



# Full wwPDB X-ray Structure Validation Report ⓘ

May 14, 2026 – 12:06 PM EDT

PDB ID : 7IOR / pdb\_00007ior  
Title : Crystal structure of A2A in complex with FU163-53  
Authors : Huang, C.-Y.; Cheng, R.Y.K.; Metz, A.  
Deposited on : 2025-08-21  
Resolution : 1.89 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity : 4-5-2 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

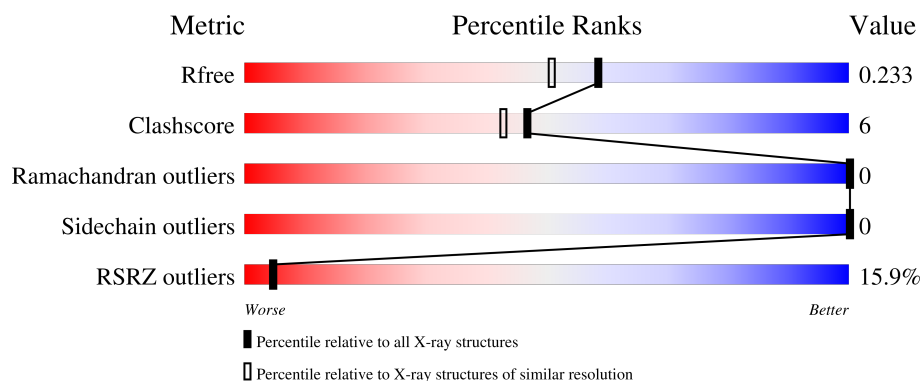
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.89 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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

Mol	Chain	Length	Quality of chain
1	A	433	<div> <div>14%</div> <div>80%</div> <div>10%</div> <div>10%</div> </div>

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	TEP	A	2401	-	X	-	-

## 2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 3763 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 Adenosine receptor A2a/Soluble cytochrome b562/Adenosine receptor A2a chimera.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	390	Total	C	N	O	S	0	13	0
			3126	2035	532	536	23			

There are 35 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-8	ASP	-	expression tag	UNP P29274
A	-7	TYR	-	expression tag	UNP P29274
A	-6	LYS	-	expression tag	UNP P29274
A	-5	ASP	-	expression tag	UNP P29274
A	-4	ASP	-	expression tag	UNP P29274
A	-3	ASP	-	expression tag	UNP P29274
A	-2	ASP	-	expression tag	UNP P29274
A	-1	GLY	-	expression tag	UNP P29274
A	0	ALA	-	expression tag	UNP P29274
A	1	PRO	-	expression tag	UNP P29274
A	54	LEU	ALA	engineered mutation	UNP P29274
A	88	ALA	THR	engineered mutation	UNP P29274
A	107	ALA	ARG	engineered mutation	UNP P29274
A	122	ALA	LYS	engineered mutation	UNP P29274
A	154	ALA	ASN	engineered mutation	UNP P29274
A	202	ALA	LEU	engineered mutation	UNP P29274
A	1007	TRP	MET	conflict	UNP P0ABE7
A	1102	ILE	-	linker	UNP P0ABE7
A	1103	GLN	-	linker	UNP P0ABE7
A	1104	LYS	-	linker	UNP P0ABE7
A	1105	TYR	-	linker	UNP P0ABE7
A	1106	LEU	-	linker	UNP P0ABE7
A	235	ALA	LEU	engineered mutation	UNP P29274
A	239	ALA	VAL	engineered mutation	UNP P29274
A	318	ALA	-	expression tag	UNP P29274
A	319	HIS	-	expression tag	UNP P29274

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Chain	Residue	Modelled	Actual	Comment	Reference
A	320	HIS	-	expression tag	UNP P29274
A	321	HIS	-	expression tag	UNP P29274
A	322	HIS	-	expression tag	UNP P29274
A	323	HIS	-	expression tag	UNP P29274
A	324	HIS	-	expression tag	UNP P29274
A	325	HIS	-	expression tag	UNP P29274
A	326	HIS	-	expression tag	UNP P29274
A	327	HIS	-	expression tag	UNP P29274
A	328	HIS	-	expression tag	UNP P29274

- Molecule 2 is SODIUM ION (CCD ID: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Na	0	0
			1	1		

- Molecule 3 is THEOPHYLLINE (CCD ID: TEP) (formula: C<sub>7</sub>H<sub>8</sub>N<sub>4</sub>O<sub>2</sub>) (labeled as "Ligand of Interest" by depositor).



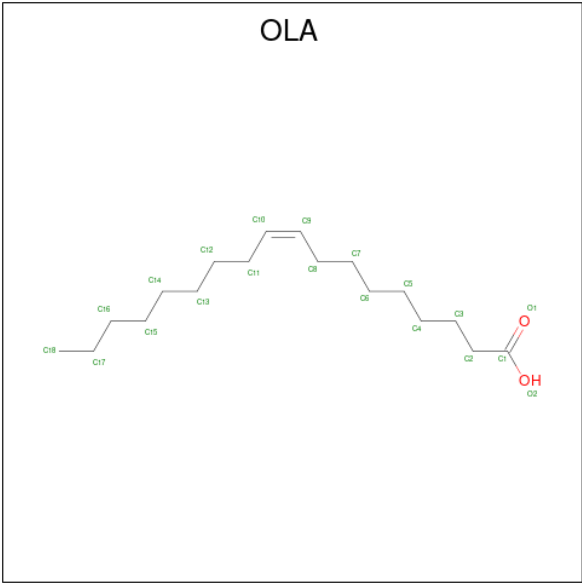
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			13	7	4	2		

- Molecule 4 is CHOLESTEROL (CCD ID: CLR) (formula: C<sub>27</sub>H<sub>46</sub>O).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			28	27	1		
4	A	1	Total	C	O	0	0
			28	27	1		
4	A	1	Total	C	O	0	0
			28	27	1		

- Molecule 5 is OLEIC ACID (CCD ID: OLA) (formula:  $C_{18}H_{34}O_2$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			20	18	2		

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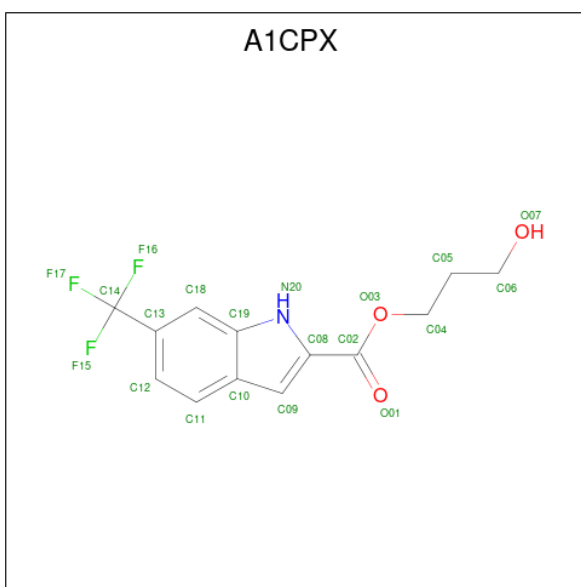
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 10 8 2	0	0
5	A	1	Total C O 9 7 2	0	0
5	A	1	Total C O 18 16 2	0	0
5	A	1	Total C O 20 18 2	0	0
5	A	1	Total C O 20 18 2	0	0
5	A	1	Total C O 12 10 2	0	0
5	A	1	Total C O 8 6 2	0	0
5	A	1	Total C O 15 13 2	0	0
5	A	1	Total C O 11 9 2	0	0
5	A	1	Total C O 19 17 2	0	0
5	A	1	Total C O 20 18 2	0	0
5	A	1	Total C 12 12	0	0
5	A	1	Total C 10 10	0	0
5	A	1	Total C 11 11	0	0
5	A	1	Total C 12 12	0	0
5	A	1	Total C 9 9	0	0
5	A	1	Total C 8 8	0	0
5	A	1	Total C O 15 13 2	0	0
5	A	1	Total C 15 15	0	0
5	A	1	Total C O 13 11 2	0	0
5	A	1	Total C O 20 18 2	0	0

- Molecule 6 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (CCD ID: OLC) (formula:  $C_{21}H_{40}O_4$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			16	12	4		
6	A	1	Total	C	O	0	0
			17	13	4		
6	A	1	Total	C	O	0	0
			23	19	4		
6	A	1	Total	C	O	0	0
			23	19	4		
6	A	1	Total	C	O	0	0
			23	19	4		
6	A	1	Total	C	O	0	0
			18	14	4		
6	A	1	Total	C	O	0	0
			16	12	4		

- Molecule 7 is 3-hydroxypropyl 6-(trifluoromethyl)-1H-indole-2-carboxylate (CCD ID: A1CPX) (formula:  $C_{13}H_{12}F_3NO_3$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	A	1	Total	C	F	N	O	0	0
			20	13	3	1	3		

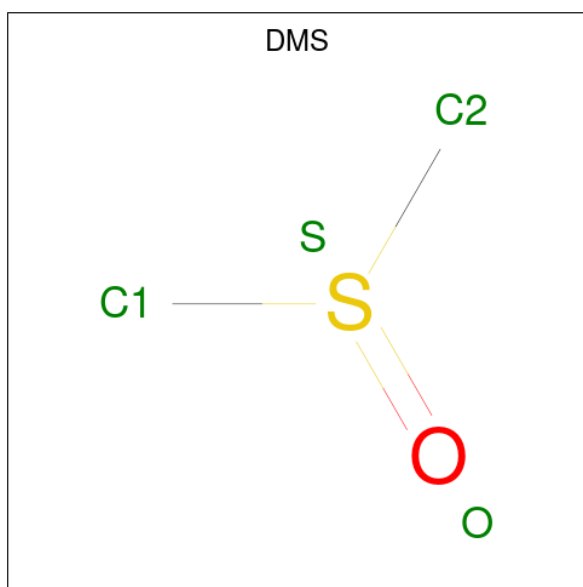
- Molecule 8 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula:  $C_4H_{10}O_3$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			7	4	3		

- Molecule 9 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula:  $C_2H_6OS$ ).





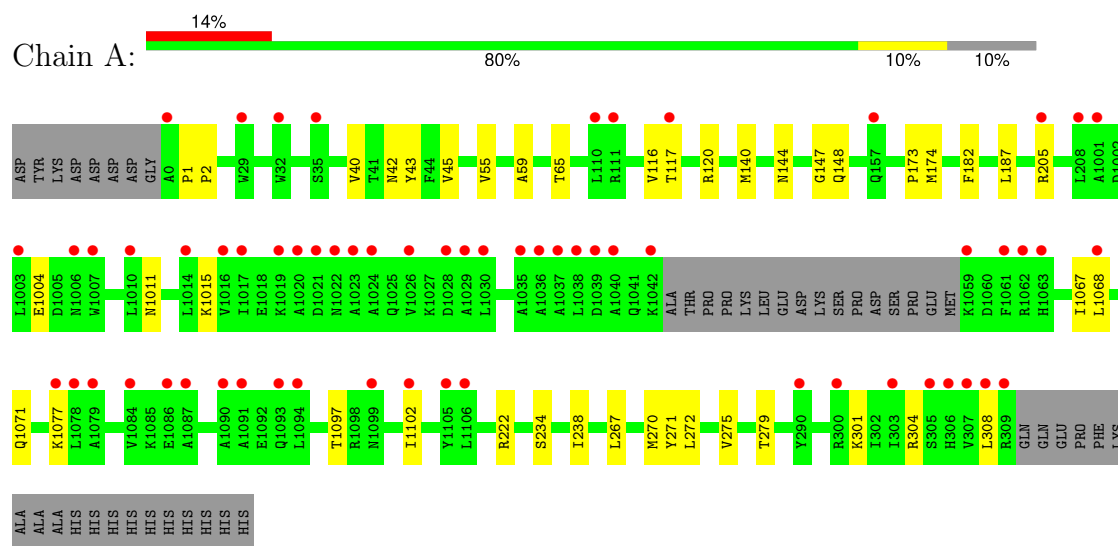
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	A	1	Total	C	O	S	0	1
			4	2	1	1		

- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	65	Total	O	0	0
			65	65		



- Molecule 1: Adenosine receptor A2a/Soluble cytochrome b562/Adenosine receptor A2a chimera



## 4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	39.30Å 178.93Å 139.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	89.47 – 1.89 89.47 – 1.89	Depositor EDS
% Data completeness (in resolution range)	76.2 (89.47-1.89) 76.5 (89.47-1.89)	Depositor EDS
$R_{merge}$	0.11	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.01 (at 1.88Å)	Xtriage
Refinement program	PHENIX (1.20_4459: ???)	Depositor
R, $R_{free}$	0.196 , 0.233 0.196 , 0.233	Depositor DCC
$R_{free}$ test set	1541 reflections (3.83%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.0	Xtriage
Anisotropy	0.061	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 46.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	3763	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.66% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: OLC, NA, TEP, A1CPX, CLR, OLA, DMS, PEG

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	0/3191	0.32	0/4338

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	3126	0	3206	34	0
2	A	1	0	0	0	0
3	A	13	0	8	0	0
4	A	84	0	138	0	0
5	A	307	0	459	12	0
6	A	136	0	186	6	0
7	A	20	0	0	0	0
8	A	7	0	10	0	0
9	A	4	0	6	0	0
10	A	65	0	0	1	0
All	All	3763	0	4013	44	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.

All (44) 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:205[B]:ARG:HH11	1:A:205[B]:ARG:HG3	1.32	0.94
1:A:117:THR:HG22	1:A:120:ARG:HD2	1.71	0.71
1:A:205[B]:ARG:HG3	1:A:205[B]:ARG:NH1	2.08	0.62
1:A:205[B]:ARG:HH11	1:A:205[B]:ARG:CG	2.11	0.62
5:A:2425:OLA:H62	6:A:2433:OLC:H3A	1.79	0.62
1:A:267:LEU:HD11	1:A:271:TYR:CZ	2.37	0.60
1:A:1068:LEU:HD22	1:A:1102:ILE:HG13	1.86	0.57
1:A:140[A]:MET:HE2	5:A:2425:OLA:H52	1.87	0.57
6:A:2429:OLC:H4A	6:A:2430:OLC:H2	1.87	0.55
5:A:2410:OLA:H183	5:A:2421:OLA:H121	1.88	0.54
1:A:275[B]:VAL:O	1:A:279[B]:THR:HG23	2.08	0.53
1:A:301:LYS:N	1:A:301:LYS:HD3	2.24	0.53
1:A:174[B]:MET:SD	10:A:2521:HOH:O	2.59	0.53
1:A:234:SER:O	1:A:238:ILE:HG12	2.11	0.50
5:A:2422:OLA:H51	5:A:2426:OLA:H131	1.93	0.50
1:A:1067:ILE:O	1:A:1071:GLN:HG3	2.11	0.49
5:A:2405:OLA:H161	6:A:2430:OLC:H14A	1.95	0.49
1:A:42:ASN:HA	1:A:45:VAL:HB	1.93	0.49
1:A:205[B]:ARG:NH1	1:A:205[B]:ARG:CG	2.72	0.48
1:A:182:PHE:CE1	1:A:187[B]:LEU:HG	2.49	0.47
1:A:270:MET:HE3	1:A:270:MET:HB3	1.55	0.47
1:A:55:VAL:HA	1:A:59:ALA:HB3	1.97	0.47
1:A:1004:GLU:OE1	1:A:222:ARG:NH1	2.42	0.46
5:A:2415:OLA:H82	5:A:2415:OLA:H111	1.80	0.46
5:A:2409:OLA:H122	5:A:2416:OLA:H151	1.98	0.46
1:A:65:THR:HG21	6:A:2430:OLC:H5	1.98	0.45
1:A:1071:GLN:OE1	1:A:1097:THR:OG1	2.30	0.45
5:A:2416:OLA:H10	5:A:2416:OLA:H132	1.85	0.44
1:A:144:ASN:HA	1:A:173:PRO:HD3	2.00	0.44
1:A:147:GLY:C	1:A:148:GLN:HG2	2.41	0.43
1:A:1011:ASN:O	1:A:1015:LYS:HG3	2.18	0.43
5:A:2408:OLA:H81	5:A:2408:OLA:H111	1.78	0.43
1:A:117:THR:HG23	1:A:120:ARG:H	1.83	0.42
1:A:279[A]:THR:HG23	5:A:2417:OLA:H131	2.00	0.42
1:A:43:TYR:HB3	6:A:2431:OLC:H2	2.00	0.42
1:A:304:ARG:O	1:A:308:LEU:HB2	2.19	0.42
1:A:43:TYR:CE2	6:A:2431:OLC:H24A	2.55	0.41
1:A:40:VAL:HG11	1:A:116:VAL:HG12	2.02	0.41
5:A:2424:OLA:H81	5:A:2424:OLA:H112	1.82	0.41
1:A:1077:LYS:HB3	1:A:1077:LYS:HE3	1.90	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:LEU:HD12	1:A:267:LEU:O	2.20	0.41
1:A:40:VAL:HG11	1:A:116:VAL:CG1	2.50	0.41
1:A:272:LEU:HD21	5:A:2410:OLA:H141	2.03	0.41
1:A:1:PRO:HA	1:A:2:PRO:HD3	1.95	0.40

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	399/433 (92%)	396 (99%)	3 (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	329/354 (93%)	329 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	75	HIS
1	A	1006	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry ⓘ

Of 37 ligands modelled in this entry, 1 is monoatomic - leaving 36 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	$\# Z  > 2$
5	OLA	A	2405	-	19,19,19	0.59	0	19,19,19	1.00	0
5	OLA	A	2423	-	14,14,19	0.68	0	14,14,19	0.98	0
7	A1CPX	A	2434	-	21,21,21	1.63	3 (14%)	30,30,30	1.94	5 (16%)
5	OLA	A	2409	-	19,19,19	0.61	0	19,19,19	0.92	0
5	OLA	A	2417	-	11,11,19	0.36	0	10,10,19	0.68	0
5	OLA	A	2419	-	10,10,19	0.44	0	9,9,19	0.79	0
5	OLA	A	2421	-	8,8,19	0.30	0	7,7,19	0.81	0
4	CLR	A	2403	-	31,31,31	0.30	0	48,48,48	0.46	0
6	OLC	A	2433	-	15,15,24	1.27	1 (6%)	16,16,25	0.93	1 (6%)
5	OLA	A	2406	-	9,9,19	0.76	0	9,9,19	1.20	0
5	OLA	A	2416	-	19,19,19	0.60	0	19,19,19	0.89	0
5	OLA	A	2411	-	11,11,19	0.75	0	11,11,19	1.12	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
8	PEG	A	2435	-	6,6,6	0.49	0	5,5,5	0.28	0
5	OLA	A	2426	-	19,19,19	0.58	0	19,19,19	1.02	0
5	OLA	A	2412	-	7,7,19	0.85	0	7,7,19	1.24	1 (14%)
6	OLC	A	2432	-	17,17,24	1.25	1 (5%)	18,18,25	1.29	2 (11%)
6	OLC	A	2430	-	22,22,24	1.08	1 (4%)	23,23,25	1.32	3 (13%)
5	OLA	A	2410	-	19,19,19	0.57	0	19,19,19	1.04	1 (5%)
4	CLR	A	2402	-	31,31,31	0.30	0	48,48,48	0.47	0
5	OLA	A	2415	-	18,18,19	0.60	0	18,18,19	1.08	1 (5%)
5	OLA	A	2424	-	14,14,19	0.36	0	13,13,19	0.67	0
6	OLC	A	2427	-	15,15,24	1.27	1 (6%)	16,16,25	1.23	1 (6%)
5	OLA	A	2407	-	8,8,19	0.80	0	8,8,19	1.30	0
5	OLA	A	2413	-	14,14,19	0.64	0	14,14,19	1.11	0
5	OLA	A	2425	-	12,12,19	0.73	0	12,12,19	1.12	0
5	OLA	A	2420	-	11,11,19	0.38	0	10,10,19	0.80	0
4	CLR	A	2404	-	31,31,31	0.29	0	48,48,48	0.39	0
6	OLC	A	2431	-	22,22,24	1.12	1 (4%)	23,23,25	1.30	2 (8%)
5	OLA	A	2418	-	9,9,19	0.41	0	8,8,19	0.69	0
6	OLC	A	2429	-	22,22,24	1.06	1 (4%)	23,23,25	1.30	3 (13%)
5	OLA	A	2414	-	10,10,19	0.71	0	10,10,19	1.25	0
6	OLC	A	2428	-	16,16,24	1.31	1 (6%)	17,17,25	1.30	2 (11%)
3	TEP	A	2401	-	14,14,14	1.79	4 (28%)	21,21,21	3.27	11 (52%)
5	OLA	A	2408	-	17,17,19	0.59	0	17,17,19	1.06	0
5	OLA	A	2422	-	7,7,19	0.30	0	6,6,19	0.83	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
5	OLA	A	2405	-	-	10/17/17/17	-
5	OLA	A	2423	-	-	5/12/12/17	-
7	A1CPX	A	2434	-	-	0/15/15/15	0/2/2/2
5	OLA	A	2409	-	-	10/17/17/17	-
5	OLA	A	2417	-	-	4/9/9/17	-
5	OLA	A	2419	-	-	6/8/8/17	-
5	OLA	A	2421	-	-	3/6/6/17	-
4	CLR	A	2403	-	-	0/10/68/68	0/4/4/4
6	OLC	A	2433	-	-	4/15/15/24	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	OLA	A	2406	-	-	4/7/7/17	-
5	OLA	A	2416	-	-	7/17/17/17	-
5	OLA	A	2411	-	-	3/9/9/17	-
8	PEG	A	2435	-	-	2/4/4/4	-
5	OLA	A	2426	-	-	10/17/17/17	-
5	OLA	A	2412	-	-	3/5/5/17	-
6	OLC	A	2432	-	-	7/17/17/24	-
6	OLC	A	2430	-	-	12/22/22/24	-
5	OLA	A	2410	-	-	8/17/17/17	-
4	CLR	A	2402	-	-	0/10/68/68	0/4/4/4
5	OLA	A	2415	-	-	4/16/16/17	-
5	OLA	A	2424	-	-	2/12/12/17	-
6	OLC	A	2427	-	-	6/15/15/24	-
5	OLA	A	2407	-	-	4/6/6/17	-
5	OLA	A	2413	-	-	4/12/12/17	-
5	OLA	A	2425	-	-	6/10/10/17	-
5	OLA	A	2420	-	-	4/9/9/17	-
4	CLR	A	2404	-	-	1/10/68/68	0/4/4/4
6	OLC	A	2431	-	-	12/22/22/24	-
5	OLA	A	2418	-	-	1/7/7/17	-
6	OLC	A	2429	-	-	12/22/22/24	-
5	OLA	A	2414	-	-	3/8/8/17	-
6	OLC	A	2428	-	-	8/16/16/24	-
3	TEP	A	2401	-	-	-	0/2/2/2
5	OLA	A	2408	-	-	7/15/15/17	-
5	OLA	A	2422	-	-	0/5/5/17	-

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	2434	A1CPX	C08-N20	-4.08	1.32	1.38
7	A	2434	A1CPX	O03-C02	3.57	1.40	1.33
3	A	2401	TEP	C6-N1	-3.52	1.33	1.40
6	A	2427	OLC	O20-C1	3.05	1.42	1.33
6	A	2433	OLC	O20-C1	3.03	1.42	1.33
6	A	2428	OLC	O20-C1	2.99	1.42	1.33
6	A	2431	OLC	O20-C1	2.95	1.42	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	2432	OLC	O20-C1	2.88	1.41	1.33
6	A	2430	OLC	O20-C1	2.84	1.41	1.33
6	A	2429	OLC	O20-C1	2.79	1.41	1.33
7	A	2434	A1CPX	C10-C19	-2.25	1.37	1.41
3	A	2401	TEP	C5-N7	-2.22	1.34	1.38
3	A	2401	TEP	C5-C4	-2.21	1.34	1.37
3	A	2401	TEP	C5-C6	-2.08	1.38	1.42

All (33) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	2434	A1CPX	O03-C02-C08	8.22	121.17	111.97
3	A	2401	TEP	C4-C5-N7	6.42	108.45	105.29
3	A	2401	TEP	C8-N9-C4	5.81	107.34	103.45
3	A	2401	TEP	C5-C6-N1	5.79	119.84	112.07
3	A	2401	TEP	C5-C4-N9	-4.80	108.26	111.63
3	A	2401	TEP	C6-C5-C4	-3.93	120.03	123.13
3	A	2401	TEP	N7-C8-N9	-3.64	108.76	112.98
3	A	2401	TEP	N3-C4-N9	3.57	131.89	126.27
3	A	2401	TEP	C6-N1-C2	-3.55	119.89	125.66
6	A	2429	OLC	C8-C9-C10	3.49	151.01	124.83
6	A	2431	OLC	C8-C9-C10	3.48	150.90	124.83
6	A	2428	OLC	C8-C9-C10	3.47	151.94	126.65
3	A	2401	TEP	O6-C6-C5	-3.45	119.59	126.64
3	A	2401	TEP	C5-C4-N3	-3.33	119.85	121.73
6	A	2430	OLC	C8-C9-C10	3.29	149.51	124.83
6	A	2427	OLC	O20-C1-C2	3.16	121.47	111.83
6	A	2432	OLC	C8-C9-C10	2.87	152.19	130.48
6	A	2431	OLC	O20-C1-C2	2.75	120.23	111.83
6	A	2432	OLC	O20-C1-C2	2.66	119.96	111.83
6	A	2430	OLC	O20-C1-C2	2.48	119.41	111.83
6	A	2428	OLC	O20-C1-C2	2.48	119.40	111.83
6	A	2433	OLC	O20-C1-C2	2.47	119.36	111.83
3	A	2401	TEP	N3-C2-N1	2.45	120.20	117.14
7	A	2434	A1CPX	C18-C19-C10	-2.31	120.30	122.59
7	A	2434	A1CPX	C04-O03-C02	-2.20	112.08	116.55
5	A	2410	OLA	C3-C2-C1	-2.19	108.80	114.51
5	A	2415	OLA	C3-C2-C1	-2.18	108.83	114.51
6	A	2430	OLC	C11-C10-C9	-2.12	108.94	124.83
7	A	2434	A1CPX	F17-C14-C13	-2.08	108.44	112.90
7	A	2434	A1CPX	O03-C02-O01	-2.05	119.68	123.37
6	A	2429	OLC	C11-C10-C9	-2.03	109.59	124.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	2429	OLC	C4-C3-C2	-2.02	105.69	113.13
5	A	2412	OLA	O2-C1-C2	2.01	120.36	114.00

There are no chirality outliers.

All (172) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	2413	OLA	C10-C11-C12-C13
5	A	2417	OLA	C6-C7-C8-C9
6	A	2429	OLC	C21-C22-C24-O25
6	A	2430	OLC	C21-C22-C24-O25
6	A	2430	OLC	O20-C21-C22-O23
6	A	2431	OLC	C10-C11-C12-C13
6	A	2431	OLC	O20-C21-C22-C24
6	A	2432	OLC	O19-C1-O20-C21
6	A	2432	OLC	C2-C1-O20-C21
6	A	2431	OLC	C2-C1-O20-C21
6	A	2431	OLC	O19-C1-O20-C21
5	A	2410	OLA	C6-C7-C8-C9
6	A	2430	OLC	O20-C21-C22-C24
6	A	2431	OLC	O20-C21-C22-O23
5	A	2425	OLA	C1-C2-C3-C4
5	A	2423	OLA	C1-C2-C3-C4
6	A	2433	OLC	C1-C2-C3-C4
6	A	2427	OLC	O20-C21-C22-O23
6	A	2428	OLC	O20-C21-C22-O23
6	A	2432	OLC	O20-C21-C22-O23
4	A	2404	CLR	C20-C22-C23-C24
6	A	2427	OLC	O20-C21-C22-C24
5	A	2415	OLA	C5-C6-C7-C8
5	A	2417	OLA	C12-C13-C14-C15
5	A	2426	OLA	C11-C12-C13-C14
5	A	2405	OLA	C14-C15-C16-C17
6	A	2429	OLC	O23-C22-C24-O25
6	A	2430	OLC	O23-C22-C24-O25
6	A	2428	OLC	C1-C2-C3-C4
6	A	2431	OLC	C1-C2-C3-C4
6	A	2430	OLC	C1-C2-C3-C4
5	A	2419	OLA	C11-C12-C13-C14
5	A	2409	OLA	C12-C13-C14-C15
8	A	2435	PEG	O1-C1-C2-O2
5	A	2423	OLA	C5-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
6	A	2432	OLC	C3-C4-C5-C6
5	A	2414	OLA	C1-C2-C3-C4
5	A	2405	OLA	C6-C7-C8-C9
5	A	2414	OLA	C4-C5-C6-C7
5	A	2405	OLA	C2-C3-C4-C5
6	A	2432	OLC	C4-C5-C6-C7
6	A	2429	OLC	C3-C4-C5-C6
6	A	2431	OLC	C4-C5-C6-C7
5	A	2423	OLA	C3-C4-C5-C6
6	A	2432	OLC	C5-C6-C7-C8
5	A	2418	OLA	C9-C10-C11-C12
6	A	2433	OLC	C2-C3-C4-C5
5	A	2410	OLA	C15-C16-C17-C18
5	A	2426	OLA	C15-C16-C17-C18
5	A	2416	OLA	C5-C6-C7-C8
6	A	2429	OLC	C10-C11-C12-C13
6	A	2431	OLC	C6-C7-C8-C9
6	A	2432	OLC	C6-C7-C8-C9
5	A	2411	OLA	C4-C5-C6-C7
6	A	2428	OLC	C3-C4-C5-C6
5	A	2410	OLA	C1-C2-C3-C4
5	A	2426	OLA	C4-C5-C6-C7
5	A	2424	OLA	C6-C7-C8-C9
5	A	2405	OLA	C12-C13-C14-C15
5	A	2407	OLA	C3-C4-C5-C6
6	A	2430	OLC	C12-C13-C14-C15
6	A	2428	OLC	C2-C1-O20-C21
5	A	2426	OLA	C6-C7-C8-C9
6	A	2430	OLC	C10-C11-C12-C13
5	A	2410	OLA	C5-C6-C7-C8
5	A	2416	OLA	C3-C4-C5-C6
6	A	2431	OLC	C5-C6-C7-C8
5	A	2419	OLA	C10-C11-C12-C13
5	A	2408	OLA	C4-C5-C6-C7
6	A	2429	OLC	C4-C5-C6-C7
5	A	2405	OLA	C10-C11-C12-C13
5	A	2408	OLA	C6-C7-C8-C9
6	A	2428	OLC	O19-C1-O20-C21
6	A	2427	OLC	C6-C7-C8-C9
5	A	2421	OLA	C12-C13-C14-C15
6	A	2429	OLC	C12-C13-C14-C15
6	A	2433	OLC	C6-C7-C8-C9

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Mol	Chain	Res	Type	Atoms
6	A	2428	OLC	C4-C5-C6-C7
5	A	2405	OLA	C15-C16-C17-C18
5	A	2416	OLA	C11-C12-C13-C14
6	A	2427	OLC	C2-C1-O20-C21
5	A	2409	OLA	C13-C14-C15-C16
5	A	2419	OLA	C12-C13-C14-C15
6	A	2430	OLC	C2-C1-O20-C21
5	A	2426	OLA	C12-C13-C14-C15
5	A	2409	OLA	C4-C5-C6-C7
5	A	2405	OLA	C11-C12-C13-C14
5	A	2423	OLA	C2-C3-C4-C5
5	A	2406	OLA	C3-C4-C5-C6
6	A	2429	OLC	C2-C3-C4-C5
5	A	2419	OLA	C15-C16-C17-C18
5	A	2426	OLA	C1-C2-C3-C4
5	A	2409	OLA	C14-C15-C16-C17
6	A	2433	OLC	C4-C5-C6-C7
5	A	2408	OLA	C1-C2-C3-C4
5	A	2420	OLA	C9-C10-C11-C12
8	A	2435	PEG	C4-C3-O2-C2
5	A	2409	OLA	C3-C4-C5-C6
5	A	2425	OLA	C5-C6-C7-C8
5	A	2421	OLA	C15-C16-C17-C18
5	A	2409	OLA	C11-C12-C13-C14
6	A	2427	OLC	O19-C1-O20-C21
6	A	2430	OLC	O19-C1-O20-C21
5	A	2409	OLA	C10-C11-C12-C13
5	A	2410	OLA	C10-C11-C12-C13
6	A	2428	OLC	C7-C8-C9-C10
5	A	2420	OLA	C4-C5-C6-C7
5	A	2425	OLA	C6-C7-C8-C9
6	A	2428	OLC	C2-C3-C4-C5
5	A	2411	OLA	C3-C4-C5-C6
5	A	2408	OLA	C12-C13-C14-C15
5	A	2426	OLA	C14-C15-C16-C17
5	A	2416	OLA	C6-C7-C8-C9
5	A	2423	OLA	C4-C5-C6-C7
5	A	2411	OLA	C5-C6-C7-C8
5	A	2417	OLA	C11-C12-C13-C14
6	A	2431	OLC	C3-C4-C5-C6
6	A	2429	OLC	C11-C12-C13-C14
6	A	2430	OLC	C4-C5-C6-C7

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Mol	Chain	Res	Type	Atoms
5	A	2419	OLA	C13-C14-C15-C16
5	A	2405	OLA	C5-C6-C7-C8
5	A	2410	OLA	C12-C13-C14-C15
5	A	2425	OLA	C4-C5-C6-C7
5	A	2426	OLA	C2-C3-C4-C5
5	A	2414	OLA	C5-C6-C7-C8
6	A	2429	OLC	C5-C6-C7-C8
5	A	2424	OLA	C12-C13-C14-C15
5	A	2420	OLA	C7-C8-C9-C10
6	A	2429	OLC	C9-C10-C11-C12
5	A	2406	OLA	C5-C6-C7-C8
5	A	2425	OLA	C2-C3-C4-C5
5	A	2408	OLA	C10-C11-C12-C13
6	A	2431	OLC	C2-C3-C4-C5
5	A	2415	OLA	C11-C12-C13-C14
6	A	2430	OLC	C5-C6-C7-C8
6	A	2429	OLC	C6-C7-C8-C9
5	A	2420	OLA	C2-C3-C4-C5
5	A	2412	OLA	C2-C3-C4-C5
6	A	2429	OLC	O20-C21-C22-O23
5	A	2408	OLA	C3-C4-C5-C6
6	A	2427	OLC	C5-C6-C7-C8
5	A	2416	OLA	C4-C5-C6-C7
6	A	2431	OLC	C9-C10-C11-C12
5	A	2405	OLA	C1-C2-C3-C4
5	A	2413	OLA	C7-C8-C9-C10
5	A	2419	OLA	C9-C10-C11-C12
5	A	2409	OLA	C9-C10-C11-C12
5	A	2410	OLA	C7-C8-C9-C10
5	A	2409	OLA	C7-C8-C9-C10
5	A	2425	OLA	C7-C8-C9-C10
5	A	2407	OLA	O2-C1-C2-C3
5	A	2413	OLA	C4-C5-C6-C7
5	A	2416	OLA	C7-C8-C9-C10
5	A	2416	OLA	C13-C14-C15-C16
5	A	2413	OLA	C9-C10-C11-C12
5	A	2421	OLA	C10-C11-C12-C13
5	A	2407	OLA	O1-C1-C2-C3
6	A	2430	OLC	C9-C10-C11-C12
5	A	2417	OLA	C14-C15-C16-C17
5	A	2408	OLA	C11-C12-C13-C14
5	A	2407	OLA	C1-C2-C3-C4

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Mol	Chain	Res	Type	Atoms
5	A	2415	OLA	C7-C8-C9-C10
5	A	2426	OLA	C7-C8-C9-C10
5	A	2409	OLA	O1-C1-C2-C3
5	A	2426	OLA	C3-C4-C5-C6
5	A	2406	OLA	O2-C1-C2-C3
5	A	2412	OLA	O2-C1-C2-C3
5	A	2412	OLA	O1-C1-C2-C3
5	A	2410	OLA	C4-C5-C6-C7
5	A	2405	OLA	O2-C1-C2-C3
5	A	2406	OLA	O1-C1-C2-C3
5	A	2415	OLA	O2-C1-C2-C3

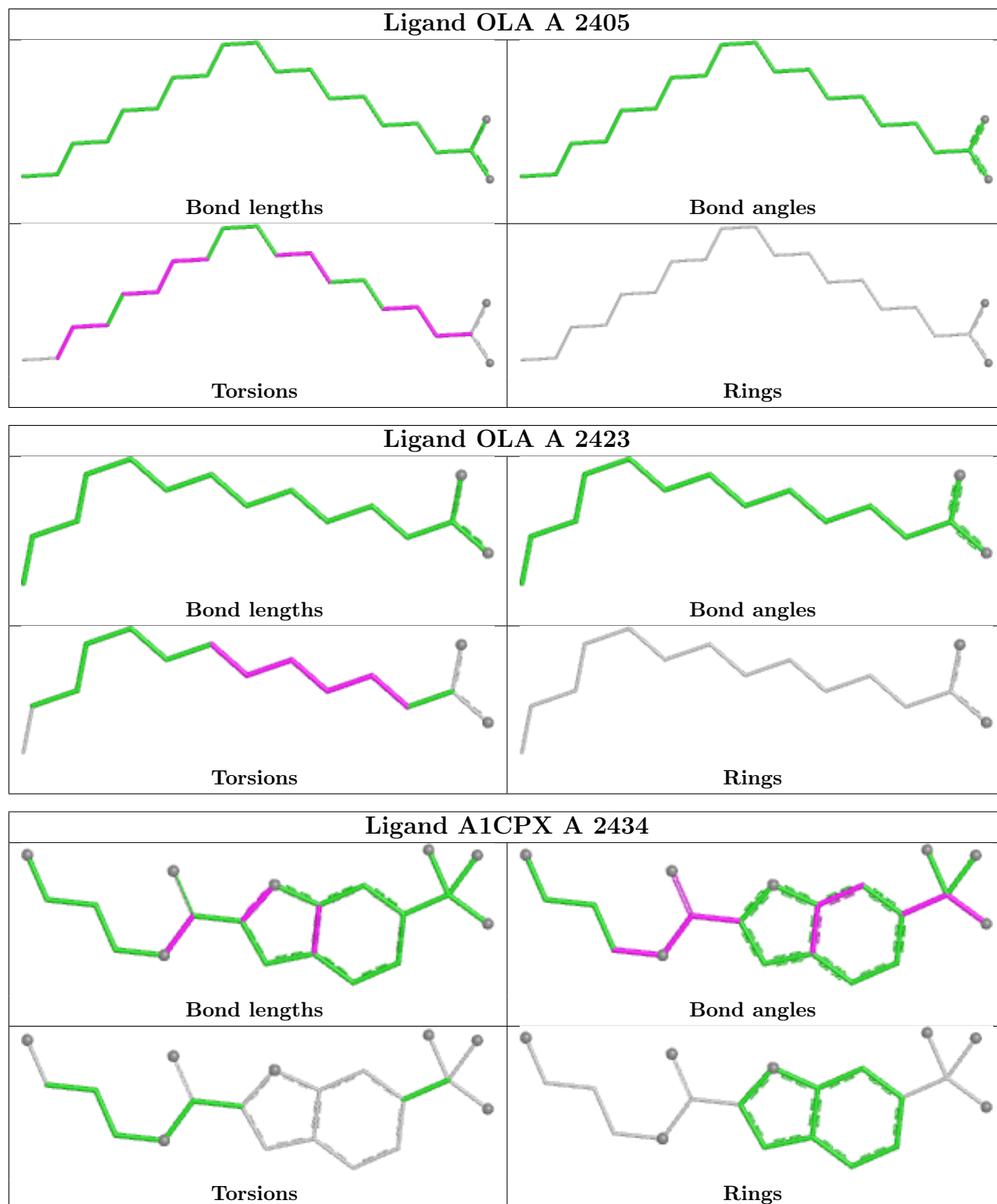
There are no ring outliers.

16 monomers are involved in 16 short contacts:

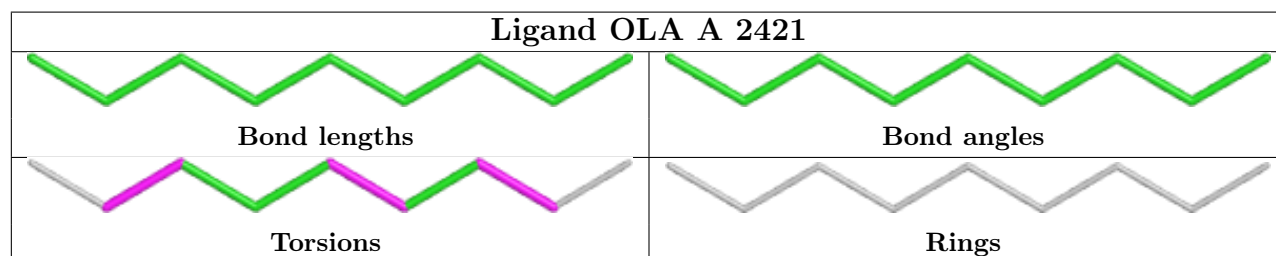
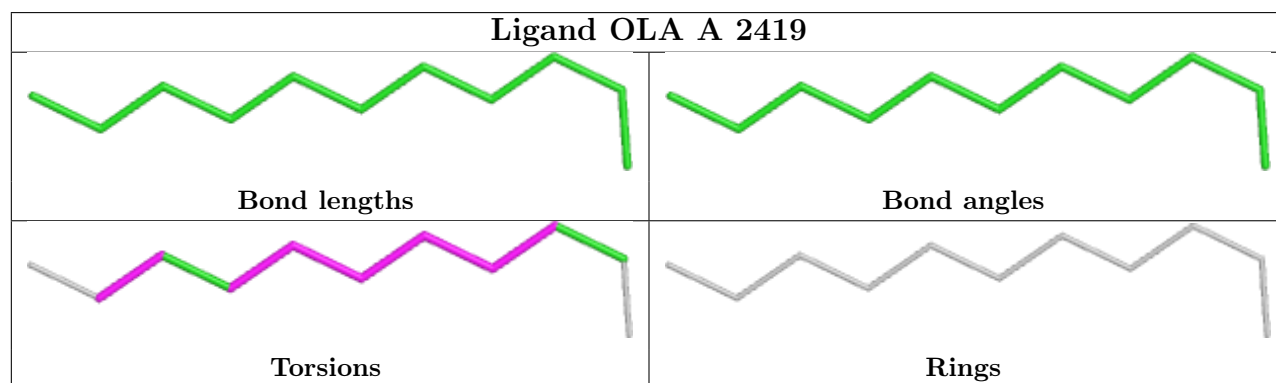
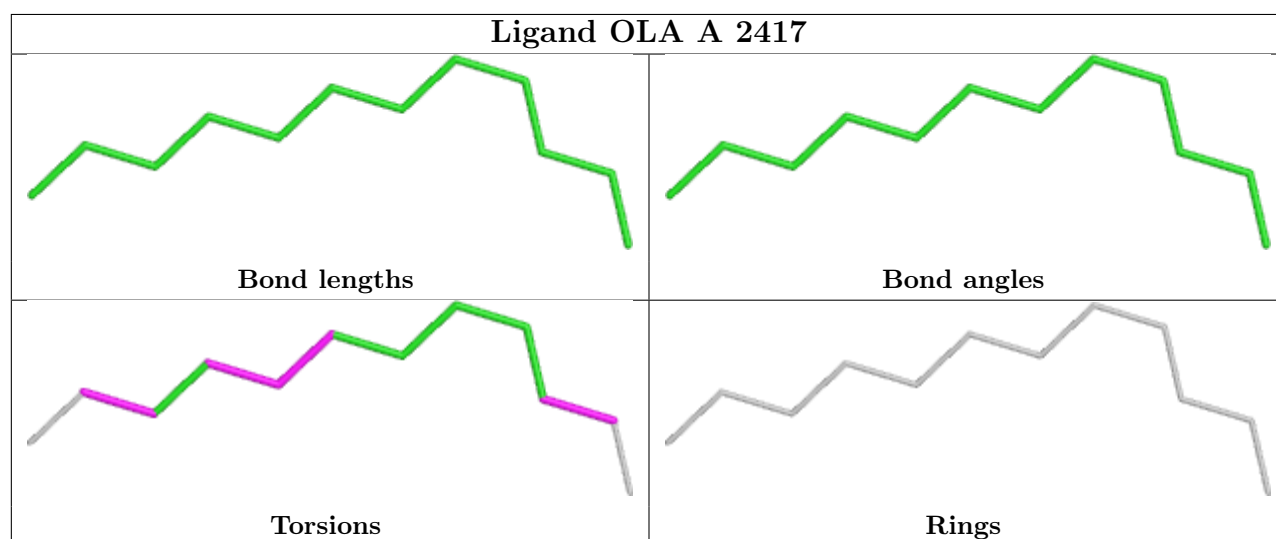
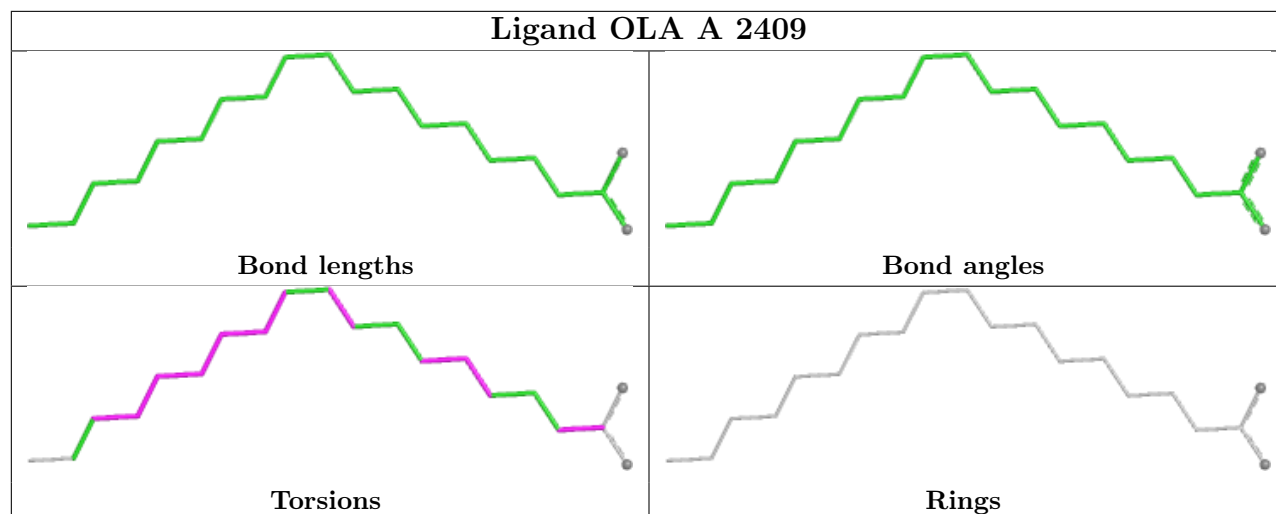
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2405	OLA	1	0
5	A	2409	OLA	1	0
5	A	2417	OLA	1	0
5	A	2421	OLA	1	0
6	A	2433	OLC	1	0
5	A	2416	OLA	2	0
5	A	2426	OLA	1	0
6	A	2430	OLC	3	0
5	A	2410	OLA	2	0
5	A	2415	OLA	1	0
5	A	2424	OLA	1	0
5	A	2425	OLA	2	0
6	A	2431	OLC	2	0
6	A	2429	OLC	1	0
5	A	2408	OLA	1	0
5	A	2422	OLA	1	0

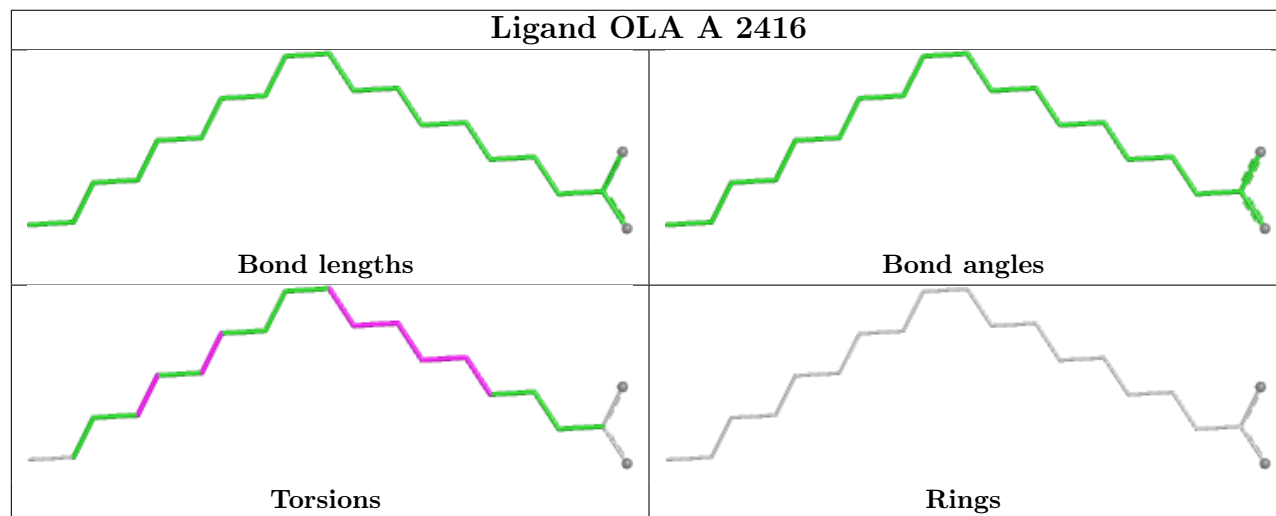
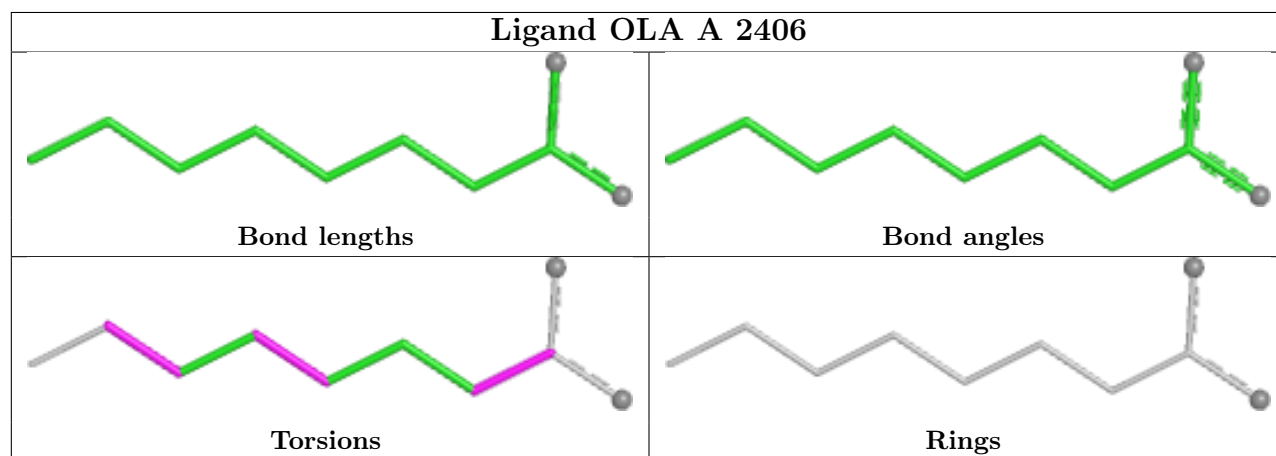
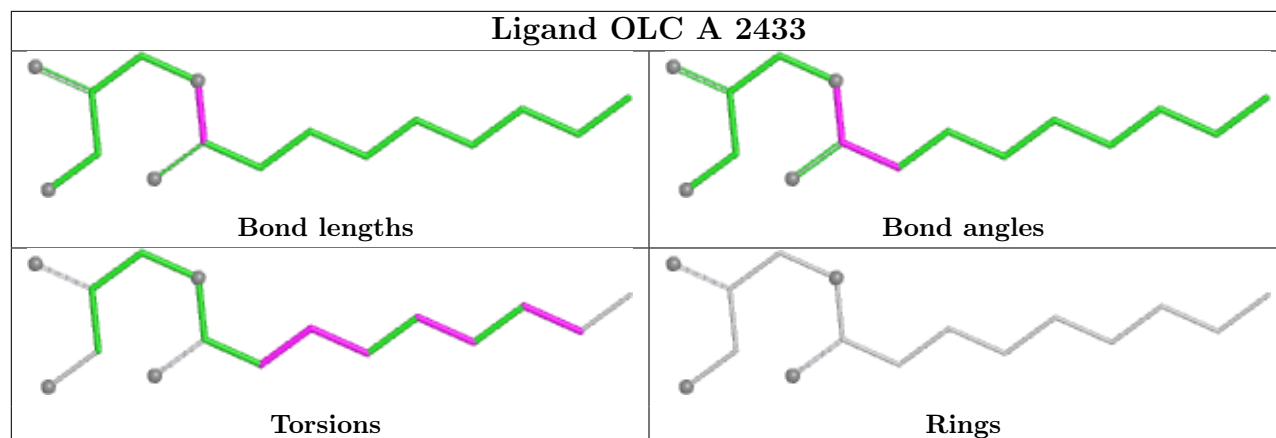
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and

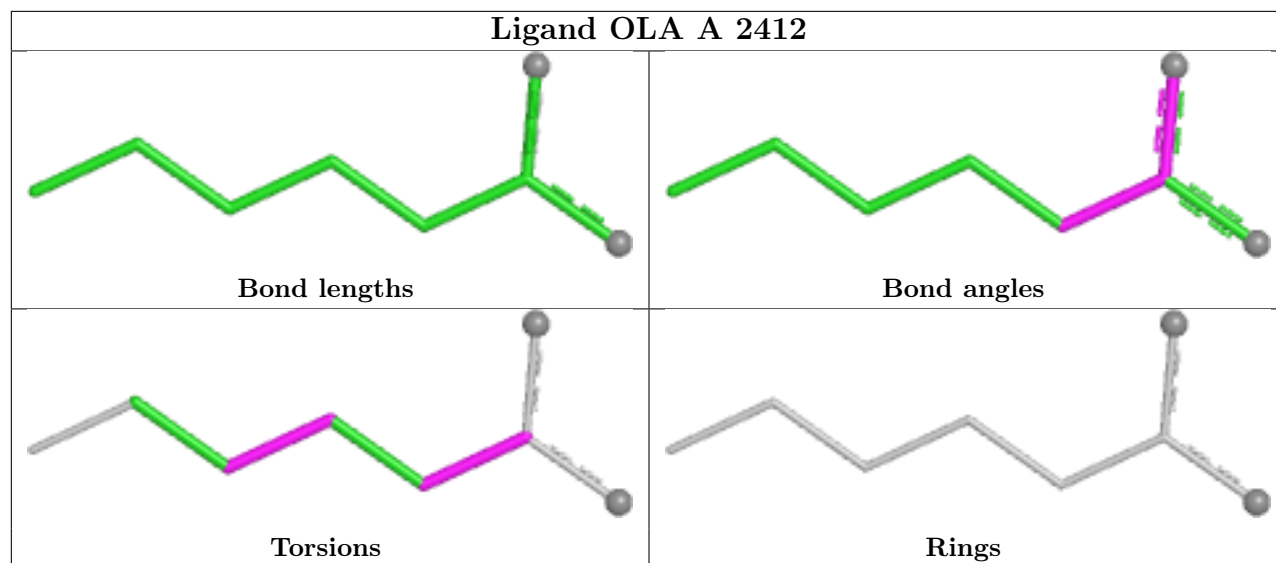
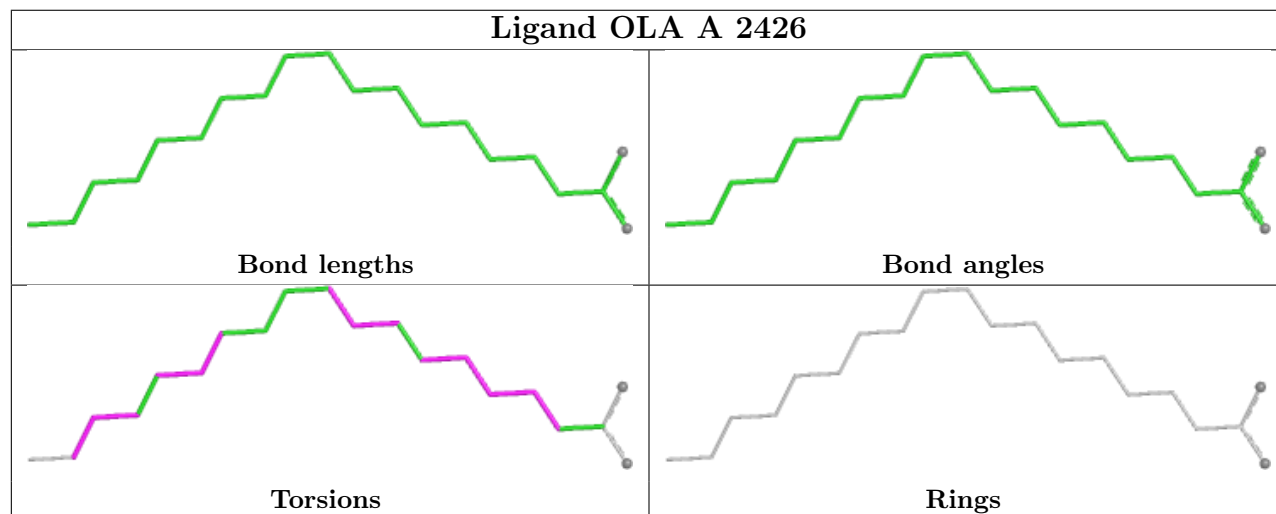
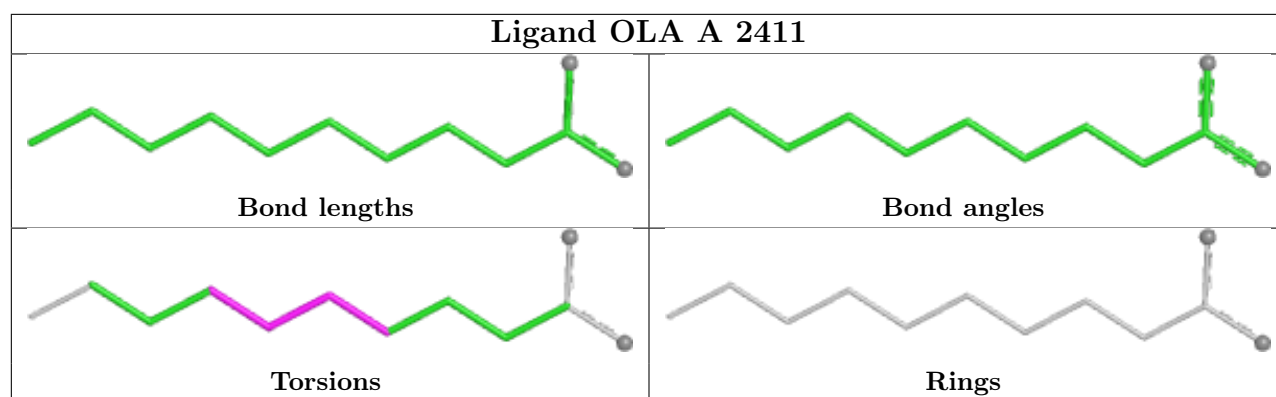
any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

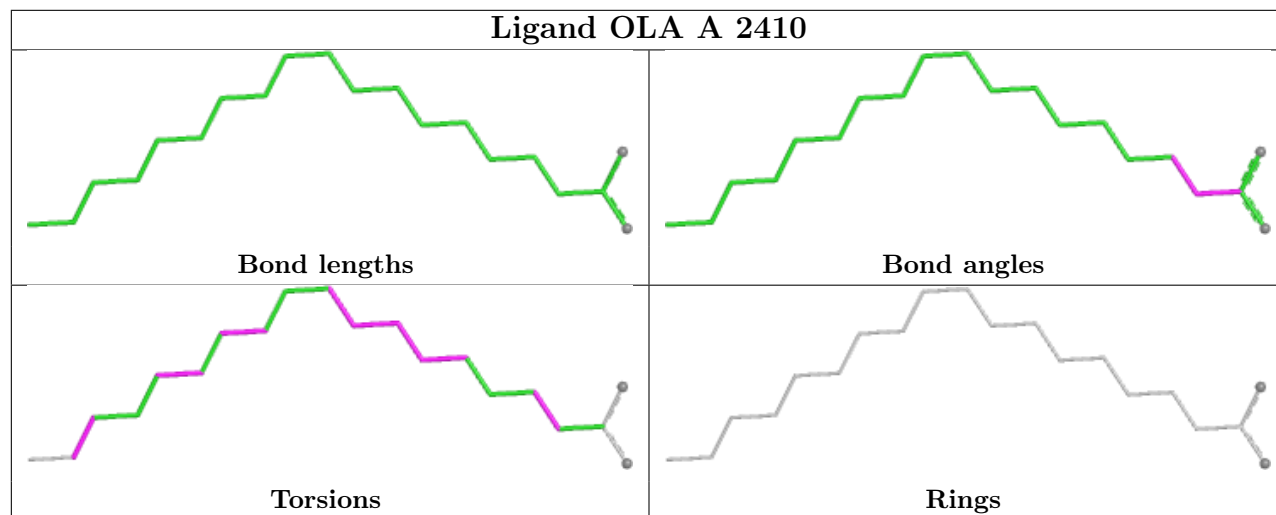
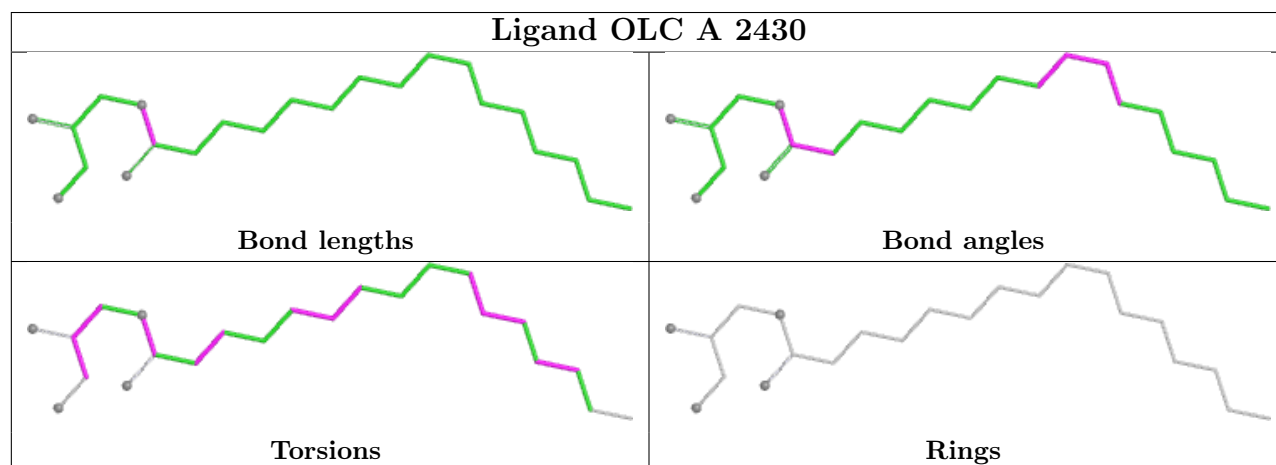
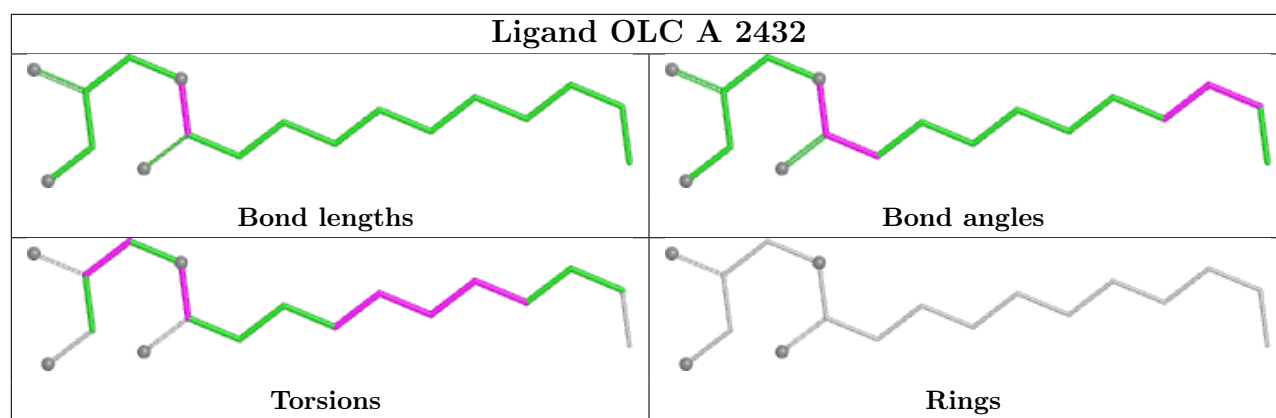


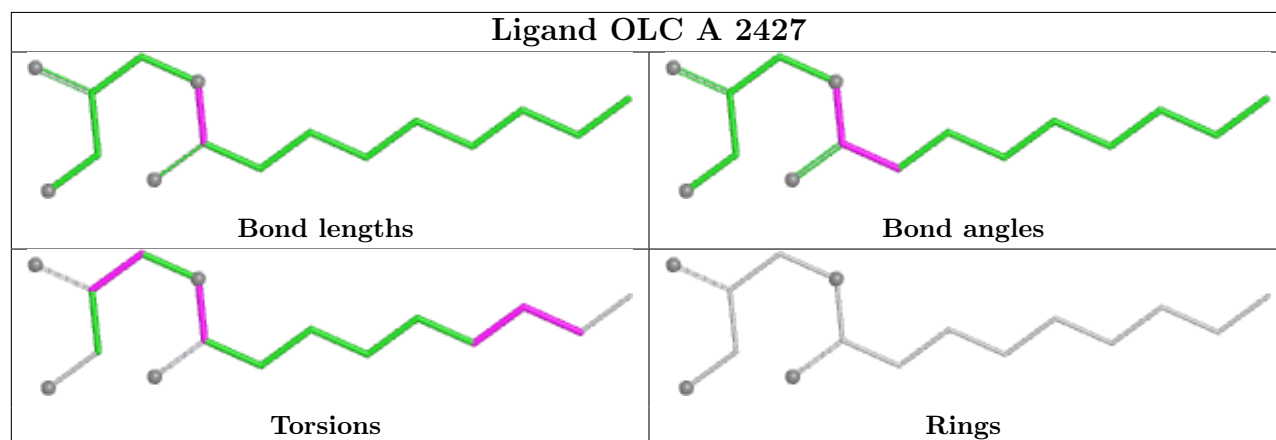
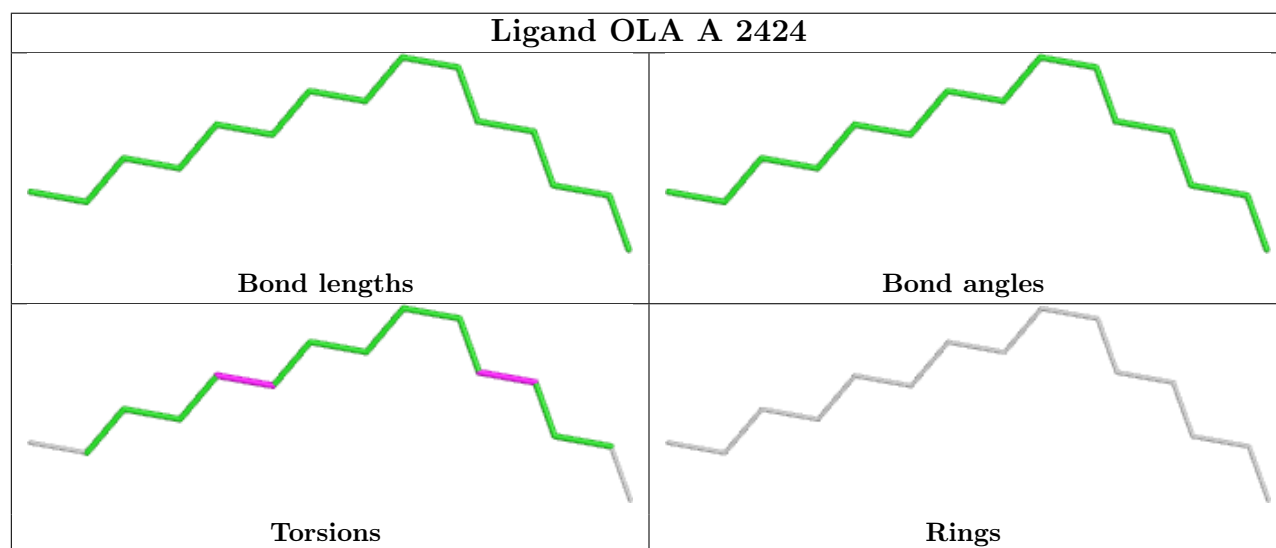
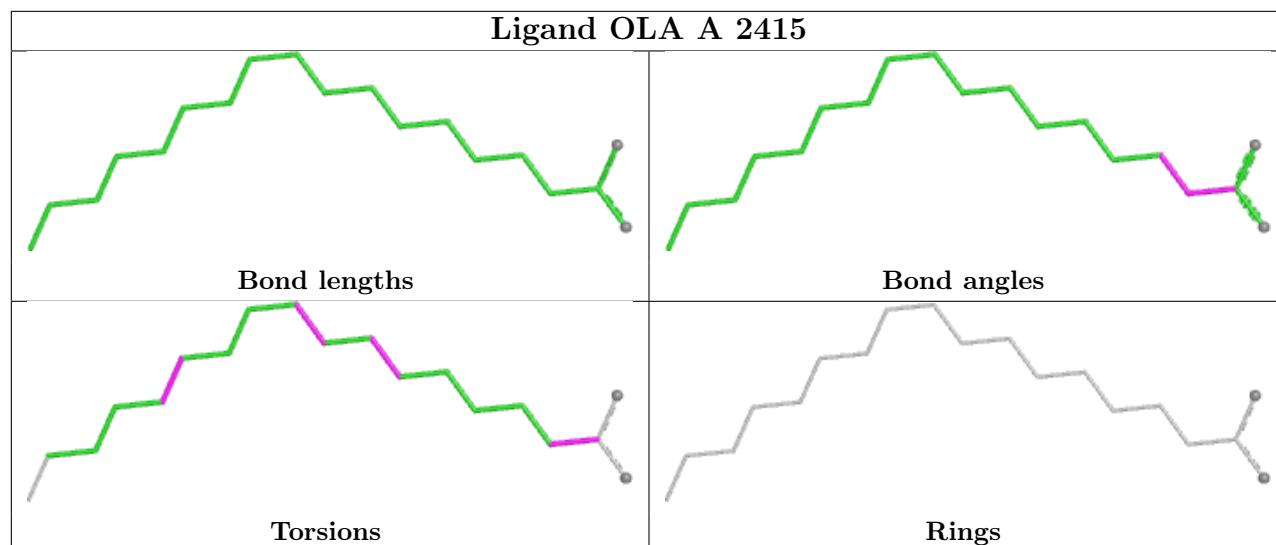


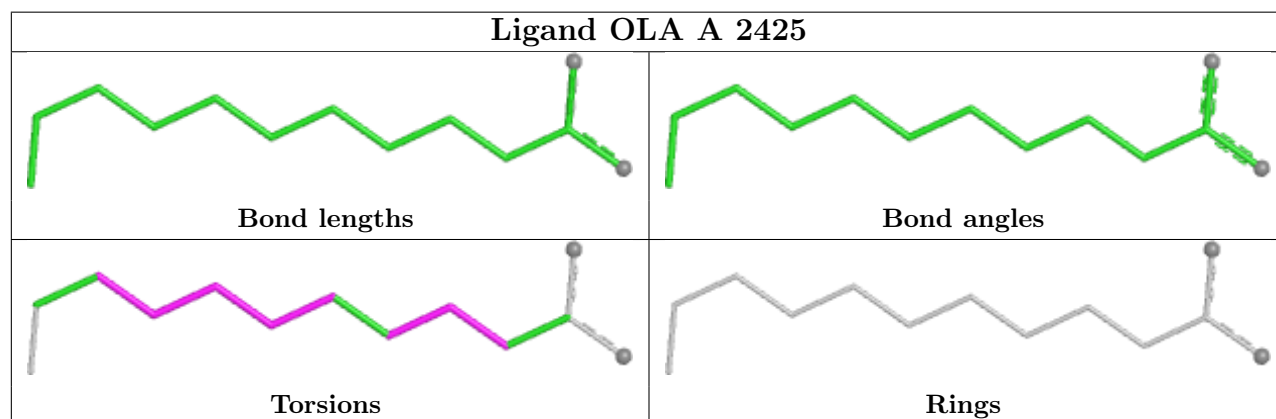
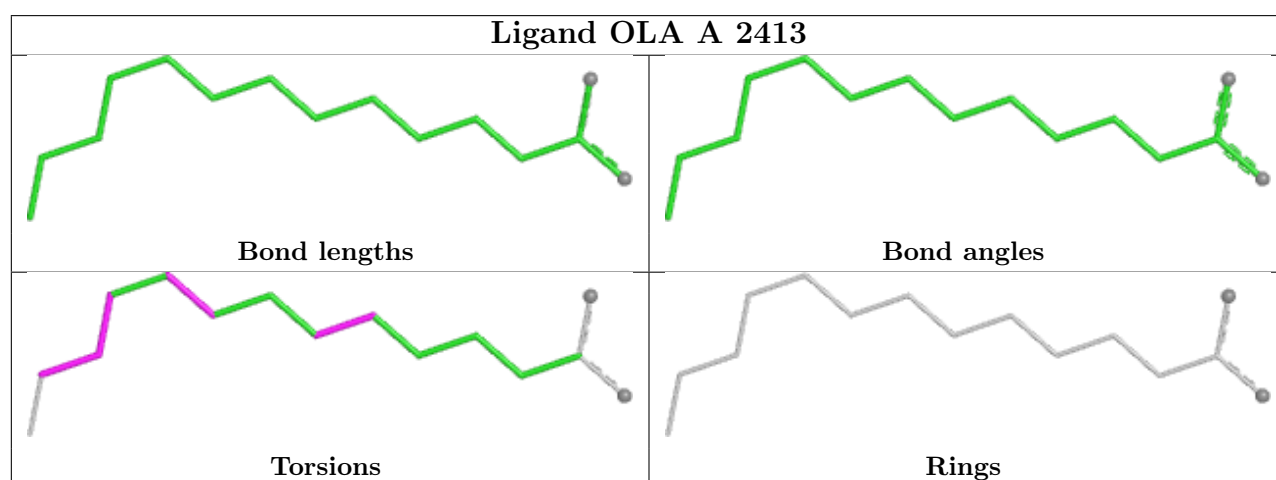
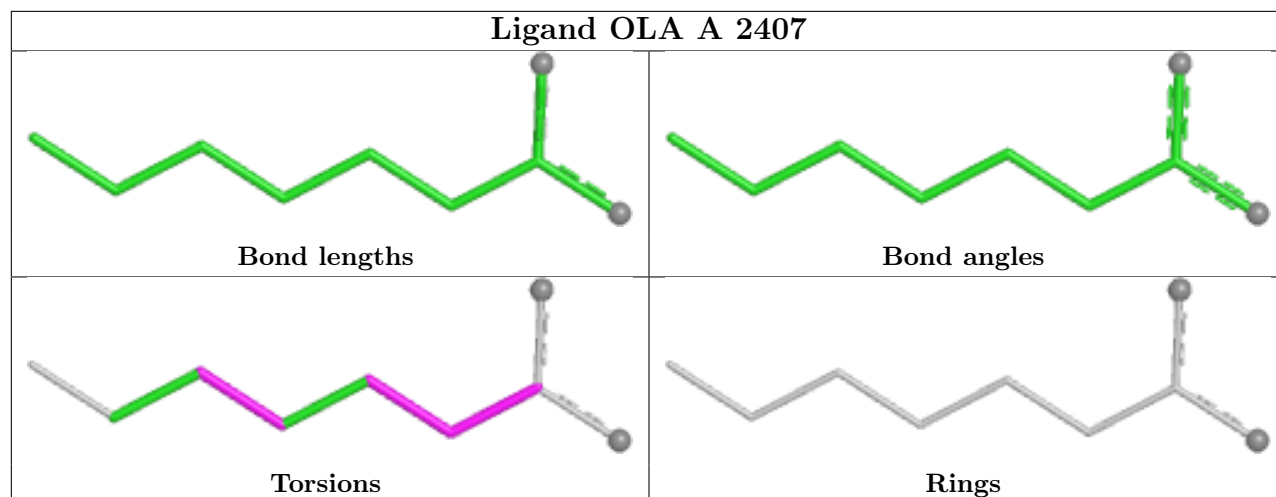


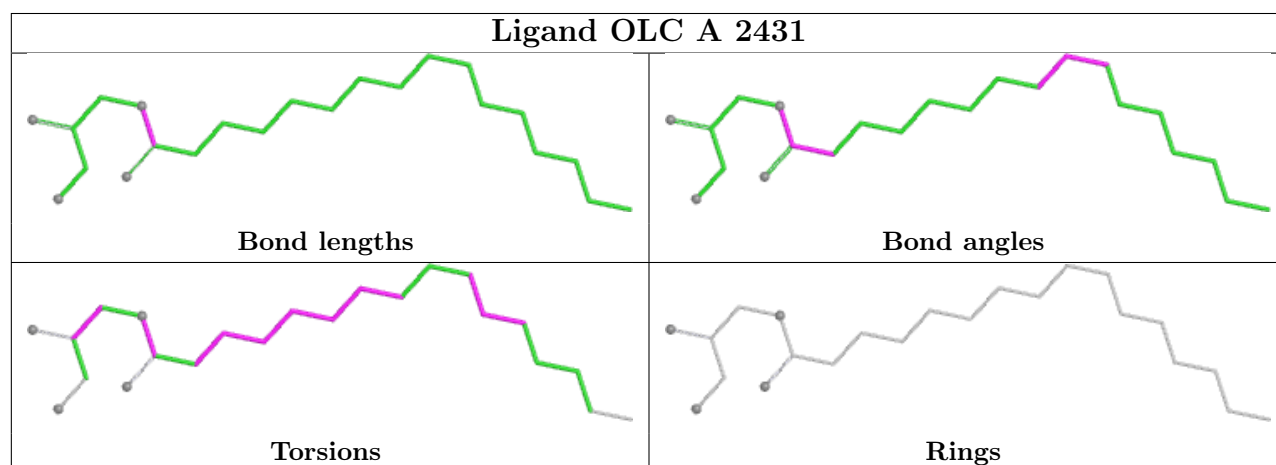
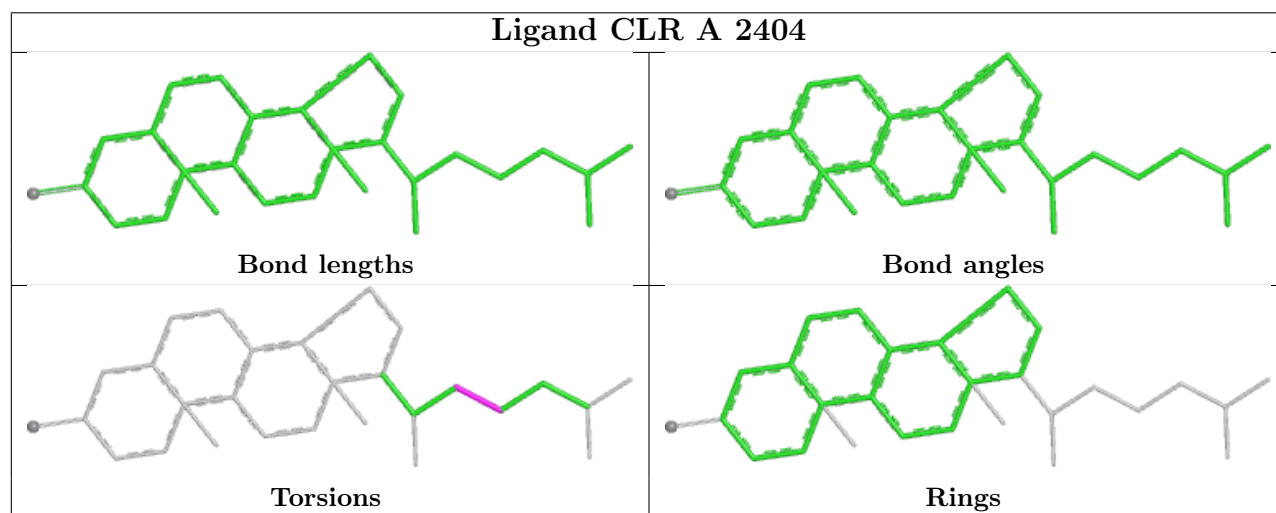
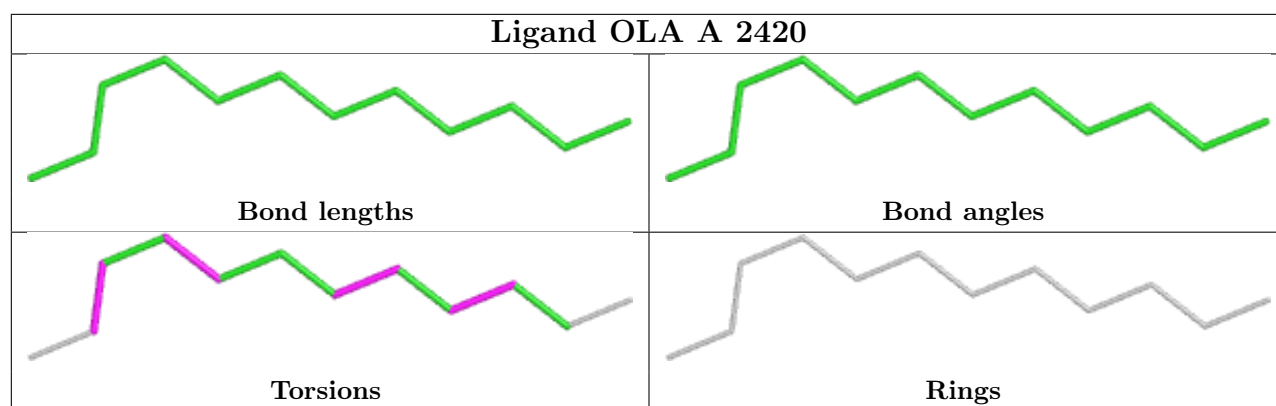


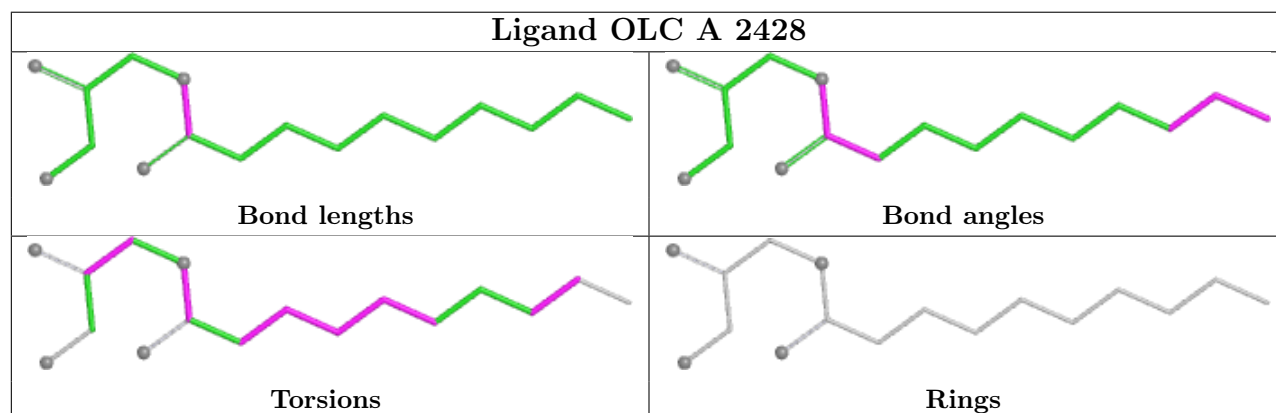
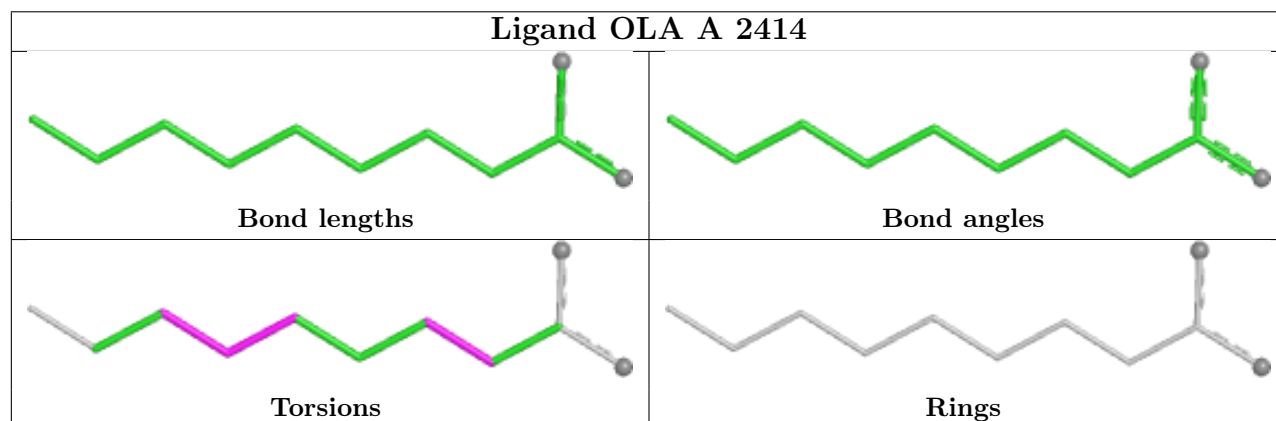
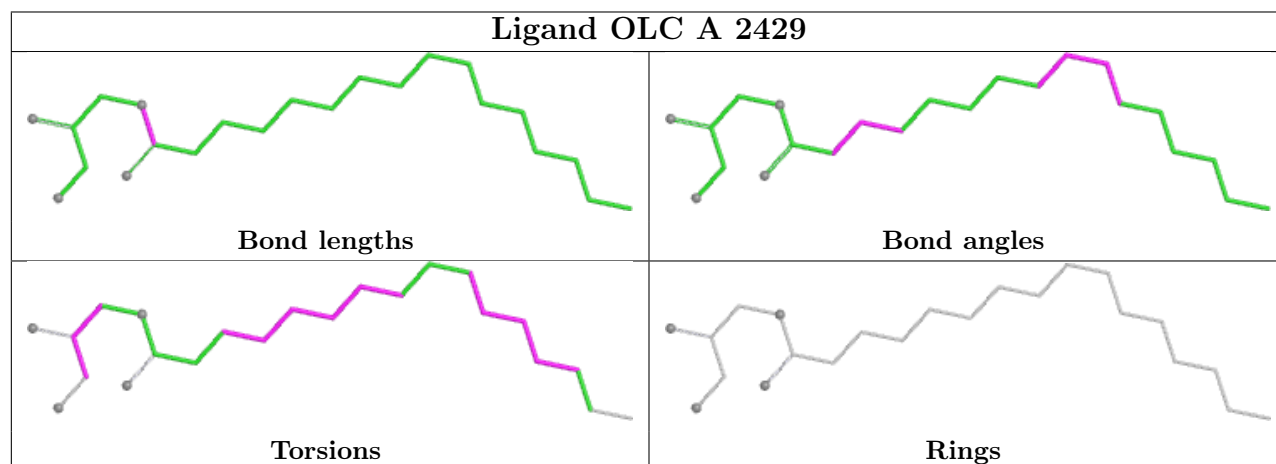
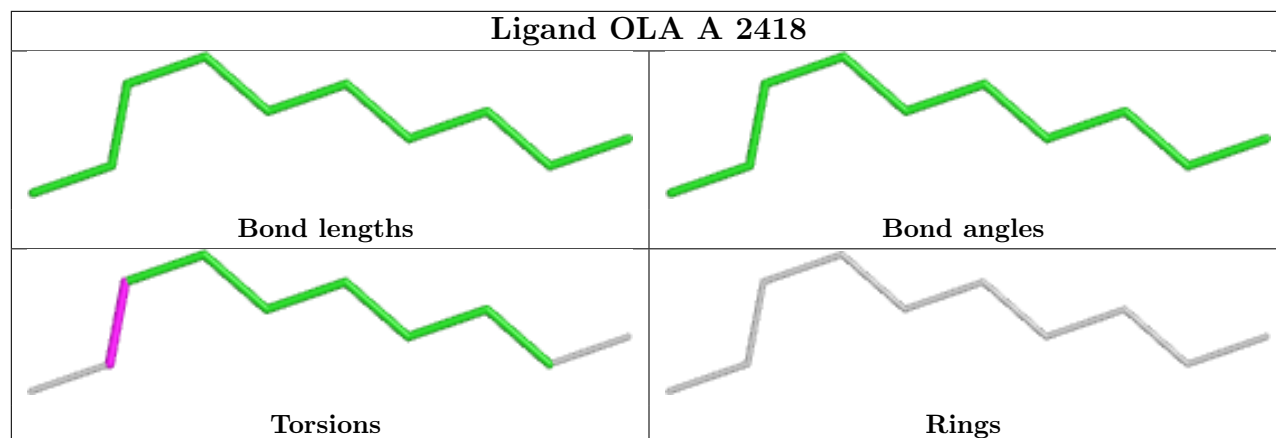






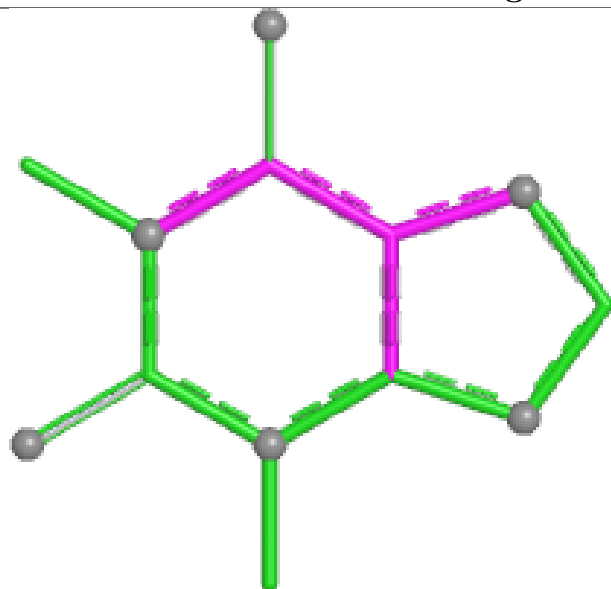




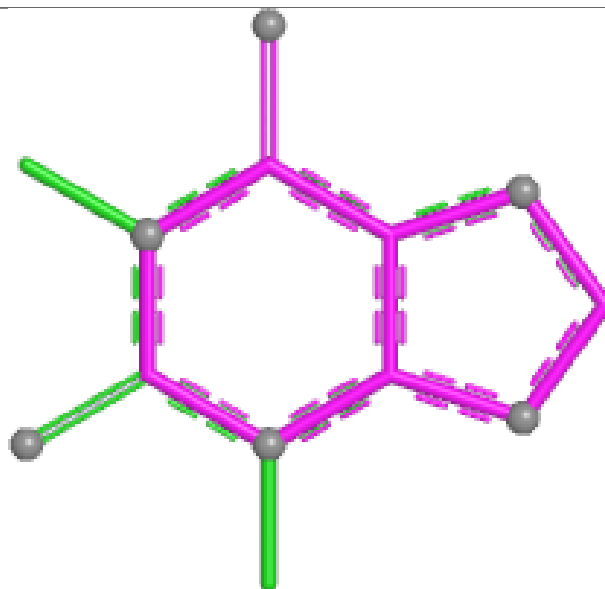




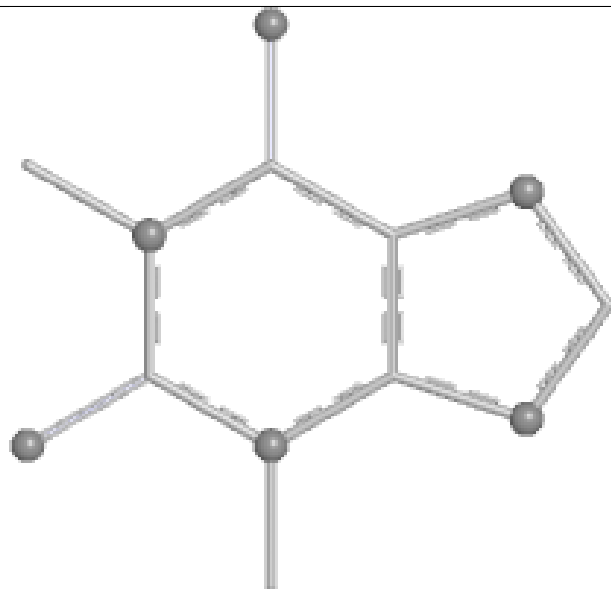
## Ligand TEP A 2401



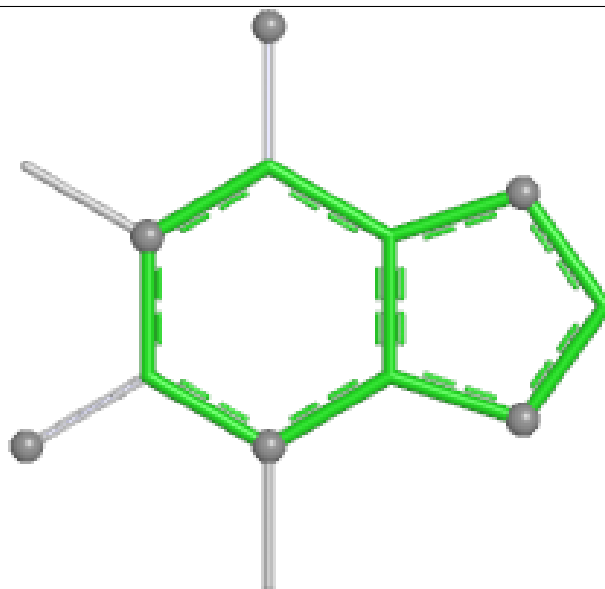
Bond lengths



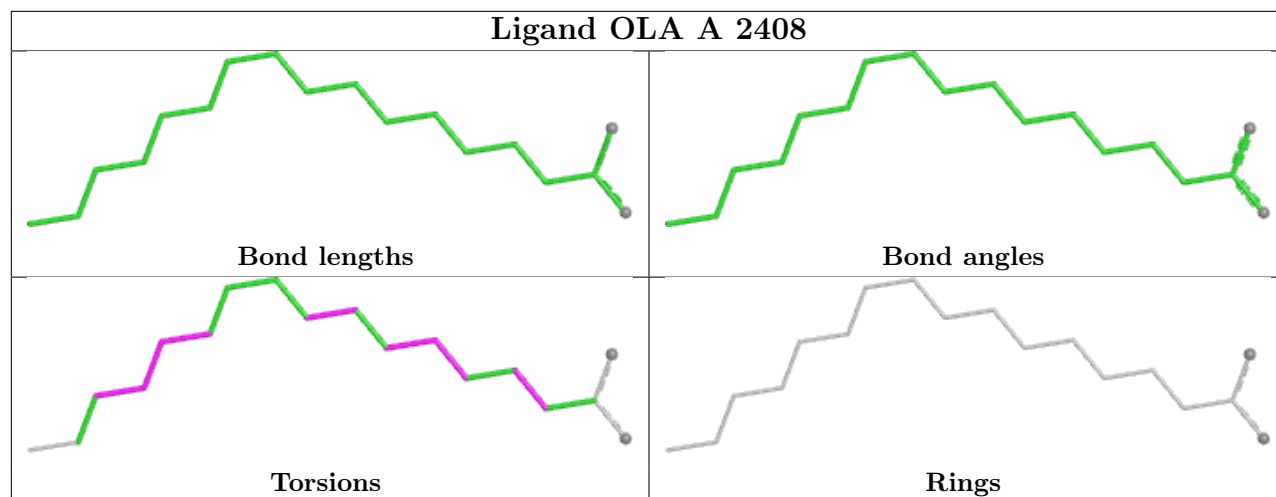
Bond angles



Torsions



Rings



## 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

### 6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	390/433 (90%)	0.79	62 (15%) <b>5</b> <b>5</b>	13, 37, 80, 110	14 (3%)

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	1061	PHE	7.1
1	A	307	VAL	5.4
1	A	308	LEU	5.0
1	A	1021	ASP	4.4
1	A	306	HIS	4.4
1	A	1106	LEU	4.1
1	A	1017	ILE	4.1
1	A	1016	VAL	4.0
1	A	29	TRP	4.0
1	A	1023	ALA	4.0
1	A	1026	VAL	3.8
1	A	1036	ALA	3.4
1	A	1059	LYS	3.4
1	A	1105	TYR	3.4
1	A	1037	ALA	3.3
1	A	1093	GLN	3.3
1	A	1084	VAL	3.2
1	A	305	SER	3.1
1	A	1022	ASN	2.9
1	A	205[A]	ARG	2.9
1	A	1028	ASP	2.9
1	A	1087	ALA	2.9
1	A	1029	ALA	2.8
1	A	303	ILE	2.8
1	A	1035	ALA	2.8
1	A	290	TYR	2.7
1	A	1010	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	A	208	LEU	2.7
1	A	1063	HIS	2.7
1	A	1003	LEU	2.6
1	A	1094	LEU	2.6
1	A	0	ALA	2.6
1	A	1040	ALA	2.6
1	A	1007	TRP	2.6
1	A	117	THR	2.5
1	A	157	GLN	2.5
1	A	1030	LEU	2.5
1	A	1079	ALA	2.5
1	A	1020	ALA	2.4
1	A	1039	ASP	2.4
1	A	111	ARG	2.4
1	A	32	TRP	2.4
1	A	1019	LYS	2.4
1	A	1086	GLU	2.3
1	A	309	ARG	2.3
1	A	1038	LEU	2.3
1	A	1062	ARG	2.3
1	A	110	LEU	2.3
1	A	1024	ALA	2.2
1	A	1102	ILE	2.2
1	A	1006	ASN	2.2
1	A	1068	LEU	2.2
1	A	1001	ALA	2.2
1	A	1042	LYS	2.2
1	A	1099	ASN	2.2
1	A	1091	ALA	2.2
1	A	1078	LEU	2.1
1	A	35	SER	2.1
1	A	300[A]	ARG	2.1
1	A	1077	LYS	2.1
1	A	1014	LEU	2.1
1	A	1090	ALA	2.0

## 6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

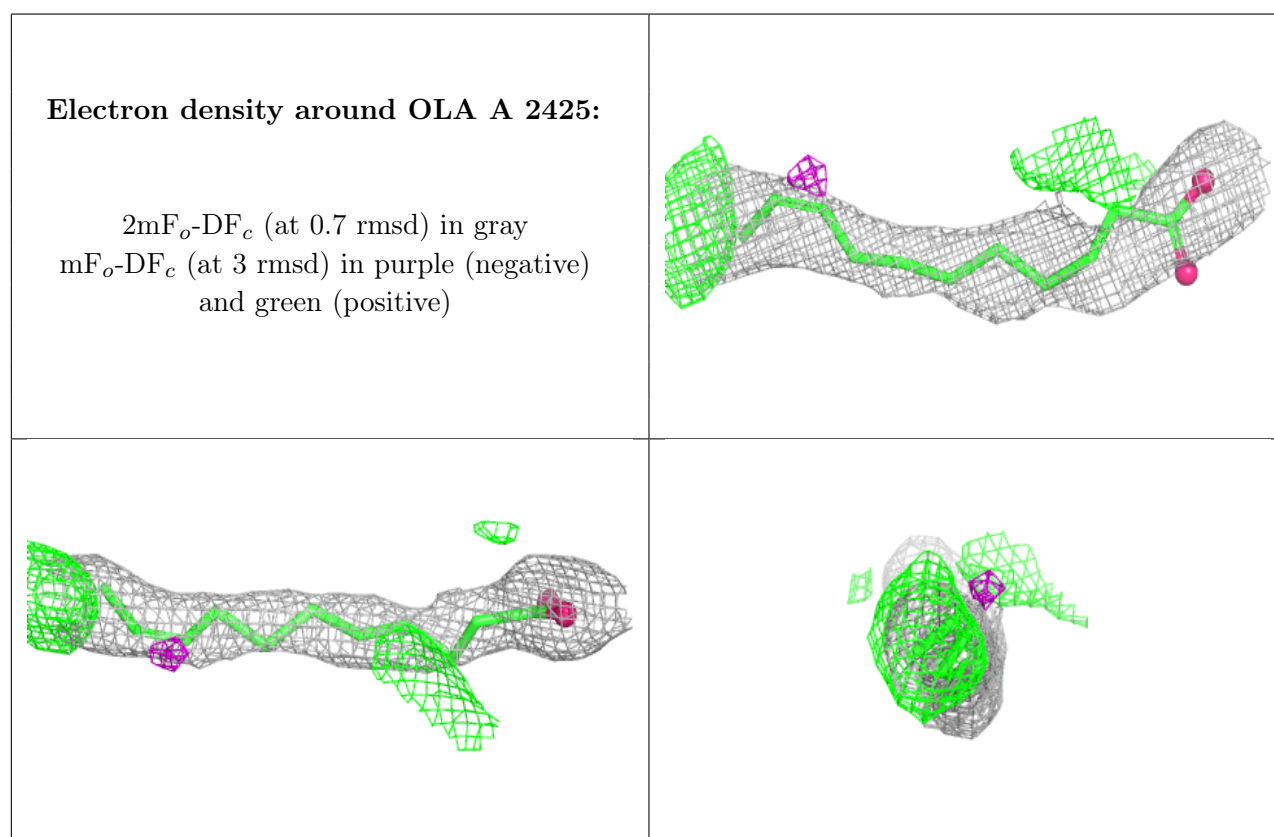
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
5	OLA	A	2425	13/20	0.66	0.20	47,54,78,89	0
5	OLA	A	2406	10/20	0.69	0.20	57,67,79,94	0
5	OLA	A	2420	12/20	0.73	0.22	51,59,62,63	0
5	OLA	A	2414	11/20	0.73	0.17	42,52,73,84	0
5	OLA	A	2412	8/20	0.74	0.18	54,55,57,58	0
5	OLA	A	2426	20/20	0.75	0.22	52,71,79,91	0
5	OLA	A	2423	15/20	0.77	0.21	48,60,76,79	0
5	OLA	A	2417	12/20	0.77	0.19	40,46,56,65	0
5	OLA	A	2413	15/20	0.77	0.20	61,66,72,75	0
6	OLC	A	2428	17/25	0.77	0.18	57,68,80,86	0
5	OLA	A	2408	18/20	0.78	0.17	41,52,73,78	0
5	OLA	A	2421	9/20	0.78	0.24	42,56,78,85	0
5	OLA	A	2409	20/20	0.78	0.20	45,60,75,82	0
5	OLA	A	2415	19/20	0.79	0.21	52,60,67,71	0
5	OLA	A	2418	10/20	0.79	0.22	47,58,65,73	0
5	OLA	A	2419	11/20	0.79	0.22	41,45,62,65	0
5	OLA	A	2416	20/20	0.80	0.20	45,62,70,71	0
3	TEP	A	2401	13/13	0.81	0.12	26,30,32,35	0
5	OLA	A	2422	8/20	0.81	0.21	53,57,59,61	0
6	OLC	A	2431	23/25	0.81	0.17	40,52,79,87	0
6	OLC	A	2429	23/25	0.82	0.18	48,57,75,79	0
5	OLA	A	2410	20/20	0.82	0.18	54,60,81,86	0
6	OLC	A	2432	18/25	0.82	0.18	47,57,74,89	0
6	OLC	A	2433	16/25	0.82	0.16	39,48,57,64	0
5	OLA	A	2407	9/20	0.83	0.17	40,43,61,63	0
5	OLA	A	2411	12/20	0.84	0.16	36,54,68,70	0
5	OLA	A	2424	15/20	0.85	0.16	43,47,55,55	0
8	PEG	A	2435	7/7	0.85	0.14	47,53,55,58	0
9	DMS	A	2436[S]	4/4	0.86	0.20	45,48,54,59	4
5	OLA	A	2405	20/20	0.87	0.17	43,54,66,66	0
6	OLC	A	2430	23/25	0.88	0.15	39,52,67,73	0
6	OLC	A	2427	16/25	0.89	0.13	42,50,63,68	0

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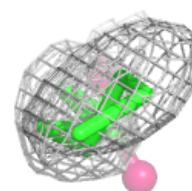
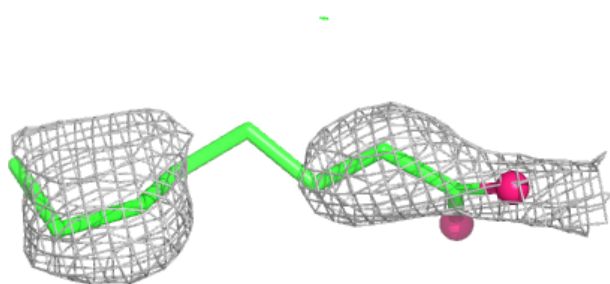
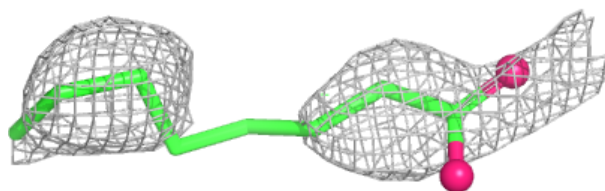
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
4	CLR	A	2402	28/28	0.92	0.11	35,39,56,65	0
4	CLR	A	2404	28/28	0.93	0.09	30,35,46,55	0
4	CLR	A	2403	28/28	0.93	0.10	36,41,49,51	0
7	A1CPX	A	2434	20/20	0.94	0.10	33,39,60,68	0
2	NA	A	2400	1/1	0.98	0.06	35,35,35,35	0

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

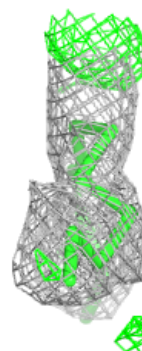
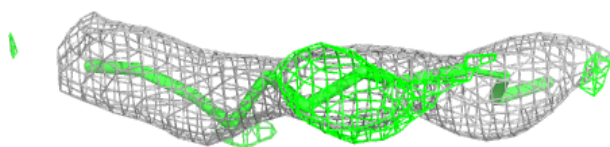
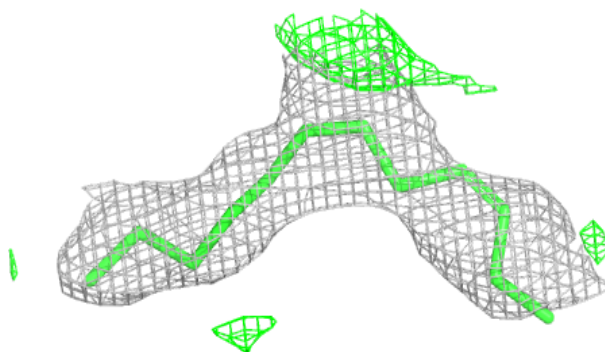


**Electron density around OLA A 2406:**

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

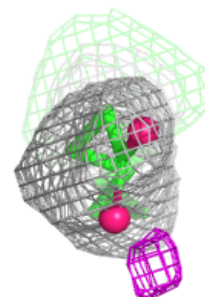
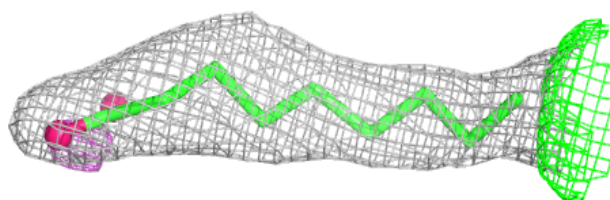
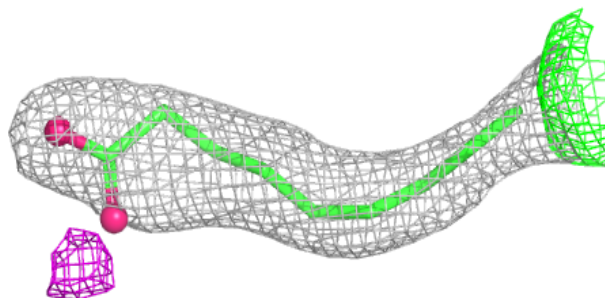
**Electron density around OLA A 2420:**

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

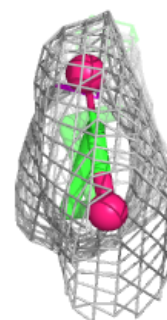
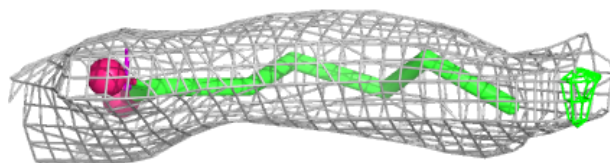
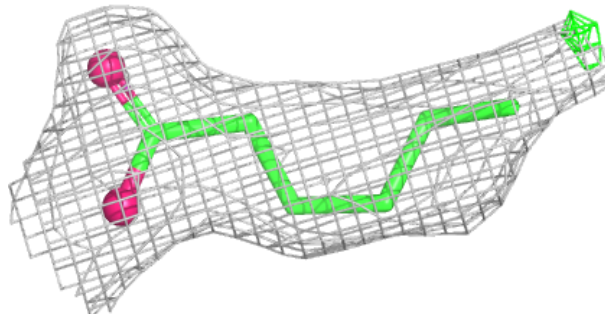


**Electron density around OLA A 2414:**

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

**Electron density around OLA A 2412:**

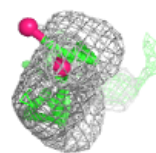
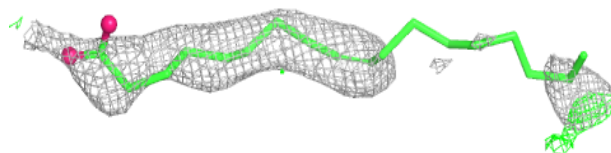
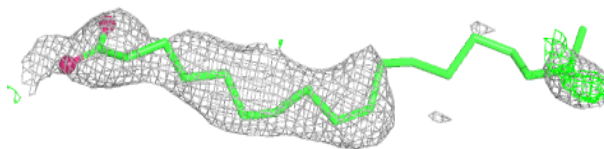
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



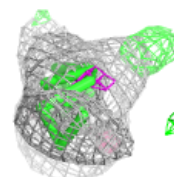
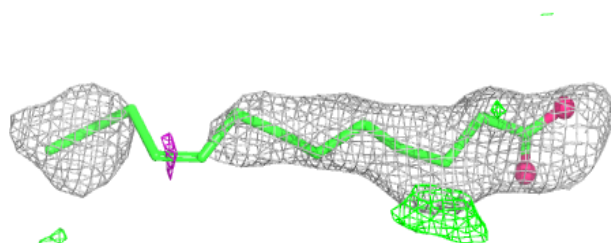
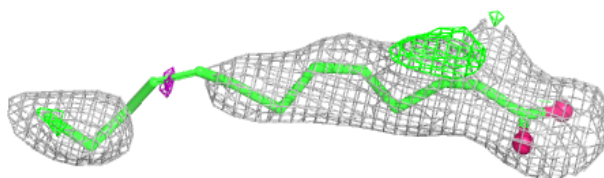


**Electron density around OLA A 2426:**

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

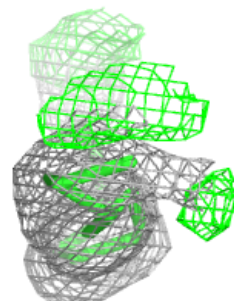
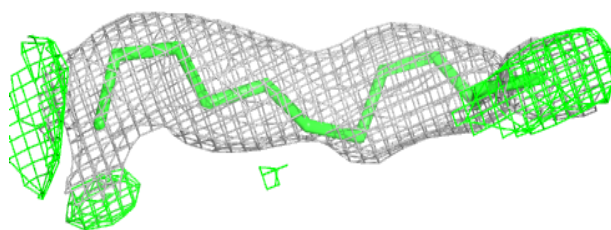
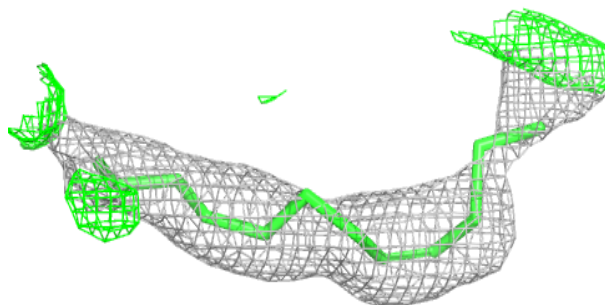
**Electron density around OLA A 2423:**

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

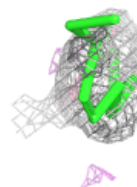
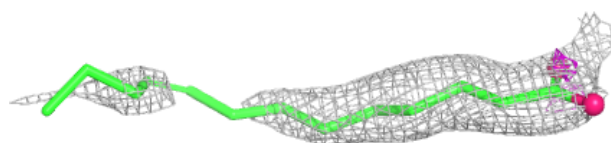
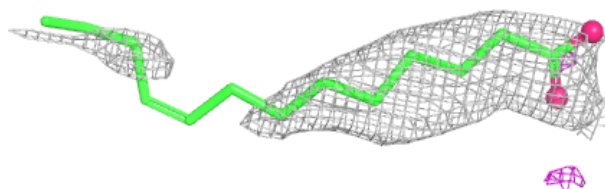


**Electron density around OLA A 2417:**

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

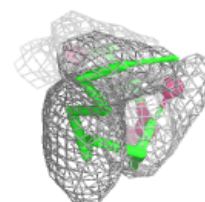
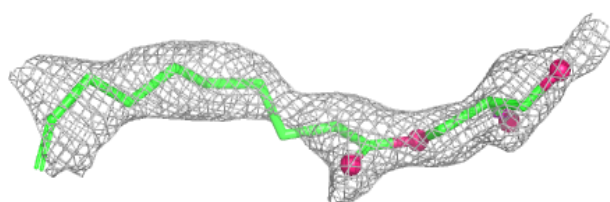
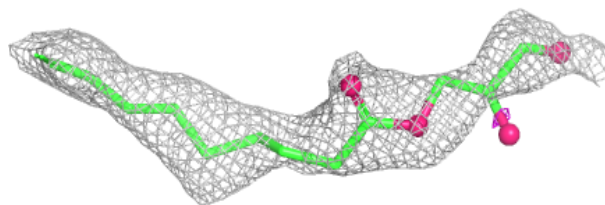
**Electron density around OLA A 2413:**

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

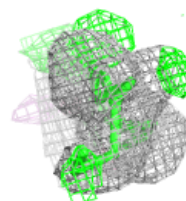
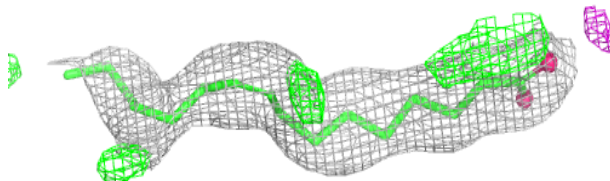
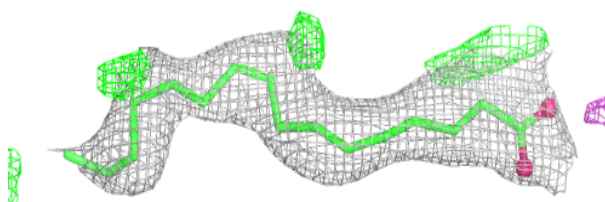


**Electron density around OLC A 2428:**

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

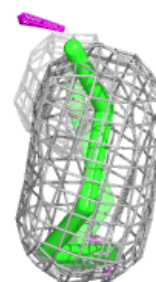
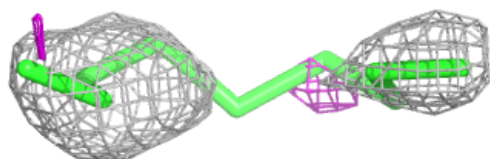
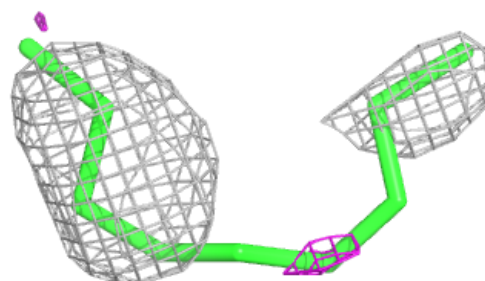
**Electron density around OLA A 2408:**

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

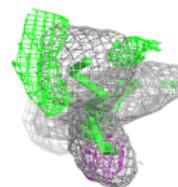
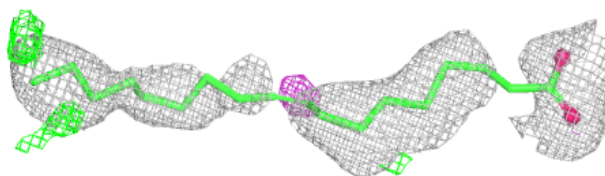
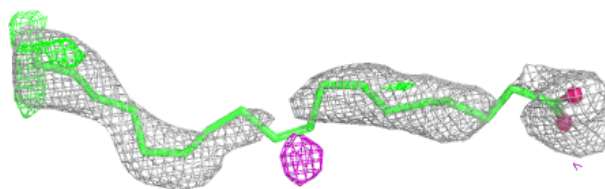


**Electron density around OLA A 2421:**

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

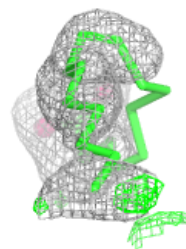
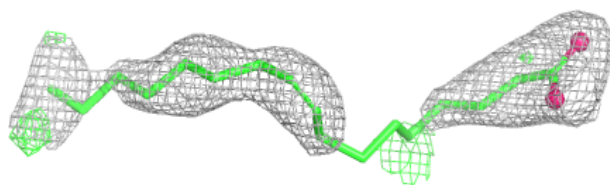
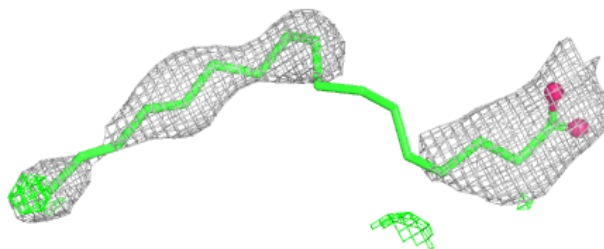
**Electron density around OLA A 2409:**

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

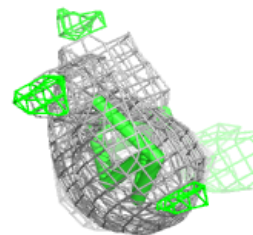
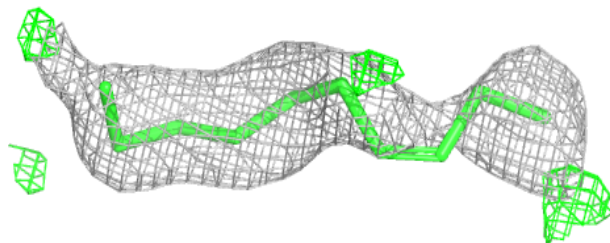
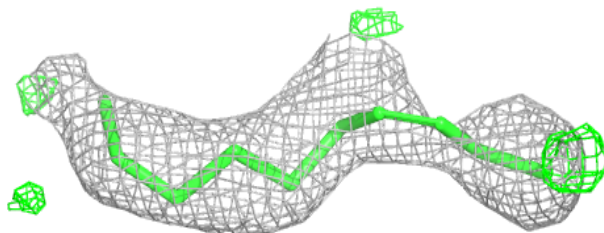


**Electron density around OLA A 2415:**

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

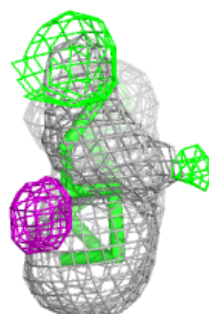
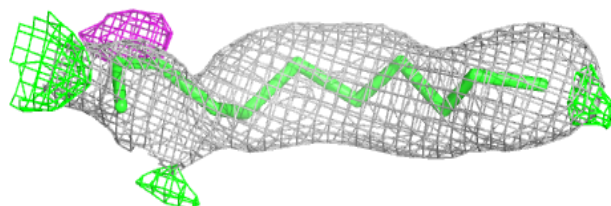
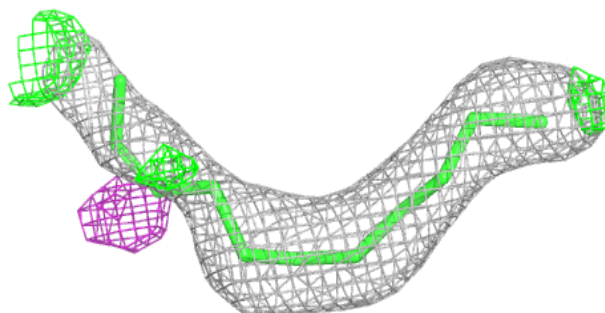
**Electron density around OLA A 2418:**

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

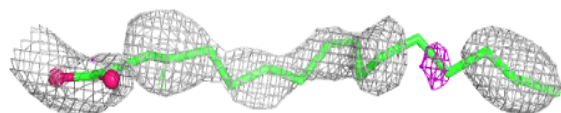
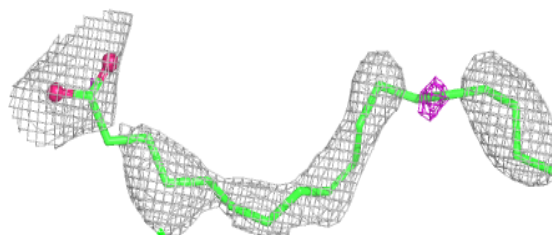


**Electron density around OLA A 2419:**

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

**Electron density around OLA A 2416:**

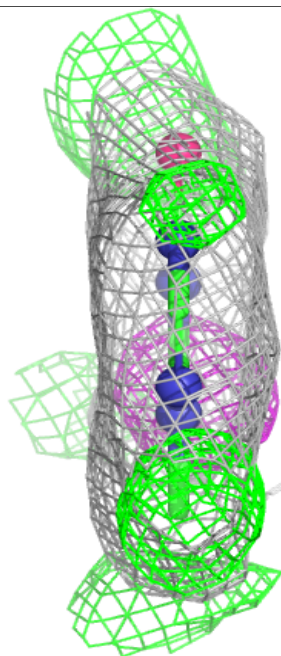
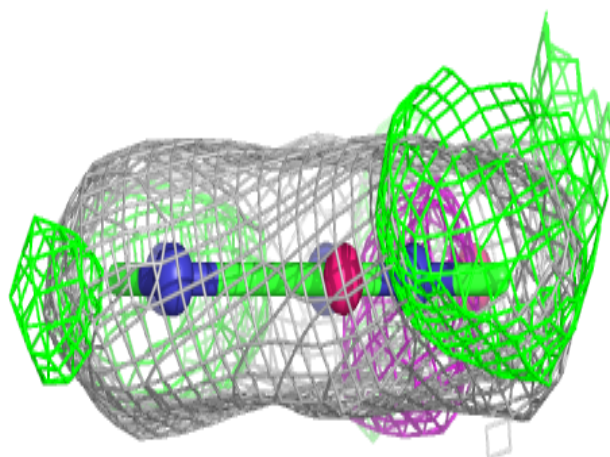
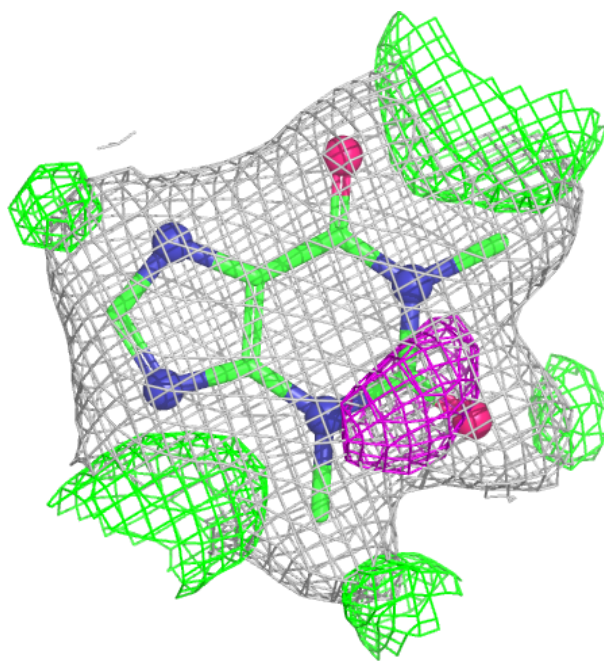
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)





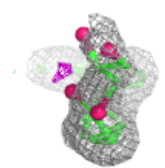
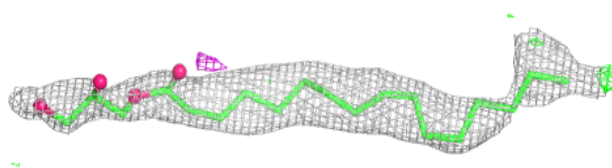
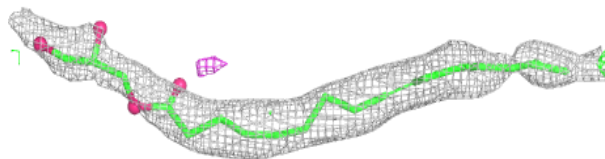
**Electron density around TEP A 2401:**

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

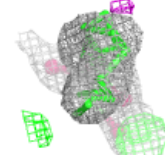
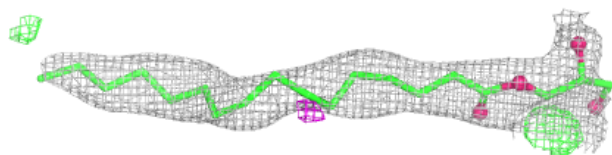
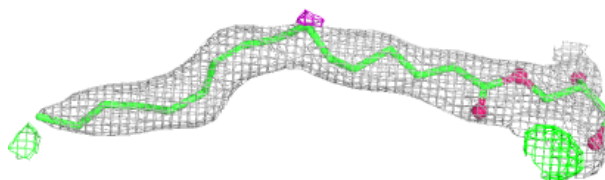


**Electron density around OLC A 2431:**

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

**Electron density around OLC A 2429:**

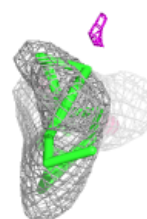
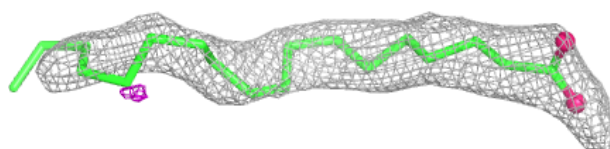
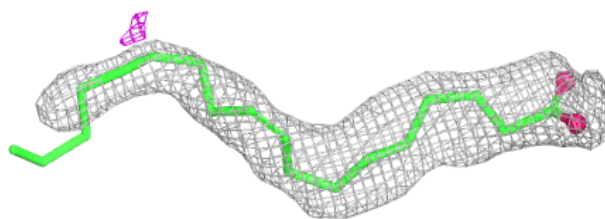
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



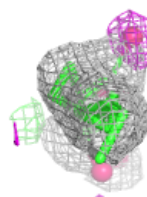
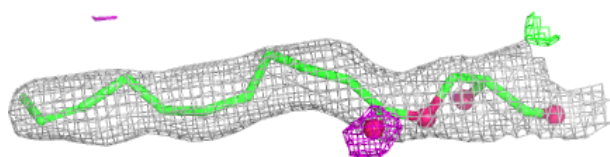
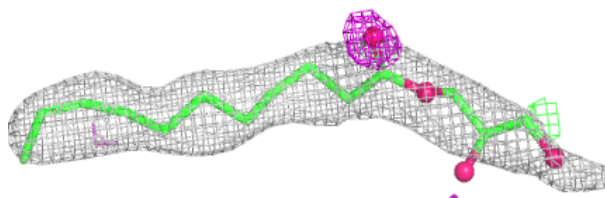


**Electron density around OLA A 2410:**

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

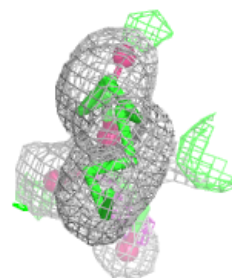
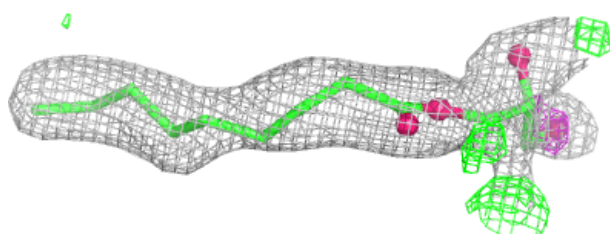
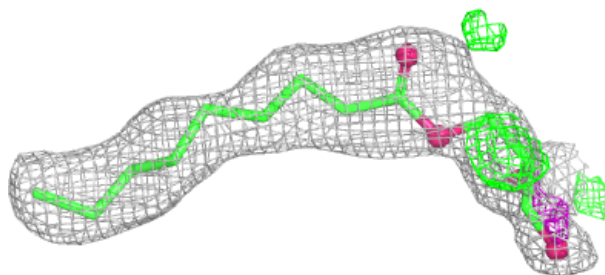
**Electron density around OLC A 2432:**

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

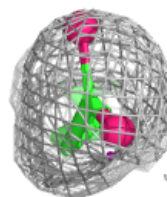
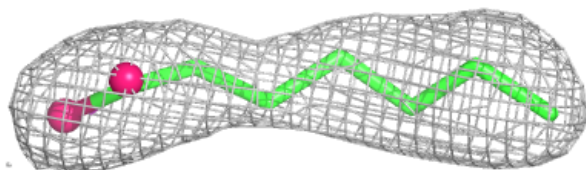
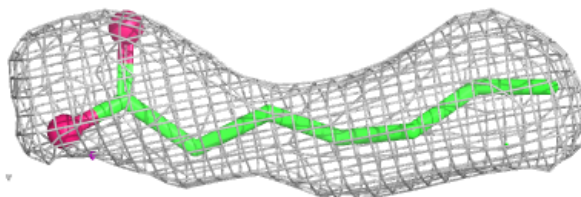


**Electron density around OLC A 2433:**

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

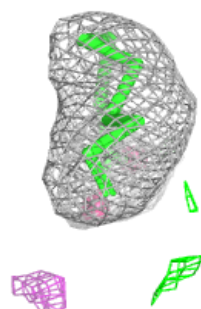
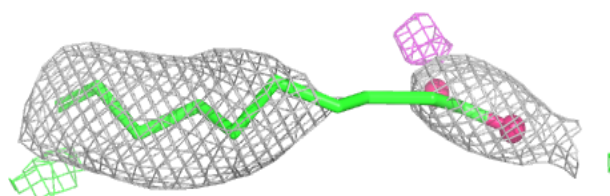
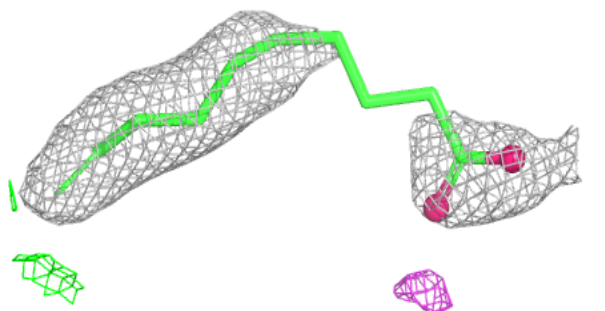
**Electron density around OLA A 2407:**

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

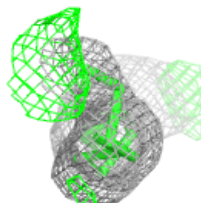
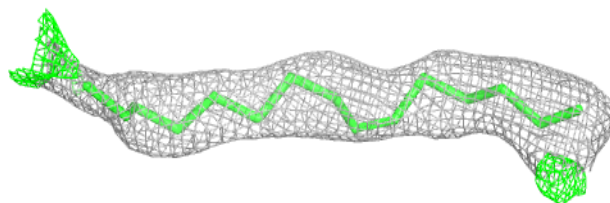
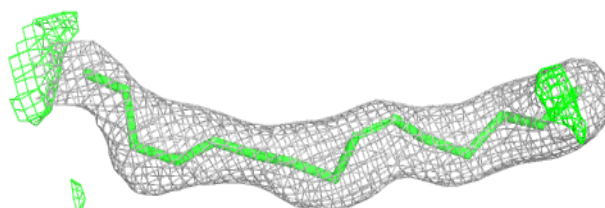


**Electron density around OLA A 2411:**

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

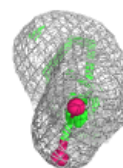
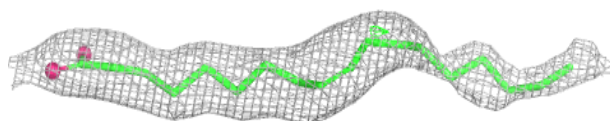
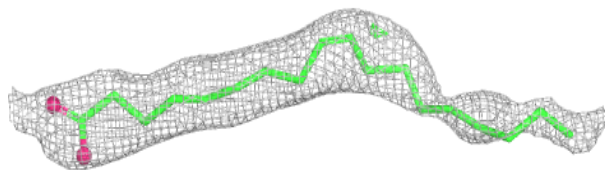
**Electron density around OLA A 2424:**

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

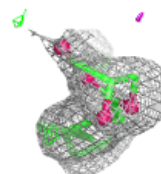
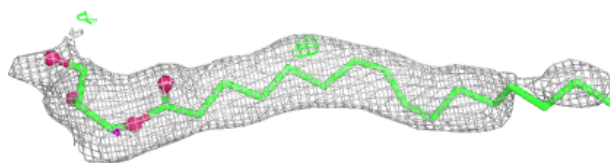
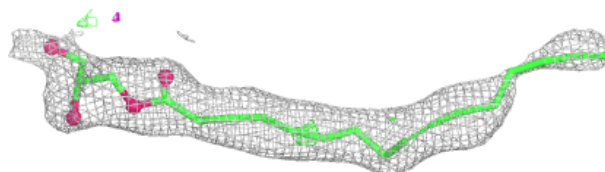


**Electron density around OLA A 2405:**

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

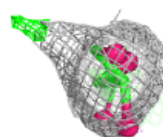
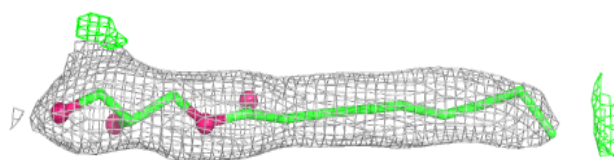
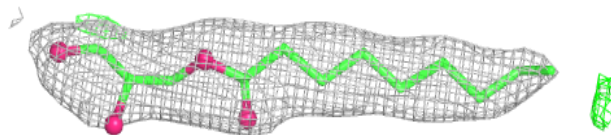
**Electron density around OLC A 2430:**

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

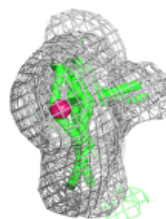
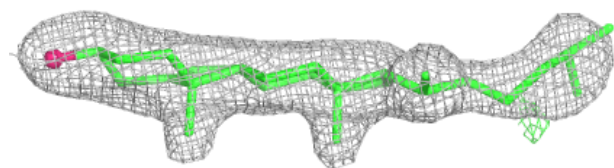
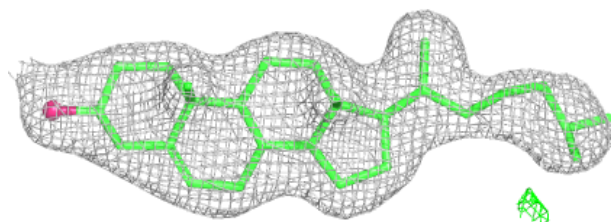


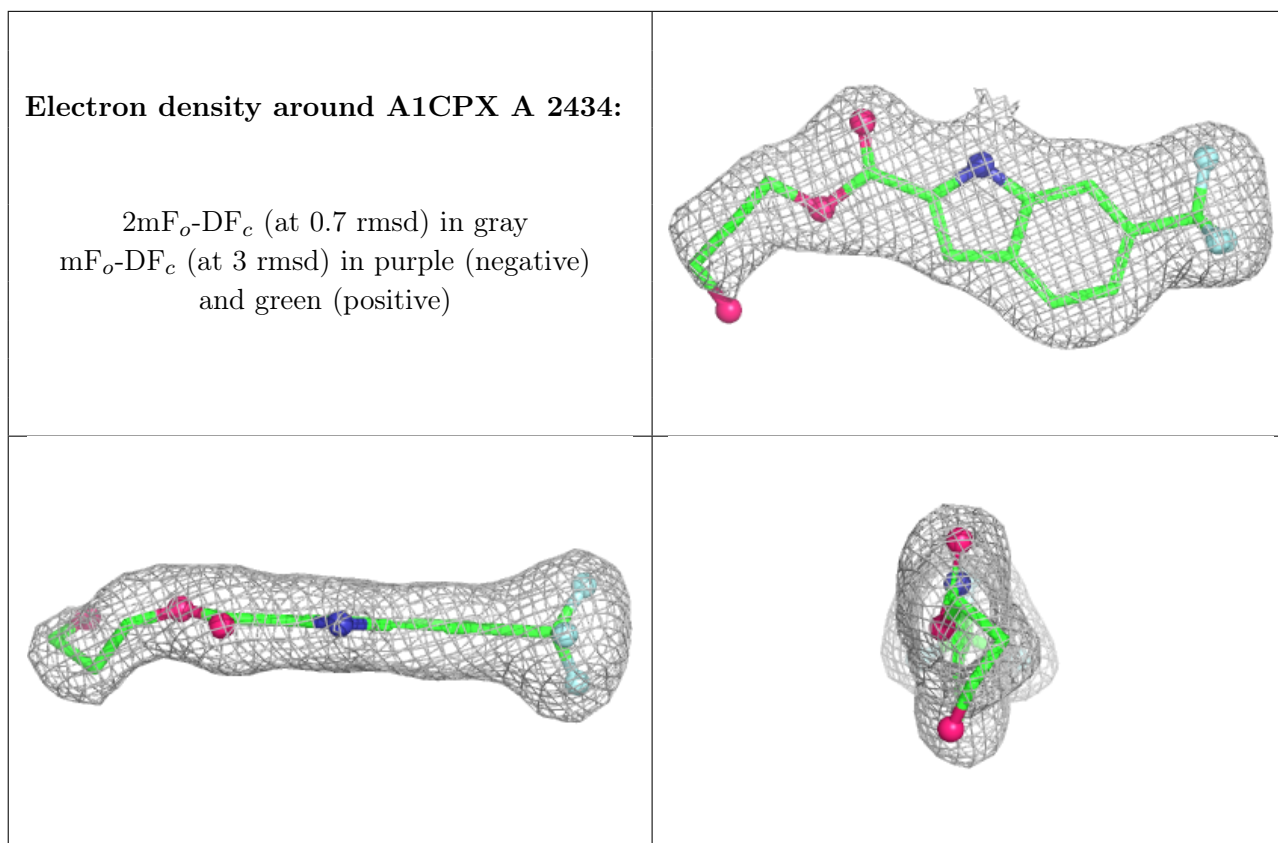
**Electron density around OLC A 2427:**

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

**Electron density around CLR A 2404:**

$2mF_o-DF_c$  (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.