



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 3, 2022 – 12:19 pm GMT

PDB ID : 7O8N  
Title : NmHR light state structure at 7.5 ms (5 - 10 ms) after photoexcitation determined by serial millisecond crystallography  
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Deposited on : 2021-04-15  
Resolution : 2.10 Å(reported)

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

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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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 : 2.26  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0267  
CCP4 : 7.1.010 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.26

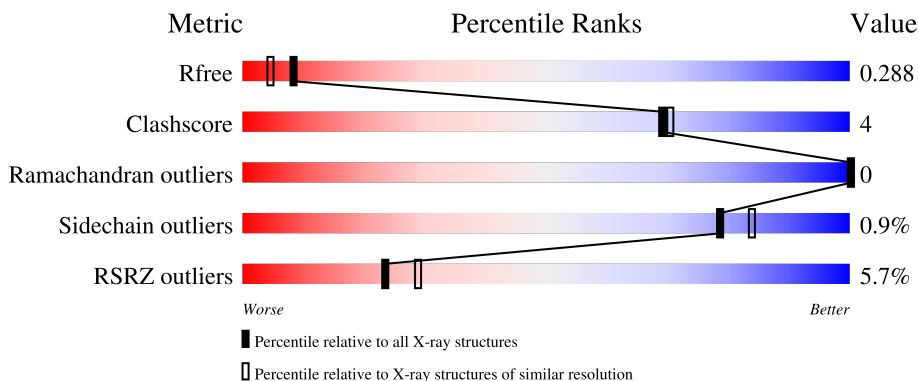
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.10 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	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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

Mol	Chain	Length	Quality of chain
1	A	296	

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
4	OLA	A	308	-	-	-	X

## 2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2312 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.

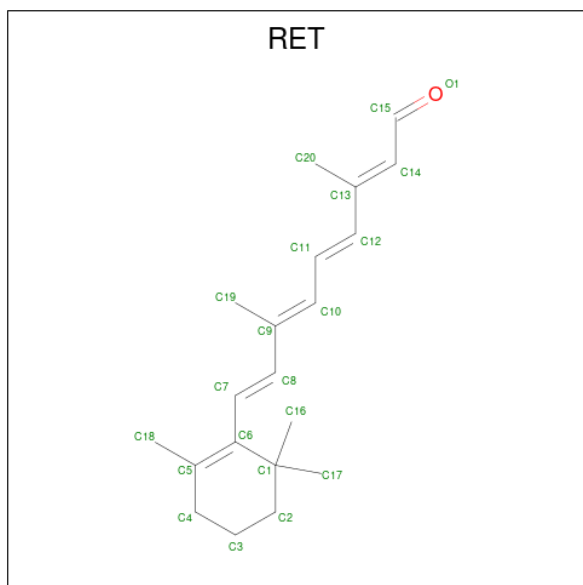
- Molecule 1 is a protein called Chloride pumping rhodopsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	265	2083	1386	320	362	15	0	1	0

There are 24 discrepancies between the modelled and reference sequences:

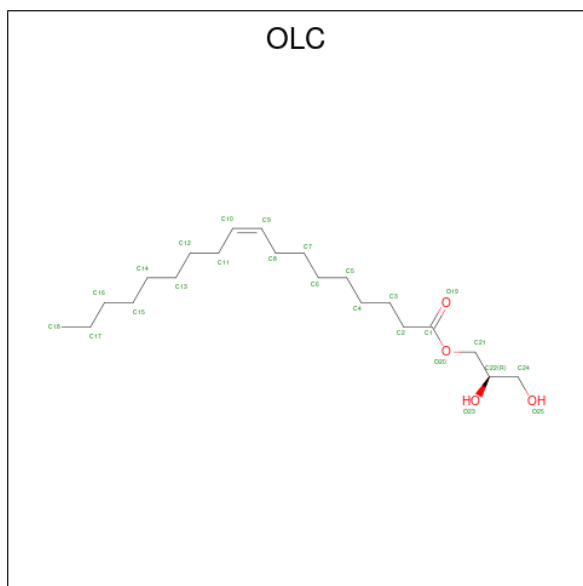
Chain	Residue	Modelled	Actual	Comment	Reference
A	-15	MET	-	initiating methionine	UNP W8VZW3
A	-14	ALA	-	expression tag	UNP W8VZW3
A	-13	SER	-	expression tag	UNP W8VZW3
A	-12	MET	-	expression tag	UNP W8VZW3
A	-11	THR	-	expression tag	UNP W8VZW3
A	-10	GLY	-	expression tag	UNP W8VZW3
A	-9	GLY	-	expression tag	UNP W8VZW3
A	-8	GLN	-	expression tag	UNP W8VZW3
A	-7	GLN	-	expression tag	UNP W8VZW3
A	-6	MET	-	expression tag	UNP W8VZW3
A	-5	GLY	-	expression tag	UNP W8VZW3
A	-4	ARG	-	expression tag	UNP W8VZW3
A	-3	ASP	-	expression tag	UNP W8VZW3
A	-2	PRO	-	expression tag	UNP W8VZW3
A	-1	ASN	-	expression tag	UNP W8VZW3
A	0	SER	-	expression tag	UNP W8VZW3
A	273	LEU	-	expression tag	UNP W8VZW3
A	274	GLU	-	expression tag	UNP W8VZW3
A	275	HIS	-	expression tag	UNP W8VZW3
A	276	HIS	-	expression tag	UNP W8VZW3
A	277	HIS	-	expression tag	UNP W8VZW3
A	278	HIS	-	expression tag	UNP W8VZW3
A	279	HIS	-	expression tag	UNP W8VZW3
A	280	HIS	-	expression tag	UNP W8VZW3

- Molecule 2 is RETINAL (three-letter code: RET) (formula: C<sub>20</sub>H<sub>28</sub>O).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C 20 20	0	0

- Molecule 3 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (three-letter code: OLC) (formula:  $C_{21}H_{40}O_4$ ).



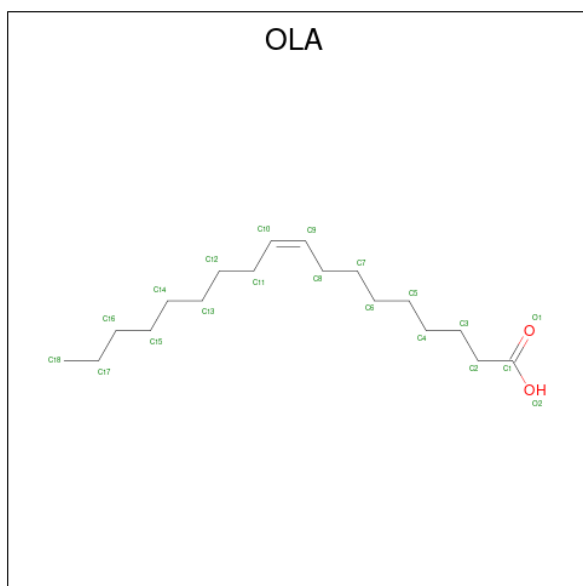
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 25 21 4	0	0
3	A	1	Total C O 16 12 4	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf		
3	A	1	Total	C	O	0	0
			10	6	4		

- Molecule 4 is OLEIC ACID (three-letter code: OLA) (formula:  $C_{18}H_{34}O_2$ ).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf		
4	A	1	Total	C	O	0	0
			20	18	2		
4	A	1	Total	C	O	0	0
			20	18	2		
4	A	1	Total	C	O	0	0
			17	15	2		
4	A	1	Total	C	O	0	0
			20	18	2		
4	A	1	Total	C	O	0	0
			14	12	2		
4	A	1	Total	C		0	0
			17	17			

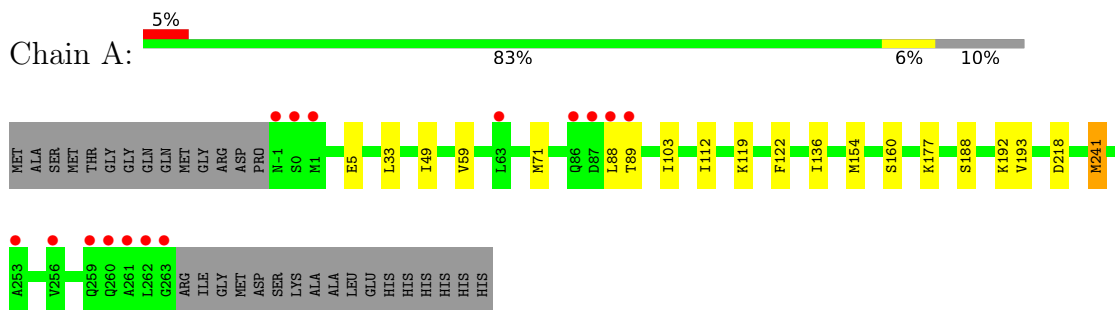
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf		
5	A	50	Total	O		0	0
			50	50			

### 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: Chloride pumping rhodopsin



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	104.48Å 51.18Å 78.33Å 90.00° 131.78° 90.00°	Depositor
Resolution (Å)	38.96 – 2.10 38.96 – 2.10	Depositor EDS
% Data completeness (in resolution range)	95.5 (38.96-2.10) 95.5 (38.96-2.10)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.50 (at 2.10Å)	Xtrriage
Refinement program	PHENIX 1.19.1_4122	Depositor
R, $R_{free}$	0.241 , 0.282 0.252 , 0.288	Depositor DCC
$R_{free}$ test set	860 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	36.4	Xtrriage
Anisotropy	0.225	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	(Not available) , (Not available)	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.52$ , $\langle L^2 \rangle = 0.36$	Xtrriage
Estimated twinning fraction	0.001 for h,-k,-h-l	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	2312	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	44.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 12.23% 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: OLC, OLA, RET

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.27	0/2139	0.41	0/2916

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2083	0	2113	15	1
2	A	20	0	27	4	0
3	A	51	0	70	0	0
4	A	108	0	169	2	0
5	A	50	0	0	0	0
All	All	2312	0	2379	17	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

All (17) 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)
2:A:301:RET:H8	2:A:301:RET:H161	1.62	0.81
1:A:103:ILE:HD12	1:A:136:ILE:HD11	1.67	0.75
2:A:301:RET:H161	2:A:301:RET:C8	2.25	0.67
1:A:103:ILE:HG12	2:A:301:RET:H202	1.79	0.65
1:A:49:ILE:HG23	1:A:112:ILE:HD13	1.80	0.62
1:A:188:SER:O	1:A:192:LYS:HE3	2.07	0.55
1:A:33:LEU:HD22	1:A:59:VAL:HG22	1.88	0.55
1:A:88:LEU:HD22	1:A:88:LEU:N	2.25	0.51
1:A:122[B]:PHE:CD1	4:A:304:OLA:H10	2.47	0.49
1:A:119:LYS:HB2	1:A:119:LYS:NZ	2.29	0.48
1:A:160:SER:OG	2:A:301:RET:H41	2.14	0.47
1:A:71:MET:HG2	1:A:89:THR:HG22	1.98	0.45
1:A:5:GLU:OE1	1:A:5:GLU:N	2.46	0.44
1:A:122[B]:PHE:HD1	4:A:304:OLA:H10	1.82	0.44
1:A:188:SER:OG	1:A:192:LYS:NZ	2.46	0.44
1:A:193:VAL:HG22	1:A:241:MET:CE	2.51	0.41
1:A:193:VAL:HG22	1:A:241:MET:HE2	2.02	0.41

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:177:LYS:NZ	1:A:218:ASP:OD2[4_445]	2.18	0.02

## 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	A	264/296 (89%)	262 (99%)	2 (1%)	0	<b>100</b> <b>100</b>

There are no Ramachandran outliers to report.

### 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	A	219/242 (90%)	217 (99%)	2 (1%)	<a href="#">78</a> <a href="#">84</a>

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

Mol	Chain	Res	Type
1	A	154	MET
1	A	241	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	-1	ASN

### 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](#)

10 ligands 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
2	RET	A	301	1	20,20,21	0.63	0	27,27,28	1.95	5 (18%)
4	OLA	A	309	-	10,13,19	0.91	1 (10%)	8,13,19	0.64	0
3	OLC	A	307	-	9,9,24	1.22	1 (11%)	10,10,25	1.27	1 (10%)
4	OLA	A	304	-	16,19,19	0.74	1 (6%)	15,19,19	0.57	0
3	OLC	A	303	-	15,15,24	0.99	1 (6%)	16,16,25	0.99	1 (6%)
3	OLC	A	302	-	24,24,24	0.81	1 (4%)	25,25,25	0.99	1 (4%)
4	OLA	A	310	-	16,16,19	0.91	1 (6%)	15,15,19	0.76	0
4	OLA	A	306	-	13,16,19	0.78	1 (7%)	12,16,19	0.61	0
4	OLA	A	305	-	16,19,19	0.73	1 (6%)	15,19,19	0.58	0
4	OLA	A	308	-	16,19,19	0.70	1 (6%)	15,19,19	0.60	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
2	RET	A	301	1	-	0/13/30/31	0/1/1/1
4	OLA	A	309	-	-	5/9/11/17	-
3	OLC	A	307	-	-	8/9/9/24	-
4	OLA	A	304	-	-	9/15/17/17	-
3	OLC	A	303	-	-	9/15/15/24	-
3	OLC	A	302	-	-	9/24/24/24	-
4	OLA	A	310	-	-	7/14/14/17	-
4	OLA	A	306	-	-	9/12/14/17	-
4	OLA	A	305	-	-	9/15/17/17	-
4	OLA	A	308	-	-	8/15/17/17	-

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	310	OLA	C10-C9	3.43	1.51	1.31
4	A	304	OLA	C10-C9	2.76	1.47	1.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	309	OLA	C10-C9	2.73	1.47	1.31
4	A	305	OLA	C10-C9	2.73	1.47	1.31
4	A	306	OLA	C10-C9	2.68	1.47	1.31
4	A	308	OLA	C10-C9	2.66	1.47	1.31
3	A	307	OLC	O20-C1	2.56	1.40	1.33
3	A	303	OLC	O20-C1	2.54	1.40	1.33
3	A	302	OLC	O20-C1	2.54	1.40	1.33

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	301	RET	C20-C13-C14	-5.16	108.48	123.71
2	A	301	RET	C3-C4-C5	-4.70	105.68	114.08
2	A	301	RET	C12-C13-C14	4.50	133.05	118.80
3	A	307	OLC	O20-C1-C2	3.46	120.46	111.38
3	A	302	OLC	O20-C1-C2	3.10	121.62	111.91
3	A	303	OLC	O20-C1-C2	2.79	120.66	111.91
2	A	301	RET	C7-C8-C9	-2.63	122.26	126.23
2	A	301	RET	C18-C5-C6	-2.33	121.91	124.53

There are no chirality outliers.

All (73) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	303	OLC	C21-C22-C24-O25
3	A	307	OLC	O20-C21-C22-C24
3	A	307	OLC	O20-C21-C22-O23
4	A	304	OLA	C6-C7-C8-C9
4	A	305	OLA	C1-C2-C3-C4
4	A	306	OLA	C1-C2-C3-C4
4	A	308	OLA	C1-C2-C3-C4
3	A	303	OLC	O20-C21-C22-C24
3	A	303	OLC	C4-C5-C6-C7
4	A	304	OLA	C3-C4-C5-C6
4	A	304	OLA	C2-C3-C4-C5
4	A	304	OLA	C11-C12-C13-C14
4	A	310	OLA	C3-C4-C5-C6
3	A	302	OLC	C2-C1-O20-C21
3	A	302	OLC	C13-C14-C15-C16
4	A	310	OLA	C2-C3-C4-C5
4	A	309	OLA	C3-C4-C5-C6
3	A	303	OLC	O20-C21-C22-O23

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Mol	Chain	Res	Type	Atoms
3	A	302	OLC	C6-C7-C8-C9
3	A	303	OLC	C1-C2-C3-C4
4	A	308	OLA	C13-C14-C15-C16
3	A	302	OLC	C5-C6-C7-C8
4	A	305	OLA	C4-C5-C6-C7
4	A	308	OLA	C3-C4-C5-C6
3	A	303	OLC	O23-C22-C24-O25
4	A	306	OLA	C10-C11-C12-C13
4	A	306	OLA	C2-C3-C4-C5
3	A	302	OLC	O19-C1-O20-C21
4	A	310	OLA	C11-C12-C13-C14
3	A	303	OLC	C6-C7-C8-C9
4	A	306	OLA	C6-C7-C8-C9
4	A	308	OLA	C6-C7-C8-C9
4	A	310	OLA	C10-C11-C12-C13
4	A	305	OLA	C10-C11-C12-C13
3	A	307	OLC	C2-C1-O20-C21
4	A	306	OLA	C11-C10-C9-C8
4	A	308	OLA	C11-C10-C9-C8
3	A	302	OLC	C4-C5-C6-C7
4	A	310	OLA	C13-C14-C15-C16
4	A	308	OLA	C15-C16-C17-C18
4	A	305	OLA	C11-C10-C9-C8
3	A	307	OLC	O19-C1-O20-C21
4	A	308	OLA	C12-C13-C14-C15
3	A	302	OLC	C15-C16-C17-C18
4	A	308	OLA	C5-C6-C7-C8
4	A	304	OLA	C5-C6-C7-C8
4	A	304	OLA	C4-C5-C6-C7
4	A	309	OLA	C5-C6-C7-C8
4	A	305	OLA	C3-C4-C5-C6
4	A	305	OLA	C14-C15-C16-C17
4	A	306	OLA	C4-C5-C6-C7
4	A	305	OLA	C11-C12-C13-C14
4	A	306	OLA	C11-C12-C13-C14
4	A	310	OLA	C6-C7-C8-C9
4	A	309	OLA	C1-C2-C3-C4
4	A	305	OLA	C12-C13-C14-C15
4	A	309	OLA	C4-C5-C6-C7
4	A	309	OLA	C7-C8-C9-C10
3	A	307	OLC	O20-C1-C2-C3
3	A	303	OLC	C2-C3-C4-C5

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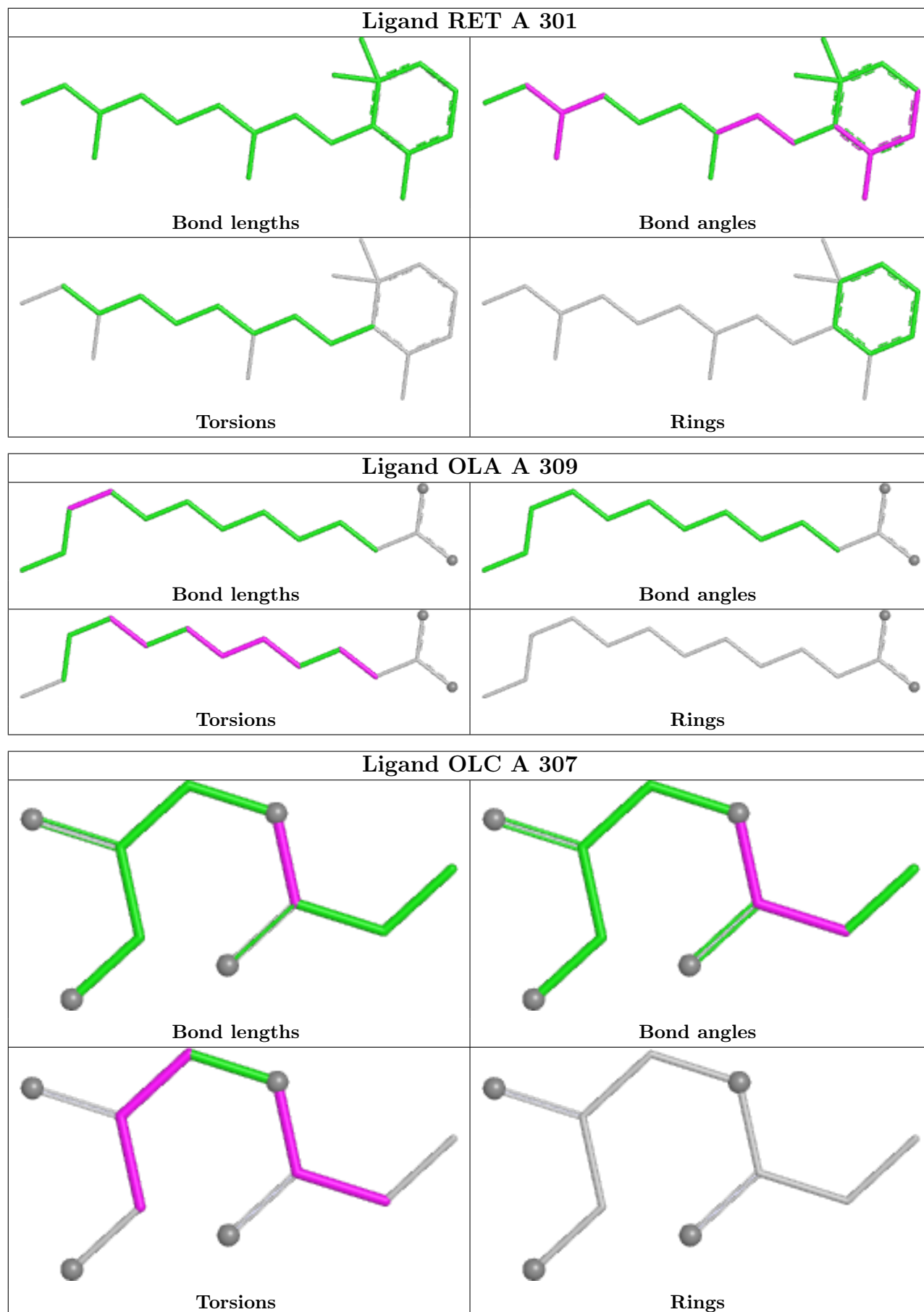
Mol	Chain	Res	Type	Atoms
3	A	307	OLC	O19-C1-C2-C3
3	A	302	OLC	C21-C22-C24-O25
3	A	307	OLC	C21-C22-C24-O25
3	A	302	OLC	O23-C22-C24-O25
3	A	307	OLC	O23-C22-C24-O25
4	A	304	OLA	C7-C8-C9-C10
4	A	306	OLA	C9-C10-C11-C12
4	A	304	OLA	C11-C10-C9-C8
4	A	305	OLA	C9-C10-C11-C12
4	A	306	OLA	C7-C8-C9-C10
4	A	304	OLA	C9-C10-C11-C12
4	A	310	OLA	C7-C8-C9-C10
3	A	303	OLC	O19-C1-O20-C21

There are no ring outliers.

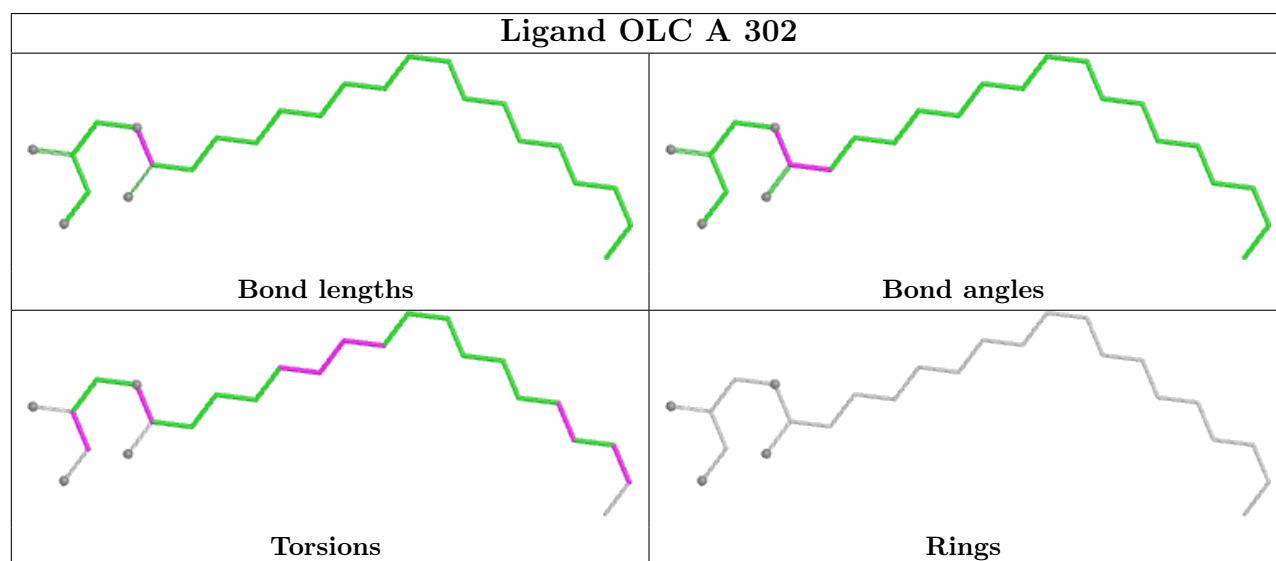
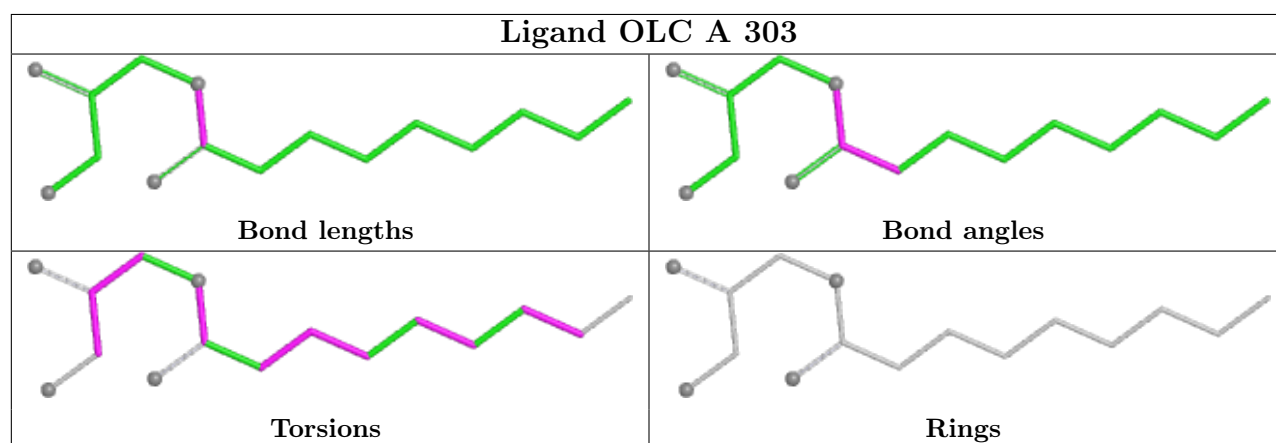
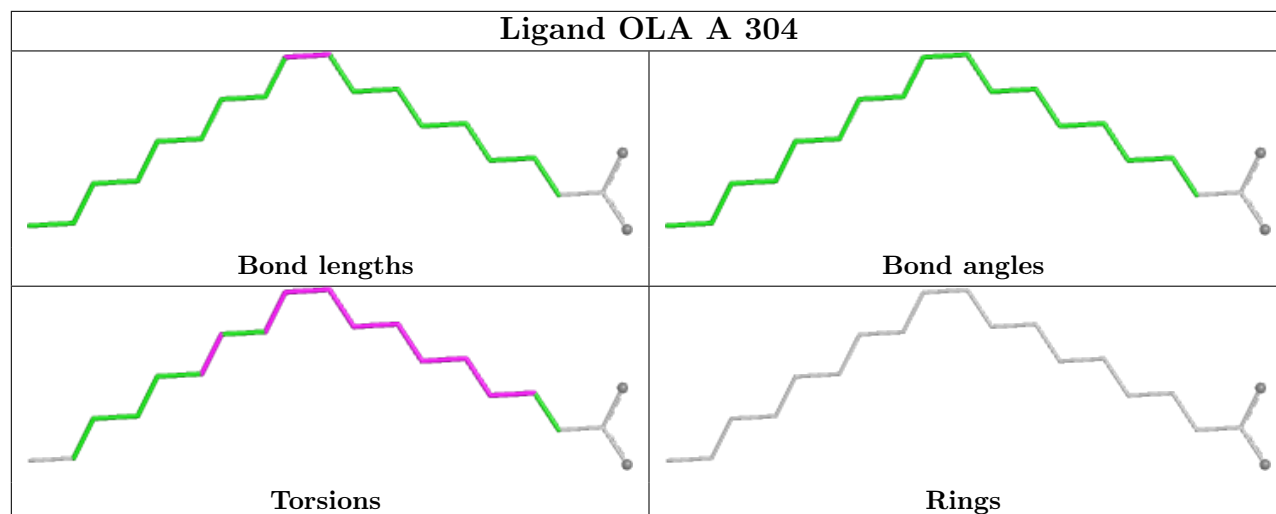
2 monomers are involved in 6 short contacts:

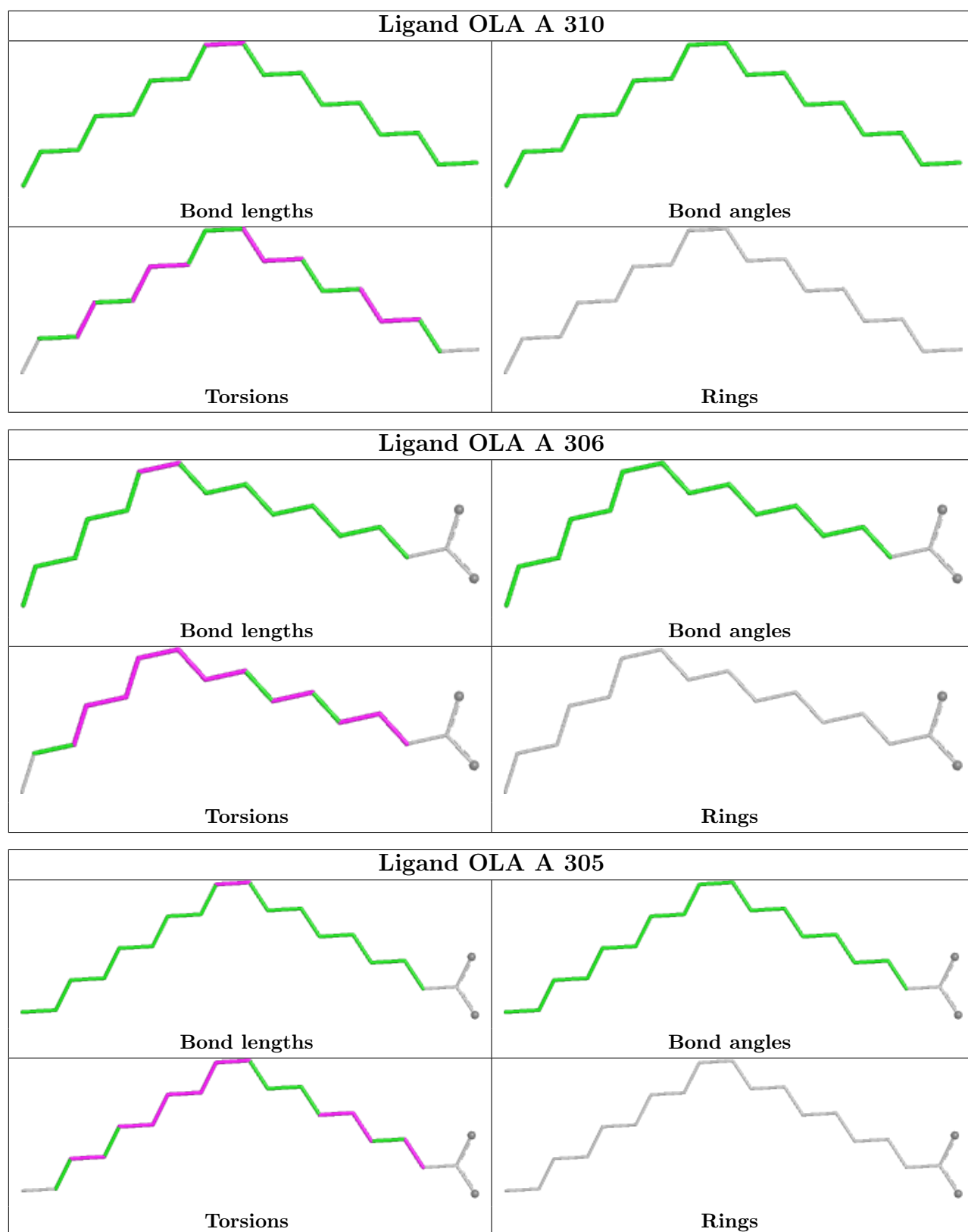
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	RET	4	0
4	A	304	OLA	2	0

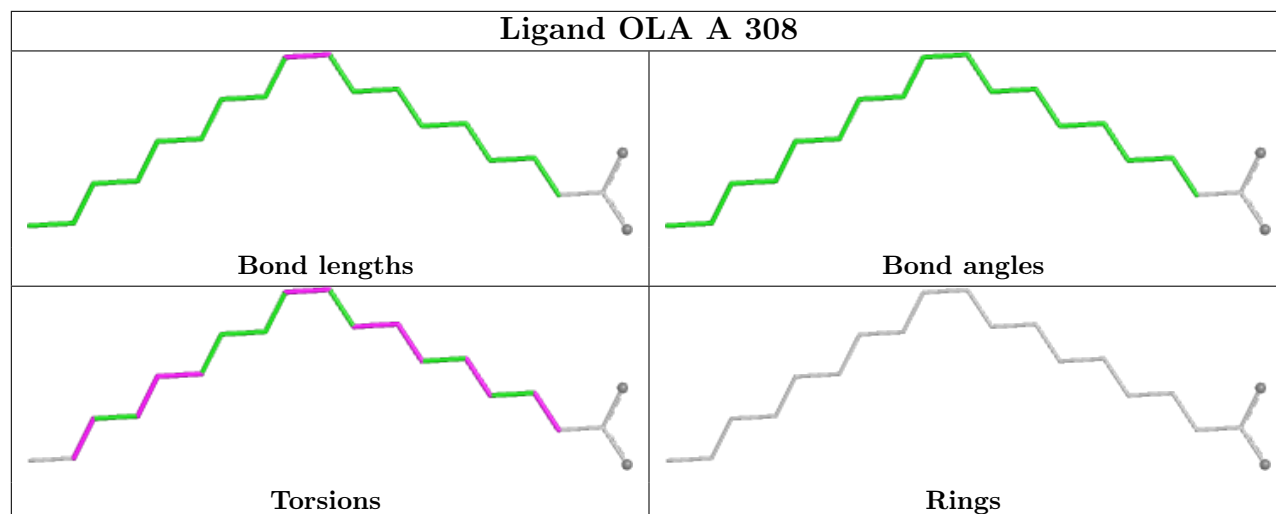
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	265/296 (89%)	0.30	15 (5%) <b>23</b> <b>29</b>	28, 39, 65, 93	0

All (15) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1	MET	7.5
1	A	262	LEU	5.2
1	A	260	GLN	5.0
1	A	263	GLY	4.6
1	A	256	VAL	3.7
1	A	-1	ASN	3.3
1	A	0	SER	3.2
1	A	253	ALA	2.7
1	A	259	GLN	2.6
1	A	63	LEU	2.6
1	A	261	ALA	2.6
1	A	86	GLN	2.6
1	A	88	LEU	2.6
1	A	87	ASP	2.5
1	A	89	THR	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](#)

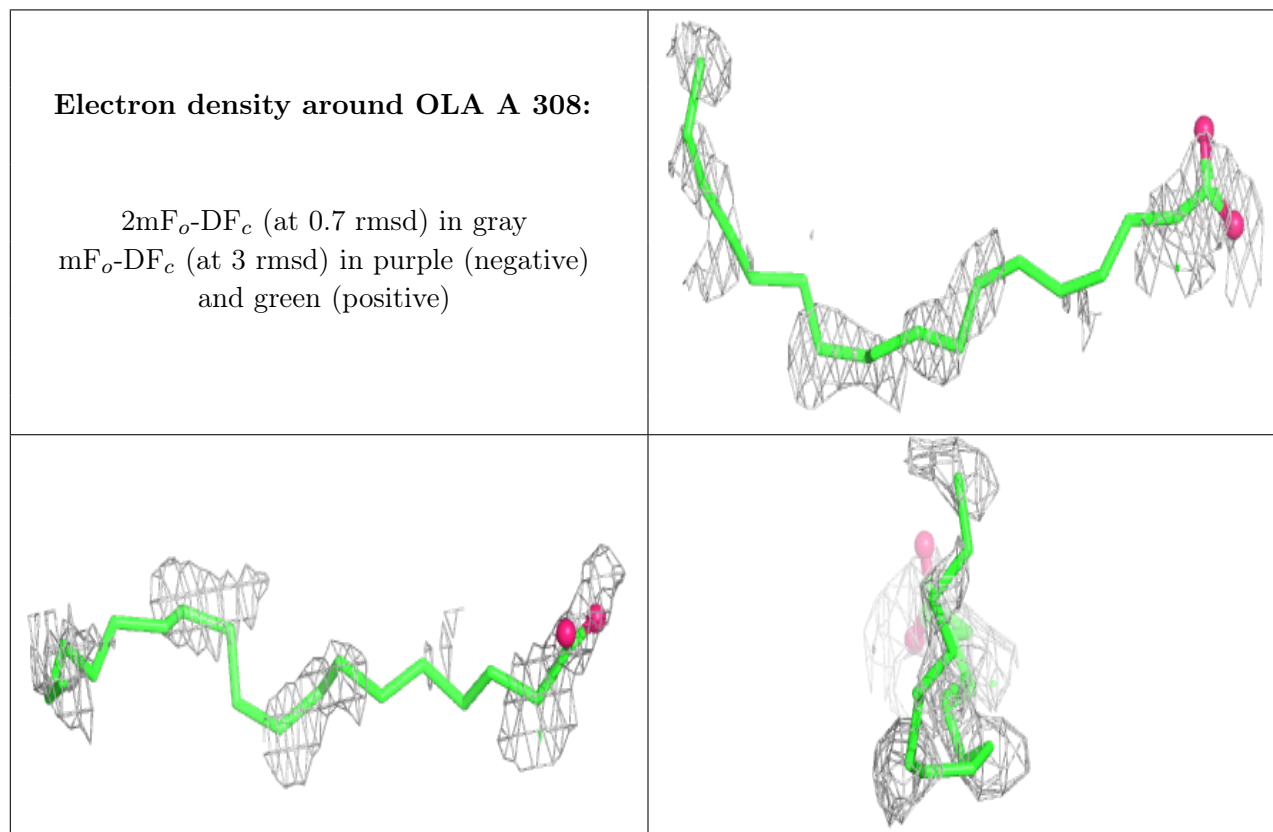
There are no monosaccharides in this entry.

## 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.

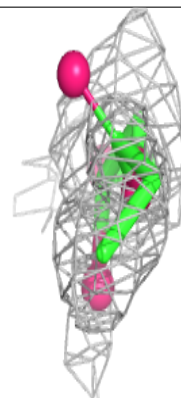
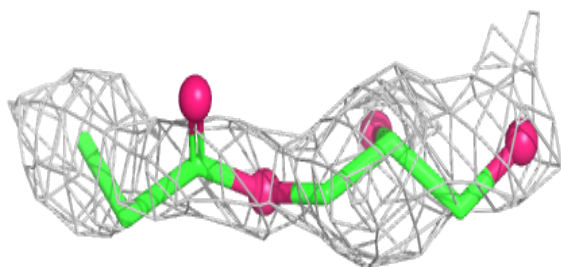
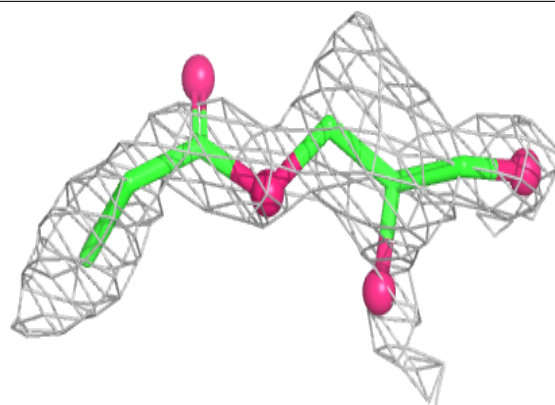
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
4	OLA	A	308	20/20	0.47	0.48	65,95,103,104	0
3	OLC	A	307	10/25	0.56	0.34	59,105,116,119	0
4	OLA	A	304	20/20	0.65	0.29	53,69,96,102	0
4	OLA	A	305	20/20	0.74	0.31	60,71,87,96	0
4	OLA	A	309	14/20	0.74	0.32	57,65,80,92	0
4	OLA	A	310	17/20	0.75	0.19	44,64,76,82	0
3	OLC	A	302	25/25	0.77	0.23	36,53,70,78	0
3	OLC	A	303	16/25	0.77	0.24	50,59,77,84	0
4	OLA	A	306	17/20	0.80	0.22	45,62,82,89	0
2	RET	A	301	20/21	0.93	0.13	27,31,38,39	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.

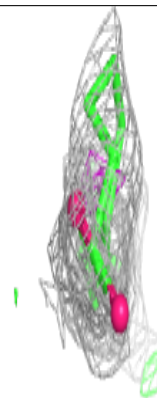
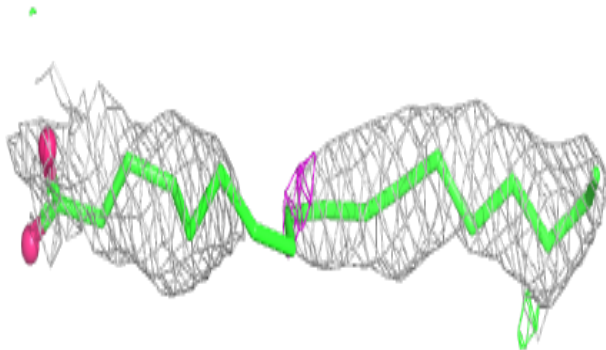
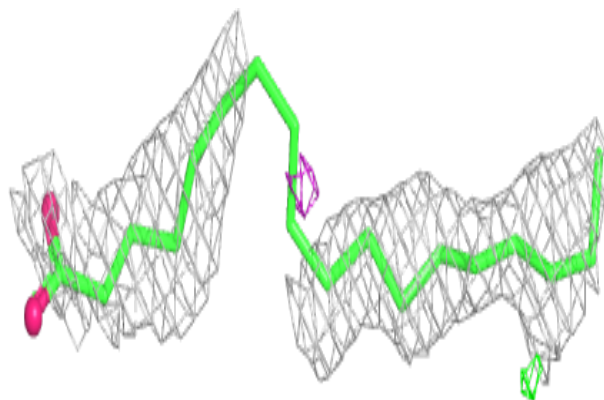


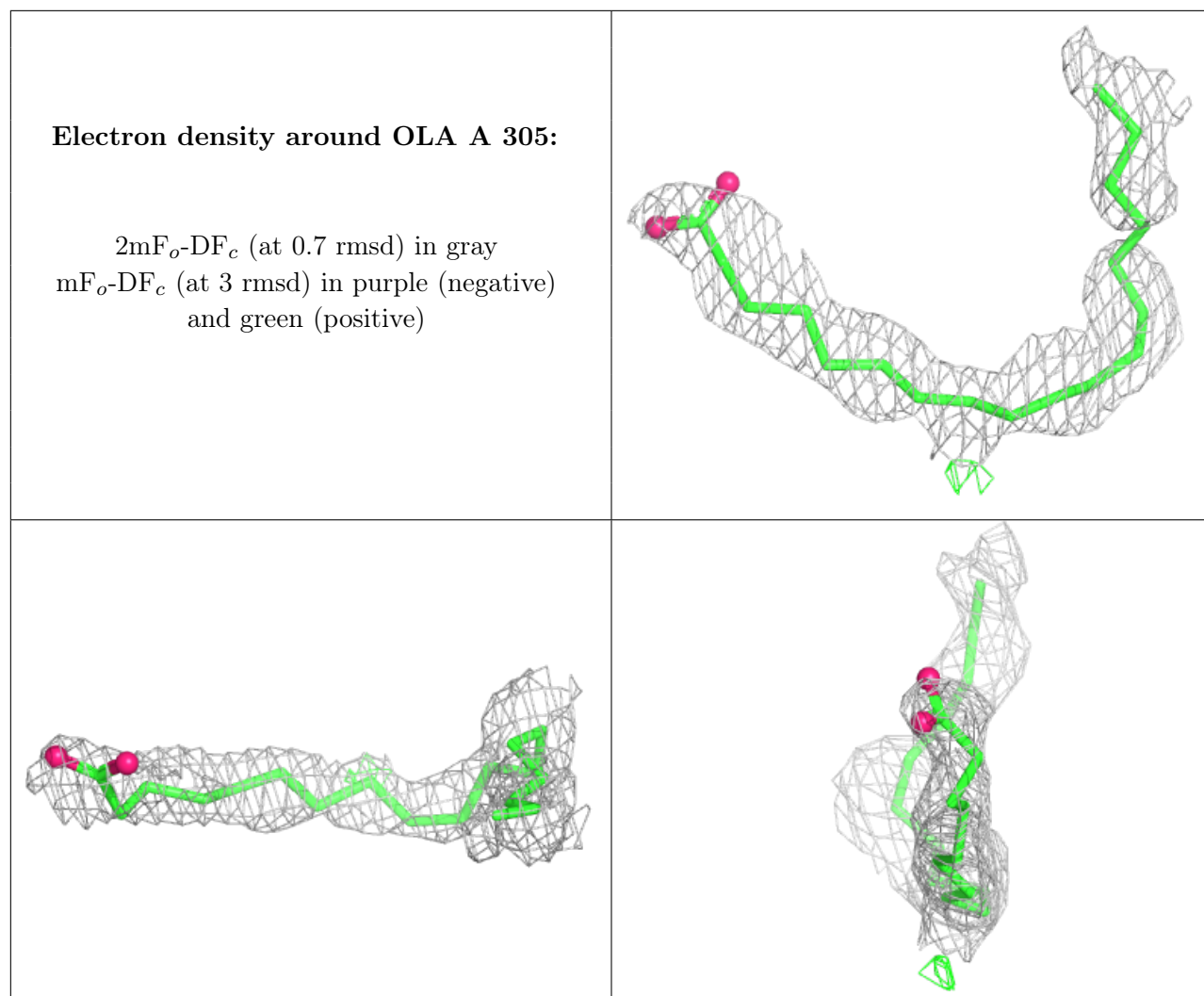
**Electron density around OLC A 307:**

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

**Electron density around OLA A 304:**

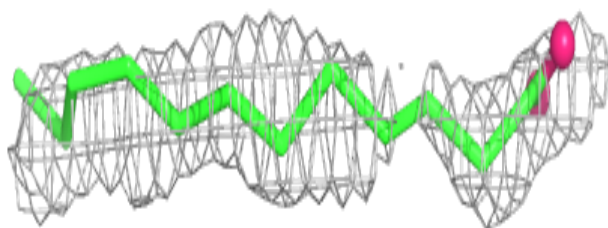
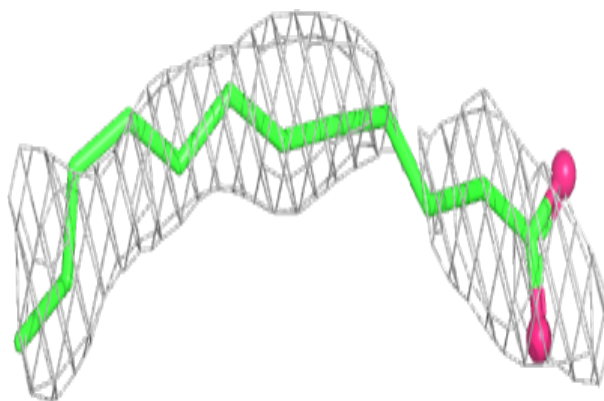
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



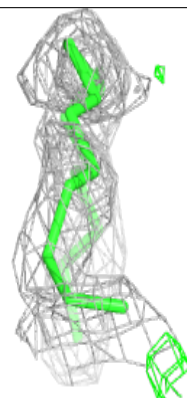
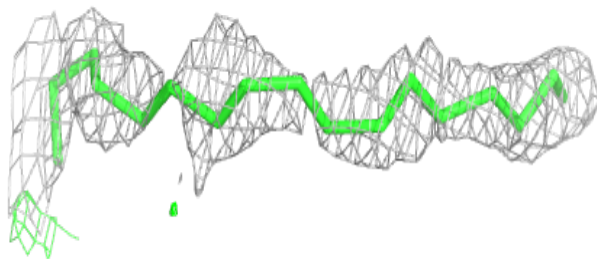
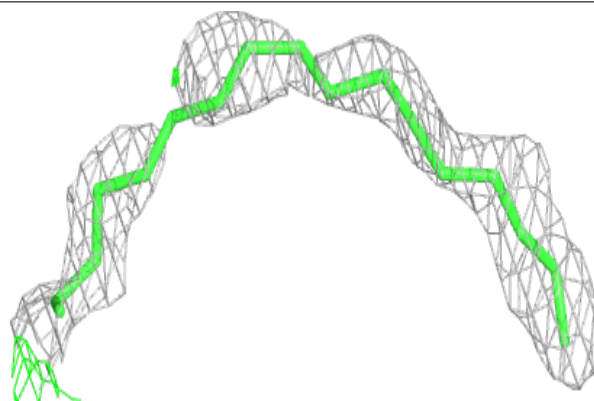


**Electron density around OLA A 309:**

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

**Electron density around OLA A 310:**

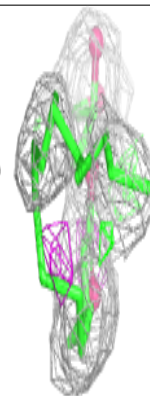
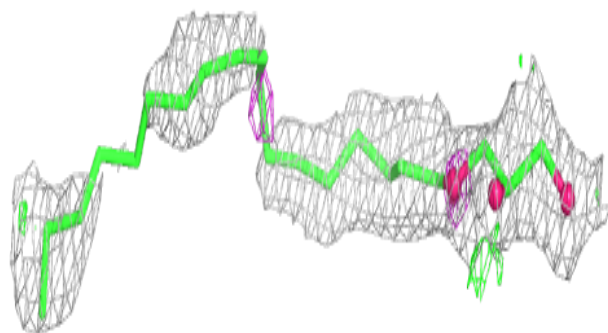
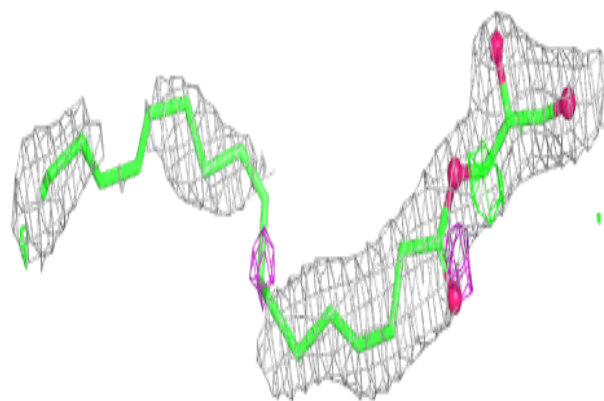
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



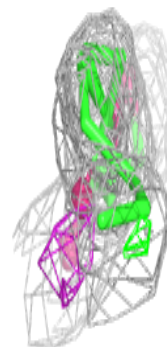
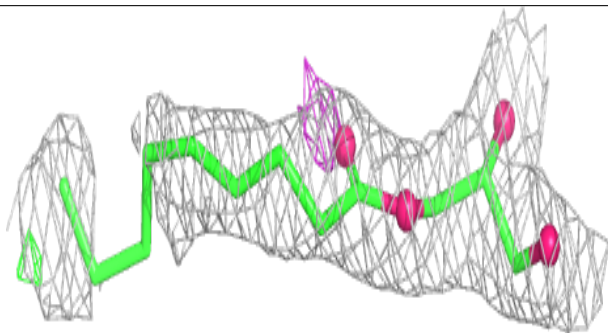
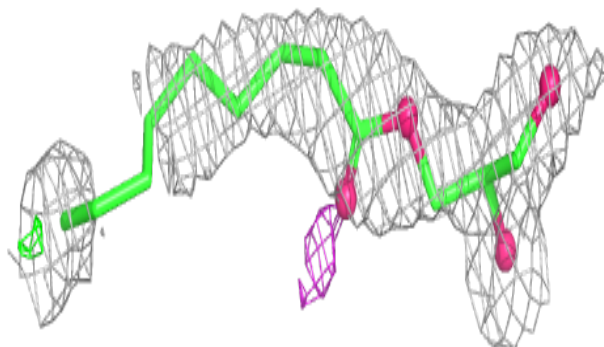


**Electron density around OLC A 302:**

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

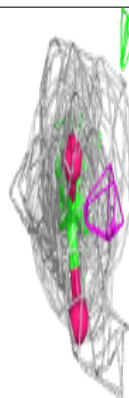
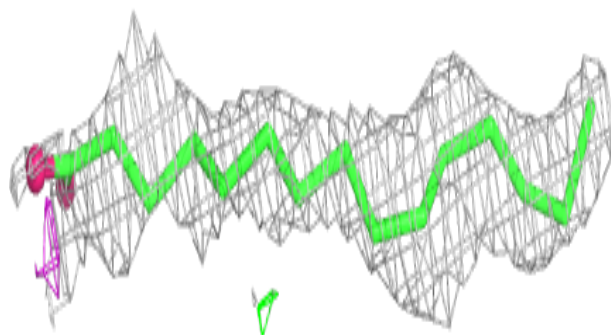
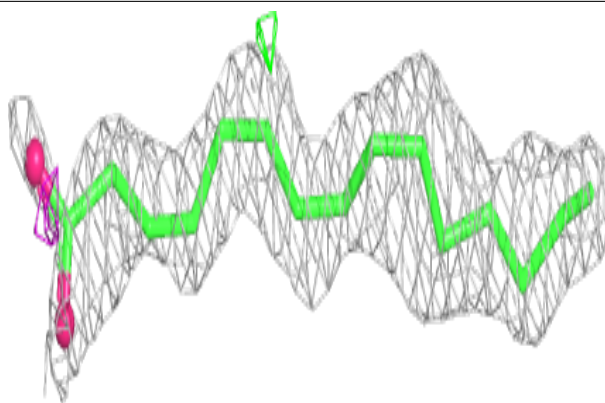
**Electron density around OLC A 303:**

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

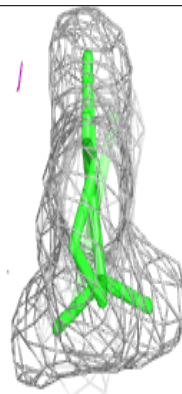
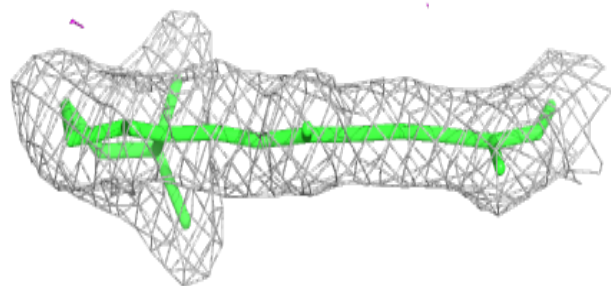
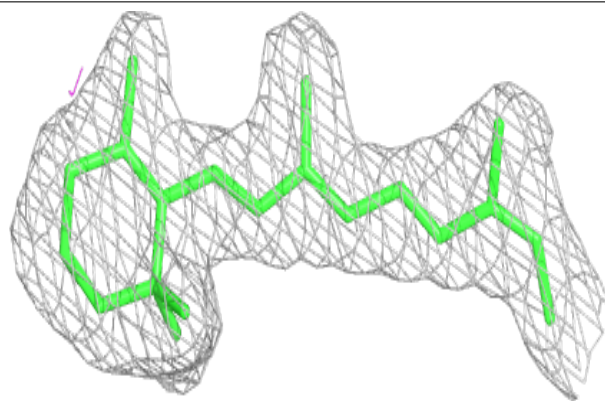


**Electron density around OLA A 306:**

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

**Electron density around RET A 301:**

$2mF_o-DF_c$  (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.