



wwPDB X-ray Structure Validation Summary Report ⓘ

Jan 15, 2024 – 11:09 am GMT

PDB ID : 6YT4
Title : Crystal structure of the N112A mutant of the light-driven sodium pump KR2 in the pentameric form, pH 8.0
Authors : Kovalev, K.; Maliar, N.; Astashkin, R.; Gordeliy, V.
Deposited on : 2020-04-23
Resolution : 2.40 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4.02b-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

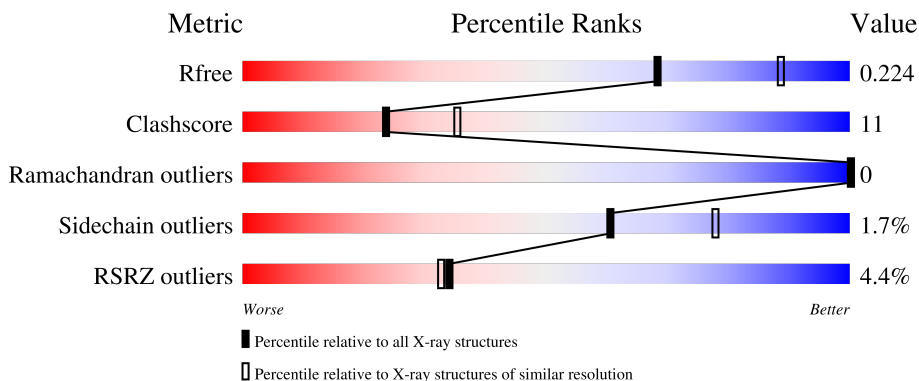
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

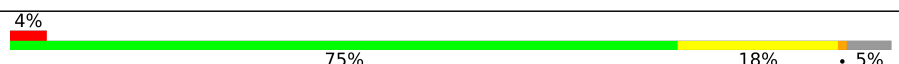
The reported resolution of this entry is 2.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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

Mol	Chain	Length	Quality of chain
1	A	288	 4% 75% 17% • 7%
1	B	288	 3% 77% 15% • 7%
1	C	288	 6% 76% 17% 7%
1	D	288	 3% 78% 15% 6%
1	E	288	 4% 75% 18% • 5%

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
3	OLA	A	307	-	-	-	X
3	OLA	E	403	-	-	-	X
5	BOG	A	309	-	-	-	X
5	BOG	B	310	-	-	-	X
5	BOG	C	1408	-	-	-	X
5	BOG	D	408	-	-	-	X
5	BOG	E	412	-	-	-	X

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 11860 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 Sodium pumping rhodopsin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	269	Total 2154	C 1437	N 327	O 381	S 9	0	2	0
1	B	269	Total 2153	C 1436	N 327	O 381	S 9	0	2	0
1	D	270	Total 2160	C 1440	N 329	O 382	S 9	0	2	0
1	E	273	Total 2179	C 1452	N 333	O 385	S 9	0	2	0
1	C	268	Total 2145	C 1431	N 325	O 380	S 9	0	2	0

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	112	ALA	ASN	engineered mutation	UNP N0DKS8
A	281	LEU	-	expression tag	UNP N0DKS8
A	282	GLU	-	expression tag	UNP N0DKS8
A	283	HIS	-	expression tag	UNP N0DKS8
A	284	HIS	-	expression tag	UNP N0DKS8
A	285	HIS	-	expression tag	UNP N0DKS8
A	286	HIS	-	expression tag	UNP N0DKS8
A	287	HIS	-	expression tag	UNP N0DKS8
A	288	HIS	-	expression tag	UNP N0DKS8
B	112	ALA	ASN	engineered mutation	UNP N0DKS8
B	281	LEU	-	expression tag	UNP N0DKS8
B	282	GLU	-	expression tag	UNP N0DKS8
B	283	HIS	-	expression tag	UNP N0DKS8
B	284	HIS	-	expression tag	UNP N0DKS8
B	285	HIS	-	expression tag	UNP N0DKS8
B	286	HIS	-	expression tag	UNP N0DKS8
B	287	HIS	-	expression tag	UNP N0DKS8
B	288	HIS	-	expression tag	UNP N0DKS8
D	112	ALA	ASN	engineered mutation	UNP N0DKS8

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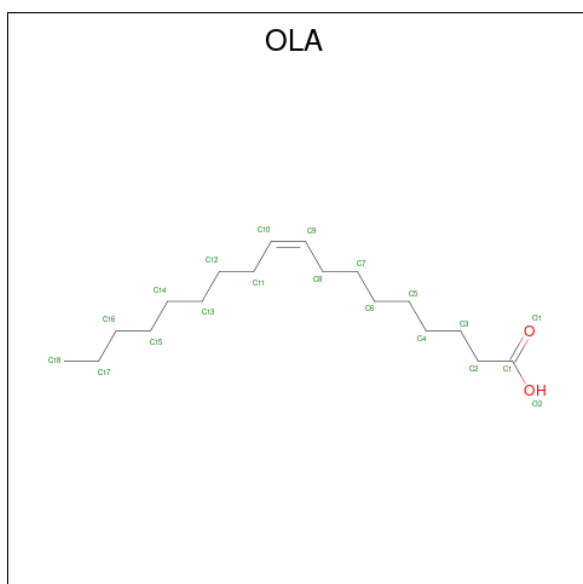
Chain	Residue	Modelled	Actual	Comment	Reference
D	281	LEU	-	expression tag	UNP N0DKS8
D	282	GLU	-	expression tag	UNP N0DKS8
D	283	HIS	-	expression tag	UNP N0DKS8
D	284	HIS	-	expression tag	UNP N0DKS8
D	285	HIS	-	expression tag	UNP N0DKS8
D	286	HIS	-	expression tag	UNP N0DKS8
D	287	HIS	-	expression tag	UNP N0DKS8
D	288	HIS	-	expression tag	UNP N0DKS8
E	112	ALA	ASN	engineered mutation	UNP N0DKS8
E	281	LEU	-	expression tag	UNP N0DKS8
E	282	GLU	-	expression tag	UNP N0DKS8
E	283	HIS	-	expression tag	UNP N0DKS8
E	284	HIS	-	expression tag	UNP N0DKS8
E	285	HIS	-	expression tag	UNP N0DKS8
E	286	HIS	-	expression tag	UNP N0DKS8
E	287	HIS	-	expression tag	UNP N0DKS8
E	288	HIS	-	expression tag	UNP N0DKS8
C	112	ALA	ASN	engineered mutation	UNP N0DKS8
C	281	LEU	-	expression tag	UNP N0DKS8
C	282	GLU	-	expression tag	UNP N0DKS8
C	283	HIS	-	expression tag	UNP N0DKS8
C	284	HIS	-	expression tag	UNP N0DKS8
C	285	HIS	-	expression tag	UNP N0DKS8
C	286	HIS	-	expression tag	UNP N0DKS8
C	287	HIS	-	expression tag	UNP N0DKS8
C	288	HIS	-	expression tag	UNP N0DKS8

- Molecule 2 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (three-letter code: OLC) (formula: C₂₁H₄₀O₄).



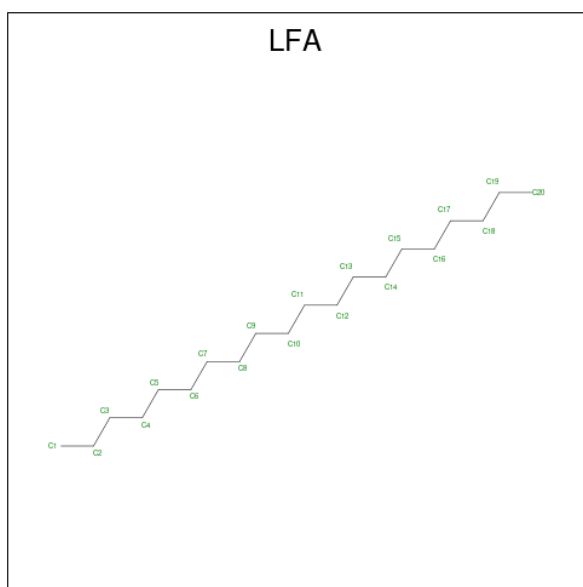
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C O	0	0
			12	8 4		
2	B	1	Total	C O	0	0
			13	10 3		
2	D	1	Total	C O	0	0
			15	12 3		
2	E	1	Total	C O	0	0
			10	6 4		
2	C	1	Total	C O	0	0
			11	8 3		

- Molecule 3 is OLEIC ACID (three-letter code: OLA) (formula: $C_{18}H_{34}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			17	15	2		
3	A	1	Total	C	O	0	0
			16	14	2		
3	A	1	Total	C	O	0	0
			10	8	2		
3	A	1	Total	C	O	0	0
			13	11	2		
3	A	1	Total	C	O	0	0
			5	3	2		
3	B	1	Total	C	O	0	0
			14	12	2		
3	B	1	Total	C	O	0	0
			4	2	2		
3	B	1	Total	C	O	0	0
			16	14	2		
3	B	1	Total	C	O	0	0
			12	10	2		
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			20	18	2		
3	E	1	Total	C	O	0	0
			4	2	2		
3	E	1	Total	C	O	0	0
			20	18	2		
3	E	1	Total	C	O	0	0
			12	10	2		
3	E	1	Total	C	O	0	0
			18	16	2		
3	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 4 is EICOSANE (three-letter code: LFA) (formula: C₂₀H₄₂).



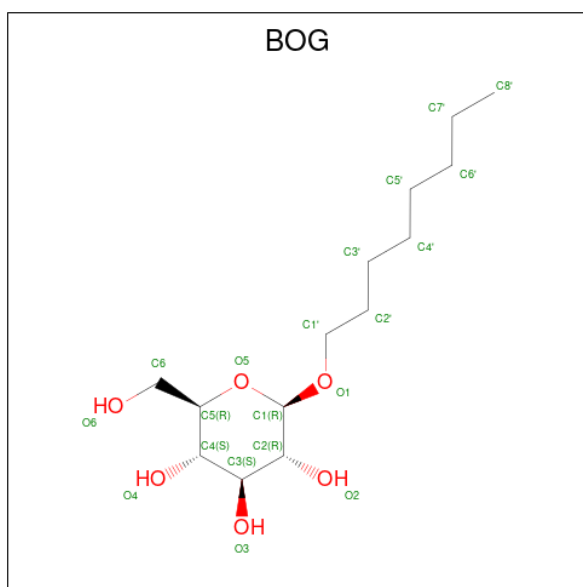
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C 20 20	0	0
4	A	1	Total C 5 5	0	0
4	A	1	Total C 8 8	0	0
4	B	1	Total C 19 19	0	0
4	B	1	Total C 11 11	0	0
4	B	1	Total C 20 20	0	0
4	B	1	Total C 6 6	0	0
4	B	1	Total C 9 9	0	0
4	B	1	Total C 20 20	0	0
4	D	1	Total C 20 20	0	0
4	D	1	Total C 18 18	0	0
4	D	1	Total C 10 10	0	0
4	D	1	Total C 15 15	0	0
4	D	1	Total C 7 7	0	0

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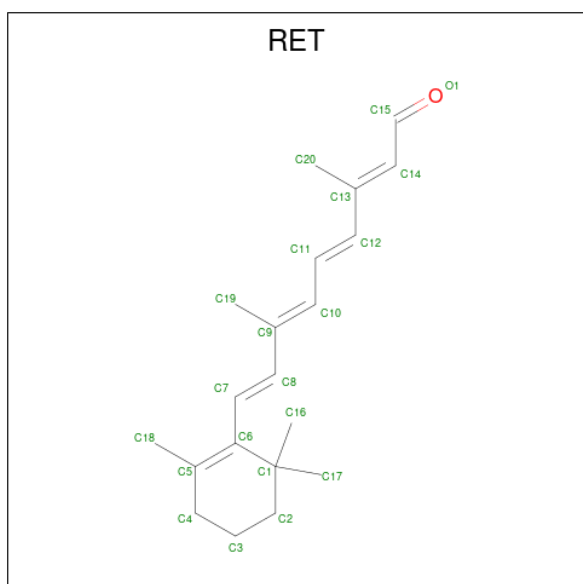
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	D	1	Total C 8 8	0	0
4	D	1	Total C 7 7	0	0
4	E	1	Total C 5 5	0	0
4	E	1	Total C 20 20	0	0
4	E	1	Total C 19 19	0	0
4	E	1	Total C 18 18	0	0
4	E	1	Total C 5 5	0	0
4	E	1	Total C 5 5	0	0
4	E	1	Total C 8 8	0	0
4	C	1	Total C 16 16	0	0
4	C	1	Total C 15 15	0	0
4	C	1	Total C 3 3	0	0
4	C	1	Total C 6 6	0	0
4	C	1	Total C 5 5	0	0

- Molecule 5 is octyl beta-D-glucopyranoside (three-letter code: BOG) (formula: C₁₄H₂₈O₆).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 20 14 6	0	0
5	B	1	Total C O 20 14 6	0	0
5	D	1	Total C O 20 14 6	0	0
5	E	1	Total C O 20 14 6	0	0
5	C	1	Total C O 20 14 6	0	0

- Molecule 6 is RETINAL (three-letter code: RET) (formula: $C_{20}H_{28}O$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	1	Total C 20 20	0	0
6	B	1	Total C 20 20	0	0
6	D	1	Total C 20 20	0	0
6	E	1	Total C 20 20	0	0
6	C	1	Total C 20 20	0	0

- Molecule 7 is SODIUM ION (three-letter code: NA) (formula: Na) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total Na 1 1	0	0
7	B	1	Total Na 1 1	0	0
7	D	1	Total Na 1 1	0	0
7	E	1	Total Na 1 1	0	0
7	C	1	Total Na 1 1	0	0

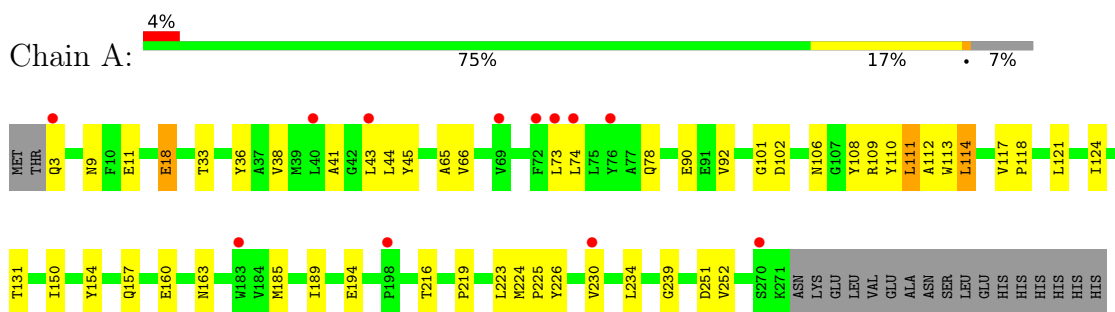
- Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	56	Total O 56 56	0	0
8	B	67	Total O 67 67	0	0
8	D	54	Total O 54 54	0	0
8	E	56	Total O 56 56	0	0
8	C	53	Total O 53 53	0	0

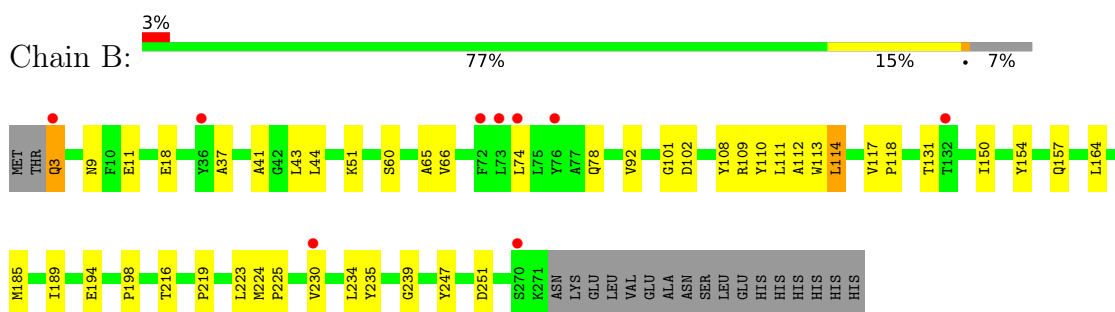
3 Residue-property plots [i](#)

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.

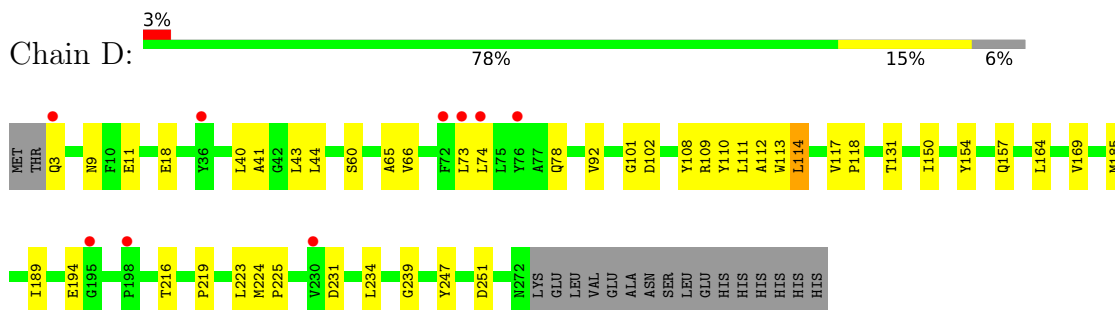
- Molecule 1: Sodium pumping rhodopsin



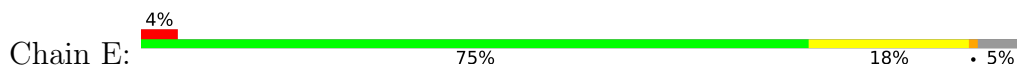
- Molecule 1: Sodium pumping rhodopsin

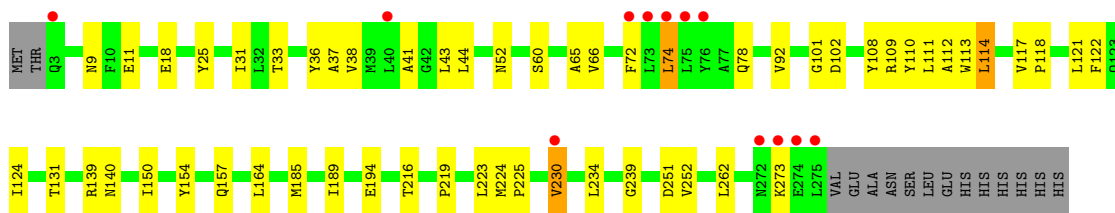


- Molecule 1: Sodium pumping rhodopsin

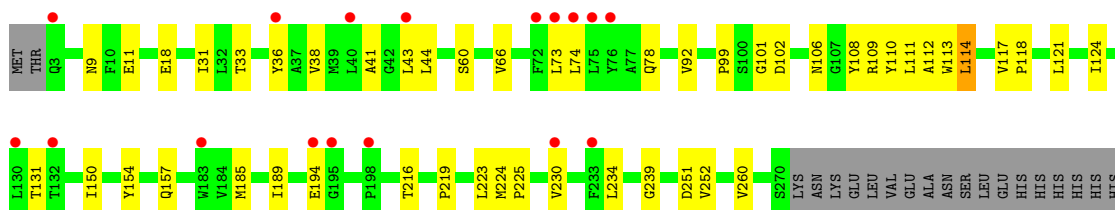
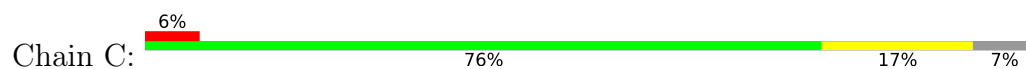


- Molecule 1: Sodium pumping rhodopsin





● Molecule 1: Sodium pumping rhodopsin



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	129.84Å 239.70Å 134.58Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.40 47.85 – 2.40	Depositor EDS
% Data completeness (in resolution range)	93.8 (20.00-2.40) 94.0 (47.85-2.40)	Depositor EDS
R_{merge}	0.22	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.39Å)	Xtrriage
Refinement program	REFMAC 5.8.0258	Depositor
R, R_{free}	0.186 , 0.217 0.198 , 0.224	Depositor DCC
R_{free} test set	3771 reflections (4.89%)	wwPDB-VP
Wilson B-factor (Å ²)	41.4	Xtrriage
Anisotropy	0.056	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.35$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11860	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.98% 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: NA, OLC, OLA, BOG, LFA, 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	0.63	0/2212	0.62	0/3008
1	B	0.63	0/2211	0.63	0/3007
1	C	0.63	0/2203	0.63	0/2997
1	D	0.63	0/2218	0.62	0/3016
1	E	0.63	0/2237	0.63	0/3041
All	All	0.63	0/11081	0.62	0/15069

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	2154	0	2133	50	0
1	B	2153	0	2131	49	0
1	C	2145	0	2120	43	0
1	D	2160	0	2140	50	0
1	E	2179	0	2157	60	0
2	A	12	0	13	0	0
2	B	13	0	14	0	0
2	C	11	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	15	0	18	0	0
2	E	10	0	9	0	0
3	A	61	0	76	6	0
3	B	46	0	55	2	0
3	C	4	0	0	2	0
3	D	24	0	33	2	0
3	E	54	0	74	4	0
4	A	33	0	63	0	0
4	B	85	0	170	8	0
4	C	45	0	82	1	0
4	D	85	0	158	6	0
4	E	80	0	150	12	0
5	A	20	0	28	1	0
5	B	20	0	28	2	0
5	C	20	0	28	0	0
5	D	20	0	28	2	0
5	E	20	0	28	4	0
6	A	20	0	27	3	0
6	B	20	0	27	4	0
6	C	20	0	27	4	0
6	D	20	0	27	4	0
6	E	20	0	27	4	0
7	A	1	0	0	0	0
7	B	1	0	0	0	0
7	C	1	0	0	0	0
7	D	1	0	0	0	0
7	E	1	0	0	0	0
8	A	56	0	0	3	0
8	B	67	0	0	8	0
8	C	53	0	0	4	0
8	D	54	0	0	3	0
8	E	56	0	0	5	0
All	All	11860	0	11881	262	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

The worst 5 of 262 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:D:401:LFA:C1	3:C:1401:OLA:C2	1.79	1.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:403:OLA:C2	4:E:406:LFA:H11	1.05	1.52
3:D:406:OLA:C2	4:E:402:LFA:H11	1.38	1.49
3:E:403:OLA:C2	4:E:406:LFA:C1	1.86	1.49
3:A:311:OLA:C3	4:B:304:LFA:H201	1.56	1.35

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	A	269/288 (93%)	263 (98%)	6 (2%)	0	100	100
1	B	269/288 (93%)	262 (97%)	7 (3%)	0	100	100
1	C	268/288 (93%)	262 (98%)	6 (2%)	0	100	100
1	D	270/288 (94%)	261 (97%)	9 (3%)	0	100	100
1	E	273/288 (95%)	265 (97%)	8 (3%)	0	100	100
All	All	1349/1440 (94%)	1313 (97%)	36 (3%)	0	100	100

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	227/247 (92%)	222 (98%)	5 (2%)	52	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	B	227/247 (92%)	222 (98%)	5 (2%)	52	71
1	C	226/247 (92%)	223 (99%)	3 (1%)	69	84
1	D	228/247 (92%)	226 (99%)	2 (1%)	78	90
1	E	229/247 (93%)	224 (98%)	5 (2%)	52	71
All	All	1137/1235 (92%)	1117 (98%)	20 (2%)	60	76

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	92	VAL
1	C	92	VAL
1	C	230	VAL
1	C	114	LEU
1	B	51	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	3	GLN
1	B	81	ASN
1	E	141	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 64 ligands modelled in this entry, 5 are monoatomic - leaving 59 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	LFA	E	406	-	18,18,19	0.08	0	17,17,18	0.06	0
4	LFA	E	409	-	4,4,19	0.15	0	3,3,18	0.24	0
2	OLC	B	301	-	11,12,24	1.36	1 (9%)	11,13,25	1.12	2 (18%)
3	OLA	D	409	-	19,19,19	0.53	0	19,19,19	0.48	0
4	LFA	D	411	-	6,6,19	0.14	0	5,5,18	0.12	0
3	OLA	B	305	-	13,13,19	0.61	0	12,13,19	0.57	0
3	OLA	B	309	-	15,15,19	0.58	0	15,15,19	0.52	0
2	OLC	C	1402	-	10,10,24	1.46	1 (10%)	10,10,25	1.18	1 (10%)
4	LFA	A	305	-	4,4,19	0.15	0	3,3,18	0.22	0
4	LFA	B	302	-	18,18,19	0.08	0	17,17,18	0.07	0
4	LFA	B	312	-	19,19,19	0.07	0	18,18,18	0.04	0
4	LFA	E	410	-	4,4,19	0.16	0	3,3,18	0.21	0
6	RET	E	414	1	20,20,21	1.64	3 (15%)	27,27,28	1.09	1 (3%)
4	LFA	D	403	-	9,9,19	0.12	0	8,8,18	0.07	0
4	LFA	E	402	-	19,19,19	0.08	0	18,18,18	0.04	0
6	RET	B	311	1	20,20,21	1.64	3 (15%)	27,27,28	1.14	2 (7%)
3	OLA	A	308	-	12,12,19	0.63	0	12,12,19	0.61	0
4	LFA	D	407	-	6,6,19	0.13	0	5,5,18	0.07	0
4	LFA	C	1403	-	15,15,19	0.11	0	14,14,18	0.07	0
4	LFA	D	404	-	14,14,19	0.08	0	13,13,18	0.06	0
4	LFA	D	401	-	19,19,19	0.08	0	18,18,18	0.05	0
5	BOG	C	1408	-	20,20,20	0.59	1 (5%)	25,25,25	0.59	0
4	LFA	E	401	-	4,4,19	0.17	0	3,3,18	0.24	0
4	LFA	B	304	-	19,19,19	0.07	0	18,18,18	0.04	0
4	LFA	D	402	-	17,17,19	0.08	0	16,16,18	0.05	0
4	LFA	B	303	-	10,10,19	0.12	0	9,9,18	0.08	0
3	OLA	A	307	-	9,9,19	0.71	0	9,9,19	0.66	0
4	LFA	A	303	-	19,19,19	0.07	0	18,18,18	0.04	0
3	OLA	B	308	-	3,3,19	0.87	0	3,3,19	0.77	0
4	LFA	A	306	-	7,7,19	0.14	0	6,6,18	0.11	0
5	BOG	B	310	-	20,20,20	0.59	1 (5%)	25,25,25	0.52	0
4	LFA	E	413	-	7,7,19	0.12	0	6,6,18	0.12	0
3	OLA	E	405	-	19,19,19	0.51	0	19,19,19	0.47	0
5	BOG	A	309	-	20,20,20	0.65	1 (5%)	25,25,25	0.59	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	OLC	E	404	-	9,9,24	1.42	1 (11%)	10,10,25	1.52	2 (20%)
4	LFA	E	408	-	17,17,19	0.08	0	16,16,18	0.07	0
3	OLA	D	406	-	3,3,19	0.88	0	3,3,19	0.80	0
6	RET	C	1409	1	20,20,21	1.64	3 (15%)	27,27,28	1.09	1 (3%)
3	OLA	B	313	-	11,11,19	0.67	0	11,11,19	0.60	0
3	OLA	E	411	-	17,17,19	0.55	0	17,17,19	0.49	0
2	OLC	A	301	-	11,11,24	1.33	1 (9%)	12,12,25	1.12	2 (16%)
3	OLA	A	304	-	15,15,19	0.58	0	15,15,19	0.54	0
6	RET	A	310	1	20,20,21	1.66	3 (15%)	27,27,28	1.09	1 (3%)
3	OLA	E	407	-	11,11,19	0.64	0	11,11,19	0.65	0
4	LFA	D	410	-	7,7,19	0.11	0	6,6,18	0.11	0
2	OLC	D	405	-	13,14,24	1.27	1 (7%)	13,15,25	1.04	1 (7%)
3	OLA	E	403	-	3,3,19	0.87	0	3,3,19	0.77	0
3	OLA	A	302	-	16,16,19	0.56	0	16,16,19	0.55	0
6	RET	D	412	1	20,20,21	1.69	3 (15%)	27,27,28	1.08	2 (7%)
5	BOG	D	408	-	20,20,20	0.60	1 (5%)	25,25,25	0.60	0
3	OLA	A	311	-	4,4,19	0.98	0	4,4,19	0.71	0
4	LFA	C	1406	-	5,5,19	0.12	0	4,4,18	0.11	0
3	OLA	C	1401	-	3,3,19	0.85	0	3,3,19	0.80	0
4	LFA	C	1405	-	2,2,19	0.06	0	0,1,18	-	-
4	LFA	B	307	-	8,8,19	0.13	0	7,7,18	0.10	0
4	LFA	C	1404	-	14,14,19	0.08	0	13,13,18	0.06	0
4	LFA	C	1407	-	4,4,19	0.15	0	3,3,18	0.23	0
4	LFA	B	306	-	5,5,19	0.14	0	4,4,18	0.09	0
5	BOG	E	412	-	20,20,20	0.62	1 (5%)	25,25,25	0.51	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	LFA	E	406	-	-	9/16/16/17	-
4	LFA	E	409	-	-	1/2/2/17	-
2	OLC	B	301	-	-	2/11/11/24	-
3	OLA	D	409	-	-	11/17/17/17	-
4	LFA	D	411	-	-	3/4/4/17	-
3	OLA	B	305	-	-	8/11/11/17	-
3	OLA	B	309	-	-	5/13/13/17	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OLC	C	1402	-	-	1/9/9/24	-
4	LFA	A	305	-	-	0/2/2/17	-
4	LFA	B	302	-	-	4/16/16/17	-
4	LFA	B	312	-	-	11/17/17/17	-
4	LFA	E	410	-	-	1/2/2/17	-
6	RET	E	414	1	-	1/13/30/31	0/1/1/1
4	LFA	D	403	-	-	5/7/7/17	-
4	LFA	E	402	-	-	7/17/17/17	-
6	RET	B	311	1	-	0/13/30/31	0/1/1/1
3	OLA	A	308	-	-	7/10/10/17	-
4	LFA	D	407	-	-	1/4/4/17	-
4	LFA	C	1403	-	-	9/13/13/17	-
4	LFA	D	404	-	-	6/12/12/17	-
4	LFA	D	401	-	-	7/17/17/17	-
5	BOG	C	1408	-	-	5/11/31/31	0/1/1/1
4	LFA	E	401	-	-	0/2/2/17	-
4	LFA	B	304	-	-	11/17/17/17	-
4	LFA	D	402	-	-	7/15/15/17	-
4	LFA	B	303	-	-	4/8/8/17	-
3	OLA	A	307	-	-	5/7/7/17	-
4	LFA	A	303	-	-	9/17/17/17	-
4	LFA	A	306	-	-	5/5/5/17	-
5	BOG	B	310	-	-	8/11/31/31	0/1/1/1
4	LFA	E	413	-	-	4/5/5/17	-
3	OLA	E	405	-	-	8/17/17/17	-
5	BOG	A	309	-	-	5/11/31/31	0/1/1/1
2	OLC	E	404	-	-	2/9/9/24	-
4	LFA	E	408	-	-	6/15/15/17	-
6	RET	C	1409	1	-	0/13/30/31	0/1/1/1
3	OLA	B	313	-	-	3/9/9/17	-
3	OLA	E	411	-	-	12/15/15/17	-
2	OLC	A	301	-	-	2/11/11/24	-
3	OLA	A	304	-	-	4/13/13/17	-
6	RET	A	310	1	-	0/13/30/31	0/1/1/1
3	OLA	E	407	-	-	8/9/9/17	-
4	LFA	D	410	-	-	1/5/5/17	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	OLC	D	405	-	-	8/13/13/24	-
3	OLA	A	302	-	-	10/14/14/17	-
6	RET	D	412	1	-	1/13/30/31	0/1/1/1
5	BOG	D	408	-	-	5/11/31/31	0/1/1/1
3	OLA	A	311	-	-	0/2/2/17	-
4	LFA	C	1406	-	-	1/3/3/17	-
4	LFA	B	307	-	-	0/6/6/17	-
4	LFA	C	1404	-	-	5/12/12/17	-
4	LFA	C	1407	-	-	0/2/2/17	-
4	LFA	B	306	-	-	0/3/3/17	-
5	BOG	E	412	-	-	5/11/31/31	0/1/1/1

The worst 5 of 25 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	310	RET	C10-C9	4.51	1.41	1.35
6	C	1409	RET	C10-C9	4.43	1.41	1.35
2	D	405	OLC	O20-C1	4.42	1.46	1.33
2	C	1402	OLC	O20-C1	4.40	1.46	1.33
6	D	412	RET	C10-C9	4.36	1.41	1.35

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	311	RET	C19-C9-C10	-4.06	117.24	122.92
2	E	404	OLC	O20-C1-C2	3.89	121.58	111.38
6	E	414	RET	C19-C9-C10	-3.83	117.55	122.92
6	C	1409	RET	C19-C9-C10	-3.82	117.58	122.92
6	A	310	RET	C19-C9-C10	-3.77	117.65	122.92

There are no chirality outliers.

5 of 243 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	308	OLA	C11-C10-C9-C8
5	B	310	BOG	C4-C5-C6-O6
2	D	405	OLC	O19-C1-O20-C21
2	D	405	OLC	C2-C1-O20-C21
5	E	412	BOG	O5-C5-C6-O6

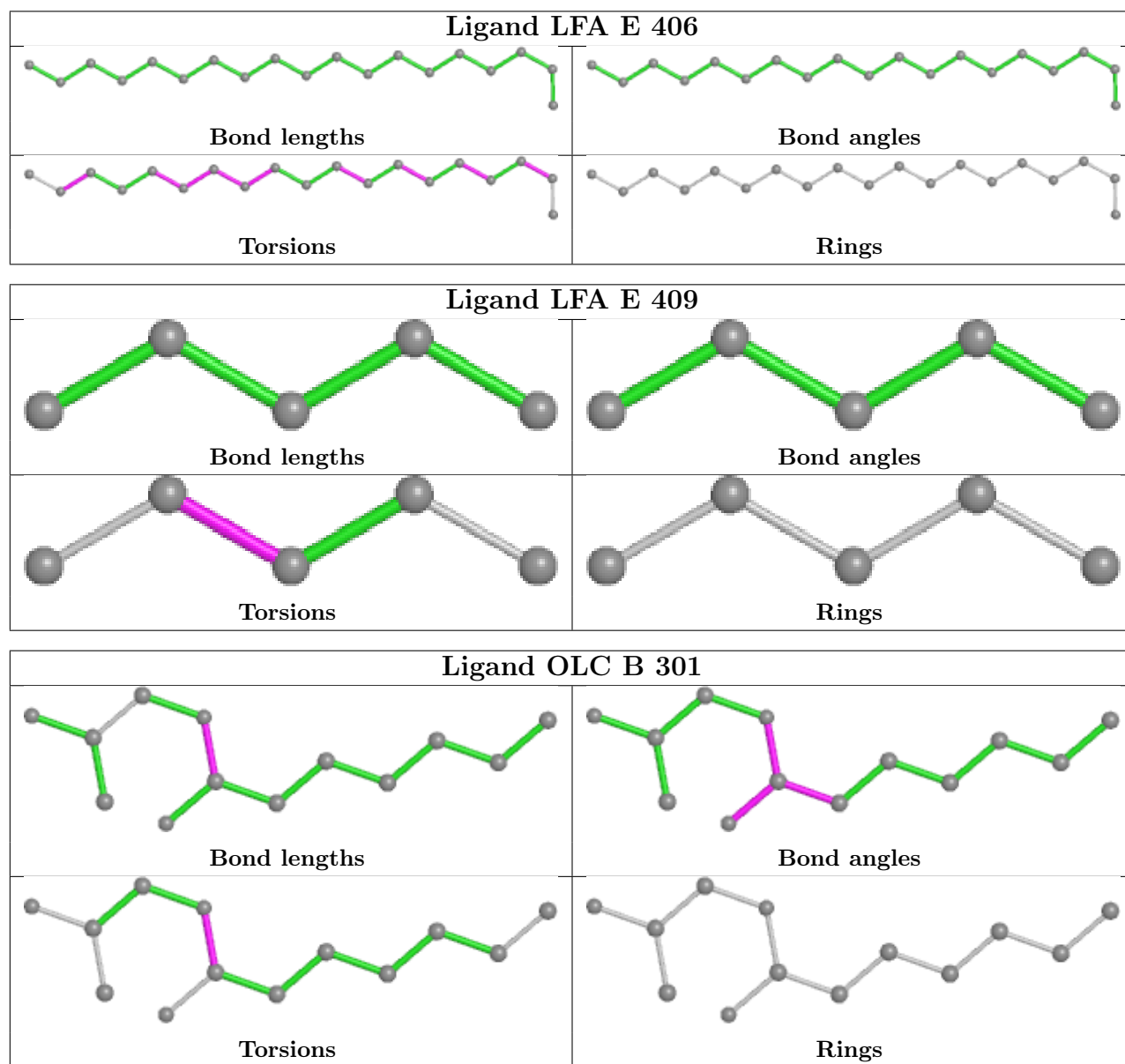
There are no ring outliers.

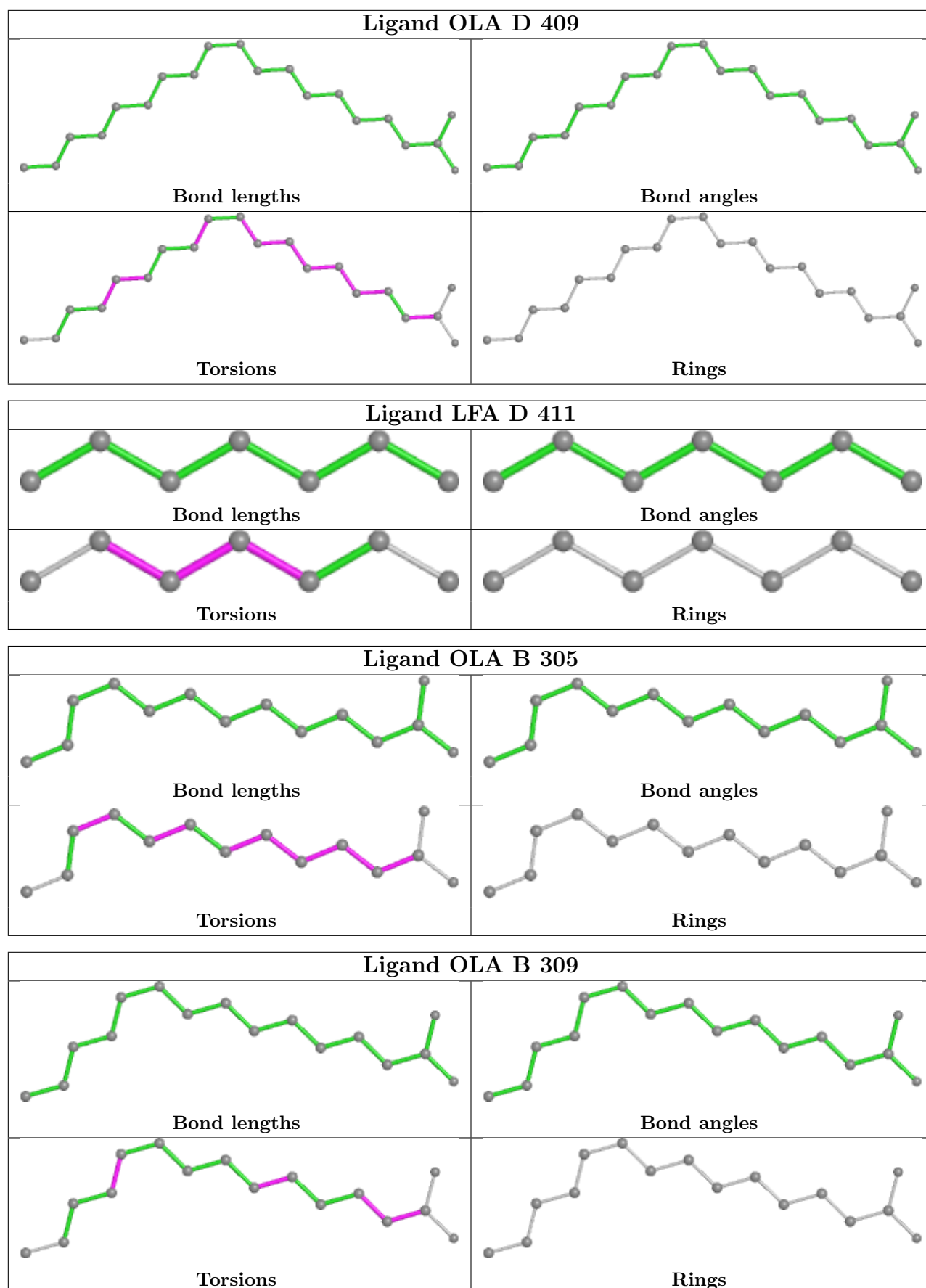
30 monomers are involved in 56 short contacts:

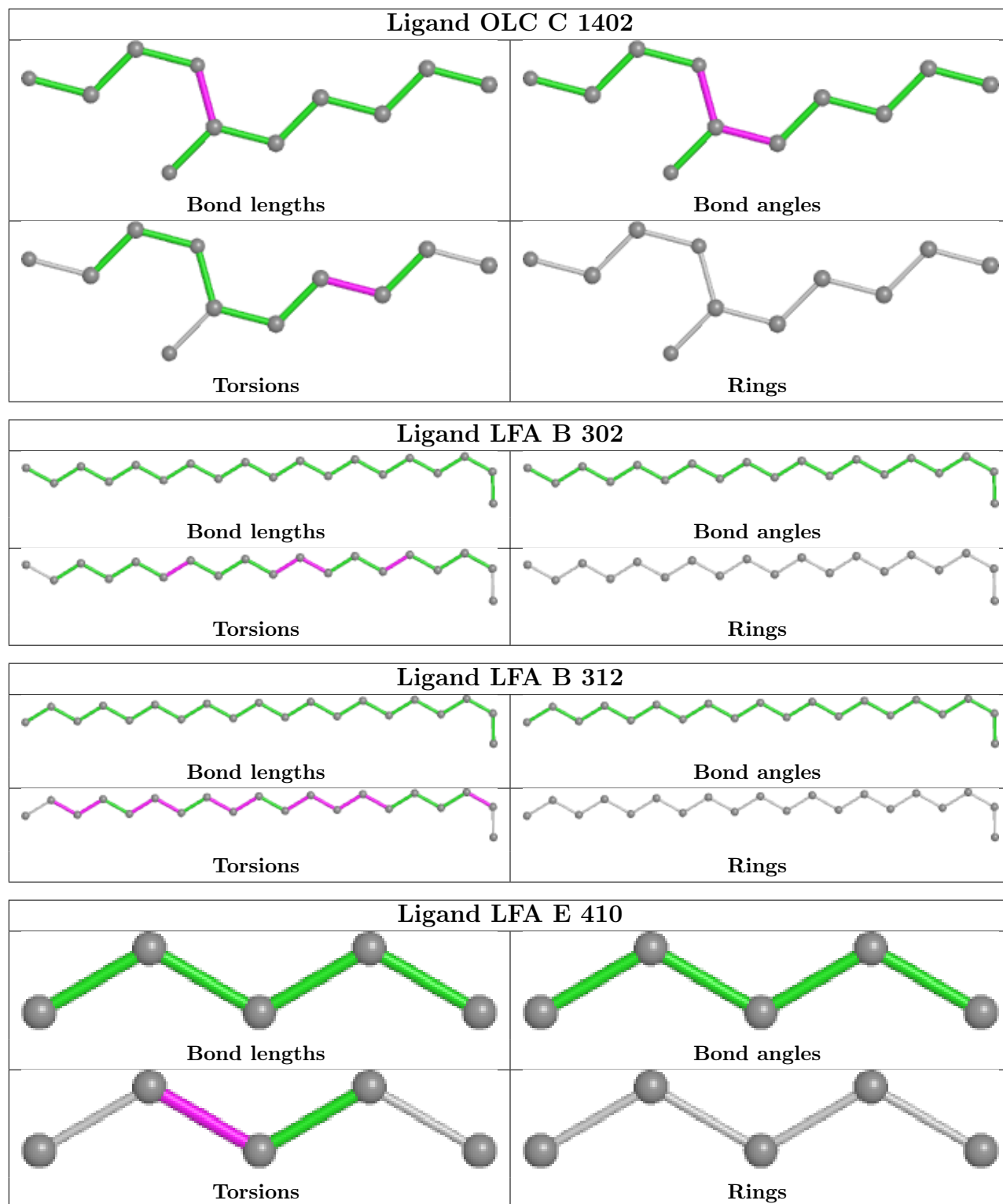
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	E	406	LFA	5	0
4	E	409	LFA	1	0
4	B	312	LFA	4	0
6	E	414	RET	4	0
4	D	403	LFA	1	0
4	E	402	LFA	3	0
6	B	311	RET	4	0
3	A	308	OLA	1	0
4	D	407	LFA	1	0
4	D	401	LFA	3	0
4	E	401	LFA	2	0
4	B	304	LFA	4	0
4	D	402	LFA	1	0
4	B	303	LFA	1	0
3	B	308	OLA	2	0
5	B	310	BOG	2	0
5	A	309	BOG	1	0
4	E	408	LFA	2	0
3	D	406	OLA	2	0
6	C	1409	RET	4	0
6	A	310	RET	3	0
3	E	407	OLA	1	0
3	E	403	OLA	3	0
3	A	302	OLA	2	0
6	D	412	RET	4	0
5	D	408	BOG	2	0
3	A	311	OLA	3	0
3	C	1401	OLA	2	0
4	C	1404	LFA	1	0
5	E	412	BOG	4	0

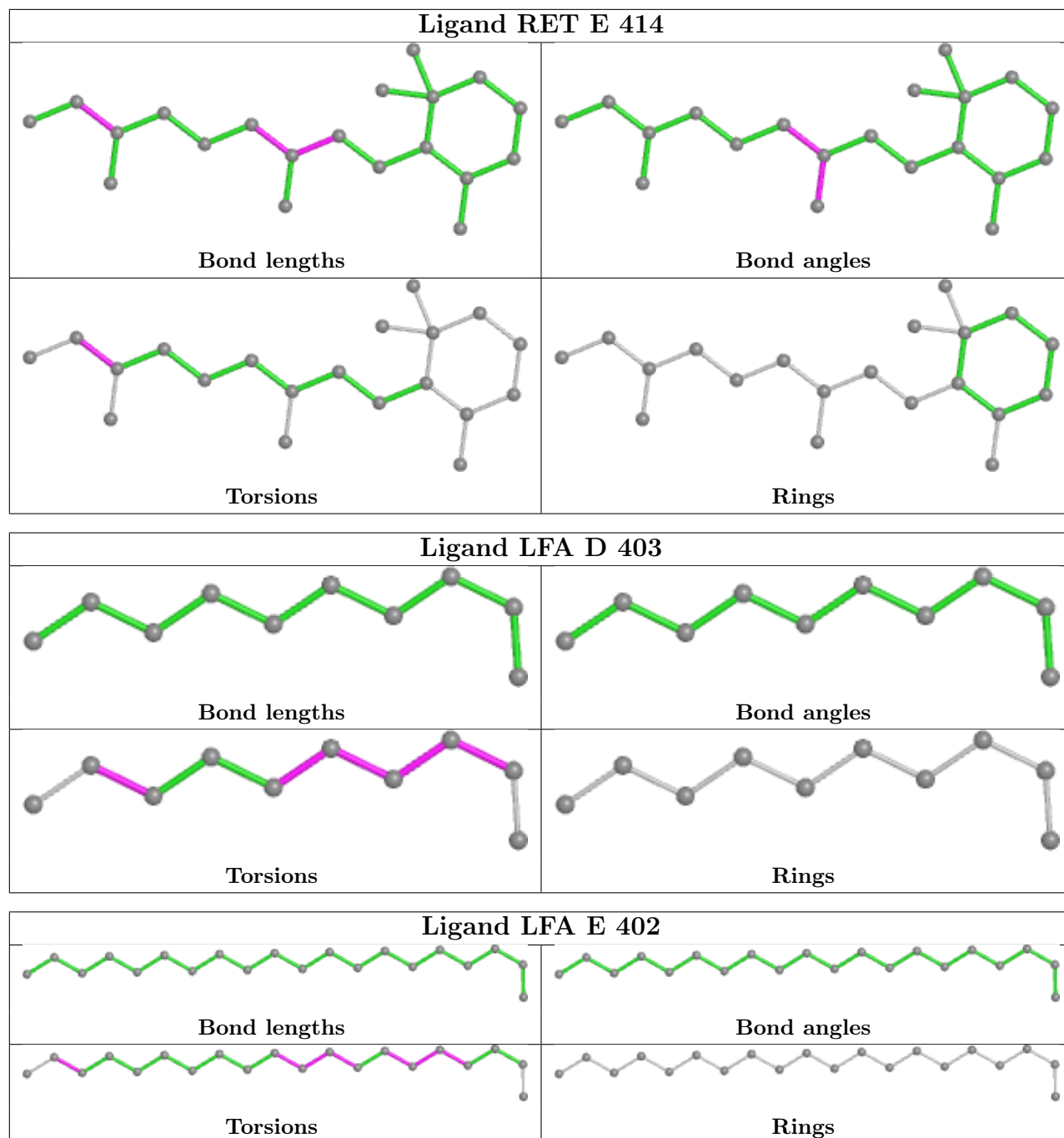
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

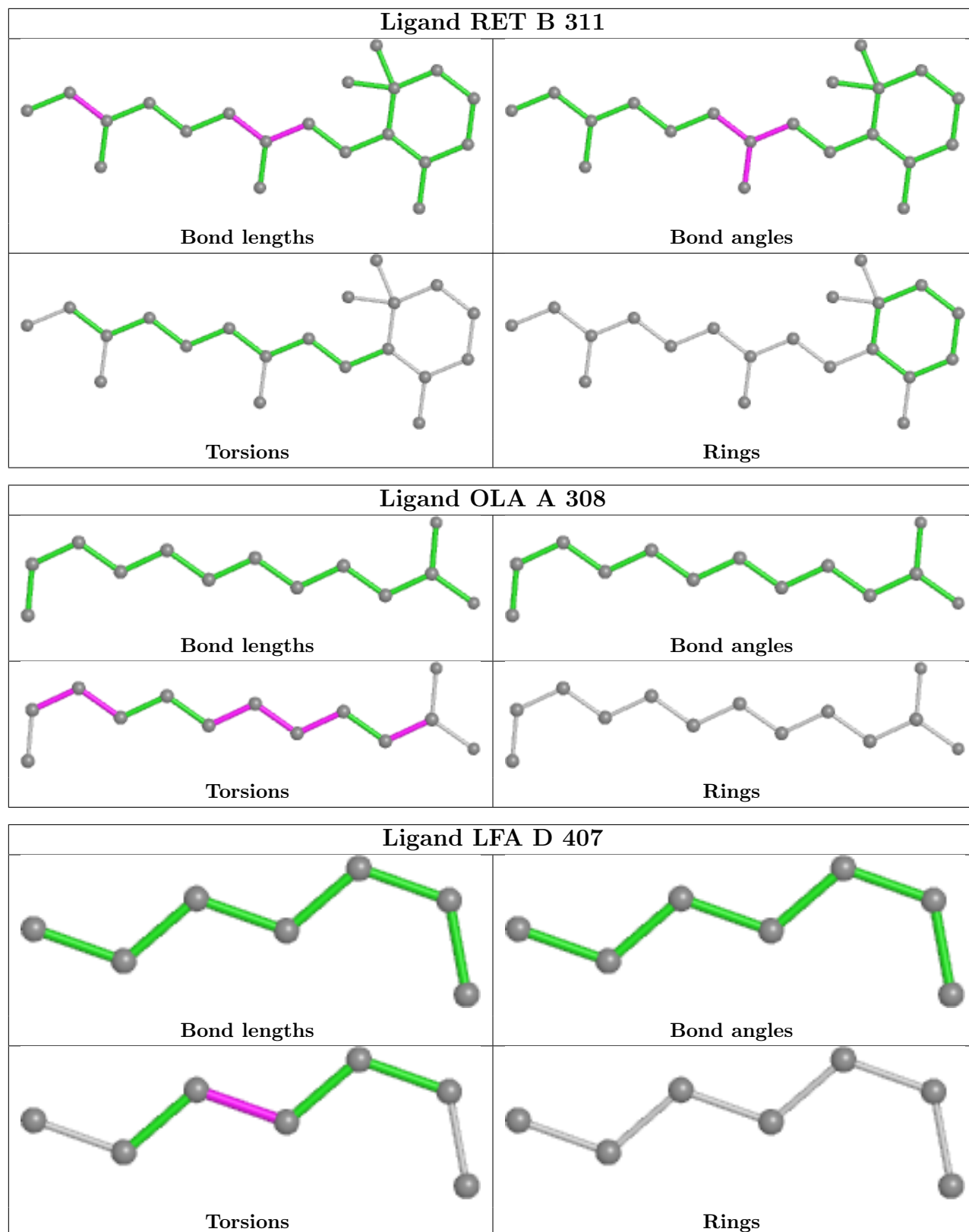
The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

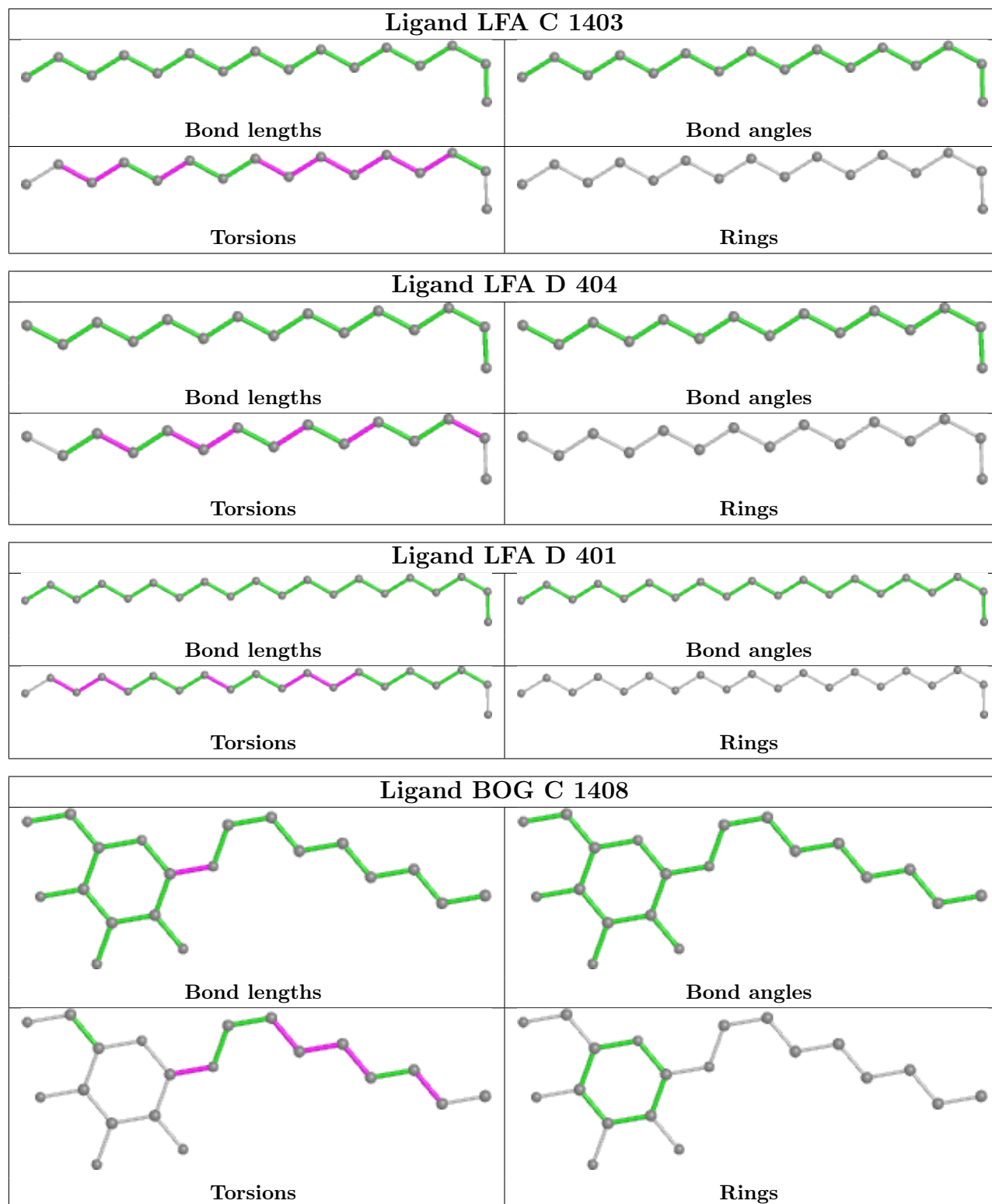


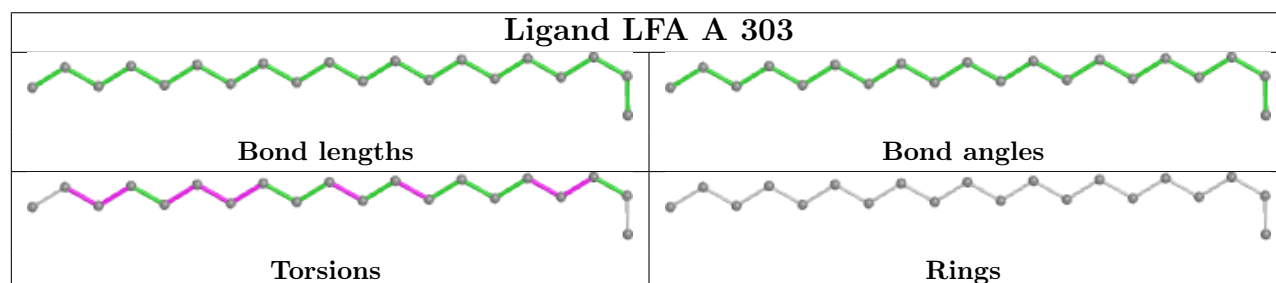
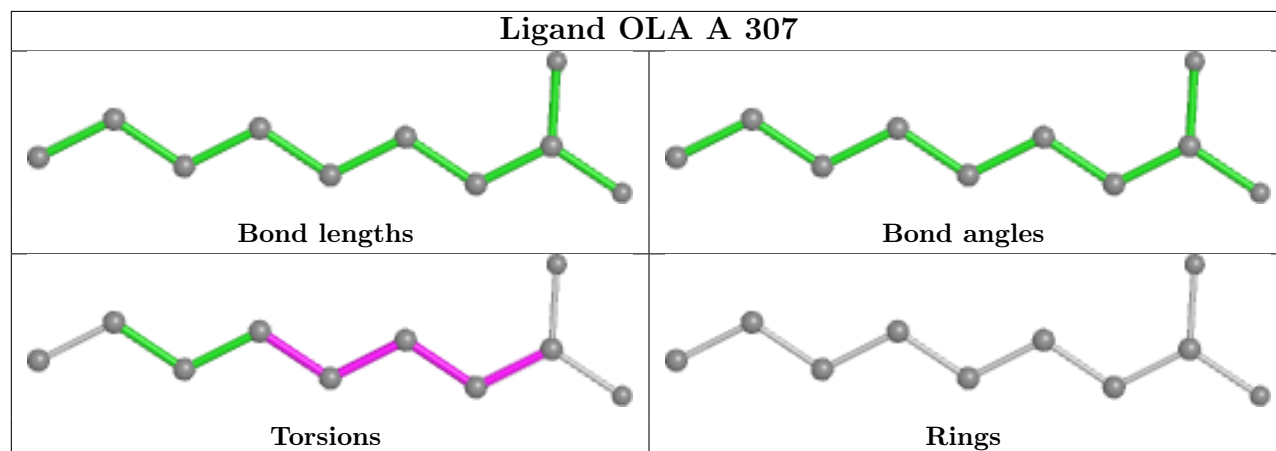
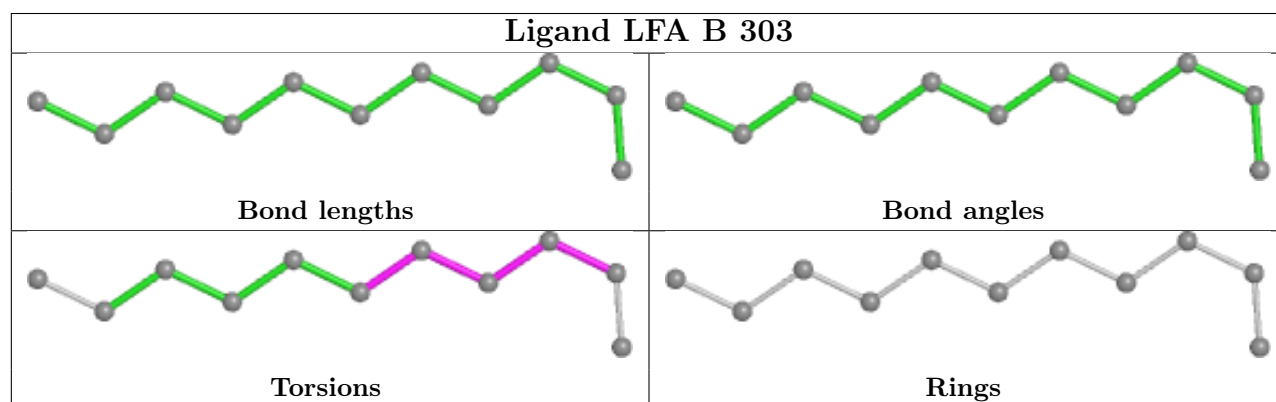
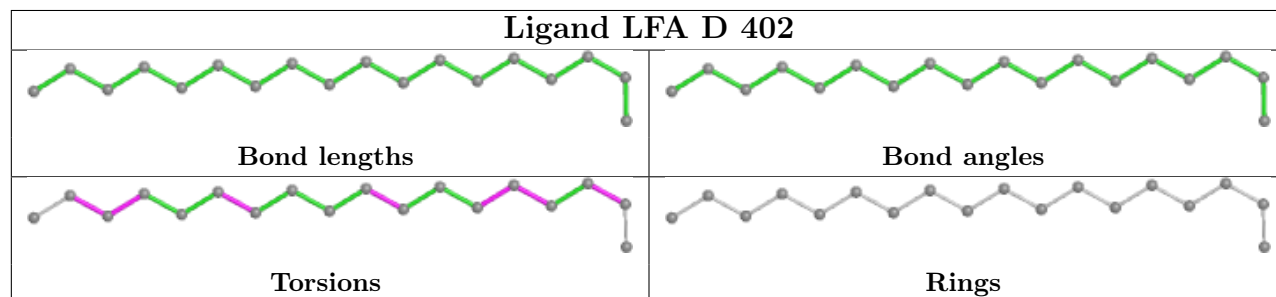
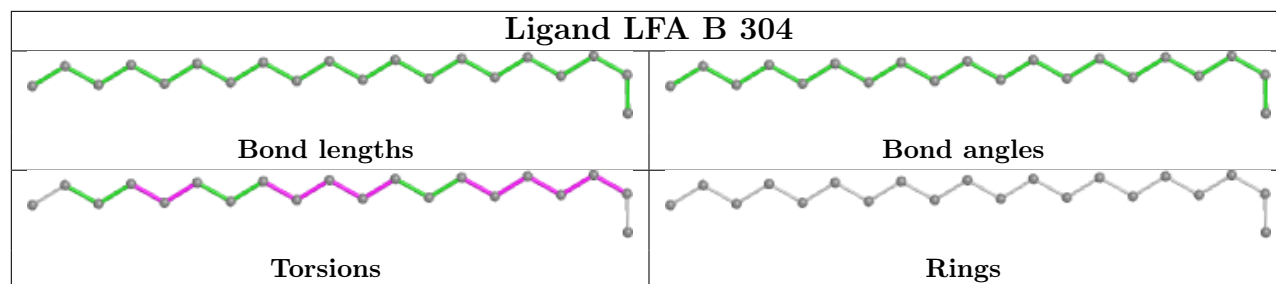


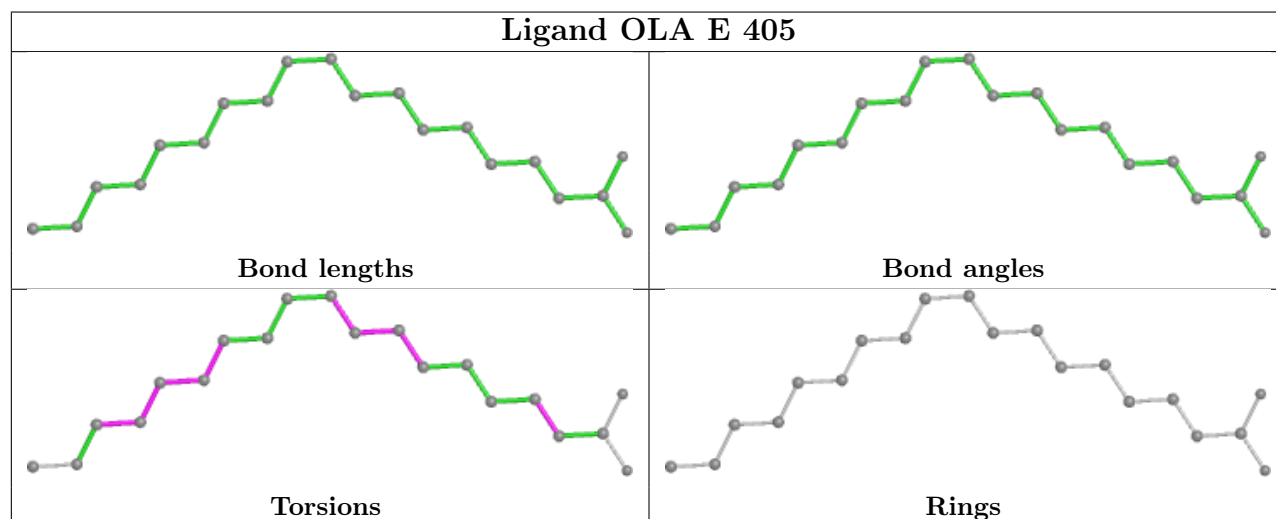
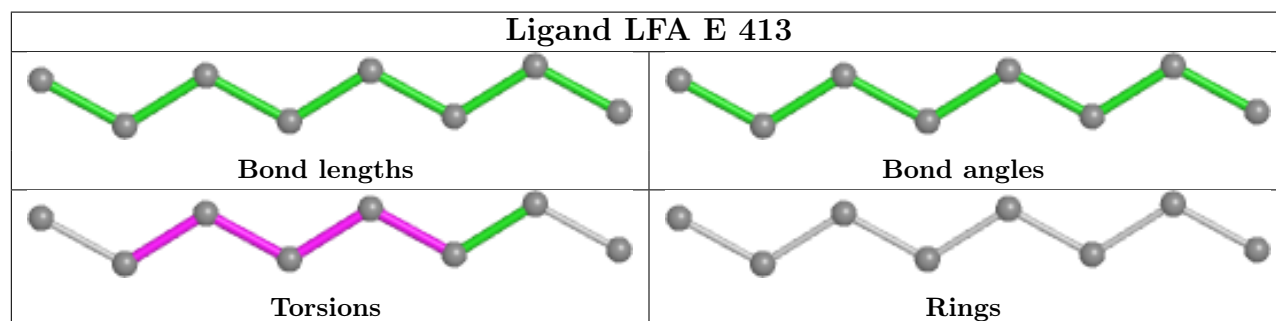
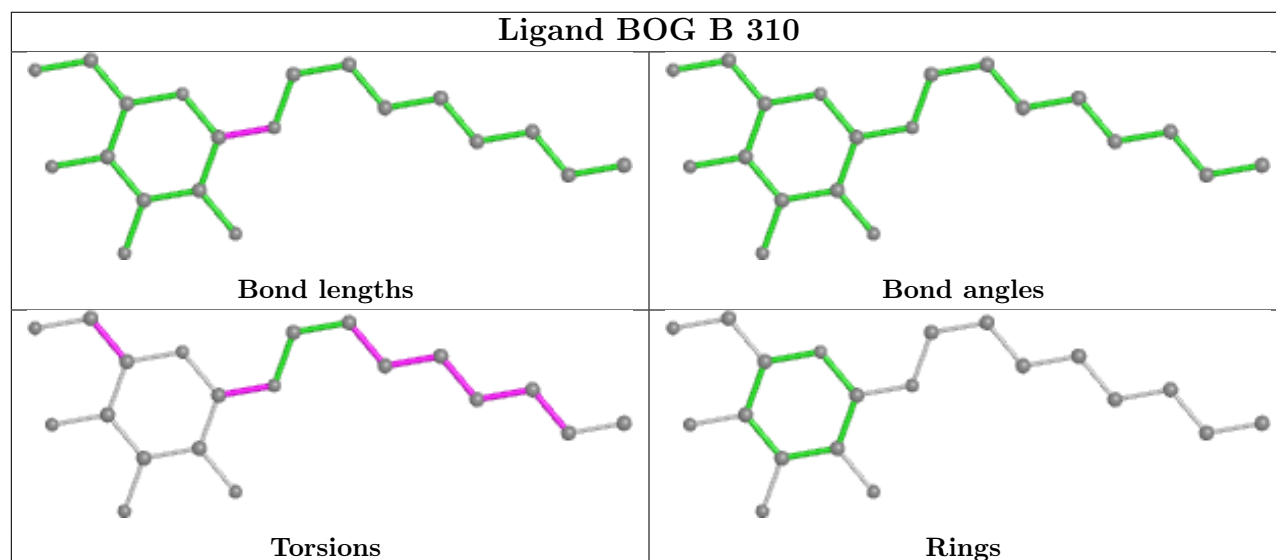
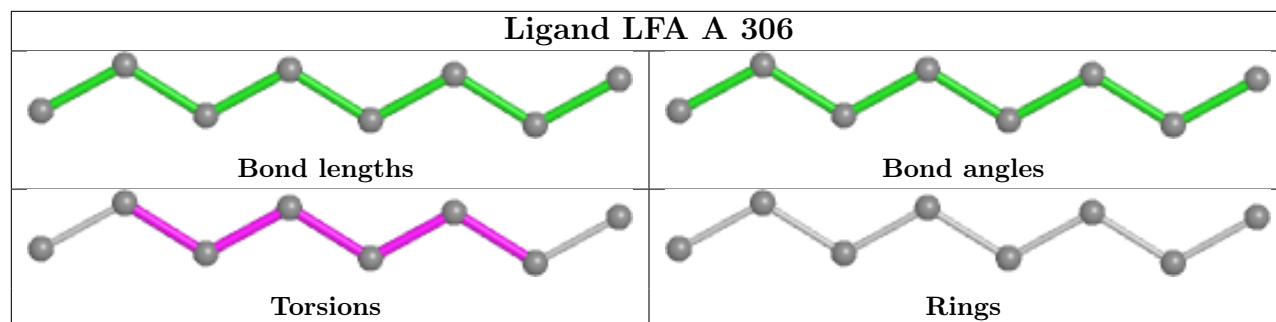


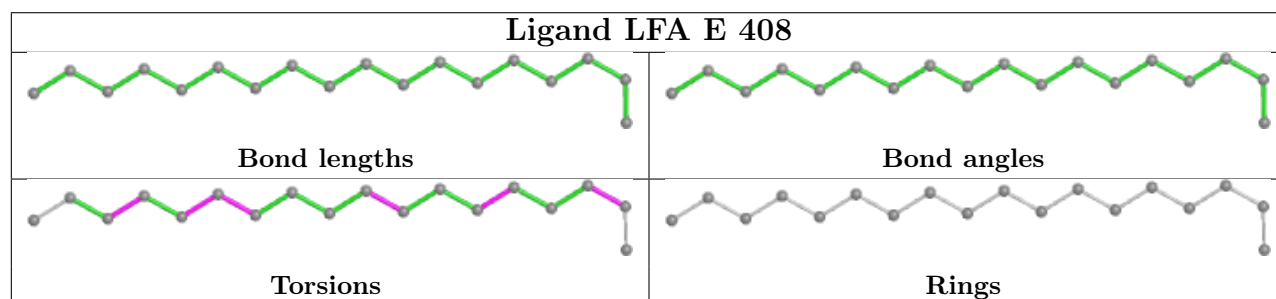
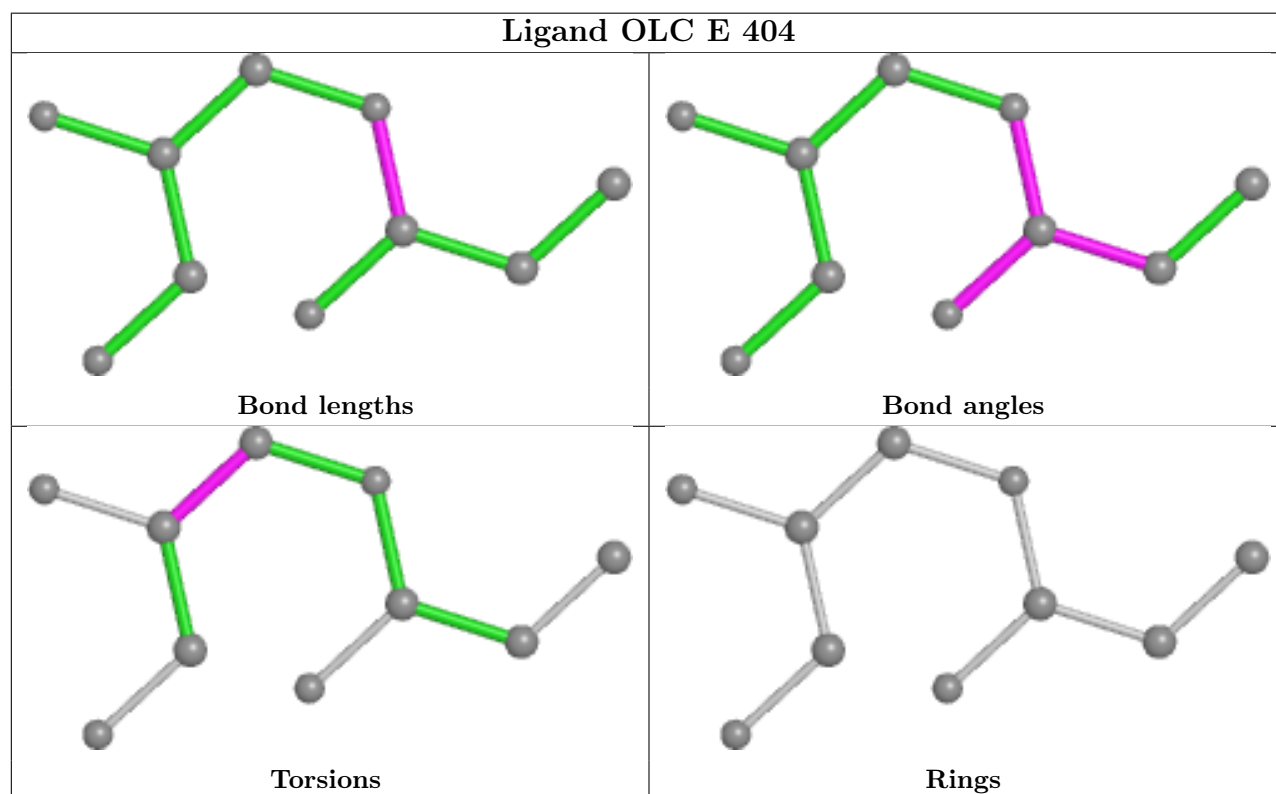
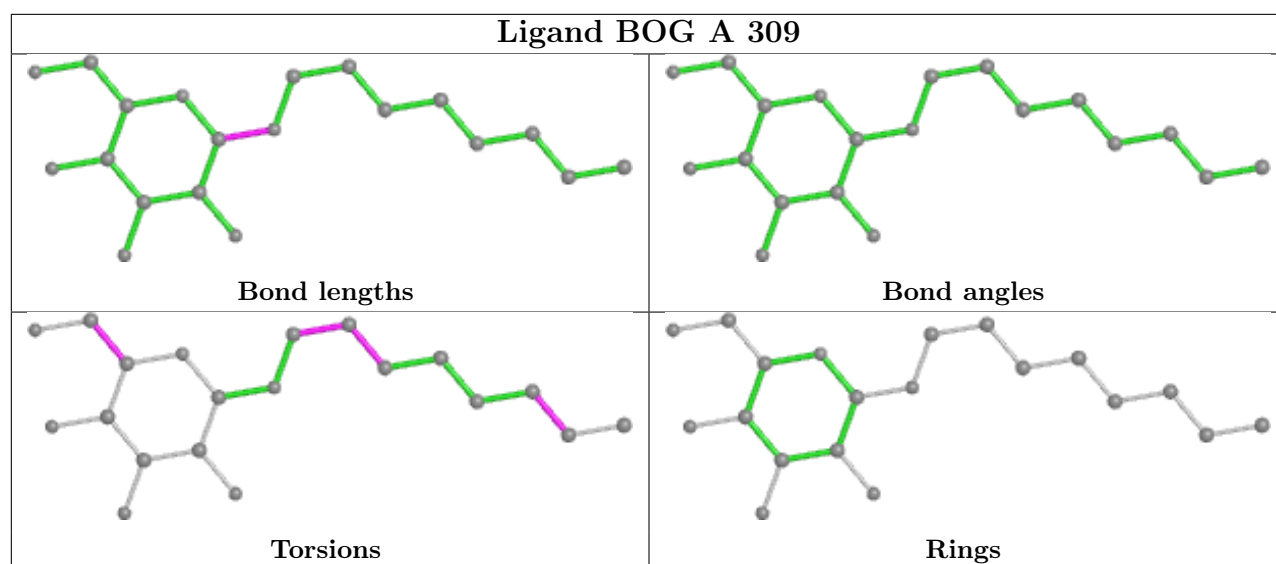


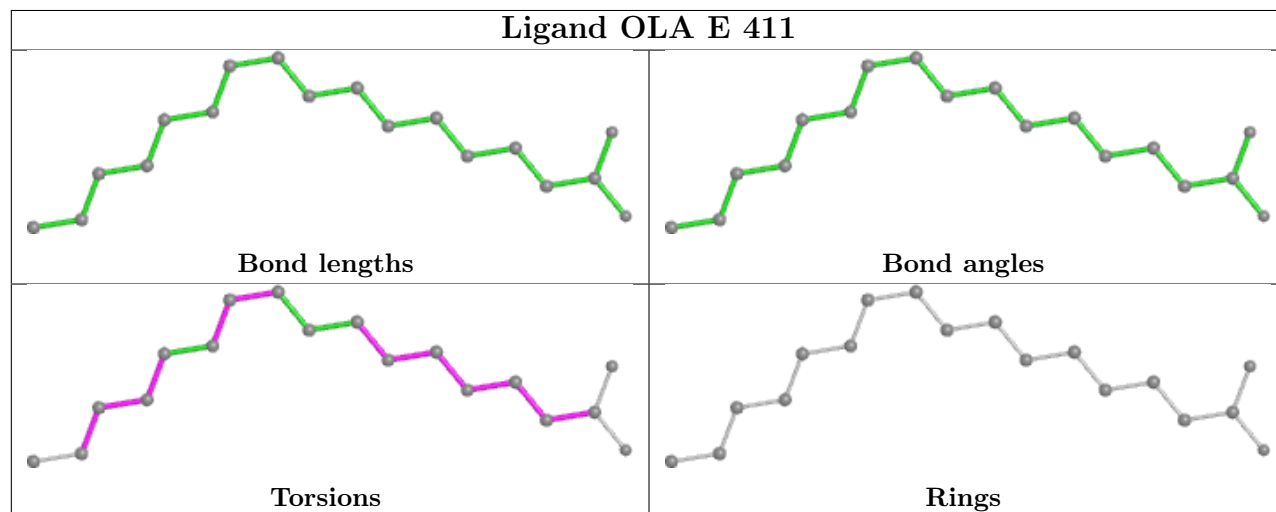
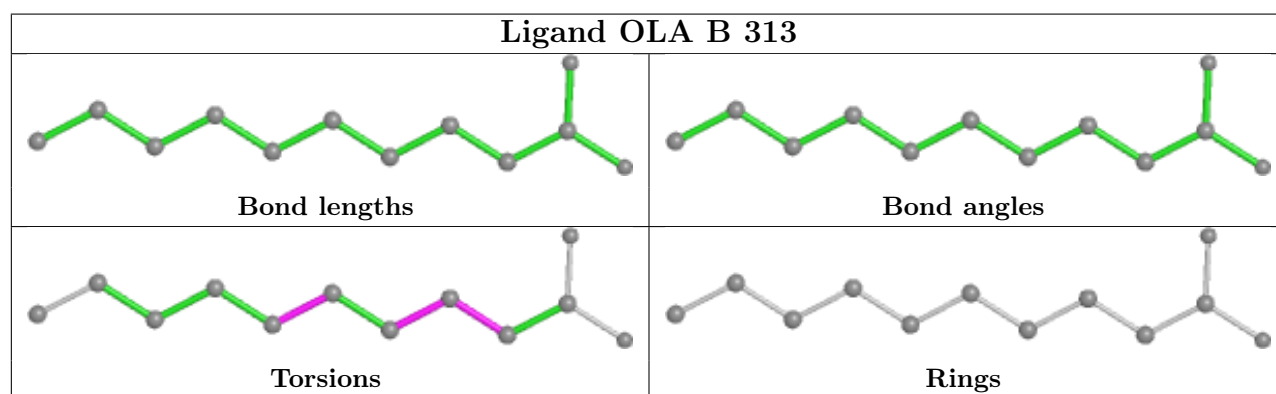
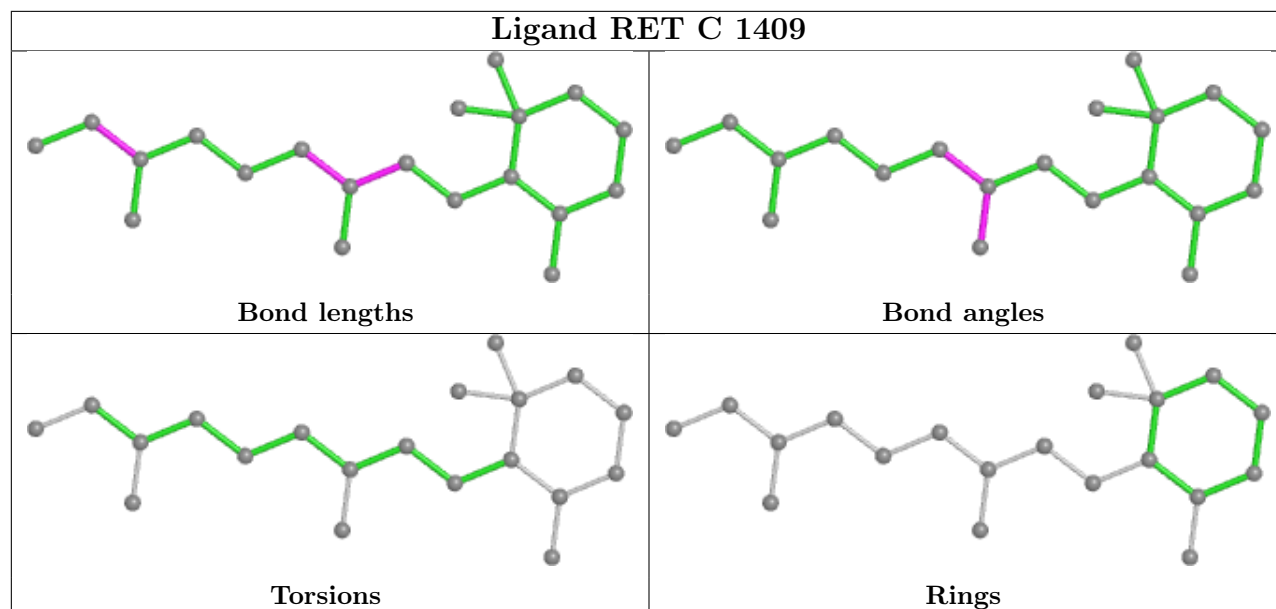


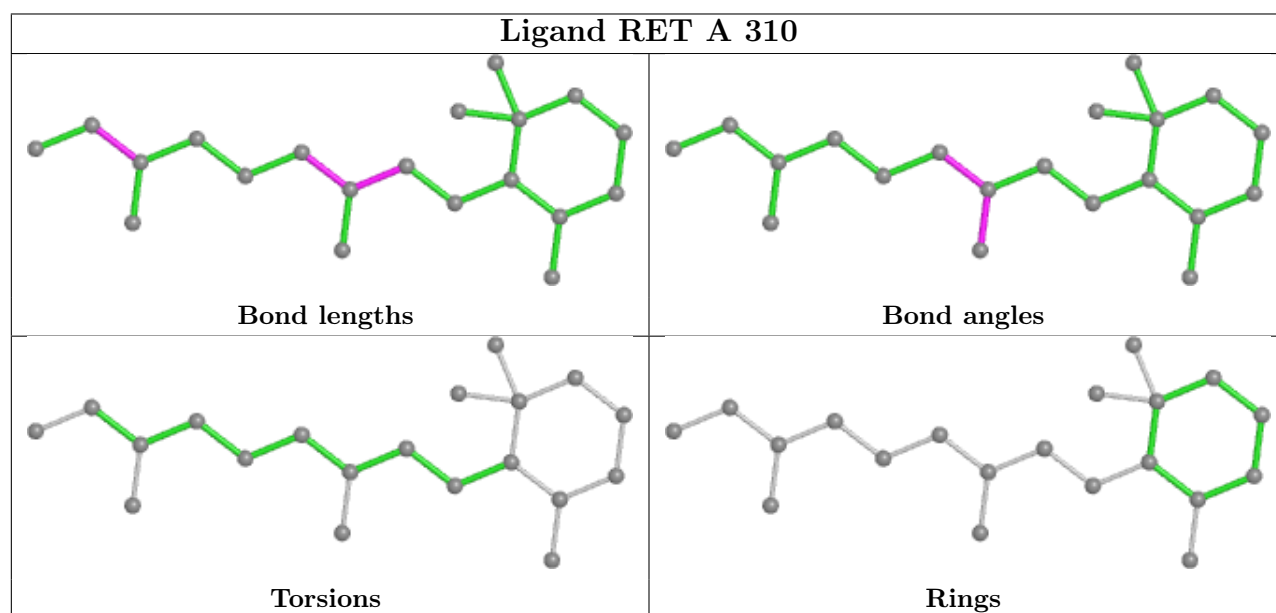
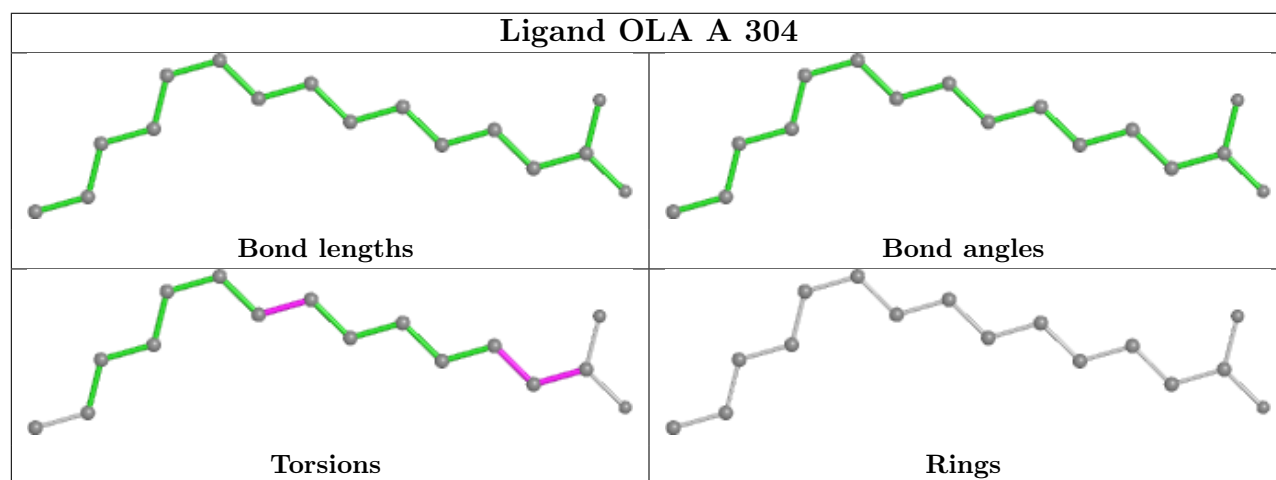
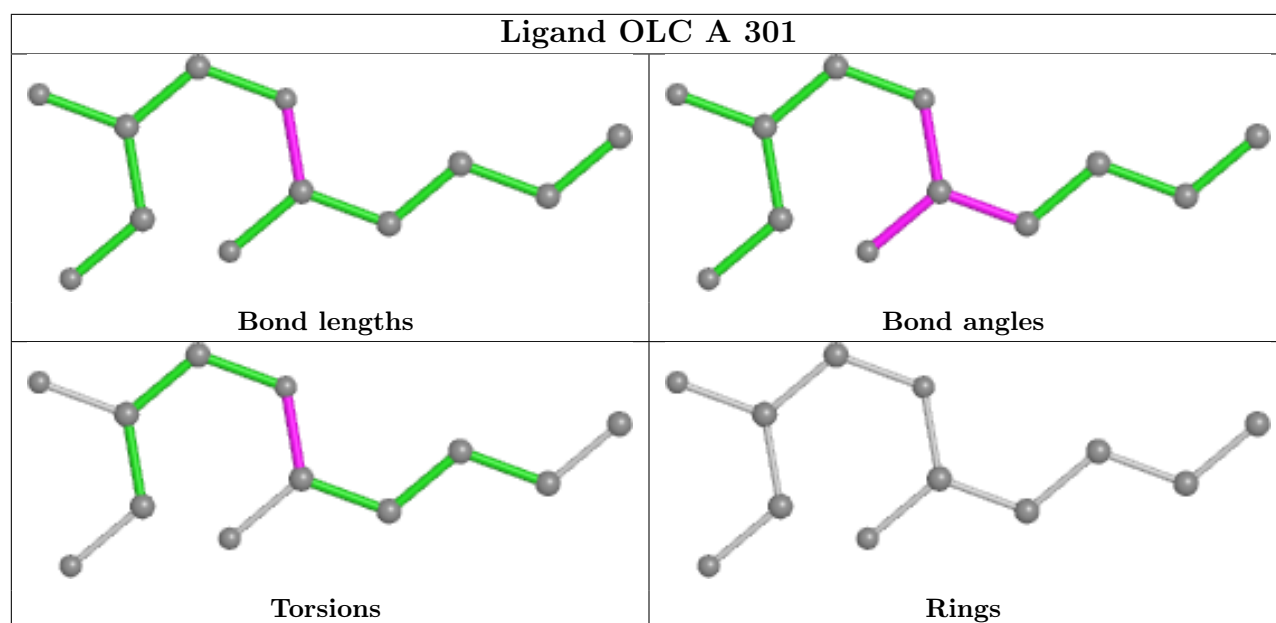


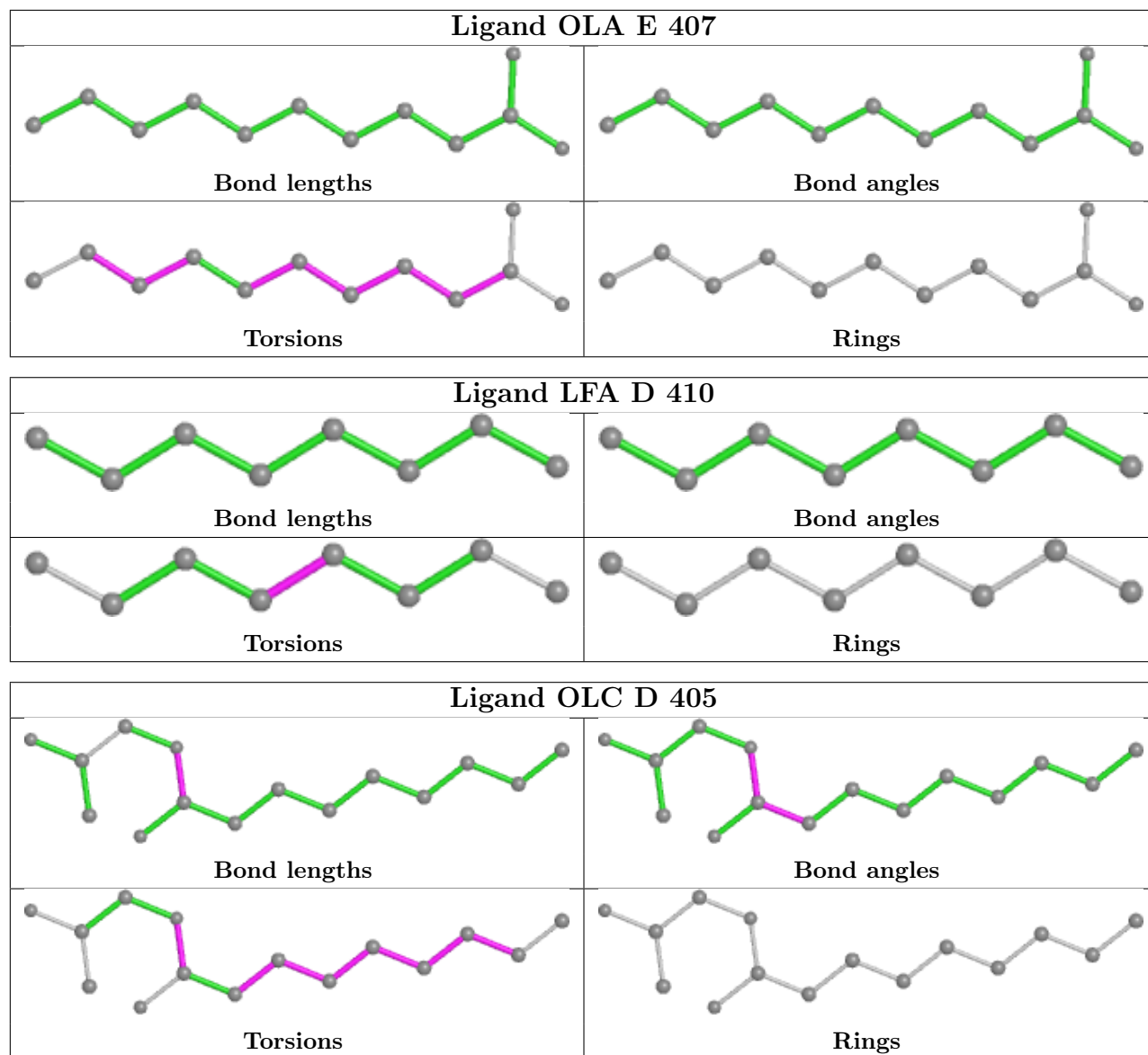


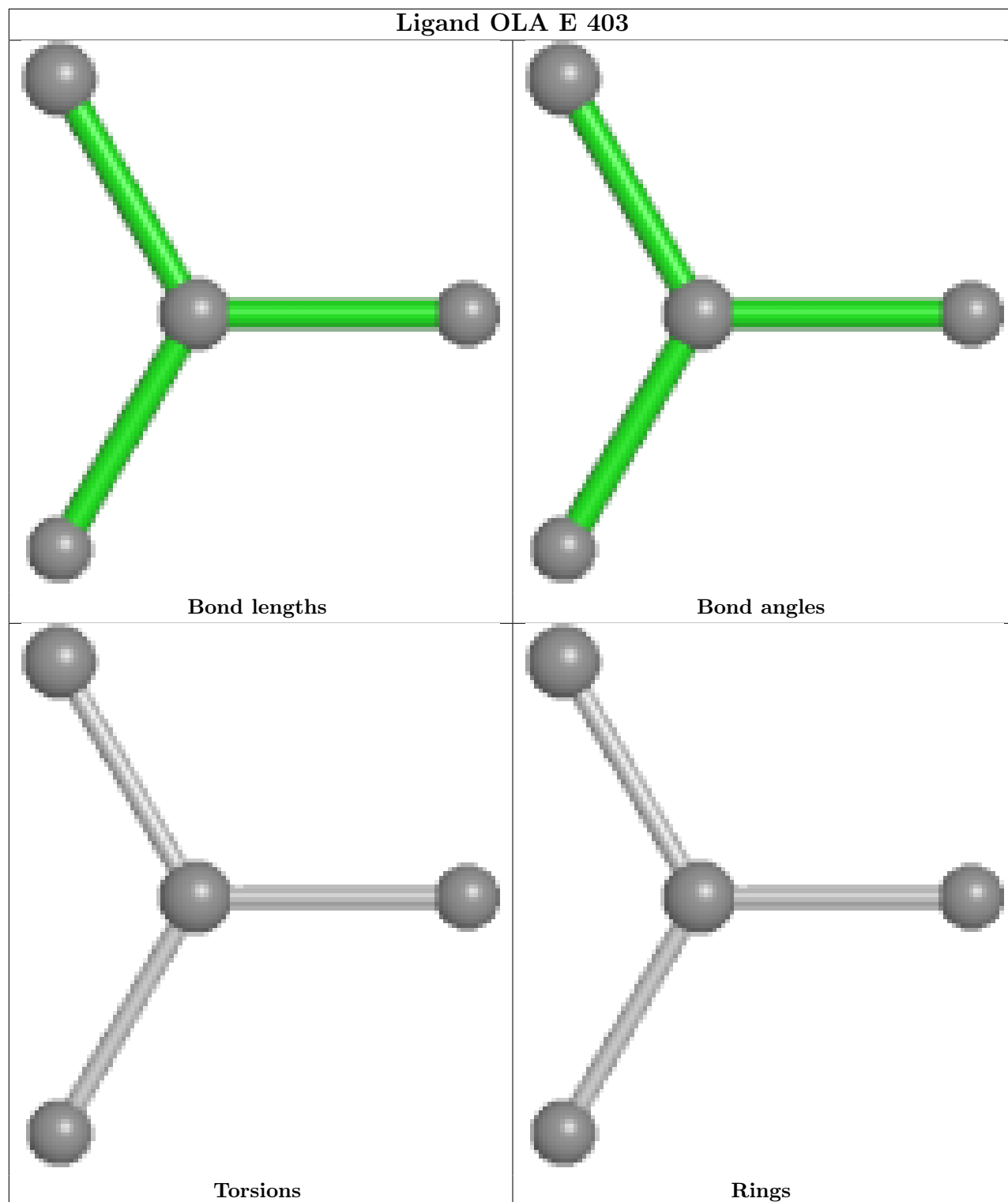


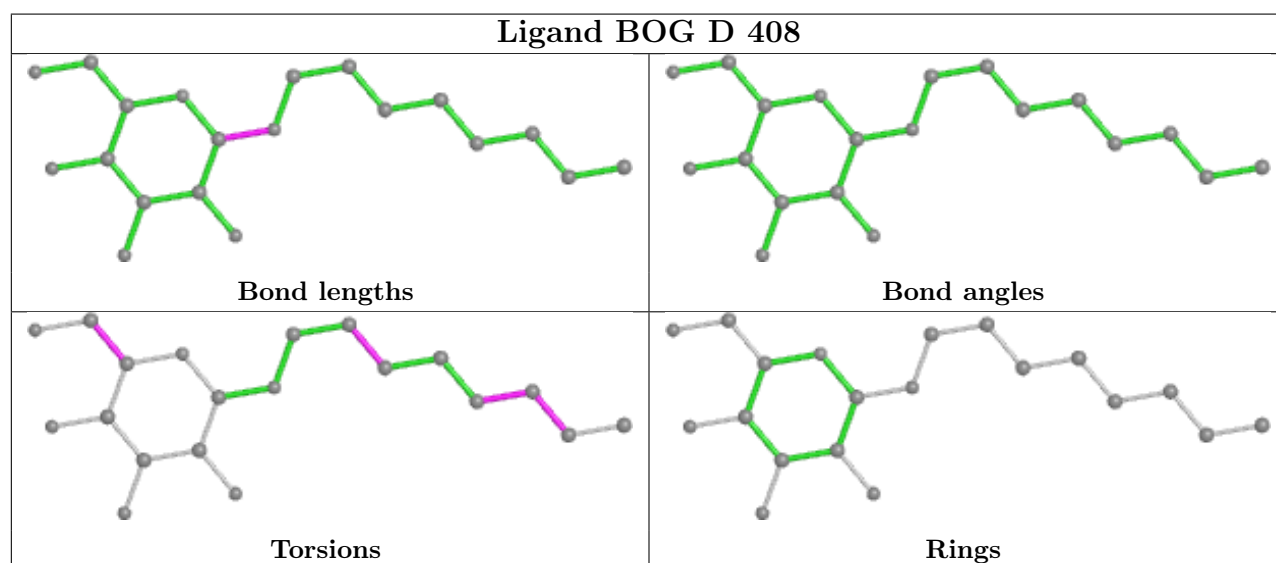
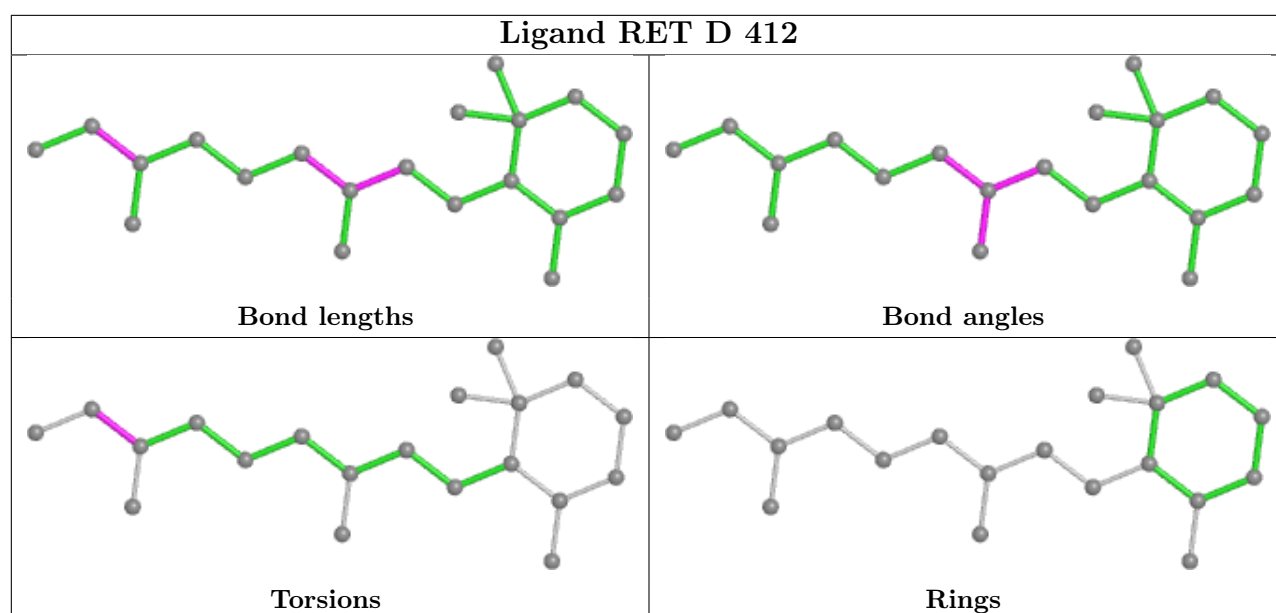
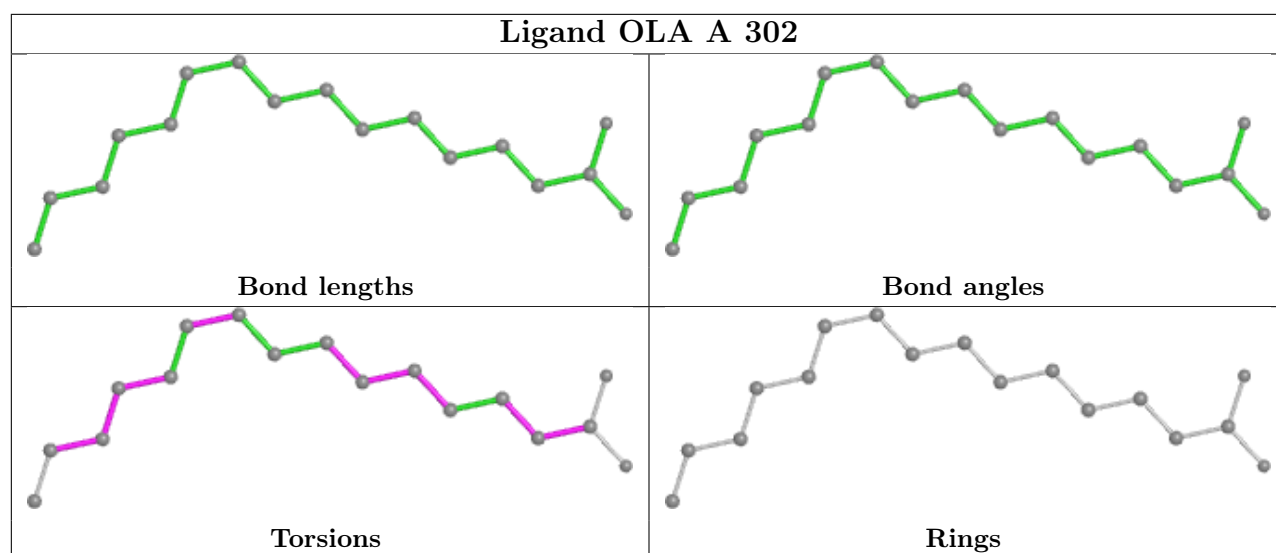


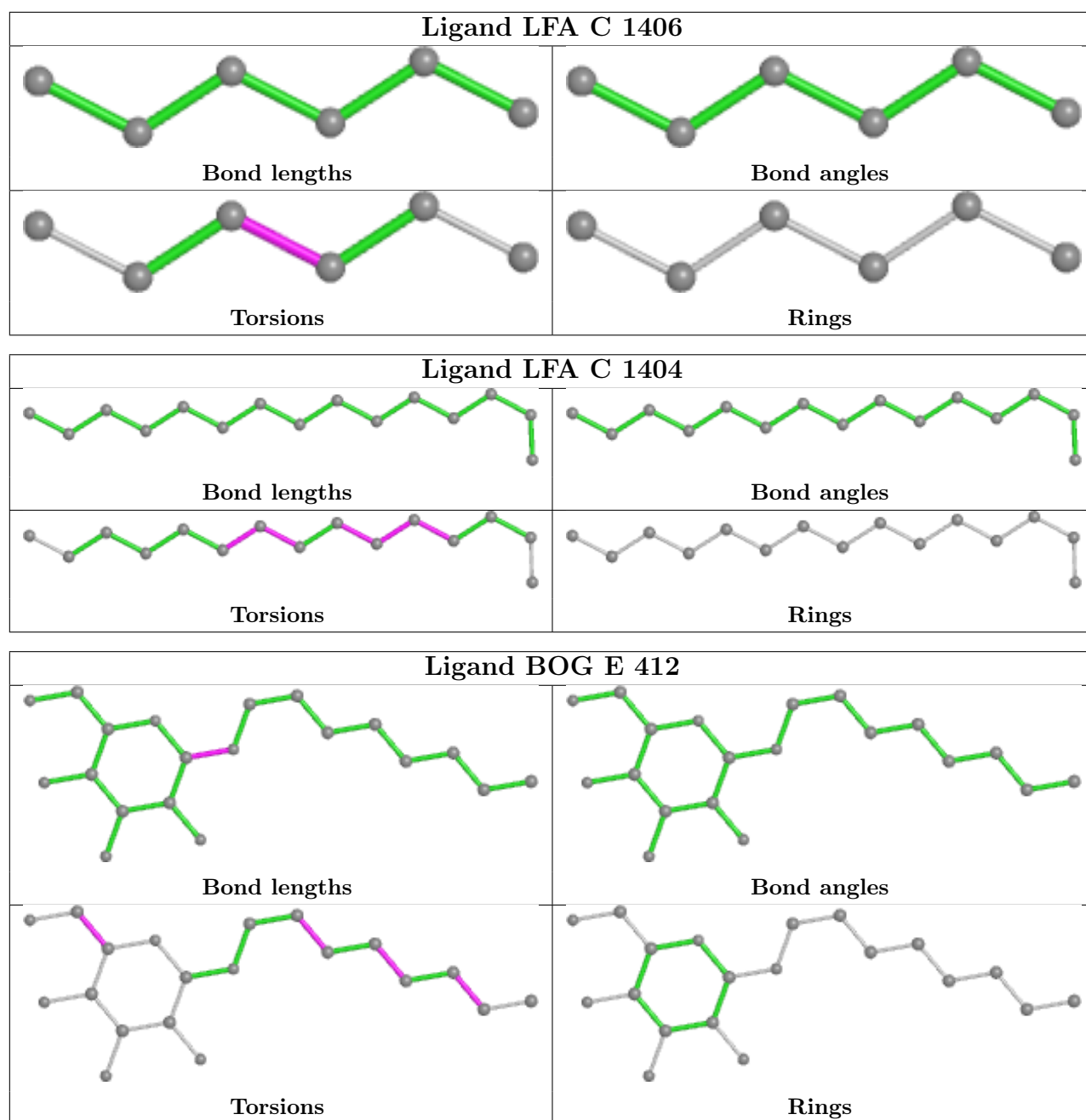












5.7 Other polymers [i](#)

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	A	269/288 (93%)	-0.09	12 (4%) 33 31	29, 40, 59, 82	0
1	B	269/288 (93%)	-0.04	9 (3%) 46 45	30, 41, 60, 86	0
1	C	268/288 (93%)	0.06	17 (6%) 20 18	30, 43, 64, 89	0
1	D	270/288 (93%)	-0.05	9 (3%) 46 45	29, 42, 62, 91	0
1	E	273/288 (94%)	0.01	12 (4%) 34 33	30, 43, 67, 130	0
All	All	1349/1440 (93%)	-0.02	59 (4%) 34 33	29, 42, 63, 130	0

The worst 5 of 59 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	274	GLU	7.7
1	E	230	VAL	4.8
1	C	230	VAL	4.1
1	D	230	VAL	4.0
1	A	73	LEU	3.8

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, 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.

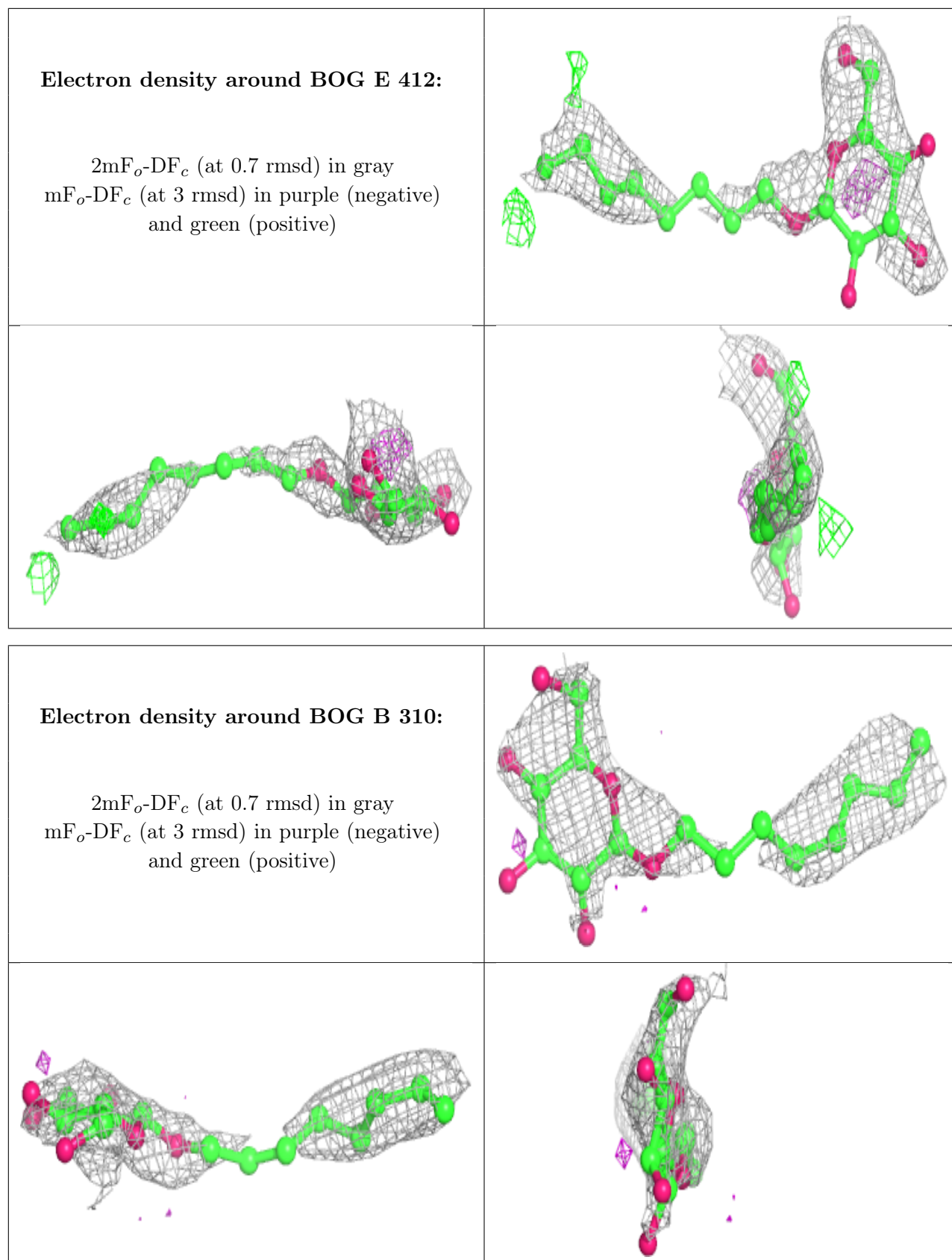
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	BOG	E	412	20/20	0.62	0.52	69,102,129,136	0
5	BOG	B	310	20/20	0.67	0.52	65,99,114,118	0
3	OLA	A	307	10/20	0.70	0.42	65,76,82,85	0
5	BOG	A	309	20/20	0.73	0.52	66,93,122,127	0
4	LFA	B	307	9/20	0.75	0.39	58,67,80,84	0
3	OLA	D	409	20/20	0.76	0.23	65,80,103,103	0
3	OLA	E	403	4/20	0.77	0.84	50,77,81,82	0
5	BOG	C	1408	20/20	0.77	0.51	71,94,116,116	0
5	BOG	D	408	20/20	0.78	0.53	65,81,109,120	0
4	LFA	D	404	15/20	0.80	0.29	65,74,88,89	0
4	LFA	D	407	7/20	0.81	0.21	59,62,67,68	0
4	LFA	E	410	5/20	0.82	0.16	57,63,72,73	0
4	LFA	D	403	10/20	0.82	0.17	59,65,69,72	0
3	OLA	E	407	12/20	0.83	0.17	72,79,85,86	0
3	OLA	A	308	13/20	0.84	0.24	55,65,84,93	0
3	OLA	B	309	16/20	0.84	0.22	64,76,84,88	0
4	LFA	A	306	8/20	0.84	0.24	55,64,76,77	0
3	OLA	A	304	16/20	0.84	0.19	59,79,88,89	0
4	LFA	C	1406	6/20	0.84	0.20	58,61,68,72	0
3	OLA	A	311	5/20	0.85	0.81	65,73,82,93	0
4	LFA	E	413	8/20	0.85	0.25	62,69,79,84	0
3	OLA	E	405	20/20	0.85	0.28	51,58,87,98	0
3	OLA	B	305	14/20	0.85	0.28	72,79,88,89	0
3	OLA	E	411	18/20	0.85	0.30	58,74,92,100	0
4	LFA	A	305	5/20	0.85	0.22	59,59,69,70	0
4	LFA	D	410	8/20	0.85	0.21	55,63,73,74	0
4	LFA	D	411	7/20	0.85	0.31	56,62,64,69	0
3	OLA	D	406	4/20	0.86	0.89	67,78,81,96	0
4	LFA	B	303	11/20	0.86	0.26	52,59,70,71	0
4	LFA	D	401	20/20	0.87	0.75	54,65,76,77	0
3	OLA	B	313	12/20	0.87	0.31	64,73,84,88	0
3	OLA	C	1401	4/20	0.87	0.67	52,73,75,86	0
3	OLA	B	308	4/20	0.87	0.98	62,78,80,92	0
4	LFA	C	1403	16/20	0.87	0.33	50,54,61,62	0
4	LFA	C	1404	15/20	0.87	0.25	57,71,87,89	0
4	LFA	E	406	19/20	0.88	0.80	51,69,76,82	0
4	LFA	E	402	20/20	0.89	0.77	54,67,83,87	0
4	LFA	E	401	5/20	0.89	0.19	48,54,59,63	0
4	LFA	C	1405	3/20	0.89	0.23	54,54,57,59	0
4	LFA	E	408	18/20	0.90	0.19	62,72,81,83	0

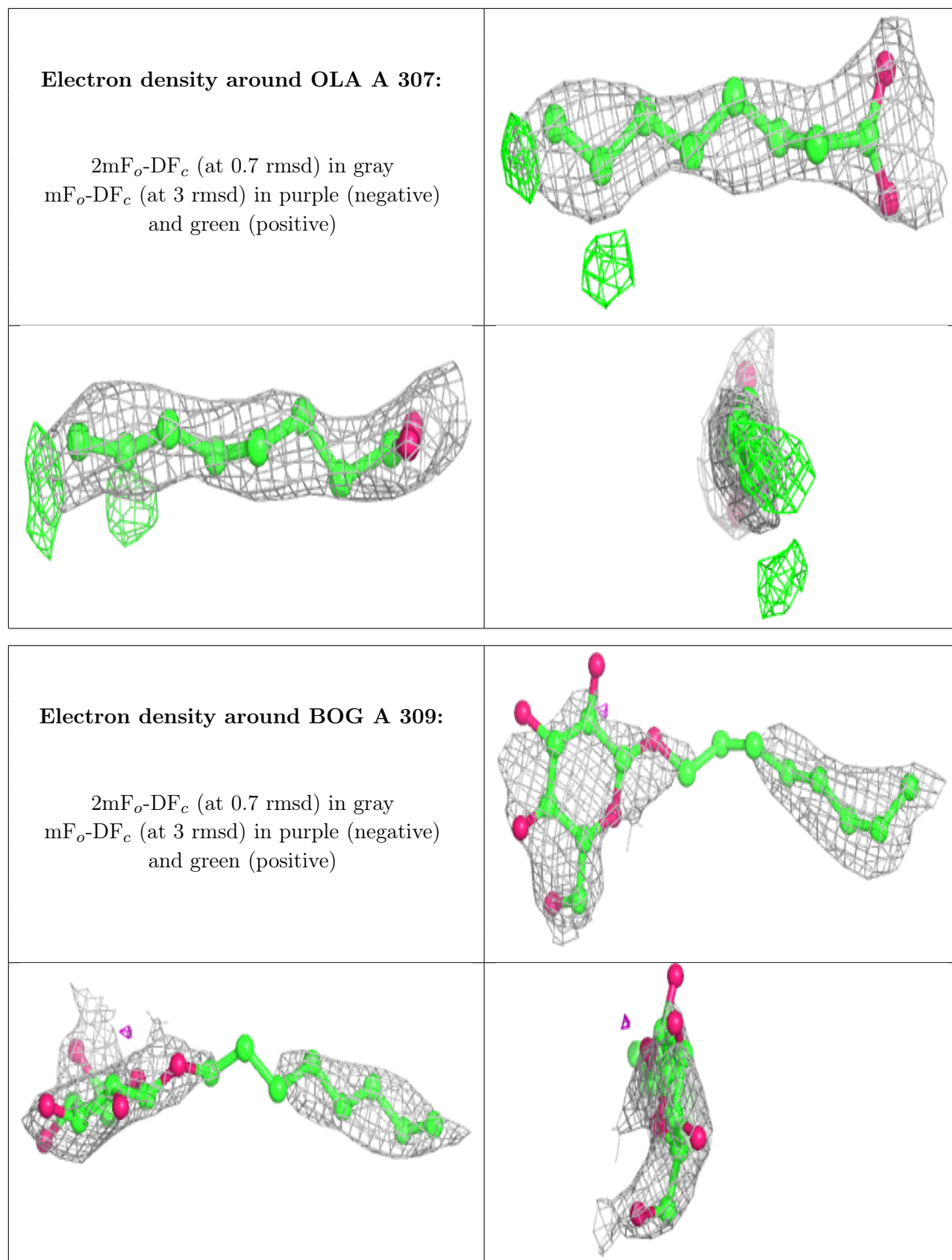
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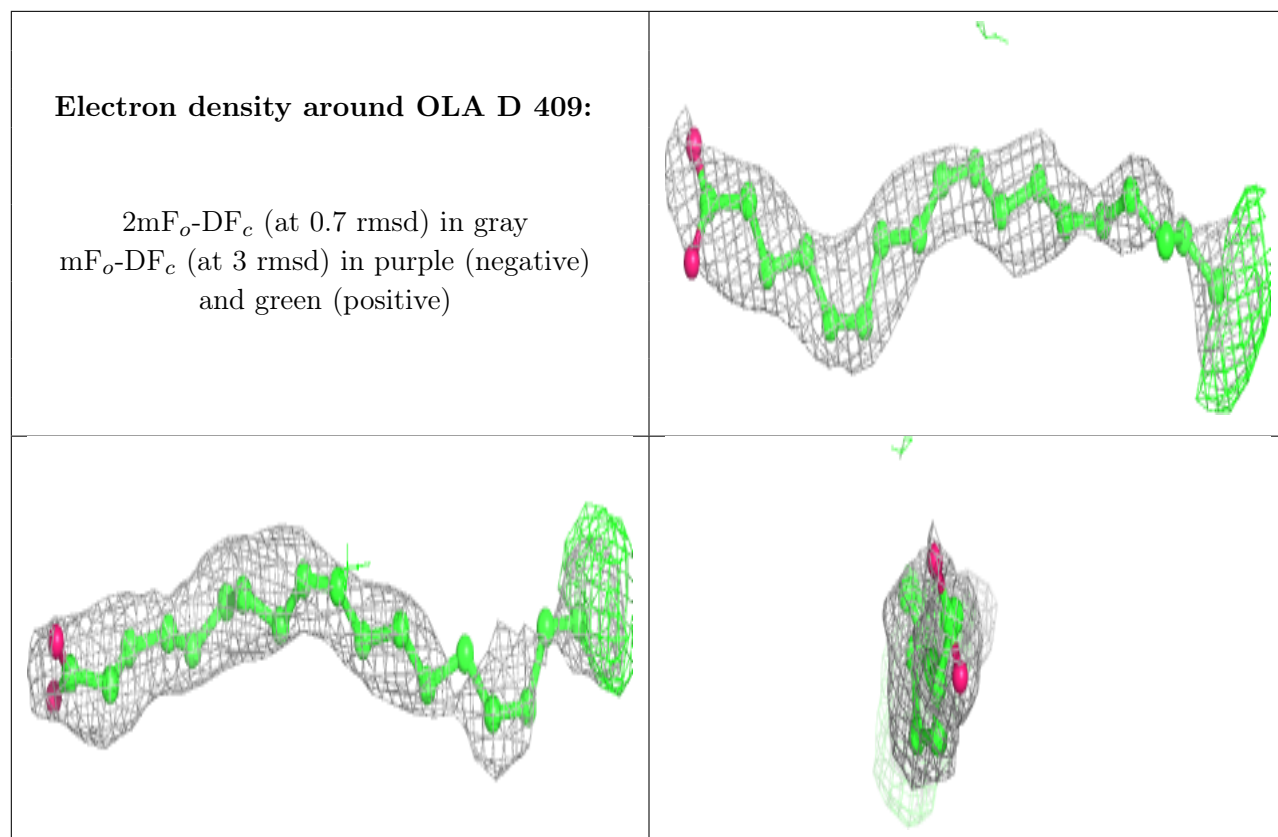
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	LFA	B	312	20/20	0.90	0.70	56,68,77,77	0
2	OLC	C	1402	11/25	0.90	0.19	57,67,70,73	0
4	LFA	B	302	19/20	0.90	0.30	47,62,70,79	0
3	OLA	A	302	17/20	0.90	0.20	45,54,70,72	0
2	OLC	E	404	10/25	0.90	0.23	64,72,78,82	0
4	LFA	C	1407	5/20	0.91	0.16	47,61,63,64	0
4	LFA	A	303	20/20	0.91	0.22	60,73,90,94	0
4	LFA	E	409	5/20	0.91	0.24	60,63,65,69	0
2	OLC	B	301	13/25	0.93	0.22	63,68,78,83	0
2	OLC	D	405	15/25	0.93	0.19	61,74,78,80	0
6	RET	A	310	20/21	0.93	0.12	34,44,49,49	0
4	LFA	D	402	18/20	0.94	0.25	46,56,66,69	0
4	LFA	B	304	20/20	0.94	0.71	61,73,76,78	0
2	OLC	A	301	12/25	0.94	0.23	58,66,74,78	0
4	LFA	B	306	6/20	0.95	0.13	50,64,71,74	0
6	RET	D	412	20/21	0.95	0.13	36,43,46,47	0
6	RET	E	414	20/21	0.95	0.12	39,44,48,52	0
6	RET	C	1409	20/21	0.95	0.12	38,42,51,52	0
6	RET	B	311	20/21	0.96	0.14	34,39,42,44	0
7	NA	B	314	1/1	0.98	0.06	33,33,33,33	0
7	NA	E	415	1/1	0.98	0.07	36,36,36,36	0
7	NA	D	413	1/1	0.99	0.05	31,31,31,31	0
7	NA	A	312	1/1	0.99	0.07	32,32,32,32	0
7	NA	C	1410	1/1	0.99	0.06	33,33,33,33	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.

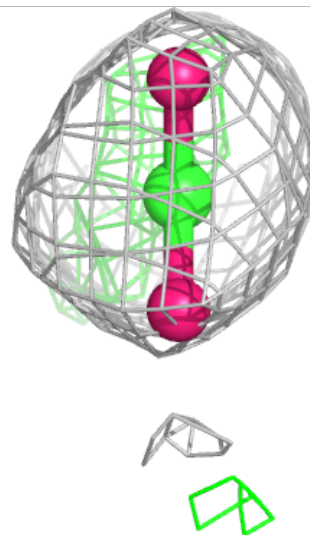
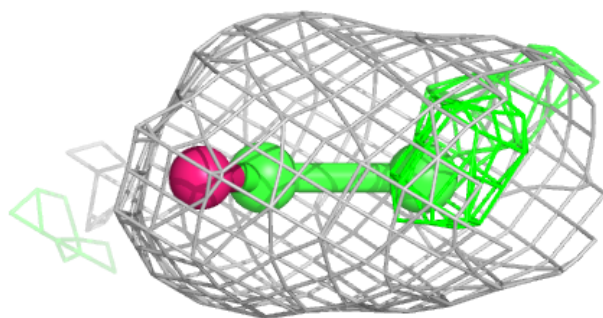
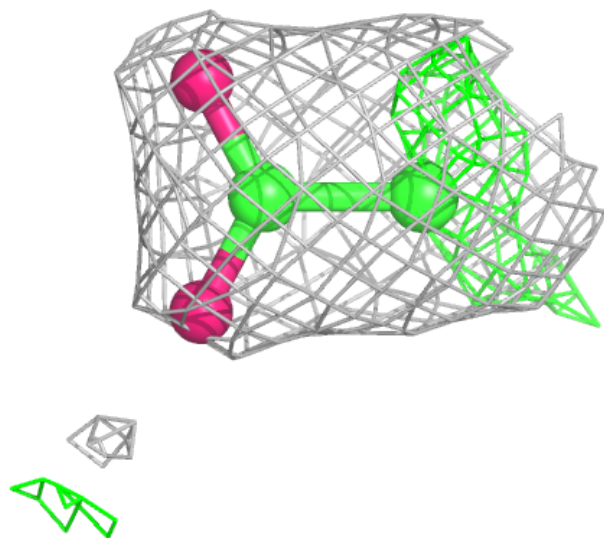






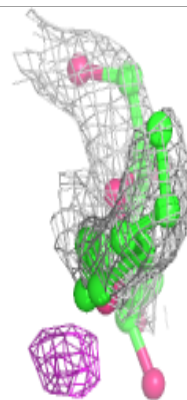
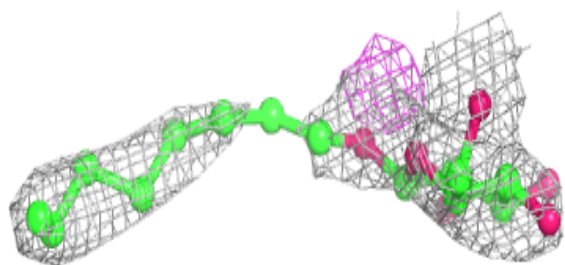
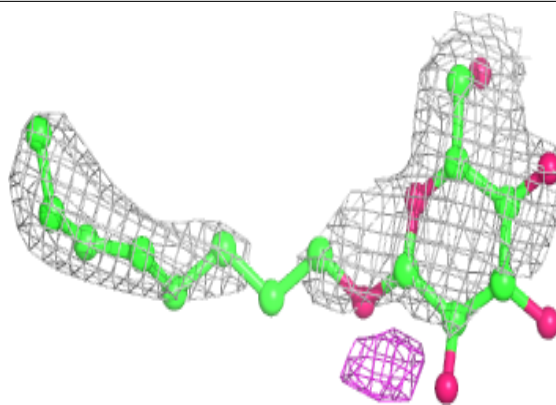
Electron density around OLA E 403:

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

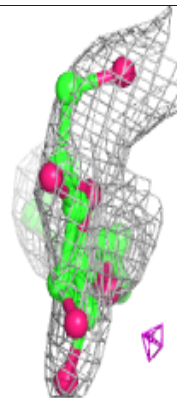
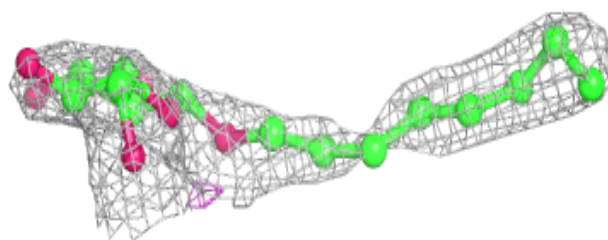
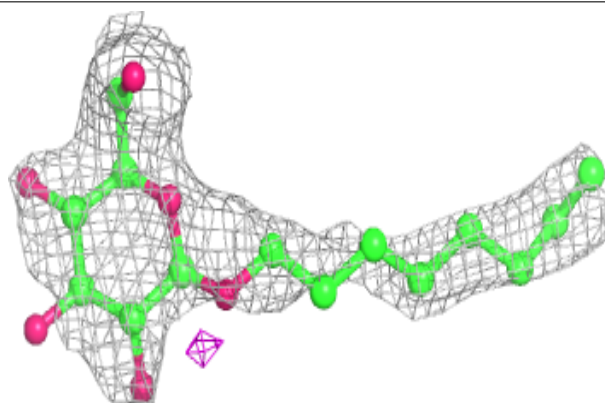


Electron density around BOG C 1408:

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

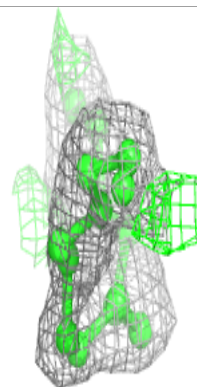
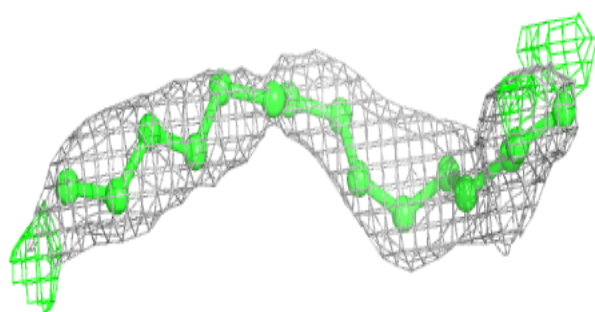
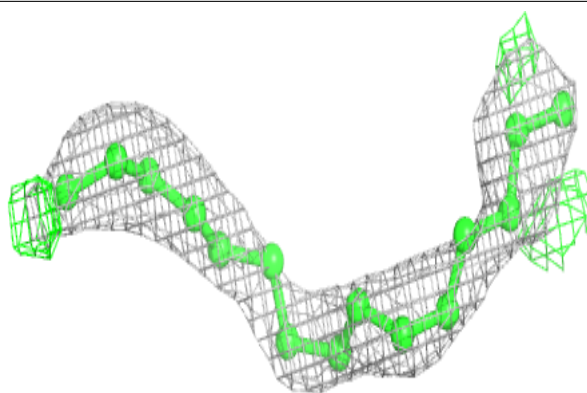
**Electron density around BOG D 408:**

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

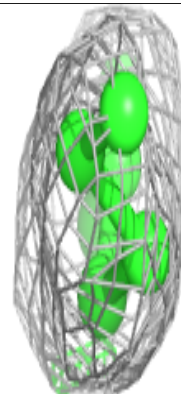
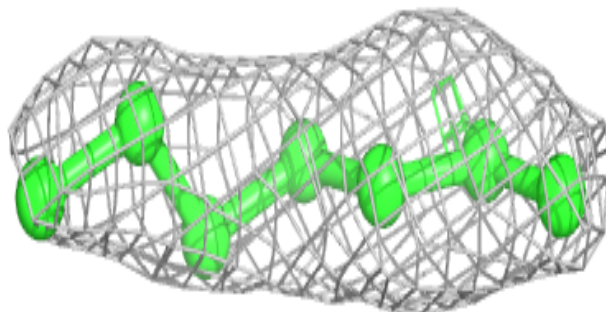
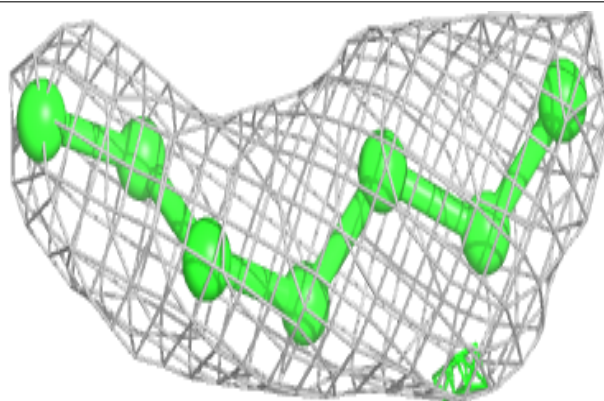


Electron density around LFA D 404:

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

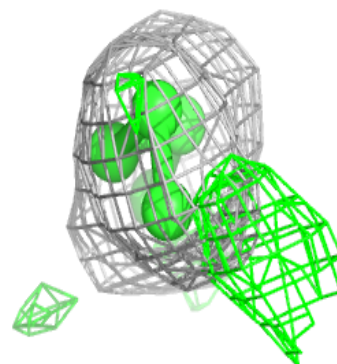
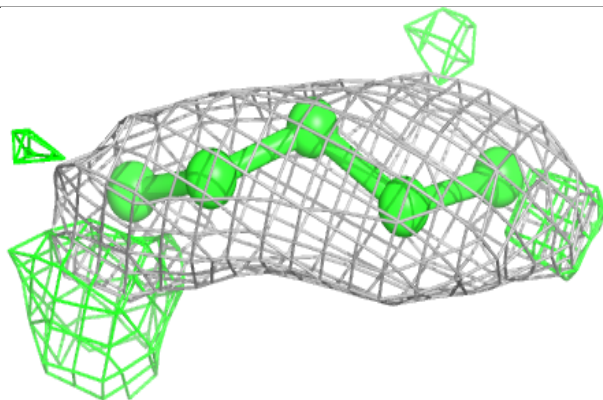
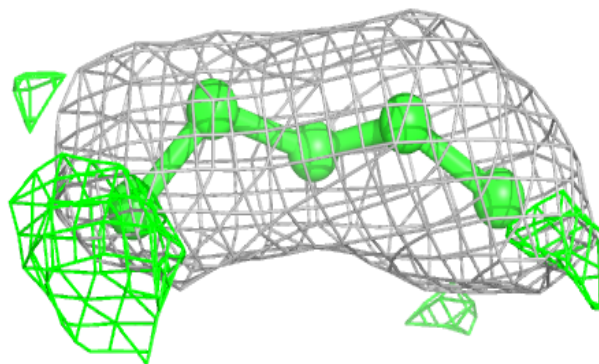
**Electron density around LFA D 407:**

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

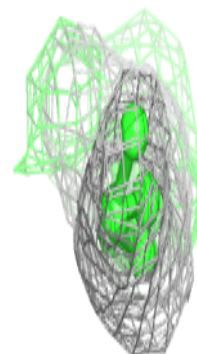
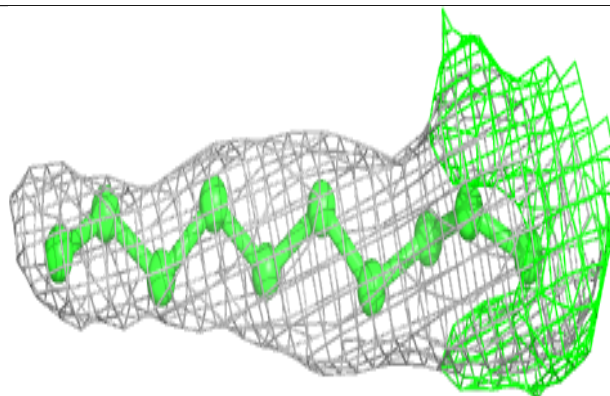
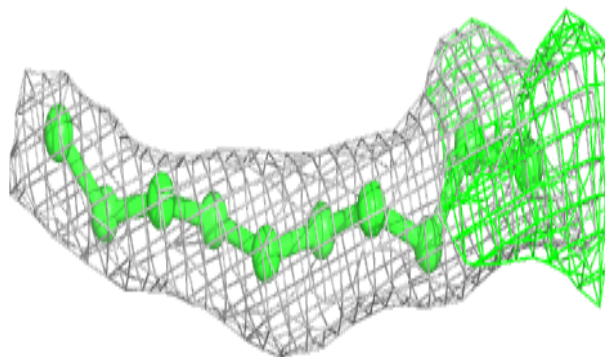


Electron density around LFA E 410:

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

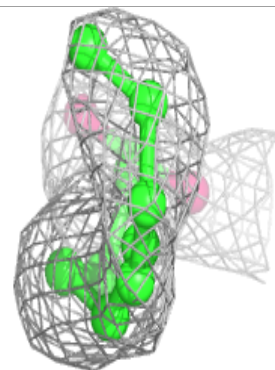
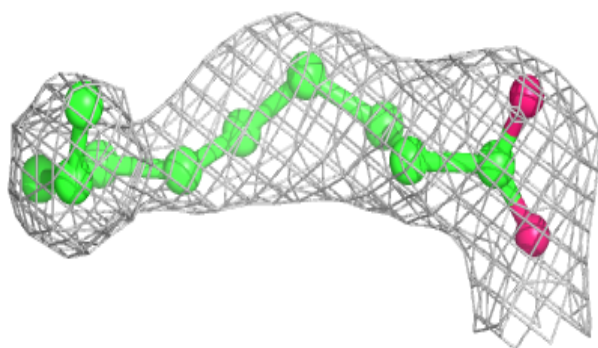
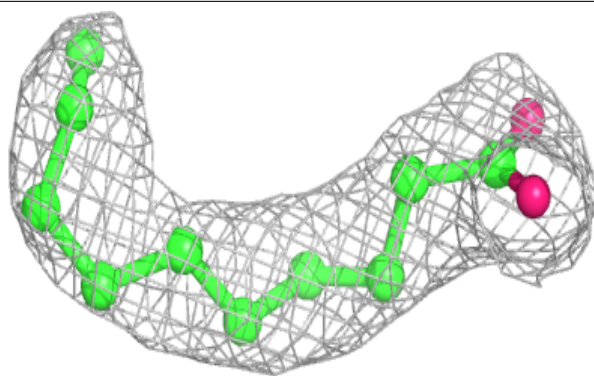
**Electron density around LFA D 403:**

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

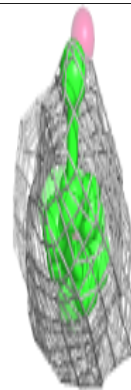
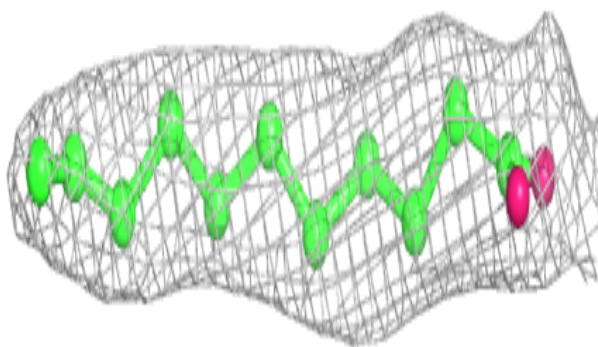
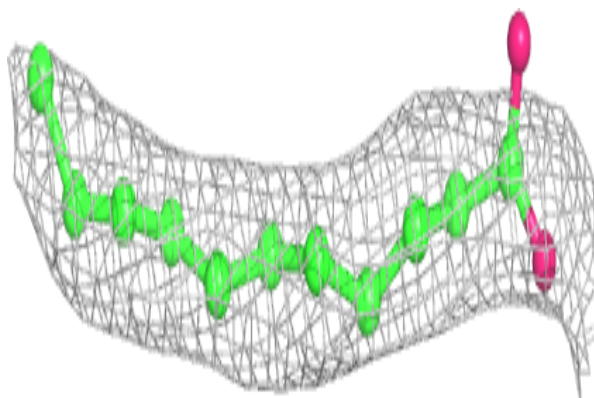


Electron density around OLA E 407:

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

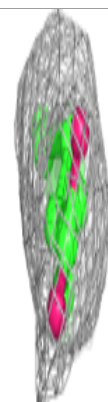
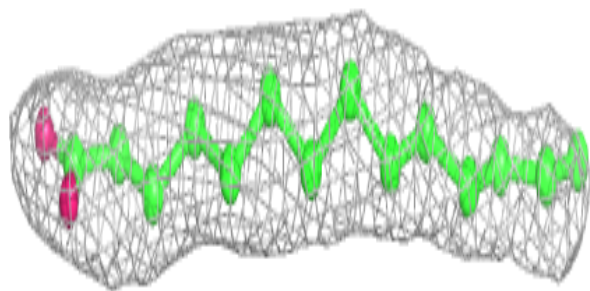
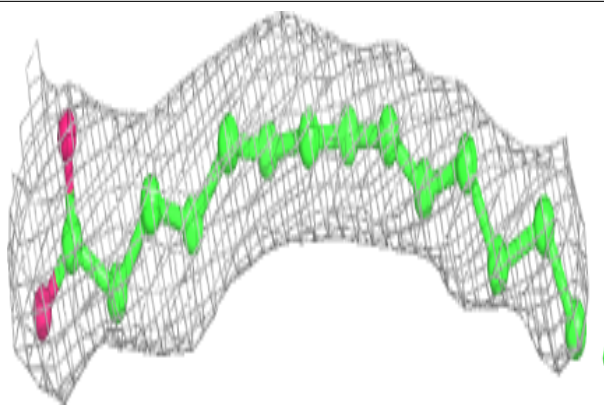
**Electron density around OLA A 308:**

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

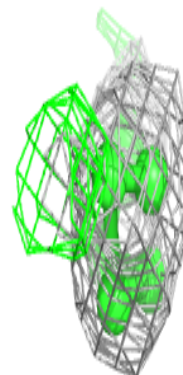
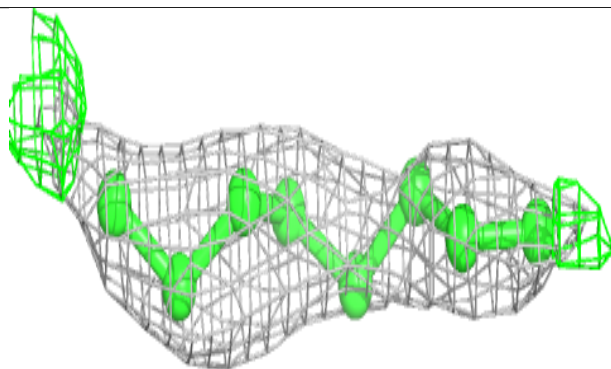
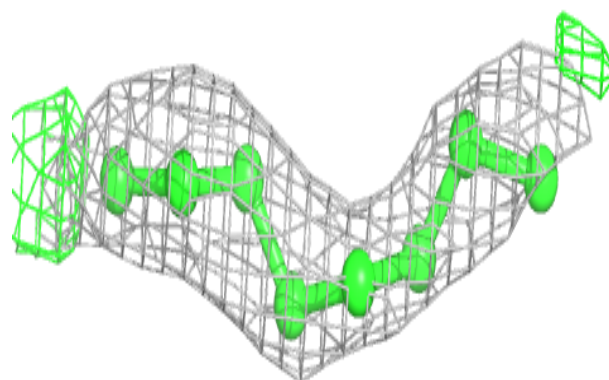


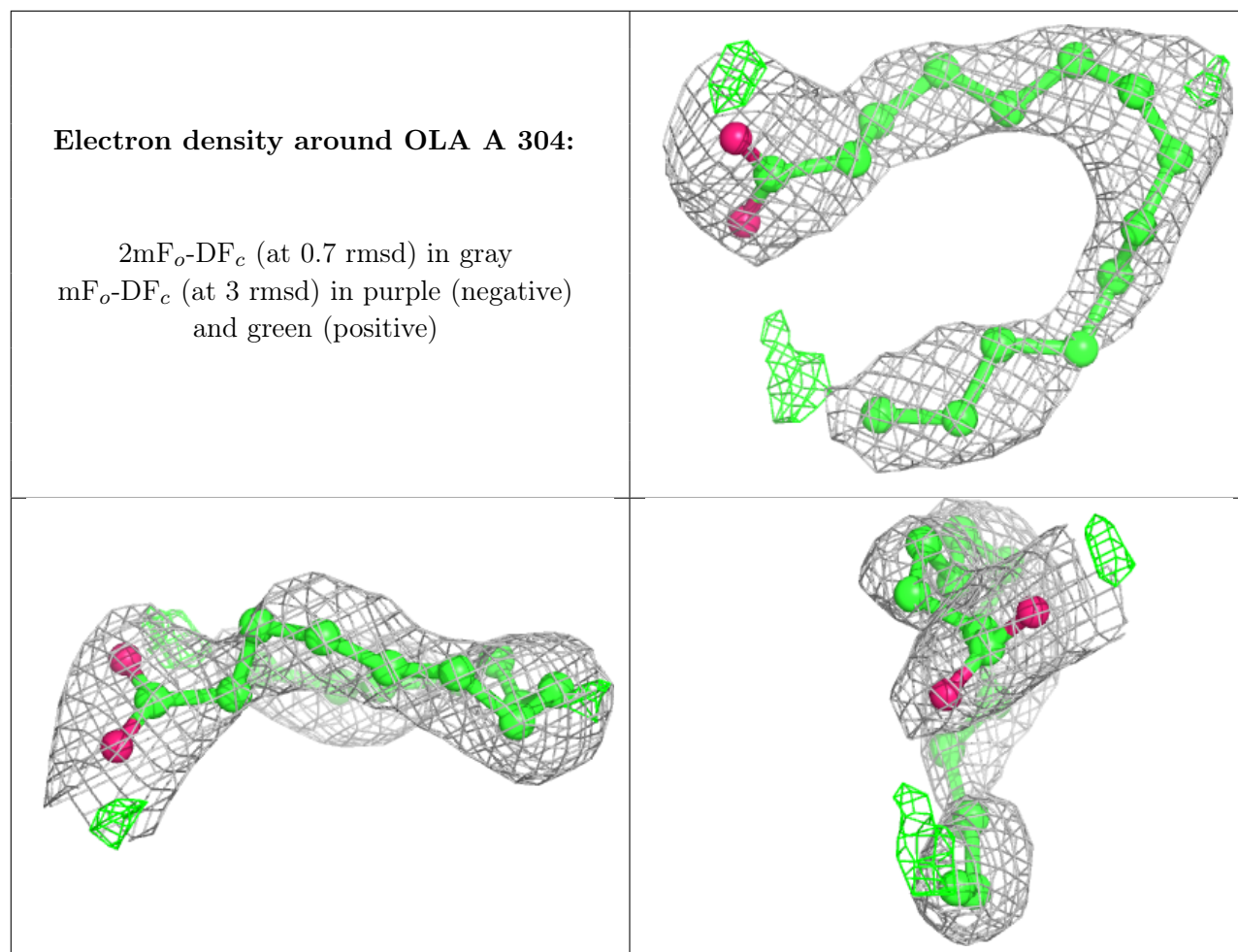
Electron density around OLA B 309:

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

**Electron density around LFA A 306:**

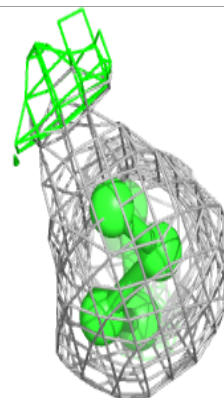
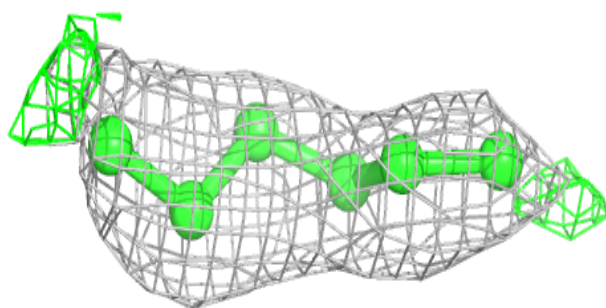
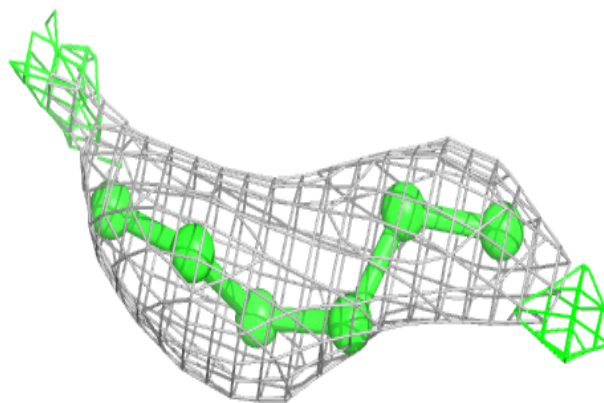
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



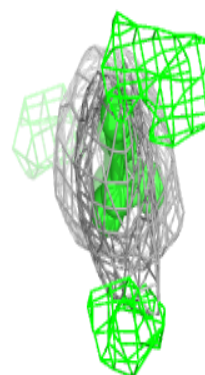
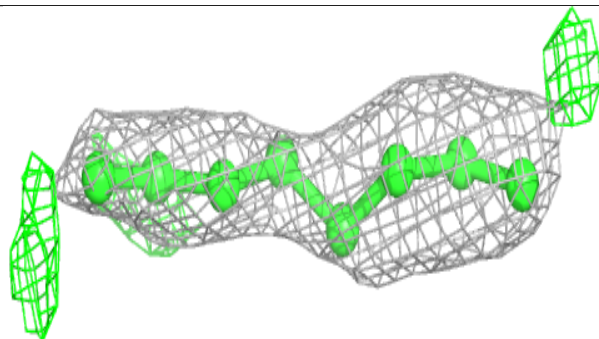
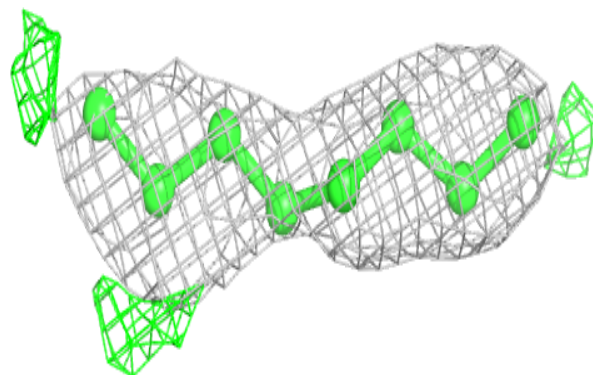


Electron density around LFA C 1406:

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

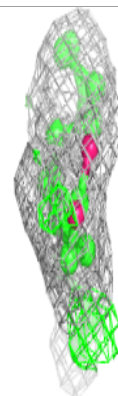
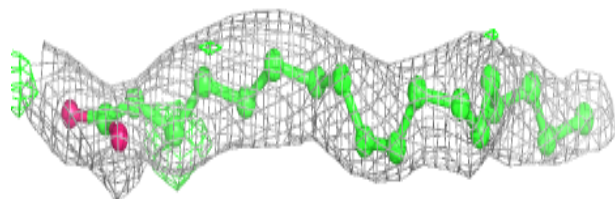
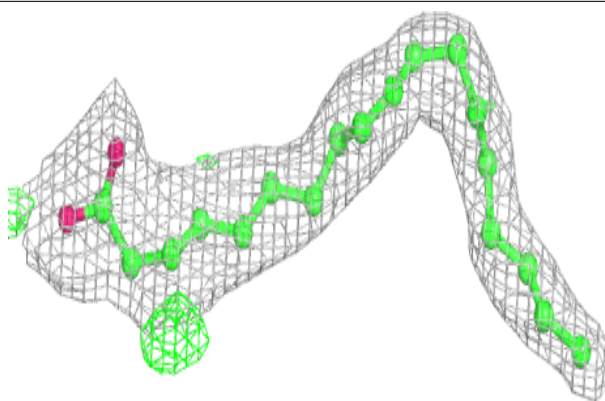
**Electron density around LFA E 413:**

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

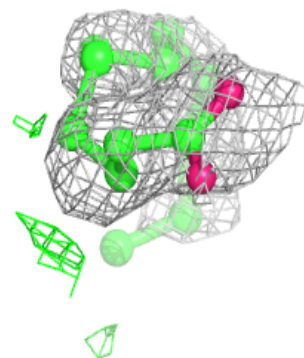
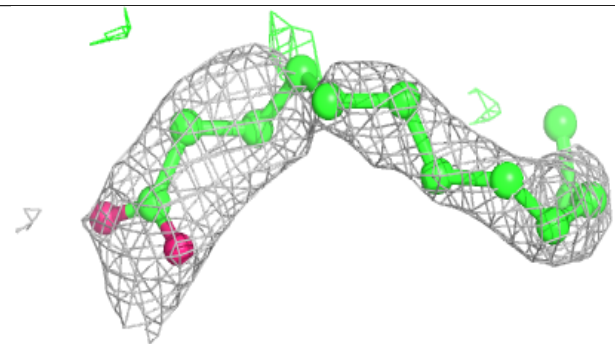
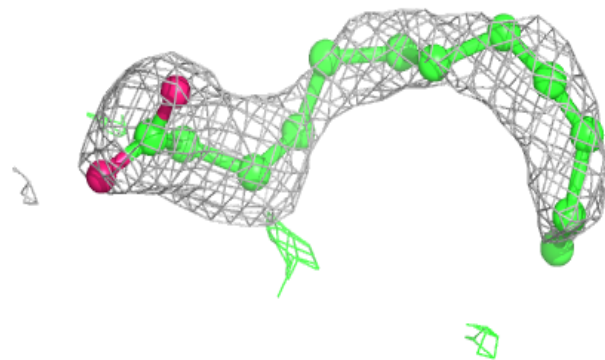


Electron density around OLA E 405:

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

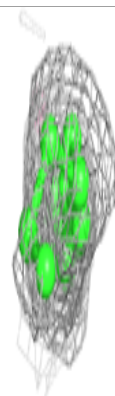
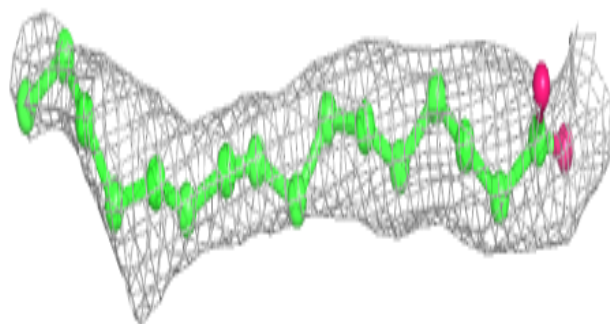
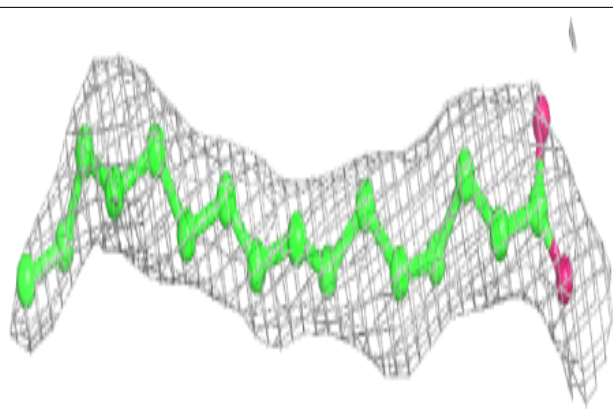
**Electron density around OLA B 305:**

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

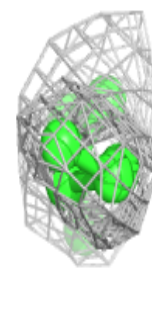
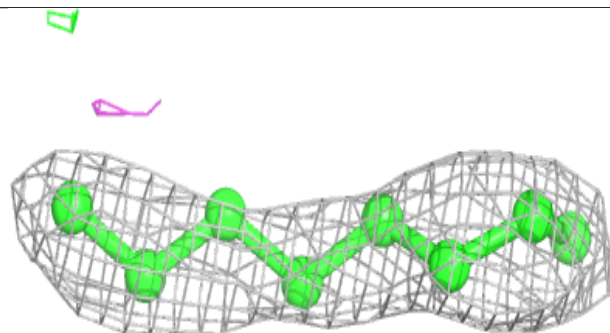
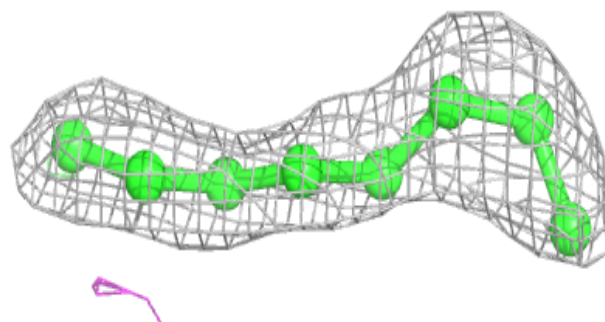


Electron density around OLA E 411:

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

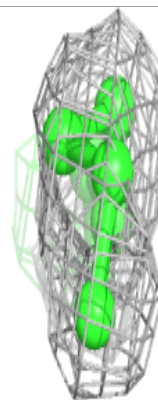
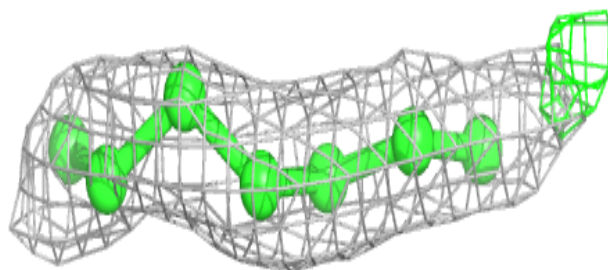
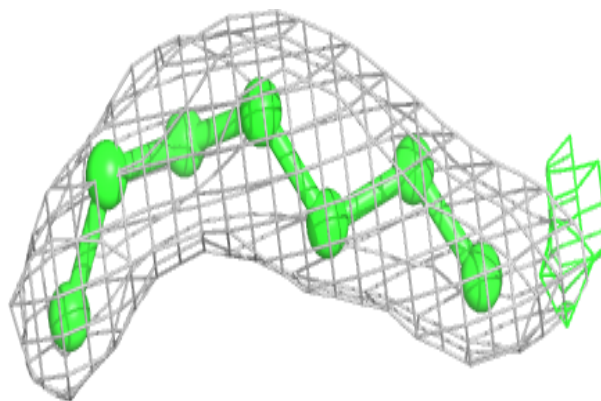
**Electron density around LFA D 410:**

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

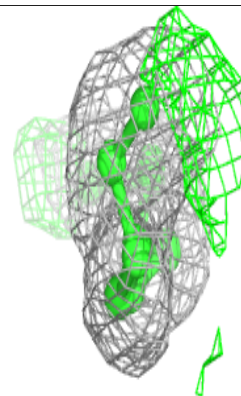
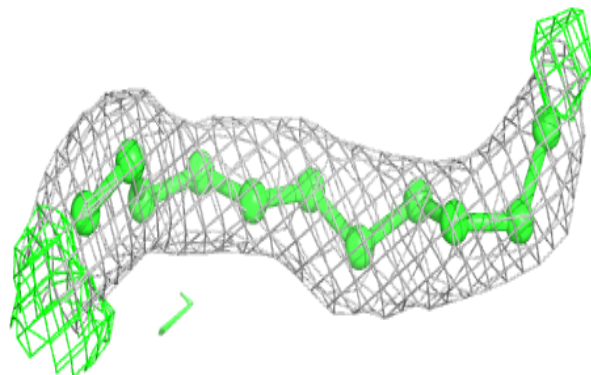
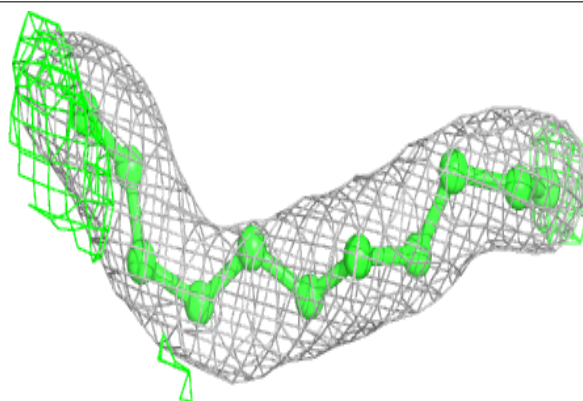


Electron density around LFA D 411:

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

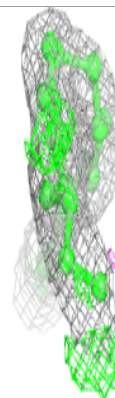
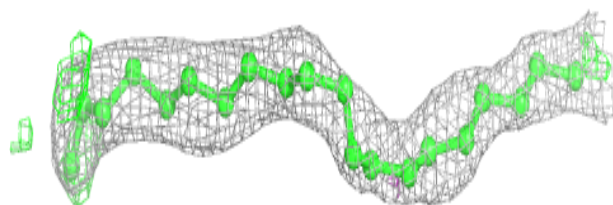
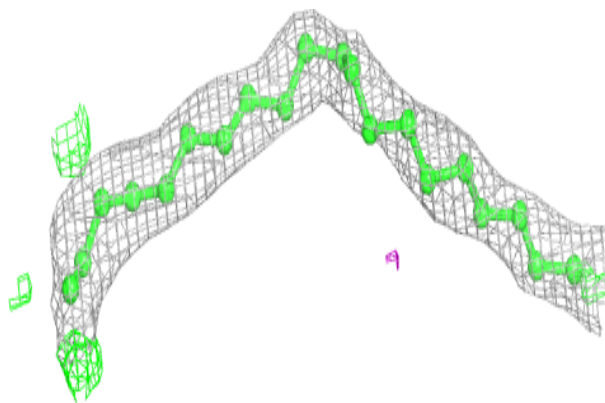
**Electron density around LFA B 303:**

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

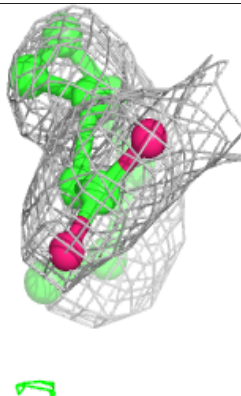
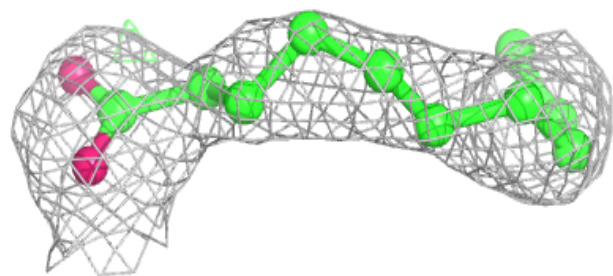
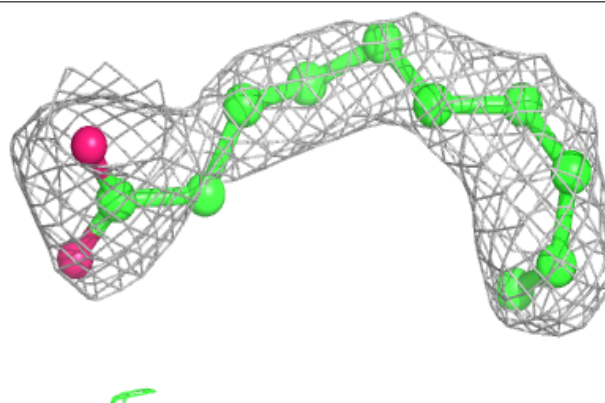


Electron density around LFA D 401:

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

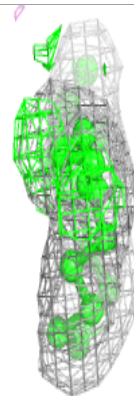
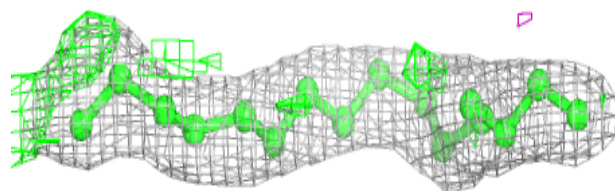
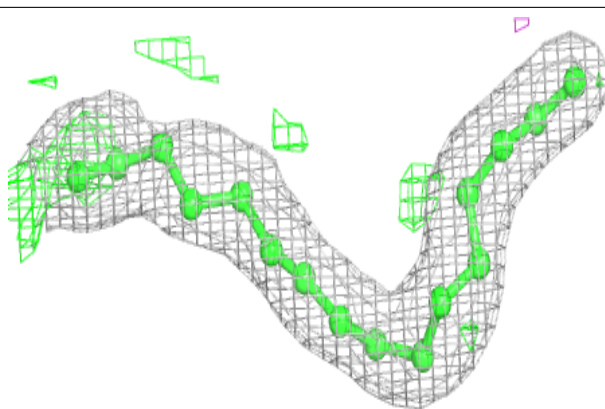
**Electron density around OLA B 313:**

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

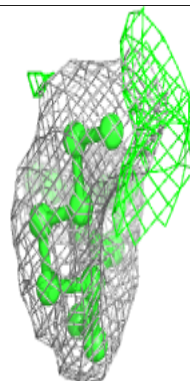
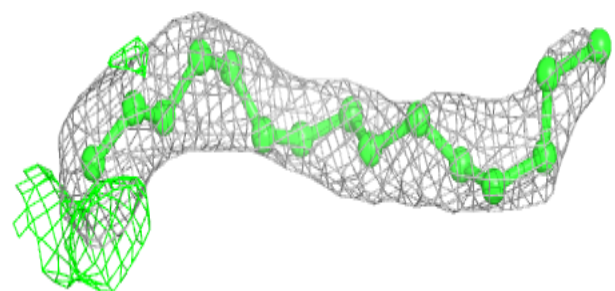
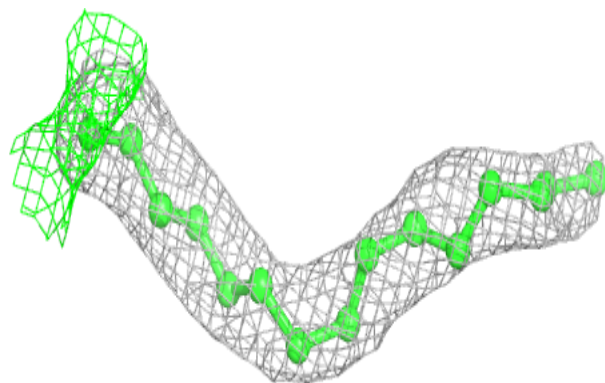


Electron density around LFA C 1403:

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

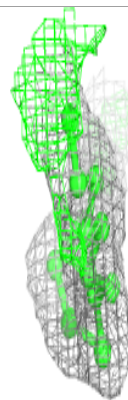
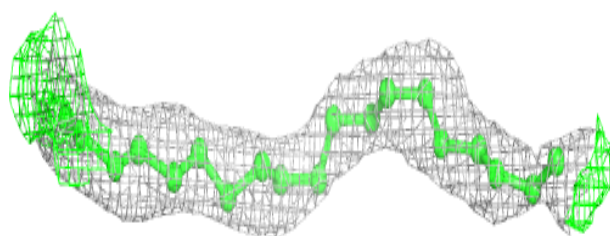
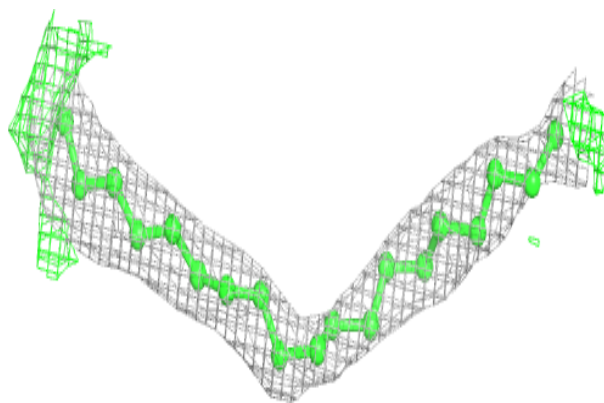
**Electron density around LFA C 1404:**

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

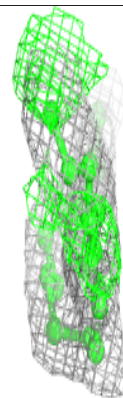
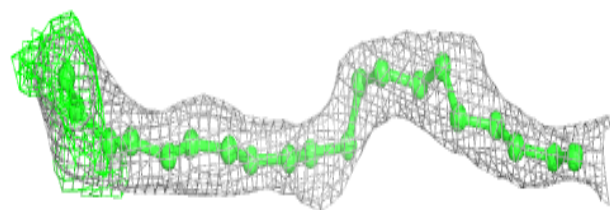
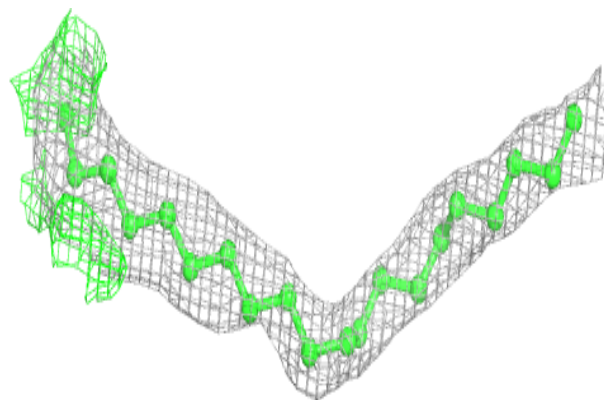


Electron density around LFA E 406:

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

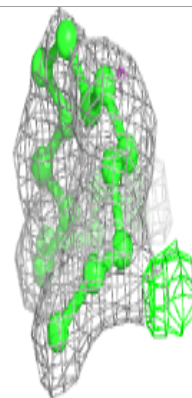
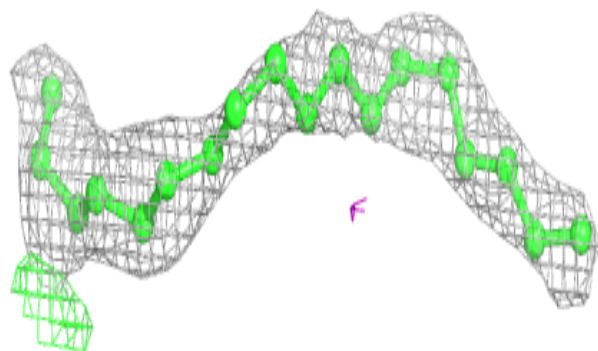
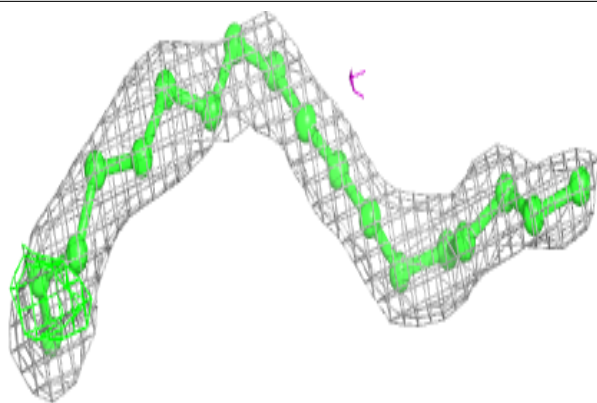
**Electron density around LFA E 402:**

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

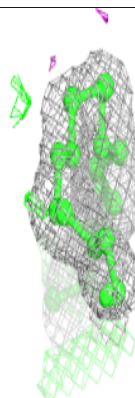
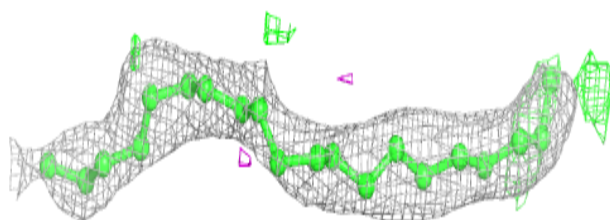
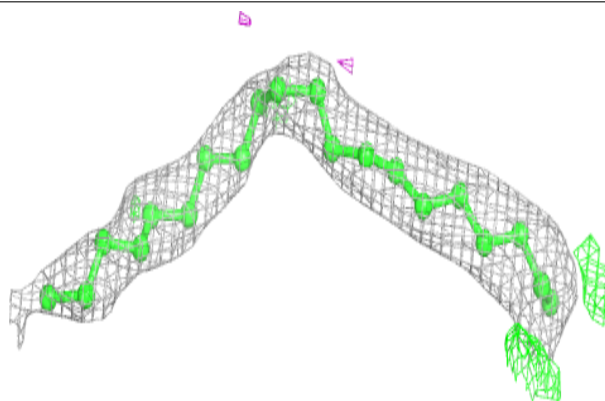


Electron density around LFA E 408:

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

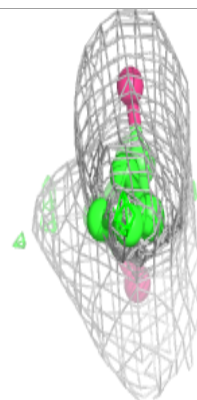
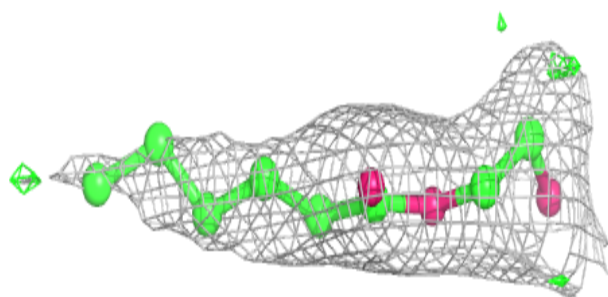
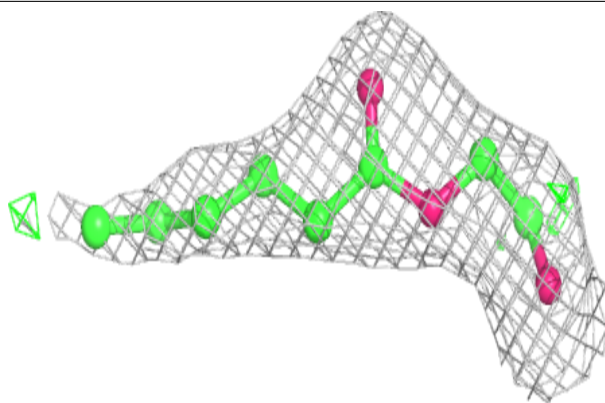
**Electron density around LFA B 312:**

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

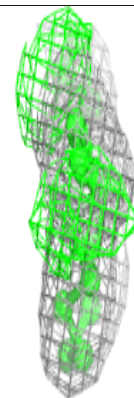
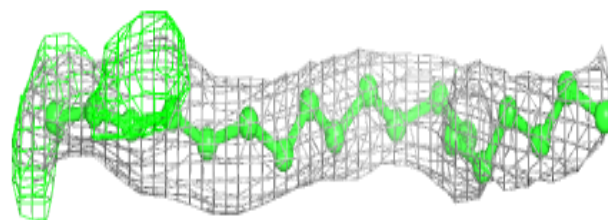
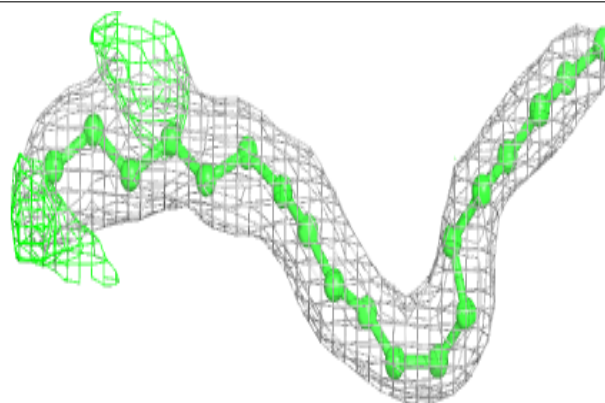


Electron density around OLC C 1402:

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

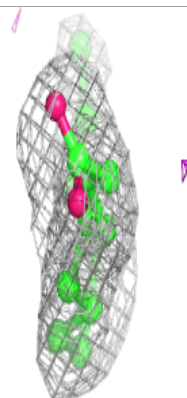
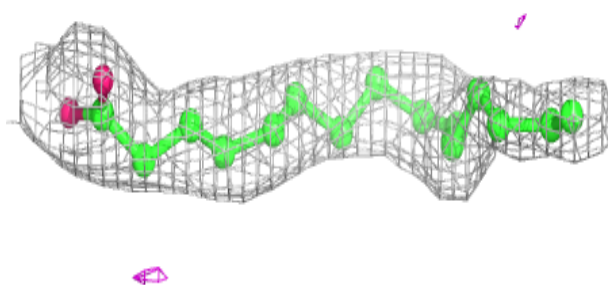
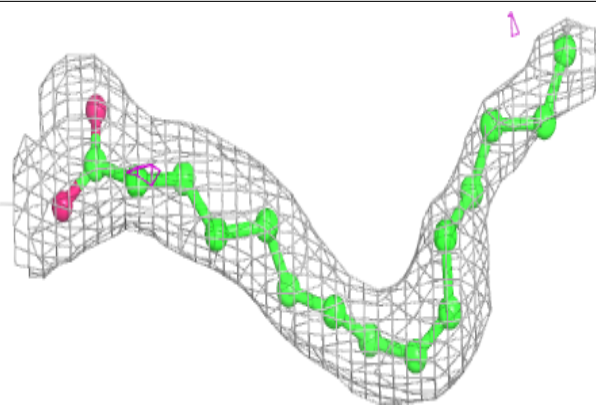
**Electron density around LFA B 302:**

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

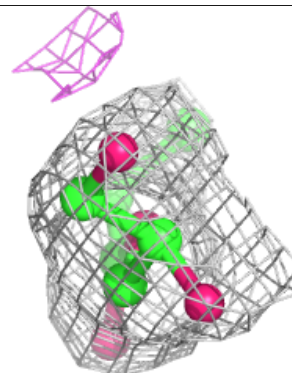
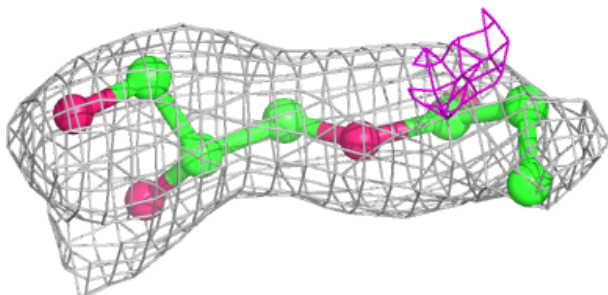
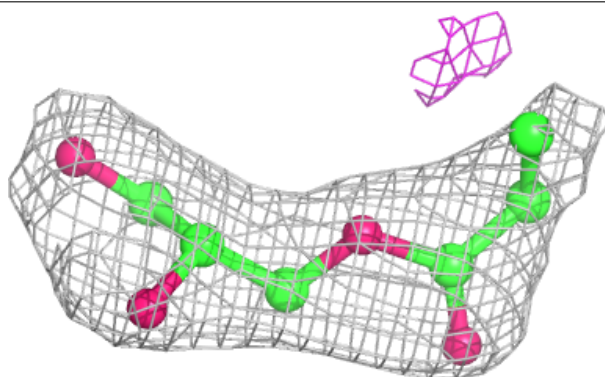


Electron density around OLA A 302:

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

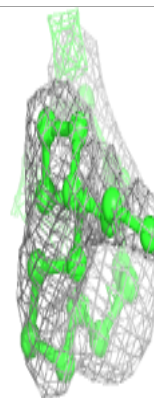
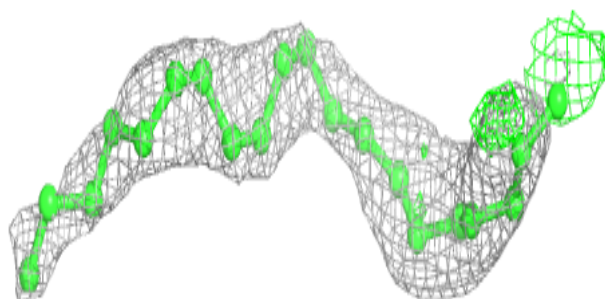
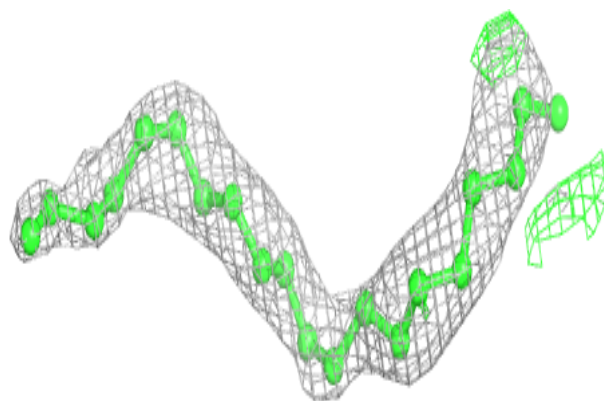
**Electron density around OLC E 404:**

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

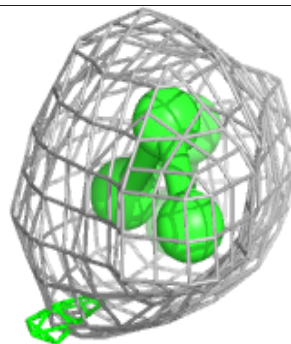
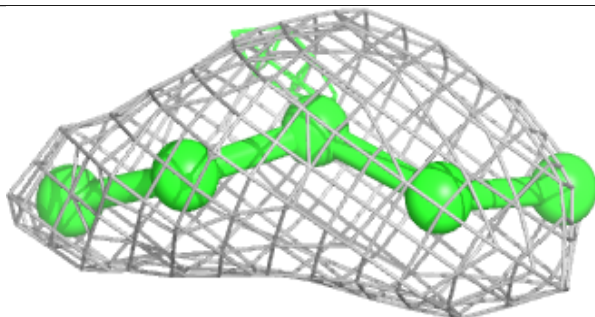
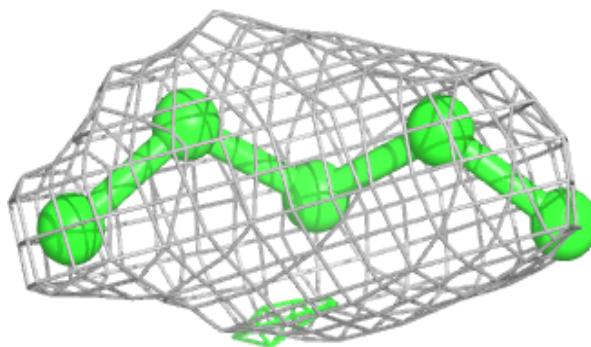


Electron density around LFA A 303:

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

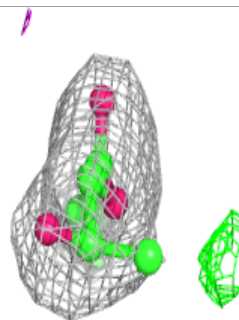
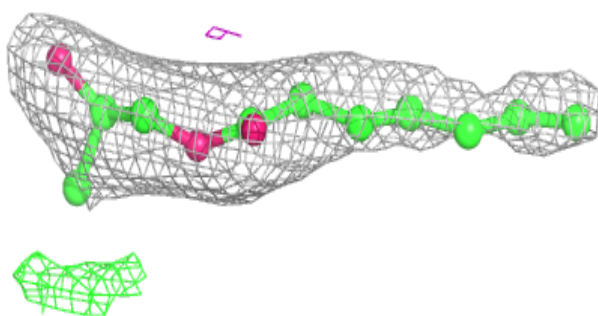
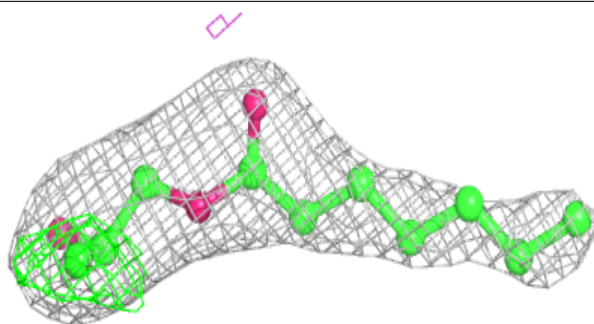
**Electron density around LFA E 409:**

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

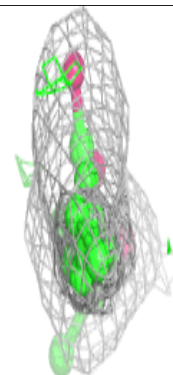
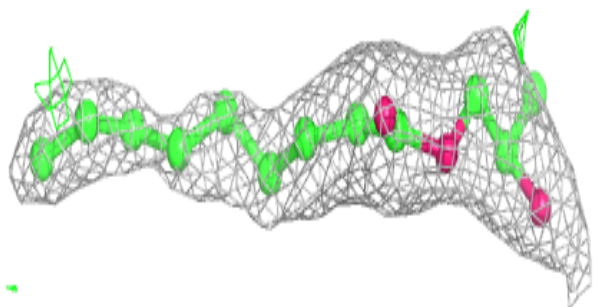
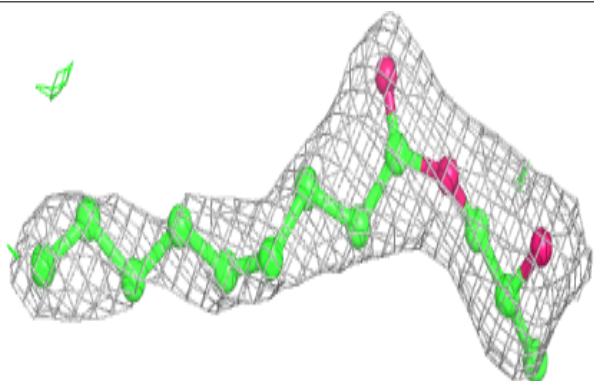


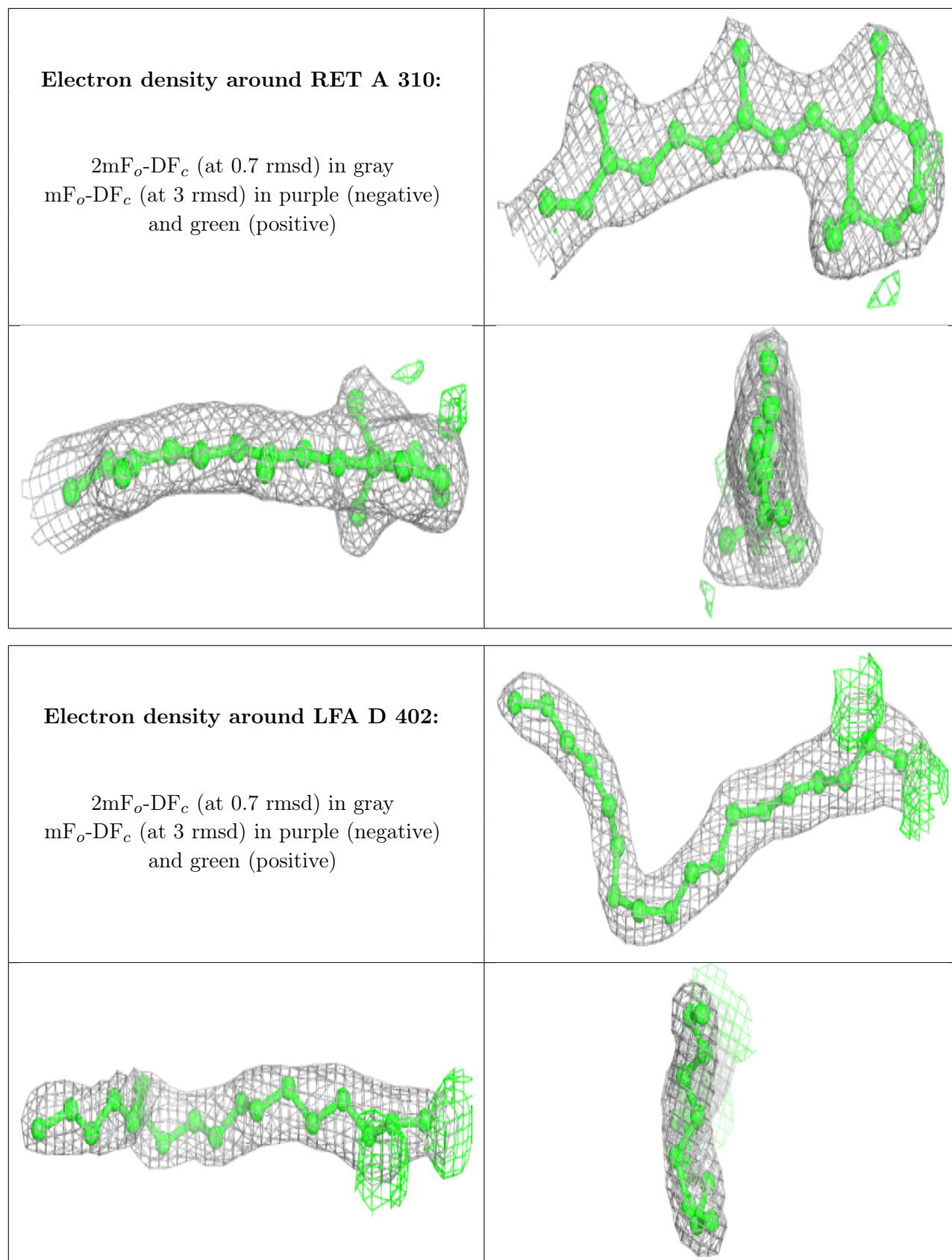
Electron density around OLC B 301:

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

**Electron density around OLC D 405:**

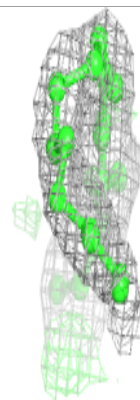
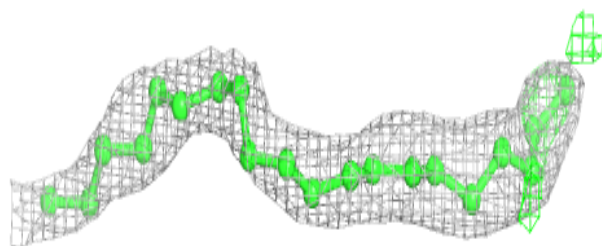
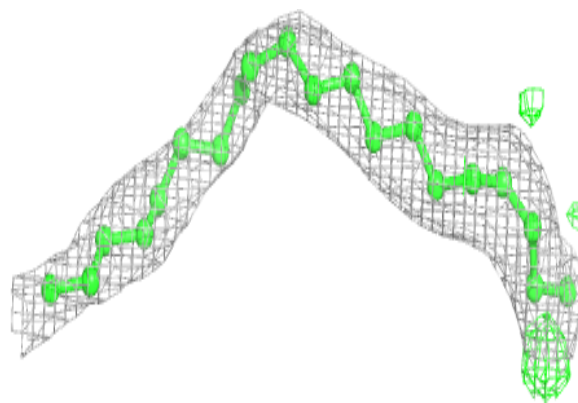
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



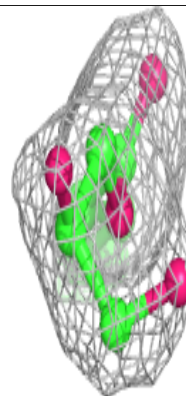
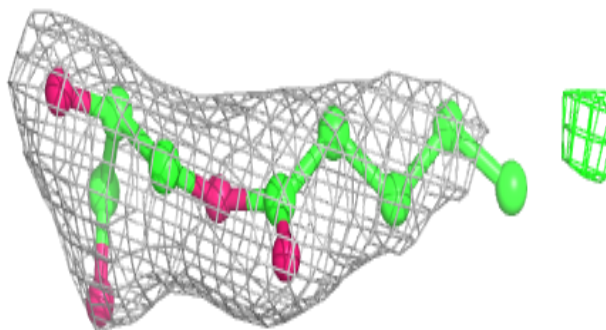
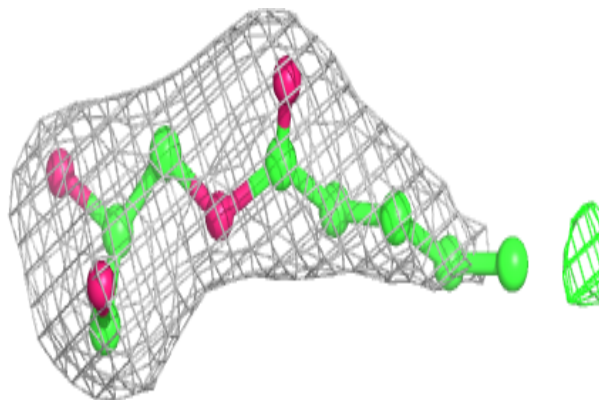


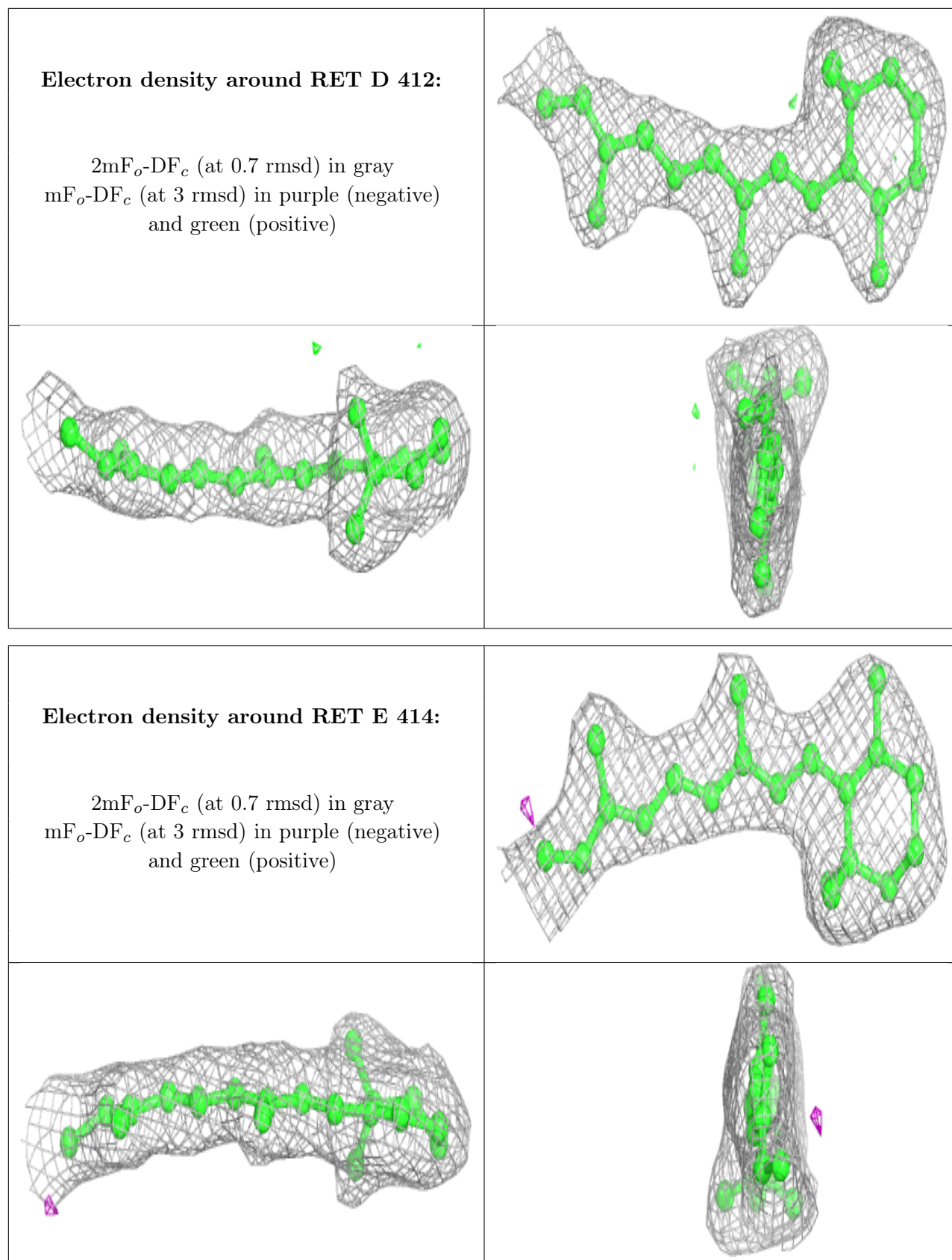
Electron density around LFA B 304:

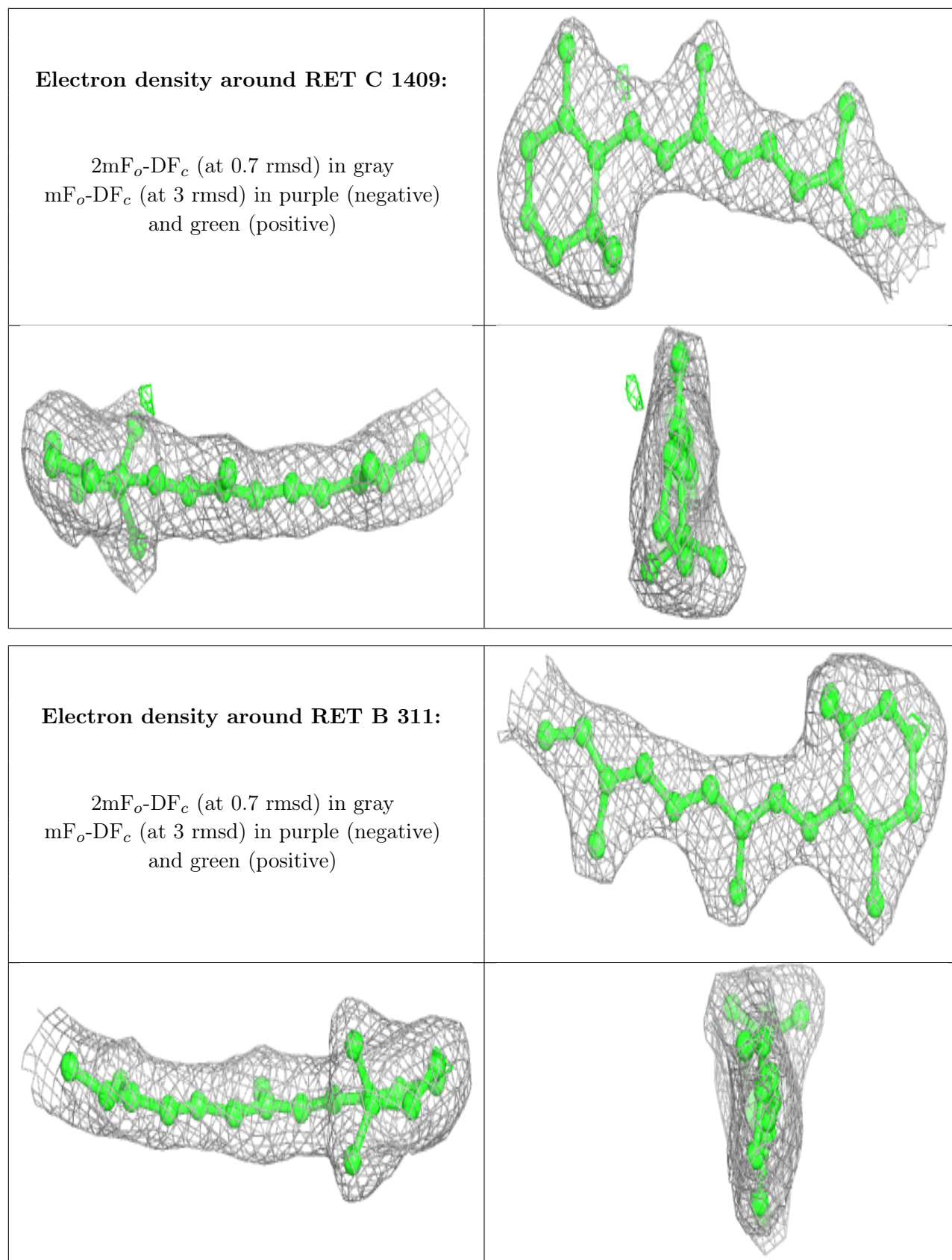
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around OLC 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)

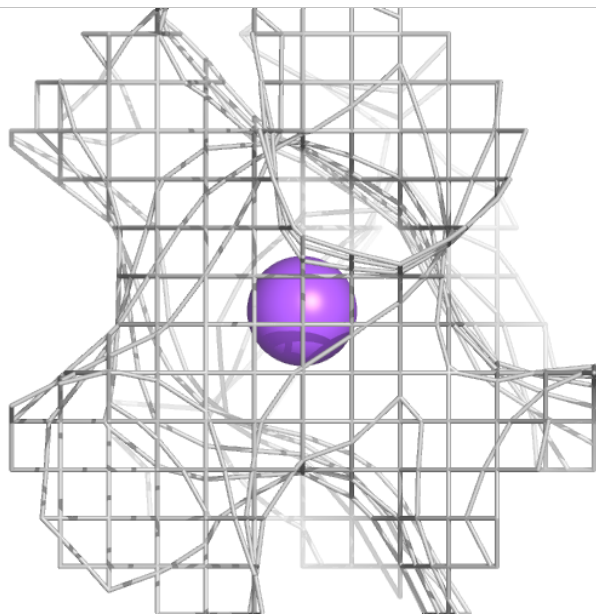
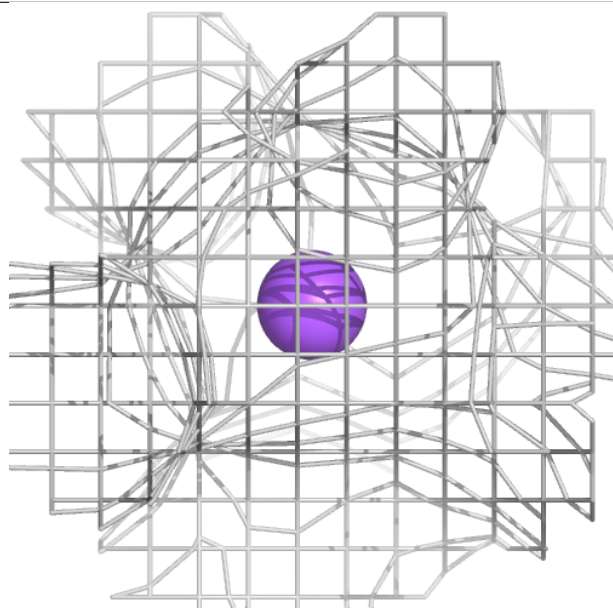
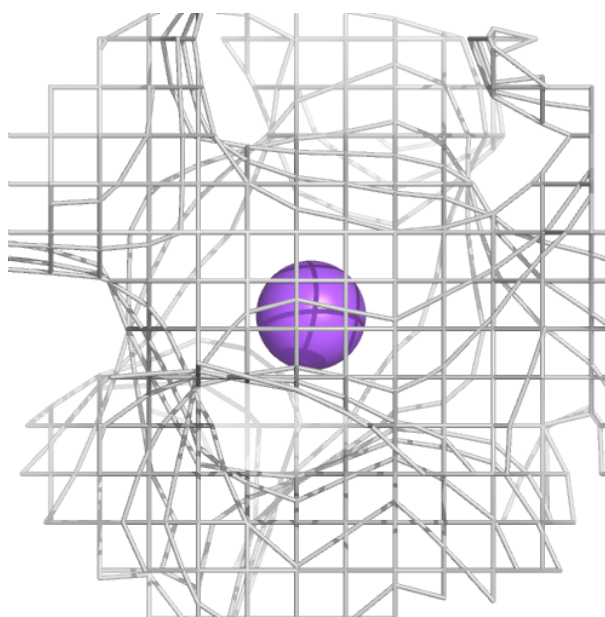






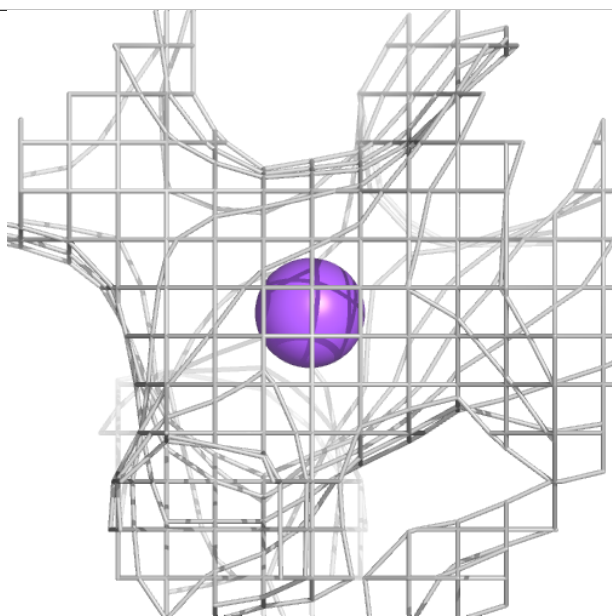
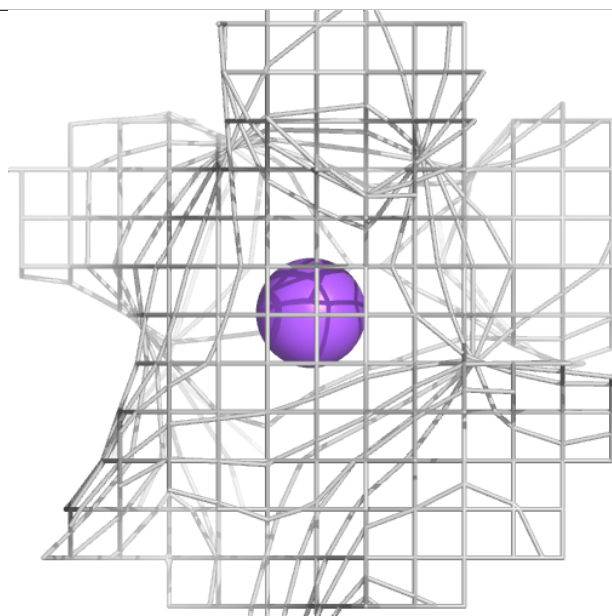
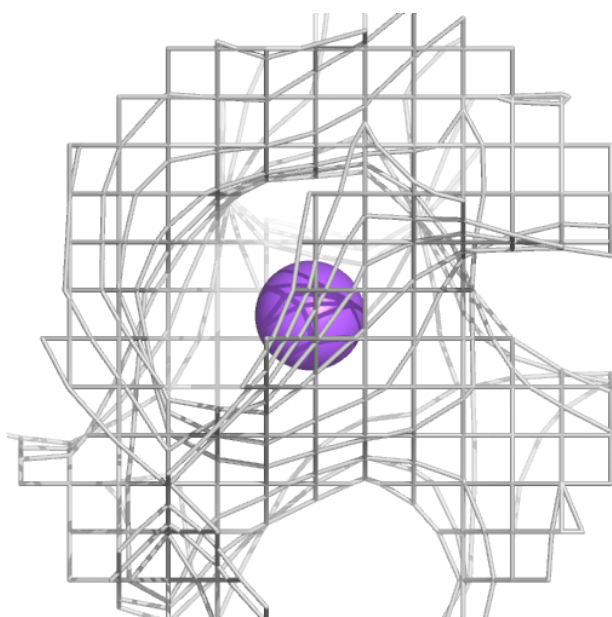
Electron density around NA B 314:

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



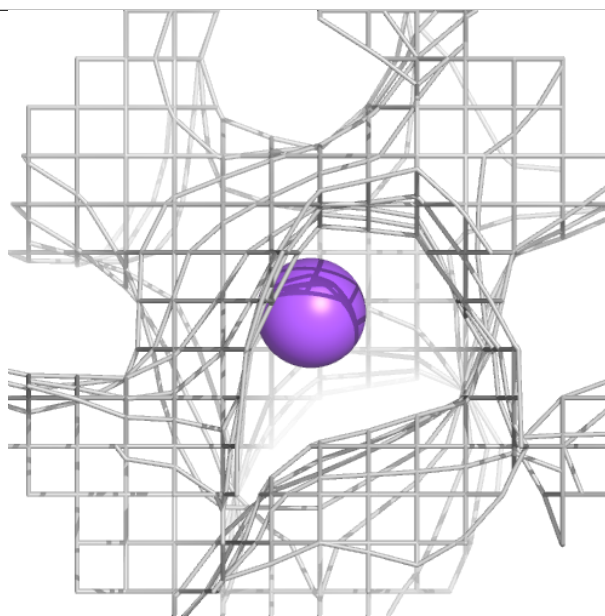
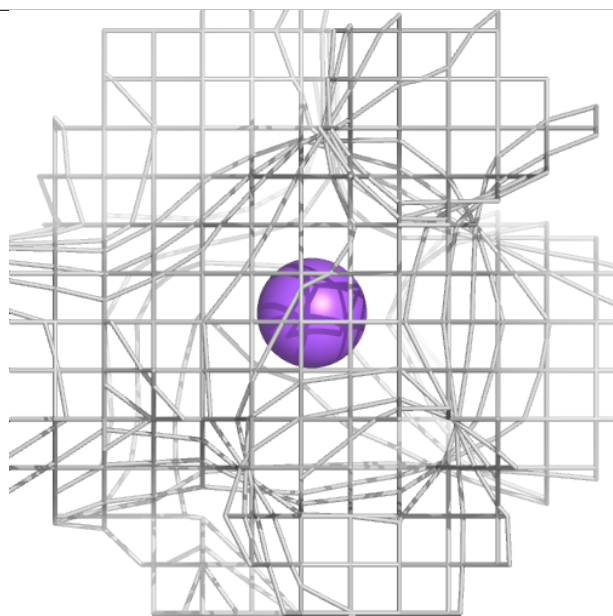
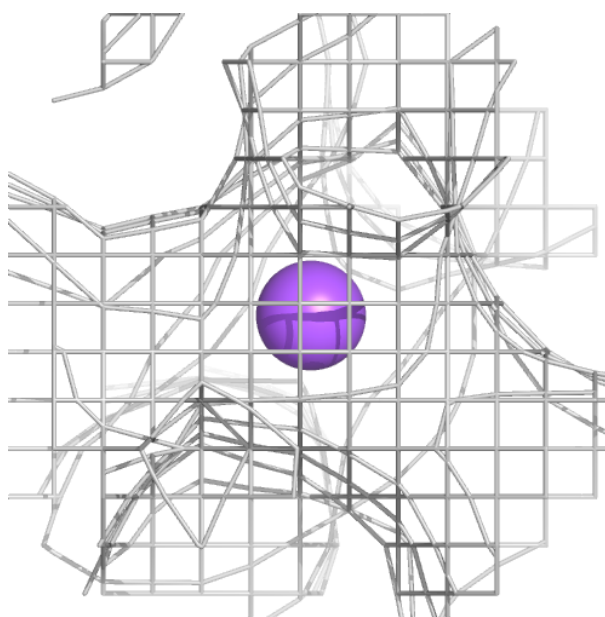
Electron density around NA E 415:

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



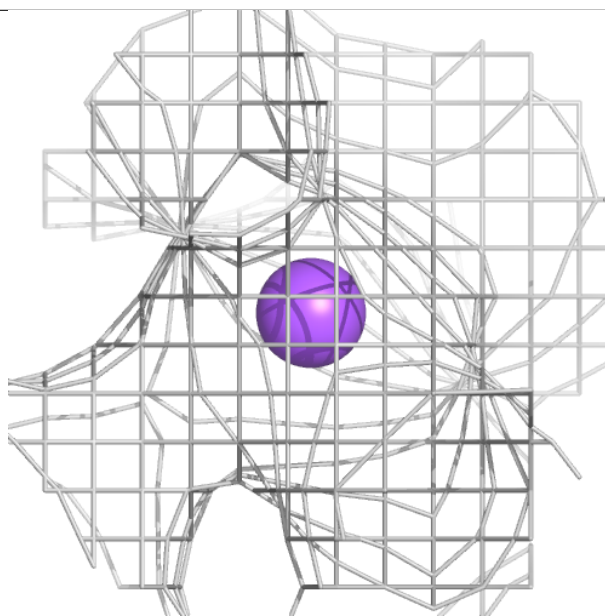
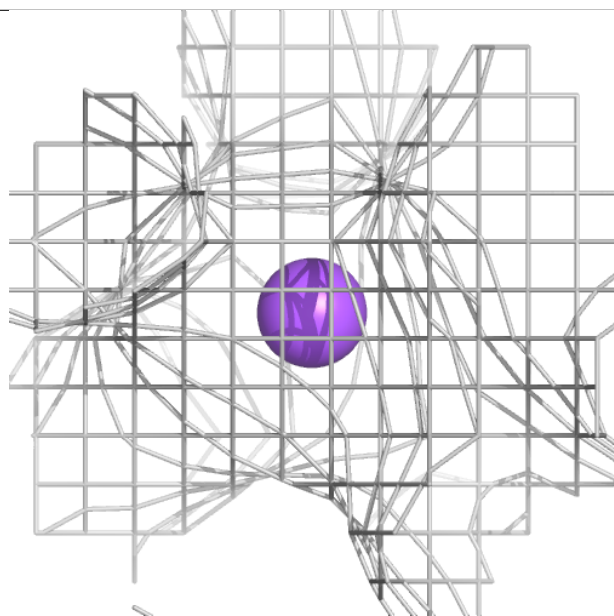
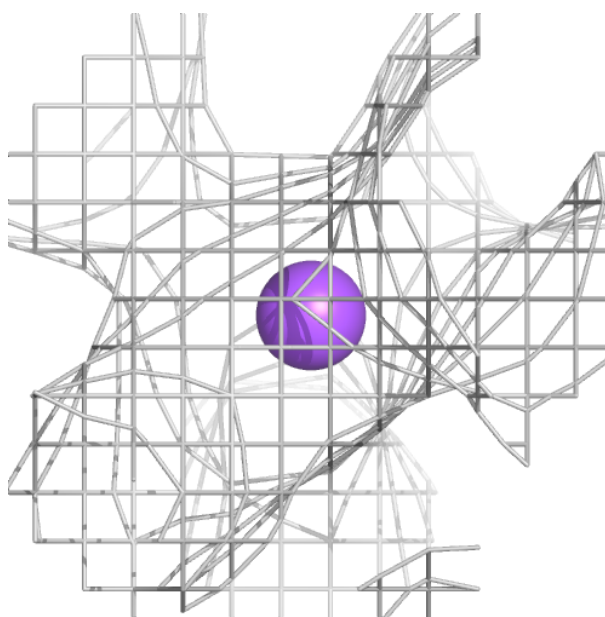
Electron density around NA D 413:

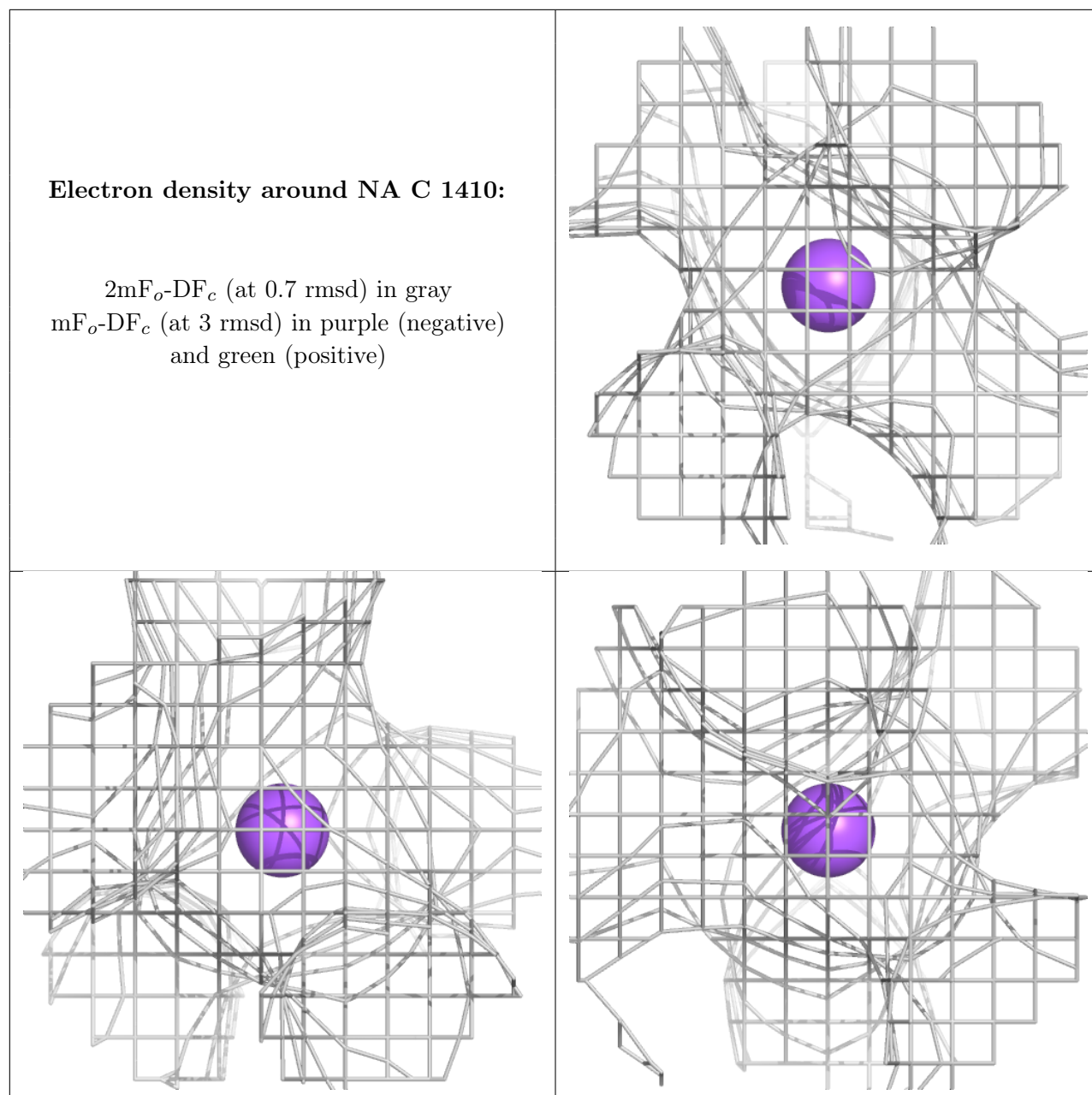
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around NA A 312:

$2mF_o-DF_c$ (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.