



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 02:22 AM UTC

PDB ID : 6YX9 / pdb_00006yx9
Title : Cryogenic human adiponectin receptor 2 (ADIPOR2) at 2.4 Å resolution determined by Serial Crystallography (SSX) using CrystalDirect
Authors : Healey, R.D.; Basu, S.; Humm, A.S.; Leyrat, C.; Dupeux, F.; Pica, A.; Granier, S.; Marquez, J.A.
Deposited on : 2020-04-30
Resolution : 2.40 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

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

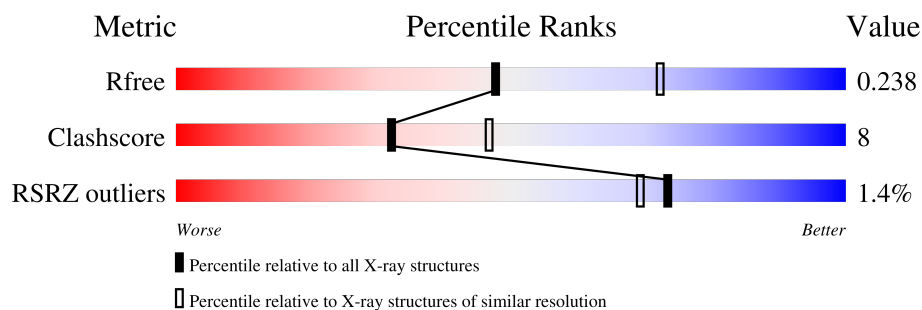
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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

Mol	Chain	Length	Quality of chain
1	A	292	2% 77% 21% .
2	H	121	91% 8% .
3	L	108	% 81% 18% .

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 4309 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 Adiponectin receptor protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	285	Total	C	N	O	S	0	0	0
			2297	1541	373	366	17			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	expression tag	UNP Q86V24
A	-3	GLY	-	expression tag	UNP Q86V24
A	-2	SER	-	expression tag	UNP Q86V24
A	-1	GLU	-	expression tag	UNP Q86V24
A	0	PHE	-	expression tag	UNP Q86V24

- Molecule 2 is a protein called V REGION HEAVY CHAIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	121	Total	C	N	O	S	0	0	0
			935	596	151	184	4			

- Molecule 3 is a protein called V REGION LIGHT CHAIN.

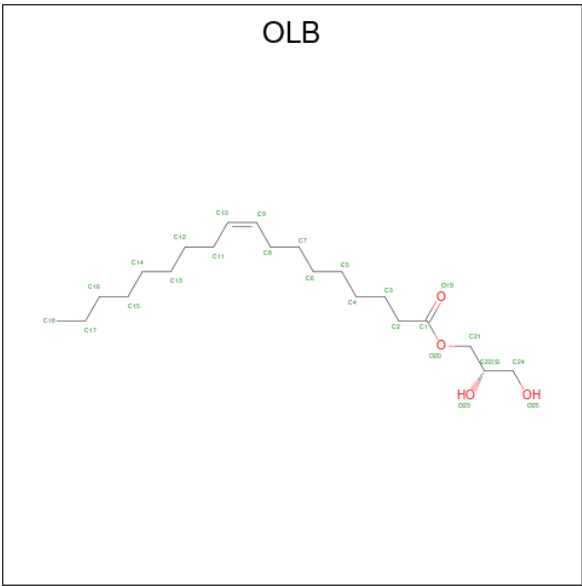
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	108	Total	C	N	O	S	0	0	0
			826	520	138	165	3			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn) (labeled as "Ligand of Interest" by depositor).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		

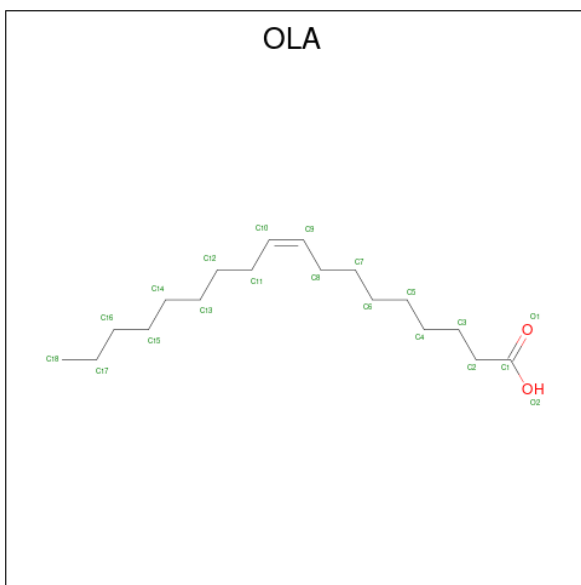
- Molecule 5 is (2S)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (CCD ID: OLB) (formula:

C₂₁H₄₀O₄) (labeled as "Ligand of Interest" by depositor).



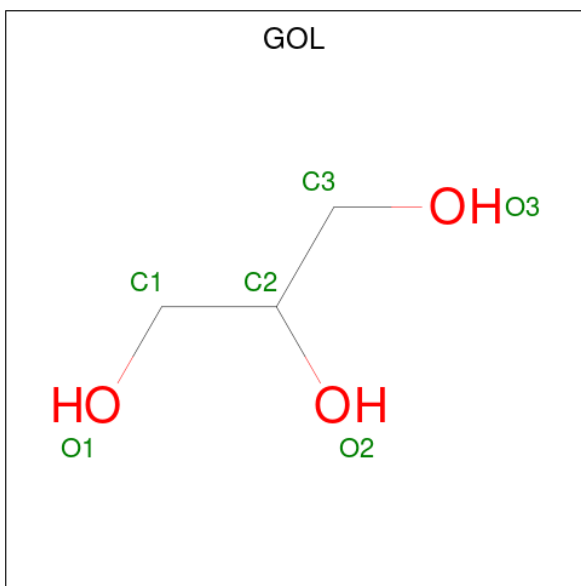
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		
5	A	1	Total	C	O	0	0
			25	21	4		

- Molecule 6 is OLEIC ACID (CCD ID: OLA) (formula: C₁₈H₃₄O₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$) (labeled as "Ligand of Interest" by depositor).



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

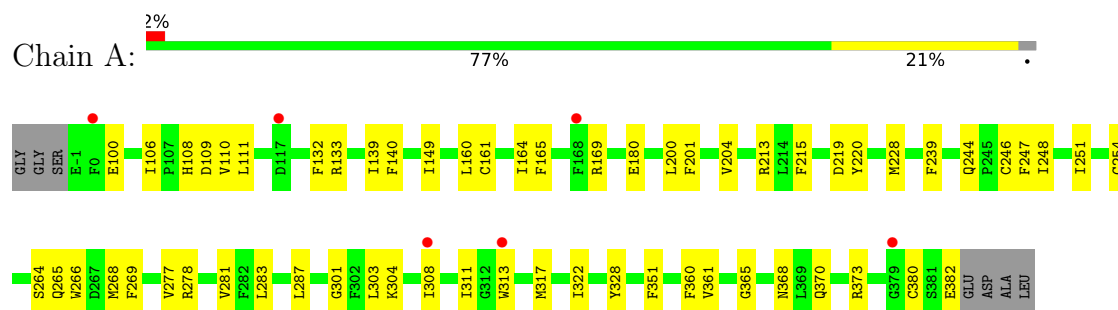
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	34	Total 34	O 34	0	0
8	H	42	Total 42	O 42	0	0
8	L	23	Total 23	O 23	0	0

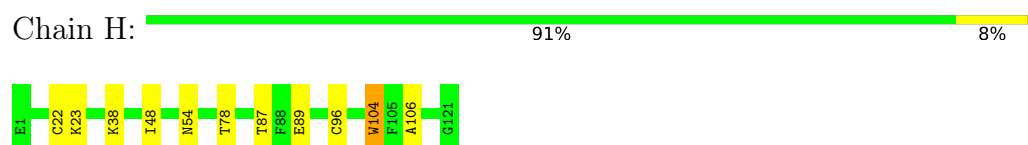
3 Residue-property plots [i](#)

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.

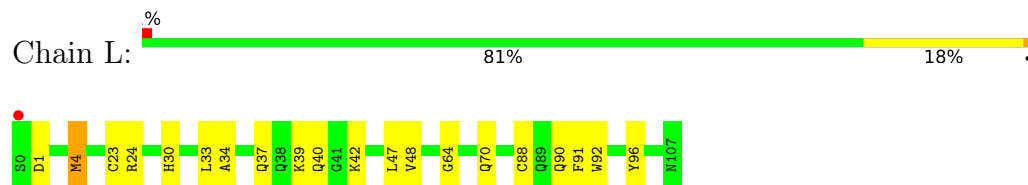
- Molecule 1: Adiponectin receptor protein 2



- Molecule 2: V REGION HEAVY CHAIN



- Molecule 3: V REGION LIGHT CHAIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2 ₁	Depositor
Cell constants a, b, c, α , β , γ	74.72Å 100.58Å 109.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	22.00 – 2.40 22.00 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.4 (22.00-2.40) 99.2 (22.00-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.02 (at 2.41Å)	Xtriage
Refinement program	BUSTER 2.10.3 (3-OCT-2019)	Depositor
R, R_{free}	0.207 , 0.248 (Not available) , 0.238	Depositor DCC
R_{free} test set	1632 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	60.7	Xtriage
Anisotropy	0.622	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 63.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4309	wwPDB-VP
Average B, all atoms (Å ²)	80.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.50% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, OLB, ZN, OLA

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	0.92	0/2381	1.23	12/3236 (0.4%)
2	H	0.84	0/959	1.09	2/1301 (0.2%)
3	L	0.87	1/846 (0.1%)	1.03	1/1148 (0.1%)
All	All	0.89	1/4186 (0.0%)	1.16	15/5685 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	4	MET	SD-CE	-5.93	1.64	1.79

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	GLU	CB-CG-CD	6.04	122.87	112.60
1	A	140	PHE	CA-CB-CG	-6.03	107.77	113.80
1	A	368	ASN	CA-CB-CG	5.83	118.43	112.60
1	A	265	GLN	CA-C-N	5.61	128.81	120.95
1	A	265	GLN	C-N-CA	5.61	128.81	120.95
1	A	360	PHE	CA-CB-CG	-5.58	108.22	113.80
1	A	109	ASP	CA-CB-CG	5.49	118.09	112.60
1	A	204	VAL	CA-C-N	5.44	127.89	120.54
1	A	204	VAL	C-N-CA	5.44	127.89	120.54
2	H	104	TRP	N-CA-C	-5.43	101.42	110.17
2	H	54	ASN	CA-CB-CG	5.34	117.94	112.60
1	A	219	ASP	CA-C-N	5.26	127.28	120.44
1	A	219	ASP	C-N-CA	5.26	127.28	120.44
3	L	1	ASP	N-CA-C	5.24	118.00	111.24
1	A	219	ASP	CA-CB-CG	5.12	117.72	112.60

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	2297	0	2271	45	0
2	H	935	0	894	6	0
3	L	826	0	791	12	0
4	A	1	0	0	0	0
5	A	125	0	200	16	0
6	A	20	0	33	4	0
7	L	6	0	8	0	0
8	A	34	0	0	3	0
8	H	42	0	0	0	0
8	L	23	0	0	0	0
All	All	4309	0	4197	66	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 8.

All (66) 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)
6:A:407:OLA:H32	8:A:502:HOH:O	1.43	1.16
1:A:373:ARG:HH21	1:A:373:ARG:HG2	1.34	0.90
2:H:22:CYS:HG	2:H:96:CYS:HG	0.92	0.84
1:A:351:PHE:HE1	6:A:407:OLA:H112	1.52	0.75
1:A:373:ARG:HG2	1:A:373:ARG:NH2	2.06	0.69
1:A:228:MET:HE3	1:A:254:CYS:HA	1.74	0.69
2:H:23:LYS:HG2	2:H:78:THR:HG22	1.73	0.67
1:A:373:ARG:HH21	1:A:373:ARG:CG	2.08	0.65
5:A:402:OLB:H2A	3:L:30:HIS:HD2	1.61	0.63
2:H:22:CYS:HG	2:H:96:CYS:CB	2.12	0.63
1:A:160:LEU:O	1:A:164:ILE:HG12	1.99	0.62
1:A:303:LEU:HB3	1:A:304:LYS:HZ2	1.66	0.61
1:A:220:TYR:HB3	1:A:264:SER:HB2	1.82	0.60
1:A:311:ILE:HD13	5:A:406:OLB:H22	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:PHE:HB2	5:A:402:OLB:H5A	1.86	0.57
1:A:133:ARG:HB2	5:A:402:OLB:H3	1.87	0.57
1:A:165:PHE:HD1	5:A:403:OLB:H3A	1.70	0.57
1:A:132:PHE:N	5:A:402:OLB:H2	2.20	0.56
1:A:161:CYS:HB2	5:A:403:OLB:H29	1.89	0.55
1:A:139:ILE:HG12	1:A:149:ILE:HG21	1.89	0.54
3:L:40:GLN:NE2	3:L:40:GLN:H	2.04	0.54
1:A:106:ILE:HD12	1:A:110:VAL:HG23	1.90	0.54
1:A:169:ARG:HH12	1:A:370:GLN:NE2	2.06	0.53
1:A:287:LEU:HD11	1:A:322:ILE:HD11	1.90	0.52
1:A:283:LEU:HD21	1:A:322:ILE:HD13	1.92	0.51
3:L:48:VAL:HG11	3:L:64:GLY:HA3	1.91	0.51
3:L:91:PHE:HA	3:L:96:TYR:CD1	2.45	0.51
1:A:328:TYR:CE1	6:A:407:OLA:H21	2.46	0.50
2:H:87:THR:HG22	2:H:89:GLU:H	1.76	0.49
1:A:266:TRP:CE2	1:A:268:MET:HB2	2.46	0.49
1:A:165:PHE:HB3	5:A:403:OLB:C3	2.43	0.48
1:A:200:LEU:HD11	5:A:402:OLB:H40	1.96	0.48
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.96	0.47
1:A:301:GLY:HA3	1:A:304:LYS:HZ3	1.80	0.47
3:L:23:CYS:SG	3:L:88:CYS:SG	3.06	0.47
1:A:317:MET:HE2	5:A:406:OLB:H40	1.97	0.47
1:A:380:CYS:SG	1:A:382:GLU:OE2	2.72	0.47
1:A:247:PHE:O	1:A:251:ILE:HG12	2.14	0.47
1:A:165:PHE:O	5:A:403:OLB:H3	2.15	0.47
1:A:165:PHE:CD1	5:A:403:OLB:H3A	2.50	0.47
1:A:100:GLU:CD	3:L:92:TRP:HE1	2.25	0.45
2:H:104:TRP:CZ3	2:H:106:ALA:HB2	2.51	0.45
1:A:269:PHE:O	1:A:278:ARG:HD3	2.16	0.45
1:A:213:ARG:NE	1:A:213:ARG:HA	2.32	0.44
1:A:277:VAL:O	1:A:281:VAL:HG23	2.18	0.44
1:A:317:MET:SD	1:A:361:VAL:HG22	2.57	0.44
1:A:382:GLU:H	1:A:382:GLU:HG3	1.58	0.44
5:A:402:OLB:H221	5:A:402:OLB:H27	1.86	0.44
1:A:322:ILE:HG23	5:A:405:OLB:H34	2.00	0.44
3:L:24:ARG:HG3	3:L:70:GLN:HE21	1.82	0.44
1:A:304:LYS:O	1:A:308:ILE:HG12	2.18	0.43
3:L:33:LEU:HG	3:L:34:ALA:N	2.33	0.43
1:A:301:GLY:HA3	1:A:304:LYS:NZ	2.33	0.43
2:H:38:LYS:HB2	2:H:48:ILE:HD11	2.00	0.43
1:A:132:PHE:H	5:A:402:OLB:H2	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:407:OLA:C3	8:A:502:HOH:O	2.27	0.42
1:A:244:GLN:O	1:A:248:ILE:HG12	2.20	0.42
1:A:201:PHE:HB2	1:A:215:PHE:HB3	2.02	0.42
1:A:239:PHE:HB2	1:A:246:CYS:HB2	2.02	0.42
3:L:4:MET:HE2	3:L:90:GLN:HB3	2.00	0.42
1:A:108:HIS:HA	1:A:111:LEU:HD12	2.02	0.41
5:A:402:OLB:O19	8:A:501:HOH:O	2.22	0.41
1:A:220:TYR:CB	1:A:264:SER:HB2	2.50	0.41
3:L:39:LYS:HB2	3:L:42:LYS:HE3	2.03	0.41
3:L:33:LEU:HD11	3:L:88:CYS:HB2	2.02	0.41
1:A:313:TRP:CE2	1:A:365:GLY:HA3	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

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

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

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 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	OLB	A	403	-	24,24,24	0.25	0	25,25,25	0.22	0
6	OLA	A	407	4	19,19,19	0.40	0	19,19,19	0.52	0
5	OLB	A	404	-	24,24,24	0.21	0	25,25,25	0.26	0
5	OLB	A	406	-	24,24,24	0.23	0	25,25,25	0.28	0
5	OLB	A	405	-	24,24,24	0.22	0	25,25,25	0.32	0
5	OLB	A	402	-	24,24,24	0.27	0	25,25,25	0.37	0
7	GOL	L	201	-	5,5,5	0.14	0	5,5,5	0.28	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	OLB	A	403	-	-	11/24/24/24	-
6	OLA	A	407	4	-	7/17/17/17	-
5	OLB	A	404	-	-	7/24/24/24	-
5	OLB	A	406	-	-	4/24/24/24	-
5	OLB	A	405	-	-	5/24/24/24	-
5	OLB	A	402	-	-	15/24/24/24	-
7	GOL	L	201	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	402	OLB	C21-C22-C24-O25
5	A	402	OLB	O23-C22-C24-O25
7	L	201	GOL	O1-C1-C2-O2
7	L	201	GOL	O1-C1-C2-C3
6	A	407	OLA	C2-C3-C4-C5
5	A	402	OLB	O20-C21-C22-O23
5	A	405	OLB	C1-C2-C3-C4
5	A	403	OLB	O20-C21-C22-O23
5	A	402	OLB	C11-C12-C13-C14
5	A	404	OLB	C2-C3-C4-C5
5	A	402	OLB	C12-C13-C14-C15
6	A	407	OLA	C13-C14-C15-C16
5	A	403	OLB	C14-C15-C16-C17
5	A	402	OLB	C2-C3-C4-C5
5	A	402	OLB	C3-C4-C5-C6
5	A	403	OLB	O20-C21-C22-C24
5	A	402	OLB	C2-C1-O20-C21
5	A	402	OLB	C10-C11-C12-C13
6	A	407	OLA	C10-C11-C12-C13
5	A	402	OLB	C1-C2-C3-C4
5	A	402	OLB	O20-C21-C22-C24
5	A	406	OLB	C10-C11-C12-C13
5	A	403	OLB	C11-C12-C13-C14
5	A	403	OLB	C12-C13-C14-C15
5	A	402	OLB	C13-C14-C15-C16
6	A	407	OLA	C4-C5-C6-C7
5	A	403	OLB	C3-C4-C5-C6
5	A	405	OLB	C21-C22-C24-O25
5	A	403	OLB	O20-C1-C2-C3
5	A	402	OLB	O19-C1-O20-C21
5	A	404	OLB	C9-C10-C11-C12
5	A	406	OLB	C9-C10-C11-C12
5	A	404	OLB	O20-C21-C22-O23
5	A	405	OLB	C7-C8-C9-C10
5	A	402	OLB	C15-C16-C17-C18
5	A	406	OLB	C5-C6-C7-C8
5	A	406	OLB	C11-C12-C13-C14
5	A	403	OLB	C9-C10-C11-C12
5	A	404	OLB	C7-C8-C9-C10
5	A	405	OLB	O20-C21-C22-O23
5	A	404	OLB	C12-C13-C14-C15
5	A	405	OLB	O20-C21-C22-C24
6	A	407	OLA	C9-C10-C11-C12

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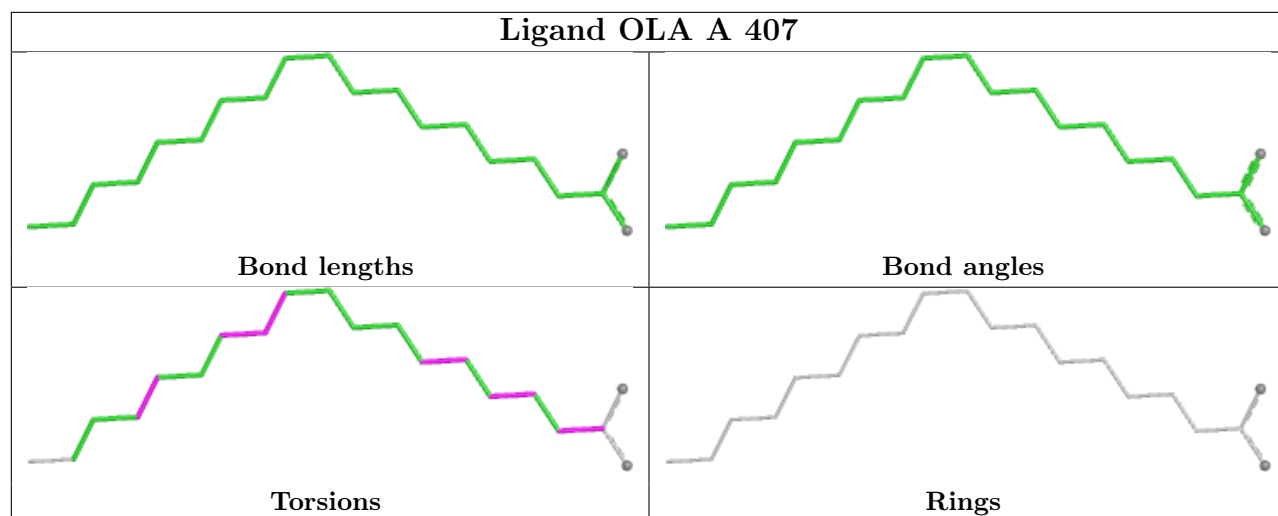
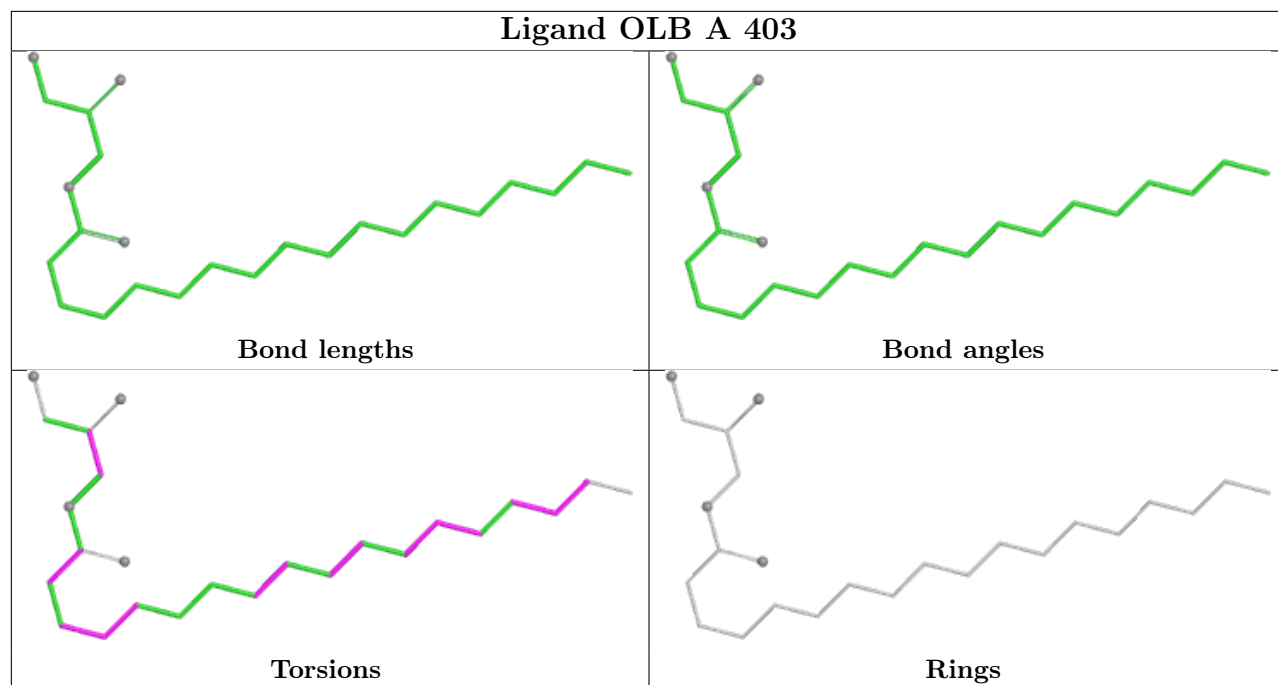
Mol	Chain	Res	Type	Atoms
5	A	402	OLB	C9-C10-C11-C12
5	A	403	OLB	C15-C16-C17-C18
5	A	404	OLB	C13-C14-C15-C16
6	A	407	OLA	O1-C1-C2-C3
6	A	407	OLA	O2-C1-C2-C3
5	A	403	OLB	C2-C3-C4-C5
5	A	404	OLB	C14-C15-C16-C17
5	A	403	OLB	C7-C8-C9-C10

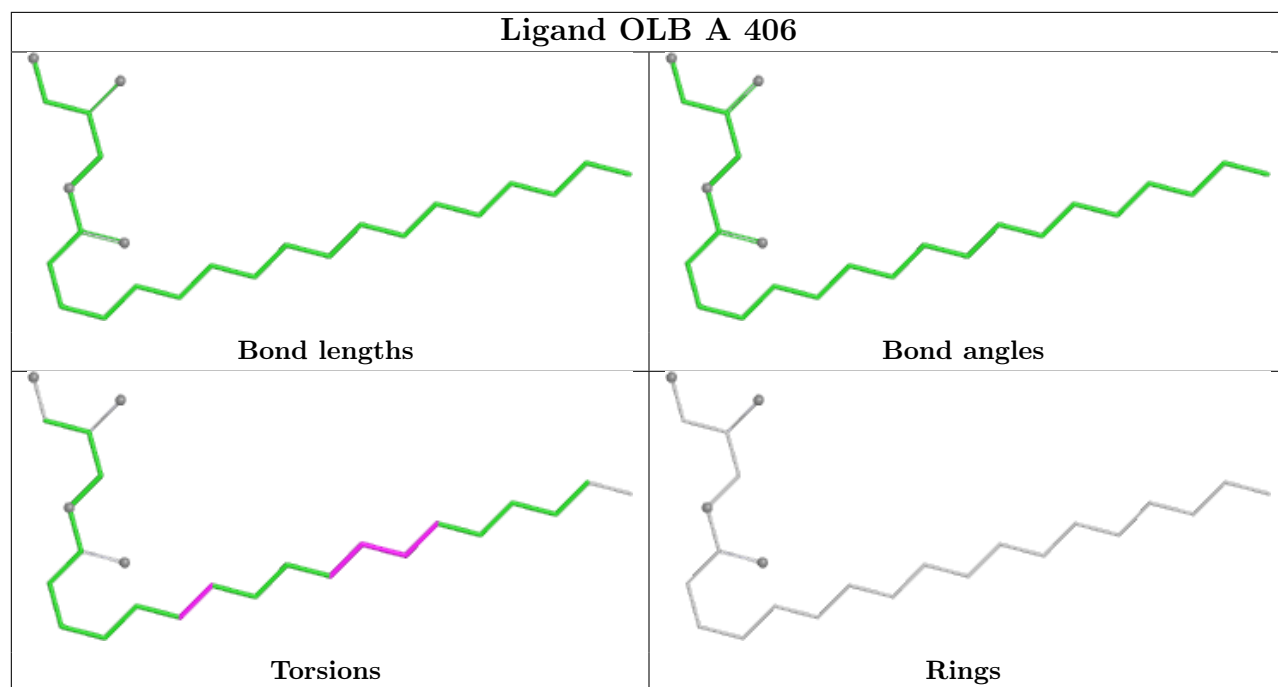
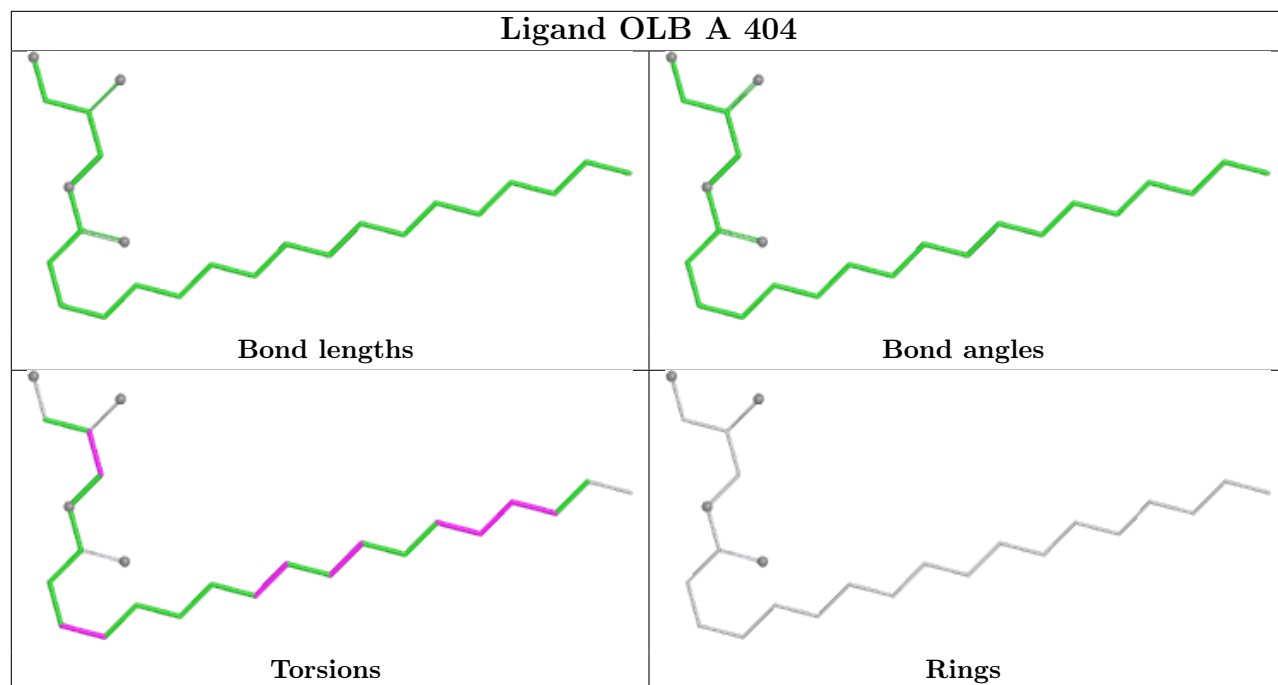
There are no ring outliers.

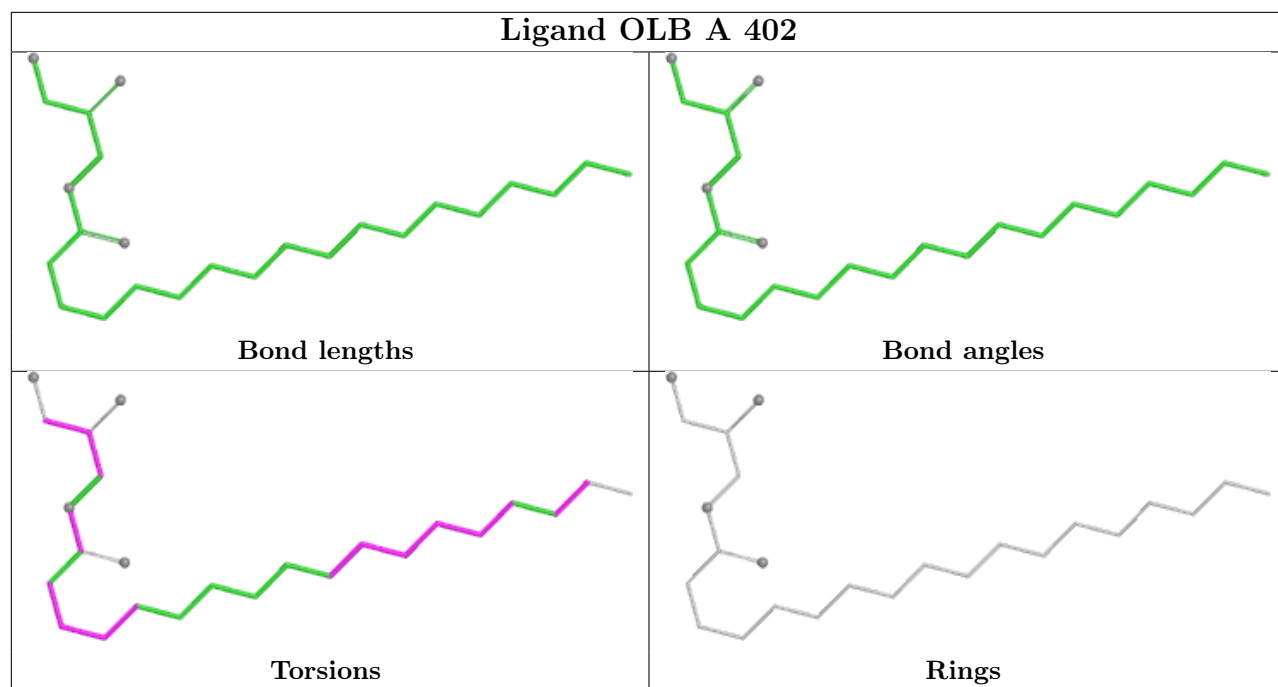
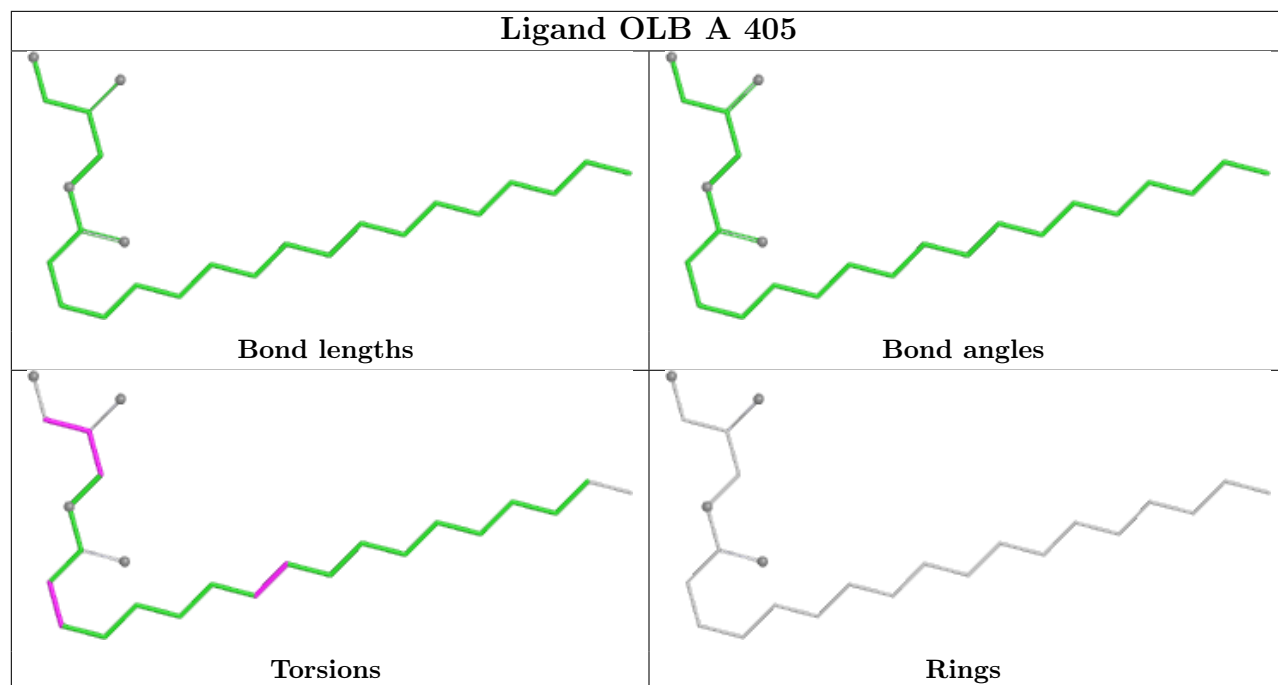
5 monomers are involved in 20 short contacts:

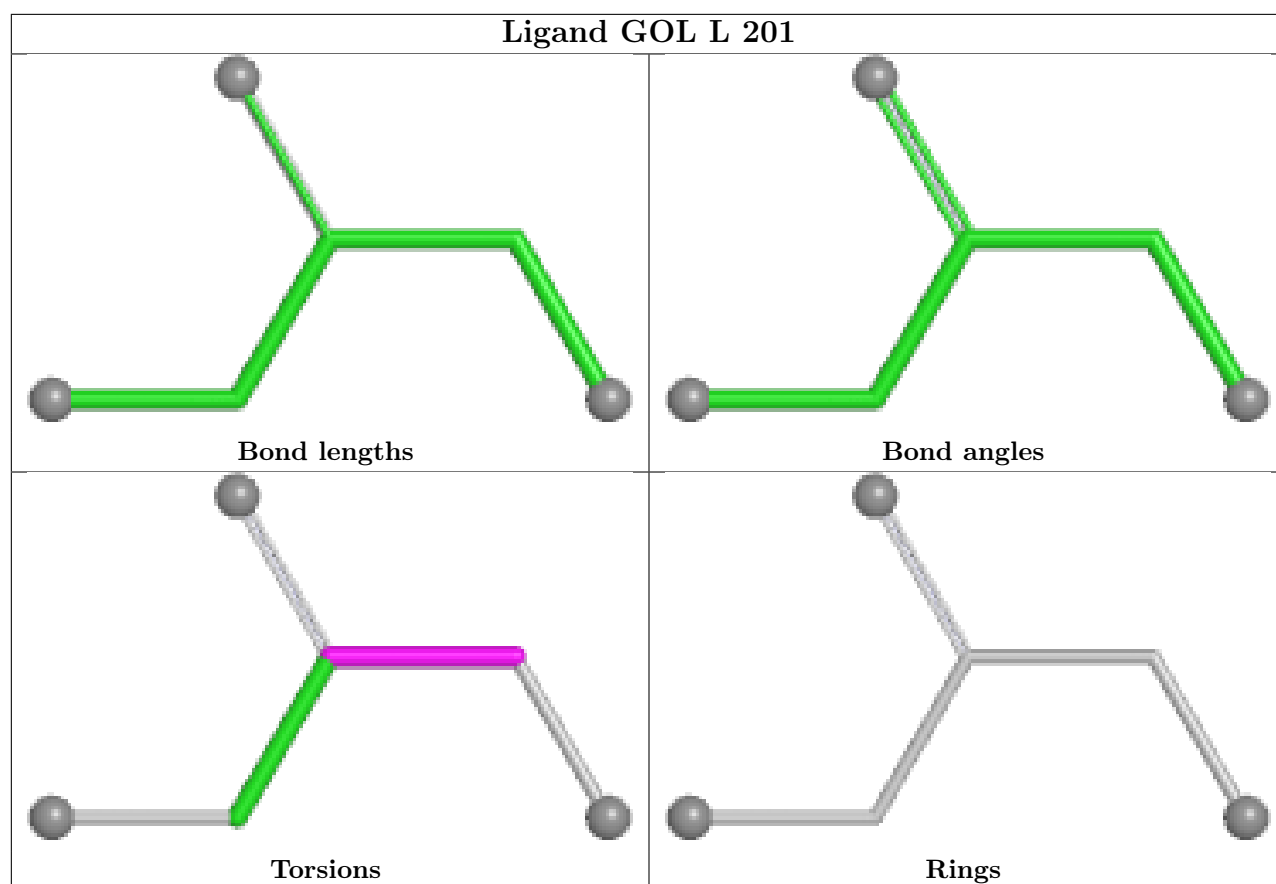
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	403	OLB	5	0
6	A	407	OLA	4	0
5	A	406	OLB	2	0
5	A	405	OLB	1	0
5	A	402	OLB	8	0

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









5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	285/292 (97%)	0.19	6 (2%) 63 59	63, 78, 111, 144	0
2	H	121/121 (100%)	-0.06	0 100 100	63, 77, 92, 97	0
3	L	108/108 (100%)	-0.14	1 (0%) 81 78	60, 76, 93, 100	0
All	All	514/521 (98%)	0.06	7 (1%) 73 69	60, 77, 104, 144	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	313	TRP	4.6
1	A	308	ILE	3.5
1	A	0	PHE	3.2
3	L	0	SER	3.0
1	A	117	ASP	2.9
1	A	168	PHE	2.4
1	A	379	GLY	2.2

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,

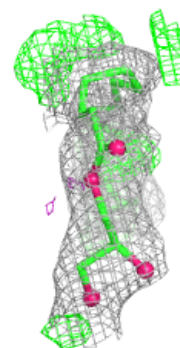
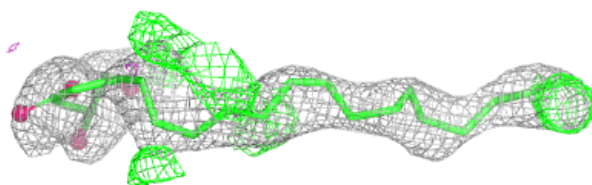
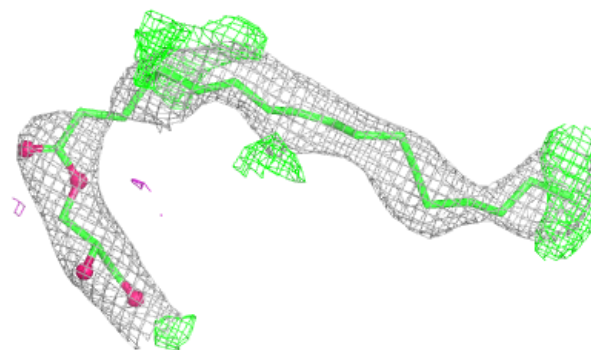
median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	OLB	A	403	25/25	0.69	0.20	97,101,103,103	0
5	OLB	A	406	25/25	0.76	0.26	110,122,132,133	0
5	OLB	A	404	25/25	0.79	0.19	107,109,116,116	0
5	OLB	A	405	25/25	0.80	0.20	97,102,114,114	0
5	OLB	A	402	25/25	0.80	0.18	90,94,102,102	0
7	GOL	L	201	6/6	0.84	0.16	113,114,114,114	0
6	OLA	A	407	20/20	0.91	0.14	70,74,81,82	0
4	ZN	A	401	1/1	0.98	0.08	72,72,72,72	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.

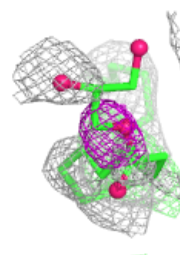
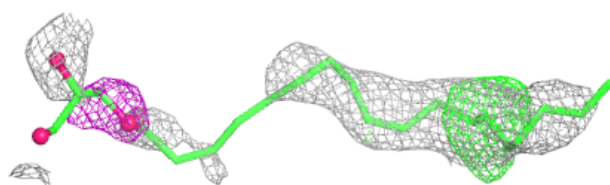
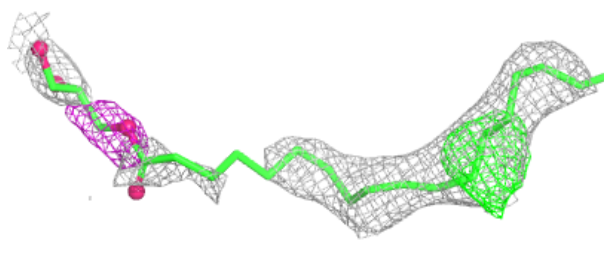
Electron density around OLB A 403:

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

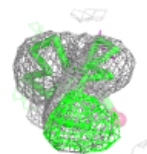
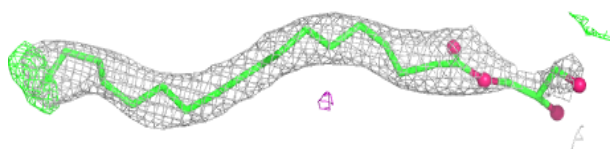
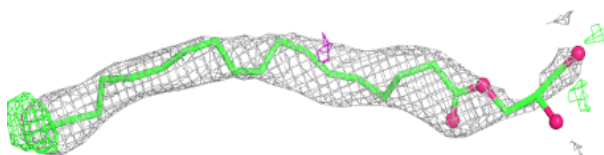


Electron density around OLB A 406:

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

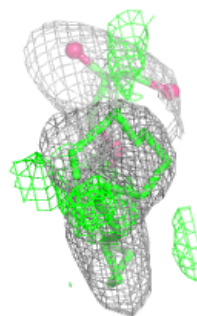
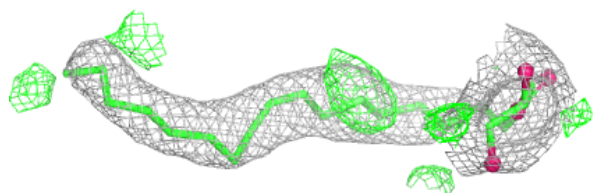
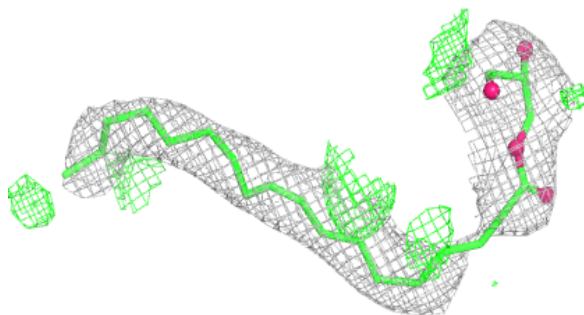
**Electron density around OLB A 404:**

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

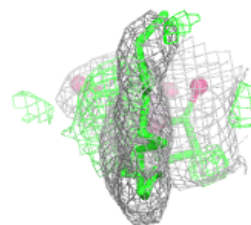
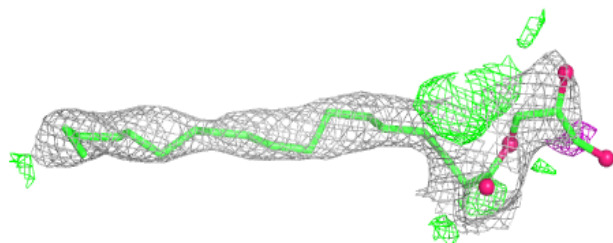
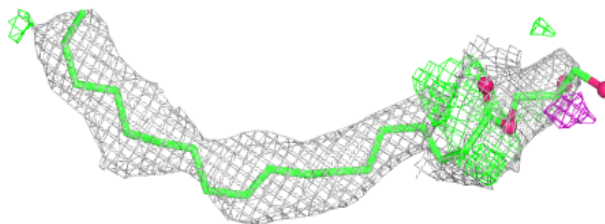


Electron density around OLB A 405:

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

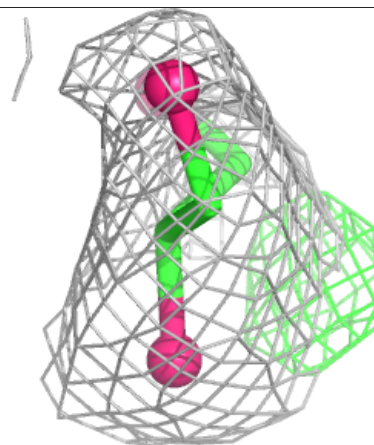
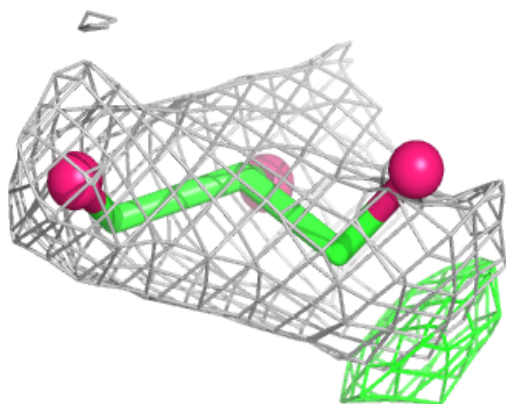
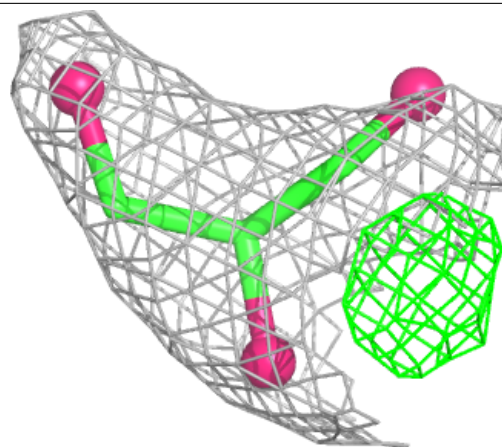
**Electron density around OLB A 402:**

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



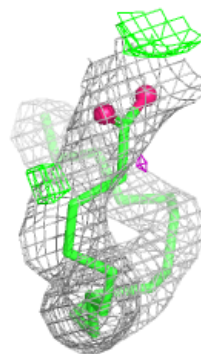
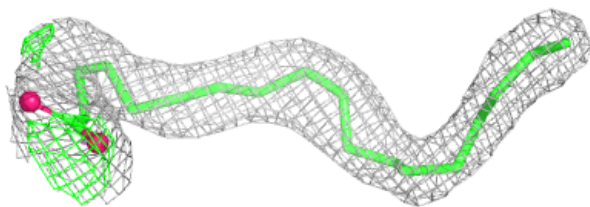
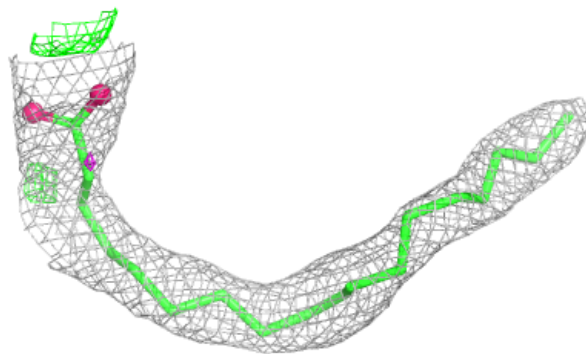
Electron density around GOL L 201:

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



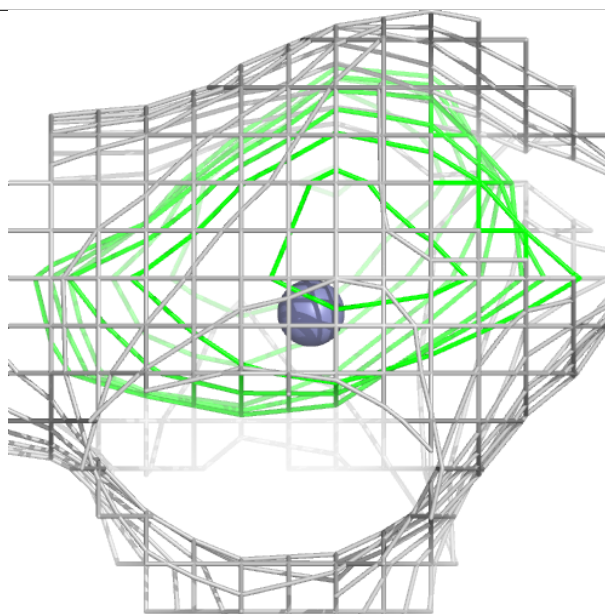
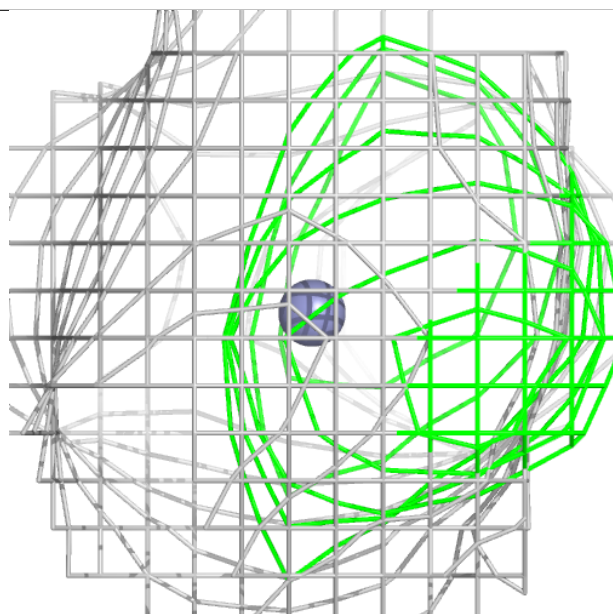
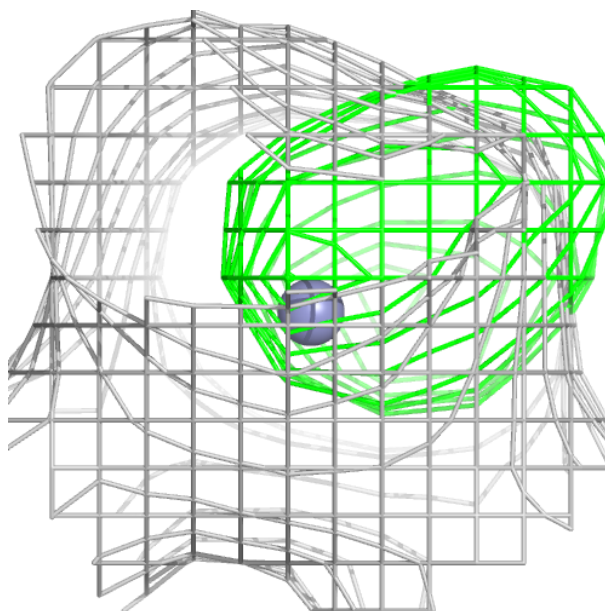
Electron density around OLA A 407:

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



Electron density around ZN A 401:

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



6.5 Other polymers ⓘ

There are no such residues in this entry.