- Molecule 11 is FE (III) ION (three-letter code: FE) (formula: Fe).

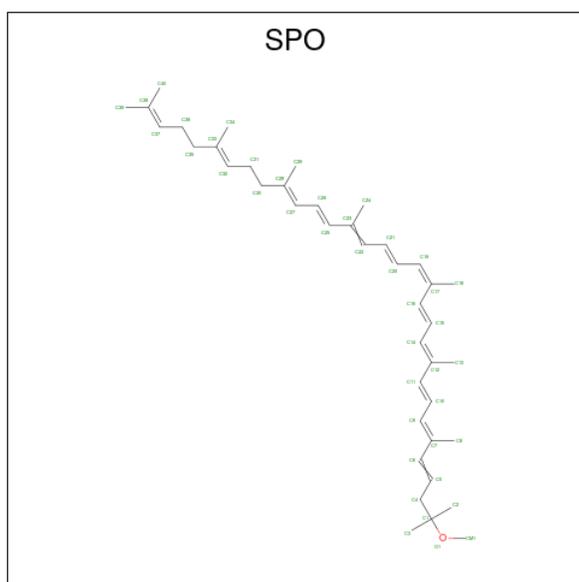
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	M	1	Total	Fe	0	0
			1	1		

- Molecule 12 is UBIQUINONE-10 (three-letter code: U10) (formula:  $C_{59}H_{90}O_4$ ).



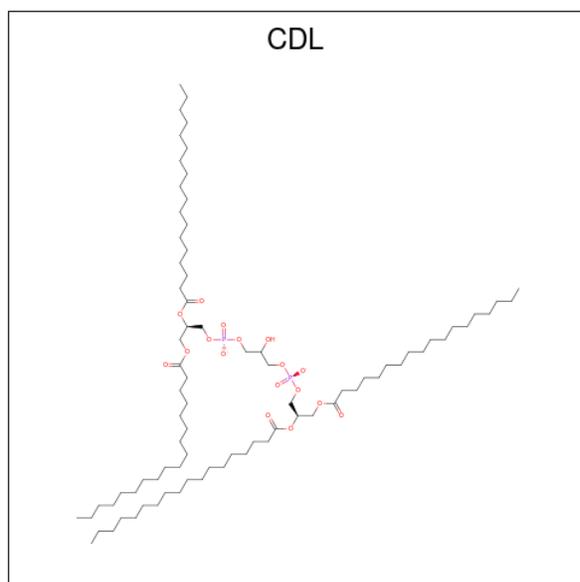
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
12	M	1	Total	C	O	0	0
			48	44	4		

- Molecule 13 is SPHEROIDENE (three-letter code: SPO) (formula:  $C_{41}H_{60}O$ ).



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

- Molecule 14 is CARDIOLIPIN (three-letter code: CDL) (formula:  $C_{81}H_{156}O_{17}P_2$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
14	M	1	81	62	17	2	0	0

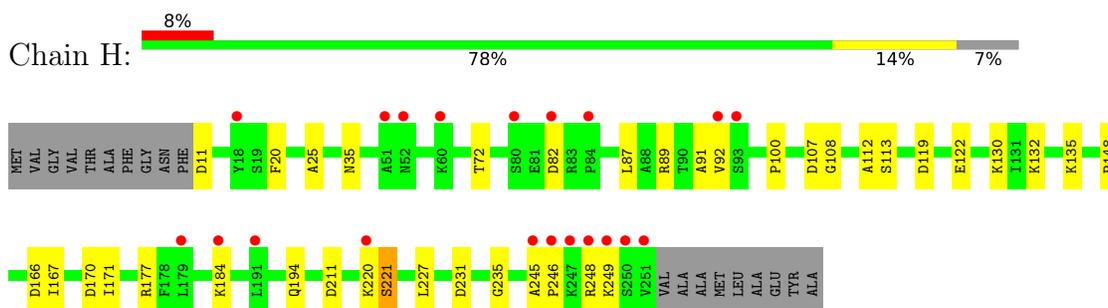
- Molecule 15 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
15	H	122	122	122	0	0
15	L	134	134	134	0	0
15	M	147	147	147	0	0

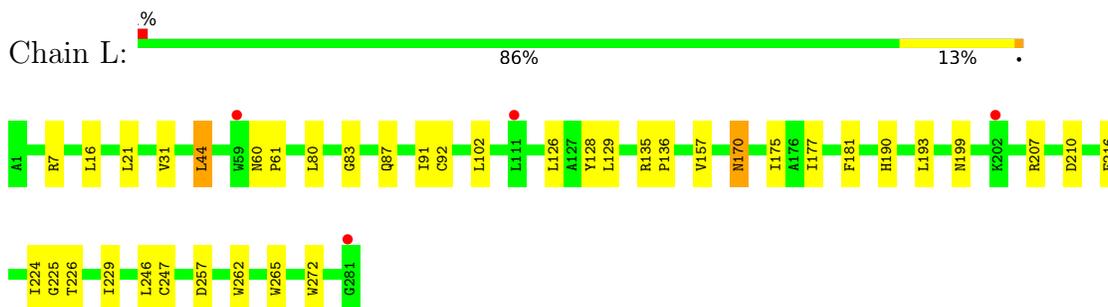
### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

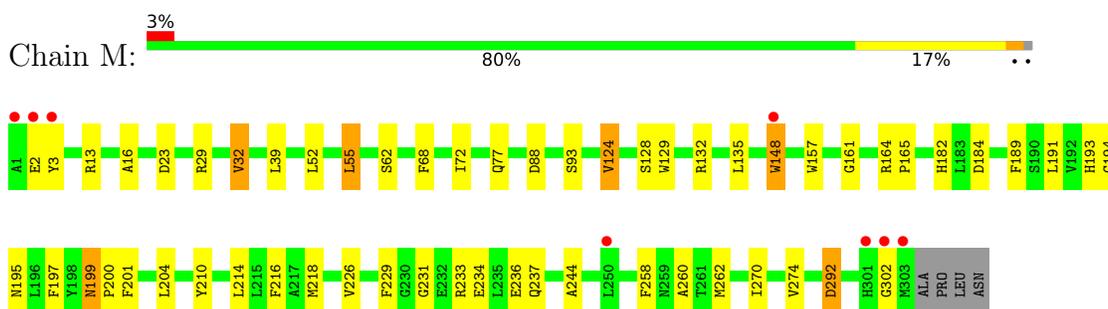
- Molecule 1: REACTION CENTER PROTEIN H CHAIN



- 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 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	139.38Å 139.38Å 235.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	119.52 – 2.25 27.12 – 2.25	Depositor EDS
% Data completeness (in resolution range)	94.0 (119.52-2.25) 90.1 (27.12-2.25)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.47 (at 2.24Å)	Xtrriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.194 , 0.225 0.192 , 0.193	Depositor DCC
$R_{free}$ test set	5205 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	39.0	Xtrriage
Anisotropy	0.158	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 57.3	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	7707	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	49.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.05% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: HTO, U10, SPO, FE, BPH, LDA, PO4, UQ2, GOL, BCL, CDL

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	0.63	0/1906	0.79	6/2591 (0.2%)
2	L	0.73	0/2320	0.71	1/3175 (0.0%)
3	M	0.69	0/2501	0.71	5/3415 (0.1%)
All	All	0.69	0/6727	0.73	12/9181 (0.1%)

There are no bond length outliers.

The worst 5 of 12 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	H	11	ASP	CB-CG-OD2	7.73	125.25	118.30
1	H	82	ASP	CB-CG-OD2	6.77	124.39	118.30
1	H	107	ASP	CB-CG-OD2	5.88	123.60	118.30
3	M	184	ASP	CB-CG-OD1	5.74	123.46	118.30
3	M	234	GLU	OE1-CD-OE2	-5.62	116.56	123.30

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	1846	0	1861	22	0
2	L	2232	0	2187	27	0
3	M	2409	0	2321	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	18	0	24	1	0
4	L	18	0	24	2	0
5	L	132	0	148	8	0
5	M	132	0	148	16	0
6	L	32	0	62	0	0
6	M	112	0	217	8	0
7	L	65	0	76	7	0
7	M	65	0	76	9	0
8	L	46	0	52	13	0
9	L	5	0	0	0	0
10	L	20	0	32	2	0
11	M	1	0	0	0	0
12	M	48	0	63	2	0
13	M	42	0	60	4	0
14	M	81	0	92	1	0
15	H	122	0	0	0	0
15	L	134	0	0	2	0
15	M	147	0	0	2	0
All	All	7707	0	7443	118	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 118 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:218:MET:SD	3:M:218:MET:CE	2.04	1.43
2:L:224:ILE:HG22	8:L:1286[B]:UQ2:C8	1.87	1.04
2:L:224:ILE:HG22	8:L:1286[B]:UQ2:H8	0.99	0.97
2:L:224:ILE:CG2	8:L:1286[B]:UQ2:H8	1.96	0.93
7:L:1285:BPH:HHC	7:L:1285:BPH:HBB3	1.51	0.91

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	242/260 (93%)	235 (97%)	7 (3%)	0	100	100
2	L	279/281 (99%)	269 (96%)	9 (3%)	1 (0%)	34	37
3	M	301/307 (98%)	288 (96%)	12 (4%)	1 (0%)	41	46
All	All	822/848 (97%)	792 (96%)	28 (3%)	2 (0%)	47	55

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	M	302	GLY
2	L	31	VAL

### 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	H	198/208 (95%)	193 (98%)	5 (2%)	47	56
2	L	220/220 (100%)	206 (94%)	14 (6%)	17	16
3	M	236/240 (98%)	222 (94%)	14 (6%)	19	19
All	All	654/668 (98%)	621 (95%)	33 (5%)	24	26

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	M	182	HIS
3	M	191	LEU
3	M	292	ASP
2	L	170	ASN
2	L	129	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 7 such sidechains are listed below:

Mol	Chain	Res	Type
3	M	77	GLN
3	M	187	ASN
3	M	199	ASN
3	M	193	HIS
2	L	264	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](#)

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

### 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

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

Of 30 ligands modelled in this entry, 1 is monoatomic - leaving 29 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$
7	BPH	L	1285	-	51,70,70	2.77	7 (13%)	52,101,101	2.10	14 (26%)
4	GOL	L	1291	-	5,5,5	0.42	0	5,5,5	0.34	0
6	LDA	L	1283	-	12,15,15	1.96	1 (8%)	14,17,17	0.63	0
5	BCL	M	1304	3	64,74,74	2.11	9 (14%)	78,115,115	2.09	18 (23%)
8	UQ2	L	1286[A]	-	23,23,23	2.74	8 (34%)	28,31,31	1.33	5 (17%)
4	GOL	H	1253	-	5,5,5	0.38	0	5,5,5	0.23	0
6	LDA	M	1307	-	12,15,15	1.94	1 (8%)	14,17,17	0.55	0
6	LDA	M	1310	-	12,15,15	2.03	1 (8%)	14,17,17	0.44	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
6	LDA	M	1311	-	12,15,15	2.08	1 (8%)	14,17,17	0.62	0
4	GOL	H	1252	-	5,5,5	0.38	0	5,5,5	0.30	0
6	LDA	L	1284	-	12,15,15	1.99	1 (8%)	14,17,17	0.48	0
4	GOL	H	1251	-	5,5,5	0.45	0	5,5,5	0.26	0
8	UQ2	L	1286[B]	-	23,23,23	2.59	9 (39%)	28,31,31	2.23	9 (32%)
6	LDA	M	1305	-	12,15,15	2.06	1 (8%)	14,17,17	0.63	0
10	HTO	L	1289	-	9,9,9	0.41	0	10,10,10	0.56	0
5	BCL	L	1287	2	64,74,74	2.01	11 (17%)	78,115,115	2.04	20 (25%)
5	BCL	M	1303	3	64,74,74	2.01	11 (17%)	78,115,115	2.05	19 (24%)
12	U10	M	1314	-	48,48,63	2.72	12 (25%)	58,61,79	1.53	12 (20%)
6	LDA	M	1309	-	12,15,15	2.07	1 (8%)	14,17,17	0.47	0
4	GOL	L	1290	-	5,5,5	0.18	0	5,5,5	0.59	0
14	CDL	M	1316	-	80,80,99	1.93	16 (20%)	86,92,111	2.97	17 (19%)
6	LDA	M	1308	-	12,15,15	2.07	1 (8%)	14,17,17	0.47	0
13	SPO	M	1315	-	40,41,41	3.94	12 (30%)	47,50,50	2.28	17 (36%)
5	BCL	L	1282	2	64,74,74	1.94	10 (15%)	78,115,115	2.24	23 (29%)
6	LDA	M	1306	-	12,15,15	2.04	1 (8%)	14,17,17	0.59	0
7	BPH	M	1313	-	51,70,70	2.62	9 (17%)	52,101,101	2.02	10 (19%)
9	PO4	L	1288	-	4,4,4	0.86	0	6,6,6	0.47	0
10	HTO	L	1904	-	9,9,9	0.34	0	10,10,10	1.06	1 (10%)
4	GOL	L	1292	-	5,5,5	0.42	0	5,5,5	0.38	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
7	BPH	L	1285	-	-	6/37/105/105	0/5/6/6
4	GOL	L	1291	-	-	2/4/4/4	-
6	LDA	L	1283	-	-	7/13/13/13	-
5	BCL	M	1304	3	2/2/21/25	5/37/137/137	-
8	UQ2	L	1286[A]	-	-	9/15/39/39	0/1/1/1
4	GOL	H	1253	-	-	3/4/4/4	-
6	LDA	M	1307	-	-	9/13/13/13	-
6	LDA	M	1310	-	-	6/13/13/13	-
6	LDA	M	1311	-	-	5/13/13/13	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	H	1252	-	-	2/4/4/4	-
6	LDA	L	1284	-	-	8/13/13/13	-
4	GOL	H	1251	-	-	2/4/4/4	-
8	UQ2	L	1286[B]	-	-	8/15/39/39	0/1/1/1
6	LDA	M	1305	-	-	5/13/13/13	-
10	HTO	L	1289	-	-	6/10/10/10	-
5	BCL	L	1287	2	2/2/21/25	13/37/137/137	-
5	BCL	M	1303	3	2/2/21/25	16/37/137/137	-
12	U10	M	1314	-	-	9/45/69/87	0/1/1/1
6	LDA	M	1309	-	-	9/13/13/13	-
4	GOL	L	1290	-	-	2/4/4/4	-
14	CDL	M	1316	-	-	43/91/91/110	-
6	LDA	M	1308	-	-	5/13/13/13	-
13	SPO	M	1315	-	-	8/47/47/47	-
5	BCL	L	1282	2	2/2/21/25	9/37/137/137	-
6	LDA	M	1306	-	-	6/13/13/13	-
7	BPH	M	1313	-	-	16/37/105/105	0/5/6/6
10	HTO	L	1904	-	-	8/10/10/10	-
4	GOL	L	1292	-	-	4/4/4/4	-

The worst 5 of 123 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	L	1285	BPH	OBD-CAD	11.77	1.38	1.22
13	M	1315	SPO	C27-C28	11.70	1.46	1.34
7	M	1313	BPH	OBD-CAD	10.53	1.37	1.22
5	L	1287	BCL	OBD-CAD	10.45	1.40	1.22
5	M	1303	BCL	OBD-CAD	10.13	1.39	1.22

The worst 5 of 165 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	M	1316	CDL	C33-C32-C31	17.31	175.40	113.19
14	M	1316	CDL	C12-C11-CA5	10.70	152.53	113.62
14	M	1316	CDL	C35-C34-C33	8.43	157.24	114.42
7	L	1285	BPH	O2D-CGD-CBD	8.39	121.63	111.00
7	M	1313	BPH	O2D-CGD-CBD	7.41	120.38	111.00

5 of 8 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
5	L	1282	BCL	C13
5	L	1282	BCL	C8
5	L	1287	BCL	C13
5	L	1287	BCL	C8
5	M	1303	BCL	C13

5 of 231 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	1251	GOL	O1-C1-C2-C3
4	H	1252	GOL	C1-C2-C3-O3
4	H	1253	GOL	O1-C1-C2-C3
4	L	1290	GOL	C1-C2-C3-O3
4	L	1290	GOL	O2-C2-C3-O3

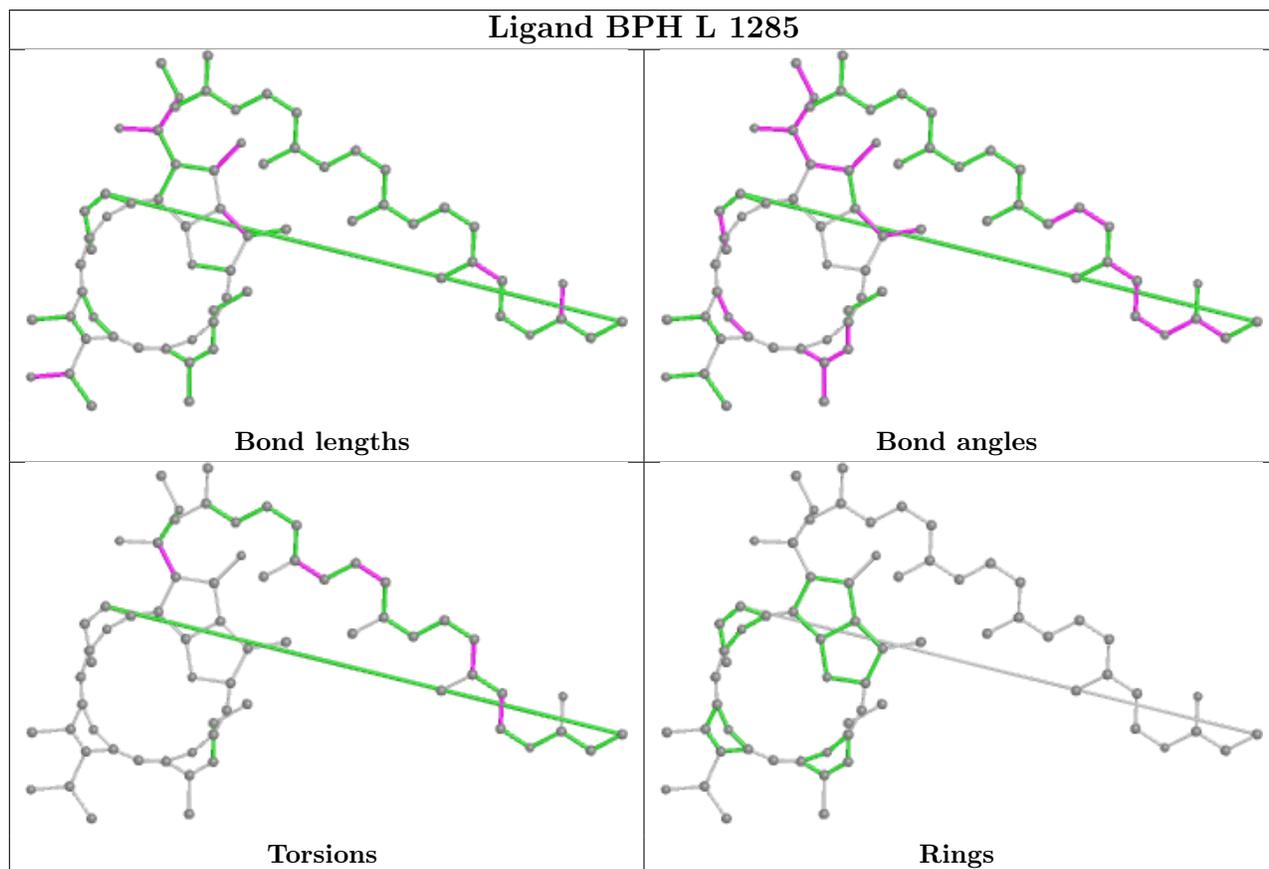
There are no ring outliers.

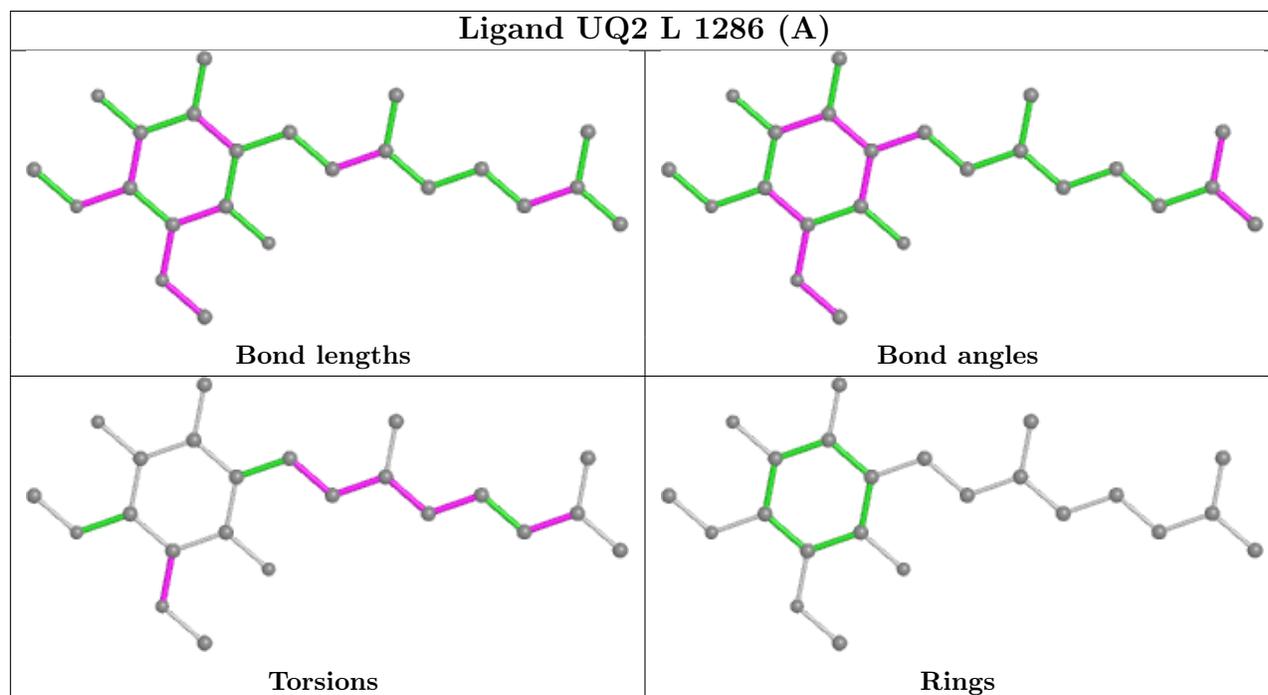
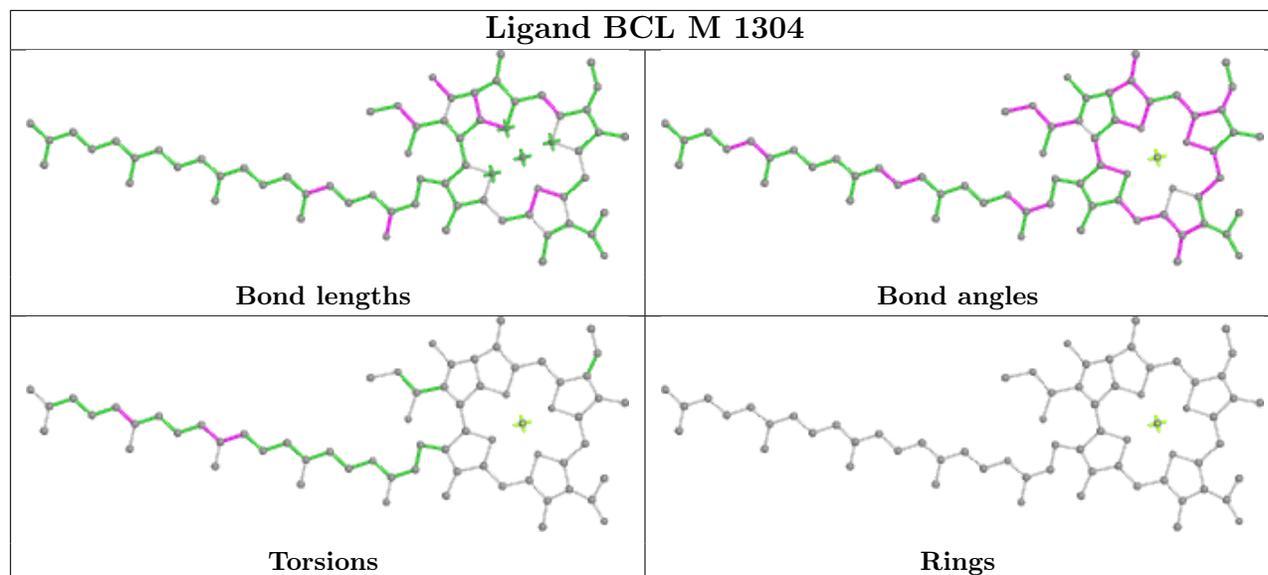
20 monomers are involved in 69 short contacts:

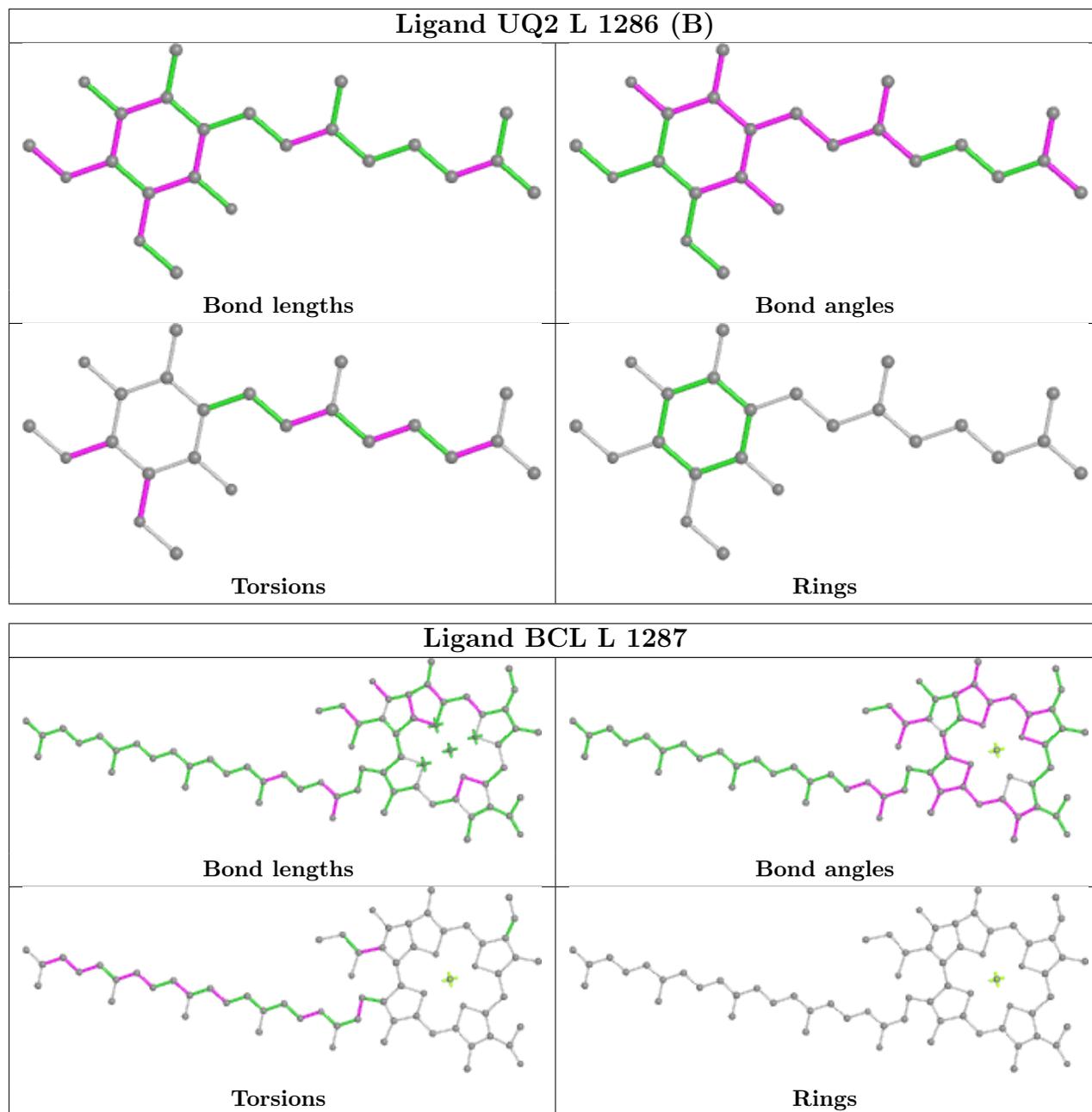
Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	L	1285	BPH	7	0
5	M	1304	BCL	11	0
8	L	1286[A]	UQ2	3	0
6	M	1307	LDA	5	0
6	M	1311	LDA	1	0
4	H	1251	GOL	1	0
8	L	1286[B]	UQ2	10	0
6	M	1305	LDA	2	0
10	L	1289	HTO	2	0
5	L	1287	BCL	4	0
5	M	1303	BCL	7	0
12	M	1314	U10	2	0
6	M	1309	LDA	1	0
4	L	1290	GOL	1	0
14	M	1316	CDL	1	0
6	M	1308	LDA	1	0
13	M	1315	SPO	4	0
5	L	1282	BCL	5	0
7	M	1313	BPH	9	0
4	L	1292	GOL	1	0

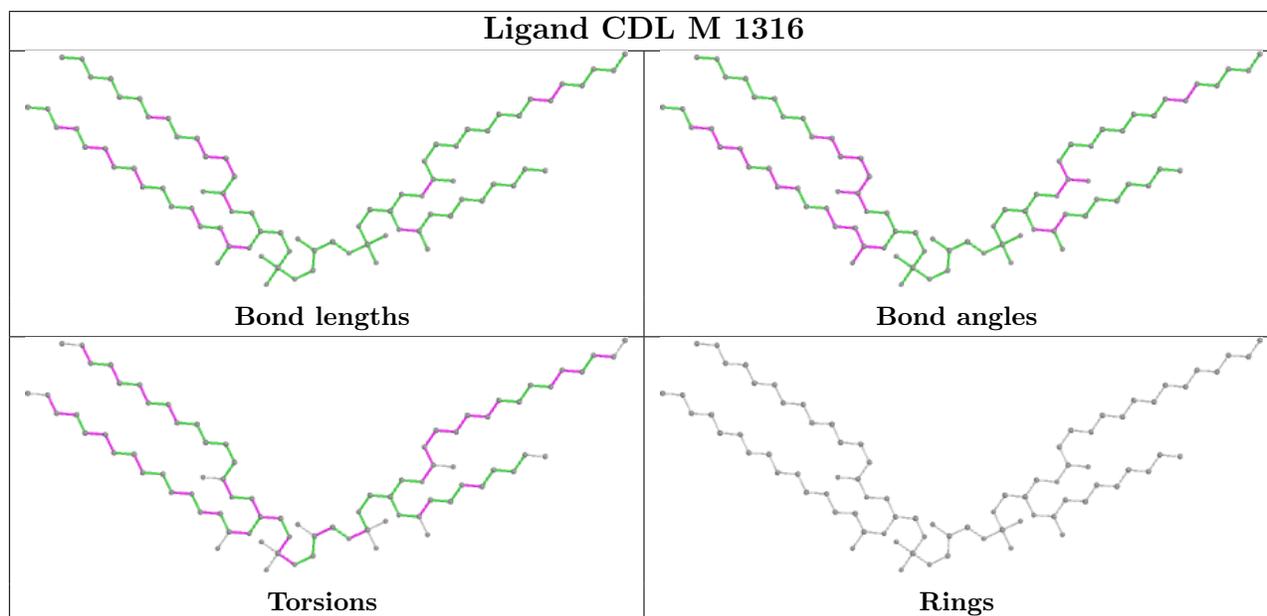
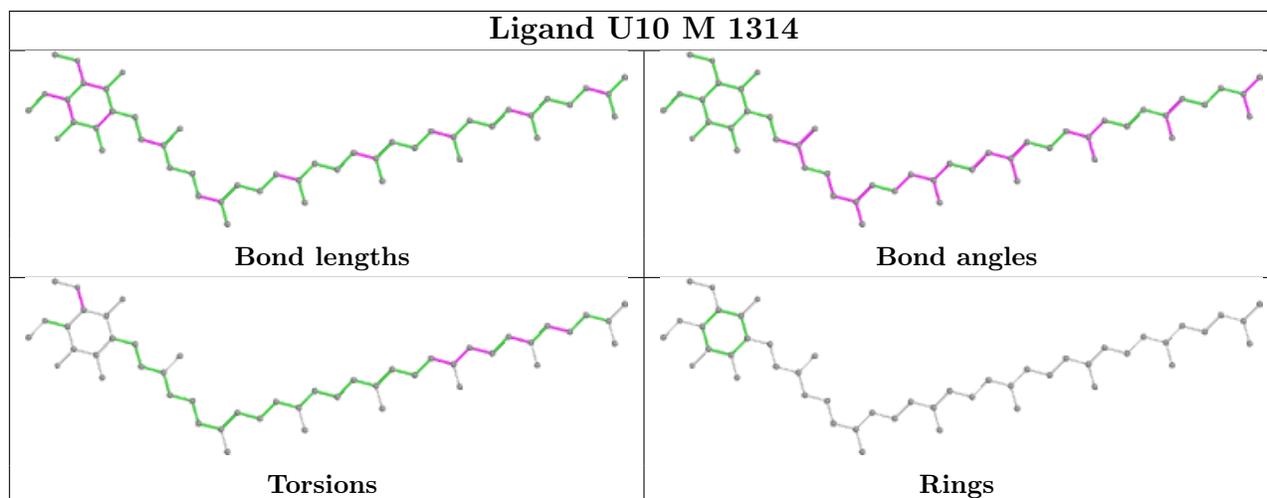
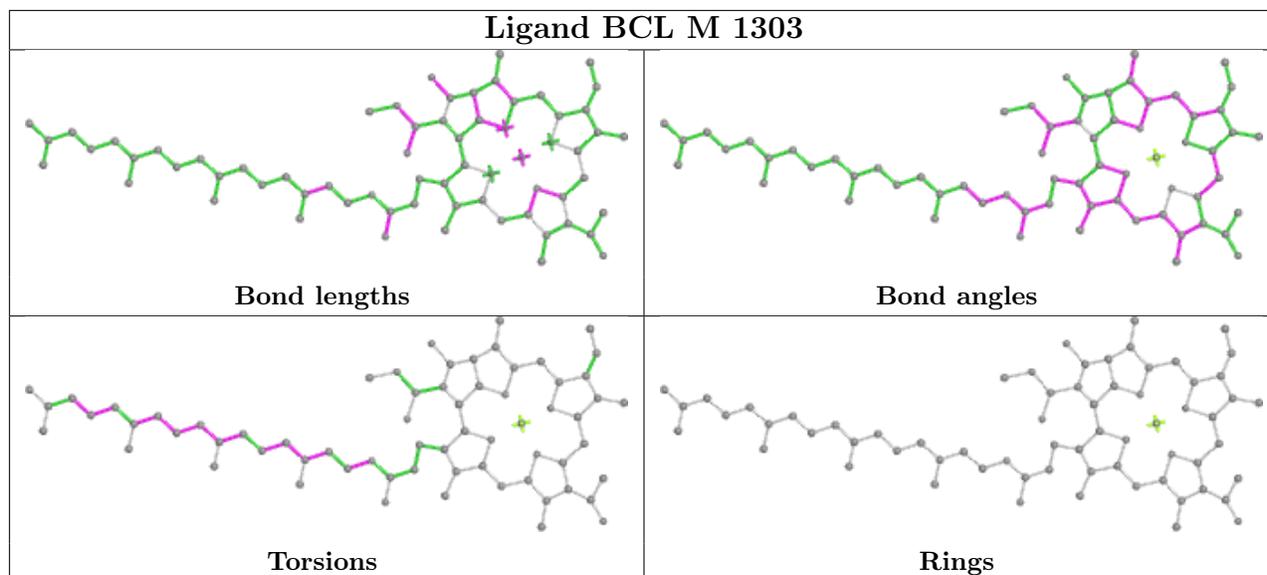
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

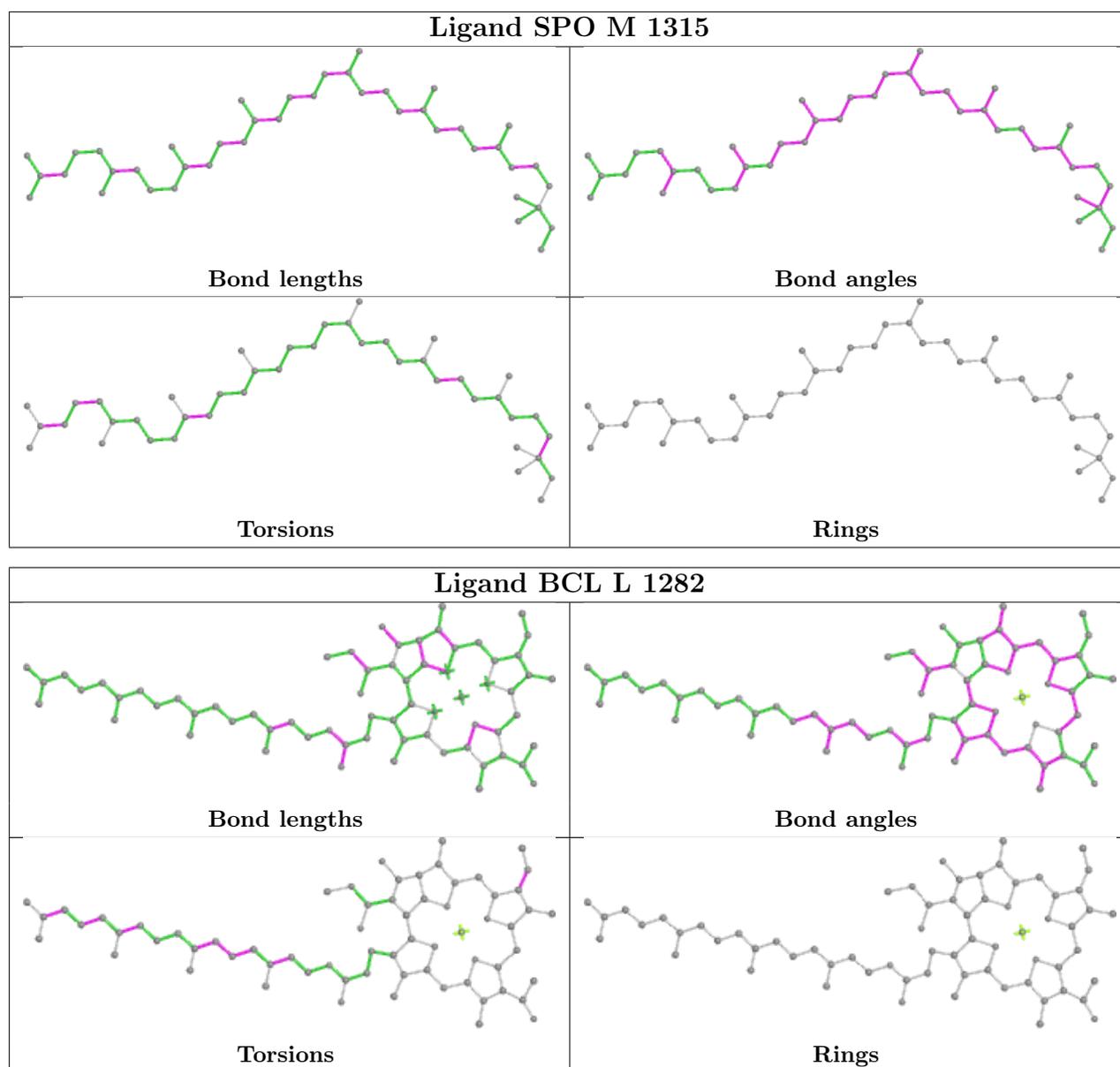
bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within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

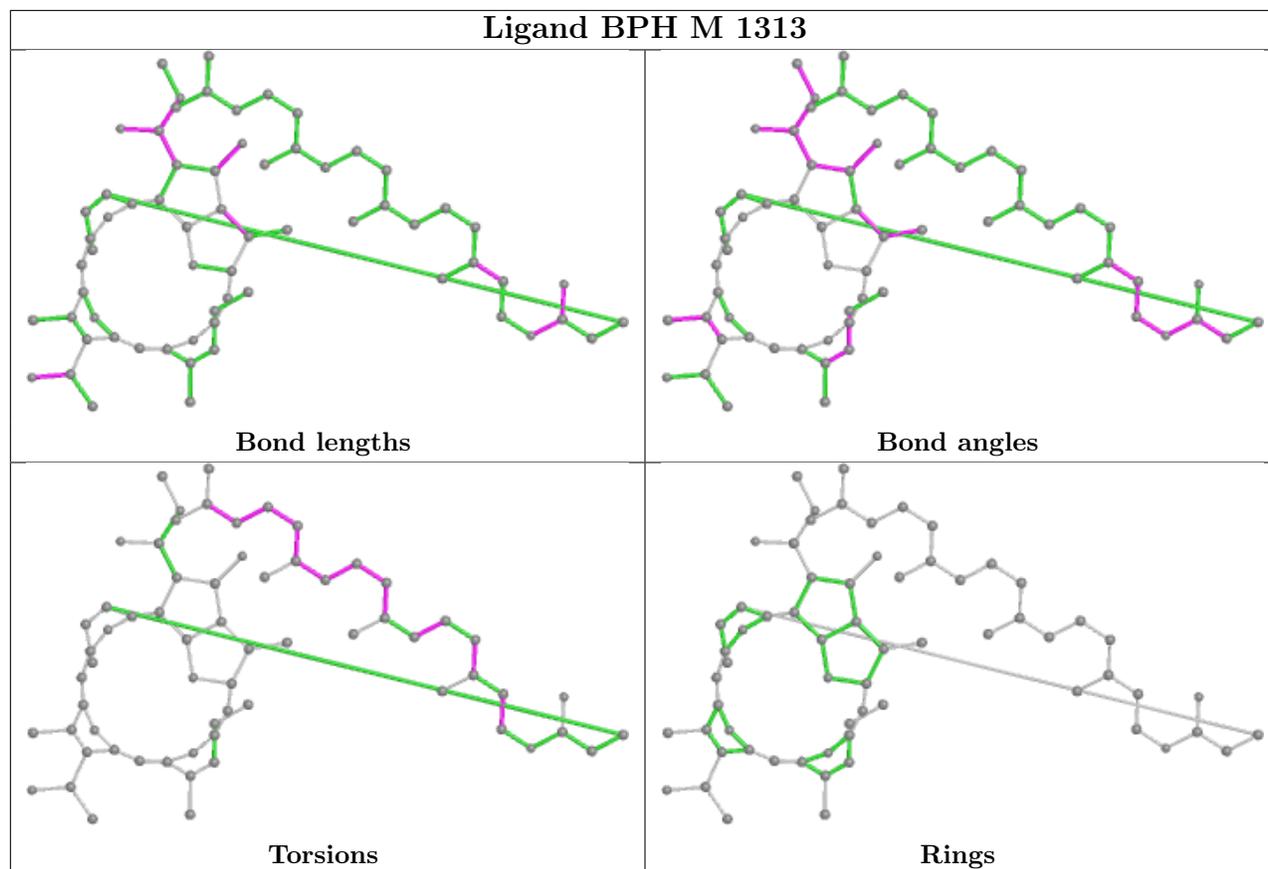












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	H	241/260 (92%)	0.31	20 (8%) 11 12	40, 51, 64, 110	0
2	L	281/281 (100%)	-0.40	4 (1%) 75 77	32, 42, 56, 65	0
3	M	303/307 (98%)	-0.20	8 (2%) 56 59	31, 43, 58, 88	0
All	All	825/848 (97%)	-0.12	32 (3%) 39 42	31, 45, 60, 110	0

The worst 5 of 32 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	251	VAL	14.3
1	H	250	SER	9.8
3	M	1	ALA	9.5
1	H	249	LYS	8.6
3	M	303	MET	8.3

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

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

### 6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

### 6.4 Ligands [i](#)

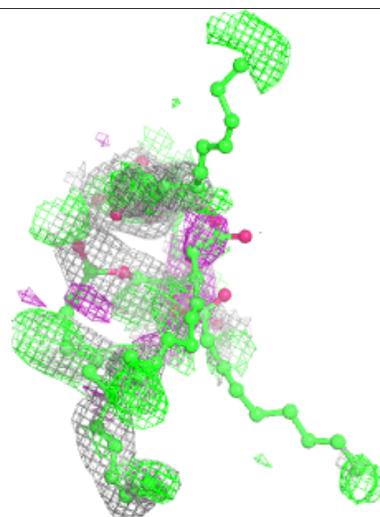
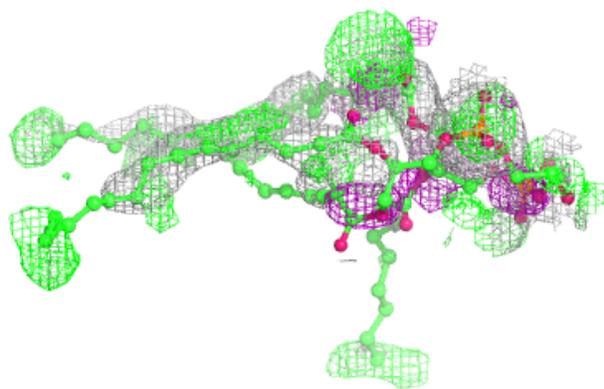
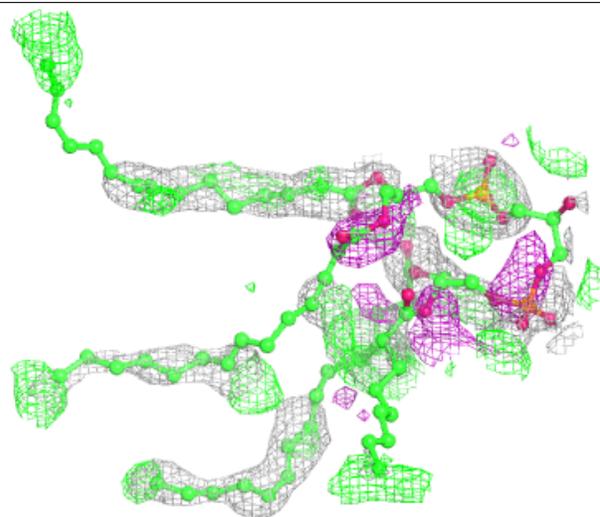
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
6	LDA	L	1284	16/16	0.08	0.43	100,108,118,118	0
6	LDA	M	1310	16/16	0.27	0.38	107,113,120,120	0
6	LDA	L	1283	16/16	0.33	0.39	76,102,115,115	0
6	LDA	M	1311	16/16	0.34	0.35	95,99,106,106	0
10	HTO	L	1904	10/10	0.37	0.40	107,110,111,112	0
14	CDL	M	1316	81/100	0.37	0.46	102,120,128,129	0
6	LDA	M	1308	16/16	0.38	0.35	124,125,128,128	0
4	GOL	L	1292	6/6	0.44	0.29	118,119,119,119	0
6	LDA	M	1309	16/16	0.61	0.29	103,105,111,112	0
4	GOL	H	1251	6/6	0.65	0.32	100,101,102,103	0
9	PO4	L	1288	5/5	0.65	0.32	166,166,167,167	0
6	LDA	M	1305	16/16	0.68	0.28	59,75,80,81	0
10	HTO	L	1289	10/10	0.75	0.29	70,73,73,74	0
6	LDA	M	1306	16/16	0.75	0.32	86,90,94,94	0
4	GOL	H	1253	6/6	0.75	0.26	119,120,120,120	0
4	GOL	H	1252	6/6	0.78	0.30	92,93,94,94	0
6	LDA	M	1307	16/16	0.81	0.21	75,76,81,82	0
13	SPO	M	1315	42/42	0.85	0.20	35,44,66,71	0
4	GOL	L	1291	6/6	0.85	0.18	80,81,82,83	0
4	GOL	L	1290	6/6	0.86	0.19	66,70,72,73	0
8	UQ2	L	1286[B]	23/23	0.90	0.27	24,31,41,42	23
8	UQ2	L	1286[A]	23/23	0.90	0.27	24,38,57,59	23
7	BPH	M	1313	65/65	0.91	0.16	35,41,98,101	0
12	U10	M	1314	48/63	0.92	0.15	36,50,76,77	0
5	BCL	M	1303	66/66	0.95	0.13	28,34,79,80	0
5	BCL	M	1304	66/66	0.96	0.16	25,34,58,64	0
5	BCL	L	1287	66/66	0.96	0.16	27,34,54,64	0
7	BPH	L	1285	65/65	0.96	0.15	25,37,46,48	0
5	BCL	L	1282	66/66	0.97	0.17	31,35,62,67	0
11	FE	M	1312	1/1	1.00	0.03	35,35,35,35	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.

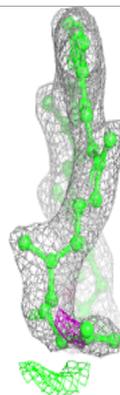
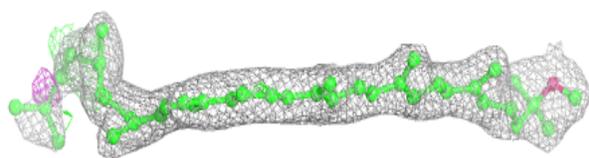
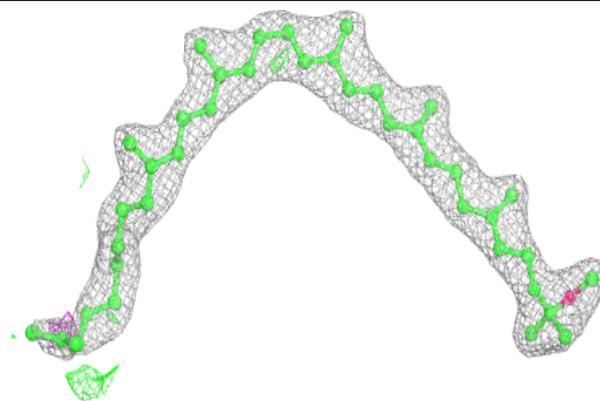
**Electron density around CDL M 1316:**

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

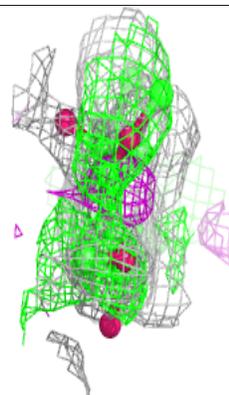
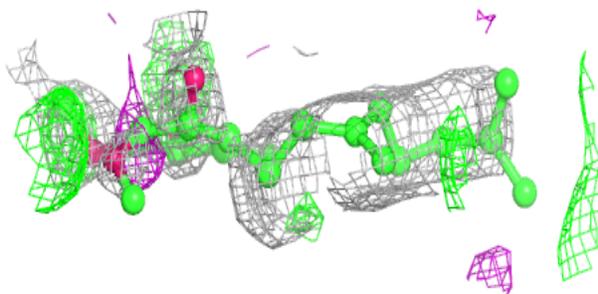
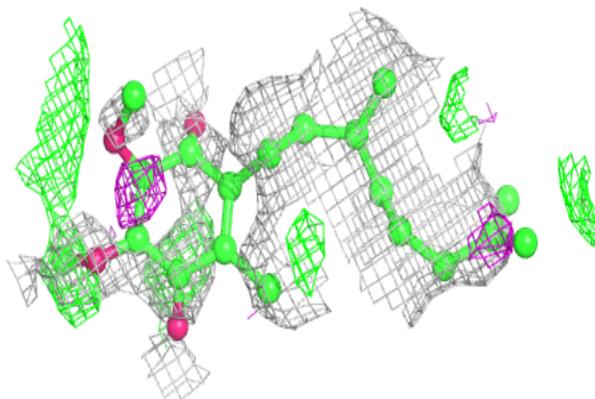


**Electron density around SPO M 1315:**

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

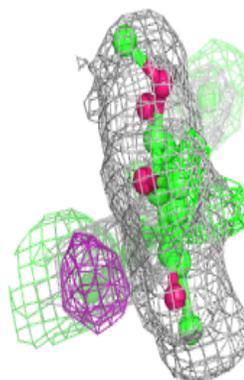
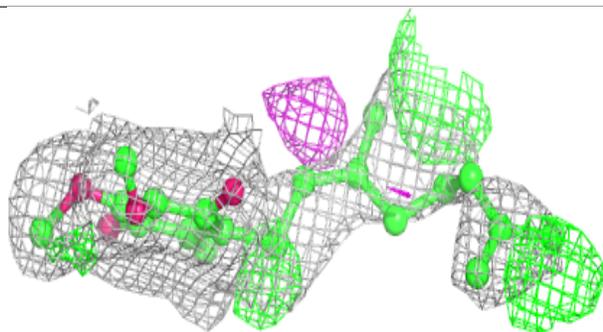
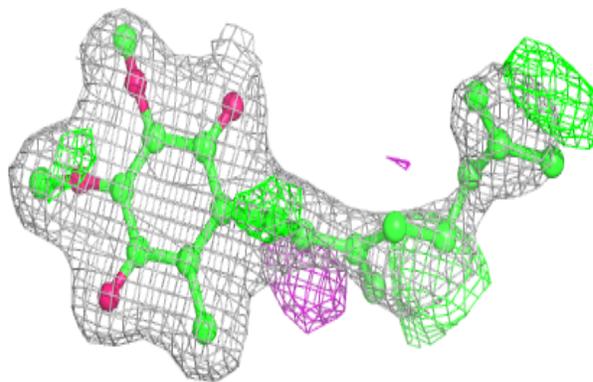
**Electron density around UQ2 L 1286 (B):**

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

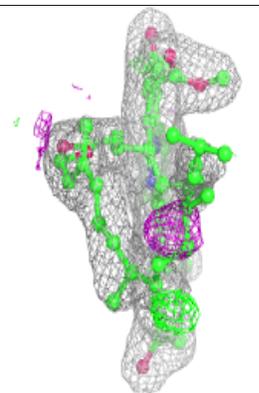
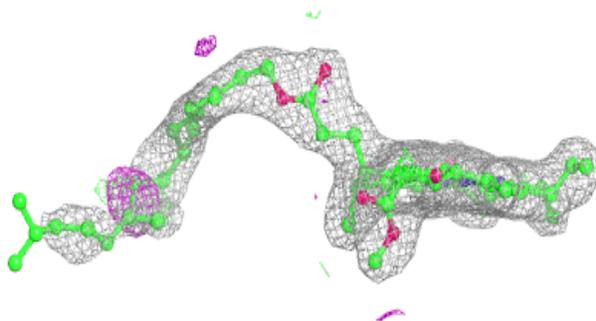
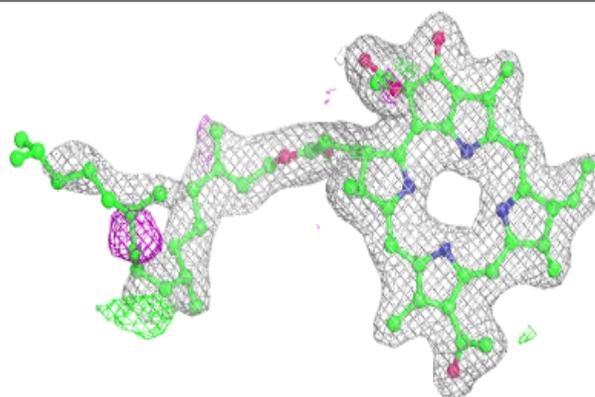


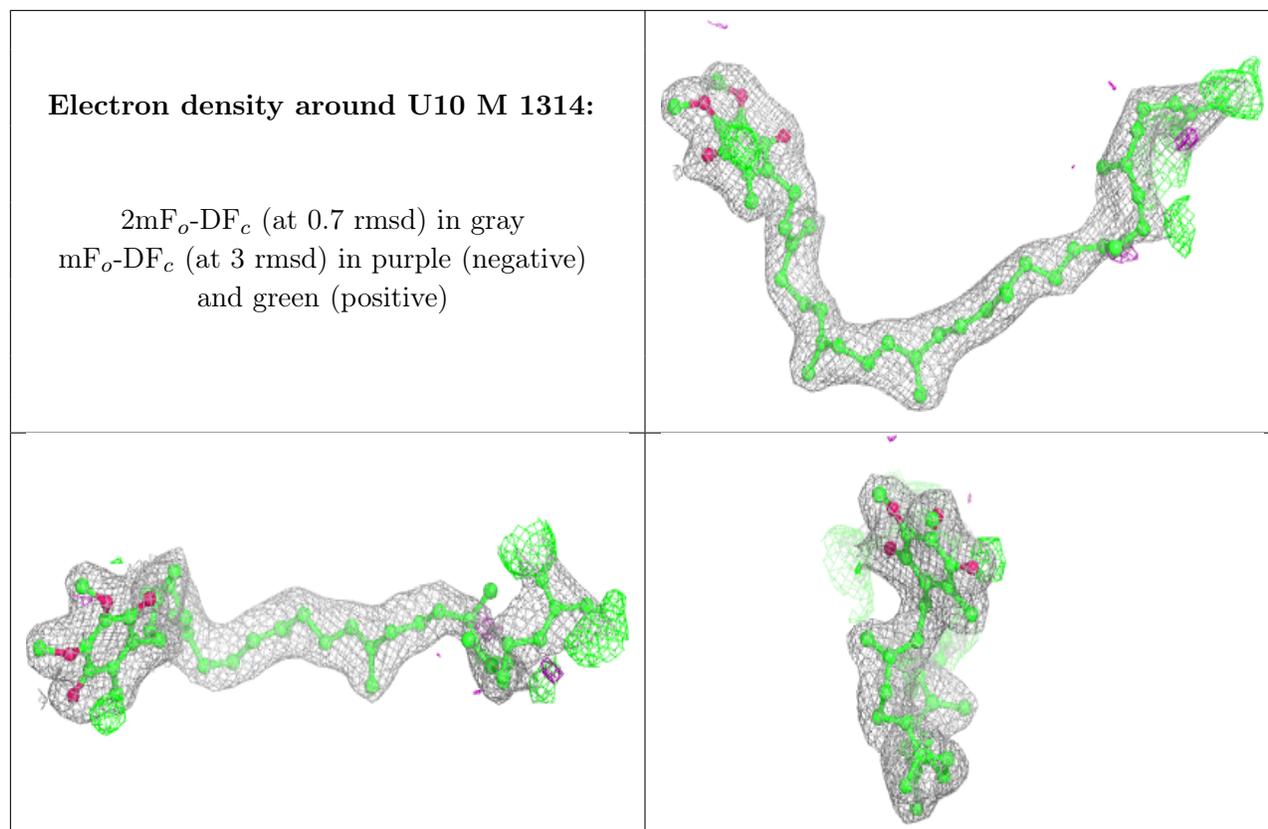
**Electron density around UQ2 L 1286 (A):**

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

**Electron density around BPH M 1313:**

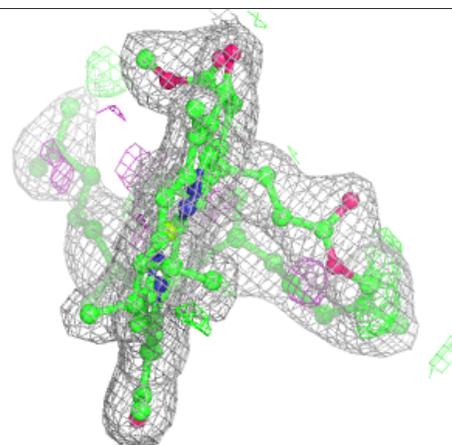
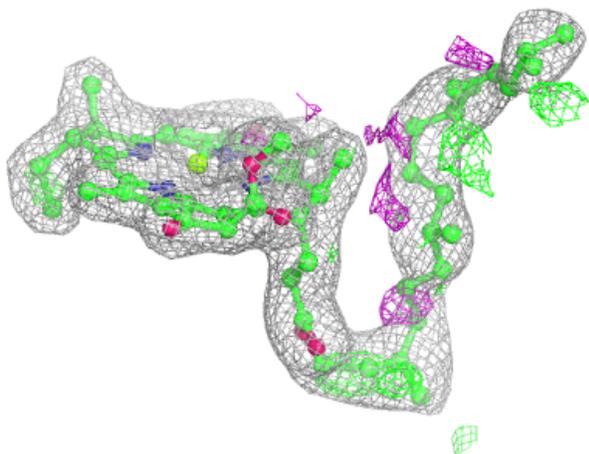
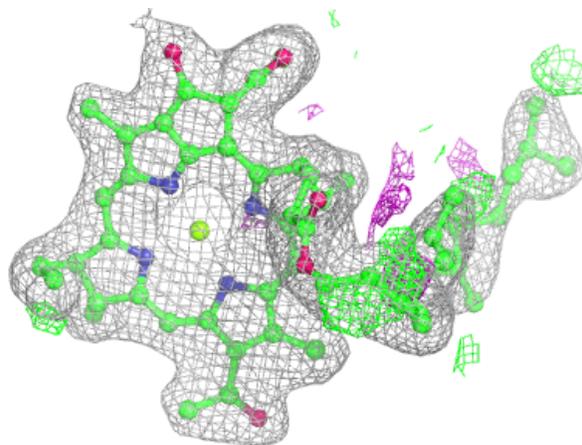
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





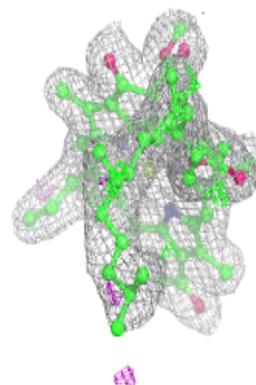
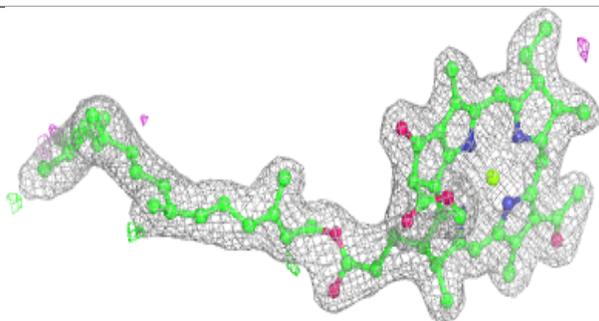
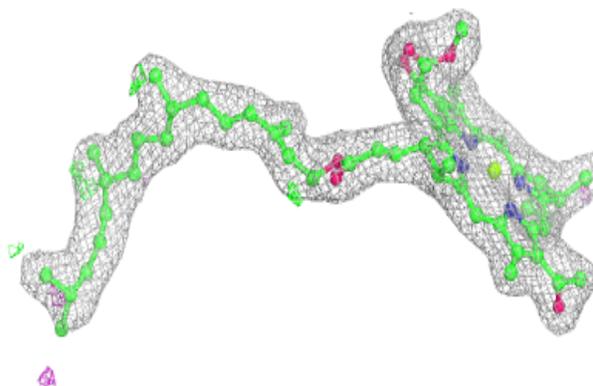
**Electron density around BCL M 1303:**

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

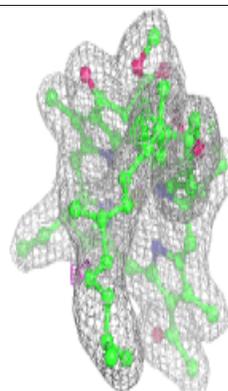
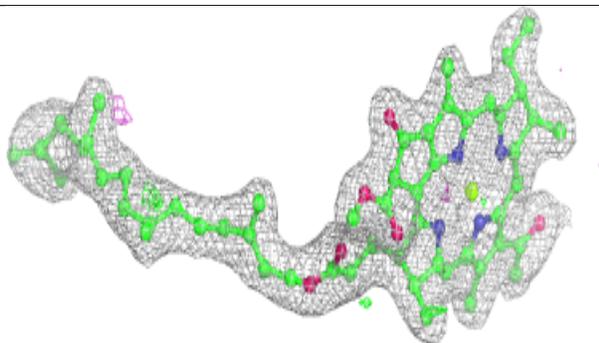
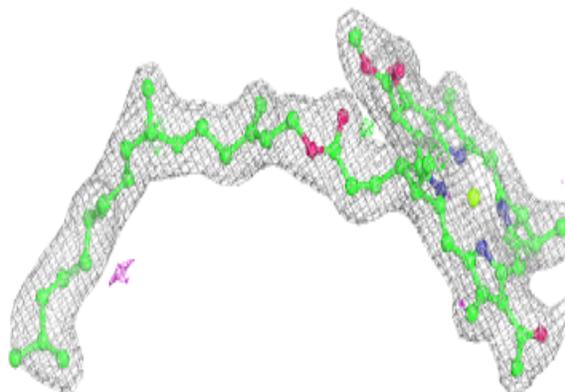


**Electron density around BCL M 1304:**

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and green (positive)

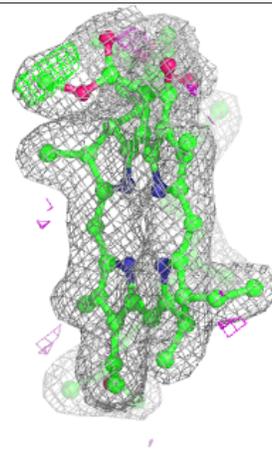
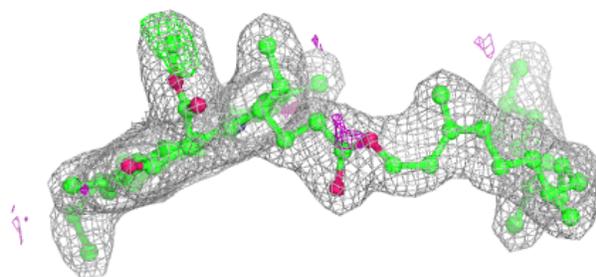
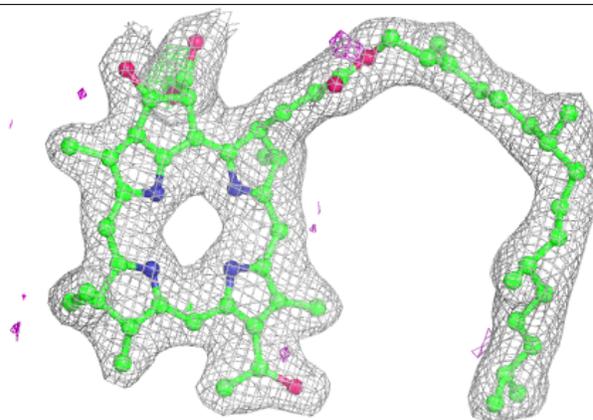
**Electron density around BCL L 1287:**

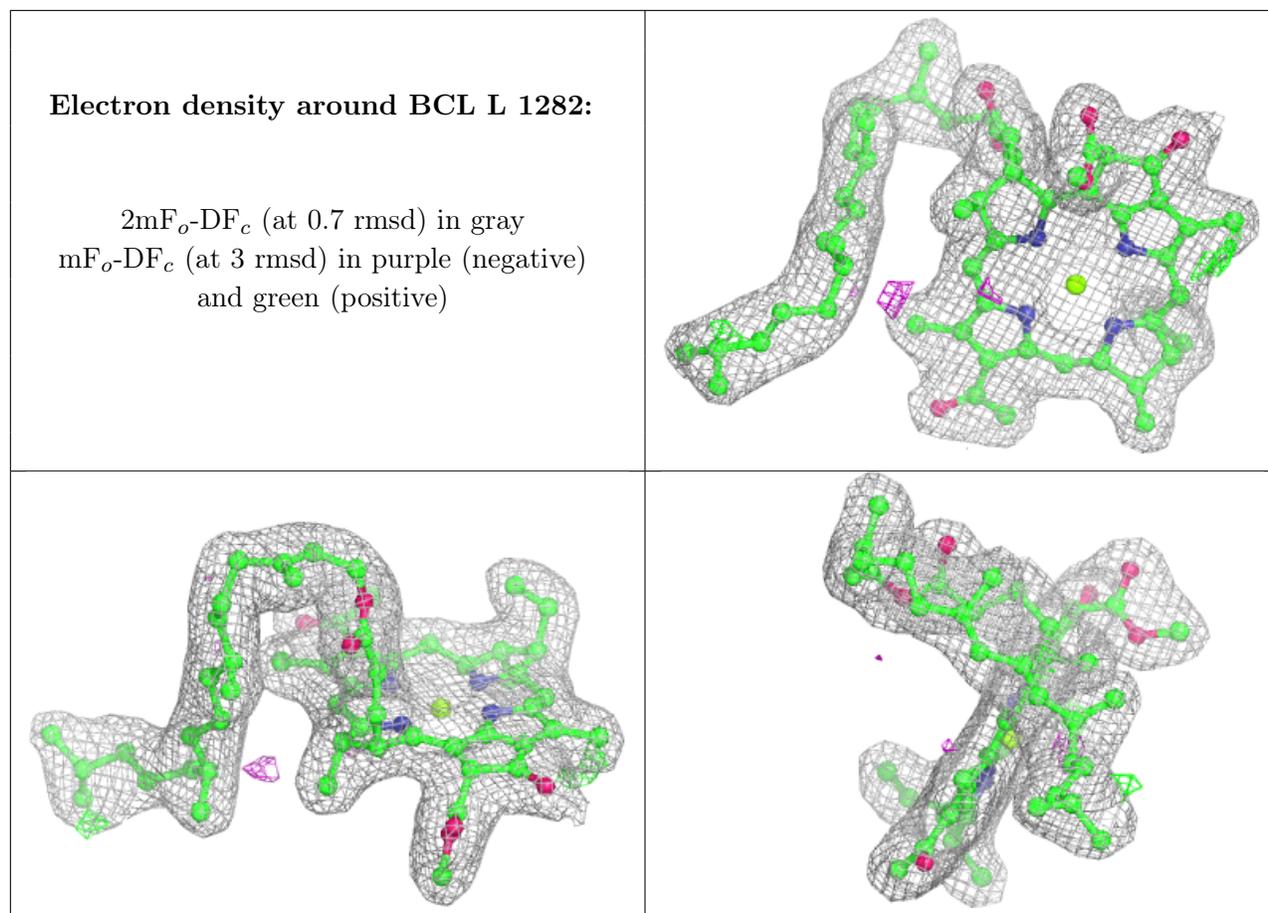
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around BPH L 1285:**

$2mF_o-DF_c$  (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.