



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 12:23 PM UTC

PDB ID : 8SJX / pdb_00008sjx
Title : Structure of lens aquaporin-0 array in sphingomyelin/cholesterol bilayer (2SM:1Chol)
Authors : Chiu, P.-L.; Walz, T.
Deposited on : 2023-04-18
Resolution : 2.50 Å (reported)
Based on initial model : 2B6O

This is a Full wwPDB EM Validation Report for a publicly released PDB/EMDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
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)
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

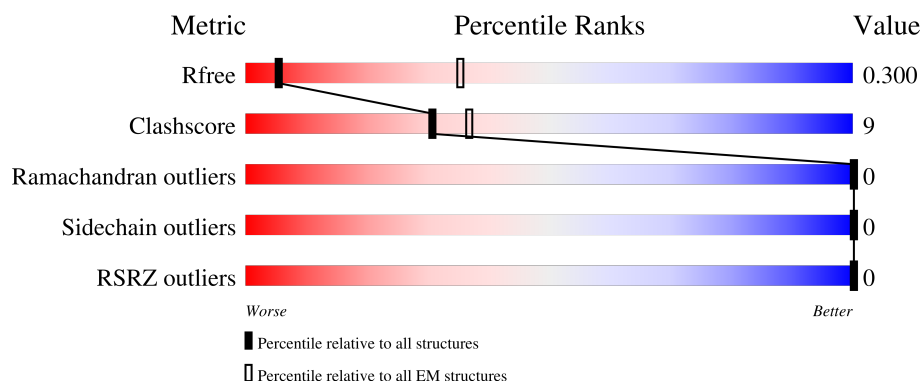
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON CRYSTALLOGRAPHY

The reported resolution of this entry is 2.50 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)	EM structures (#Entries)
R_{free}	180332	208
Clashscore	229148	23984
Ramachandran outliers	224038	23583
Sidechain outliers	223484	23102
RSRZ outliers	180361	209

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the 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 $\leq 5\%$.

Mol	Chain	Length	Quality of chain
1	A	263	 67% 16% 16%

2 Entry composition ⓘ

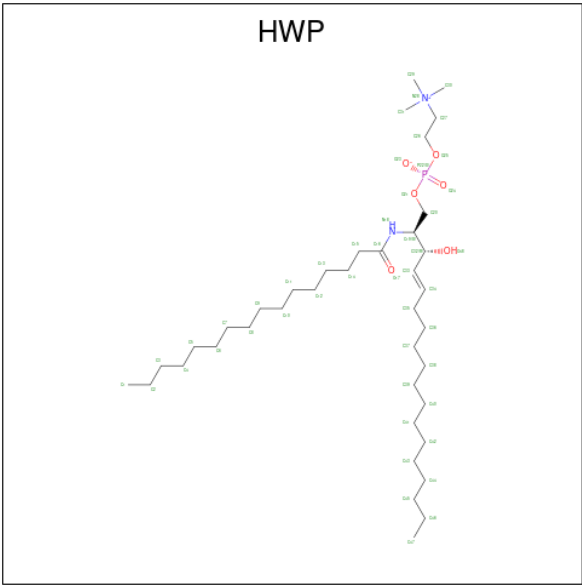
There are 4 unique types of molecules in this entry. The entry contains 2010 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, 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 Lens fiber major intrinsic protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	220	1663	1107	276	275	5	0	0

- Molecule 2 is [(E,2S,3R)-2-(hexadecanoylamino)-3-oxidanyl-octadec-4-enyl] 2-(trimethylazaniumyl)ethyl phosphate (CCD ID: HWP) (formula: C₃₉H₇₉N₂O₆P) (labeled as "Ligand of Interest" by depositor).



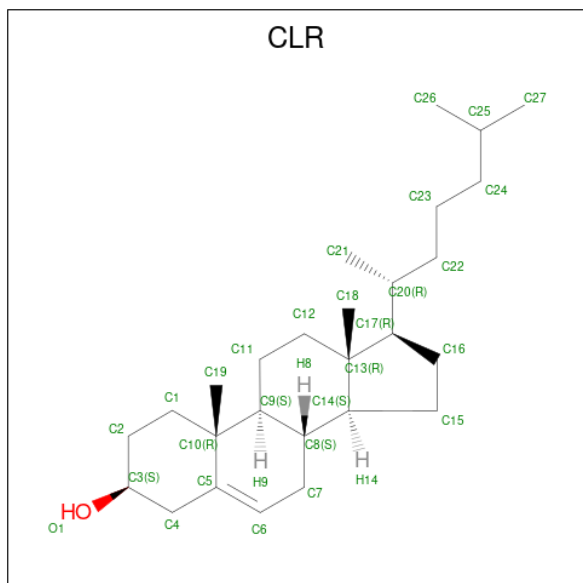
Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
2	A	1	Total 43	34	2	6	1	0
2	A	1	Total 46	37	2	6	1	0
2	A	1	Total 48	39	2	6	1	0
2	A	1	Total 45	36	2	6	1	0
2	A	1	Total 43	34	2	6	1	0

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Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	P	0
			41	32	2	6	1	
2	A	1	Total	C	N	O	P	0
			42	33	2	6	1	

- Molecule 3 is CHOLESTEROL (CCD ID: CLR) (formula: $C_{27}H_{46}O$) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		AltConf
4	A	11	Total	O	0
			11	11	

- Molecule 1: Lens fiber major intrinsic protein



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 2 2	Depositor
Cell constants a, b, c, α , β , γ	65.50Å 65.50Å 200.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	2.50 – 2.50 2.50 – 2.50	Depositor EDS
% Data completeness (in resolution range)	86.7 (2.50-2.50) 0.0 (2.50-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.35 (at 2.29Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.260 , 0.286 0.273 , 0.300	Depositor DCC
R_{free} test set	1752 reflections (8.89%)	wwPDB-VP
Wilson B-factor (Å ²)	47.5	Xtriage
Anisotropy	0.306	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 100.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.00	EDS
Total number of atoms	2010	wwPDB-VP
Average B, all atoms (Å ²)	66.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 13.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: CLR, HWP

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	0.18	0/1713	0.35	0/2344

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1663	0	1679	29	0
2	A	308	0	0	2	0
3	A	28	0	46	3	0
4	A	11	0	0	0	0
All	All	2010	0	1725	33	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 9.

All (33) 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:87:ILE:HG23	3:A:308:CLR:H25	1.70	0.74
1:A:137:LEU:HD11	1:A:175:GLY:HA3	1.78	0.64
1:A:84:LEU:HA	1:A:87:ILE:HD12	1.83	0.60
2:A:302:HWP:O21	2:A:302:HWP:O48	2.18	0.60
1:A:134:GLU:OE2	1:A:179:THR:OG1	2.25	0.53
1:A:68:ASN:HD21	1:A:184:ASN:HA	1.76	0.50
1:A:137:LEU:HD21	1:A:174:PHE:HD2	1.77	0.50
1:A:13:ILE:HD12	1:A:88:CYS:HB3	1.92	0.50
1:A:28:LEU:HG	1:A:117:ALA:HB3	1.94	0.49
1:A:171:GLY:HA3	1:A:183:MET:HE1	1.95	0.48
1:A:38:PRO:HG2	1:A:40:HIS:HD2	1.79	0.48
1:A:68:ASN:ND2	1:A:184:ASN:HA	2.30	0.47
1:A:41:VAL:HG21	1:A:177:TYR:CZ	2.49	0.47
2:A:301:HWP:O21	2:A:301:HWP:O48	2.31	0.47
3:A:308:CLR:H232	3:A:308:CLR:H211	1.71	0.46
1:A:70:ALA:HB1	1:A:212:ALA:HB2	1.97	0.46
1:A:184:ASN:HB3	1:A:187:ARG:HB3	1.98	0.46
1:A:142:VAL:HG21	1:A:212:ALA:HA	1.98	0.46
1:A:148:THR:HB	1:A:160:VAL:HB	1.97	0.46
1:A:119:ASN:HB3	1:A:204:TYR:CE1	2.51	0.45
1:A:110:PRO:HA	1:A:113:ARG:HB2	1.98	0.45
1:A:12:ALA:HA	1:A:62:ILE:HG22	1.99	0.45
1:A:220:ASP:O	1:A:225:PRO:HA	2.17	0.45
1:A:24:VAL:HG13	1:A:28:LEU:HD22	1.97	0.44
1:A:24:VAL:O	1:A:28:LEU:HB2	2.17	0.44
1:A:15:ALA:HB1	1:A:59:VAL:HB	2.00	0.43
1:A:141:PHE:CZ	1:A:145:ILE:HD11	2.52	0.43
1:A:97:ALA:HB1	1:A:190:ALA:HB2	1.99	0.43
3:A:308:CLR:H262	3:A:308:CLR:H231	1.76	0.43
1:A:35:ALA:HB3	1:A:38:PRO:HG3	2.01	0.43
1:A:68:ASN:ND2	1:A:185:PRO:HD2	2.36	0.41
1:A:9:PHE:O	1:A:13:ILE:HG12	2.20	0.41
1:A:119:ASN:HB3	1:A:204:TYR:CZ	2.56	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 PDB entries followed by that with respect to all EM

entries.

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	218/263 (83%)	208 (95%)	10 (5%)	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 PDB entries followed by that with respect to all EM entries.

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	166/205 (81%)	166 (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 (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	68	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	HWP	A	301	-	41,42,47	1.40	5 (12%)	47,50,55	1.39	6 (12%)
2	HWP	A	303	-	46,47,47	1.28	5 (10%)	52,55,55	1.13	1 (1%)
2	HWP	A	304	-	43,44,47	1.20	6 (13%)	49,52,55	1.47	2 (4%)
2	HWP	A	307	-	40,41,47	1.48	8 (20%)	46,49,55	1.07	2 (4%)
2	HWP	A	305	-	41,42,47	1.47	7 (17%)	47,50,55	1.58	8 (17%)
2	HWP	A	306	-	39,40,47	1.46	6 (15%)	45,48,55	1.24	5 (11%)
2	HWP	A	302	-	44,45,47	1.43	5 (11%)	50,53,55	1.07	2 (4%)
3	CLR	A	308	-	31,31,31	0.41	0	48,48,48	0.78	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HWP	A	301	-	-	26/47/47/52	-
2	HWP	A	303	-	-	18/52/52/52	-
2	HWP	A	304	-	-	11/49/49/52	-
2	HWP	A	307	-	-	14/46/46/52	-
2	HWP	A	305	-	-	22/47/47/52	-
2	HWP	A	306	-	-	14/45/45/52	-
2	HWP	A	302	-	-	19/50/50/52	-
3	CLR	A	308	-	-	9/10/68/68	0/4/4/4

All (42) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	302	HWP	C20-C19	4.95	1.60	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	305	HWP	P22-O25	4.56	1.77	1.59
2	A	306	HWP	P22-O25	4.56	1.77	1.59
2	A	303	HWP	P22-O25	4.51	1.77	1.59
2	A	302	HWP	P22-O25	4.46	1.76	1.59
2	A	301	HWP	C20-C19	4.46	1.59	1.51
2	A	303	HWP	C20-C19	4.46	1.59	1.51
2	A	301	HWP	P22-O25	4.42	1.76	1.59
2	A	307	HWP	P22-O25	4.39	1.76	1.59
2	A	304	HWP	P22-O25	4.29	1.76	1.59
2	A	306	HWP	C20-C19	4.09	1.59	1.51
2	A	305	HWP	C20-C19	3.83	1.58	1.51
2	A	307	HWP	C20-C19	3.47	1.58	1.51
2	A	307	HWP	C16-N18	3.36	1.41	1.34
2	A	301	HWP	P22-O21	3.15	1.71	1.59
2	A	303	HWP	P22-O21	3.13	1.71	1.59
2	A	305	HWP	C19-N18	3.12	1.51	1.46
2	A	307	HWP	C19-N18	3.10	1.51	1.46
2	A	302	HWP	P22-O21	3.09	1.71	1.59
2	A	306	HWP	P22-O21	3.02	1.71	1.59
2	A	305	HWP	P22-O21	2.95	1.70	1.59
2	A	304	HWP	P22-O21	2.87	1.70	1.59
2	A	306	HWP	C16-N18	2.76	1.39	1.34
2	A	301	HWP	C16-N18	2.70	1.39	1.34
2	A	304	HWP	C20-C19	2.68	1.56	1.51
2	A	305	HWP	C16-N18	2.65	1.39	1.34
2	A	307	HWP	P22-O21	2.59	1.69	1.59
2	A	307	HWP	O25-C26	-2.56	1.34	1.44
2	A	302	HWP	C16-N18	2.43	1.39	1.34
2	A	305	HWP	O25-C26	-2.43	1.35	1.44
2	A	306	HWP	C19-N18	2.42	1.50	1.46
2	A	303	HWP	O25-C26	-2.39	1.35	1.44
2	A	301	HWP	O25-C26	-2.39	1.35	1.44
2	A	304	HWP	O25-C26	-2.35	1.35	1.44
2	A	302	HWP	O25-C26	-2.34	1.35	1.44
2	A	306	HWP	O25-C26	-2.33	1.35	1.44
2	A	303	HWP	C16-N18	2.22	1.38	1.34
2	A	305	HWP	O48-C32	-2.08	1.39	1.43
2	A	307	HWP	O21-C20	-2.08	1.36	1.44
2	A	304	HWP	O48-C32	-2.05	1.39	1.43
2	A	307	HWP	C15-C16	2.04	1.55	1.51
2	A	304	HWP	O21-C20	-2.01	1.37	1.44

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	304	HWP	C32-C33-C34	-7.22	109.73	124.69
2	A	305	HWP	C11-C10-C9	-5.15	88.35	114.37
2	A	305	HWP	C13-C14-C15	-4.43	96.86	113.13
2	A	301	HWP	C32-C33-C34	-4.30	115.78	124.69
2	A	306	HWP	C35-C34-C33	-3.42	110.23	125.47
2	A	303	HWP	C32-C33-C34	-3.31	117.84	124.69
2	A	306	HWP	C32-C33-C34	-3.06	118.36	124.69
2	A	306	HWP	C19-N18-C16	2.86	128.18	123.40
2	A	302	HWP	C32-C33-C34	-2.75	119.00	124.69
2	A	305	HWP	C11-C12-C13	-2.61	101.15	114.37
2	A	305	HWP	C19-N18-C16	2.60	127.74	123.40
2	A	305	HWP	C9-C8-C7	-2.59	101.26	114.37
2	A	301	HWP	C13-C14-C15	-2.58	103.66	113.13
2	A	306	HWP	C14-C15-C16	-2.54	106.13	113.19
2	A	302	HWP	C19-N18-C16	2.54	127.64	123.40
2	A	301	HWP	C20-C19-C32	-2.47	107.38	112.90
2	A	305	HWP	C32-C19-N18	2.42	113.76	109.66
2	A	305	HWP	C32-C33-C34	-2.41	119.69	124.69
2	A	301	HWP	C35-C34-C33	-2.27	115.34	125.47
2	A	301	HWP	C40-C39-C38	-2.26	102.94	114.37
2	A	305	HWP	C20-C19-C32	-2.26	107.86	112.90
2	A	301	HWP	C15-C16-N18	2.21	119.77	115.86
2	A	307	HWP	C32-C19-N18	2.21	113.41	109.66
2	A	306	HWP	C20-C19-C32	-2.19	108.02	112.90
2	A	304	HWP	C14-C15-C16	-2.18	107.15	113.19
2	A	307	HWP	C15-C16-N18	2.04	119.45	115.86

There are no chirality outliers.

All (133) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	HWP	C19-C20-O21-P22
2	A	301	HWP	O25-C26-C27-N28
2	A	301	HWP	C20-O21-P22-O23
2	A	301	HWP	C20-O21-P22-O25
2	A	301	HWP	C15-C16-N18-C19
2	A	301	HWP	O17-C16-N18-C19
2	A	301	HWP	C32-C19-N18-C16
2	A	301	HWP	C20-C19-C32-C33
2	A	301	HWP	C20-C19-C32-O48
2	A	301	HWP	N18-C19-C32-C33
2	A	301	HWP	N18-C19-C32-O48
2	A	302	HWP	N18-C19-C20-O21

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Mol	Chain	Res	Type	Atoms
2	A	302	HWP	C32-C19-C20-O21
2	A	302	HWP	C26-O25-P22-O21
2	A	302	HWP	C15-C16-N18-C19
2	A	302	HWP	C32-C19-N18-C16
2	A	302	HWP	C20-C19-C32-O48
2	A	302	HWP	N18-C19-C32-C33
2	A	302	HWP	C19-C32-C33-C34
2	A	302	HWP	O48-C32-C33-C34
2	A	303	HWP	C27-C26-O25-P22
2	A	303	HWP	C20-O21-P22-O23
2	A	303	HWP	C20-O21-P22-O25
2	A	303	HWP	C15-C16-N18-C19
2	A	303	HWP	O17-C16-N18-C19
2	A	303	HWP	C20-C19-C32-C33
2	A	303	HWP	C20-C19-C32-O48
2	A	303	HWP	N18-C19-C32-C33
2	A	303	HWP	N18-C19-C32-O48
2	A	303	HWP	C19-C32-C33-C34
2	A	304	HWP	C20-O21-P22-O23
2	A	304	HWP	C15-C16-N18-C19
2	A	304	HWP	O17-C16-N18-C19
2	A	304	HWP	C19-C32-C33-C34
2	A	305	HWP	C19-C20-O21-P22
2	A	305	HWP	C27-C26-O25-P22
2	A	305	HWP	C20-O21-P22-O25
2	A	305	HWP	C15-C16-N18-C19
2	A	305	HWP	O17-C16-N18-C19
2	A	305	HWP	C32-C19-N18-C16
2	A	305	HWP	C20-C19-C32-C33
2	A	305	HWP	C20-C19-C32-O48
2	A	305	HWP	N18-C19-C32-C33
2	A	305	HWP	N18-C19-C32-O48
2	A	306	HWP	N18-C19-C20-O21
2	A	306	HWP	C32-C19-C20-O21
2	A	306	HWP	O25-C26-C27-N28
2	A	306	HWP	C15-C16-N18-C19
2	A	306	HWP	O17-C16-N18-C19
2	A	306	HWP	C20-C19-N18-C16
2	A	306	HWP	C32-C19-N18-C16
2	A	306	HWP	N18-C19-C32-O48
2	A	306	HWP	C19-C32-C33-C34
2	A	307	HWP	N18-C19-C20-O21

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Mol	Chain	Res	Type	Atoms
2	A	307	HWP	C27-C26-O25-P22
2	A	307	HWP	C20-C19-C32-C33
2	A	307	HWP	C20-C19-C32-O48
2	A	307	HWP	N18-C19-C32-C33
2	A	307	HWP	N18-C19-C32-O48
2	A	307	HWP	C19-C32-C33-C34
2	A	307	HWP	O48-C32-C33-C34
2	A	302	HWP	O17-C16-N18-C19
2	A	303	HWP	C14-C15-C16-N18
2	A	303	HWP	C14-C15-C16-O17
2	A	307	HWP	C14-C15-C16-O17
2	A	304	HWP	C14-C15-C16-N18
2	A	307	HWP	C14-C15-C16-N18
3	A	308	CLR	C13-C17-C20-C22
2	A	304	HWP	C14-C15-C16-O17
2	A	303	HWP	O48-C32-C33-C34
2	A	306	HWP	O48-C32-C33-C34
2	A	305	HWP	C11-C12-C13-C14
3	A	308	CLR	C20-C22-C23-C24
2	A	305	HWP	C11-C10-C9-C8
2	A	305	HWP	C10-C11-C12-C13
3	A	308	CLR	C22-C23-C24-C25
2	A	301	HWP	C10-C11-C12-C13
3	A	308	CLR	C13-C17-C20-C21
3	A	308	CLR	C16-C17-C20-C21
2	A	306	HWP	C33-C34-C35-C36
2	A	306	HWP	C13-C14-C15-C16
2	A	305	HWP	C12-C13-C14-C15
2	A	304	HWP	O48-C32-C33-C34
2	A	307	HWP	C34-C35-C36-C37
2	A	305	HWP	C7-C8-C9-C10
3	A	308	CLR	C16-C17-C20-C22
2	A	301	HWP	C12-C13-C14-C15
2	A	307	HWP	C36-C37-C38-C39
2	A	301	HWP	C6-C7-C8-C9
2	A	302	HWP	N18-C19-C32-O48
2	A	301	HWP	C9-C10-C11-C12
3	A	308	CLR	C21-C20-C22-C23
2	A	301	HWP	C11-C10-C9-C8
2	A	301	HWP	C27-C26-O25-P22
2	A	302	HWP	C20-C19-C32-C33
2	A	306	HWP	C20-C19-C32-O48

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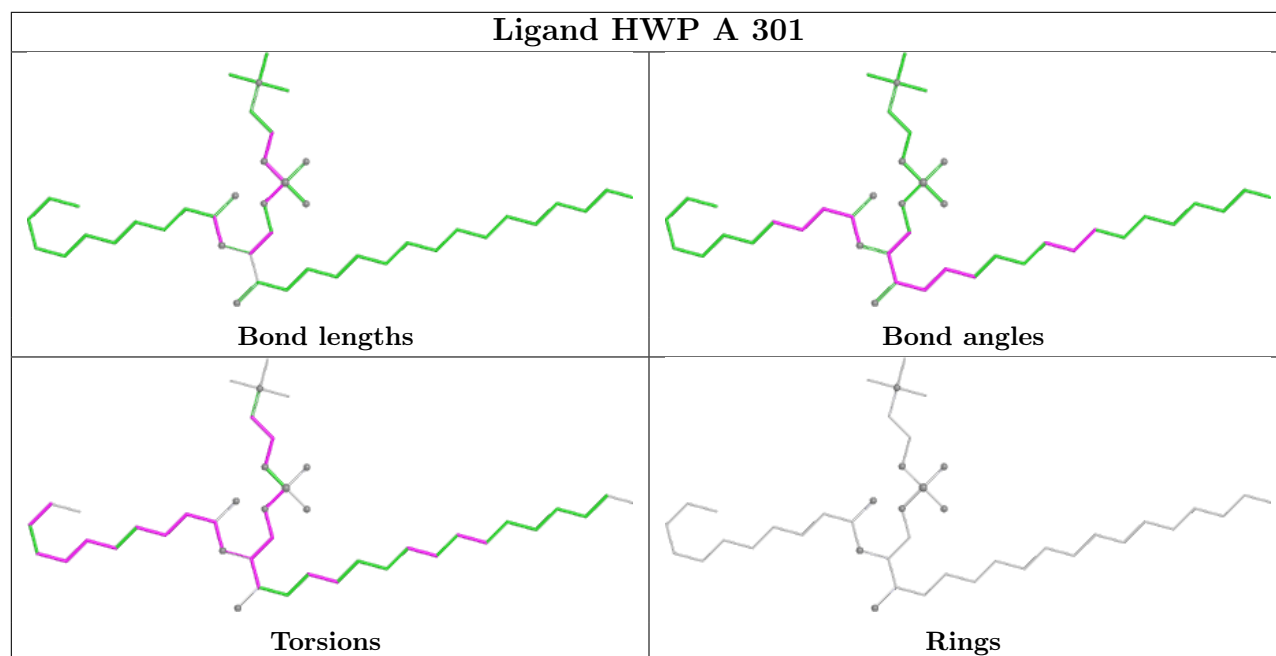
Mol	Chain	Res	Type	Atoms
2	A	301	HWP	N18-C19-C20-O21
2	A	302	HWP	O25-C26-C27-N28
2	A	303	HWP	O25-C26-C27-N28
2	A	304	HWP	O25-C26-C27-N28
2	A	305	HWP	O25-C26-C27-N28
2	A	302	HWP	C19-C20-O21-P22
2	A	304	HWP	C19-C20-O21-P22
2	A	301	HWP	C14-C15-C16-O17
2	A	302	HWP	C14-C15-C16-O17
2	A	307	HWP	C37-C38-C39-C40
2	A	301	HWP	C32-C19-C20-O21
2	A	301	HWP	C20-O21-P22-O24
2	A	302	HWP	C20-O21-P22-O24
2	A	302	HWP	C26-O25-P22-O24
2	A	303	HWP	C26-O25-P22-O24
2	A	304	HWP	C26-O25-P22-O21
2	A	305	HWP	C20-O21-P22-O24
2	A	305	HWP	C26-O25-P22-O24
2	A	301	HWP	C33-C34-C35-C36
3	A	308	CLR	C23-C24-C25-C26
2	A	301	HWP	C14-C15-C16-N18
2	A	304	HWP	C33-C34-C35-C36
2	A	302	HWP	C14-C15-C16-N18
3	A	308	CLR	C23-C24-C25-C27
2	A	302	HWP	C41-C42-C43-C44
2	A	301	HWP	C13-C14-C15-C16
2	A	305	HWP	C13-C14-C15-C16
2	A	305	HWP	C37-C38-C39-C40
2	A	305	HWP	C33-C34-C35-C36
2	A	305	HWP	C5-C6-C7-C8
2	A	303	HWP	C36-C37-C38-C39
2	A	301	HWP	C37-C38-C39-C40
2	A	306	HWP	C20-C19-C32-C33
2	A	301	HWP	C39-C40-C41-C42
2	A	303	HWP	C19-C20-O21-P22
2	A	303	HWP	C11-C10-C9-C8
2	A	307	HWP	C32-C19-C20-O21

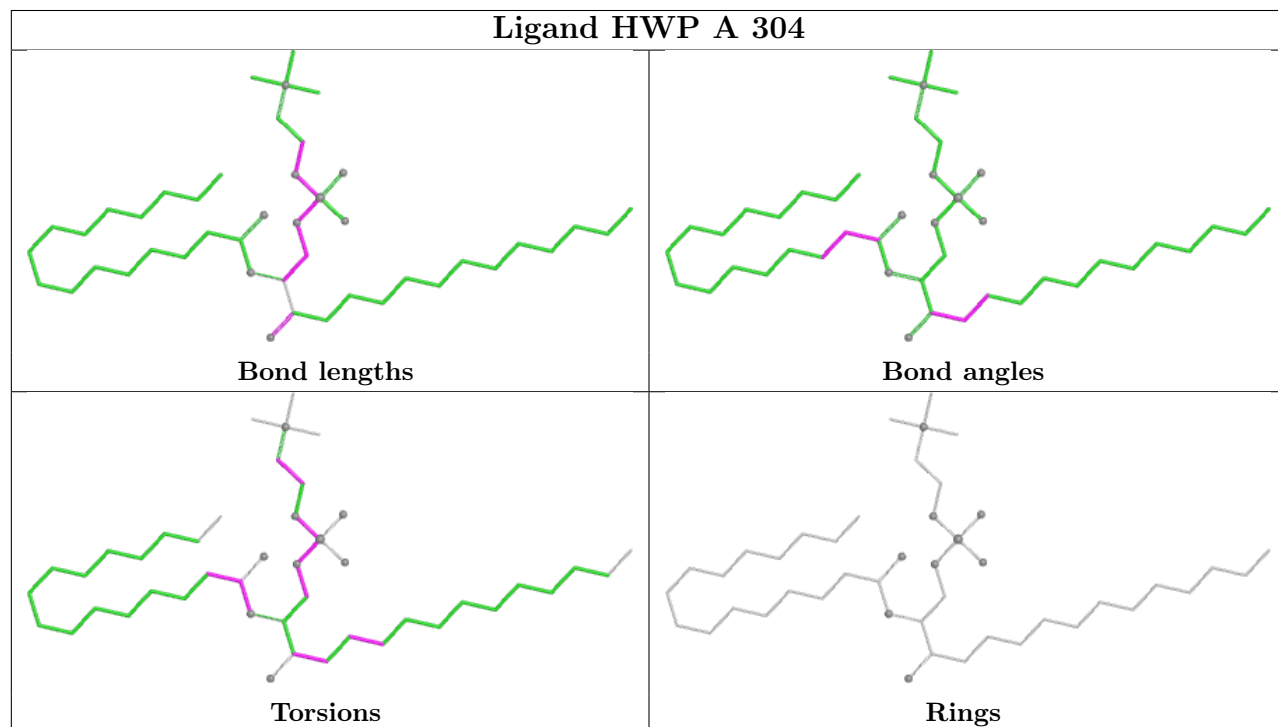
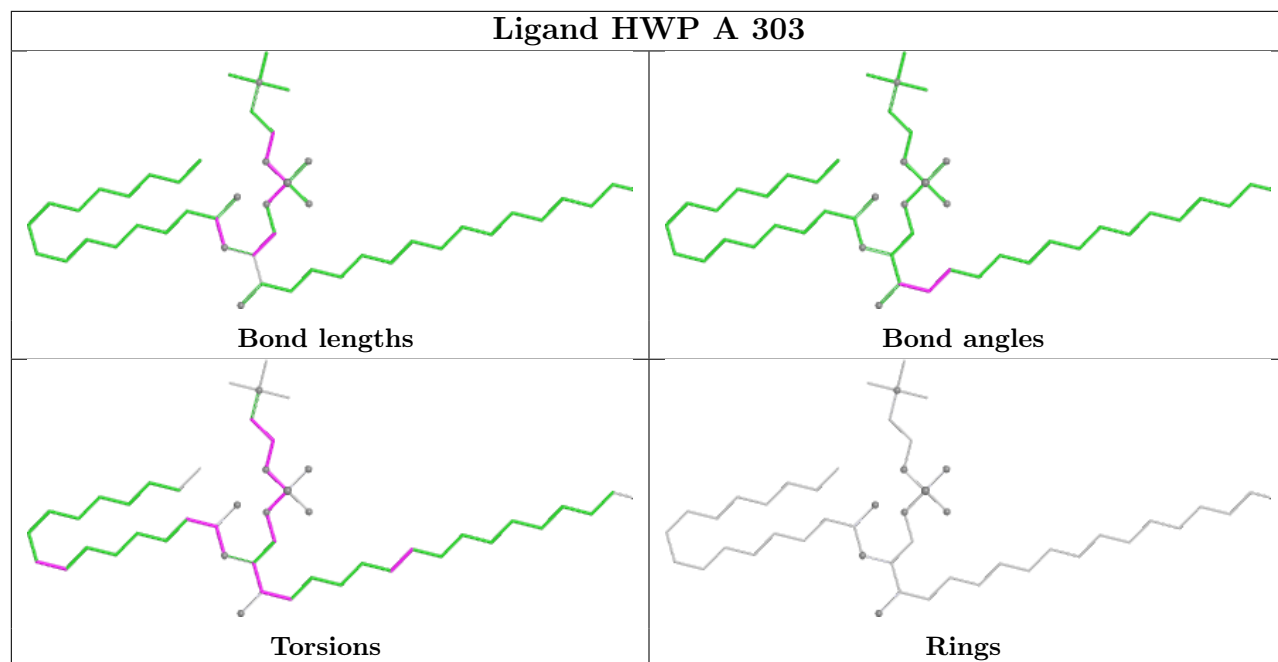
There are no ring outliers.

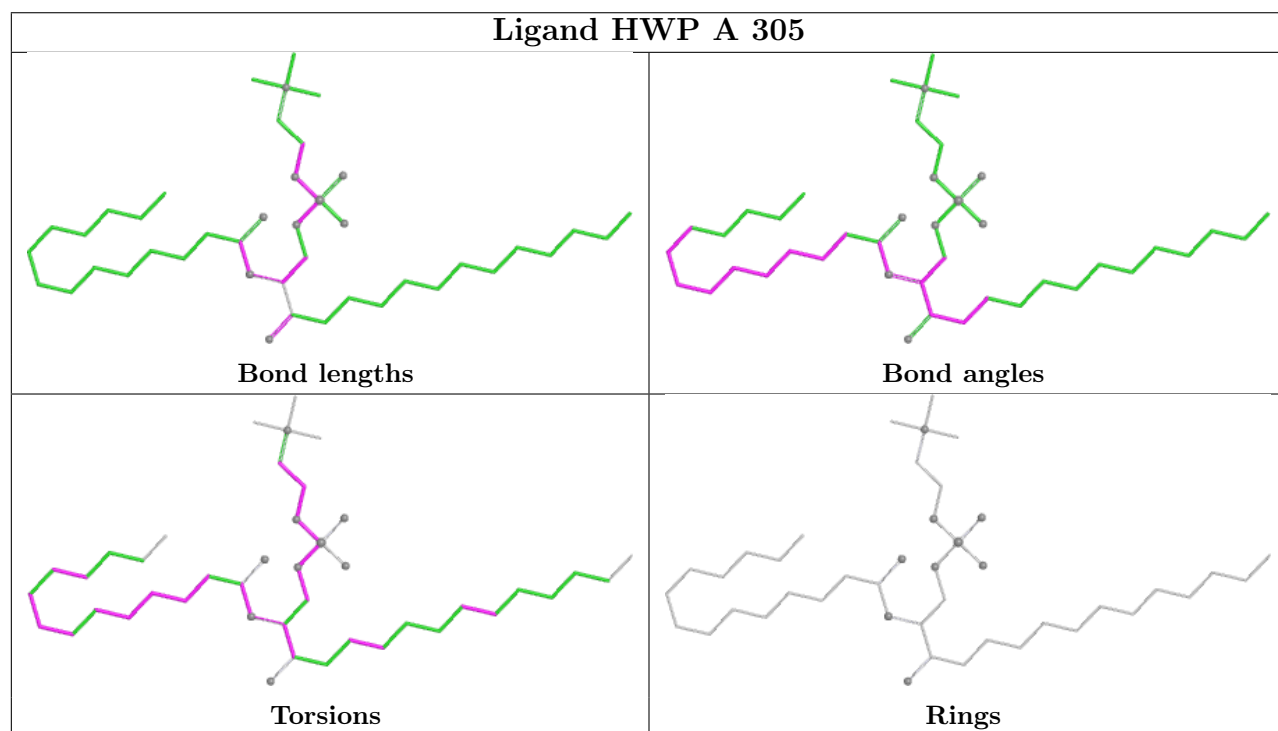
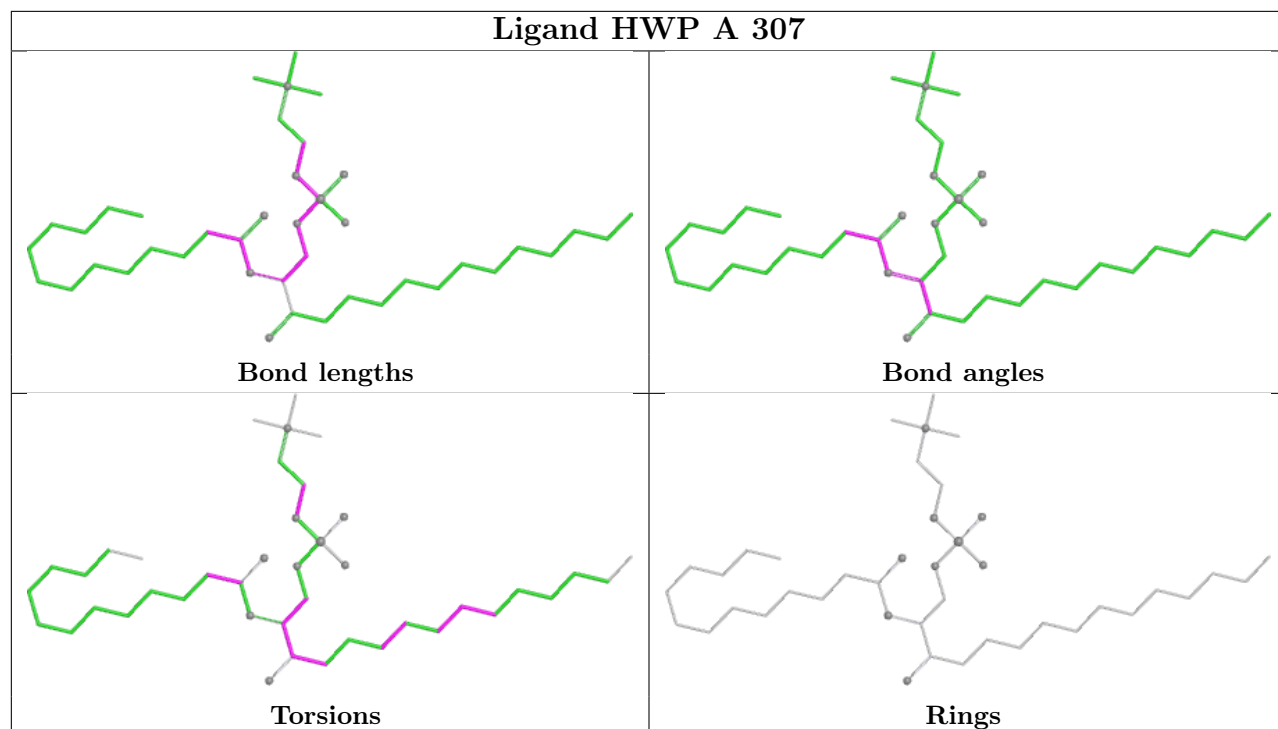
3 monomers are involved in 5 short contacts:

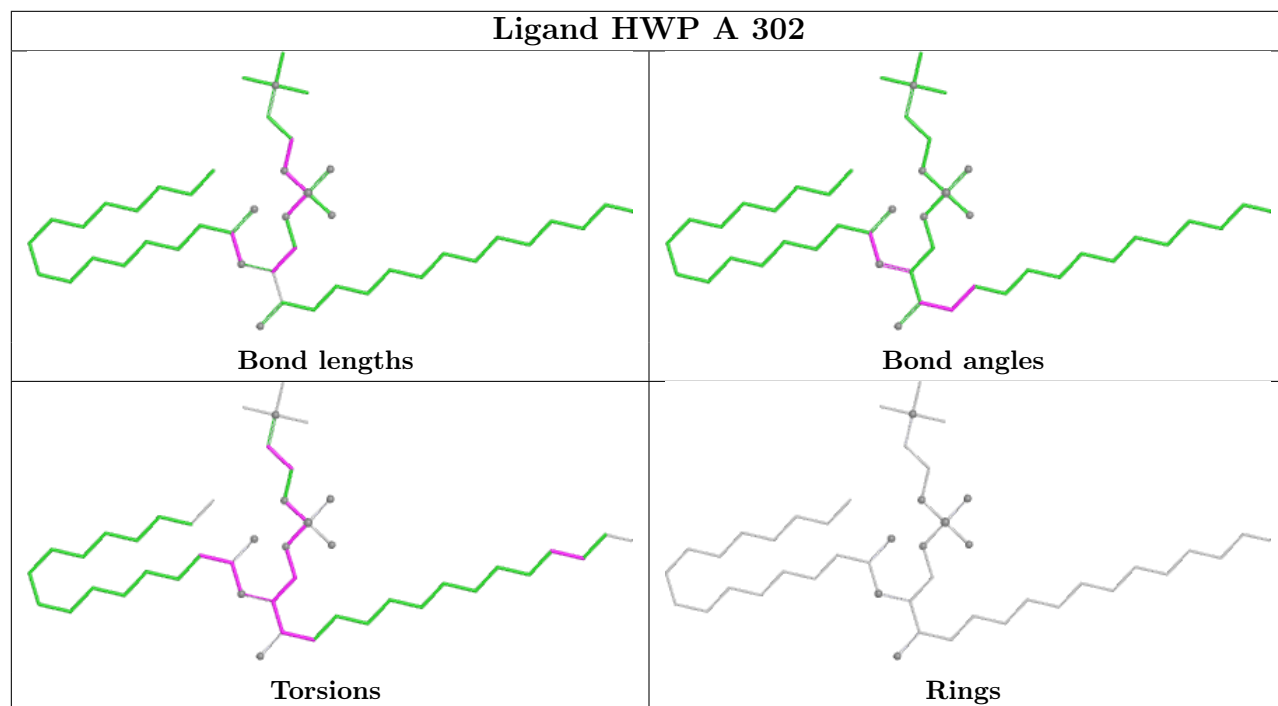
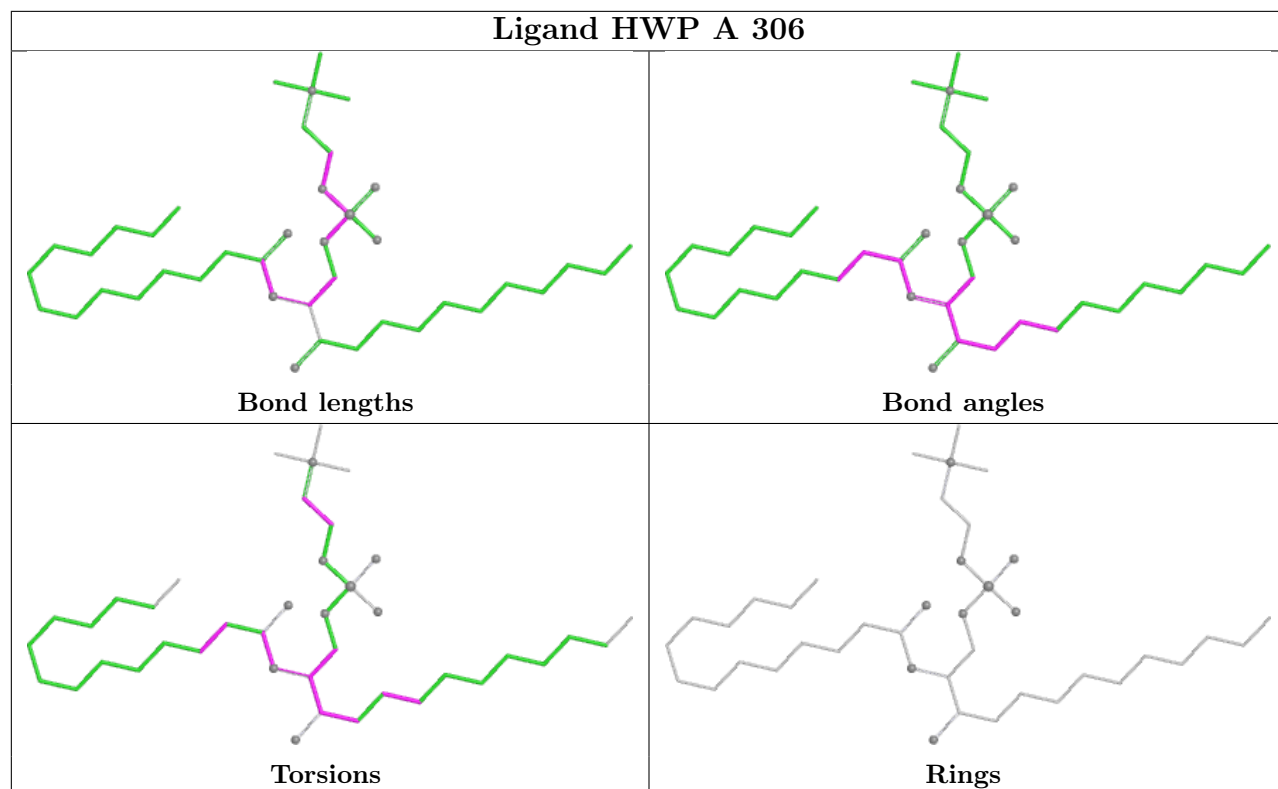
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	301	HWP	1	0
2	A	302	HWP	1	0
3	A	308	CLR	3	0

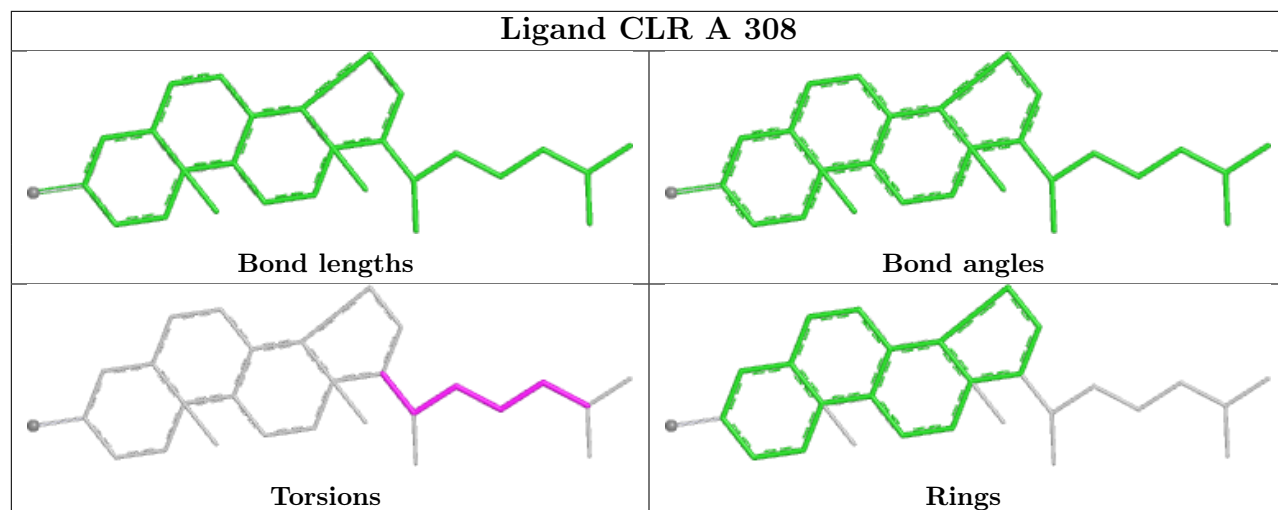
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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











5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.