



Full wwPDB EM Validation Report ⓘ

Jul 24, 2025 – 12:12 PM JST

PDB ID : 9JOU / pdb_00009jou
EMDB ID : EMD-61685
Title : Cryo-EM structure of the zeaxanthin-bound light-driven proton pumping rhodopsin, NM-R1
Authors : Hosaka, T.; Shirouzu, M.
Deposited on : 2024-09-25
Resolution : 2.46 Å(reported)

This is a Full wwPDB EM 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/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:

EMDB validation analysis : 0.0.1.dev118
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0rc1
buster-report : 1.1.7 (2018)
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.44

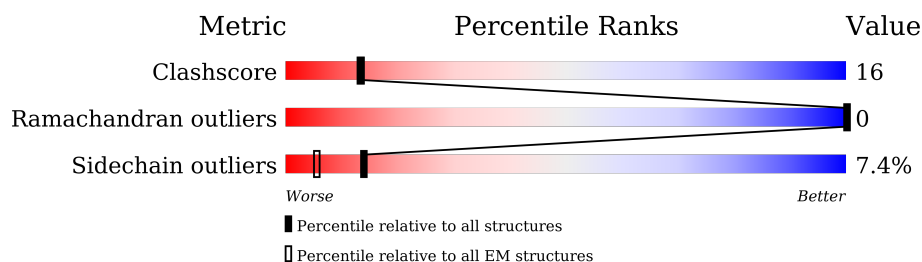
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.46 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)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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 $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	260	
1	B	260	
1	C	260	
1	D	260	
1	E	260	

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
2	RET	E	302	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 9913 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 Proteorhodopsin.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	230	Total	C	N	O	S	0	0
			1840	1247	283	298	12		
1	B	230	Total	C	N	O	S	0	0
			1840	1247	283	298	12		
1	C	230	Total	C	N	O	S	0	0
			1840	1247	283	298	12		
1	D	230	Total	C	N	O	S	0	0
			1840	1247	283	298	12		
1	E	230	Total	C	N	O	S	0	0
			1840	1247	283	298	12		

There are 70 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	247	LEU	-	expression tag	UNP W8VZ92
A	248	GLU	-	expression tag	UNP W8VZ92
A	249	GLU	-	expression tag	UNP W8VZ92
A	250	ASN	-	expression tag	UNP W8VZ92
A	251	LEU	-	expression tag	UNP W8VZ92
A	252	TYR	-	expression tag	UNP W8VZ92
A	253	PHE	-	expression tag	UNP W8VZ92
A	254	GLN	-	expression tag	UNP W8VZ92
A	255	GLY	-	expression tag	UNP W8VZ92
A	256	TRP	-	expression tag	UNP W8VZ92
A	257	SER	-	expression tag	UNP W8VZ92
A	258	HIS	-	expression tag	UNP W8VZ92
A	259	PRO	-	expression tag	UNP W8VZ92
A	260	GLN	-	expression tag	UNP W8VZ92
B	247	LEU	-	expression tag	UNP W8VZ92
B	248	GLU	-	expression tag	UNP W8VZ92
B	249	GLU	-	expression tag	UNP W8VZ92
B	250	ASN	-	expression tag	UNP W8VZ92
B	251	LEU	-	expression tag	UNP W8VZ92
B	252	TYR	-	expression tag	UNP W8VZ92

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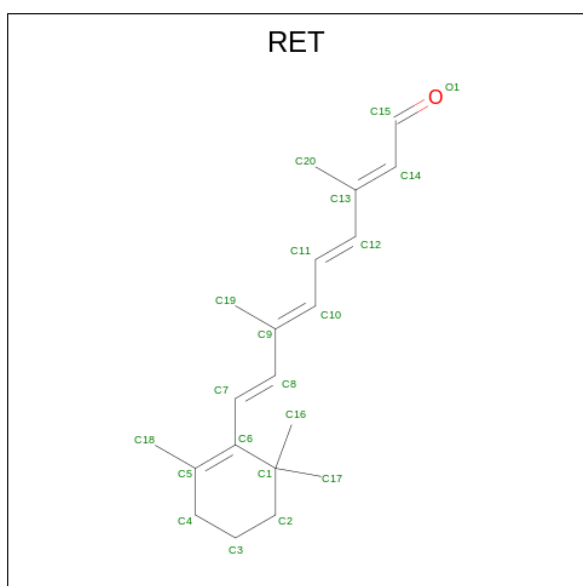
Chain	Residue	Modelled	Actual	Comment	Reference
B	253	PHE	-	expression tag	UNP W8VZ92
B	254	GLN	-	expression tag	UNP W8VZ92
B	255	GLY	-	expression tag	UNP W8VZ92
B	256	TRP	-	expression tag	UNP W8VZ92
B	257	SER	-	expression tag	UNP W8VZ92
B	258	HIS	-	expression tag	UNP W8VZ92
B	259	PRO	-	expression tag	UNP W8VZ92
B	260	GLN	-	expression tag	UNP W8VZ92
C	247	LEU	-	expression tag	UNP W8VZ92
C	248	GLU	-	expression tag	UNP W8VZ92
C	249	GLU	-	expression tag	UNP W8VZ92
C	250	ASN	-	expression tag	UNP W8VZ92
C	251	LEU	-	expression tag	UNP W8VZ92
C	252	TYR	-	expression tag	UNP W8VZ92
C	253	PHE	-	expression tag	UNP W8VZ92
C	254	GLN	-	expression tag	UNP W8VZ92
C	255	GLY	-	expression tag	UNP W8VZ92
C	256	TRP	-	expression tag	UNP W8VZ92
C	257	SER	-	expression tag	UNP W8VZ92
C	258	HIS	-	expression tag	UNP W8VZ92
C	259	PRO	-	expression tag	UNP W8VZ92
C	260	GLN	-	expression tag	UNP W8VZ92
D	247	LEU	-	expression tag	UNP W8VZ92
D	248	GLU	-	expression tag	UNP W8VZ92
D	249	GLU	-	expression tag	UNP W8VZ92
D	250	ASN	-	expression tag	UNP W8VZ92
D	251	LEU	-	expression tag	UNP W8VZ92
D	252	TYR	-	expression tag	UNP W8VZ92
D	253	PHE	-	expression tag	UNP W8VZ92
D	254	GLN	-	expression tag	UNP W8VZ92
D	255	GLY	-	expression tag	UNP W8VZ92
D	256	TRP	-	expression tag	UNP W8VZ92
D	257	SER	-	expression tag	UNP W8VZ92
D	258	HIS	-	expression tag	UNP W8VZ92
D	259	PRO	-	expression tag	UNP W8VZ92
D	260	GLN	-	expression tag	UNP W8VZ92
E	247	LEU	-	expression tag	UNP W8VZ92
E	248	GLU	-	expression tag	UNP W8VZ92
E	249	GLU	-	expression tag	UNP W8VZ92
E	250	ASN	-	expression tag	UNP W8VZ92
E	251	LEU	-	expression tag	UNP W8VZ92
E	252	TYR	-	expression tag	UNP W8VZ92

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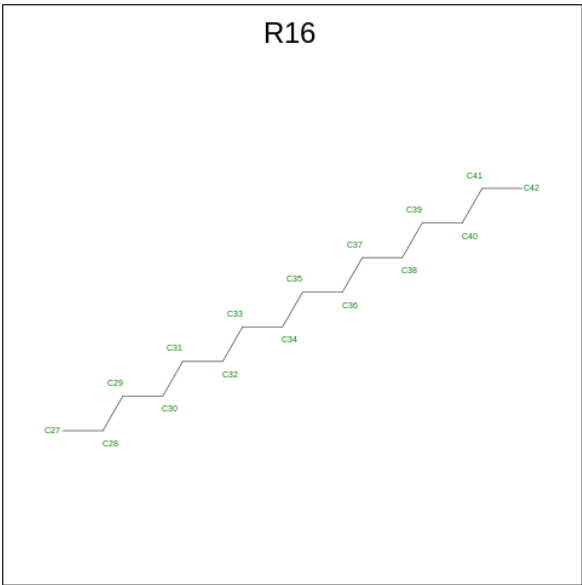
Chain	Residue	Modelled	Actual	Comment	Reference
E	253	PHE	-	expression tag	UNP W8VZ92
E	254	GLN	-	expression tag	UNP W8VZ92
E	255	GLY	-	expression tag	UNP W8VZ92
E	256	TRP	-	expression tag	UNP W8VZ92
E	257	SER	-	expression tag	UNP W8VZ92
E	258	HIS	-	expression tag	UNP W8VZ92
E	259	PRO	-	expression tag	UNP W8VZ92
E	260	GLN	-	expression tag	UNP W8VZ92

- Molecule 2 is RETINAL (CCD ID: RET) (formula: $C_{20}H_{28}O$) (labeled as "Ligand of Interest" by depositor).



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

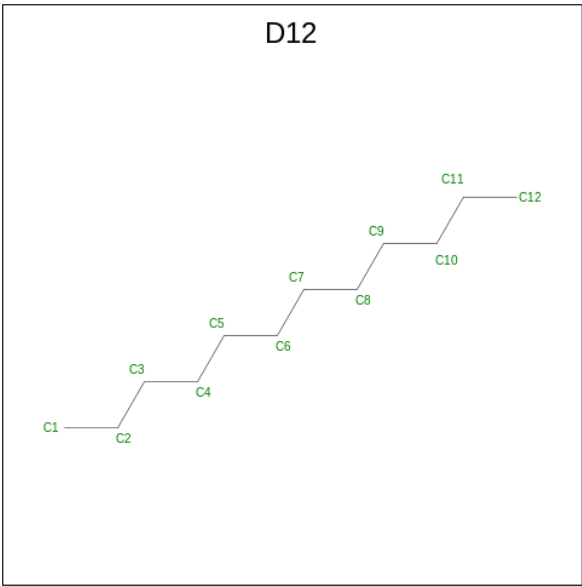
- Molecule 3 is HEXADECANE (CCD ID: R16) (formula: $C_{16}H_{34}$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	AltConf
3	A	1	Total C 16 16	0
3	A	1	Total C 16 16	0
3	B	1	Total C 16 16	0
3	C	1	Total C 16 16	0
3	C	1	Total C 16 16	0
3	C	1	Total C 16 16	0
3	C	1	Total C 16 16	0
3	D	1	Total C 16 16	0
3	D	1	Total C 16 16	0
3	D	1	Total C 16 16	0
3	D	1	Total C 16 16	0
3	E	1	Total C 16 16	0
3	E	1	Total C 16 16	0

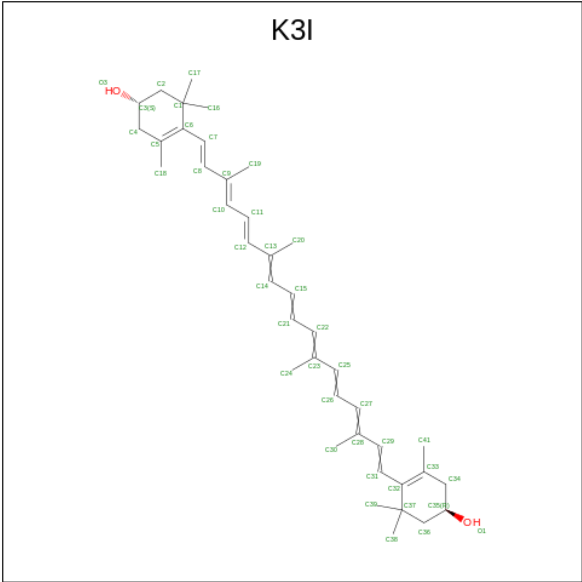
- Molecule 4 is DODECANE (CCD ID: D12) (formula: C₁₂H₂₆) (labeled as "Ligand of Interest")

by depositor).



Mol	Chain	Residues	Atoms	AltConf
4	A	1	Total C 12 12	0
4	A	1	Total C 12 12	0
4	A	1	Total C 12 12	0
4	B	1	Total C 12 12	0
4	B	1	Total C 12 12	0
4	B	1	Total C 12 12	0
4	C	1	Total C 12 12	0
4	C	1	Total C 12 12	0
4	D	1	Total C 12 12	0
4	E	1	Total C 12 12	0
4	E	1	Total C 12 12	0
4	E	1	Total C 12 12	0

- Molecule 5 is Zeaxanthin (CCD ID: K3I) (formula: C₄₀H₅₆O₂).



Mol	Chain	Residues	Atoms			AltConf
5	A	1	Total	C	O	0
			42	40	2	
5	B	1	Total	C	O	0
			42	40	2	
5	C	1	Total	C	O	0
			42	40	2	
5	D	1	Total	C	O	0
			42	40	2	
5	E	1	Total	C	O	0
			42	40	2	

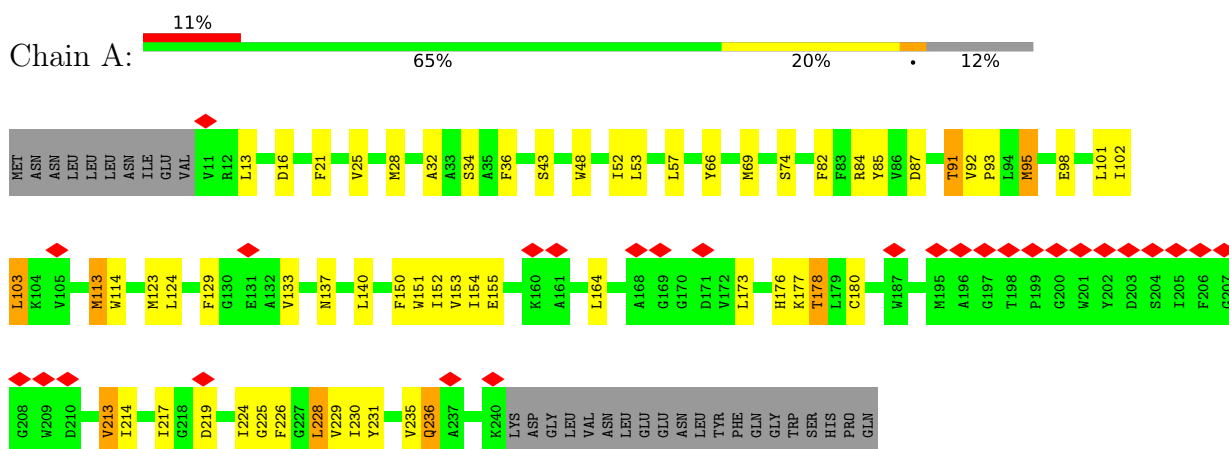
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		AltConf
6	A	8	Total	O	0
			8	8	
6	B	12	Total	O	0
			12	12	
6	C	10	Total	O	0
			10	10	
6	D	11	Total	O	0
			11	11	
6	E	10	Total	O	0
			10	10	

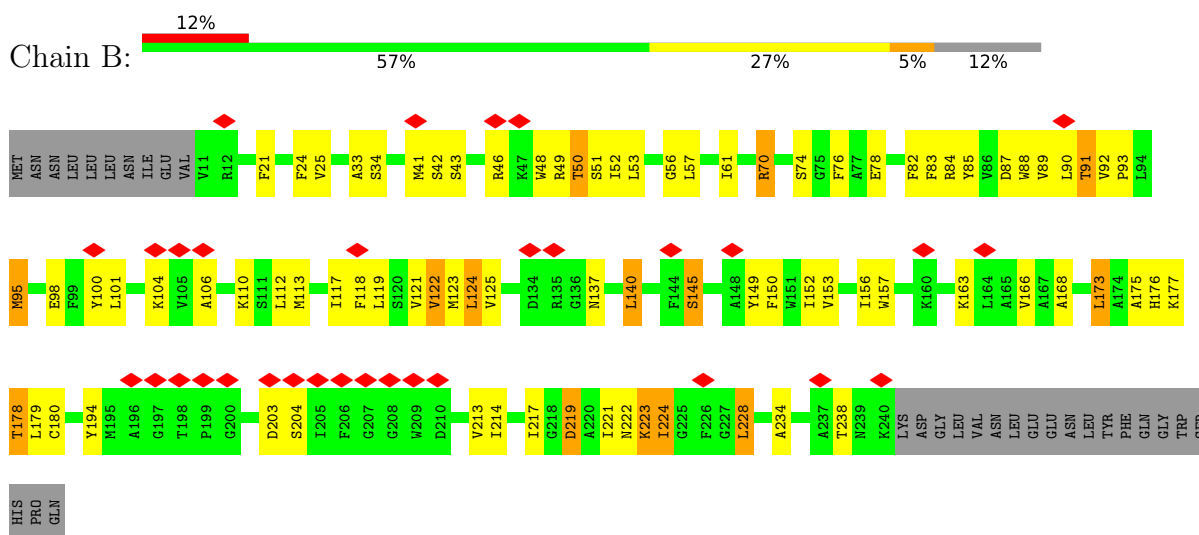
3 Residue-property plots

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 atom inclusion in map 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 diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). 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: Proteorhodopsin

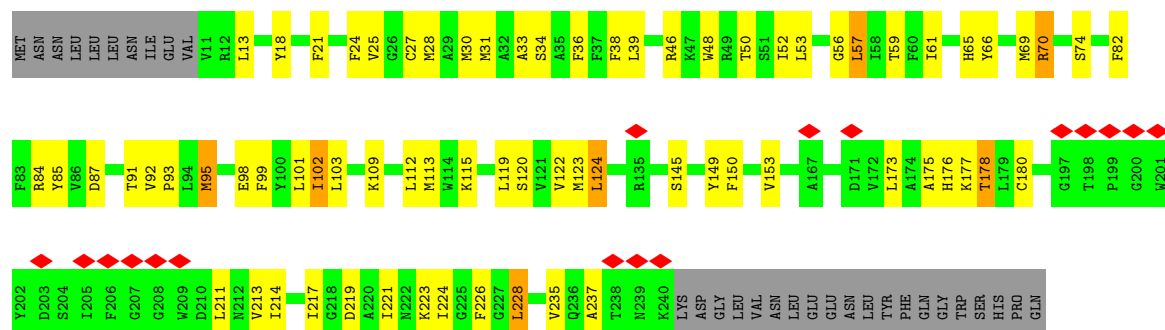


• Molecule 1: Proteorhodopsin

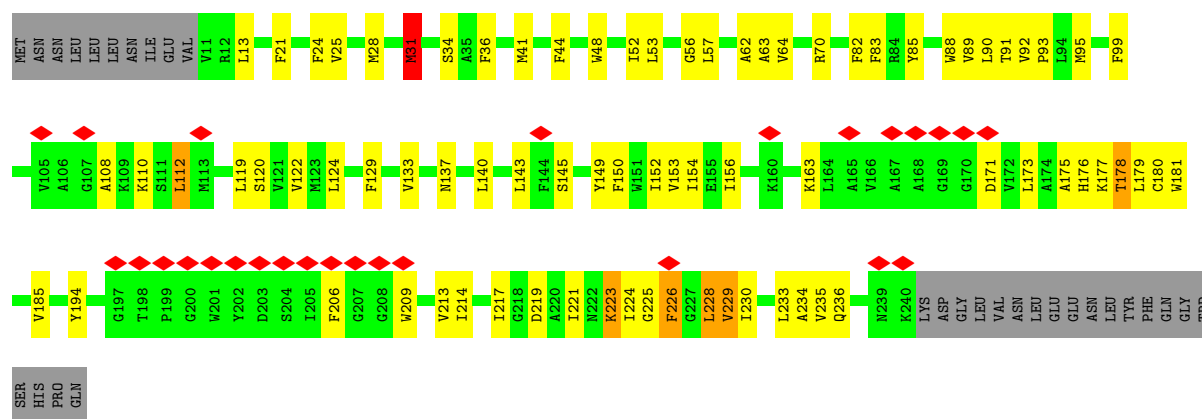


• Molecule 1: Proteorhodopsin

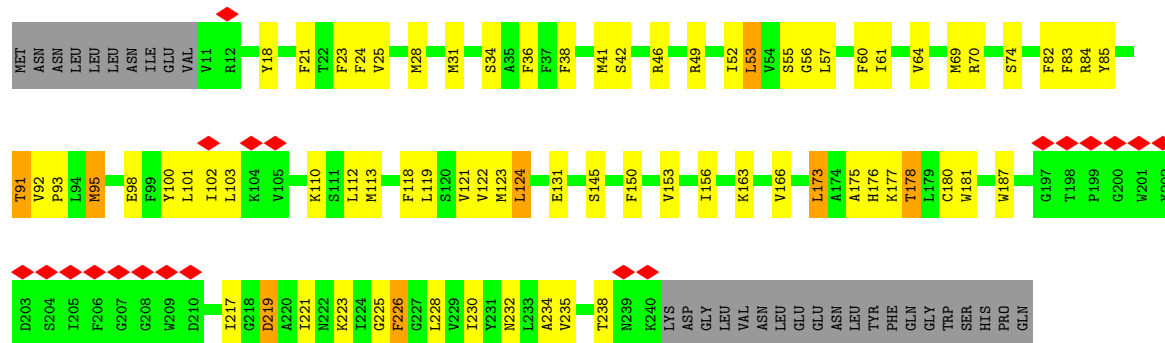




• Molecule 1: Proteorhodopsin



• Molecule 1: Proteorhodopsin



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	1814697	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	54	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.134	Depositor
Minimum map value	-0.076	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.035	Depositor
Map size (Å)	182.59999, 182.59999, 182.59999	wwPDB
Map dimensions	220, 220, 220	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: K3I, D12, R16, RET

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.55	0/1904	1.12	4/2592 (0.2%)
1	B	0.56	0/1904	1.08	3/2592 (0.1%)
1	C	0.57	0/1904	1.07	2/2592 (0.1%)
1	D	0.58	0/1904	1.12	3/2592 (0.1%)
1	E	0.57	0/1904	1.09	4/2592 (0.2%)
All	All	0.56	0/9520	1.10	16/12960 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1
1	C	0	1
All	All	0	2

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	236	GLN	N-CA-CB	9.03	125.05	110.42
1	D	31	MET	CG-SD-CE	-8.27	82.71	100.90
1	C	178	THR	CA-CB-OG1	-7.96	97.67	109.60
1	E	178	THR	OG1-CB-CG2	7.88	125.07	109.30
1	D	178	THR	CA-CB-OG1	-7.78	97.93	109.60
1	B	178	THR	CA-CB-OG1	-6.47	99.89	109.60
1	E	31	MET	CG-SD-CE	6.31	114.79	100.90
1	A	178	THR	OG1-CB-CG2	6.21	121.71	109.30
1	B	91	THR	CA-CB-OG1	-5.50	101.36	109.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	16	ASP	CA-CB-CG	5.34	117.94	112.60
1	D	229	VAL	N-CA-CB	5.26	117.30	110.57
1	E	178	THR	CA-CB-OG1	-5.22	101.77	109.60
1	A	91	THR	CA-CB-OG1	-5.12	101.92	109.60
1	B	122	VAL	N-CA-CB	5.10	116.52	110.55
1	E	91	THR	CA-CB-OG1	-5.05	102.02	109.60
1	C	50	THR	CA-CB-OG1	-5.00	102.10	109.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	70	ARG	Sidechain
1	C	70	ARG	Sidechain

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	1840	0	1823	45	0
1	B	1840	0	1823	74	0
1	C	1840	0	1823	61	0
1	D	1840	0	1823	74	0
1	E	1840	0	1823	72	0
2	A	20	0	27	1	0
2	B	20	0	27	7	0
2	C	20	0	27	4	0
2	D	20	0	27	7	0
2	E	20	0	27	15	0
3	A	32	0	68	3	0
3	B	16	0	34	1	0
3	C	64	0	136	5	0
3	D	64	0	136	1	0
3	E	32	0	68	7	0
4	A	36	0	78	2	0
4	B	36	0	78	1	0
4	C	24	0	52	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	12	0	26	1	0
4	E	36	0	78	7	0
5	A	42	0	0	0	0
5	B	42	0	0	0	0
5	C	42	0	0	0	0
5	D	42	0	0	0	0
5	E	42	0	0	0	0
6	A	8	0	0	6	0
6	B	12	0	0	5	0
6	C	10	0	0	6	0
6	D	11	0	0	4	0
6	E	10	0	0	6	0
All	All	9913	0	10004	316	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 16.

All (316) 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:B:113:MET:O	1:B:117:ILE:HD12	1.63	0.98
1:C:84:ARG:NH1	6:C:401:HOH:O	1.97	0.96
1:D:95:MET:HE1	2:D:302:RET:H201	1.53	0.91
1:D:31:MET:HE1	1:E:60:PHE:HZ	1.35	0.90
1:E:219:ASP:OD1	1:E:223:LYS:HE2	1.74	0.88
1:C:123:MET:HE3	1:C:149:TYR:HB2	1.55	0.87
2:B:302:RET:H8	2:B:302:RET:H161	1.56	0.86
1:E:123:MET:CE	2:E:302:RET:H192	2.06	0.85
1:A:32:ALA:HB2	1:B:61:ILE:HD11	1.61	0.82
1:C:30:MET:HG3	1:C:59:THR:HG23	1.62	0.81
1:C:57:LEU:O	1:C:61:ILE:HD12	1.80	0.80
1:E:123:MET:HE1	2:E:302:RET:H192	1.63	0.80
1:D:108:ALA:HB1	1:D:112:LEU:HD21	1.62	0.80
1:B:87:ASP:OD2	6:B:401:HOH:O	2.00	0.79
1:C:28:MET:HE1	6:D:407:HOH:O	1.84	0.77
1:C:211:LEU:O	1:C:214:ILE:HG13	1.85	0.76
1:E:95:MET:HE1	2:E:302:RET:H201	1.65	0.76
1:D:149:TYR:O	1:D:152:ILE:HG22	1.87	0.75
1:D:156:ILE:HG13	1:D:180:CYS:SG	2.25	0.75
1:C:66:TYR:HD1	6:C:404:HOH:O	1.70	0.73
2:E:302:RET:H161	2:E:302:RET:H8	1.69	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:219:ASP:OD2	6:D:401:HOH:O	2.08	0.72
1:E:57:LEU:O	1:E:61:ILE:HD13	1.90	0.72
1:D:95:MET:HE1	2:D:302:RET:C20	2.19	0.72
1:B:113:MET:O	1:B:117:ILE:CD1	2.36	0.72
1:B:219:ASP:O	1:B:223:LYS:HG2	1.91	0.70
1:D:31:MET:HE1	1:E:60:PHE:CZ	2.22	0.70
1:C:224:ILE:O	1:C:228:LEU:HD22	1.93	0.69
1:D:224:ILE:O	1:D:228:LEU:HD22	1.93	0.67
1:E:219:ASP:OD2	6:E:401:HOH:O	2.12	0.67
1:C:87:ASP:OD2	6:C:402:HOH:O	2.12	0.67
1:C:27:CYS:O	1:C:31:MET:HG3	1.95	0.66
1:B:234:ALA:O	1:B:238:THR:HG23	1.94	0.66
1:D:213:VAL:O	1:D:217:ILE:HD12	1.94	0.66
1:D:31:MET:CE	1:E:60:PHE:HZ	2.07	0.66
1:C:66:TYR:CD1	6:C:404:HOH:O	2.46	0.66
1:B:123:MET:CE	1:B:152:ILE:HD12	2.25	0.65
1:A:219:ASP:OD2	6:A:401:HOH:O	2.13	0.65
2:C:303:RET:H8	2:C:303:RET:H161	1.78	0.65
1:B:213:VAL:O	1:B:217:ILE:HD12	1.98	0.64
1:A:123:MET:CE	1:A:152:ILE:HD12	2.28	0.64
1:E:123:MET:HE3	2:E:302:RET:H7	1.79	0.64
1:A:87:ASP:OD2	6:A:402:HOH:O	2.15	0.63
1:D:28:MET:HE2	1:D:28:MET:HA	1.81	0.63
1:A:102:ILE:HG13	1:A:103:LEU:HD22	1.80	0.63
1:E:234:ALA:O	1:E:238:THR:HG23	1.98	0.63
1:D:31:MET:CE	1:E:60:PHE:CZ	2.82	0.63
1:B:76:PHE:O	1:B:78:GLU:OE2	2.17	0.62
1:D:150:PHE:CZ	1:D:154:ILE:HD11	2.34	0.62
1:E:123:MET:HE2	2:E:302:RET:H192	1.81	0.62
1:E:98:GLU:O	1:E:102:ILE:HG12	2.00	0.62
1:B:224:ILE:O	1:B:228:LEU:HD22	2.00	0.61
1:B:43:SER:HA	1:C:46:ARG:NH1	2.15	0.61
1:A:224:ILE:HG22	1:A:228:LEU:CD2	2.30	0.61
1:B:84:ARG:NH1	6:B:402:HOH:O	2.30	0.61
1:B:222:ASN:HB3	6:B:407:HOH:O	2.00	0.60
1:B:224:ILE:HG22	1:B:228:LEU:CD2	2.31	0.60
1:C:98:GLU:O	1:C:102:ILE:HG13	2.01	0.60
1:A:137:ASN:OD1	1:A:140:LEU:HB3	2.01	0.60
1:C:24:PHE:CE1	1:C:70:ARG:HD2	2.37	0.60
1:C:123:MET:HE2	2:C:303:RET:H7	1.84	0.59
1:B:221:ILE:HD13	3:B:303:R16:H351	1.82	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:95:MET:O	1:A:98:GLU:HB2	2.03	0.59
2:E:302:RET:H161	2:E:302:RET:C8	2.32	0.59
1:A:224:ILE:O	1:A:228:LEU:HD22	2.03	0.59
1:C:95:MET:HE1	2:C:303:RET:H201	1.84	0.58
1:E:24:PHE:CZ	1:E:70:ARG:HD3	2.38	0.58
1:D:150:PHE:CE2	1:D:154:ILE:HD11	2.38	0.58
1:E:123:MET:HE3	2:E:302:RET:H181	1.84	0.58
1:D:24:PHE:CZ	1:D:70:ARG:HD3	2.38	0.58
1:D:224:ILE:HG22	1:D:228:LEU:CD2	2.33	0.58
1:C:224:ILE:HG22	1:C:228:LEU:CD2	2.33	0.57
1:E:84:ARG:NH1	6:E:401:HOH:O	2.22	0.57
1:E:219:ASP:OD1	6:E:402:HOH:O	2.18	0.57
1:B:95:MET:O	1:B:98:GLU:HB2	2.04	0.57
1:C:221:ILE:HD13	3:C:304:R16:H342	1.87	0.57
1:A:173:LEU:HD13	1:A:177:LYS:HD2	1.87	0.57
1:D:44:PHE:CE2	1:D:234:ALA:HB1	2.40	0.56
1:B:123:MET:HE1	1:B:152:ILE:HD12	1.87	0.56
3:A:303:R16:H272	1:B:50:THR:HG21	1.86	0.56
1:B:24:PHE:CE1	1:B:70:ARG:HD2	2.41	0.56
1:B:106:ALA:HB2	1:B:168:ALA:HB2	1.86	0.56
1:C:28:MET:CE	6:D:407:HOH:O	2.46	0.56
1:B:95:MET:HE1	2:B:302:RET:H201	1.87	0.56
1:D:62:ALA:HB2	1:D:223:LYS:HE3	1.87	0.56
1:E:100:TYR:CD1	1:E:113:MET:HG2	2.41	0.56
2:B:302:RET:H161	2:B:302:RET:C8	2.31	0.56
1:E:235:VAL:HG11	4:E:305:D12:H121	1.88	0.56
1:B:51:SER:OG	1:B:101:LEU:HB3	2.06	0.55
1:A:226:PHE:HB3	6:A:407:HOH:O	2.07	0.55
1:E:123:MET:HE3	2:E:302:RET:C7	2.38	0.54
1:A:69:MET:HE1	6:A:405:HOH:O	2.06	0.54
1:B:113:MET:HE3	1:B:117:ILE:HD11	1.88	0.54
1:C:213:VAL:O	1:C:217:ILE:HD12	2.07	0.54
1:D:171:ASP:OD2	1:D:236:GLN:OE1	2.26	0.54
1:B:48:TRP:CZ3	1:B:238:THR:HG22	2.43	0.54
1:B:219:ASP:CG	6:B:402:HOH:O	2.50	0.54
1:E:41:MET:SD	1:E:53:LEU:HD23	2.48	0.53
1:D:185:VAL:HG23	4:D:306:D12:H92	1.89	0.53
1:C:150:PHE:O	1:C:153:VAL:HG12	2.08	0.53
3:E:304:R16:H332	4:E:305:D12:H102	1.91	0.53
1:A:173:LEU:CD1	1:A:177:LYS:HD2	2.39	0.53
1:E:119:LEU:O	1:E:122:VAL:HG12	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:150:PHE:O	1:E:153:VAL:HG12	2.08	0.53
1:A:150:PHE:O	1:A:153:VAL:HG12	2.09	0.52
1:B:89:VAL:HG12	1:B:124:LEU:HB3	1.90	0.52
1:C:30:MET:O	1:C:59:THR:CG2	2.57	0.52
1:D:150:PHE:O	1:D:153:VAL:HG12	2.08	0.52
1:D:206:PHE:CG	1:D:206:PHE:O	2.63	0.52
1:E:52:ILE:N	1:E:52:ILE:HD12	2.24	0.52
1:A:66:TYR:OH	1:A:84:ARG:NH2	2.42	0.52
1:A:57:LEU:HD11	1:E:36:PHE:HB2	1.91	0.52
1:D:221:ILE:HD13	3:D:303:R16:H351	1.92	0.52
2:D:302:RET:H161	2:D:302:RET:H8	1.92	0.52
1:D:181:TRP:O	1:D:185:VAL:HG22	2.10	0.51
1:B:150:PHE:HA	1:B:153:VAL:HG12	1.92	0.51
1:B:41:MET:HG2	1:B:49:ARG:HG3	1.93	0.51
1:C:119:LEU:O	1:C:122:VAL:HG12	2.10	0.51
1:D:119:LEU:O	1:D:122:VAL:HG12	2.11	0.51
1:A:13:LEU:HD11	1:B:83:PHE:CD1	2.46	0.51
1:B:119:LEU:O	1:B:122:VAL:HG12	2.11	0.51
1:D:88:TRP:C	1:D:124:LEU:HD23	2.36	0.51
1:E:38:PHE:O	6:E:403:HOH:O	2.19	0.51
1:E:123:MET:HE1	2:E:302:RET:C19	2.39	0.51
1:B:150:PHE:O	1:B:153:VAL:HG12	2.11	0.51
1:D:44:PHE:HE2	1:D:234:ALA:C	2.18	0.50
1:E:150:PHE:HA	1:E:153:VAL:HG12	1.92	0.50
1:C:219:ASP:CG	6:C:402:HOH:O	2.54	0.50
1:E:110:LYS:O	1:E:113:MET:HB2	2.10	0.50
1:B:222:ASN:CB	6:B:407:HOH:O	2.59	0.50
1:A:43:SER:C	1:B:46:ARG:HH12	2.19	0.50
1:B:137:ASN:OD1	1:B:140:LEU:HD23	2.11	0.50
1:D:108:ALA:HB1	1:D:112:LEU:CD2	2.35	0.50
1:C:99:PHE:O	1:C:102:ILE:HD12	2.12	0.50
1:E:21:PHE:O	1:E:25:VAL:HG23	2.12	0.50
1:B:21:PHE:O	1:B:25:VAL:HG23	2.12	0.49
1:B:100:TYR:CD1	1:B:113:MET:HG2	2.47	0.49
1:C:123:MET:HE1	1:C:149:TYR:HD2	1.76	0.49
1:E:41:MET:CE	1:E:49:ARG:HB3	2.41	0.49
1:D:149:TYR:O	1:D:152:ILE:CG2	2.59	0.49
1:E:98:GLU:HG2	1:E:230:ILE:HD11	1.94	0.49
1:A:129:PHE:HA	1:A:133:VAL:HG12	1.95	0.49
1:A:150:PHE:HA	1:A:153:VAL:HG12	1.93	0.49
1:C:24:PHE:CZ	1:C:70:ARG:HD2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:24:PHE:CE1	1:E:70:ARG:HD3	2.48	0.49
1:A:25:VAL:HG13	1:B:90:LEU:HD21	1.95	0.49
1:B:24:PHE:CZ	1:B:70:ARG:HD2	2.47	0.49
1:D:88:TRP:C	1:D:124:LEU:CD2	2.85	0.49
1:E:228:LEU:HD12	4:E:305:D12:H91	1.95	0.49
1:A:219:ASP:CG	6:A:402:HOH:O	2.55	0.49
1:B:33:ALA:CB	1:B:224:ILE:HD13	2.43	0.49
1:A:36:PHE:CD1	1:B:53:LEU:HD21	2.48	0.48
1:C:38:PHE:O	6:C:403:HOH:O	2.20	0.48
1:D:44:PHE:CE2	1:D:234:ALA:CB	2.95	0.48
1:E:187:TRP:NE1	6:E:404:HOH:O	2.36	0.48
1:D:236:GLN:OE1	1:D:236:GLN:C	2.56	0.48
1:D:99:PHE:CD1	1:D:156:ILE:HD11	2.48	0.48
1:B:52:ILE:O	1:B:53:LEU:C	2.56	0.48
1:C:92:VAL:HG21	1:C:124:LEU:HD21	1.96	0.48
1:B:113:MET:CE	1:B:117:ILE:HD11	2.44	0.48
1:C:175:ALA:O	1:C:178:THR:OG1	2.31	0.48
1:D:175:ALA:O	1:D:178:THR:OG1	2.32	0.47
1:E:221:ILE:HD13	3:E:303:R16:H361	1.96	0.47
1:B:156:ILE:HD13	1:B:180:CYS:SG	2.54	0.47
1:D:95:MET:CE	2:D:302:RET:C20	2.89	0.47
1:B:82:PHE:HA	1:B:85:TYR:HD2	1.80	0.47
1:D:21:PHE:O	1:D:25:VAL:HG23	2.15	0.47
4:C:306:D12:H111	1:D:110:LYS:HE3	1.96	0.47
1:D:150:PHE:HA	1:D:153:VAL:HG12	1.95	0.47
1:D:13:LEU:HD22	1:E:83:PHE:CE1	2.50	0.47
1:D:44:PHE:CD2	1:D:234:ALA:HB1	2.50	0.47
1:A:48:TRP:HD1	6:A:406:HOH:O	1.95	0.47
1:B:42:SER:O	1:C:46:ARG:NH1	2.41	0.47
1:C:18:TYR:CE1	1:D:133:VAL:HG12	2.50	0.47
1:C:150:PHE:HA	1:C:153:VAL:HG12	1.96	0.47
1:D:194:TYR:HD2	2:D:302:RET:H22	1.79	0.47
1:E:225:GLY:HA2	3:E:303:R16:H312	1.97	0.47
1:C:101:LEU:HD21	3:C:302:R16:H301	1.96	0.47
3:C:305:R16:H331	4:C:306:D12:H112	1.97	0.47
1:B:110:LYS:O	1:B:113:MET:HB2	2.15	0.47
1:C:21:PHE:O	1:C:25:VAL:HG23	2.14	0.47
1:C:52:ILE:O	1:C:53:LEU:C	2.57	0.47
1:E:123:MET:CE	2:E:302:RET:C19	2.87	0.47
1:A:28:MET:HE2	1:B:61:ILE:HG23	1.98	0.46
1:B:156:ILE:HD12	1:B:157:TRP:CD1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:176:HIS:O	1:C:177:LYS:C	2.58	0.46
1:C:25:VAL:HG13	1:D:90:LEU:HD21	1.96	0.46
1:C:36:PHE:HB2	1:D:57:LEU:HD11	1.97	0.46
1:E:95:MET:CE	2:E:302:RET:H201	2.38	0.46
1:C:102:ILE:HD12	1:C:103:LEU:HG	1.98	0.46
1:D:88:TRP:O	1:D:124:LEU:CD2	2.64	0.46
1:E:70:ARG:O	1:E:70:ARG:HG2	2.13	0.46
1:D:173:LEU:O	1:D:176:HIS:HB3	2.16	0.46
1:D:92:VAL:HG21	1:D:124:LEU:HD21	1.98	0.46
1:D:163:LYS:HD2	1:D:163:LYS:N	2.31	0.46
1:B:33:ALA:HB2	1:B:224:ILE:HD13	1.96	0.46
1:D:52:ILE:O	1:D:53:LEU:C	2.59	0.46
1:C:92:VAL:HB	1:C:93:PRO:HD3	1.99	0.45
1:B:149:TYR:HB2	2:B:302:RET:C18	2.46	0.45
1:D:92:VAL:HB	1:D:93:PRO:HD3	1.98	0.45
1:B:92:VAL:HB	1:B:93:PRO:HD3	1.97	0.45
1:C:235:VAL:HG11	4:C:306:D12:C12	2.46	0.45
1:E:52:ILE:O	1:E:53:LEU:C	2.58	0.45
1:A:13:LEU:HD23	1:A:13:LEU:H	1.81	0.45
1:A:52:ILE:HD13	1:A:230:ILE:HG22	1.98	0.45
1:C:13:LEU:HD22	1:D:83:PHE:CE1	2.52	0.45
3:E:304:R16:H312	4:E:305:D12:H123	1.98	0.45
3:A:303:R16:H332	4:A:304:D12:H101	1.99	0.45
1:D:176:HIS:O	1:D:177:LYS:C	2.60	0.45
1:A:21:PHE:O	1:A:25:VAL:HG23	2.16	0.45
1:C:56:GLY:O	1:C:57:LEU:C	2.60	0.45
1:D:235:VAL:HG11	4:E:301:D12:H121	1.99	0.45
1:D:36:PHE:HB2	1:E:57:LEU:HD21	1.98	0.44
1:E:92:VAL:HG21	1:E:124:LEU:HD21	1.99	0.44
1:A:176:HIS:O	1:A:177:LYS:C	2.60	0.44
1:B:176:HIS:O	1:B:177:LYS:C	2.58	0.44
1:B:33:ALA:HB2	1:B:224:ILE:CD1	2.47	0.44
1:E:181:TRP:HB3	4:E:306:D12:H112	2.00	0.44
1:A:92:VAL:HG21	1:A:124:LEU:HD21	1.98	0.44
1:A:95:MET:HE1	2:A:301:RET:H201	1.99	0.44
1:D:48:TRP:CZ3	1:D:233:LEU:HD12	2.53	0.44
1:A:151:TRP:O	1:A:154:ILE:HG13	2.17	0.44
1:C:123:MET:HE3	1:C:149:TYR:CB	2.39	0.44
1:E:92:VAL:HB	1:E:93:PRO:HD3	1.98	0.44
1:A:92:VAL:HB	1:A:93:PRO:HD3	1.98	0.44
1:A:213:VAL:O	1:A:217:ILE:HD12	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:41:MET:HE2	1:B:53:LEU:HA	2.00	0.44
2:C:303:RET:C8	2:C:303:RET:H171	2.47	0.44
1:A:52:ILE:O	1:A:53:LEU:C	2.60	0.44
1:D:92:VAL:CG2	1:D:124:LEU:HD21	2.48	0.44
1:E:175:ALA:HB2	6:E:407:HOH:O	2.18	0.44
1:A:231:TYR:O	1:A:235:VAL:HG23	2.17	0.44
1:B:137:ASN:OD1	1:B:137:ASN:O	2.36	0.44
1:C:82:PHE:HA	1:C:85:TYR:HD2	1.83	0.43
1:C:173:LEU:O	1:C:176:HIS:HB3	2.18	0.43
1:D:129:PHE:HA	1:D:133:VAL:HG22	2.00	0.43
1:E:232:ASN:HA	1:E:235:VAL:HG12	2.00	0.43
1:E:176:HIS:O	1:E:177:LYS:C	2.58	0.43
1:C:31:MET:HE1	1:D:64:VAL:HG11	2.00	0.43
1:C:219:ASP:O	1:C:223:LYS:HG2	2.19	0.43
1:D:28:MET:HE1	1:E:64:VAL:HG11	2.00	0.43
1:E:123:MET:CE	2:E:302:RET:H7	2.47	0.43
4:A:307:D12:H112	4:E:305:D12:H11	2.00	0.43
1:C:217:ILE:HA	3:C:304:R16:H411	2.00	0.43
1:D:56:GLY:O	1:D:57:LEU:C	2.61	0.43
1:B:88:TRP:CD1	2:B:302:RET:H14	2.54	0.43
1:D:88:TRP:CD1	2:D:302:RET:H14	2.53	0.43
1:E:82:PHE:HA	1:E:85:TYR:HD2	1.83	0.43
1:B:122:VAL:HA	1:B:125:VAL:HG12	2.00	0.43
1:C:33:ALA:HB3	1:C:59:THR:HG21	2.00	0.43
1:D:41:MET:HE2	1:D:41:MET:HB3	1.93	0.43
1:E:118:PHE:O	1:E:121:VAL:HG22	2.19	0.43
1:E:173:LEU:O	1:E:176:HIS:HB3	2.19	0.43
1:D:24:PHE:HD1	6:D:410:HOH:O	2.01	0.43
1:E:24:PHE:O	1:E:28:MET:HG2	2.19	0.43
1:B:56:GLY:O	1:B:57:LEU:C	2.61	0.43
1:B:92:VAL:HG21	1:B:124:LEU:HD21	2.00	0.43
1:E:217:ILE:HD13	3:E:303:R16:H412	2.01	0.43
1:B:156:ILE:CD1	1:B:157:TRP:CD1	3.01	0.42
1:A:32:ALA:CB	1:B:61:ILE:HD11	2.42	0.42
1:A:224:ILE:HG22	1:A:228:LEU:HD22	2.01	0.42
1:E:46:ARG:HD2	1:E:46:ARG:HA	1.81	0.42
1:A:82:PHE:HA	1:A:85:TYR:HD2	1.83	0.42
1:B:173:LEU:O	1:B:176:HIS:HB3	2.19	0.42
1:C:123:MET:HE1	1:C:149:TYR:CD2	2.53	0.42
1:E:145:SER:OG	2:E:302:RET:H41	2.19	0.42
1:E:52:ILE:O	1:E:55:SER:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:ILE:CD1	1:C:103:LEU:HG	2.49	0.42
1:D:82:PHE:HA	1:D:85:TYR:HD2	1.84	0.42
1:C:24:PHE:CD1	1:C:70:ARG:HD2	2.55	0.42
1:C:211:LEU:HA	1:C:214:ILE:HG13	2.02	0.42
1:D:89:VAL:HA	1:D:124:LEU:CD2	2.49	0.42
1:D:137:ASN:OD1	1:D:140:LEU:HB3	2.20	0.42
1:A:213:VAL:HG13	1:A:214:ILE:HD12	2.02	0.42
1:B:214:ILE:HG23	4:B:305:D12:H11	2.01	0.42
1:D:209:TRP:HB3	1:D:214:ILE:HD11	2.00	0.42
1:D:225:GLY:O	1:D:226:PHE:C	2.63	0.42
1:A:133:VAL:HG23	1:E:18:TYR:OH	2.20	0.41
1:B:145:SER:HB2	2:B:302:RET:H41	2.02	0.41
1:E:56:GLY:O	1:E:57:LEU:C	2.62	0.41
1:A:229:VAL:O	1:A:230:ILE:C	2.63	0.41
1:B:163:LYS:HA	1:B:166:VAL:HG22	2.01	0.41
1:B:177:LYS:O	1:B:180:CYS:N	2.53	0.41
1:B:118:PHE:O	1:B:121:VAL:HG22	2.21	0.41
1:C:65:HIS:ND1	1:C:69:MET:HE2	2.36	0.41
1:C:113:MET:HE2	1:C:113:MET:HB3	1.94	0.41
1:E:156:ILE:HD12	1:E:180:CYS:SG	2.60	0.41
1:E:163:LYS:HA	1:E:166:VAL:HG22	2.02	0.41
3:A:303:R16:H421	1:B:90:LEU:HD12	2.03	0.41
1:B:194:TYR:HD2	2:B:302:RET:H22	1.86	0.41
1:C:211:LEU:HA	1:C:214:ILE:CG1	2.49	0.41
1:E:225:GLY:O	1:E:226:PHE:C	2.63	0.41
1:C:25:VAL:HB	3:C:304:R16:H422	2.03	0.41
1:C:48:TRP:CZ2	1:C:237:ALA:HB3	2.55	0.41
1:A:225:GLY:O	1:A:226:PHE:C	2.63	0.41
1:C:177:LYS:O	1:C:180:CYS:N	2.53	0.41
1:D:70:ARG:O	1:D:70:ARG:HG2	2.21	0.41
1:E:177:LYS:O	1:E:180:CYS:N	2.53	0.41
1:E:217:ILE:HD13	3:E:303:R16:H421	2.02	0.41
1:D:177:LYS:O	1:D:180:CYS:N	2.54	0.41
1:B:175:ALA:O	1:B:179:LEU:HG	2.21	0.40
1:E:23:PHE:HE1	1:E:69:MET:HE3	1.86	0.40
1:E:217:ILE:HD13	3:E:303:R16:C42	2.52	0.40
1:A:113:MET:HG3	1:A:114:TRP:N	2.36	0.40
1:D:219:ASP:O	1:D:223:LYS:HG3	2.21	0.40
1:B:48:TRP:O	1:B:49:ARG:C	2.64	0.40
1:B:175:ALA:O	1:B:178:THR:OG1	2.39	0.40
1:A:177:LYS:O	1:A:180:CYS:N	2.55	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:62:ALA:O	1:D:63:ALA:C	2.63	0.40
1:E:42:SER:HA	1:E:49:ARG:HH12	1.87	0.40
1:E:124:LEU:HD13	1:E:124:LEU:HA	1.96	0.40
1:B:124:LEU:HD13	1:B:124:LEU:HA	1.95	0.40
1:B:203:ASP:OD1	1:B:204:SER:N	2.55	0.40
1:D:229:VAL:O	1:D:230:ILE:C	2.64	0.40
2:D:302:RET:H7	2:D:302:RET:H181	1.81	0.40
1:E:95:MET:HE1	2:E:302:RET:C15	2.51	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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	228/260 (88%)	221 (97%)	7 (3%)	0	100	100
1	B	228/260 (88%)	224 (98%)	4 (2%)	0	100	100
1	C	228/260 (88%)	223 (98%)	5 (2%)	0	100	100
1	D	228/260 (88%)	221 (97%)	7 (3%)	0	100	100
1	E	228/260 (88%)	222 (97%)	6 (3%)	0	100	100
All	All	1140/1300 (88%)	1111 (98%)	29 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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	183/211 (87%)	170 (93%)	13 (7%)	12	15
1	B	183/211 (87%)	168 (92%)	15 (8%)	9	10
1	C	183/211 (87%)	168 (92%)	15 (8%)	9	10
1	D	183/211 (87%)	172 (94%)	11 (6%)	16	22
1	E	183/211 (87%)	169 (92%)	14 (8%)	10	12
All	All	915/1055 (87%)	847 (93%)	68 (7%)	14	13

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	34	SER
1	A	74	SER
1	A	91	THR
1	A	95	MET
1	A	101	LEU
1	A	103	LEU
1	A	113	MET
1	A	155	GLU
1	A	164	LEU
1	A	178	THR
1	A	213	VAL
1	A	228	LEU
1	A	236	GLN
1	B	34	SER
1	B	50	THR
1	B	74	SER
1	B	91	THR
1	B	95	MET
1	B	104	LYS
1	B	112	LEU
1	B	124	LEU
1	B	140	LEU
1	B	145	SER
1	B	173	LEU
1	B	219	ASP
1	B	223	LYS
1	B	224	ILE
1	B	228	LEU
1	C	34	SER

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Mol	Chain	Res	Type
1	C	39	LEU
1	C	57	LEU
1	C	74	SER
1	C	91	THR
1	C	95	MET
1	C	102	ILE
1	C	109	LYS
1	C	112	LEU
1	C	115	LYS
1	C	120	SER
1	C	124	LEU
1	C	145	SER
1	C	226	PHE
1	C	228	LEU
1	D	31	MET
1	D	34	SER
1	D	91	THR
1	D	112	LEU
1	D	120	SER
1	D	143	LEU
1	D	145	SER
1	D	179	LEU
1	D	223	LYS
1	D	226	PHE
1	D	228	LEU
1	E	34	SER
1	E	53	LEU
1	E	74	SER
1	E	91	THR
1	E	95	MET
1	E	101	LEU
1	E	103	LEU
1	E	112	LEU
1	E	124	LEU
1	E	131	GLU
1	E	173	LEU
1	E	178	THR
1	E	219	ASP
1	E	226	PHE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	239	ASN
1	D	176	HIS
1	D	232	ASN
1	D	236	GLN
1	E	176	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

35 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	RET	D	302	1	20,20,21	2.59	3 (15%)	27,27,28	1.45	6 (22%)
3	R16	D	304	-	15,15,15	0.09	0	14,14,14	0.09	0
4	D12	E	305	-	11,11,11	0.20	0	10,10,10	0.10	0
5	K3I	C	308	-	43,43,43	1.62	6 (13%)	58,60,60	1.07	5 (8%)
3	R16	D	303	-	15,15,15	0.19	0	14,14,14	0.16	0
4	D12	A	305	-	11,11,11	0.14	0	10,10,10	0.12	0
2	RET	C	303	1	20,20,21	2.87	5 (25%)	27,27,28	1.38	3 (11%)
4	D12	A	304	-	11,11,11	0.14	0	10,10,10	0.15	0
3	R16	B	303	-	15,15,15	0.25	0	14,14,14	0.20	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	R16	A	303	-	15,15,15	0.19	0	14,14,14	0.23	0
4	D12	B	301	-	11,11,11	0.11	0	10,10,10	0.07	0
4	D12	B	304	-	11,11,11	0.18	0	10,10,10	0.08	0
3	R16	E	303	-	15,15,15	0.16	0	14,14,14	0.14	0
4	D12	E	301	-	11,11,11	0.15	0	10,10,10	0.09	0
4	D12	D	306	-	11,11,11	0.21	0	10,10,10	0.20	0
2	RET	A	301	1	20,20,21	2.22	5 (25%)	27,27,28	1.35	5 (18%)
3	R16	C	302	-	15,15,15	0.26	0	14,14,14	0.22	0
4	D12	B	305	-	11,11,11	0.17	0	10,10,10	0.17	0
3	R16	C	305	-	15,15,15	0.19	0	14,14,14	0.19	0
4	D12	E	306	-	11,11,11	0.13	0	10,10,10	0.14	0
3	R16	D	305	-	15,15,15	0.16	0	14,14,14	0.15	0
4	D12	A	307	-	11,11,11	0.13	0	10,10,10	0.14	0
5	K3I	D	307	-	43,43,43	1.80	6 (13%)	58,60,60	1.31	5 (8%)
2	RET	E	302	1	20,20,21	2.64	4 (20%)	27,27,28	1.41	3 (11%)
3	R16	C	304	-	15,15,15	0.14	0	14,14,14	0.28	0
3	R16	A	302	-	15,15,15	0.21	0	14,14,14	0.11	0
3	R16	E	304	-	15,15,15	0.21	0	14,14,14	0.12	0
4	D12	C	306	-	11,11,11	0.18	0	10,10,10	0.22	0
5	K3I	E	307	-	43,43,43	1.75	6 (13%)	58,60,60	0.97	4 (6%)
2	RET	B	302	1	20,20,21	2.77	4 (20%)	27,27,28	1.55	7 (25%)
4	D12	C	307	-	11,11,11	0.14	0	10,10,10	0.07	0
5	K3I	A	306	-	43,43,43	1.61	6 (13%)	58,60,60	1.27	6 (10%)
3	R16	C	301	-	15,15,15	0.11	0	14,14,14	0.10	0
3	R16	D	301	-	15,15,15	0.16	0	14,14,14	0.11	0
5	K3I	B	306	-	43,43,43	1.63	7 (16%)	58,60,60	1.77	11 (18%)

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	RET	D	302	1	-	2/13/30/31	0/1/1/1
3	R16	D	304	-	-	4/13/13/13	-
4	D12	E	305	-	-	1/9/9/9	-
5	K3I	C	308	-	-	1/29/67/67	0/2/2/2
3	R16	D	303	-	-	4/13/13/13	-
4	D12	A	305	-	-	1/9/9/9	-
2	RET	C	303	1	-	2/13/30/31	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	D12	A	304	-	-	4/9/9/9	-
3	R16	B	303	-	-	2/13/13/13	-
3	R16	A	303	-	-	2/13/13/13	-
4	D12	B	301	-	-	1/9/9/9	-
4	D12	B	304	-	-	1/9/9/9	-
3	R16	E	303	-	-	7/13/13/13	-
4	D12	E	301	-	-	2/9/9/9	-
4	D12	D	306	-	-	3/9/9/9	-
2	RET	A	301	1	-	6/13/30/31	0/1/1/1
3	R16	C	302	-	-	3/13/13/13	-
4	D12	B	305	-	-	2/9/9/9	-
3	R16	C	305	-	-	4/13/13/13	-
4	D12	E	306	-	-	1/9/9/9	-
3	R16	D	305	-	-	4/13/13/13	-
4	D12	A	307	-	-	2/9/9/9	-
5	K3I	D	307	-	-	5/29/67/67	0/2/2/2
2	RET	E	302	1	-	0/13/30/31	0/1/1/1
3	R16	C	304	-	-	2/13/13/13	-
3	R16	A	302	-	-	4/13/13/13	-
3	R16	E	304	-	-	5/13/13/13	-
4	D12	C	306	-	-	4/9/9/9	-
5	K3I	E	307	-	-	10/29/67/67	0/2/2/2
2	RET	B	302	1	-	2/13/30/31	0/1/1/1
4	D12	C	307	-	-	3/9/9/9	-
5	K3I	A	306	-	-	6/29/67/67	0/2/2/2
3	R16	C	301	-	-	2/13/13/13	-
3	R16	D	301	-	-	1/13/13/13	-
5	K3I	B	306	-	-	8/29/67/67	0/2/2/2

All (52) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	303	RET	C14-C13	10.46	1.41	1.33
2	B	302	RET	C14-C13	9.80	1.41	1.33
2	E	302	RET	C14-C13	9.63	1.41	1.33
2	D	302	RET	C14-C13	8.86	1.40	1.33
2	A	301	RET	C14-C13	7.24	1.39	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	302	RET	C10-C9	5.84	1.43	1.35
5	E	307	K3I	C27-C28	5.52	1.43	1.35
2	B	302	RET	C10-C9	5.44	1.43	1.35
5	D	307	K3I	C27-C28	5.27	1.42	1.35
5	D	307	K3I	C10-C9	5.24	1.42	1.35
5	E	307	K3I	C22-C23	5.22	1.42	1.35
5	D	307	K3I	C22-C23	5.14	1.42	1.35
5	D	307	K3I	C14-C13	5.12	1.42	1.35
2	E	302	RET	C10-C9	4.95	1.42	1.35
5	B	306	K3I	C27-C28	4.88	1.42	1.35
2	C	303	RET	C10-C9	4.74	1.42	1.35
5	E	307	K3I	C14-C13	4.69	1.42	1.35
5	A	306	K3I	C10-C9	4.69	1.42	1.35
5	E	307	K3I	C10-C9	4.63	1.41	1.35
5	C	308	K3I	C10-C9	4.63	1.41	1.35
5	A	306	K3I	C27-C28	4.60	1.41	1.35
5	B	306	K3I	C14-C13	4.58	1.41	1.35
5	C	308	K3I	C14-C13	4.57	1.41	1.35
5	B	306	K3I	C10-C9	4.51	1.41	1.35
5	C	308	K3I	C27-C28	4.48	1.41	1.35
5	C	308	K3I	C22-C23	4.46	1.41	1.35
5	A	306	K3I	C22-C23	4.37	1.41	1.35
5	A	306	K3I	C14-C13	4.35	1.41	1.35
5	B	306	K3I	C22-C23	4.06	1.41	1.35
2	A	301	RET	C10-C9	3.70	1.40	1.35
2	A	301	RET	C8-C9	-3.31	1.38	1.45
5	D	307	K3I	C26-C25	3.22	1.42	1.34
5	D	307	K3I	C11-C12	3.10	1.42	1.34
5	A	306	K3I	C26-C25	2.85	1.41	1.34
5	E	307	K3I	C26-C25	2.83	1.41	1.34
2	A	301	RET	C12-C13	-2.83	1.39	1.45
5	B	306	K3I	C11-C12	2.80	1.41	1.34
5	A	306	K3I	C11-C12	2.79	1.41	1.34
5	E	307	K3I	C11-C12	2.77	1.41	1.34
5	C	308	K3I	C11-C12	2.77	1.41	1.34
5	B	306	K3I	C26-C25	2.74	1.41	1.34
5	C	308	K3I	C26-C25	2.72	1.41	1.34
2	C	303	RET	C8-C9	-2.65	1.40	1.45
2	C	303	RET	C12-C13	-2.63	1.40	1.45
2	B	302	RET	C8-C7	2.59	1.40	1.33
2	B	302	RET	C11-C12	2.41	1.40	1.34
2	D	302	RET	C12-C13	-2.30	1.41	1.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	301	RET	C11-C12	2.25	1.40	1.34
2	E	302	RET	C11-C12	2.17	1.40	1.34
5	B	306	K3I	C12-C13	-2.08	1.41	1.45
2	E	302	RET	C8-C7	2.06	1.39	1.33
2	C	303	RET	C11-C12	2.05	1.39	1.34

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	306	K3I	C21-C15-C14	5.81	135.38	123.47
5	B	306	K3I	C24-C23-C22	-5.61	115.06	122.92
5	B	306	K3I	C25-C23-C22	5.52	127.41	118.94
5	D	307	K3I	C21-C15-C14	4.89	133.50	123.47
5	A	306	K3I	C24-C23-C25	4.72	125.51	118.08
2	E	302	RET	C19-C9-C10	-4.31	116.88	122.92
2	B	302	RET	C19-C9-C10	-3.88	117.49	122.92
2	C	303	RET	C19-C9-C10	-3.78	117.63	122.92
5	D	307	K3I	C20-C13-C14	-3.77	117.65	122.92
5	D	307	K3I	C24-C23-C25	3.76	123.99	118.08
5	D	307	K3I	C12-C13-C14	3.67	124.58	118.94
2	D	302	RET	C19-C9-C10	-3.62	117.85	122.92
5	C	308	K3I	C21-C15-C14	3.54	130.74	123.47
5	A	306	K3I	C21-C15-C14	3.53	130.70	123.47
5	A	306	K3I	C25-C23-C22	-3.53	113.53	118.94
5	B	306	K3I	C30-C28-C29	3.48	123.55	118.08
2	A	301	RET	C19-C9-C10	-3.32	118.28	122.92
5	B	306	K3I	C30-C28-C27	-3.31	118.29	122.92
5	E	307	K3I	C21-C15-C14	3.10	129.83	123.47
5	B	306	K3I	C20-C13-C14	-3.09	118.59	122.92
5	A	306	K3I	C20-C13-C14	-3.02	118.70	122.92
2	A	301	RET	C19-C9-C8	2.99	122.79	118.08
5	C	308	K3I	C20-C13-C14	-2.96	118.78	122.92
2	C	303	RET	C19-C9-C8	2.89	122.63	118.08
5	B	306	K3I	C3-C4-C5	-2.80	106.27	111.85
5	C	308	K3I	C30-C28-C27	-2.73	119.10	122.92
2	D	302	RET	C2-C1-C6	2.71	114.65	110.48
2	D	302	RET	C11-C10-C9	2.67	131.12	127.31
5	B	306	K3I	C12-C13-C14	2.67	123.03	118.94
5	B	306	K3I	C31-C29-C28	2.64	130.22	126.23
2	E	302	RET	C19-C9-C8	2.62	122.20	118.08
5	E	307	K3I	C15-C21-C22	2.61	128.81	123.47
5	C	308	K3I	C30-C28-C29	2.58	122.14	118.08

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	302	RET	C11-C10-C9	2.55	130.95	127.31
2	A	301	RET	C18-C5-C6	2.51	127.35	124.53
2	B	302	RET	C18-C5-C4	-2.44	108.92	113.62
5	B	306	K3I	C15-C21-C22	-2.43	118.51	123.47
2	A	301	RET	C1-C6-C7	-2.41	108.97	115.78
2	C	303	RET	C11-C10-C9	2.41	130.74	127.31
2	B	302	RET	C19-C9-C8	2.38	121.83	118.08
5	D	307	K3I	C25-C23-C22	-2.35	115.34	118.94
2	A	301	RET	C11-C10-C9	2.33	130.64	127.31
2	B	302	RET	C1-C6-C5	-2.25	119.44	122.61
5	A	306	K3I	C30-C28-C27	-2.20	119.85	122.92
5	E	307	K3I	C20-C13-C14	-2.19	119.86	122.92
2	B	302	RET	C18-C5-C6	2.18	126.98	124.53
2	D	302	RET	C1-C6-C7	2.16	121.88	115.78
5	E	307	K3I	C37-C32-C33	-2.13	119.61	122.61
2	B	302	RET	C2-C1-C6	2.12	113.75	110.48
5	A	306	K3I	C30-C28-C29	2.11	121.41	118.08
5	C	308	K3I	C12-C13-C14	2.10	122.17	118.94
2	E	302	RET	C1-C6-C5	-2.09	119.66	122.61
5	B	306	K3I	C21-C22-C23	2.08	130.28	127.31
2	D	302	RET	C19-C9-C8	2.05	121.31	118.08
2	D	302	RET	C1-C6-C5	-2.03	119.75	122.61

There are no chirality outliers.

All (111) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	RET	C20-C13-C14-C15
2	B	302	RET	C20-C13-C14-C15
2	D	302	RET	C20-C13-C14-C15
5	B	306	K3I	C5-C6-C7-C8
5	D	307	K3I	C27-C28-C29-C31
5	D	307	K3I	C30-C28-C29-C31
5	E	307	K3I	C15-C21-C22-C23
5	E	307	K3I	C25-C26-C27-C28
2	A	301	RET	C7-C8-C9-C19
5	E	307	K3I	C24-C23-C25-C26
5	E	307	K3I	C30-C28-C29-C31
2	A	301	RET	C7-C8-C9-C10
5	E	307	K3I	C22-C23-C25-C26
5	E	307	K3I	C27-C28-C29-C31
3	E	304	R16	C35-C36-C37-C38

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Mol	Chain	Res	Type	Atoms
3	A	302	R16	C32-C33-C34-C35
3	D	305	R16	C34-C35-C36-C37
3	D	305	R16	C35-C36-C37-C38
3	C	302	R16	C35-C36-C37-C38
3	C	301	R16	C31-C32-C33-C34
3	D	303	R16	C34-C35-C36-C37
3	E	303	R16	C31-C32-C33-C34
4	A	304	D12	C5-C6-C7-C8
3	E	303	R16	C28-C29-C30-C31
4	A	307	D12	C3-C4-C5-C6
3	C	305	R16	C31-C32-C33-C34
3	E	304	R16	C34-C35-C36-C37
4	D	306	D12	C6-C7-C8-C9
4	C	307	D12	C4-C5-C6-C7
4	D	306	D12	C2-C3-C4-C5
3	B	303	R16	C34-C35-C36-C37
4	E	305	D12	C6-C7-C8-C9
2	A	301	RET	C5-C6-C7-C8
5	B	306	K3I	C1-C6-C7-C8
5	D	307	K3I	C29-C31-C32-C33
5	D	307	K3I	C29-C31-C32-C37
5	E	307	K3I	C1-C6-C7-C8
5	E	307	K3I	C5-C6-C7-C8
3	C	302	R16	C32-C33-C34-C35
4	C	306	D12	C11-C10-C9-C8
5	D	307	K3I	C23-C25-C26-C27
3	A	303	R16	C36-C37-C38-C39
3	D	301	R16	C30-C31-C32-C33
3	E	303	R16	C36-C37-C38-C39
4	A	305	D12	C2-C3-C4-C5
3	A	303	R16	C39-C40-C41-C42
4	C	306	D12	C9-C10-C11-C12
3	C	305	R16	C32-C33-C34-C35
5	E	307	K3I	C26-C27-C28-C30
4	B	305	D12	C1-C2-C3-C4
3	E	303	R16	C32-C33-C34-C35
5	E	307	K3I	C26-C27-C28-C29
3	C	302	R16	C31-C32-C33-C34
2	A	301	RET	C12-C13-C14-C15
2	B	302	RET	C12-C13-C14-C15
2	C	303	RET	C12-C13-C14-C15
2	D	302	RET	C12-C13-C14-C15

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Mol	Chain	Res	Type	Atoms
4	C	306	D12	C7-C8-C9-C10
3	E	303	R16	C27-C28-C29-C30
3	B	303	R16	C35-C36-C37-C38
3	E	304	R16	C36-C37-C38-C39
4	B	301	D12	C9-C10-C11-C12
3	A	302	R16	C37-C38-C39-C40
2	C	303	RET	C20-C13-C14-C15
3	D	303	R16	C35-C36-C37-C38
3	E	304	R16	C27-C28-C29-C30
2	A	301	RET	C1-C6-C7-C8
4	E	306	D12	C3-C4-C5-C6
3	C	304	R16	C31-C32-C33-C34
4	A	307	D12	C2-C3-C4-C5
3	D	304	R16	C35-C36-C37-C38
5	B	306	K3I	C21-C22-C23-C25
4	E	301	D12	C9-C10-C11-C12
3	C	301	R16	C38-C39-C40-C41
4	E	301	D12	C5-C6-C7-C8
4	C	307	D12	C5-C6-C7-C8
3	C	305	R16	C33-C34-C35-C36
5	A	306	K3I	C23-C25-C26-C27
4	D	306	D12	C3-C4-C5-C6
3	A	302	R16	C29-C30-C31-C32
5	A	306	K3I	C1-C6-C7-C8
3	C	305	R16	C30-C31-C32-C33
5	B	306	K3I	C12-C13-C14-C15
4	B	305	D12	C2-C3-C4-C5
3	E	304	R16	C32-C33-C34-C35
5	A	306	K3I	C26-C27-C28-C30
5	B	306	K3I	C20-C13-C14-C15
5	B	306	K3I	C21-C22-C23-C24
5	B	306	K3I	C26-C27-C28-C30
4	C	306	D12	C4-C5-C6-C7
4	B	304	D12	C1-C2-C3-C4
5	A	306	K3I	C26-C27-C28-C29
3	D	305	R16	C37-C38-C39-C40
3	E	303	R16	C38-C39-C40-C41
5	A	306	K3I	C5-C6-C7-C8
5	A	306	K3I	C29-C31-C32-C37
3	D	304	R16	C37-C38-C39-C40
4	A	304	D12	C1-C2-C3-C4
4	A	304	D12	C11-C10-C9-C8

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Mol	Chain	Res	Type	Atoms
3	C	304	R16	C38-C39-C40-C41
5	C	308	K3I	C22-C23-C25-C26
5	B	306	K3I	C26-C27-C28-C29
3	D	304	R16	C38-C39-C40-C41
3	A	302	R16	C31-C32-C33-C34
3	D	305	R16	C31-C32-C33-C34
3	D	304	R16	C36-C37-C38-C39
3	E	303	R16	C30-C31-C32-C33
4	C	307	D12	C6-C7-C8-C9
4	A	304	D12	C6-C7-C8-C9
3	D	303	R16	C27-C28-C29-C30
3	D	303	R16	C32-C33-C34-C35

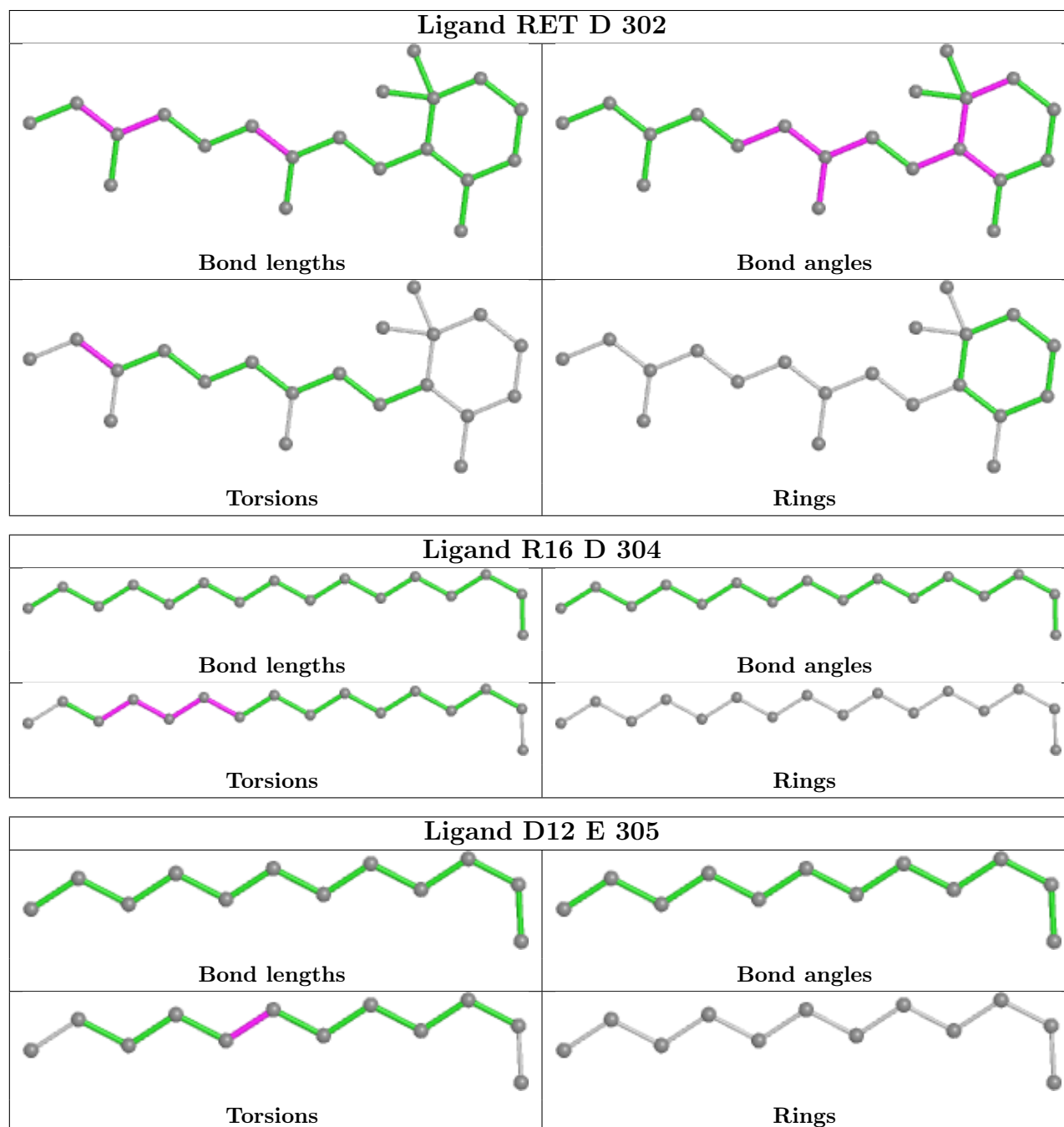
There are no ring outliers.

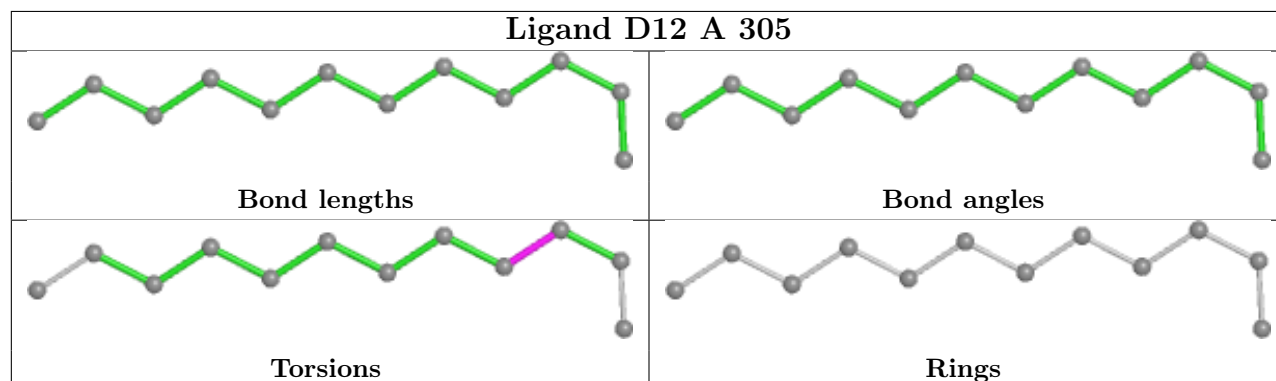
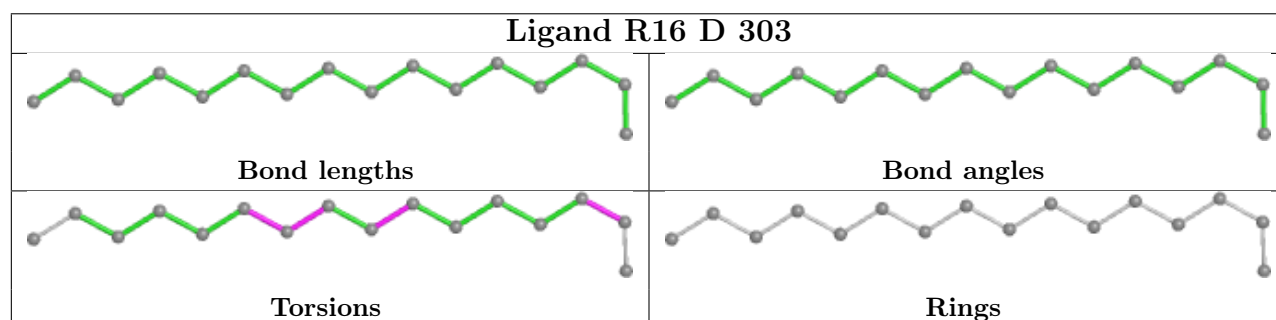
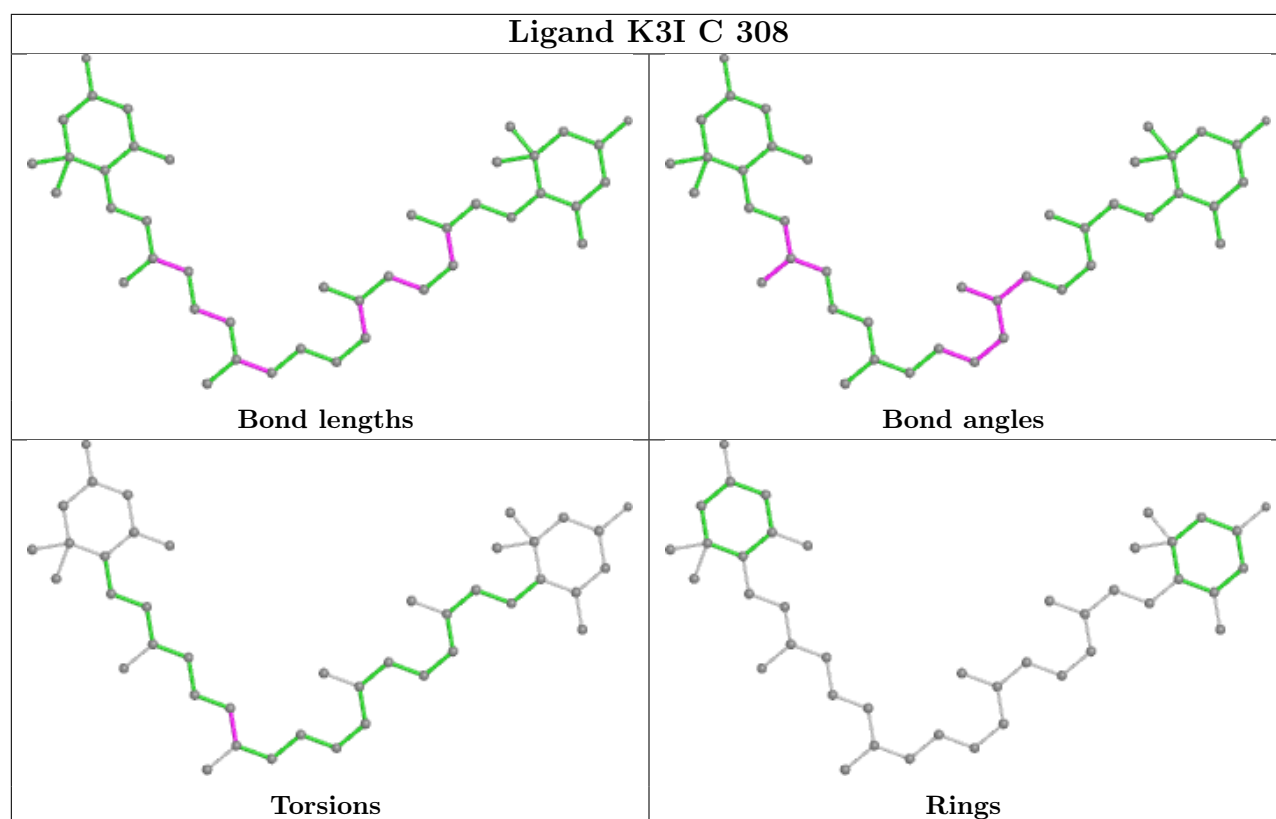
21 monomers are involved in 60 short contacts:

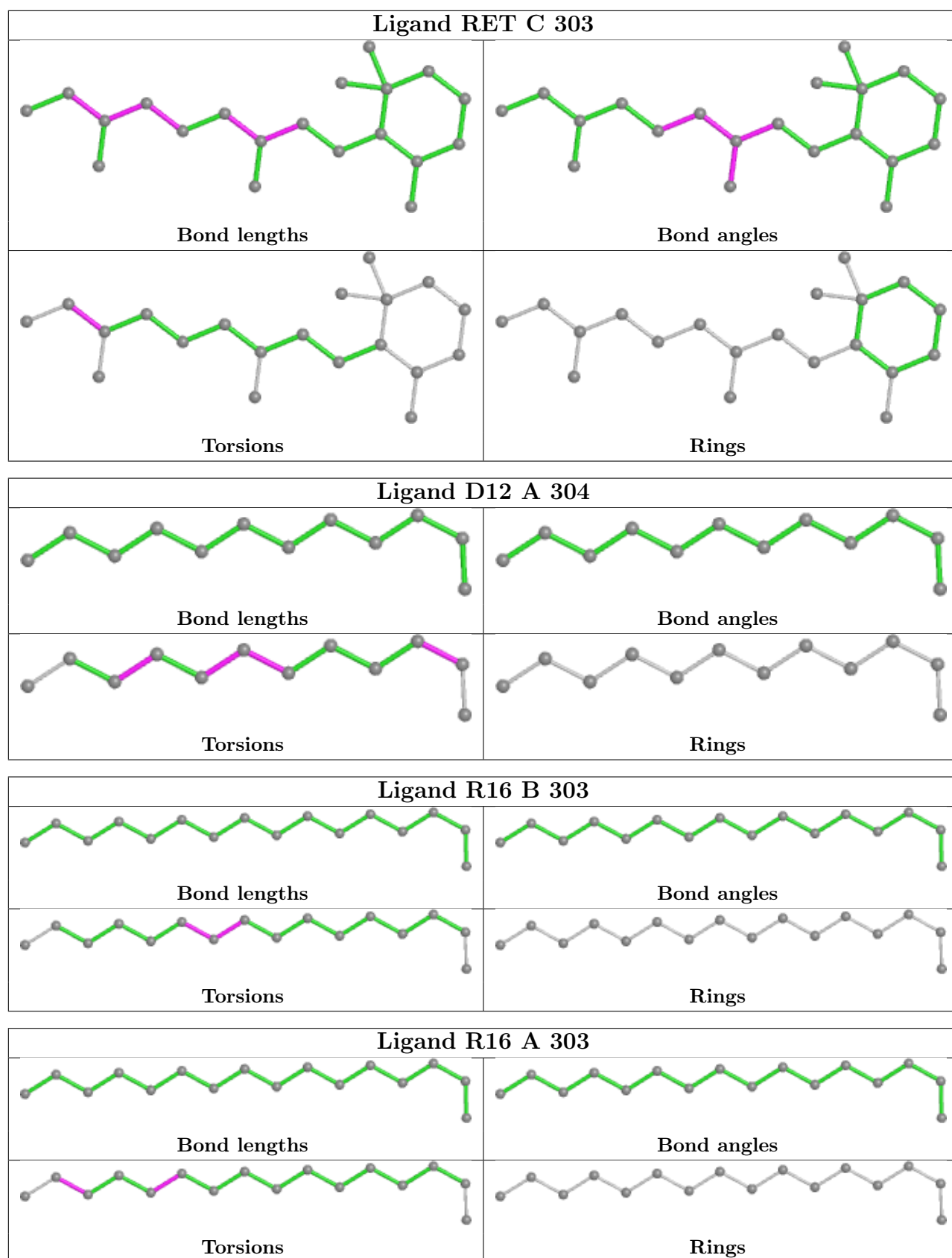
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	302	RET	7	0
4	E	305	D12	5	0
3	D	303	R16	1	0
2	C	303	RET	4	0
4	A	304	D12	1	0
3	B	303	R16	1	0
3	A	303	R16	3	0
3	E	303	R16	5	0
4	E	301	D12	1	0
4	D	306	D12	1	0
2	A	301	RET	1	0
3	C	302	R16	1	0
4	B	305	D12	1	0
3	C	305	R16	1	0
4	E	306	D12	1	0
4	A	307	D12	1	0
2	E	302	RET	15	0
3	C	304	R16	3	0
3	E	304	R16	2	0
4	C	306	D12	3	0
2	B	302	RET	7	0

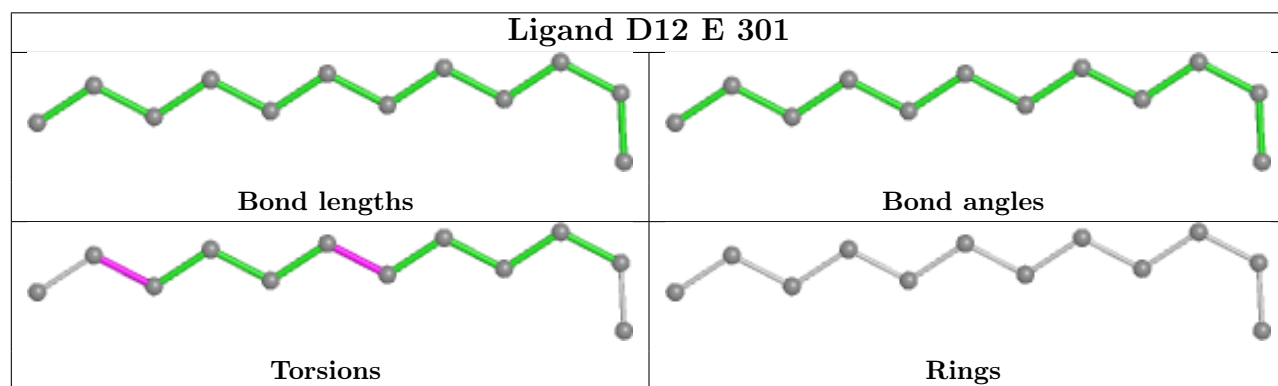
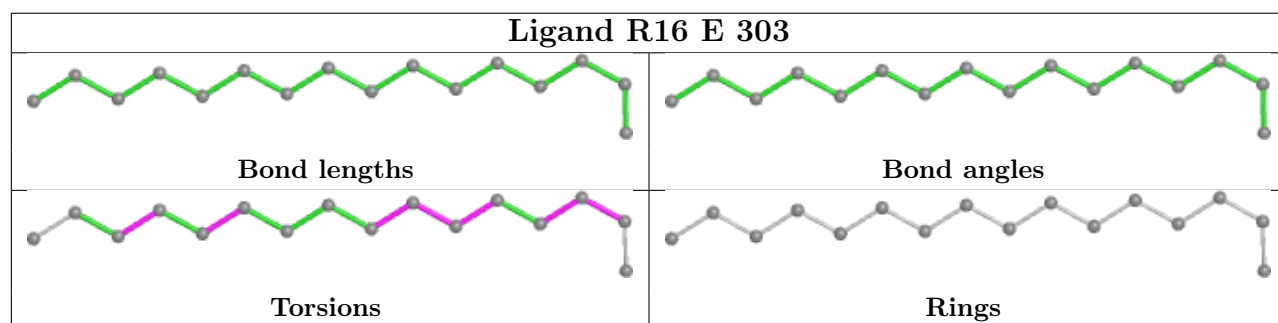
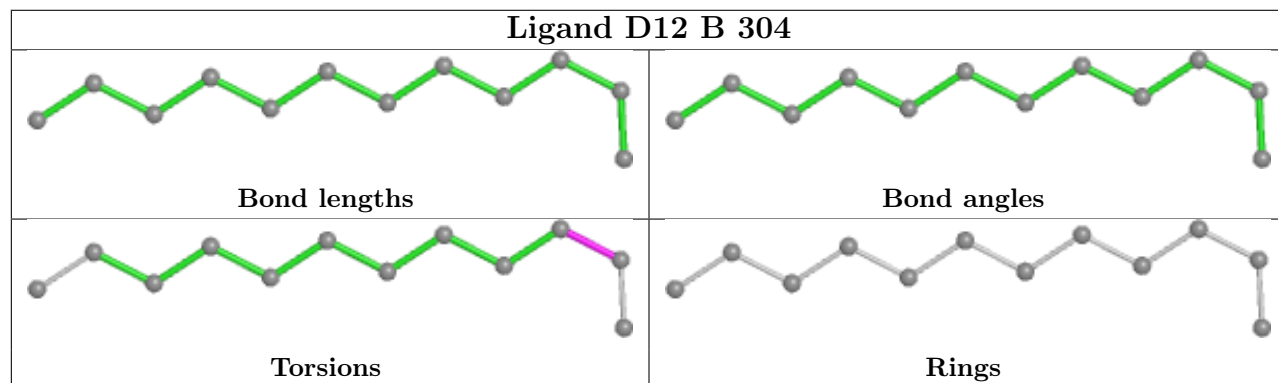
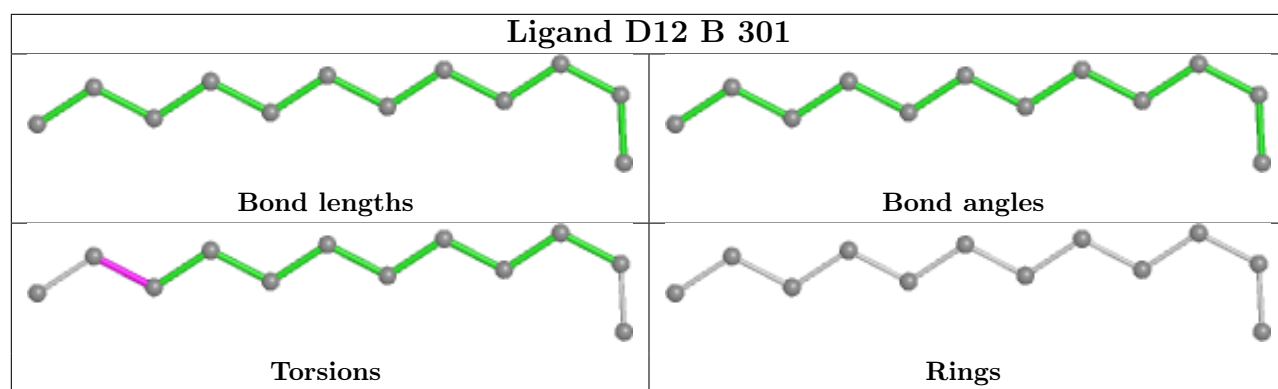
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

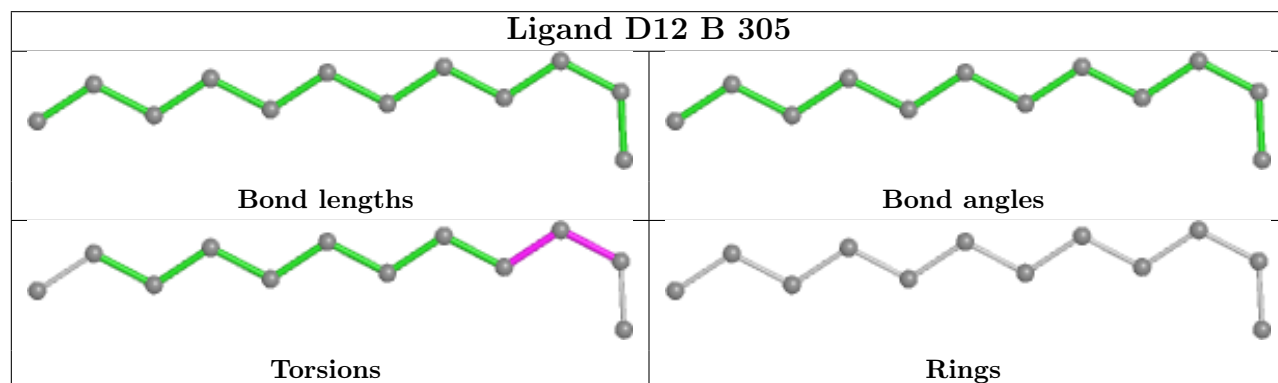
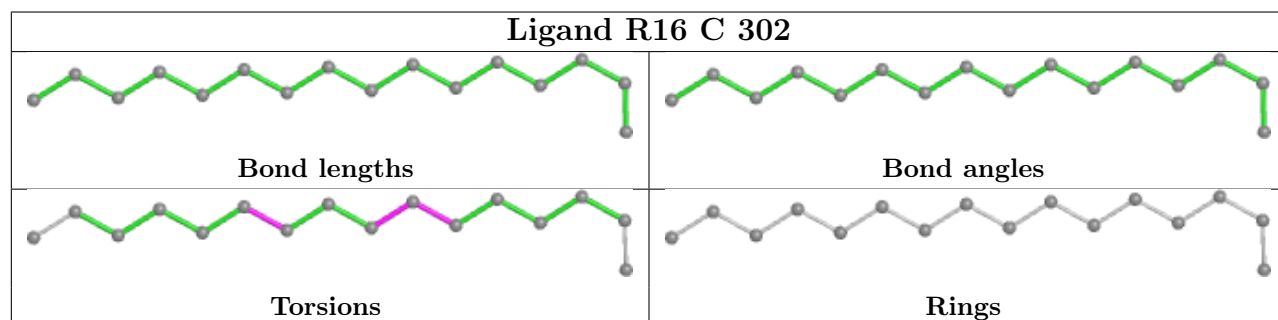
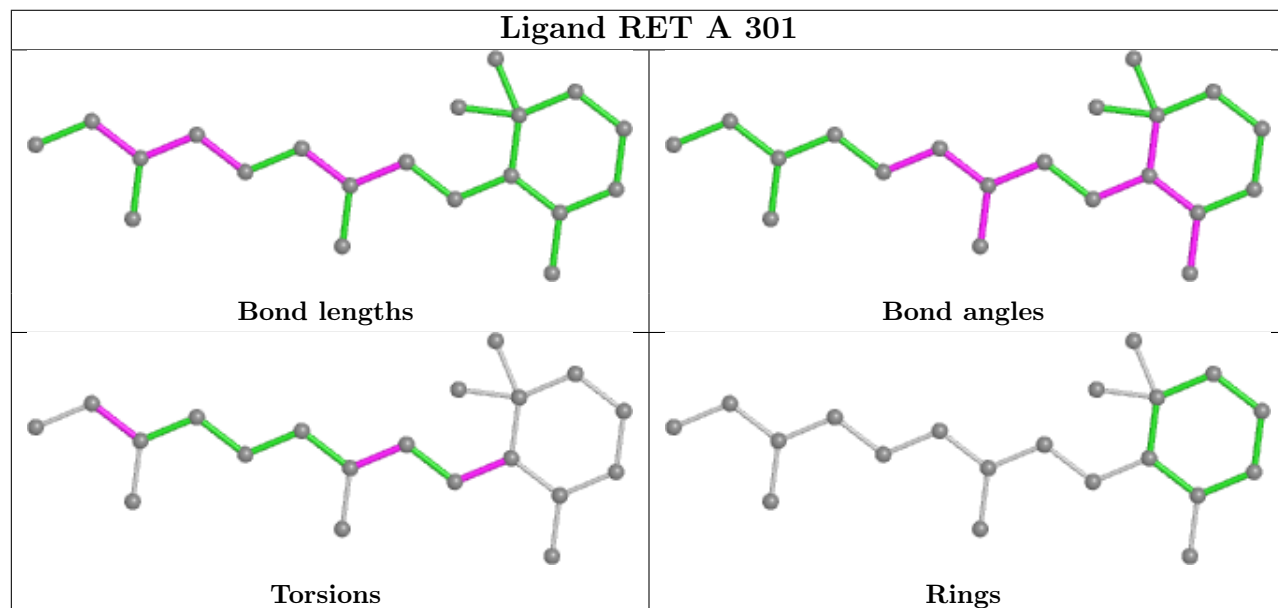
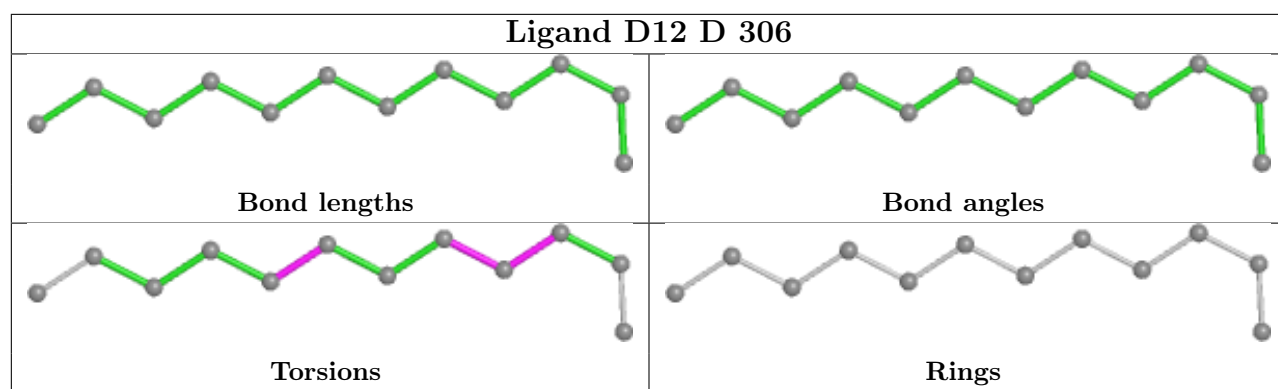
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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.

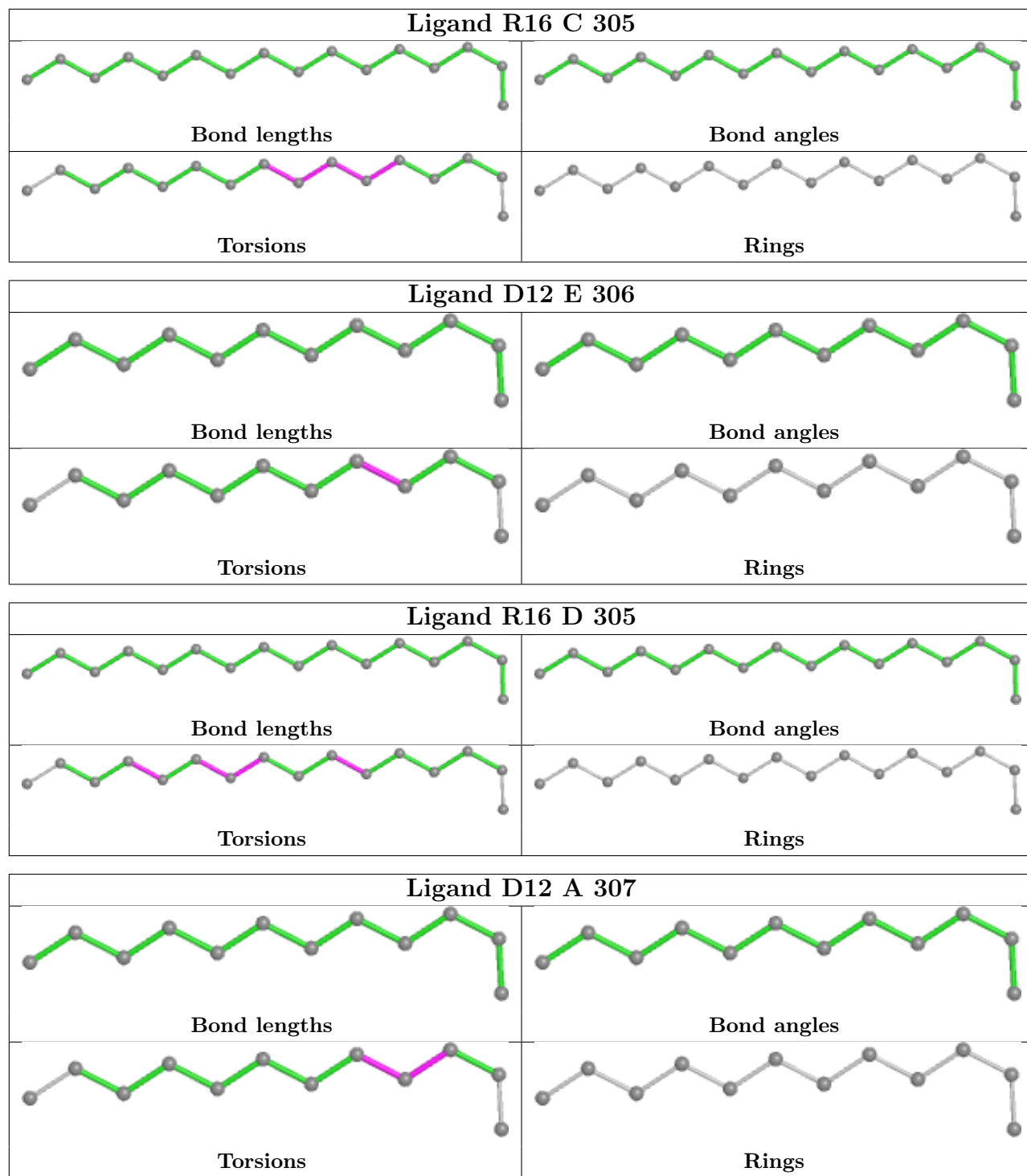




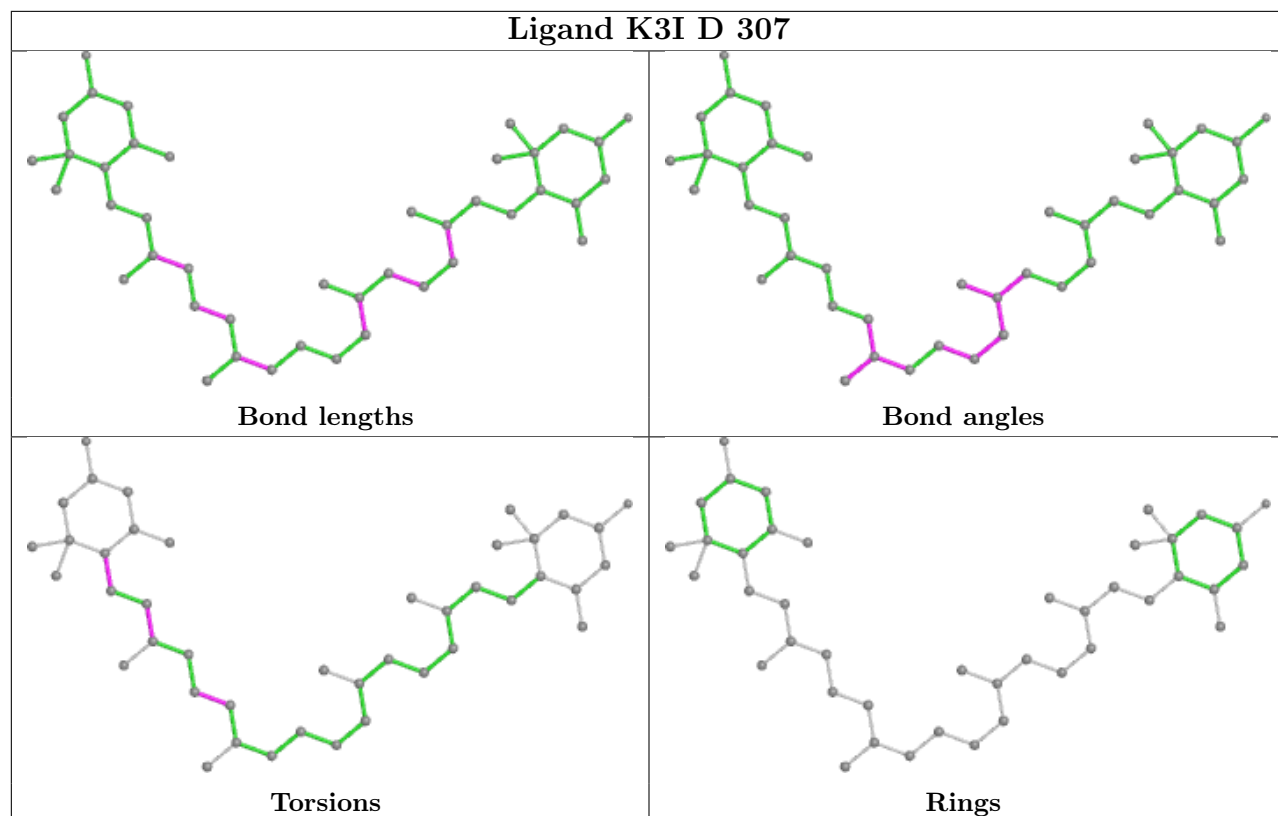




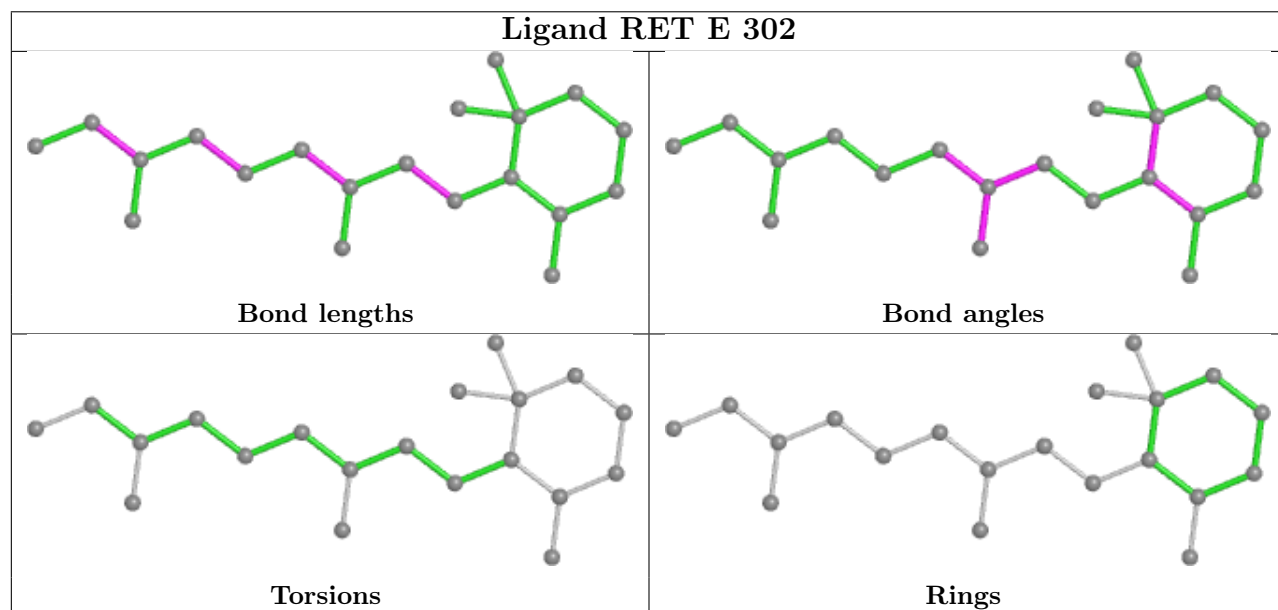




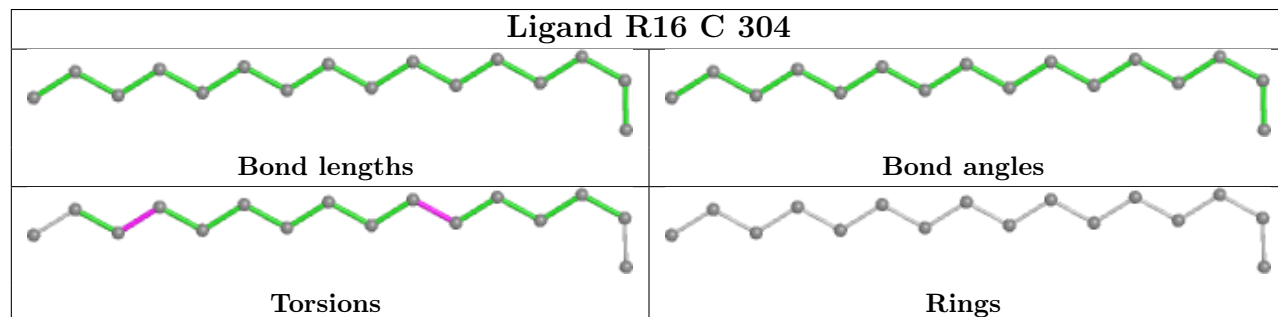
Ligand K3I D 307

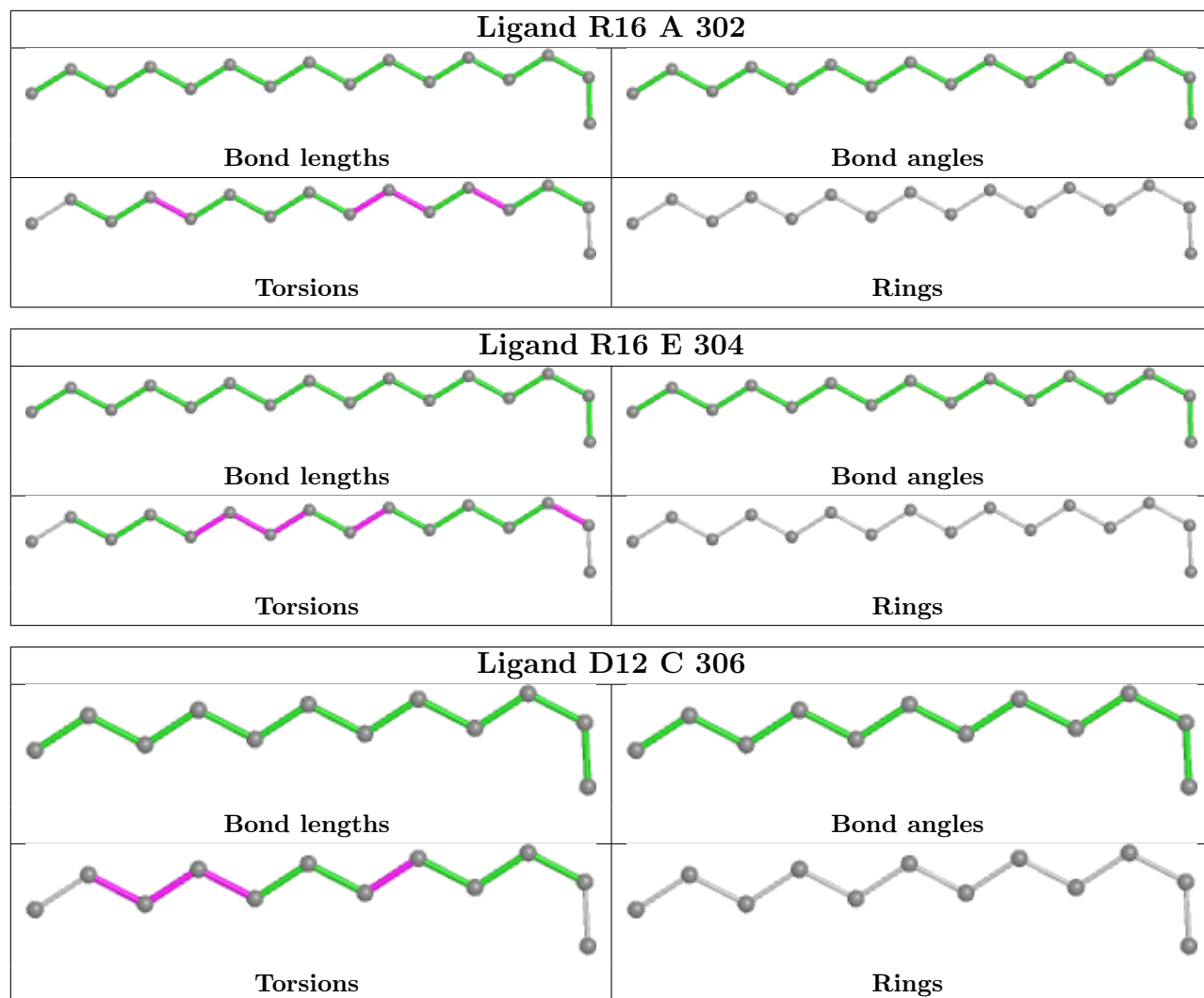


Ligand RET E 302

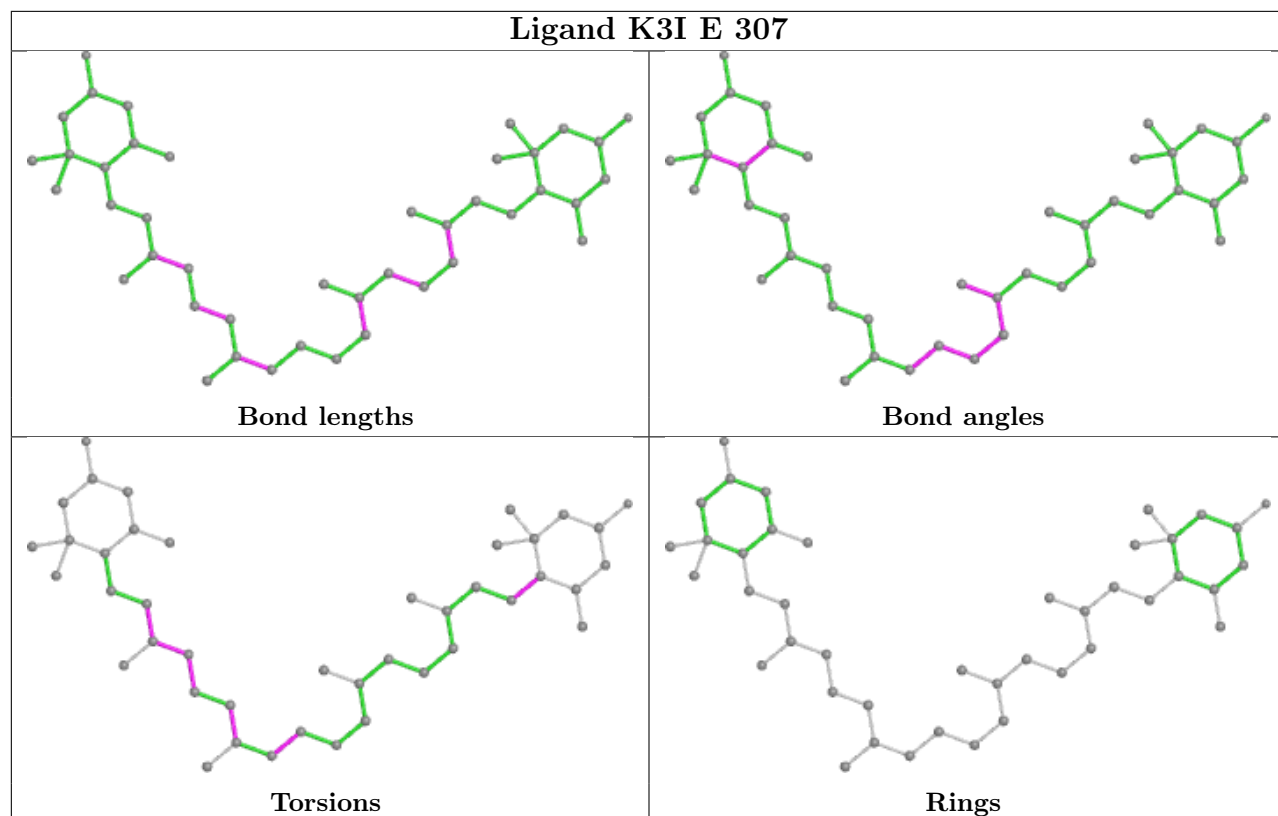


Ligand R16 C 304

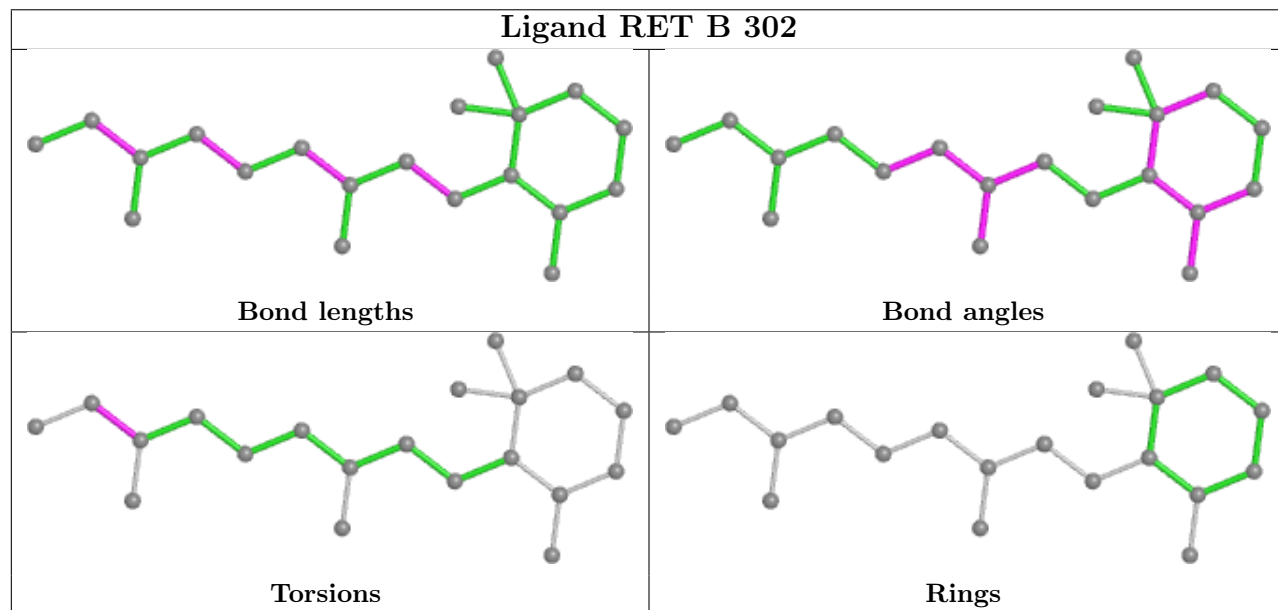


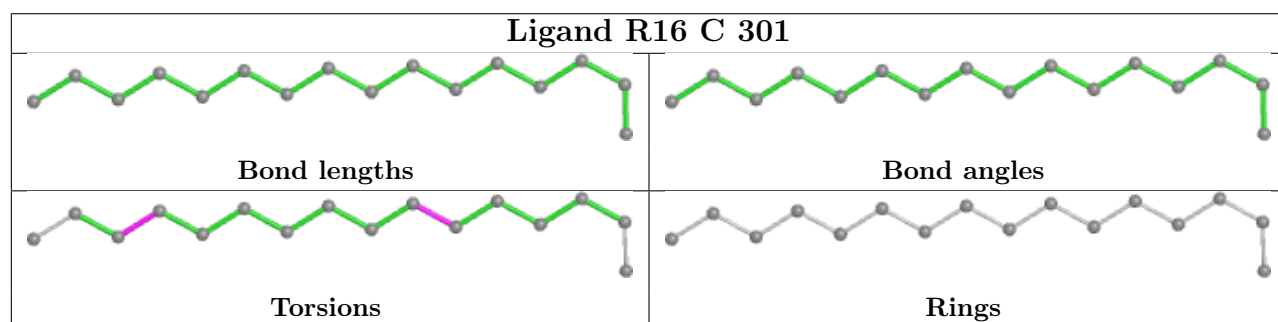
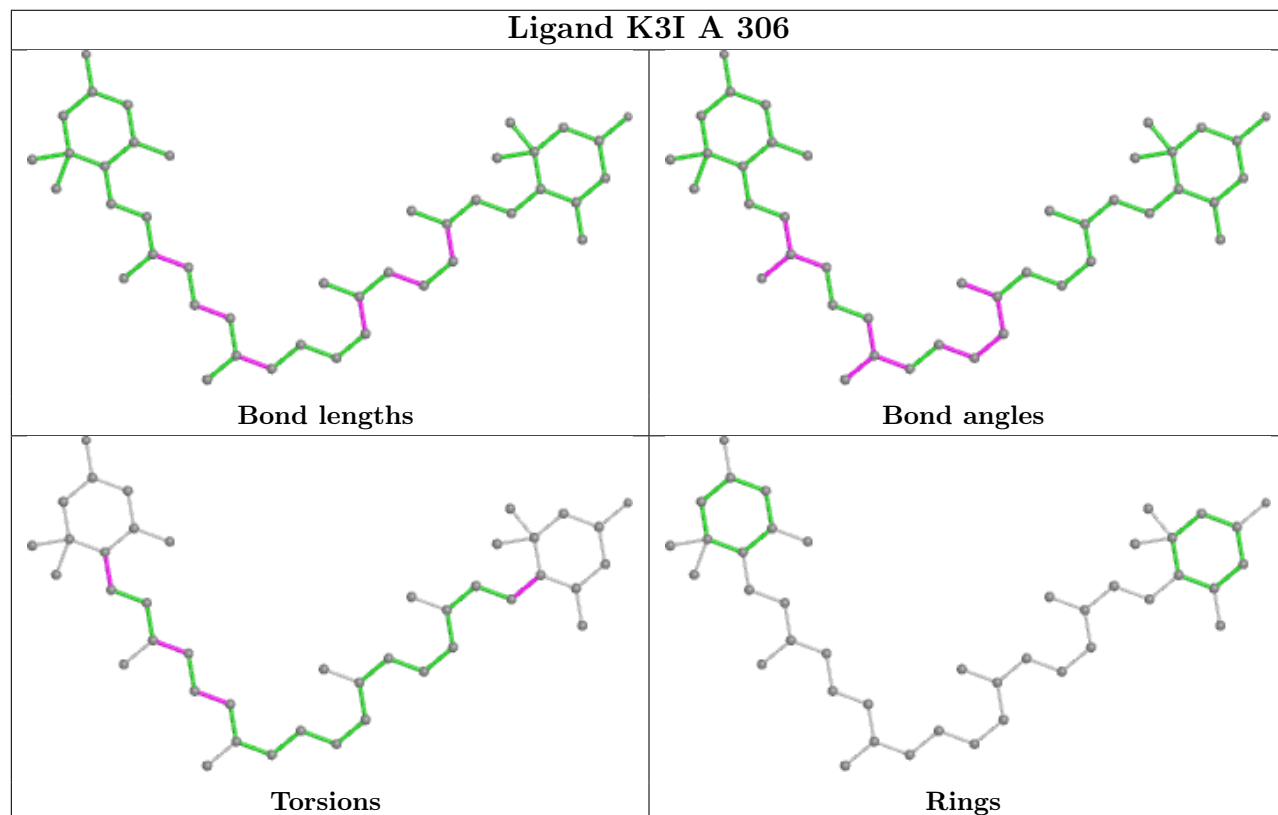
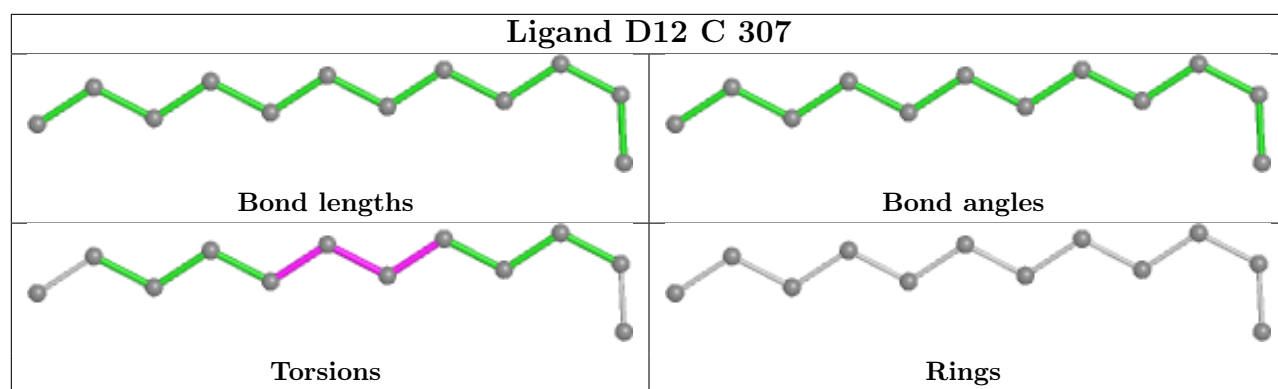


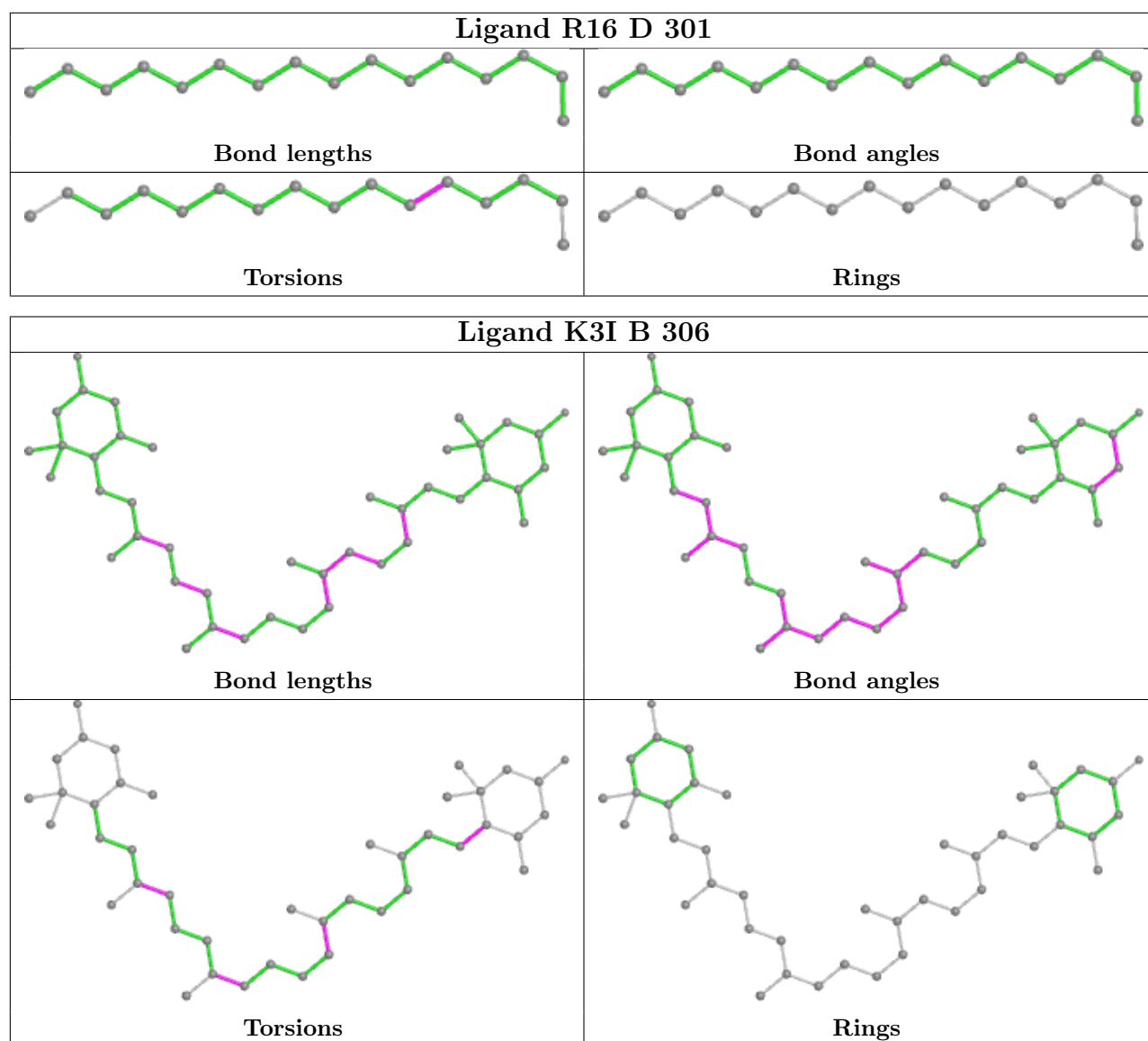
Ligand K3I E 307



Ligand RET B 302







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.

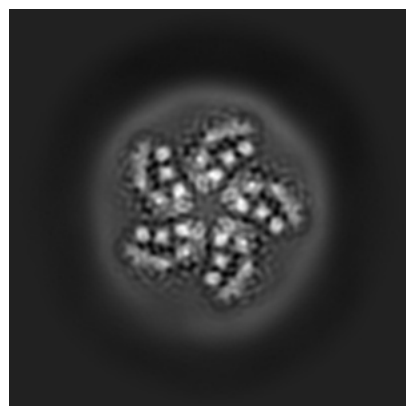
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-61685. These allow visual inspection of the internal detail of the map and identification of artifacts.

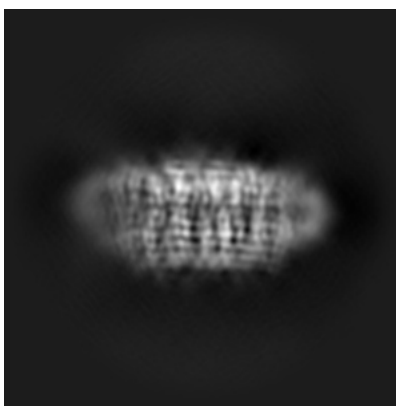
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

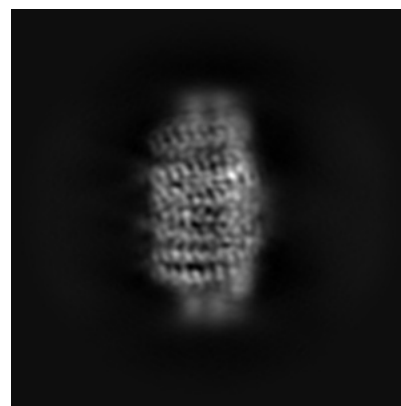
6.1.1 Primary map



X

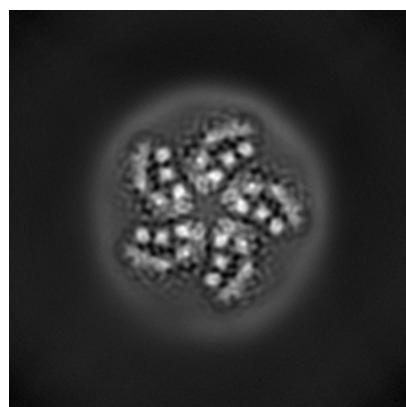


Y

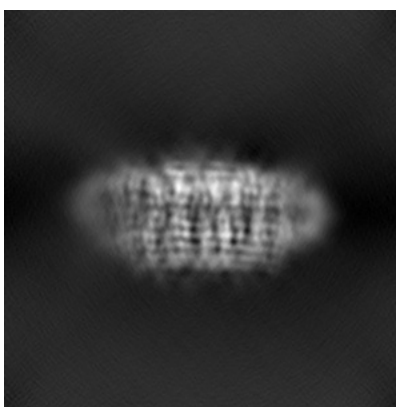


Z

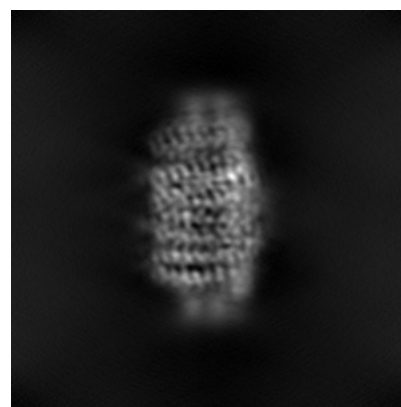
6.1.2 Raw map



X



Y

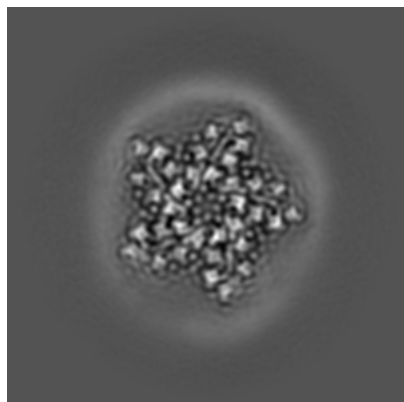


Z

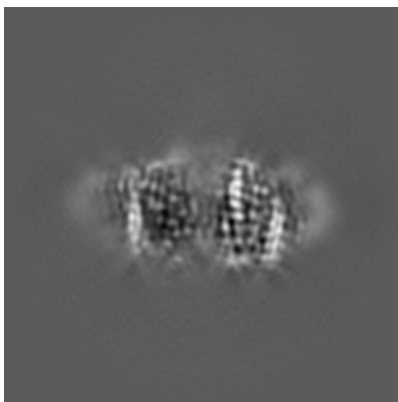
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

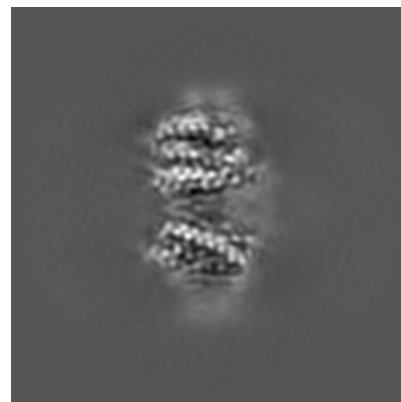
6.2.1 Primary map



X Index: 110

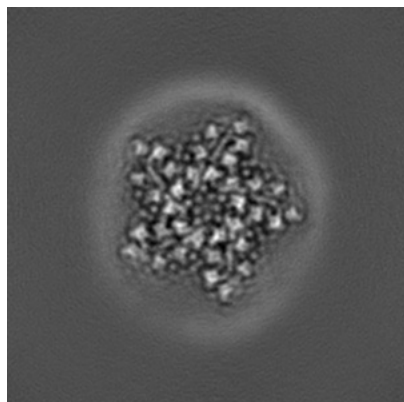


Y Index: 110

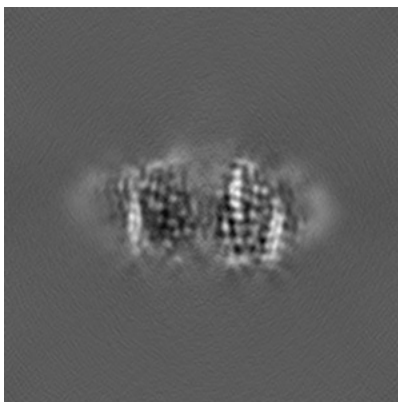


Z Index: 110

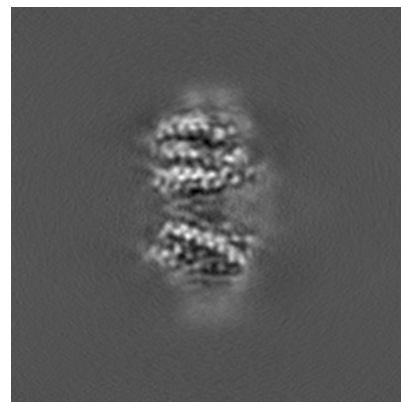
6.2.2 Raw map



X Index: 110



Y Index: 110

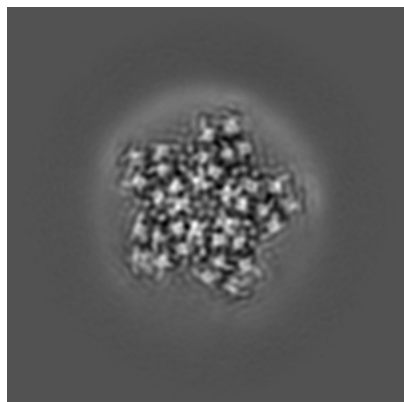


Z Index: 110

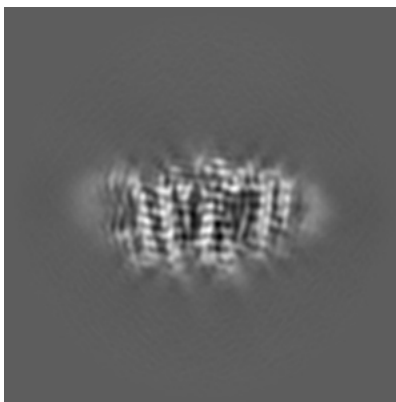
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

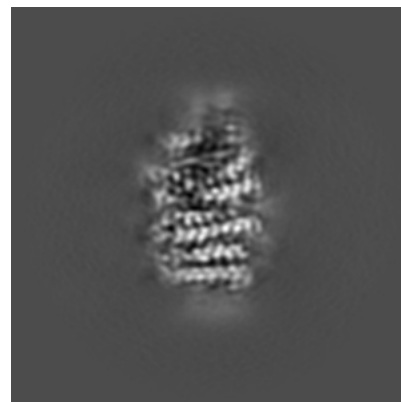
6.3.1 Primary map



X Index: 97

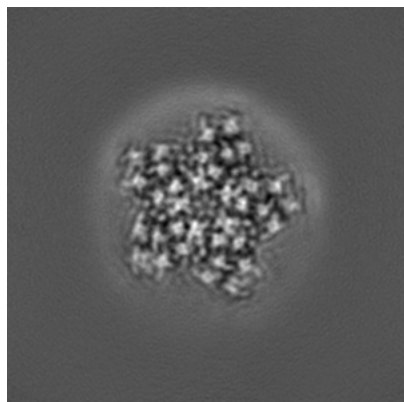


Y Index: 129

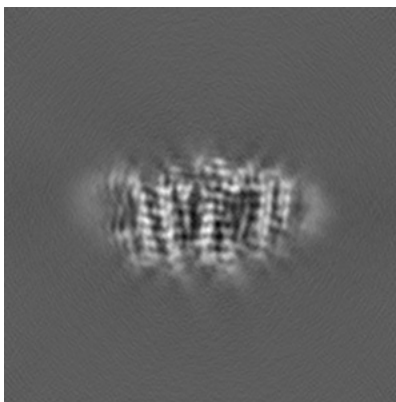


Z Index: 96

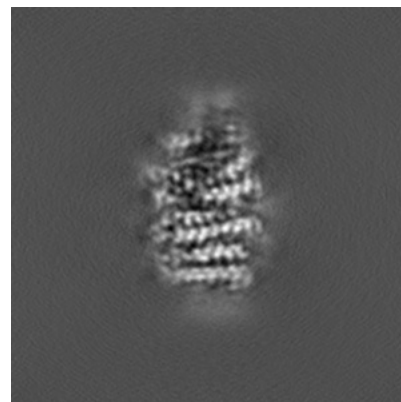
6.3.2 Raw map



X Index: 97



Y Index: 129

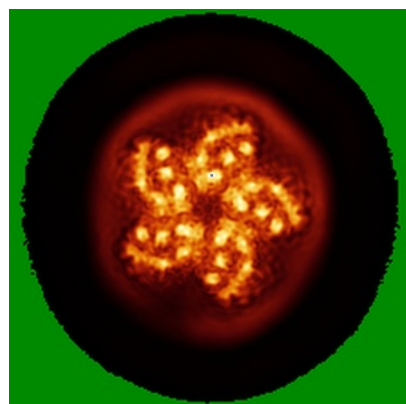


Z Index: 96

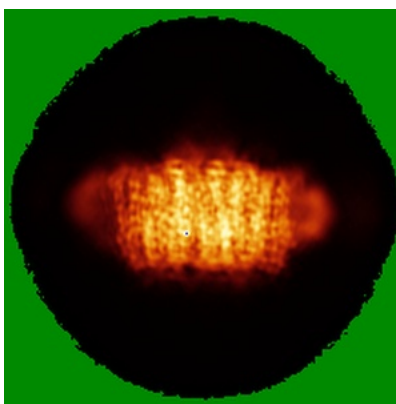
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) ⓘ

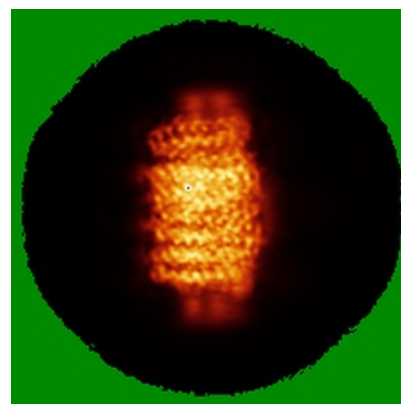
6.4.1 Primary map



X

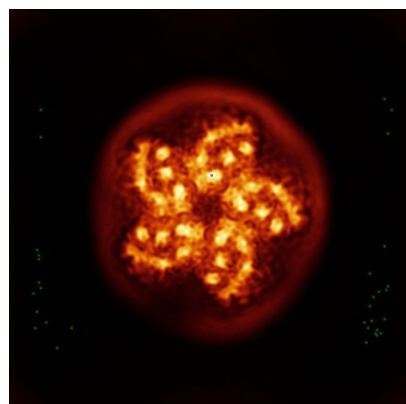


Y

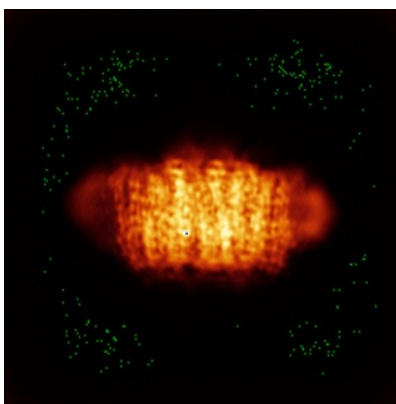


Z

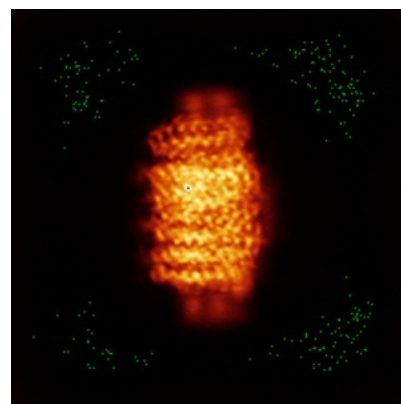
6.4.2 Raw map



X



Y

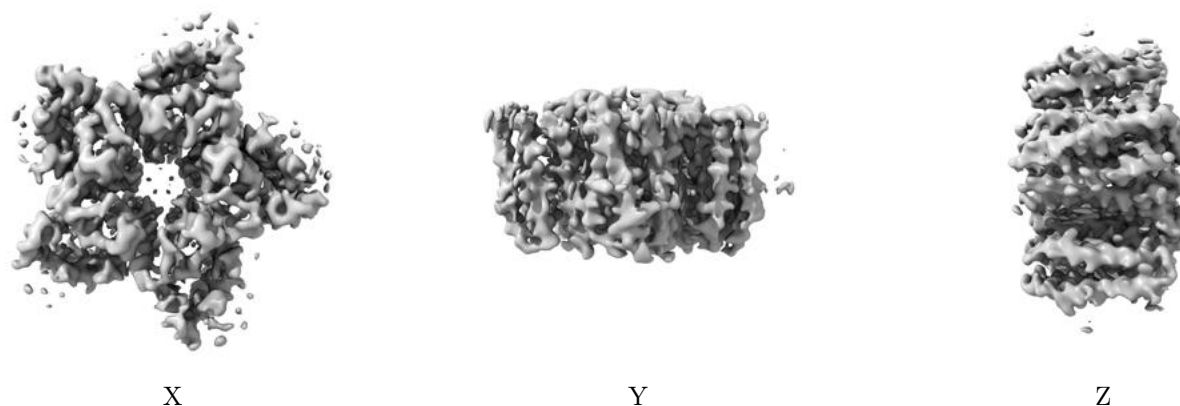


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

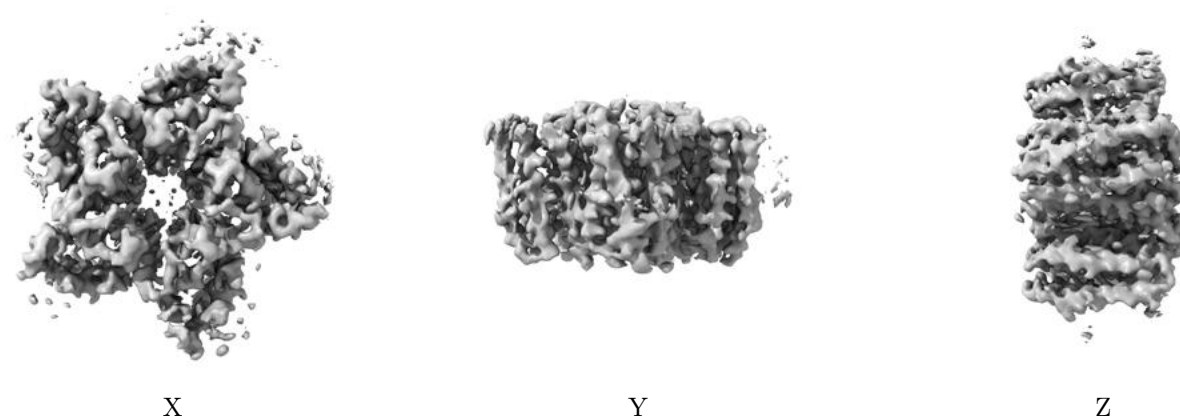
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.035. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

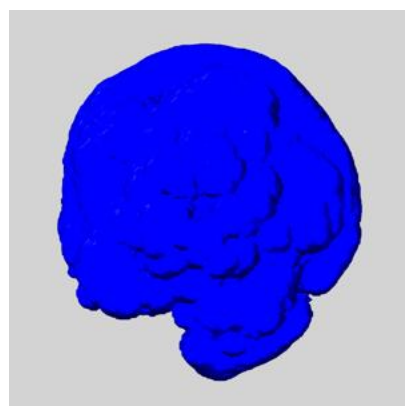
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

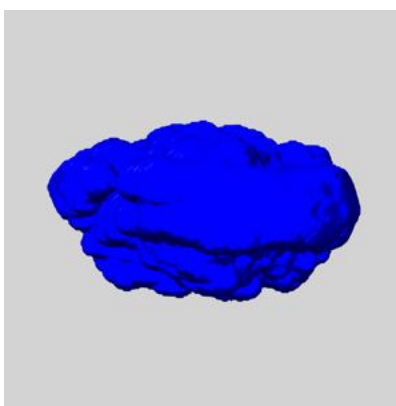
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

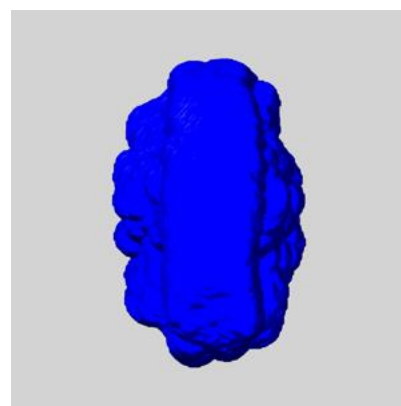
6.6.1 emd_61685_msk_1.map [i](#)



X



Y

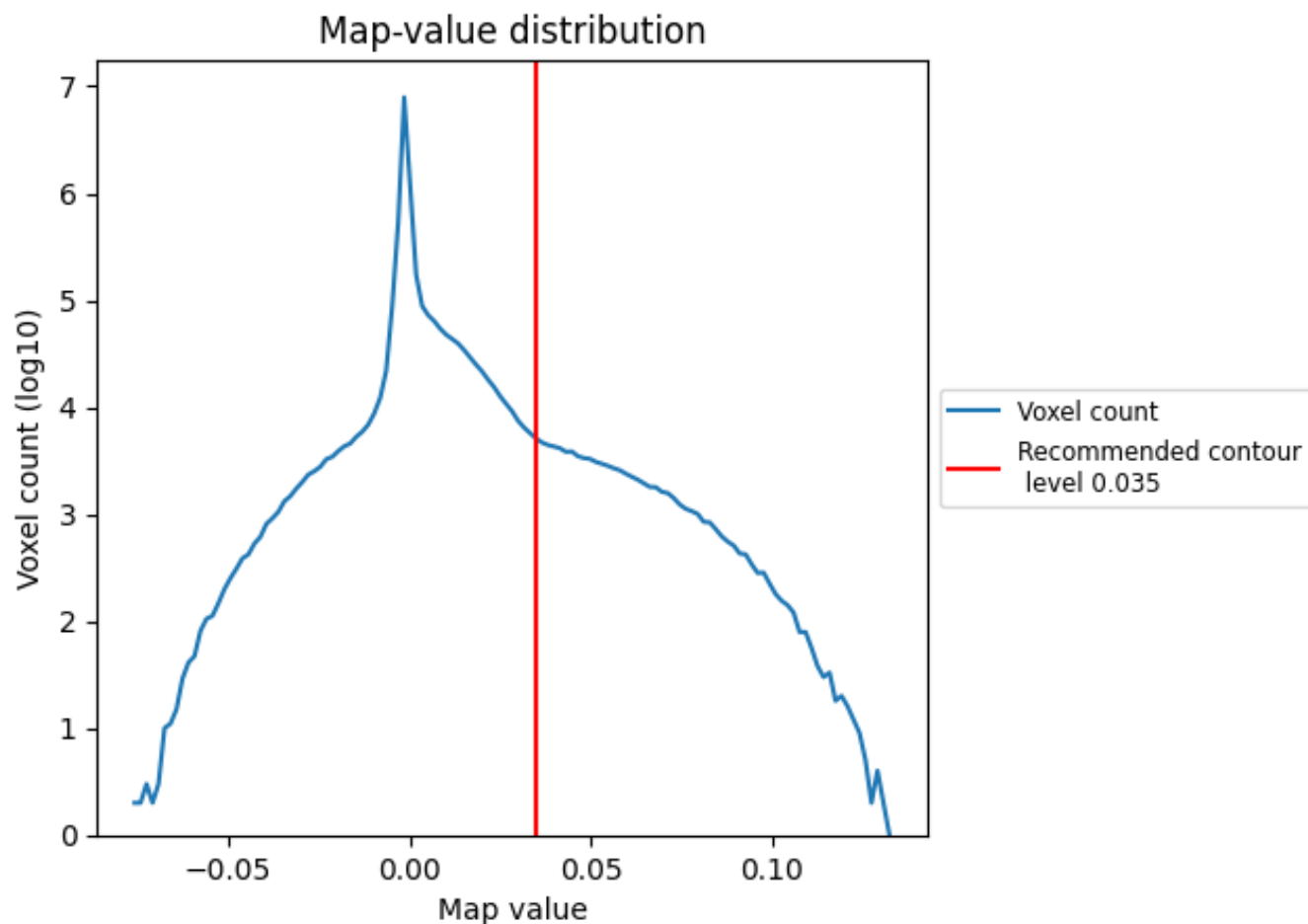


Z

7 Map analysis [i](#)

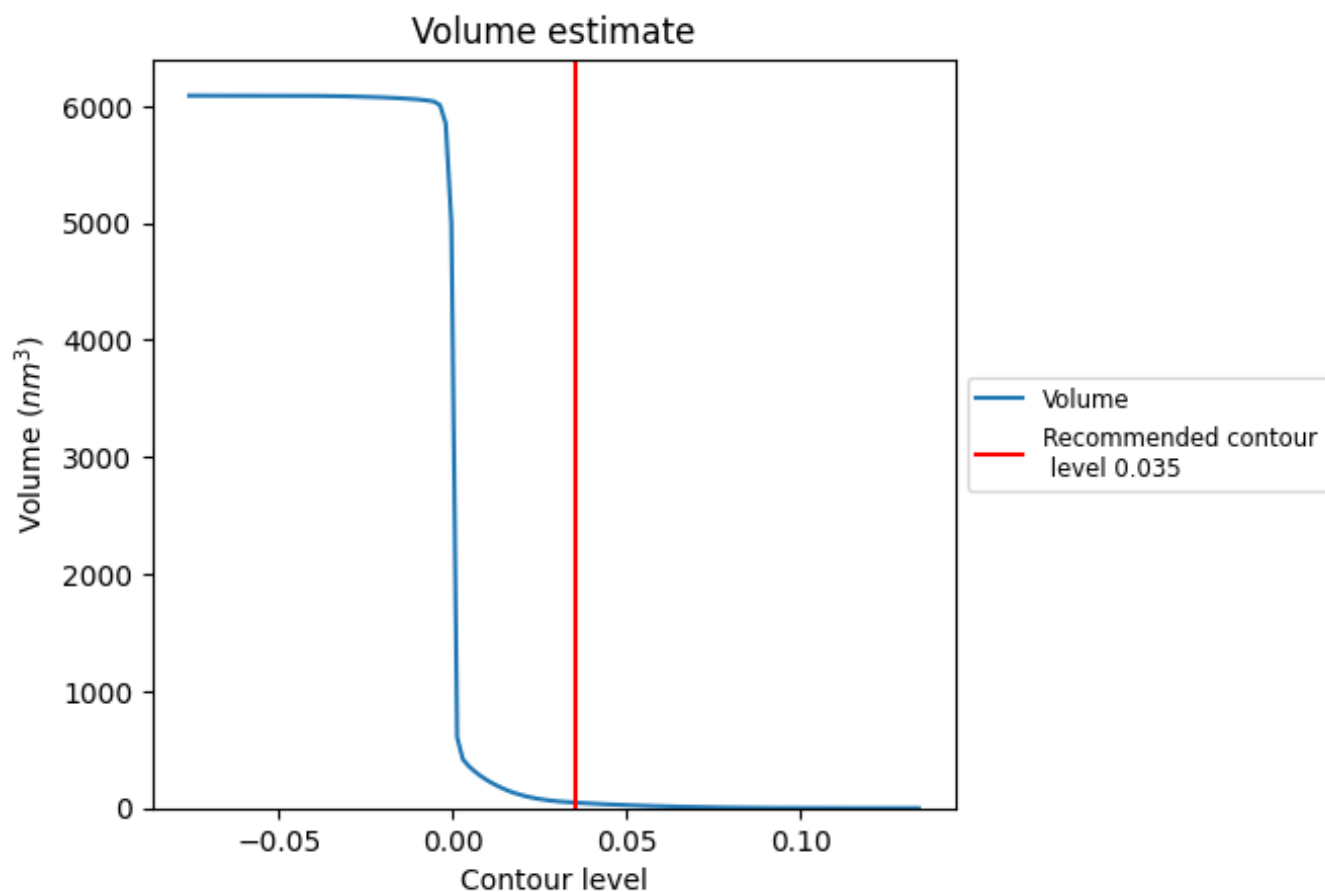
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

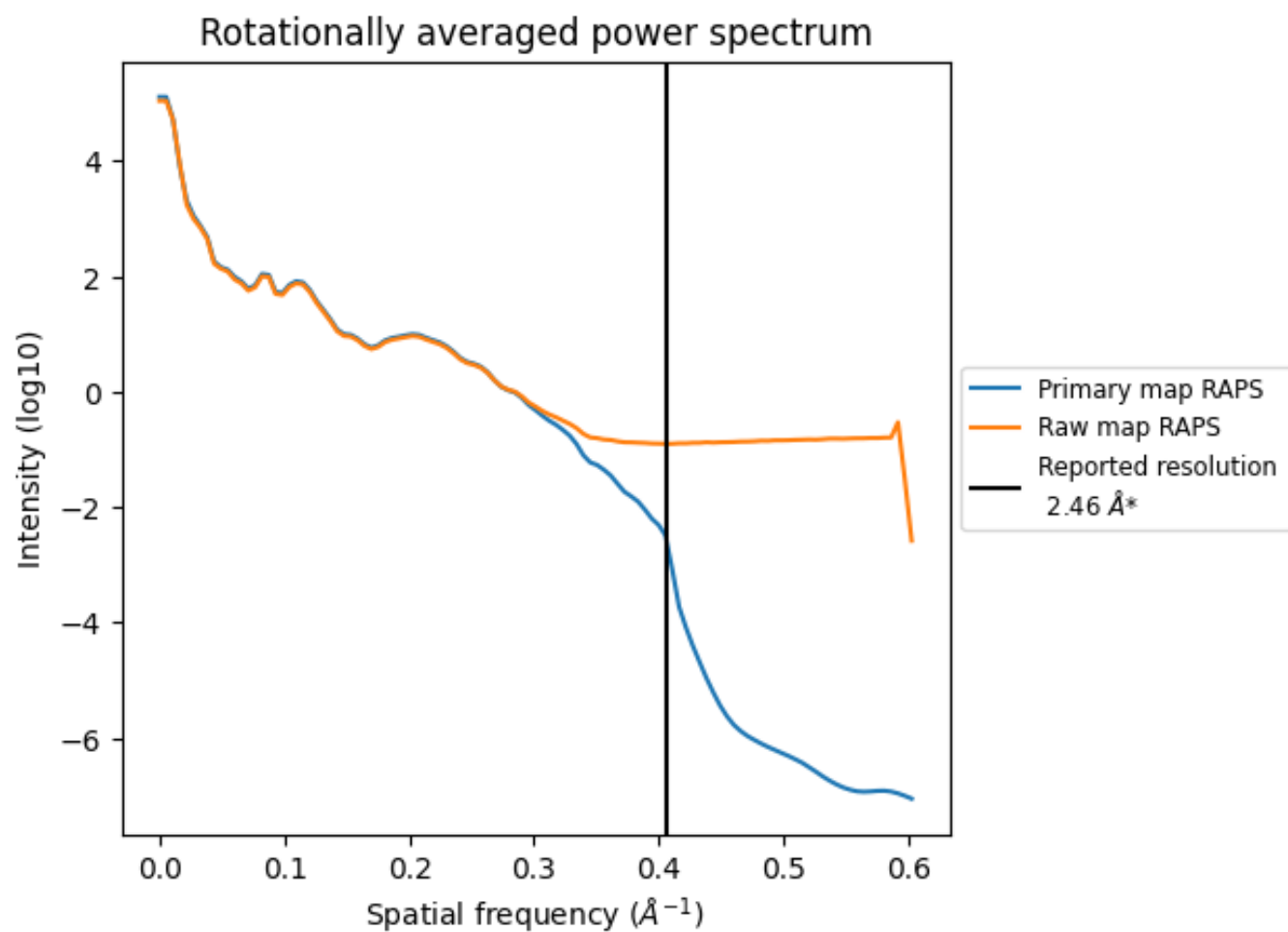
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 47 nm^3 ; this corresponds to an approximate mass of 43 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum ⓘ

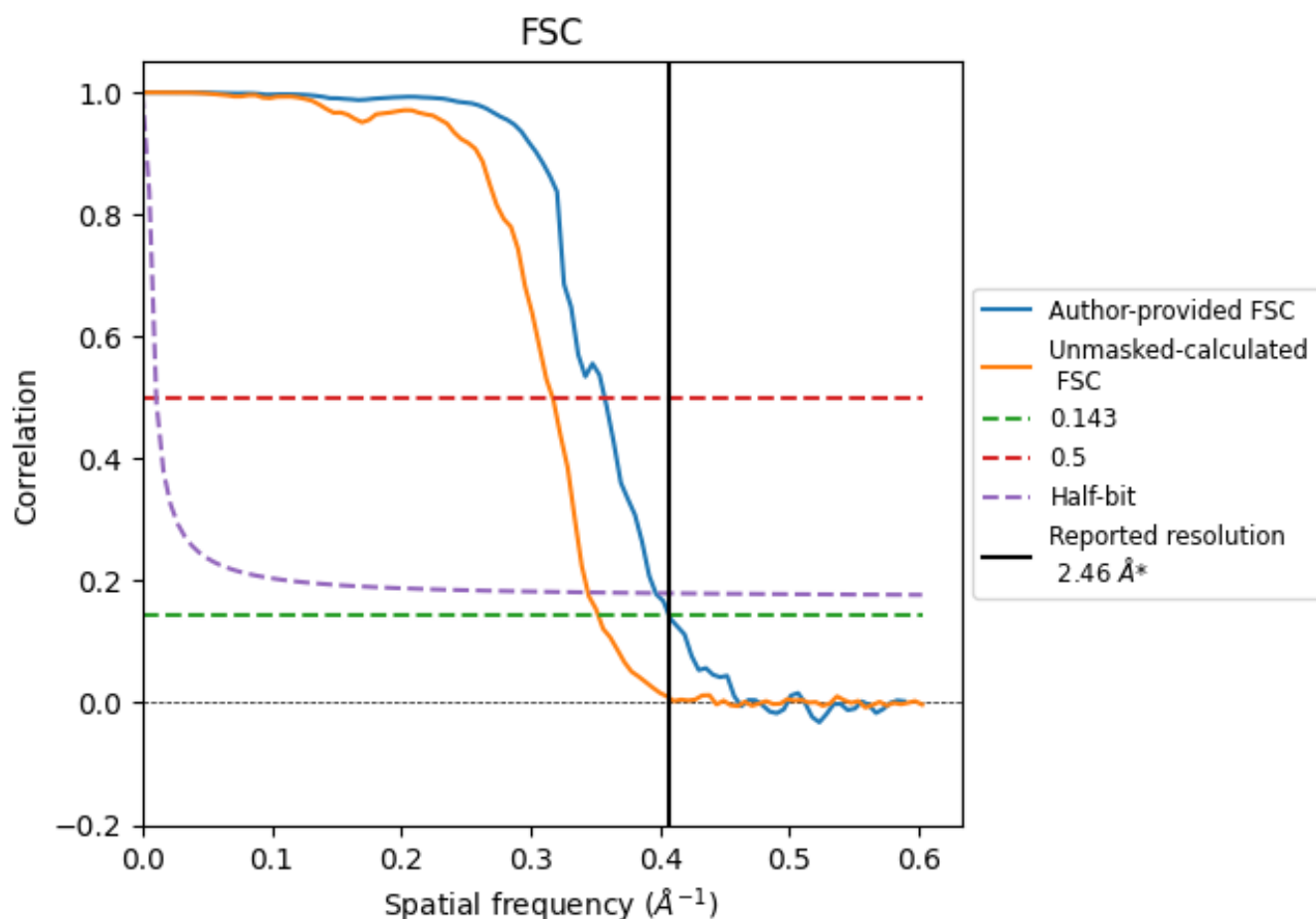


*Reported resolution corresponds to spatial frequency of 0.407 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.407 \AA^{-1}

8.2 Resolution estimates [i](#)

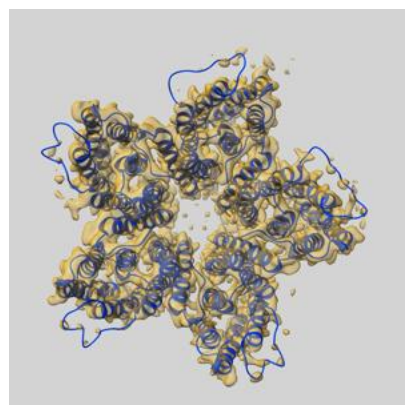
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.46	-	-
Author-provided FSC curve	2.46	2.80	2.52
Unmasked-calculated*	2.84	3.15	2.90

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 2.84 differs from the reported value 2.46 by more than 10 %

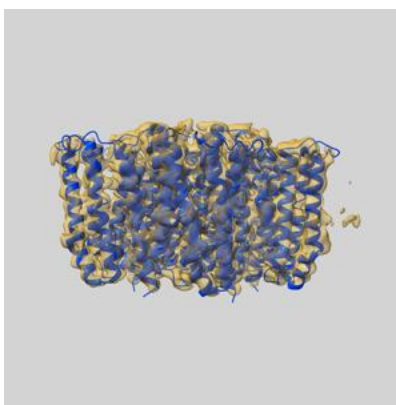
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-61685 and PDB model 9JOU. Per-residue inclusion information can be found in section 3 on page 10.

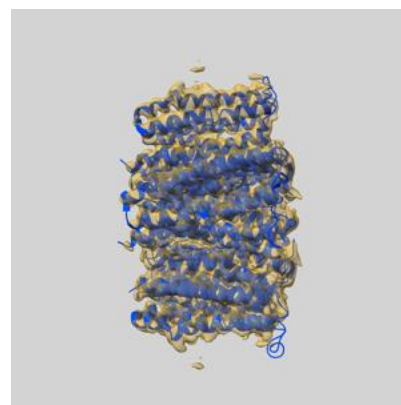
9.1 Map-model overlay [i](#)



X



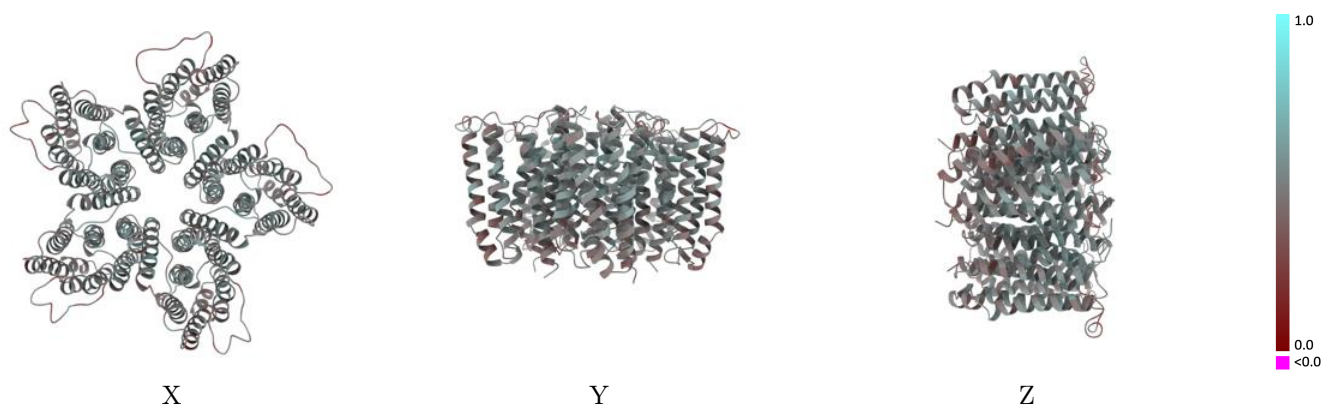
Y



Z

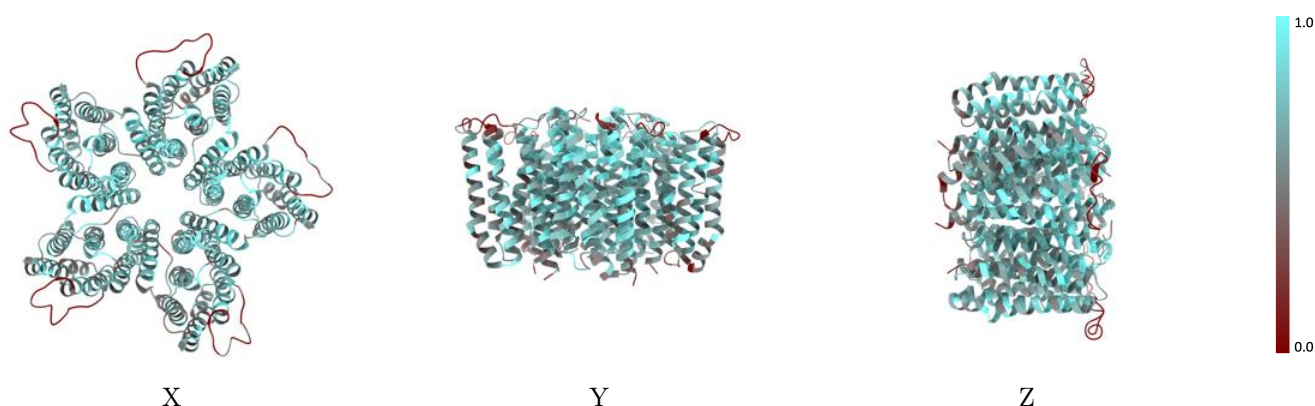
The images above show the 3D surface view of the map at the recommended contour level 0.035 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

9.2 Q-score mapped to coordinate model [i](#)



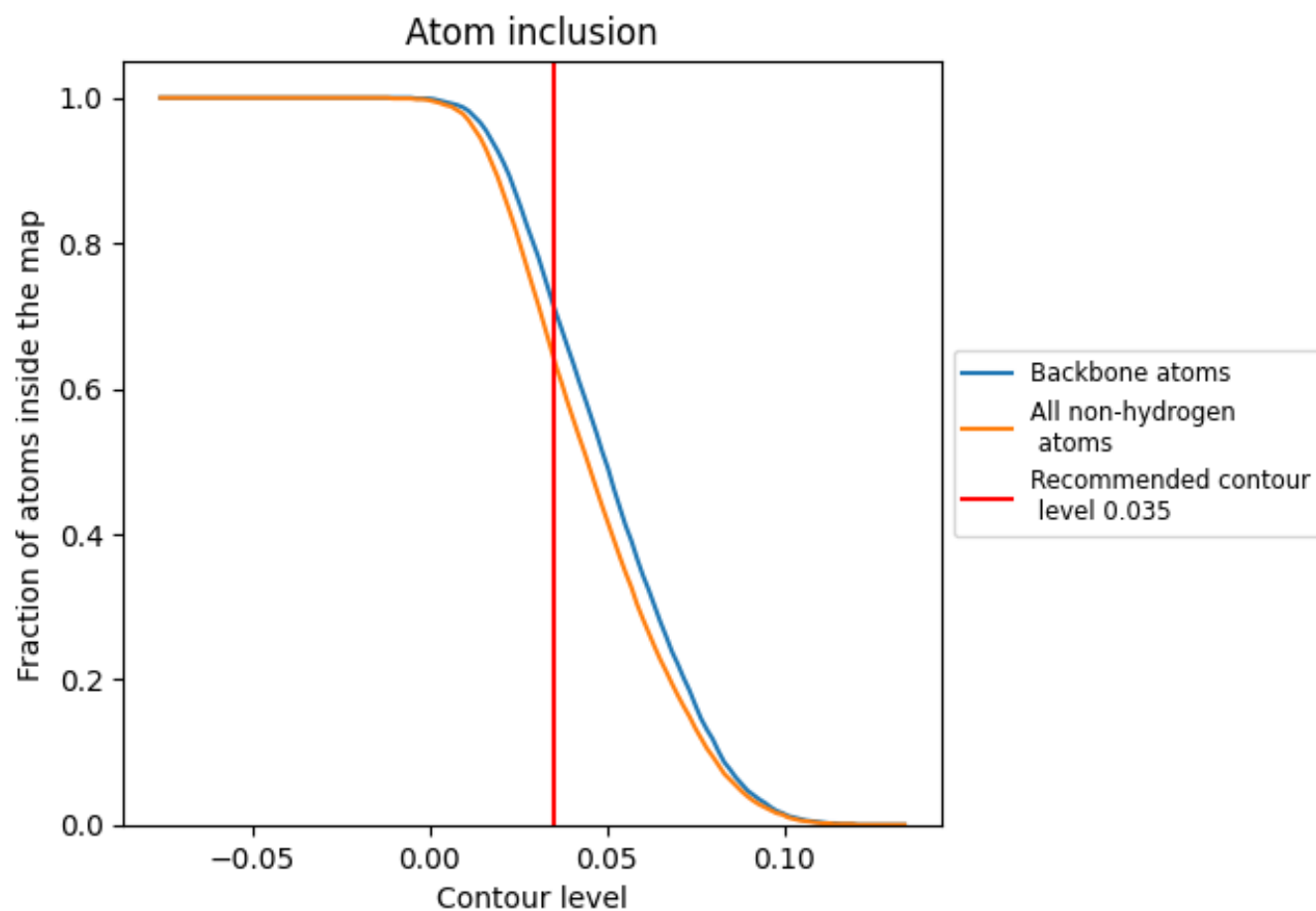
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.035).

9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 64% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.035) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.6420	<div></div> 0.4920
A	<div></div> 0.6250	<div></div> 0.4910
B	<div></div> 0.6270	<div></div> 0.4850
C	<div></div> 0.6860	<div></div> 0.4960
D	<div></div> 0.6380	<div></div> 0.4870
E	<div></div> 0.6750	<div></div> 0.4990

