



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 05:24 PM UTC

PDB ID : 5Y79 / pdb_00005y79
Title : Crystal structure of the triose-phosphate/phosphate translocator in complex with 3-phosphoglycerate
Authors : Lee, Y.; Nishizawa, T.; Takemoto, M.; Kumazaki, K.; Yamashita, K.; Hirata, K.; Minoda, A.; Nagatoishi, S.; Tsumoto, K.; Ishitani, R.; Nureki, O.
Deposited on : 2017-08-16
Resolution : 2.20 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

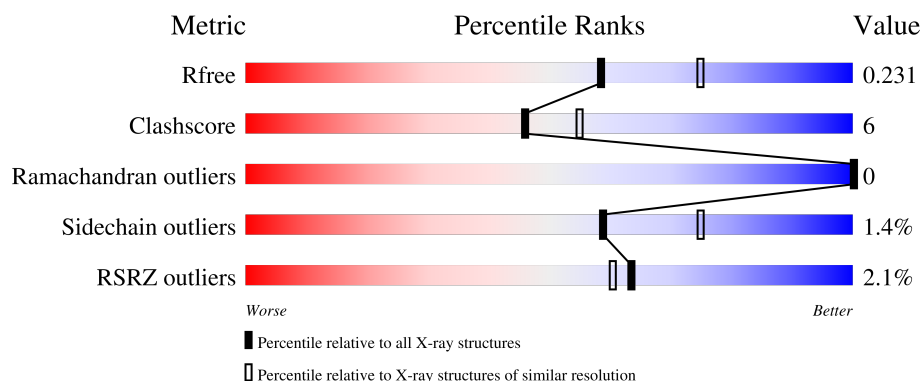
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

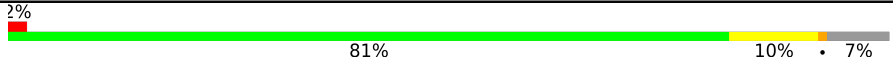
The reported resolution of this entry is 2.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	329	 2% 81% 10% 7%
1	B	329	 2% 82% 11% 7%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 5396 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 Putative hexose phosphate translocator.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	305	Total	C	N	O	S	0	0	0
			2354	1580	364	398	12			
1	B	305	Total	C	N	O	S	0	0	0
			2359	1586	364	397	12			

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	90	MET	-	initiating methionine	UNP B5AJT1
A	411	GLY	-	expression tag	UNP B5AJT1
A	412	THR	-	expression tag	UNP B5AJT1
A	413	GLU	-	expression tag	UNP B5AJT1
A	414	ASN	-	expression tag	UNP B5AJT1
A	415	LEU	-	expression tag	UNP B5AJT1
A	416	TYR	-	expression tag	UNP B5AJT1
A	417	PHE	-	expression tag	UNP B5AJT1
A	418	GLN	-	expression tag	UNP B5AJT1
B	90	MET	-	initiating methionine	UNP B5AJT1
B	411	GLY	-	expression tag	UNP B5AJT1
B	412	THR	-	expression tag	UNP B5AJT1
B	413	GLU	-	expression tag	UNP B5AJT1
B	414	ASN	-	expression tag	UNP B5AJT1
B	415	LEU	-	expression tag	UNP B5AJT1
B	416	TYR	-	expression tag	UNP B5AJT1
B	417	PHE	-	expression tag	UNP B5AJT1
B	418	GLN	-	expression tag	UNP B5AJT1

- Molecule 2 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (CCD ID: OLC) (formula: C₂₁H₄₀O₄).



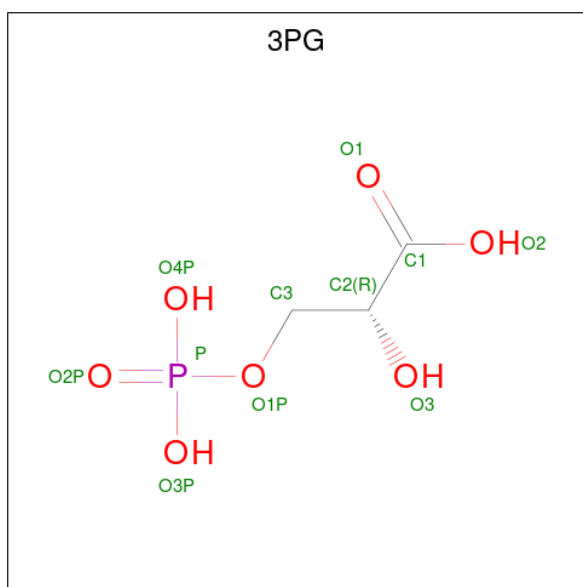
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 25 21 4	0	0
2	A	1	Total C 14 14	0	0
2	A	1	Total C 6 6	0	0
2	A	1	Total C 8 8	0	0
2	A	1	Total C O 19 15 4	0	0
2	A	1	Total C 12 12	0	0
2	A	1	Total C 13 13	0	0
2	A	1	Total C 14 14	0	0
2	A	1	Total C 8 8	0	0
2	A	1	Total C 12 12	0	0
2	A	1	Total C 7 7	0	0
2	A	1	Total C 11 11	0	0
2	A	1	Total C 8 8	0	0
2	A	1	Total C O 11 7 4	0	0

Continued on next page...

Continued from previous page...

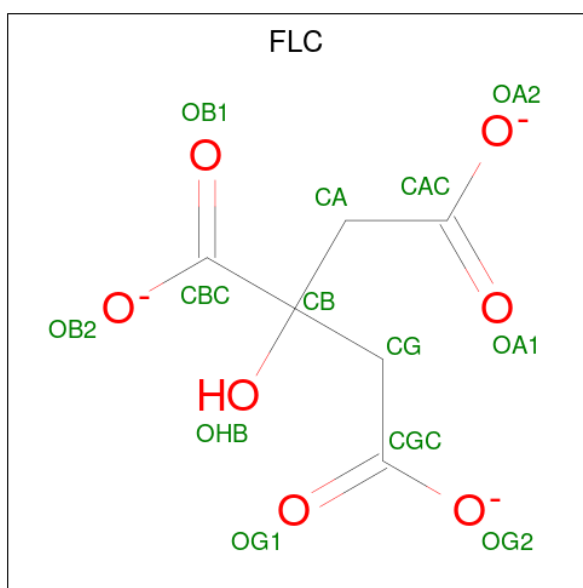
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C 13 13	0	0
2	B	1	Total C O 25 21 4	0	0
2	B	1	Total C O 25 21 4	0	0
2	B	1	Total C O 25 21 4	0	0
2	B	1	Total C 14 14	0	0
2	B	1	Total C O 19 15 4	0	0
2	B	1	Total C 8 8	0	0
2	B	1	Total C 13 13	0	0
2	B	1	Total C 8 8	0	0
2	B	1	Total C 8 8	0	0
2	B	1	Total C 16 16	0	0
2	B	1	Total C 8 8	0	0
2	B	1	Total C O 20 18 2	0	0
2	B	1	Total C 17 17	0	0
2	B	1	Total C 9 9	0	0
2	B	1	Total C 6 6	0	0
2	B	1	Total C O 12 8 4	0	0
2	B	1	Total C 12 12	0	0

- Molecule 3 is 3-PHOSPHOGLYCERIC ACID (CCD ID: 3PG) (formula: C₃H₇O₇P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			11	3	7	1		
3	B	1	Total	C	O	P	0	0
			11	3	7	1		

- Molecule 4 is CITRATE ANION (CCD ID: FLC) (formula: $C_6H_5O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			13	6	7		

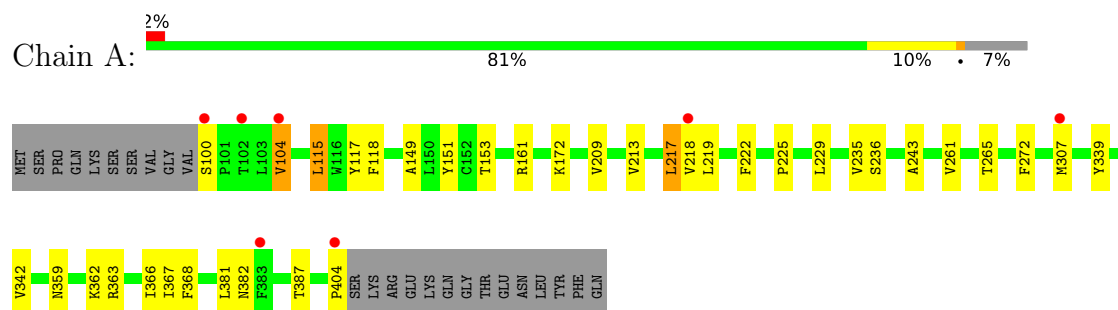
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	95	Total 95	O 95	0	0
5	B	127	Total 127	O 127	0	0

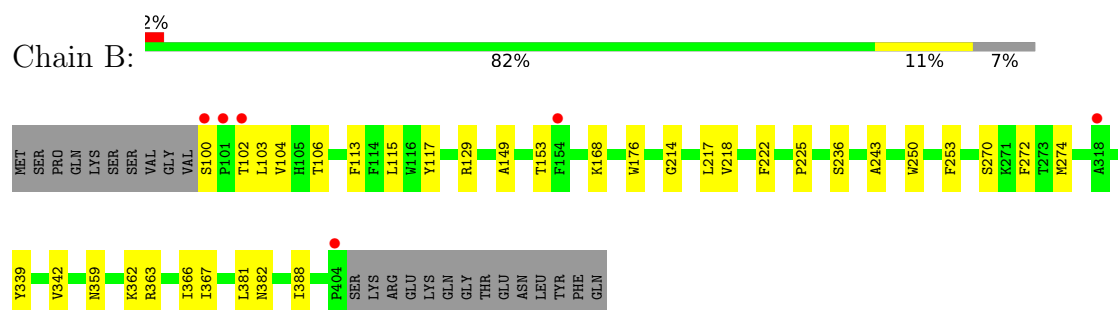
3 Residue-property plots [i](#)

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: Putative hexose phosphate translocator



- Molecule 1: Putative hexose phosphate translocator



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	107.05Å 165.33Å 41.14Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.00 – 2.20 49.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.00-2.20) 92.2 (49.00-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$	-	Xtriage
Refinement program	PHENIX (1.10_2155: ???)	Depositor
R, R_{free}	0.191 , 0.229 0.193 , 0.231	Depositor DCC
R_{free} test set	2059 reflections (5.39%)	wwPDB-VP
Wilson B-factor (Å ²)	(Not available)	Xtriage
Anisotropy	(Not available)	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.8	EDS
L-test for twinning ¹	$\langle L \rangle =$ (Not available), $\langle L^2 \rangle =$ (Not available)	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	5396	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *(Not available)*

¹Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 3PG, OLC, FLC

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	0.15	0/2423	0.32	0/3313
1	B	0.14	0/2428	0.30	0/3320
All	All	0.14	0/4851	0.31	0/6633

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

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	2354	0	2375	30	0
1	B	2359	0	2397	31	0
2	A	181	0	272	15	0
2	B	245	0	381	20	0
3	A	11	0	4	0	0
3	B	11	0	4	0	0
4	A	13	0	5	1	0
5	A	95	0	0	0	0
5	B	127	0	0	0	0
All	All	5396	0	5438	62	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 6.

The worst 5 of 62 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)
1:A:218:VAL:HG11	1:A:272:PHE:HZ	1.49	0.77
1:B:113:PHE:HB3	2:B:517:OLC:H9	1.71	0.73
4:A:517:FLC:OA2	4:A:517:FLC:OHB	2.10	0.69
2:A:512:OLC:H5	2:A:515:OLC:H10	1.73	0.69
1:B:129:ARG:HH21	2:B:514:OLC:H5	1.55	0.68

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	303/329 (92%)	300 (99%)	3 (1%)	0	100	100
1	B	303/329 (92%)	302 (100%)	1 (0%)	0	100	100
All	All	606/658 (92%)	602 (99%)	4 (1%)	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	251/286 (88%)	246 (98%)	5 (2%)	48	64
1	B	254/286 (89%)	252 (99%)	2 (1%)	73	85
All	All	505/572 (88%)	498 (99%)	7 (1%)	59	75

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	229	LEU
1	A	339	TYR
1	B	339	TYR
1	B	104	VAL
1	A	217	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

35 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	OLC	A	510	-	11,11,24	0.26	0	10,10,25	0.83	0
2	OLC	B	514	-	8,8,24	0.12	0	7,7,25	0.73	0
2	OLC	A	502	-	13,13,24	0.24	0	12,12,25	0.71	0
2	OLC	A	515	-	12,12,24	0.26	0	11,11,25	0.73	0
2	OLC	B	507	-	12,12,24	0.24	0	11,11,25	0.73	0
2	OLC	B	516	-	11,11,24	0.98	1 (9%)	12,12,25	1.18	1 (8%)
2	OLC	A	505	-	18,18,24	0.80	1 (5%)	19,19,25	1.09	1 (5%)
2	OLC	A	508	-	13,13,24	0.24	0	12,12,25	0.74	0
2	OLC	B	505	-	18,18,24	0.78	1 (5%)	19,19,25	1.16	1 (5%)
2	OLC	B	503	2	24,24,24	0.70	1 (4%)	25,25,25	0.86	1 (4%)
2	OLC	B	501	-	24,24,24	0.68	1 (4%)	25,25,25	0.99	1 (4%)
3	3PG	B	518	-	9,10,10	0.92	0	11,14,14	1.41	2 (18%)
2	OLC	A	503	-	5,5,24	0.13	0	4,4,25	0.62	0
2	OLC	B	504	-	13,13,24	0.23	0	12,12,25	0.70	0
4	FLC	A	517	-	12,12,12	1.13	0	17,17,17	1.22	1 (5%)
2	OLC	A	514	-	10,10,24	1.06	1 (10%)	11,11,25	1.33	1 (9%)
2	OLC	A	501	-	24,24,24	0.70	1 (4%)	25,25,25	1.07	1 (4%)
2	OLC	A	507	2	12,12,24	0.25	0	11,11,25	0.77	0
2	OLC	A	513	-	7,7,24	0.11	0	6,6,25	0.75	0
2	OLC	A	511	-	6,6,24	0.13	0	5,5,25	0.66	0
2	OLC	B	517	-	11,11,24	0.27	0	10,10,25	0.82	0
2	OLC	A	509	-	7,7,24	0.10	0	6,6,25	0.79	0
2	OLC	B	515	-	5,5,24	0.13	0	4,4,25	0.61	0
2	OLC	A	504	-	7,7,24	0.11	0	6,6,25	0.76	0
2	OLC	B	502	2	24,24,24	0.70	1 (4%)	25,25,25	0.91	1 (4%)
2	OLC	B	512	-	19,19,24	0.76	1 (5%)	19,19,25	0.97	0
2	OLC	B	511	-	7,7,24	0.11	0	6,6,25	0.75	0
2	OLC	B	508	-	7,7,24	0.12	0	6,6,25	0.72	0
2	OLC	A	506	-	11,11,24	0.25	0	10,10,25	0.67	0
3	3PG	A	516	-	9,10,10	0.94	0	11,14,14	1.33	1 (9%)
2	OLC	B	509	-	7,7,24	0.12	0	6,6,25	0.72	0
2	OLC	B	506	-	7,7,24	0.10	0	6,6,25	0.80	0
2	OLC	B	510	-	15,15,24	0.22	0	14,14,25	0.74	0
2	OLC	A	512	-	10,10,24	0.33	0	9,9,25	0.92	0
2	OLC	B	513	2	16,16,24	0.24	0	15,15,25	0.66	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	OLC	A	510	-	-	4/9/9/24	-
2	OLC	B	514	-	-	2/6/6/24	-
2	OLC	A	502	-	-	4/11/11/24	-
2	OLC	A	515	-	-	4/10/10/24	-
2	OLC	B	507	-	-	3/10/10/24	-
2	OLC	B	516	-	-	6/11/11/24	-
2	OLC	A	505	-	-	9/18/18/24	-
2	OLC	A	508	-	-	6/11/11/24	-
2	OLC	B	505	-	-	11/18/18/24	-
2	OLC	B	503	2	-	11/24/24/24	-
2	OLC	B	501	-	-	8/24/24/24	-
3	3PG	B	518	-	-	5/10/10/10	-
2	OLC	A	503	-	-	1/3/3/24	-
2	OLC	B	504	-	-	6/11/11/24	-
4	FLC	A	517	-	-	14/16/16/16	-
2	OLC	A	514	-	-	6/10/10/24	-
2	OLC	A	501	-	-	11/24/24/24	-
2	OLC	A	507	2	-	3/10/10/24	-
2	OLC	A	513	-	-	4/5/5/24	-
2	OLC	A	511	-	-	2/4/4/24	-
2	OLC	B	517	-	-	5/9/9/24	-
2	OLC	A	509	-	-	2/5/5/24	-
2	OLC	B	515	-	-	2/3/3/24	-
2	OLC	A	504	-	-	4/5/5/24	-
2	OLC	B	502	2	-	8/24/24/24	-
2	OLC	B	512	-	-	6/17/17/24	-
2	OLC	B	511	-	-	3/5/5/24	-
2	OLC	B	508	-	-	3/5/5/24	-
2	OLC	A	506	-	-	2/9/9/24	-
3	3PG	A	516	-	-	7/10/10/10	-
2	OLC	B	509	-	-	4/5/5/24	-
2	OLC	B	506	-	-	3/5/5/24	-
2	OLC	B	510	-	-	8/13/13/24	-
2	OLC	A	512	-	-	2/8/8/24	-
2	OLC	B	513	2	-	6/14/14/24	-

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	512	OLC	O20-C1	3.14	1.41	1.30
2	A	514	OLC	O20-C1	2.79	1.41	1.33
2	A	501	OLC	O20-C1	2.74	1.41	1.33
2	B	502	OLC	O20-C1	2.71	1.41	1.33
2	B	503	OLC	O20-C1	2.70	1.41	1.33

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	514	OLC	O20-C1-C2	3.24	121.71	111.83
2	B	516	OLC	O20-C1-C2	2.79	120.35	111.83
2	A	501	OLC	O20-C1-C2	2.78	120.30	111.83
3	B	518	3PG	O2-C1-C2	2.72	118.48	112.74
2	B	502	OLC	O20-C1-C2	2.68	120.02	111.83

There are no chirality outliers.

5 of 185 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	OLC	C21-C22-C24-O25
2	A	514	OLC	O20-C21-C22-O23
2	A	515	OLC	C6-C7-C8-C9
2	B	505	OLC	C21-C22-C24-O25
3	A	516	3PG	C3-O1P-P-O3P

There are no ring outliers.

20 monomers are involved in 36 short contacts:

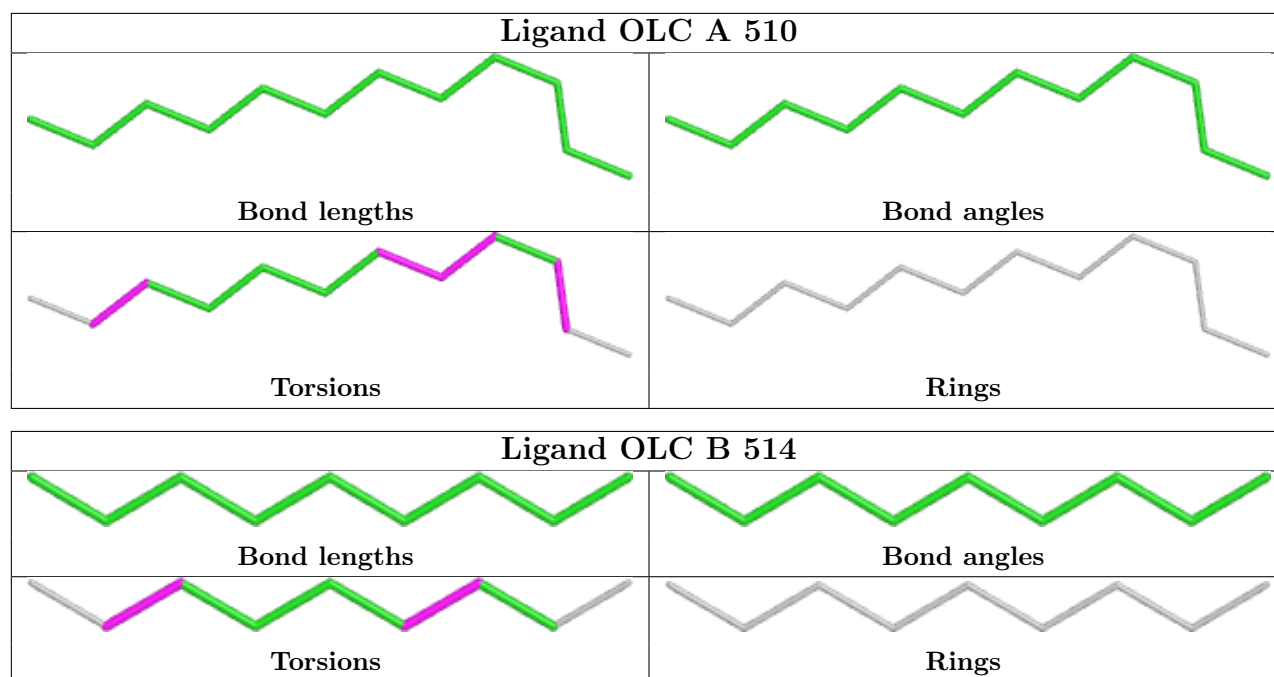
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	510	OLC	1	0
2	B	514	OLC	2	0
2	A	502	OLC	1	0
2	A	515	OLC	1	0
2	A	505	OLC	4	0
2	B	505	OLC	2	0
2	B	503	OLC	3	0
2	B	501	OLC	2	0
2	B	504	OLC	3	0
4	A	517	FLC	1	0
2	A	514	OLC	2	0

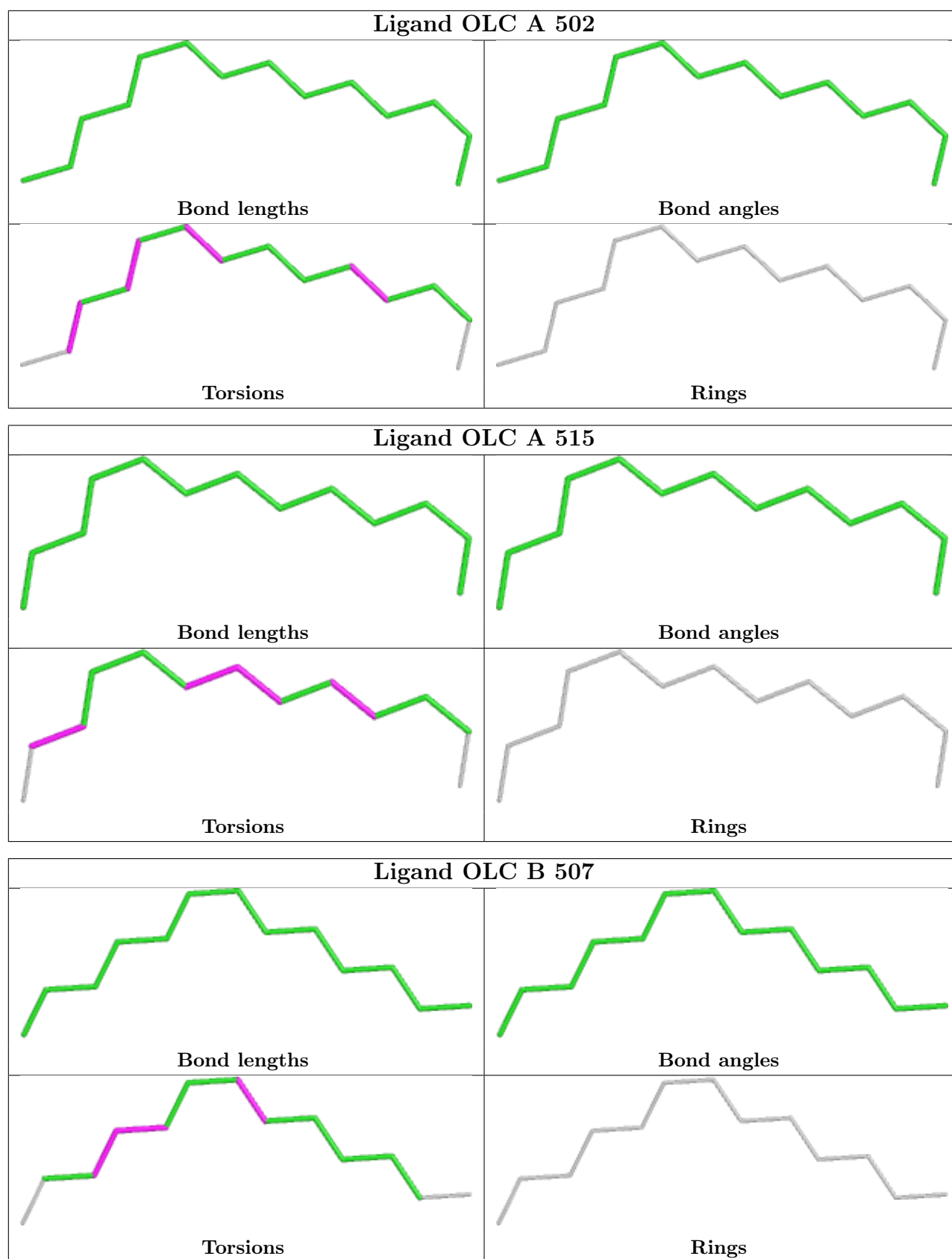
Continued on next page...

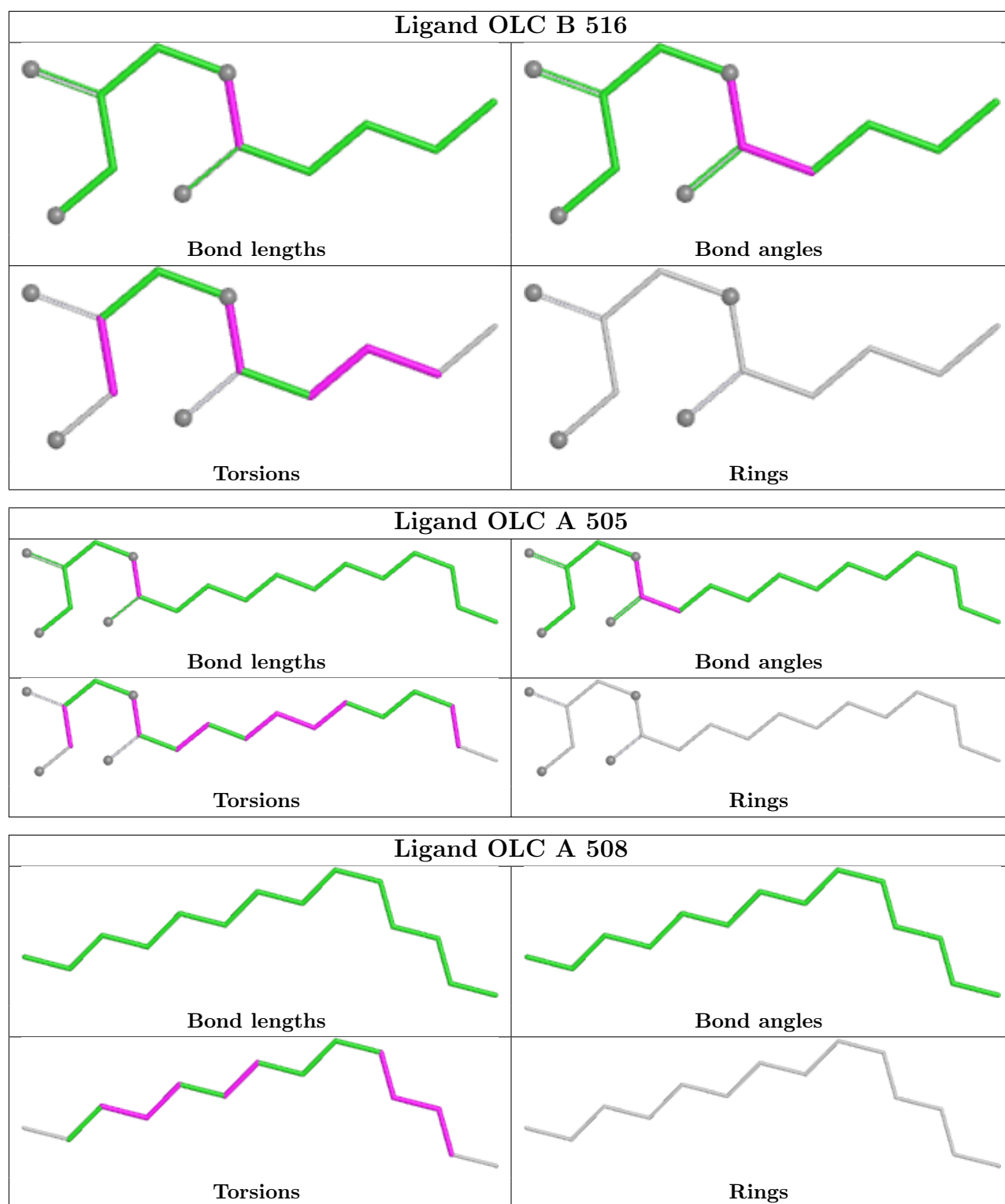
Continued from previous page...

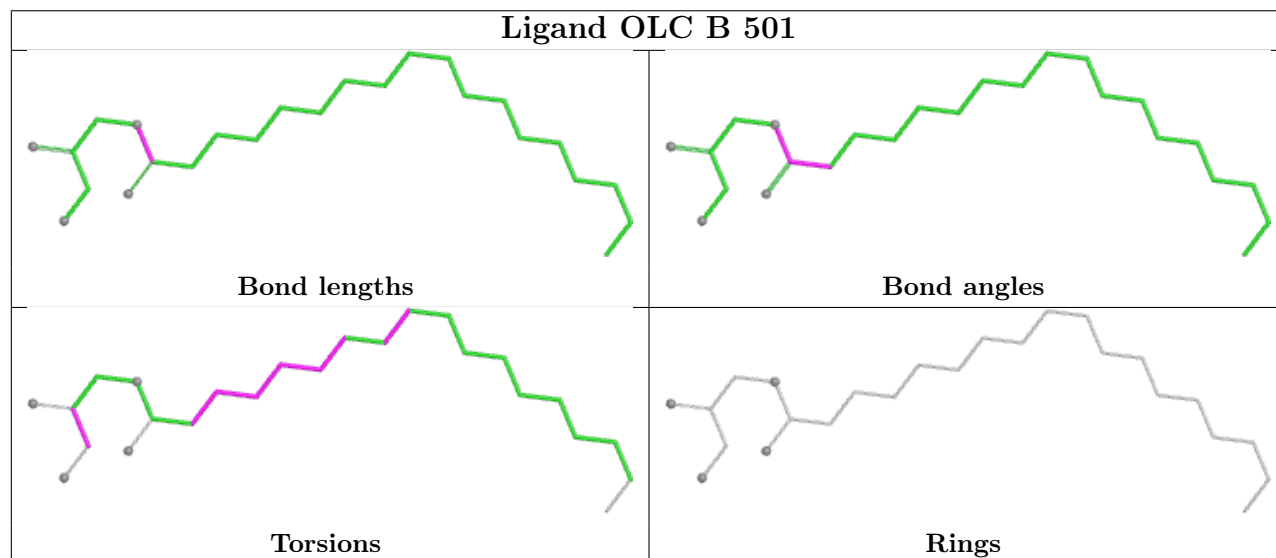
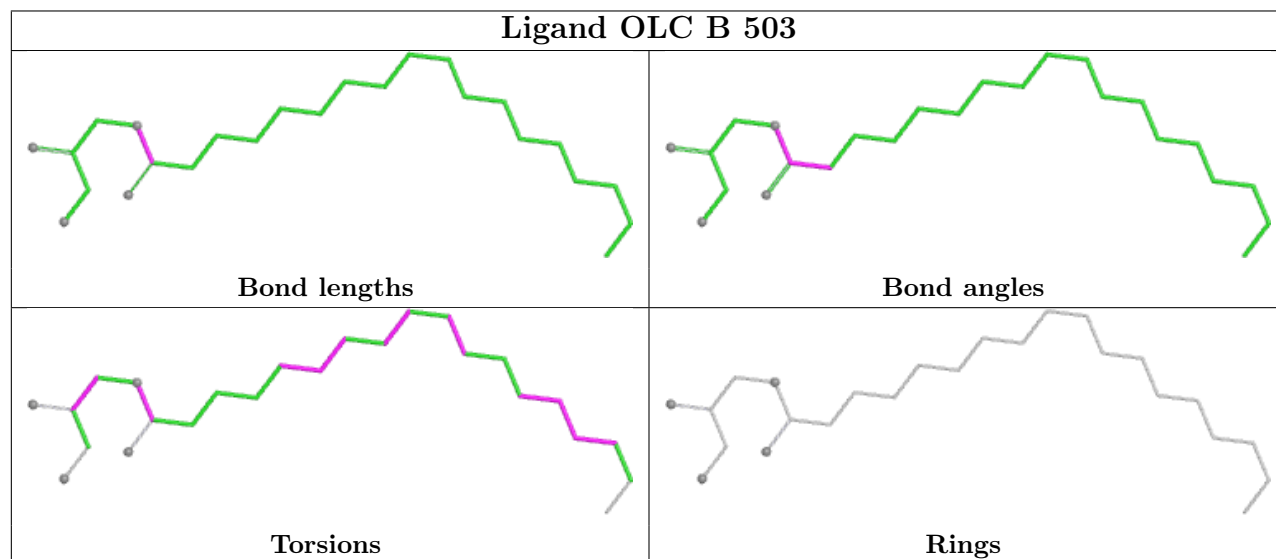
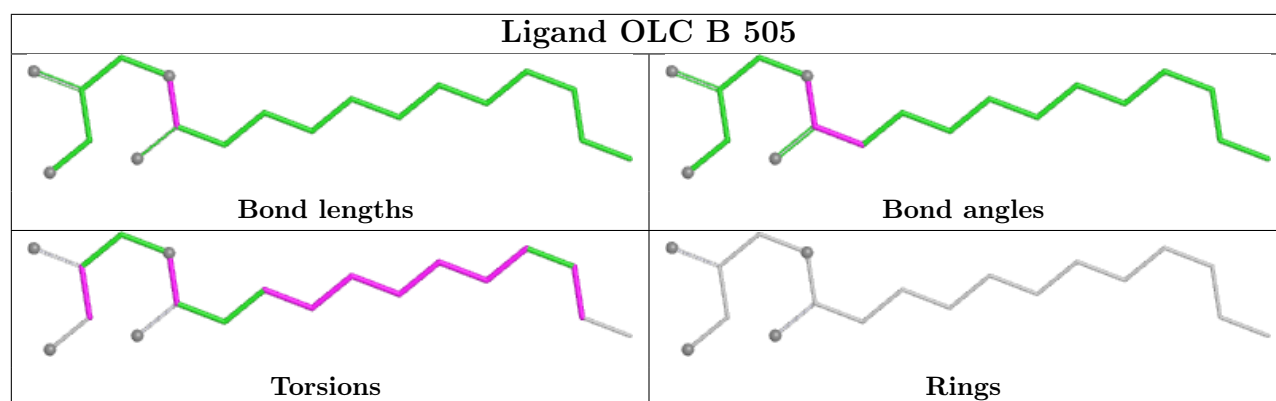
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	OLC	3	0
2	A	507	OLC	1	0
2	B	517	OLC	2	0
2	A	504	OLC	1	0
2	B	502	OLC	1	0
2	B	512	OLC	1	0
2	B	508	OLC	2	0
2	A	512	OLC	2	0
2	B	513	OLC	3	0

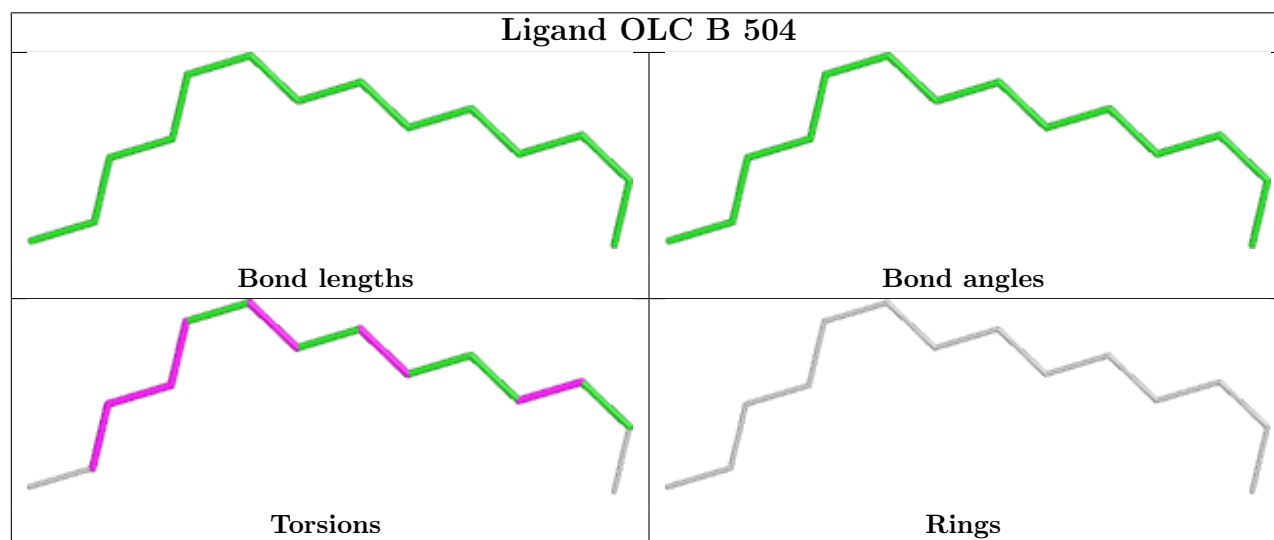
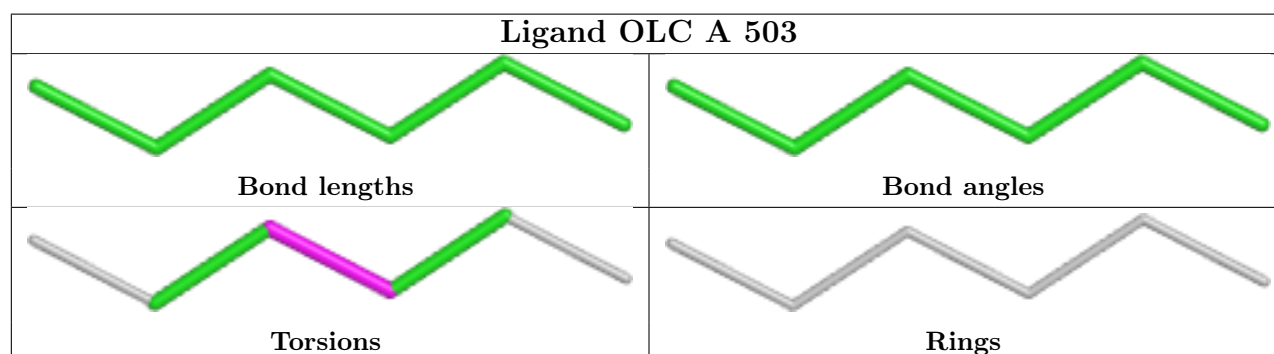
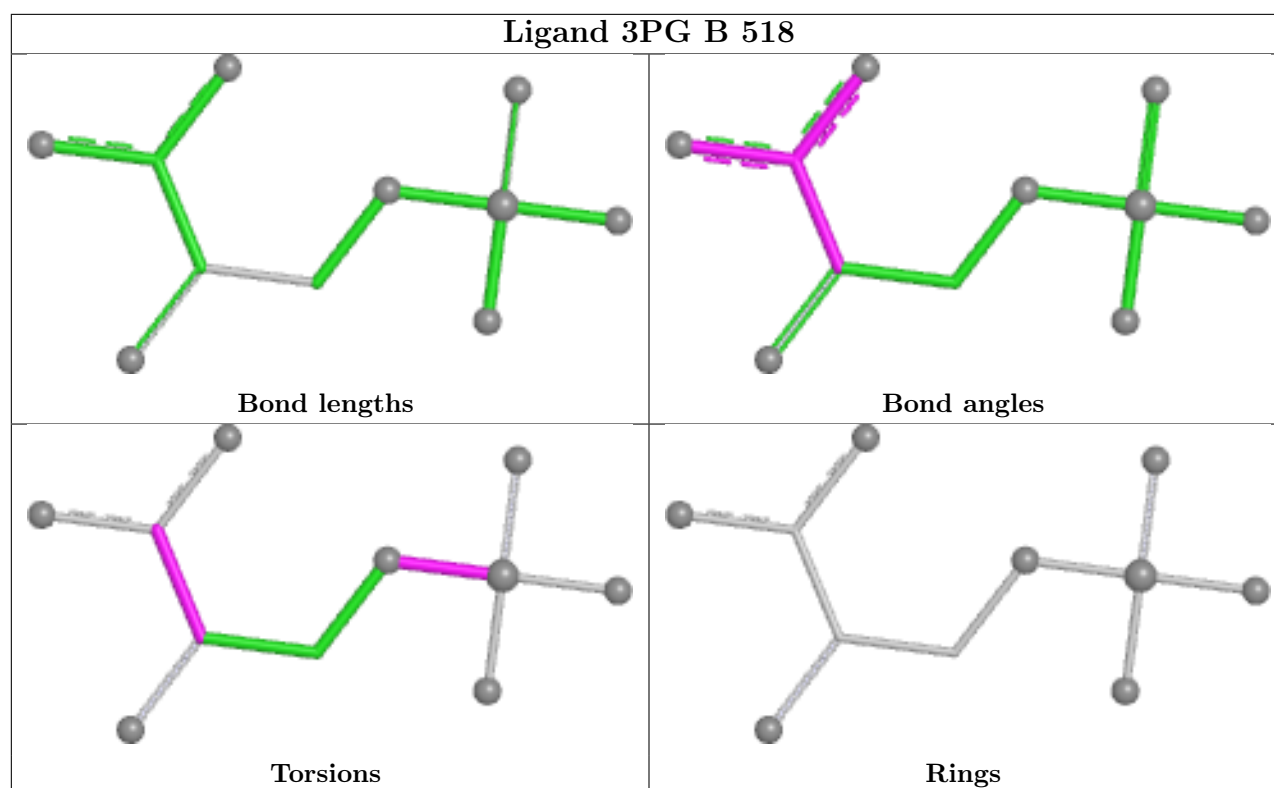
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

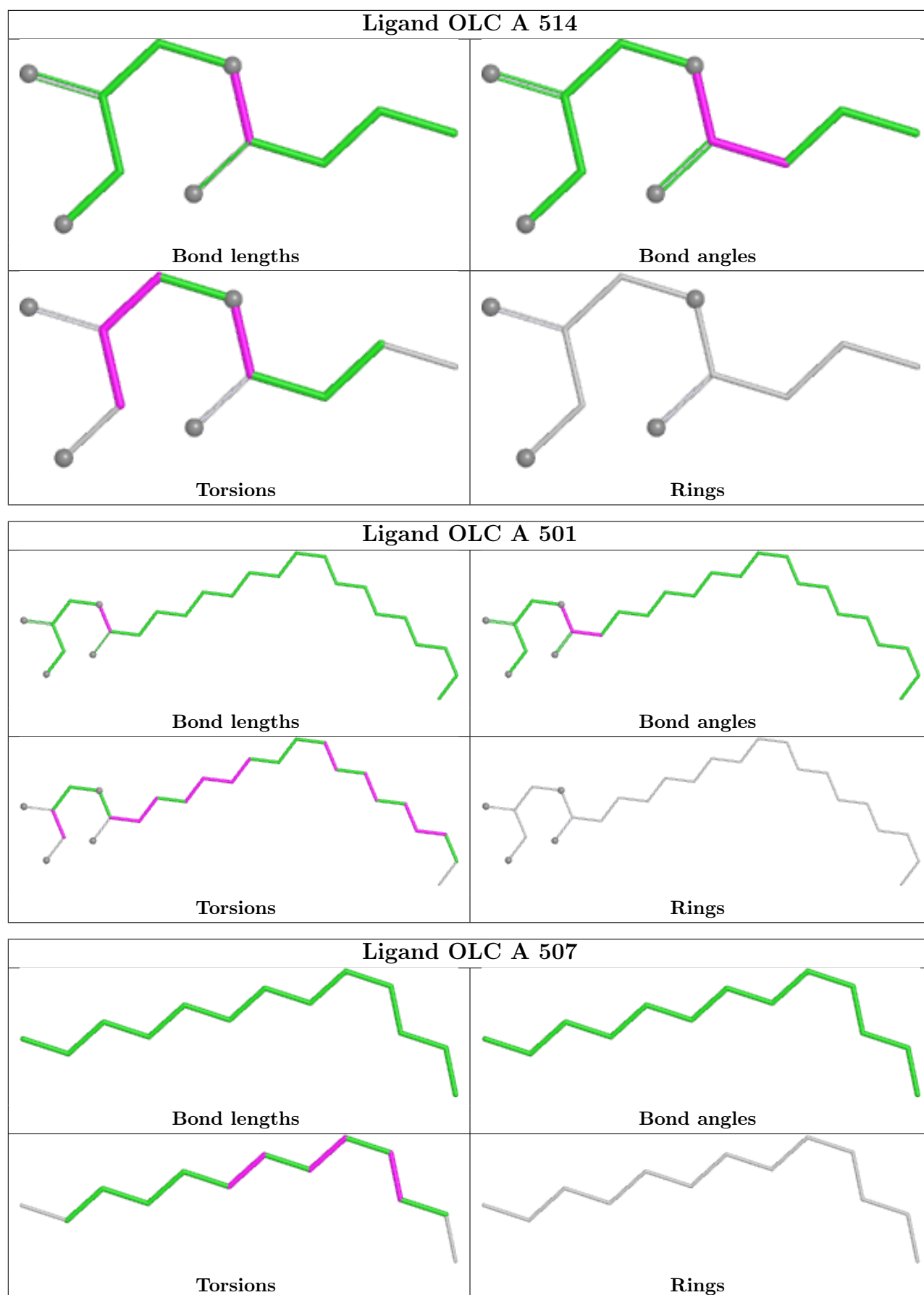


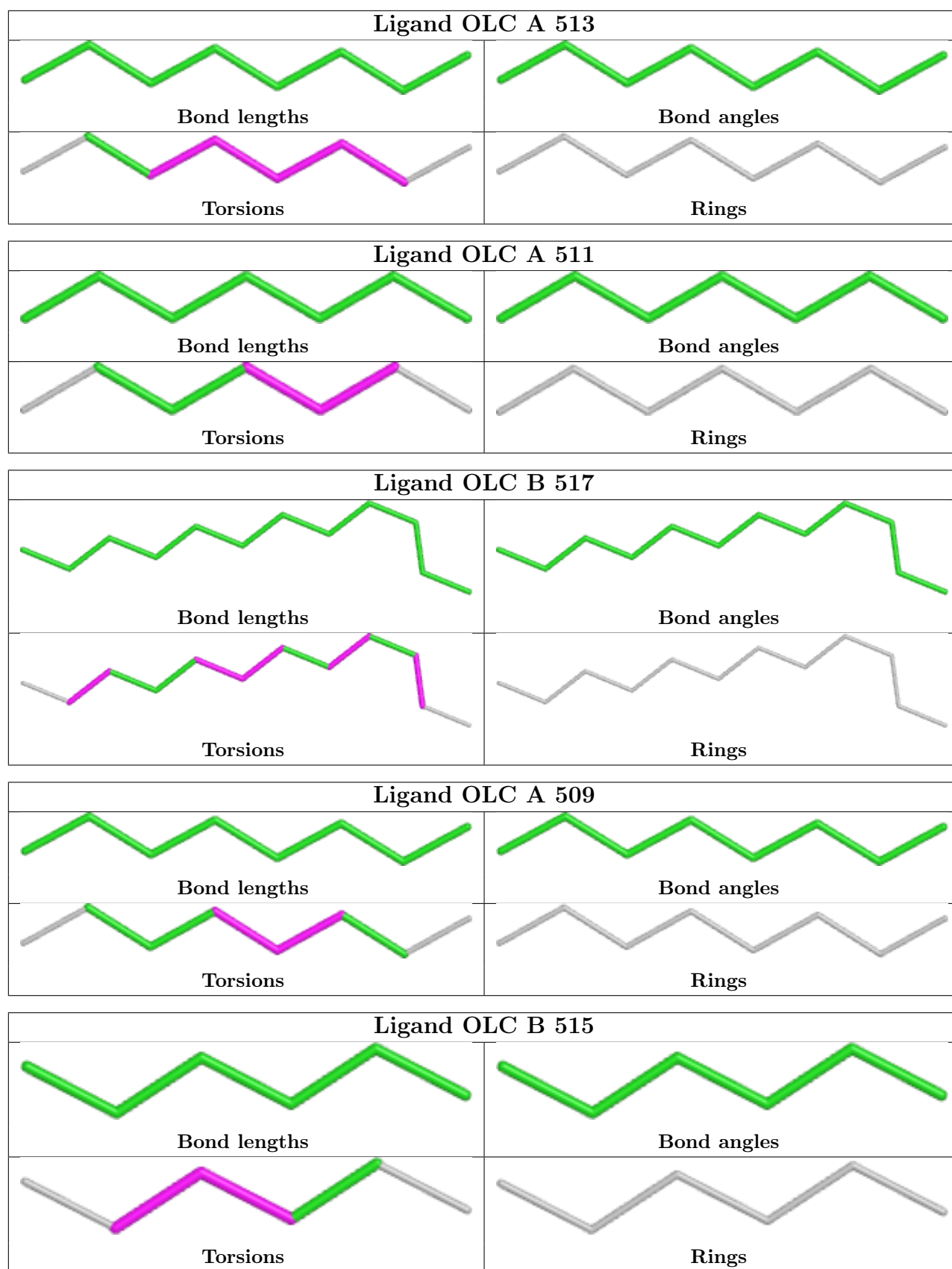


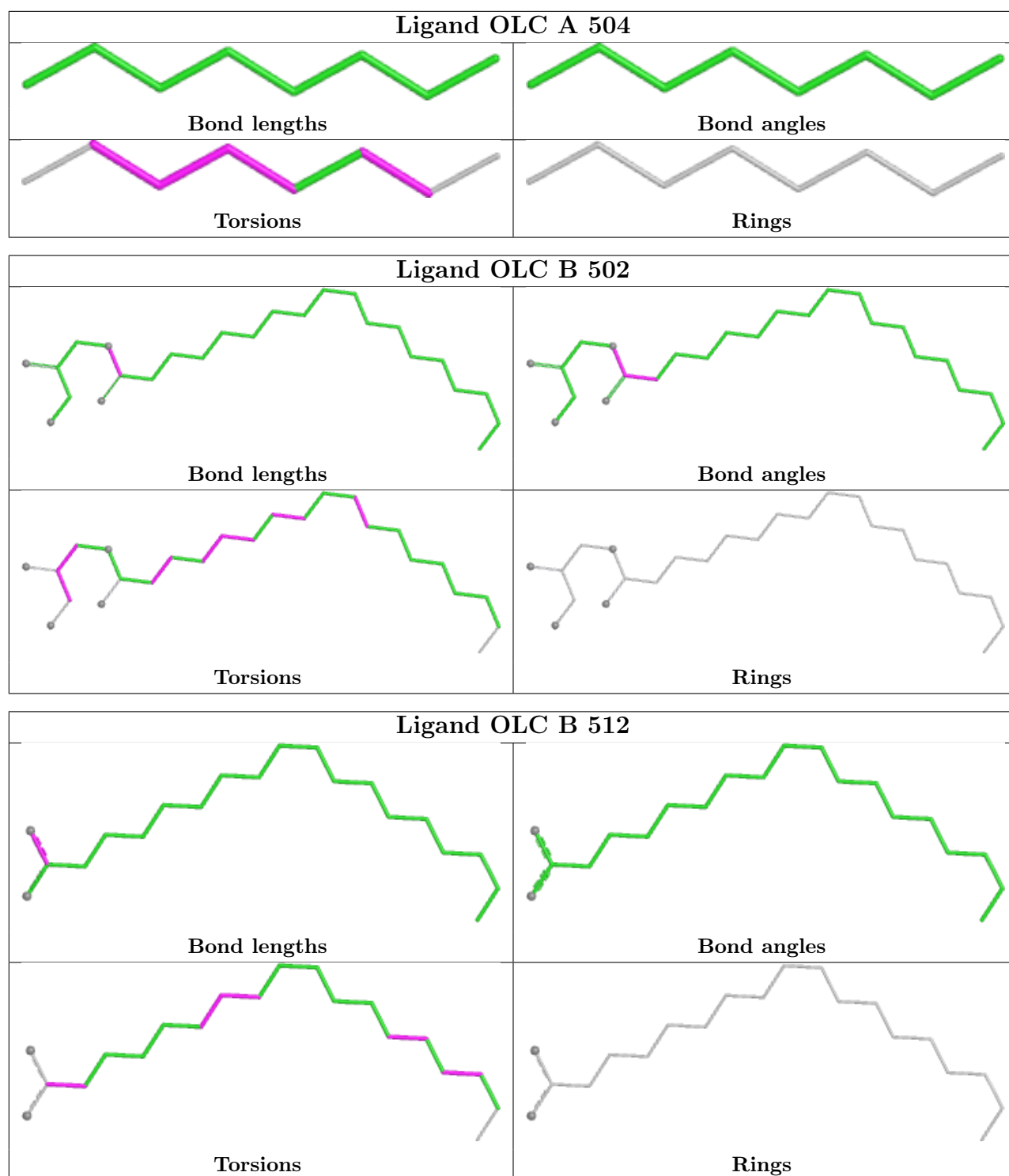


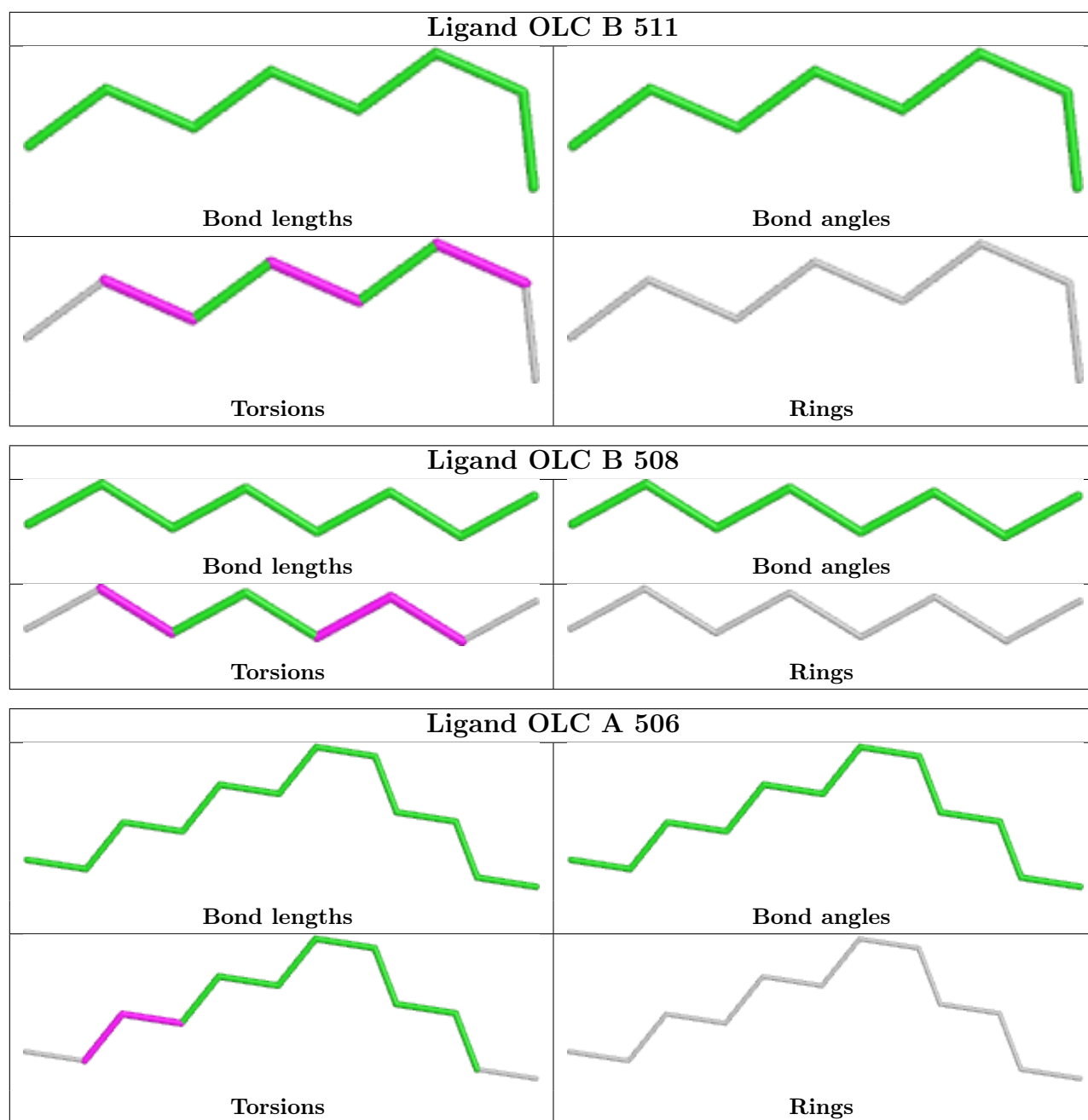


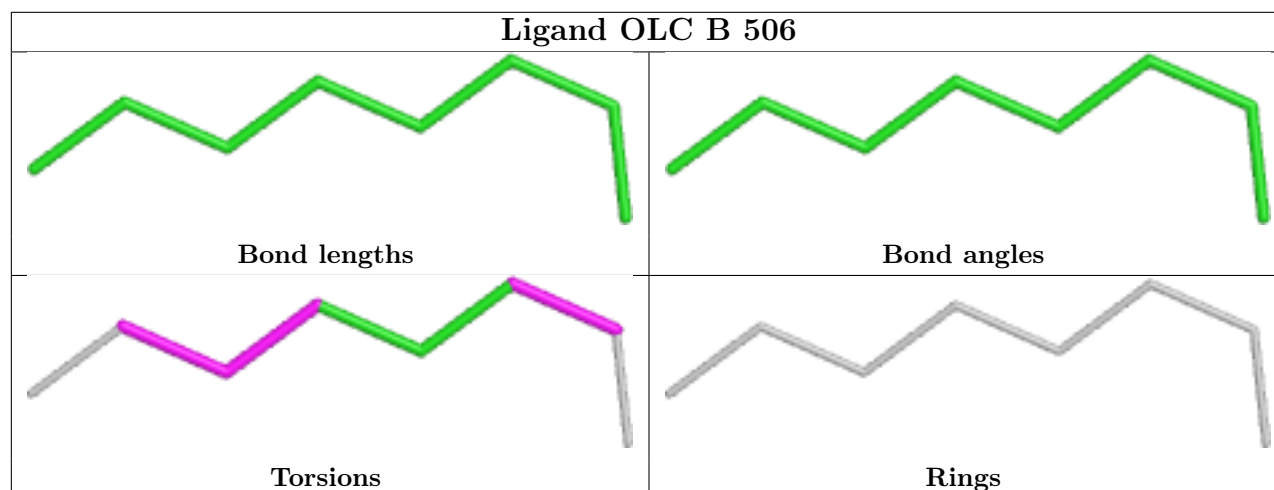
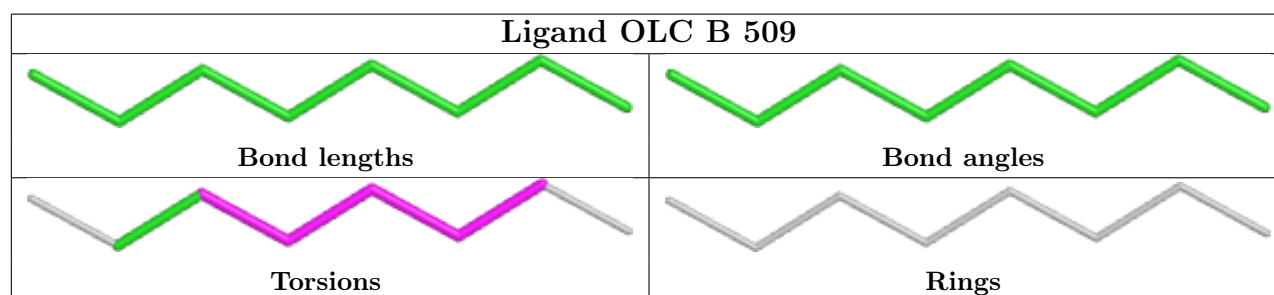
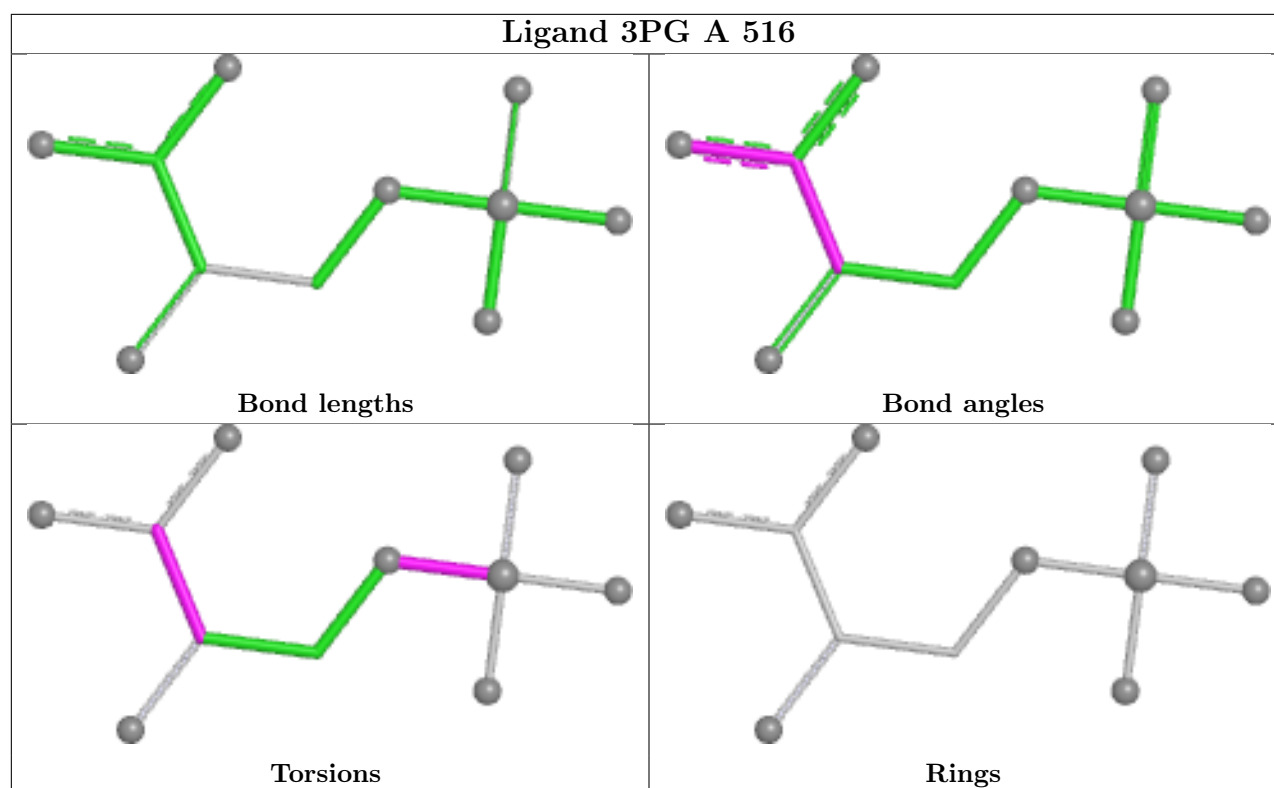


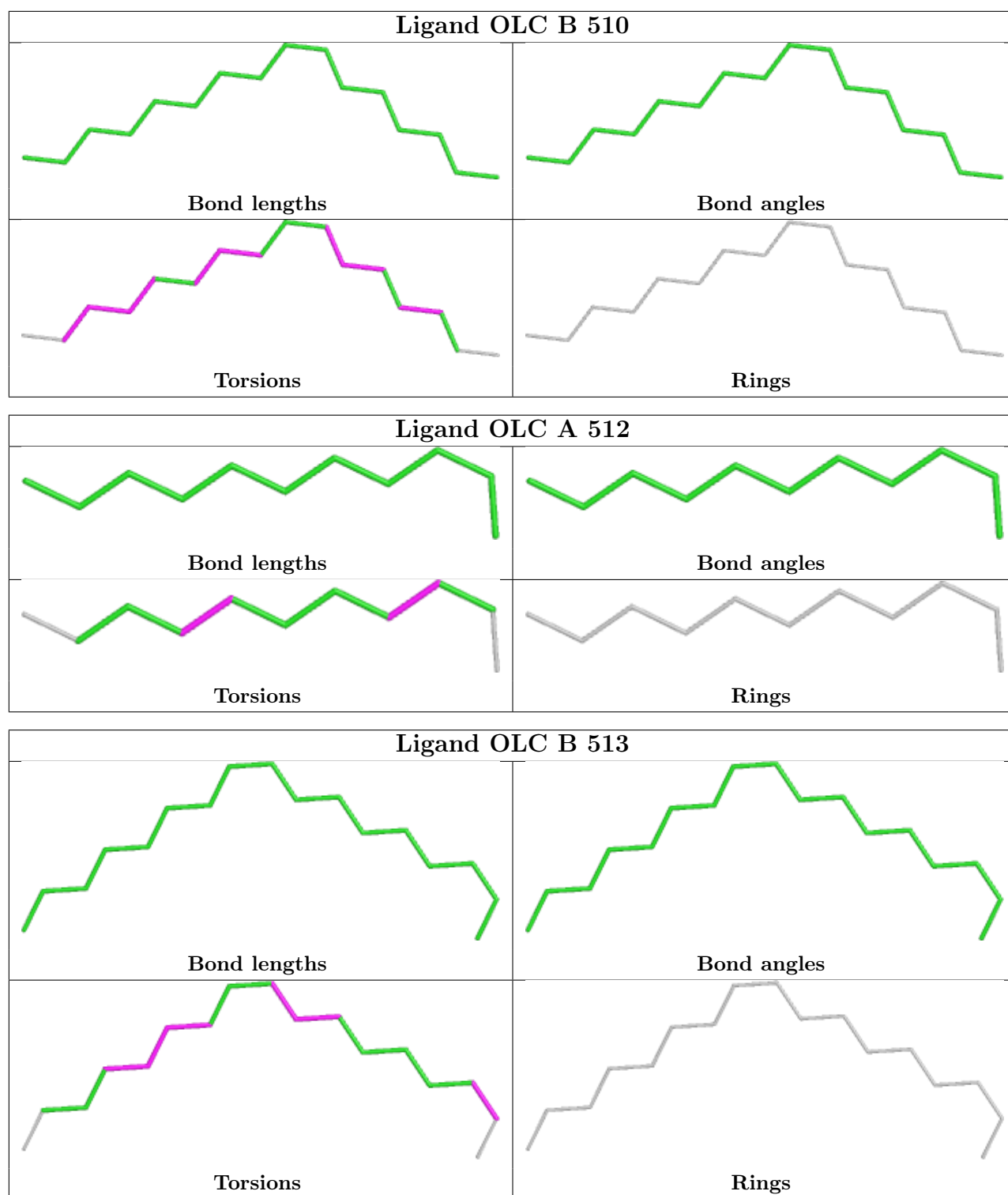












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	305/329 (92%)	0.04	7 (2%) 61 58	16, 27, 47, 66	0
1	B	305/329 (92%)	-0.10	6 (1%) 65 62	18, 26, 43, 70	0
All	All	610/658 (92%)	-0.03	13 (2%) 63 60	16, 26, 46, 70	0

The worst 5 of 13 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	SER	5.3
1	B	100	SER	3.5
1	A	404	PRO	3.4
1	A	383	PHE	3.0
1	B	101	PRO	2.9

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

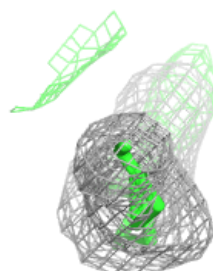
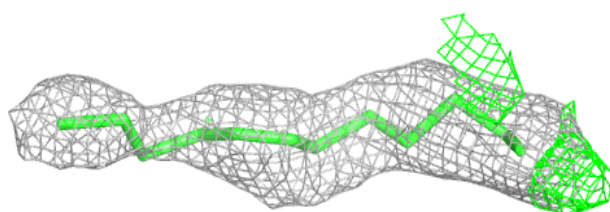
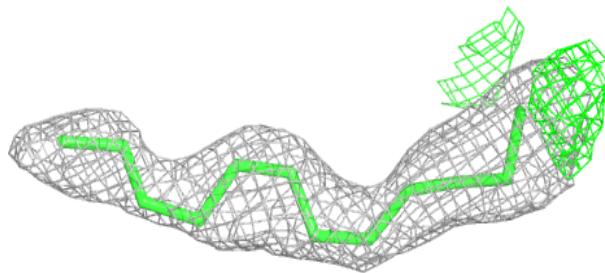
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	OLC	A	506	12/25	0.72	0.19	38,42,51,51	0
2	OLC	A	514	11/25	0.72	0.18	35,49,56,59	0
2	OLC	B	515	6/25	0.74	0.21	36,42,42,44	0
2	OLC	B	514	9/25	0.75	0.16	31,38,42,44	0
2	OLC	A	512	11/25	0.75	0.17	37,40,45,46	0
2	OLC	B	504	14/25	0.76	0.19	35,43,48,48	0
2	OLC	A	511	7/25	0.77	0.19	33,39,42,44	0
2	OLC	A	502	14/25	0.77	0.18	36,44,47,49	0
2	OLC	B	510	16/25	0.78	0.18	43,48,51,52	0
2	OLC	B	516	12/25	0.78	0.15	52,54,61,63	0
2	OLC	B	517	12/25	0.78	0.16	38,43,47,48	0
2	OLC	B	503	25/25	0.79	0.15	34,41,47,58	0
2	OLC	B	513	17/25	0.79	0.15	33,38,45,45	0
2	OLC	B	507	13/25	0.79	0.16	31,43,48,51	0
2	OLC	B	508	8/25	0.80	0.15	41,43,45,52	0
2	OLC	B	509	8/25	0.80	0.16	34,38,42,42	0
2	OLC	B	511	8/25	0.81	0.17	29,37,42,43	0
4	FLC	A	517	13/13	0.81	0.11	35,42,50,54	0
2	OLC	A	507	13/25	0.82	0.14	34,38,43,46	0
2	OLC	A	513	8/25	0.82	0.16	32,43,49,51	0
2	OLC	A	510	12/25	0.82	0.14	35,41,46,49	0
2	OLC	A	515	13/25	0.82	0.16	36,42,47,47	0
2	OLC	B	502	25/25	0.82	0.13	28,40,48,55	0
2	OLC	A	504	8/25	0.82	0.14	33,36,39,43	0
2	OLC	A	503	6/25	0.83	0.14	30,36,40,49	0
2	OLC	B	512	20/25	0.83	0.14	23,42,51,52	0
2	OLC	A	508	14/25	0.84	0.15	29,39,46,47	0
2	OLC	B	501	25/25	0.85	0.13	27,37,57,62	0
2	OLC	A	501	25/25	0.85	0.14	30,38,49,51	0
2	OLC	B	506	8/25	0.86	0.15	30,39,46,47	0
2	OLC	B	505	19/25	0.88	0.12	29,32,42,48	0
2	OLC	A	509	8/25	0.88	0.12	32,36,41,43	0
2	OLC	A	505	19/25	0.91	0.11	26,32,42,47	0
3	3PG	B	518	11/11	0.97	0.06	19,21,27,28	0
3	3PG	A	516	11/11	0.98	0.05	19,22,27,27	0

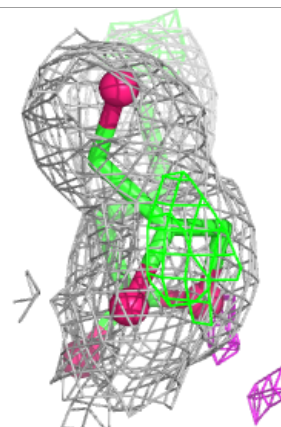
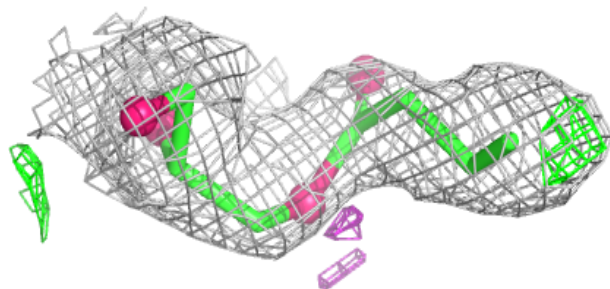
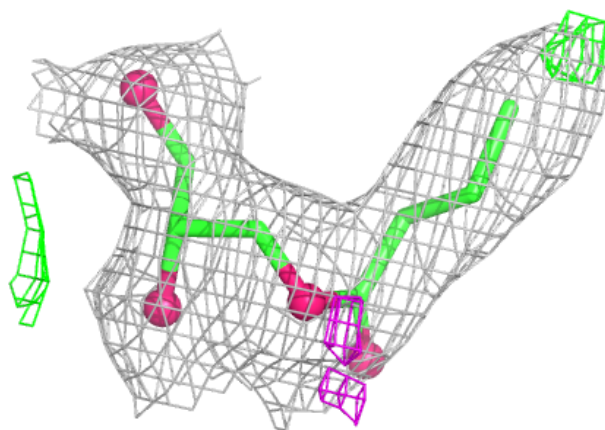
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around OLC A 506:

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

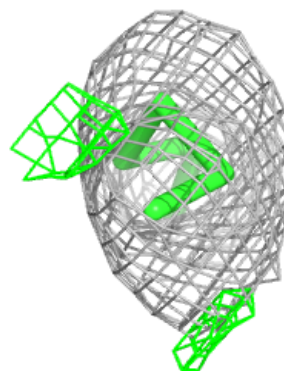
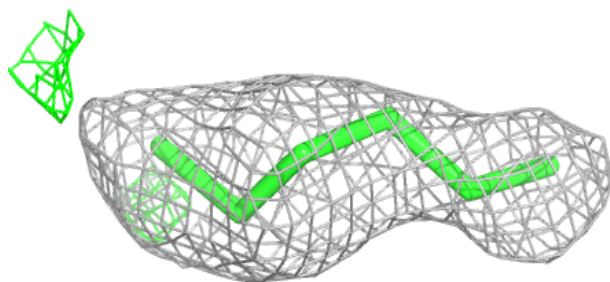
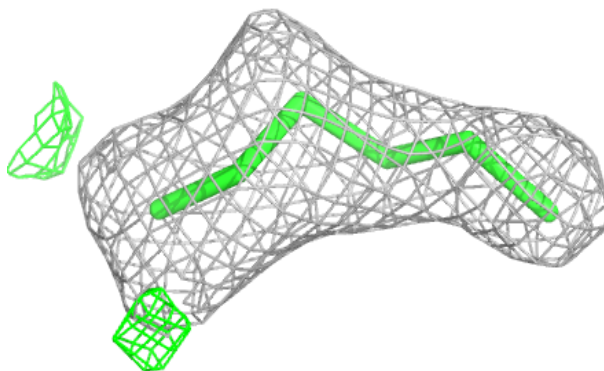
**Electron density around OLC A 514:**

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

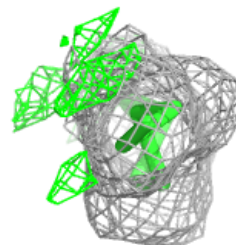
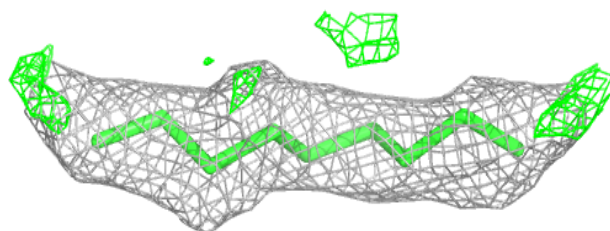
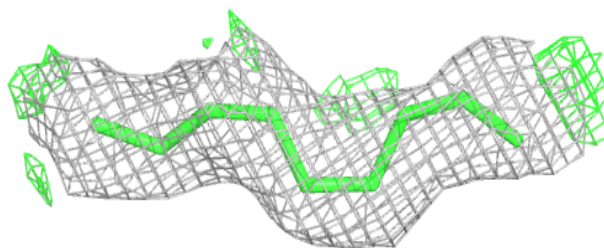


Electron density around OLC B 515:

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

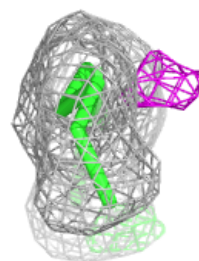
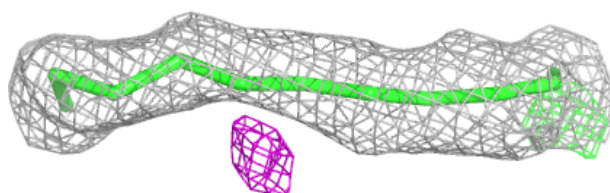
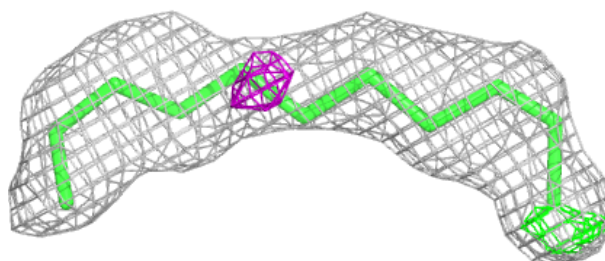
**Electron density around OLC B 514:**

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

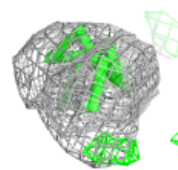
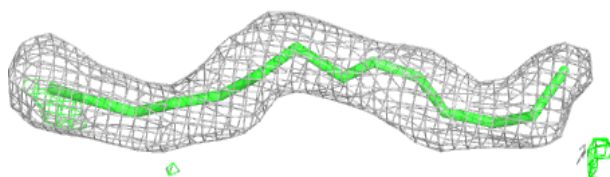
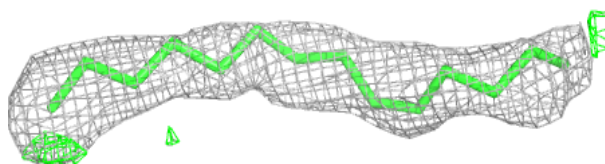


Electron density around OLC A 512:

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

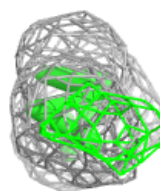
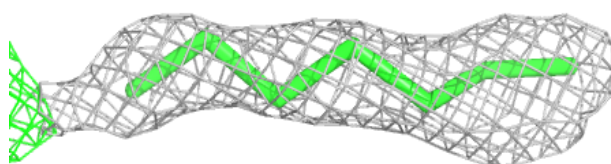
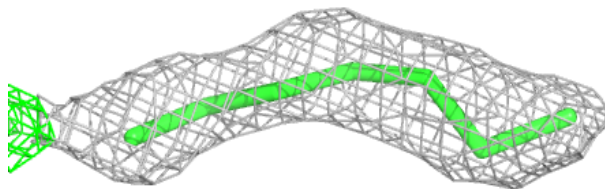
**Electron density around OLC B 504:**

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

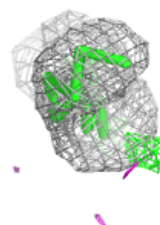
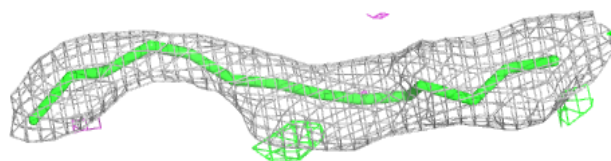
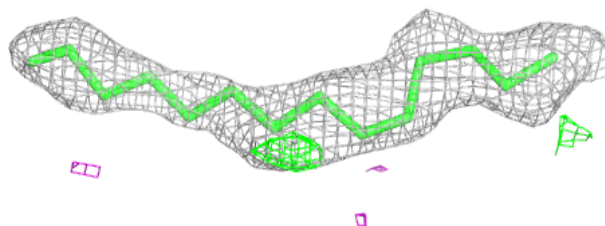


Electron density around OLC A 511:

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

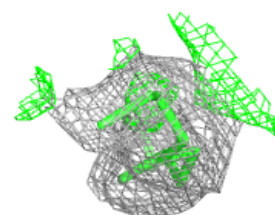
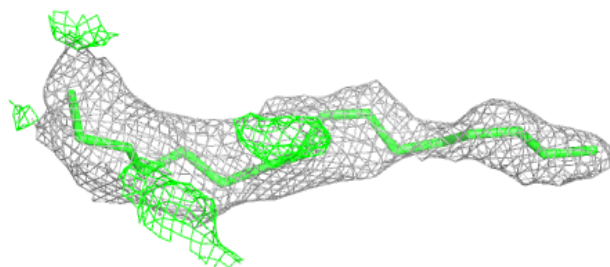
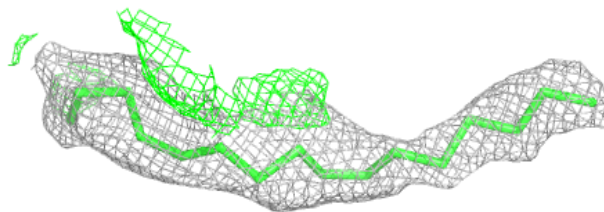
**Electron density around OLC A 502:**

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

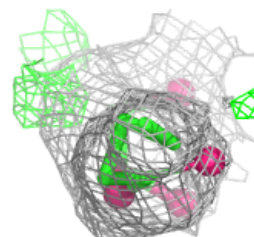
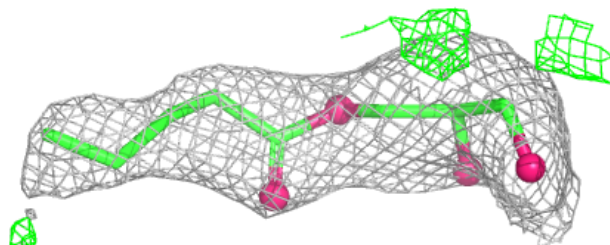
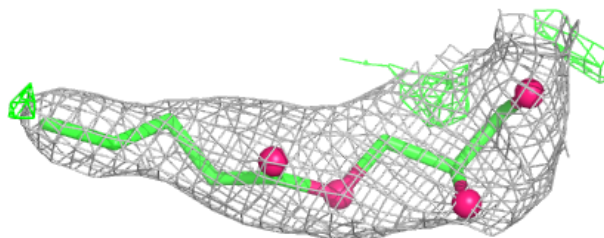


Electron density around OLC B 510:

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

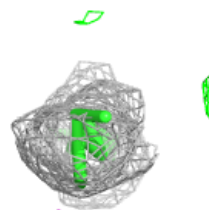
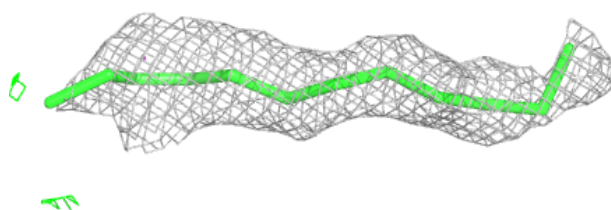
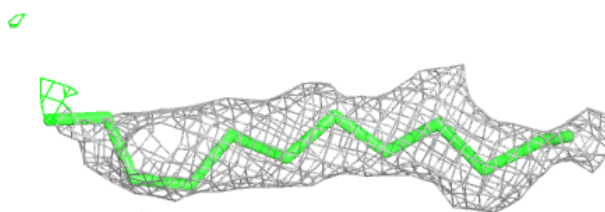
**Electron density around OLC B 516:**

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

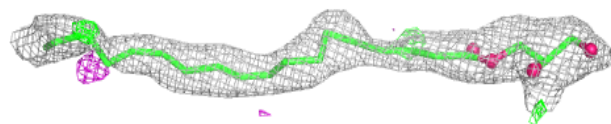
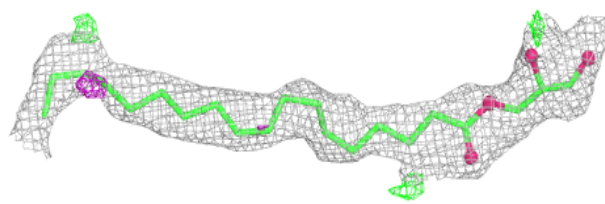


Electron density around OLC B 517:

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

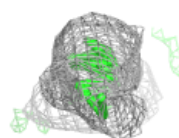
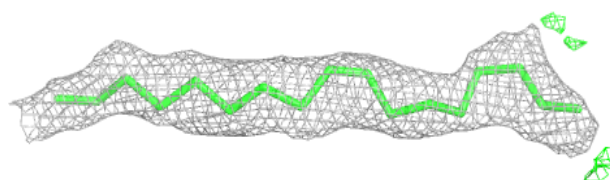
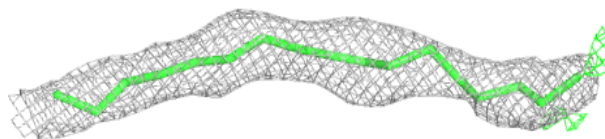
**Electron density around OLC B 503:**

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

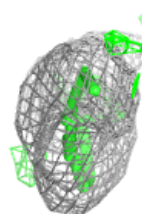
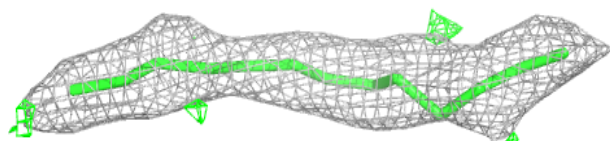
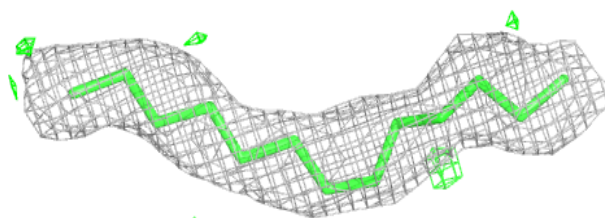


Electron density around OLC B 513:

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

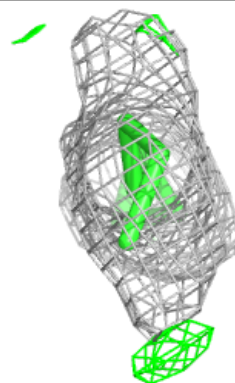
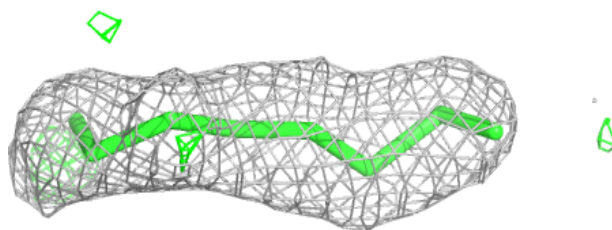
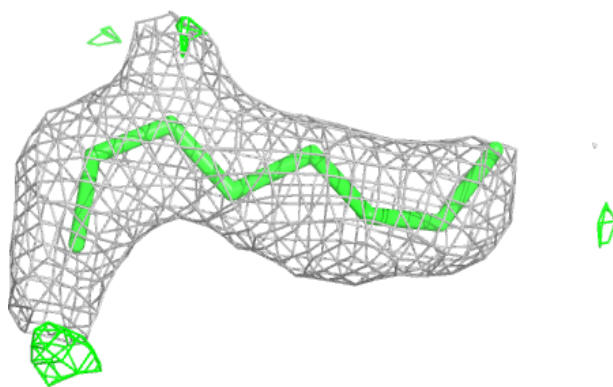
**Electron density around OLC B 507:**

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

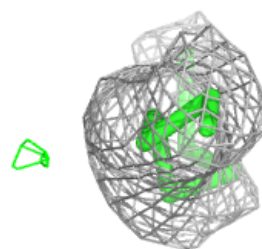
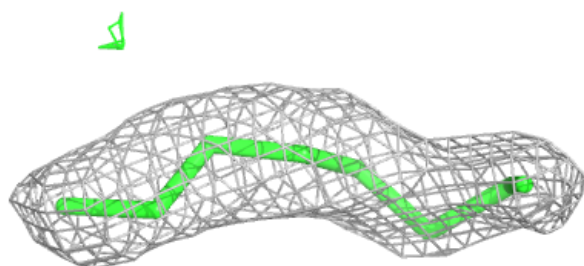
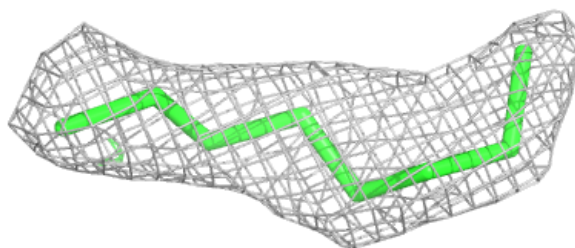


Electron density around OLC B 508:

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

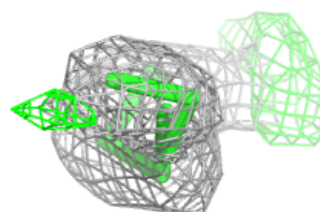
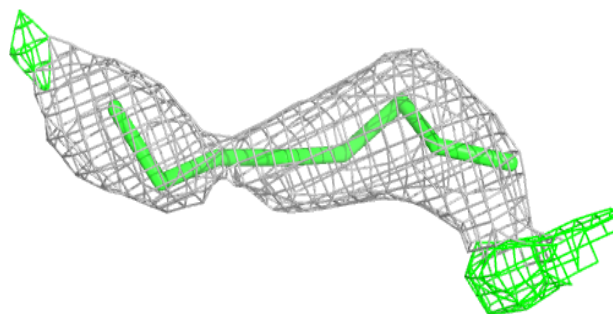
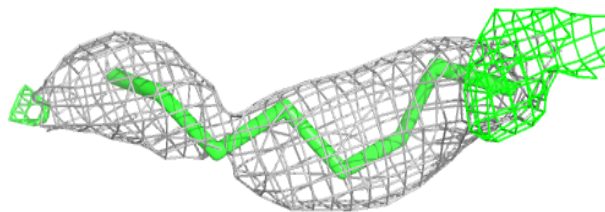
**Electron density around OLC B 509:**

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

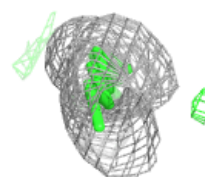
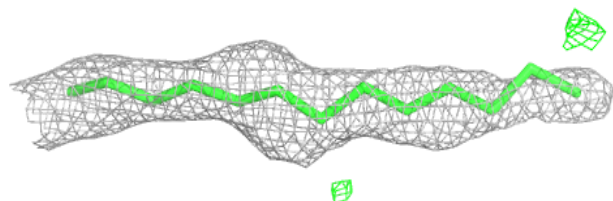
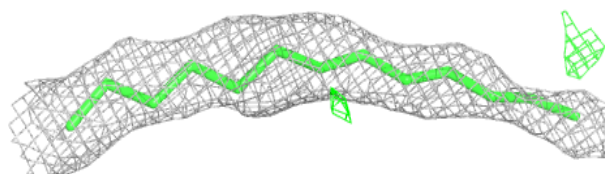


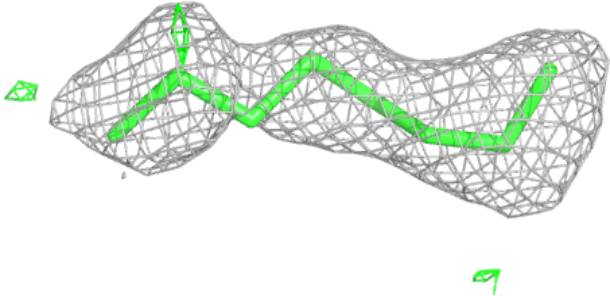
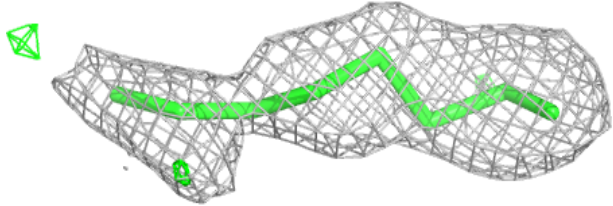
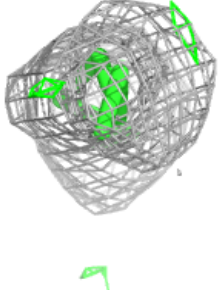
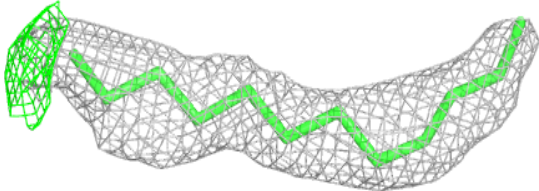
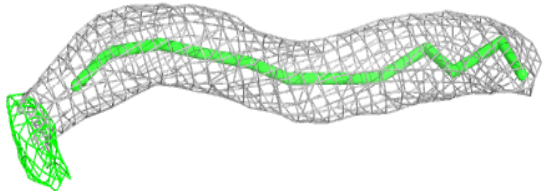
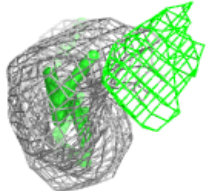
Electron density around OLC B 511:

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

**Electron density around OLC A 507:**

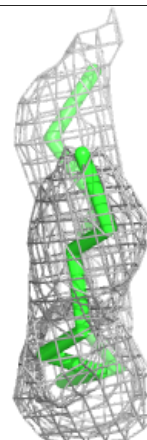
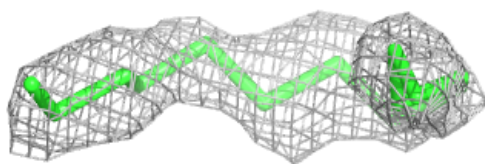
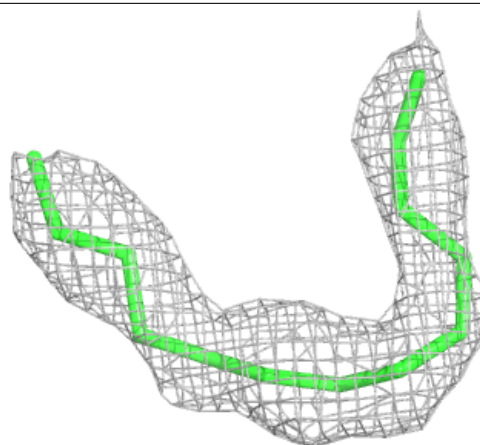
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



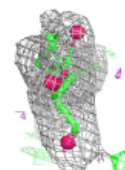
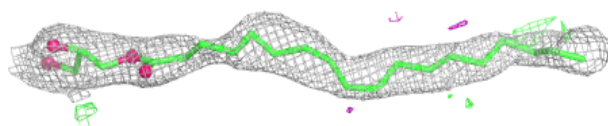
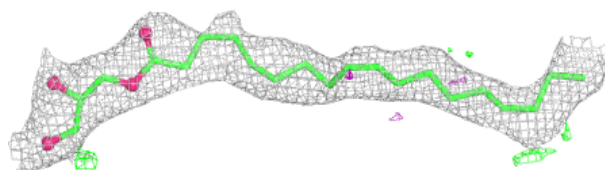
Electron density around OLC A 513: $2mF_o-DF_c$ (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)	 A 3D electron density map of OLC A 513. The map is shown as a gray wireframe mesh. A green stick model of the molecule is overlaid on the map. A small green arrow points to a specific feature on the molecule.
 A 3D electron density map of OLC A 513, showing a different view or a different part of the molecule. The map is shown as a gray wireframe mesh. A green stick model of the molecule is overlaid on the map. A small green arrow points to a specific feature on the molecule.	 A 3D electron density map of OLC A 513, showing a different view or a different part of the molecule. The map is shown as a gray wireframe mesh. A green stick model of the molecule is overlaid on the map. A small green arrow points to a specific feature on the molecule.
Electron density around OLC A 510: $2mF_o-DF_c$ (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)	 A 3D electron density map of OLC A 510. The map is shown as a gray wireframe mesh. A green stick model of the molecule is overlaid on the map. A small green arrow points to a specific feature on the molecule.
 A 3D electron density map of OLC A 510, showing a different view or a different part of the molecule. The map is shown as a gray wireframe mesh. A green stick model of the molecule is overlaid on the map. A small green arrow points to a specific feature on the molecule.	 A 3D electron density map of OLC A 510, showing a different view or a different part of the molecule. The map is shown as a gray wireframe mesh. A green stick model of the molecule is overlaid on the map. A small green arrow points to a specific feature on the molecule.

Electron density around OLC A 515:

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

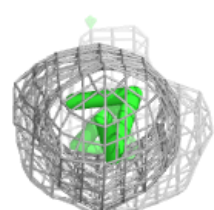
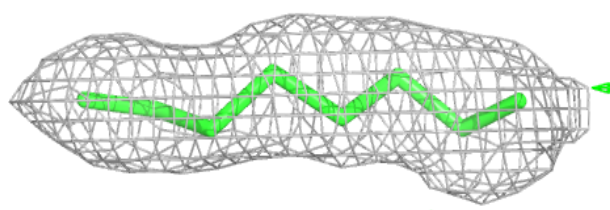
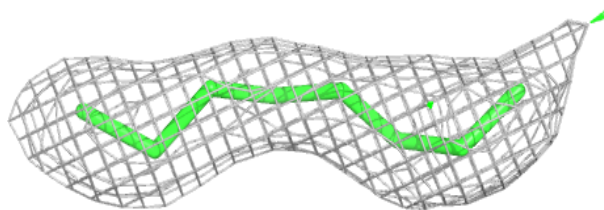
**Electron density around OLC B 502:**

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

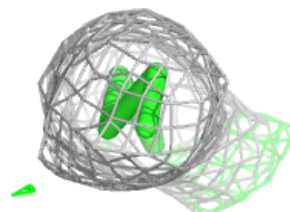
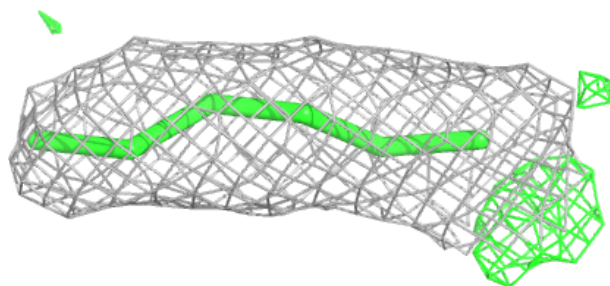
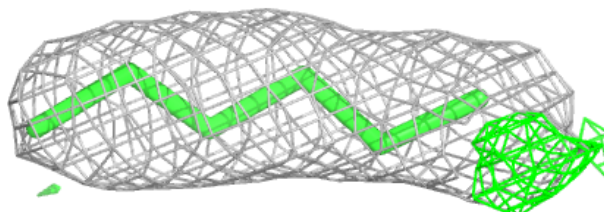


Electron density around OLC A 504:

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

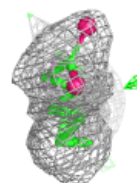
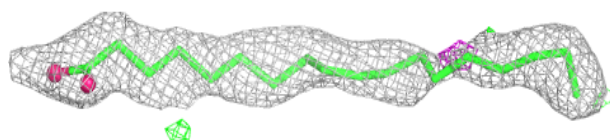
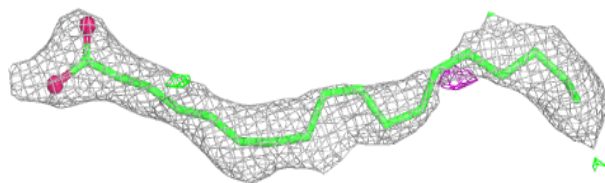
**Electron density around OLC A 503:**

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

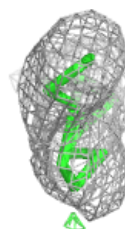
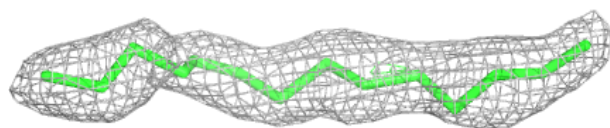
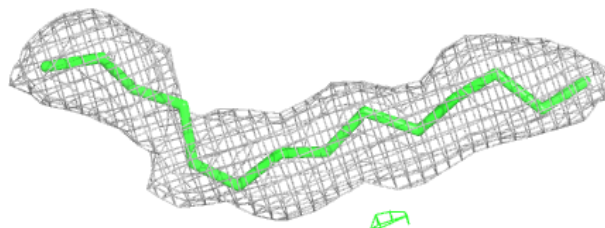


Electron density around OLC B 512:

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

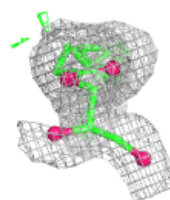
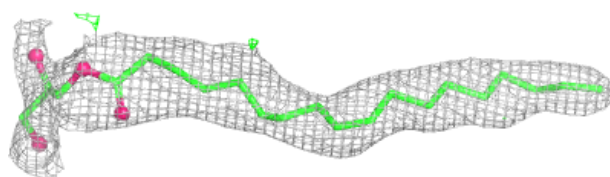
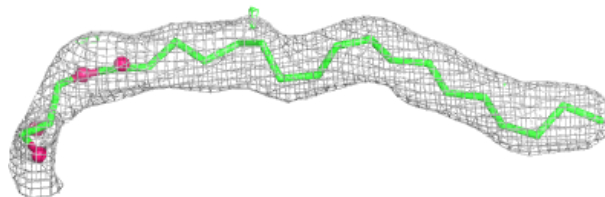
**Electron density around OLC A 508:**

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

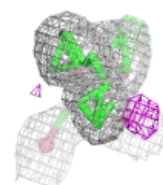
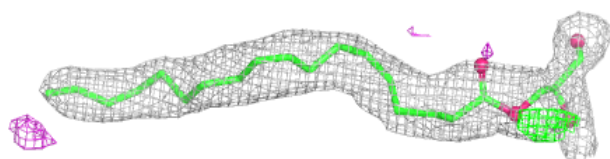
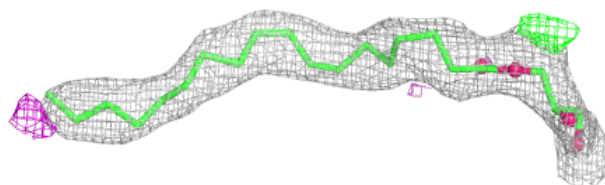


Electron density around OLC B 501:

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

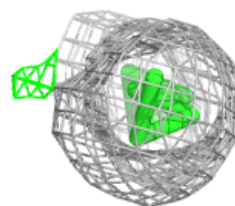
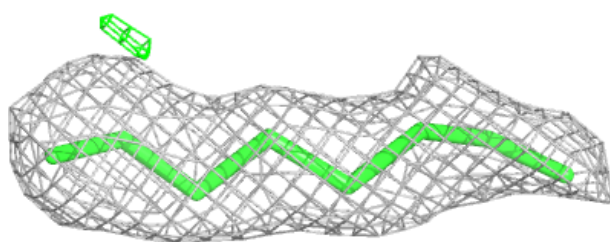
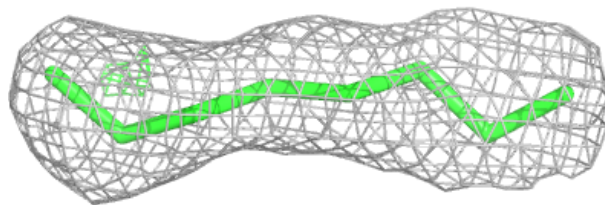
**Electron density around OLC A 501:**

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

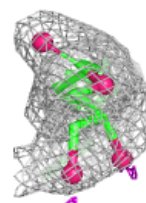
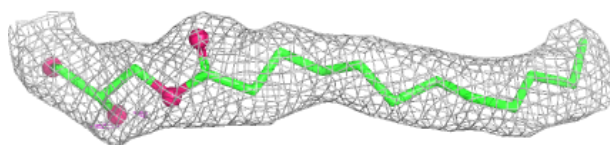
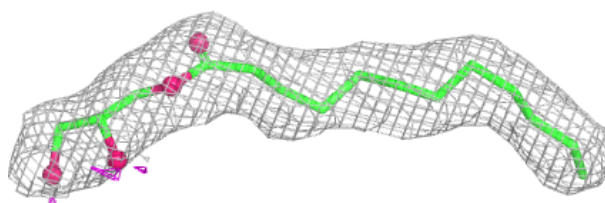


Electron density around OLC B 506:

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

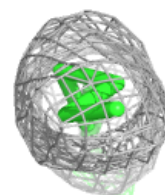
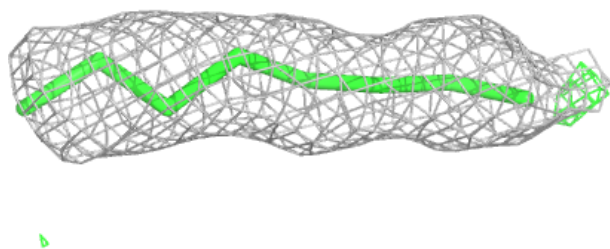
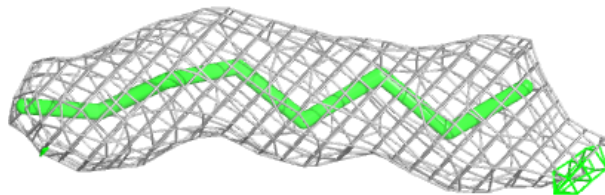
**Electron density around OLC B 505:**

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

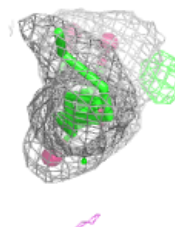
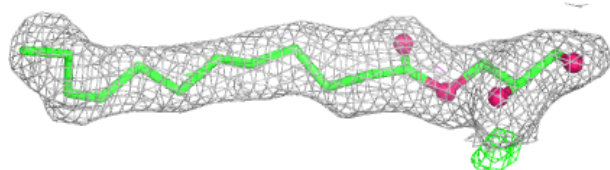
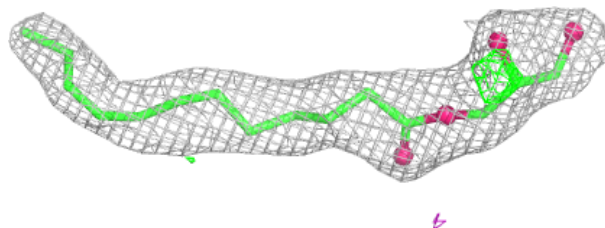


Electron density around OLC A 509:

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

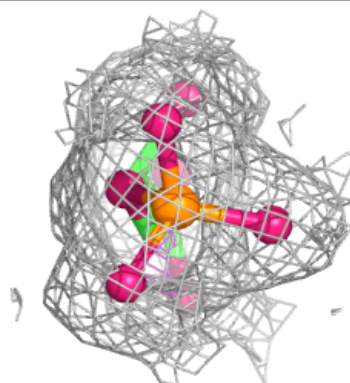
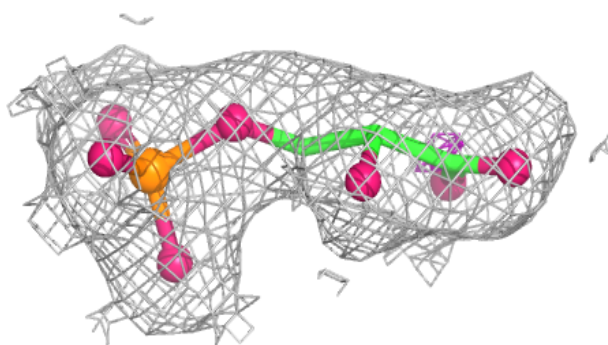
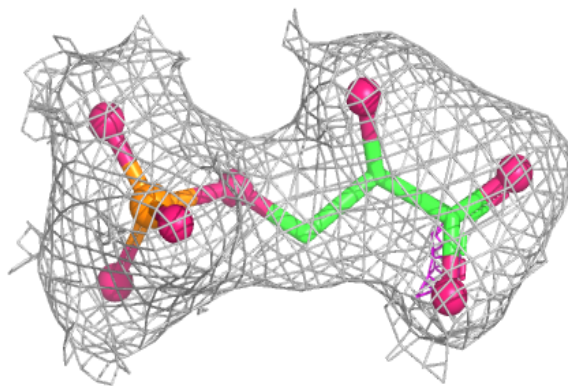
**Electron density around OLC A 505:**

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

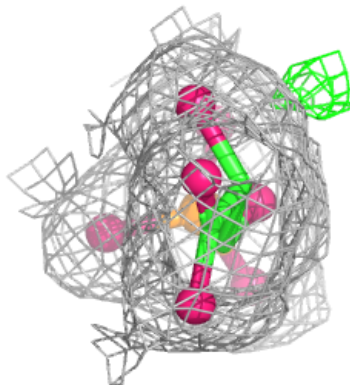
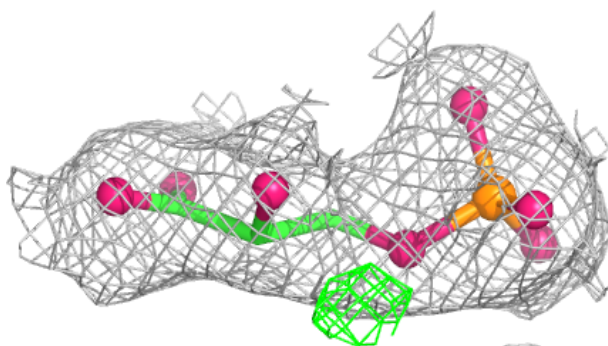
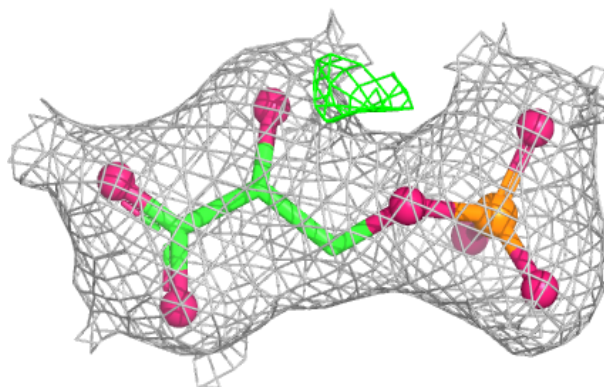


Electron density around 3PG B 518:

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

**Electron density around 3PG A 516:**

$2mF_o-DF_c$ (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.