



wwPDB EM Validation Summary Report ⓘ

Mar 29, 2026 – 03:48 AM UTC

PDB ID : 7WLI / pdb_00007wli
EMDB ID : EMD-32584
Title : CryoEM structure of human low-voltage activated T-type calcium channel CaV3.3 (apo)
Authors : He, L.; Yu, Z.; Dong, Y.; Chen, Q.; Zhao, Y.
Deposited on : 2022-01-13
Resolution : 3.30 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/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.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

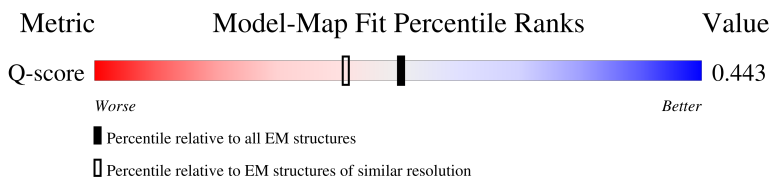
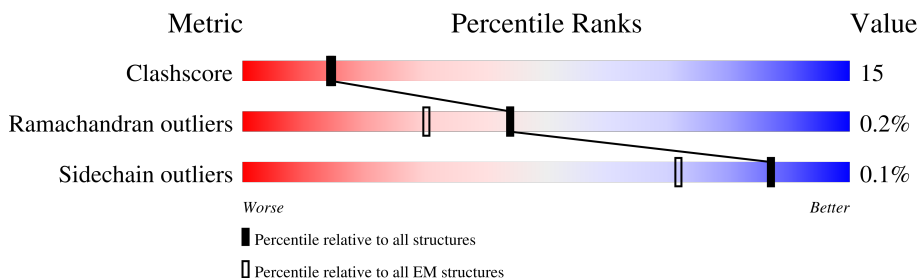
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.30 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)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15087 (2.80 - 3.80)

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	2223	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 9493 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.

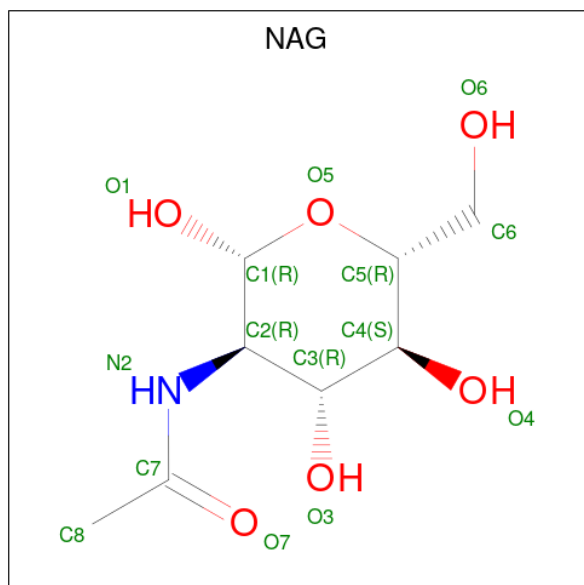
- Molecule 1 is a protein called Voltage-dependent T-type calcium channel subunit alpha-1I.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	1135	9111	5979	1499	1555	78	1	0

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

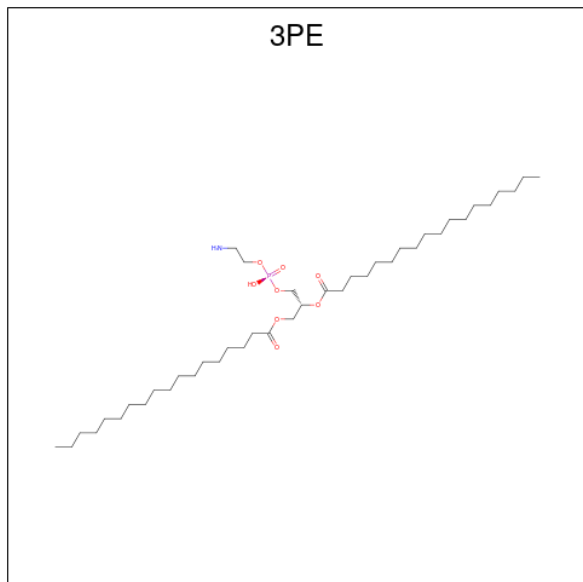
Mol	Chain	Residues	Atoms		AltConf
2	A	2	Total	Ca	0
			2	2	

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



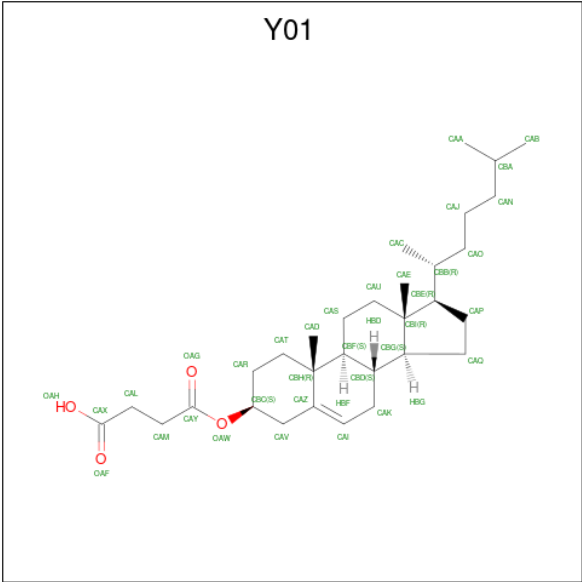
Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
3	A	1	14	8	1	5	0
3	A	1	14	8	1	5	0

- Molecule 4 is 1,2-Distearoyl-sn-glycerophosphoethanolamine (CCD ID: 3PE) (formula: $C_{41}H_{82}NO_8P$).



Mol	Chain	Residues	Atoms				AltConf
4	A	1	Total	C	O	P	0
			32	23	8	1	
4	A	1	Total	C			0
			11	11			
4	A	1	Total	C			0
			12	12			
4	A	1	Total	C			0
			12	12			
4	A	1	Total	C			0
			12	12			
4	A	1	Total	C			0
			13	13			
4	A	1	Total	C			0
			10	10			
4	A	1	Total	C			0
			13	13			
4	A	1	Total	C			0
			16	16			
4	A	1	Total	C			0
			10	10			
4	A	1	Total	C			0
			15	15			

- Molecule 5 is CHOLESTEROL HEMISUCCINATE (CCD ID: Y01) (formula: $C_{31}H_{50}O_4$).

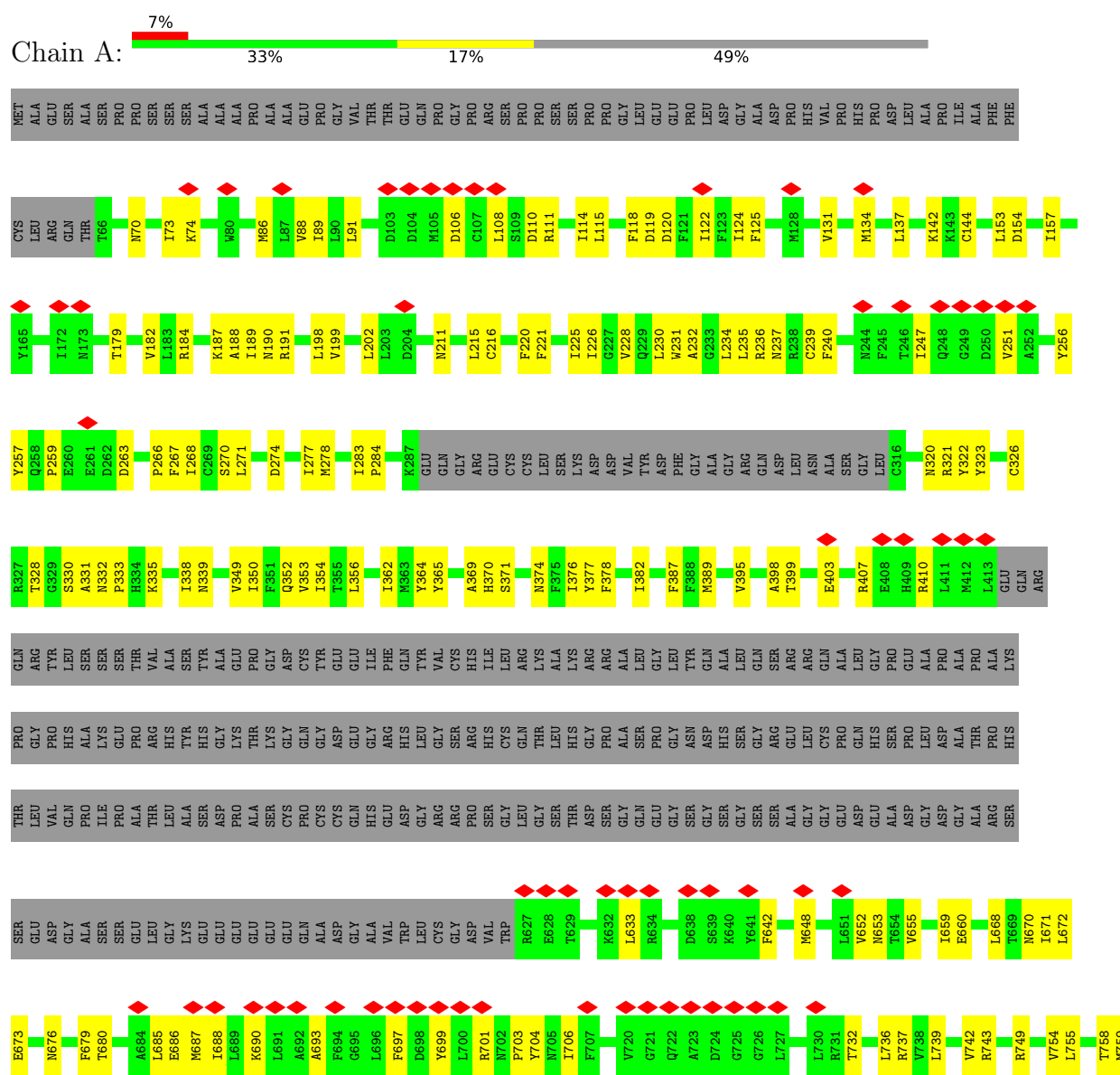


Mol	Chain	Residues	Atoms			AltConf
5	A	1	Total	C	O	0
			31	29	2	
5	A	1	Total	C	O	0
			31	29	2	
5	A	1	Total	C	O	0
			35	31	4	
5	A	1	Total	C	O	0
			35	31	4	
5	A	1	Total	C	O	0
			29	28	1	
5	A	1	Total	C	O	0
			35	31	4	

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: Voltage-dependent T-type calcium channel subunit alpha-1I





TRP	GLY	SER	SER	PRO	ARG	PRO	ASP
ALA	SER	PRO	ALA	ARG	ARG	ASP	
ALA	GLU	ALA	VAL	ASN	GLN	SER	
ASP	HIS	ALA	ALA	ARG	CYS	ALA	
SER	GLU	ARG	THR	ARG	THR	PRO	
LYS	THR	ARG	LEU	ARG	LEU	PRO	
ASP	LEU	LEU	LEU	LEU	LEU	SER	
PRO	SER	SER	ARG	SER	GLN	PRO	
PRO	SER	LEU	GLN	ARG	ALA	SER	
GLY	LEU	ARG	THR	GLY	THR	PRO	
ARG	LEU	ARG	GLY	ARG	ASP	ASP	
PRO	THR	GLY	SER	GLY	SER	ALA	
PRO	THR	GLY	SER	SER	ASP	ALA	
LEU	SER	LEU	ASP	THR	SER	SER	
GLY	PRO	GLY	ALA	LEU	LEU	LEU	
PRO	PRO	PRO	ASP	LEU	ASP	LEU	
PRO	PRO	ARG	SER	SER	PRO	PRO	
ALA	PRO	ALA	SER	ALA	SER	ALA	
PRO	PRO	HIS	SER	SER	GLU	GLU	
PRO	ALA	GLN	SER	SER	SER	PHE	
PRO	PRO	ARG	ALA	ALA	ALA	HIS	
GLY	GLY	ARG	GLY	GLY	SER	PRO	
GLU	PHE	THR	GLU	GLU	GLN	LYS	
LEU	SER	SER	ASP	ASP	LYS	GLY	
PRO	SER	PRO	SER	SER	PRO	PRO	
GLY	SER	GLY	LEU	LEU	PRO	GLY	
ASP	THR	GLY	LEU	LEU	PRO	GLU	
ALA	SER	GLY	THR	THR	THR	LYS	
ALA	SER	ASP	SER	SER	GLY	THR	
ARG	PRO	ALA	ALA	ASP	GLY	THR	
ARG	PRO	GLY	GLY	MET	ARG	THR	
	ARG	ARG	ARG	ASP	ARG	LEU	
		PRO	PRO	ALA	ALA	PRO	
		HIS	SER	LEU	GLY	LYS	
		ALA	ASP	GLY	LEU	LYS	
		ALA	GLU	PRO	PRO	ILE	
		ALA	GLU	PRO	ALA	ALA	
		LEU	GLY	ALA	LEU	GLN	
		ALA	ARG	PRO	GLN	GLY	
		HIS	GLY	ALA	ALA	GLY	
		GLY	GLY	PRO	PRO	SER	
		LEU	ALA	GLY	THR	TRP	
		ALA	GLY	ALA	ARG	ALA	
		ARG	GLY	ALA	LEU	SER	
		SER	GLY	ALA	GLY	LEU	
		PRO	GLY	LEU	THR	ASP	
		SER	ALA	THR	SER	GLY	

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	93988	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	9.6	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	3.327	Depositor
Minimum map value	-2.203	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.068	Depositor
Recommended contour level	0.35	Depositor
Map size ($\text{\AA}$)	266.24, 266.24, 266.24	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.04, 1.04, 1.04	Depositor

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: 3PE, CA, NAG, Y01

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	0.33	4/9314 (0.0%)	0.51	10/12625 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1443	ALA	C-O	10.41	1.36	1.24
1	A	1439	GLN	C-O	7.98	1.33	1.24
1	A	1459	LYS	C-O	7.97	1.33	1.24
1	A	1450	GLU	C-O	6.83	1.32	1.24

The worst 5 of 10 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1443	ALA	N-CA-C	11.92	123.96	110.97
1	A	1463	LYS	N-CA-C	9.79	123.19	108.31
1	A	1446	ALA	CA-C-N	-7.75	109.12	120.28
1	A	1446	ALA	C-N-CA	-7.75	109.12	120.28
1	A	1446	ALA	N-CA-C	-7.46	103.08	111.07

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	9111	0	9291	280	0
2	A	2	0	0	0	0
3	A	28	0	26	2	0
4	A	156	0	245	8	0
5	A	196	0	282	20	0
All	All	9493	0	9844	293	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 15.

The worst 5 of 293 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:1460:LYS:O	1:A:1463:LYS:HG3	1.47	1.12
1:A:1441:GLN:OE1	1:A:1734:ASP:OD2	1.89	0.89
1:A:1376:LEU:HD23	1:A:1382:TRP:HB2	1.56	0.87
1:A:1441:GLN:HE21	1:A:1445:GLU:CB	1.93	0.81
1:A:1441:GLN:OE1	1:A:1734:ASP:CG	2.25	0.80

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	1128/2223 (51%)	1044 (93%)	82 (7%)	2 (0%)	43 71

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1442	GLU
1	A	1441	GLN

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	1006/1879 (54%)	1005 (100%)	1 (0%)	88 90

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1444	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1481	HIS
1	A	1511	GLN
1	A	1552	ASN
1	A	1523	ASN
1	A	816	GLN

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](#)

Of 21 ligands modelled in this entry, 2 are monoatomic - leaving 19 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
4	3PE	A	2307	-	11,11,50	1.69	0	10,10,55	0.88	0
5	Y01	A	2313	-	32,32,38	0.82	2 (6%)	49,49,57	1.57	14 (28%)
4	3PE	A	2318	-	12,12,50	1.69	0	11,11,55	0.89	0
4	3PE	A	2308	-	11,11,50	1.68	0	10,10,55	0.89	0
5	Y01	A	2310	-	34,34,38	0.68	1 (2%)	52,52,57	1.74	9 (17%)
3	NAG	A	2303	1	14,14,15	0.41	0	17,19,21	1.32	2 (11%)
4	3PE	A	2305	-	31,31,50	2.48	11 (35%)	34,36,55	1.25	2 (5%)
5	Y01	A	2311	-	38,38,38	0.74	1 (2%)	57,57,57	1.64	9 (15%)
3	NAG	A	2304	1	14,14,15	0.22	0	17,19,21	0.57	0
5	Y01	A	2309	-	34,34,38	0.66	0	52,52,57	1.66	8 (15%)
4	3PE	A	2316	-	12,12,50	1.70	0	11,11,55	0.83	0
4	3PE	A	2319	-	15,15,50	1.70	0	14,14,55	0.85	0
4	3PE	A	2306	-	10,10,50	1.68	0	9,9,55	0.87	0
5	Y01	A	2315	-	38,38,38	0.69	1 (2%)	57,57,57	1.66	10 (17%)
4	3PE	A	2314	-	11,11,50	1.70	0	10,10,55	0.85	0
4	3PE	A	2320	-	9,9,50	1.67	0	8,8,55	0.81	0
4	3PE	A	2321	-	14,14,50	1.70	0	13,13,55	0.85	0
5	Y01	A	2312	-	38,38,38	0.64	0	57,57,57	1.82	13 (22%)
4	3PE	A	2317	-	9,9,50	1.66	0	8,8,55	0.80	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
4	3PE	A	2307	-	-	1/9/9/54	-
5	Y01	A	2313	-	-	6/12/70/77	0/4/4/4
4	3PE	A	2318	-	-	1/10/10/54	-
4	3PE	A	2308	-	-	0/9/9/54	-
5	Y01	A	2310	-	-	12/14/72/77	0/4/4/4
3	NAG	A	2303	1	-	6/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	3PE	A	2305	-	-	11/33/33/54	-
5	Y01	A	2311	-	-	11/19/77/77	0/4/4/4
3	NAG	A	2304	1	-	4/6/23/26	0/1/1/1
5	Y01	A	2309	-	-	5/14/72/77	0/4/4/4
4	3PE	A	2316	-	-	1/10/10/54	-
4	3PE	A	2319	-	-	5/13/13/54	-
4	3PE	A	2306	-	-	1/8/8/54	-
5	Y01	A	2315	-	-	6/19/77/77	0/4/4/4
4	3PE	A	2314	-	-	3/9/9/54	-
4	3PE	A	2320	-	-	4/7/7/54	-
4	3PE	A	2321	-	-	5/12/12/54	-
5	Y01	A	2312	-	-	11/19/77/77	0/4/4/4
4	3PE	A	2317	-	-	2/7/7/54	-

The worst 5 of 16 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	2305	3PE	P-O13	5.04	1.73	1.54
4	A	2305	3PE	O31-C31	4.69	1.47	1.33
4	A	2305	3PE	O21-C21	4.60	1.47	1.34
4	A	2305	3PE	P-O11	4.15	1.73	1.60
4	A	2305	3PE	C22-C21	3.42	1.60	1.50

The worst 5 of 67 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	2315	Y01	CBI-CBE-CBB	-5.59	110.87	119.50
5	A	2310	Y01	OAW-CAY-CAM	5.12	120.21	111.09
5	A	2309	Y01	OAW-CAY-CAM	4.85	119.75	111.09
5	A	2312	Y01	CBI-CBE-CBB	-4.62	112.36	119.50
3	A	2303	NAG	C2-N2-C7	4.56	129.00	122.90

There are no chirality outliers.

5 of 95 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	2310	Y01	OAG-CAY-OAW-CBC
5	A	2311	Y01	CAM-CAY-OAW-CBC
5	A	2312	Y01	CAR-CBC-OAW-CAY

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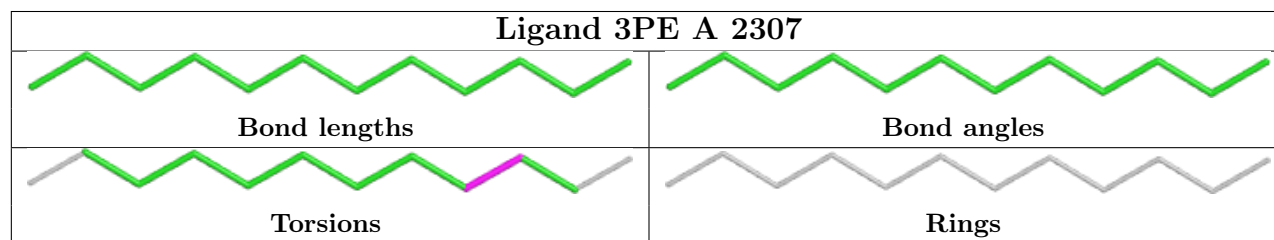
Mol	Chain	Res	Type	Atoms
5	A	2312	Y01	CAM-CAY-OAW-CBC
5	A	2313	Y01	CAV-CBC-OAW-CAY

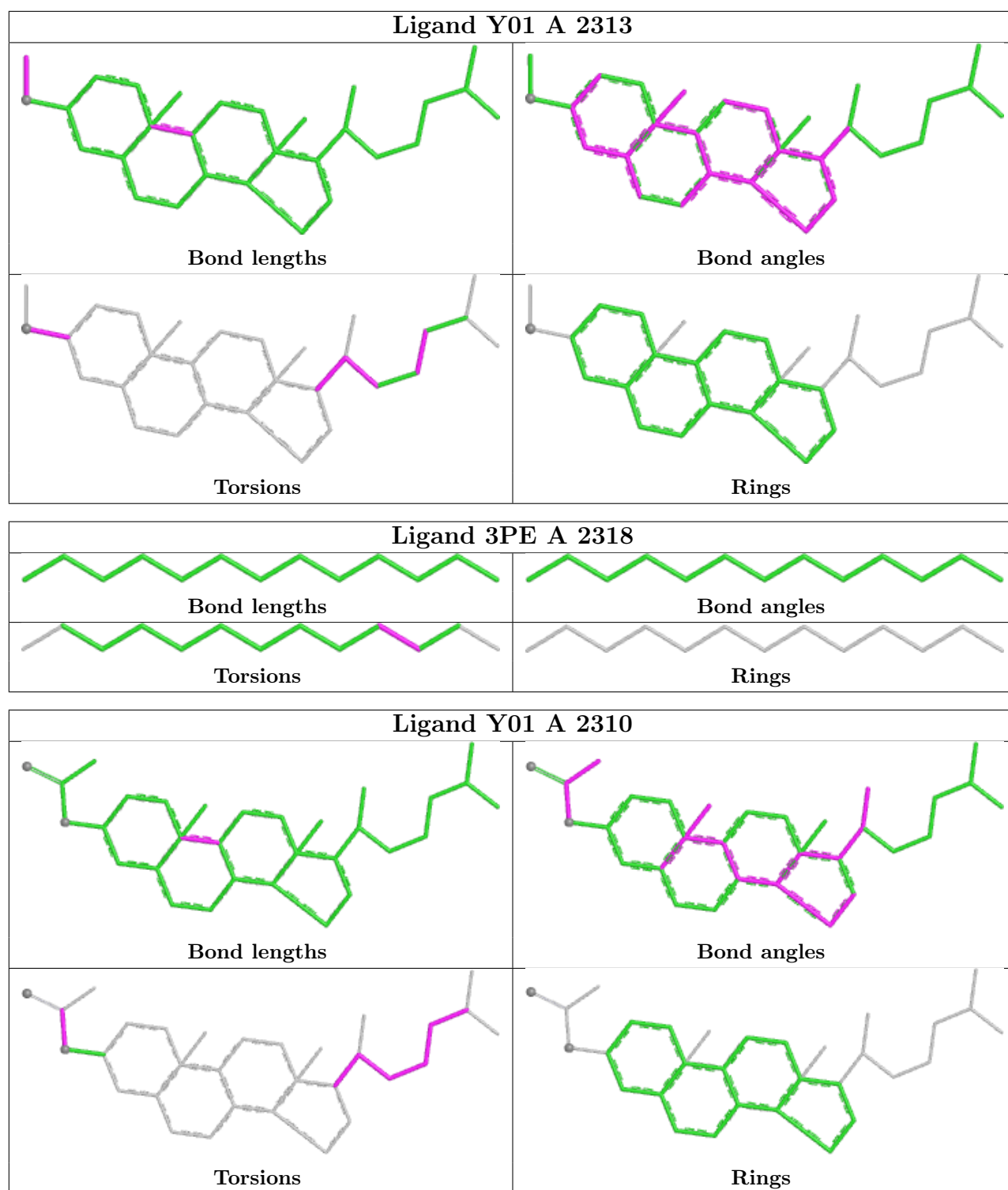
There are no ring outliers.

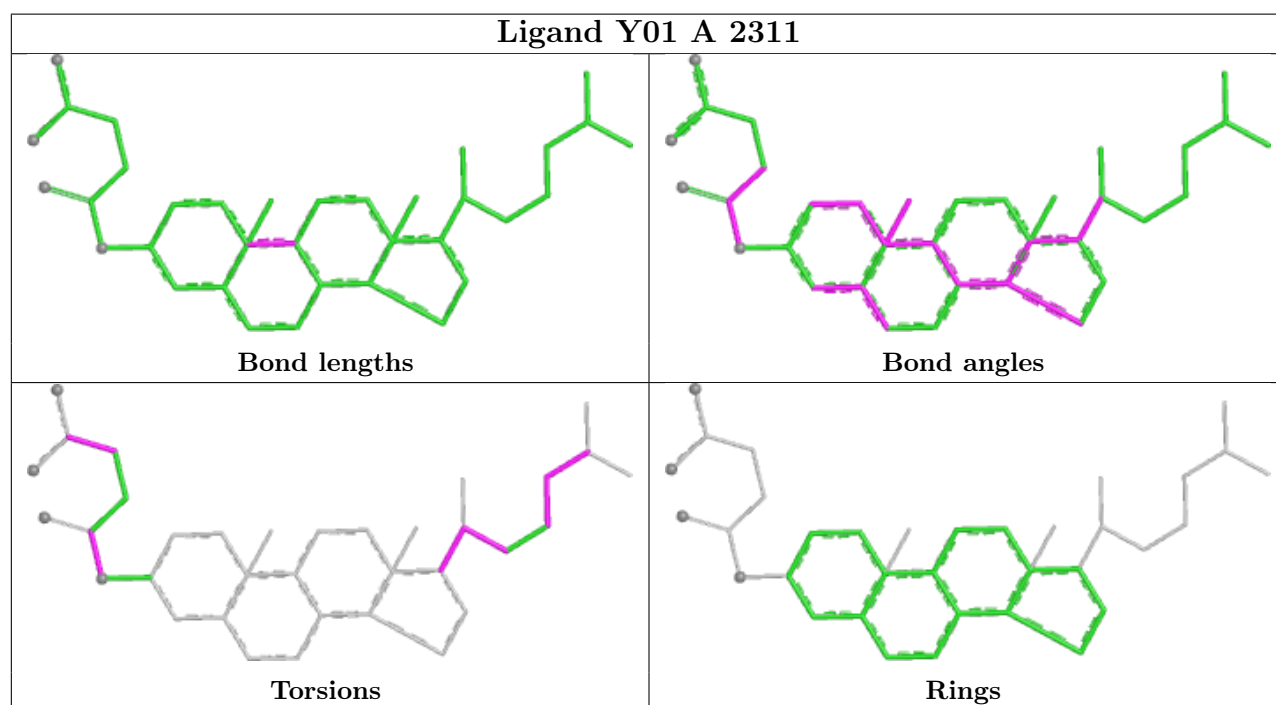
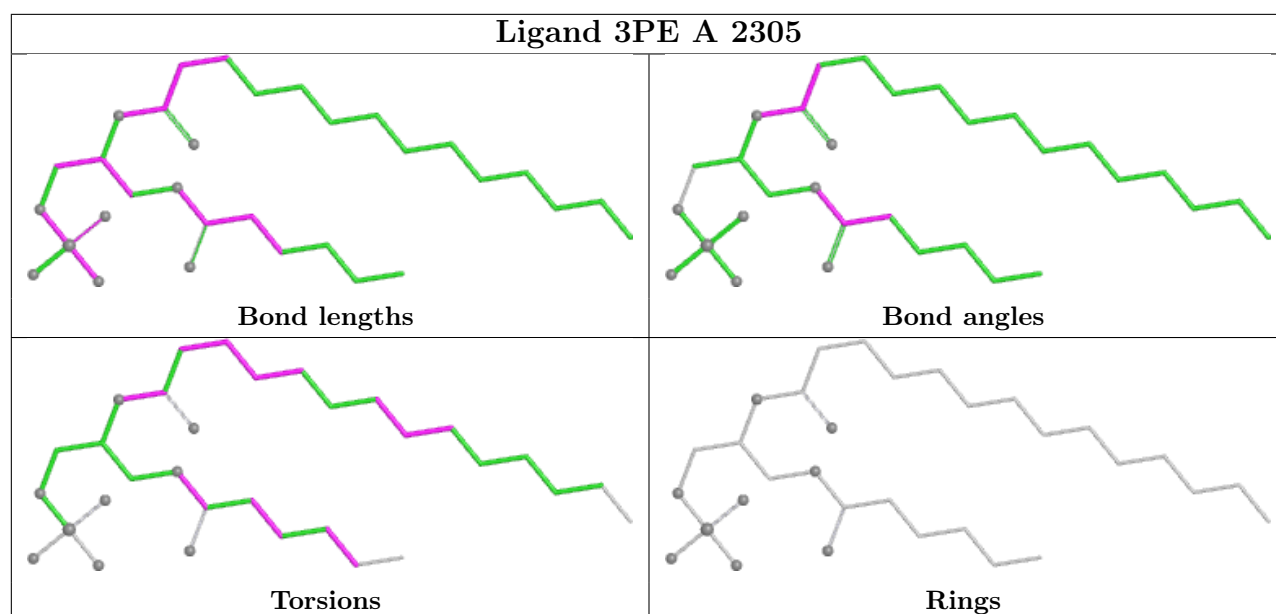
14 monomers are involved in 30 short contacts:

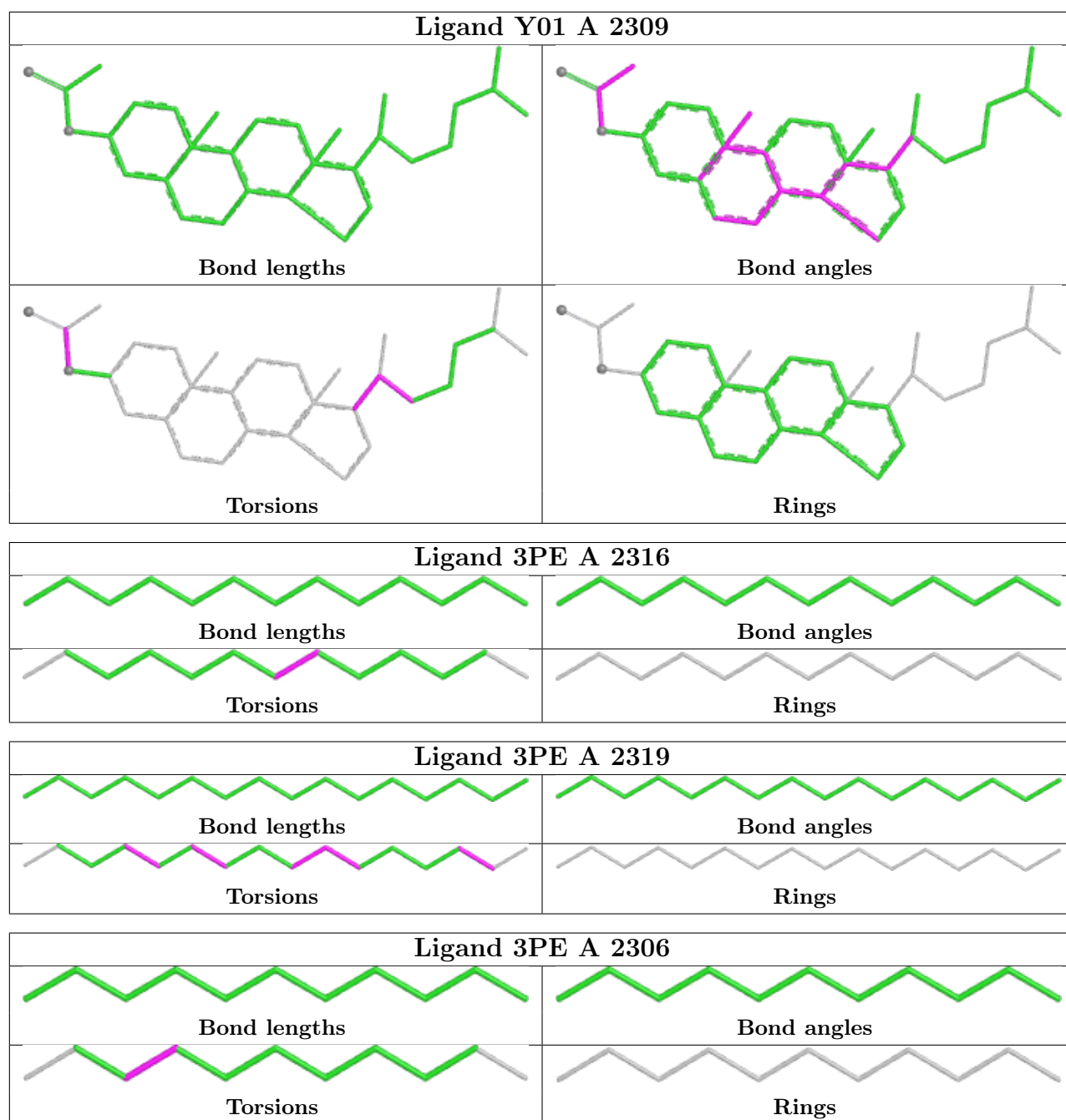
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	2313	Y01	5	0
4	A	2308	3PE	2	0
5	A	2310	Y01	1	0
3	A	2303	NAG	1	0
4	A	2305	3PE	1	0
5	A	2311	Y01	8	0
3	A	2304	NAG	1	0
5	A	2309	Y01	1	0
4	A	2316	3PE	2	0
5	A	2315	Y01	2	0
4	A	2314	3PE	1	0
4	A	2321	3PE	1	0
5	A	2312	Y01	4	0
4	A	2317	3PE	1	0

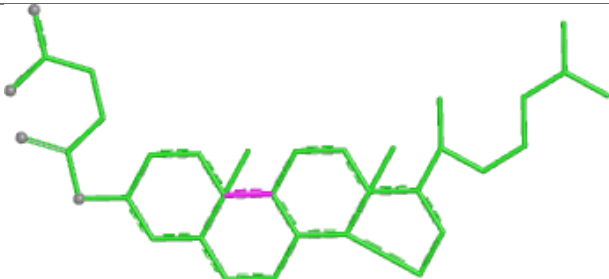
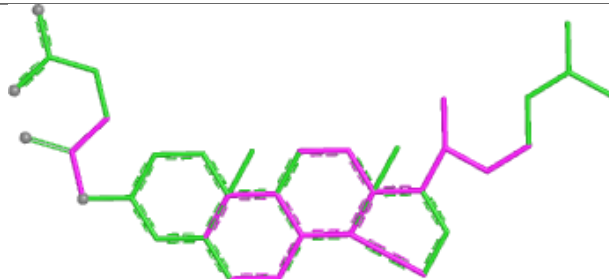
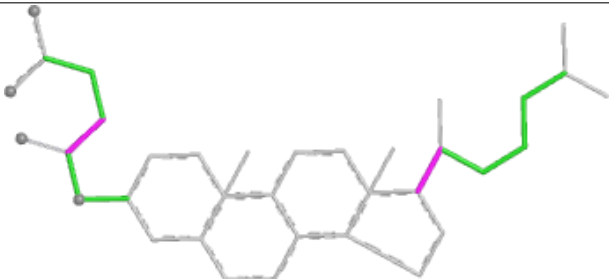
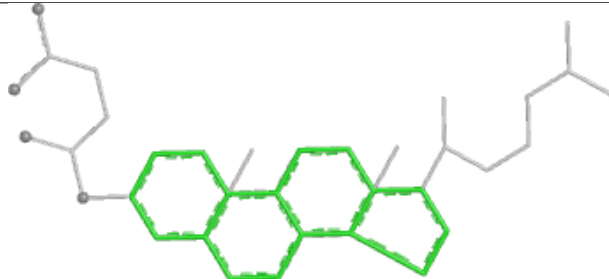
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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





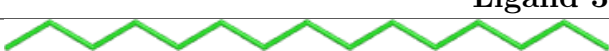
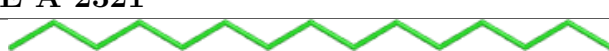


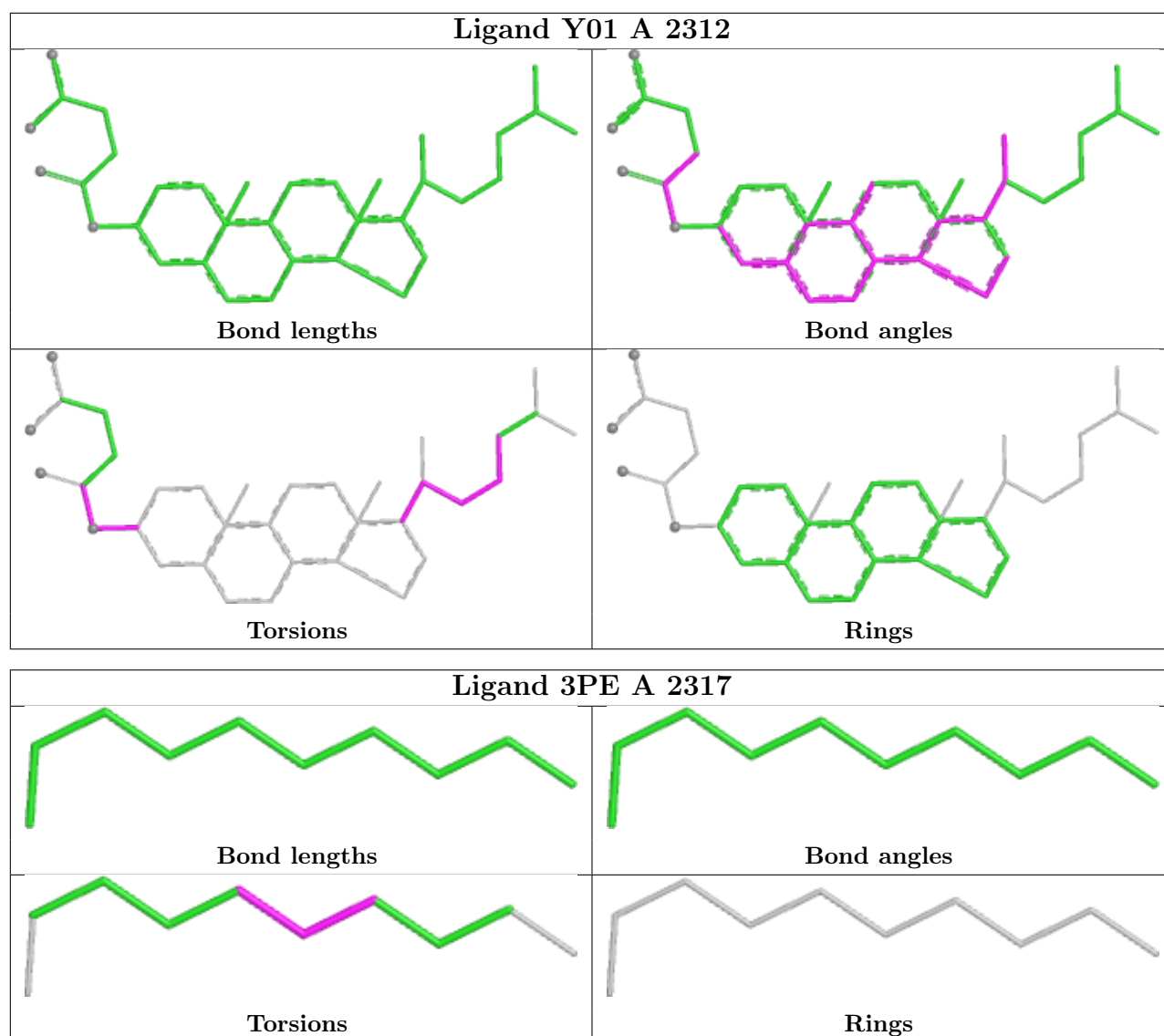


Ligand Y01 A 2315	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand 3PE A 2314	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand 3PE A 2320	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand 3PE A 2321	
	
Bond lengths	Bond angles
	
Torsions	Rings



5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

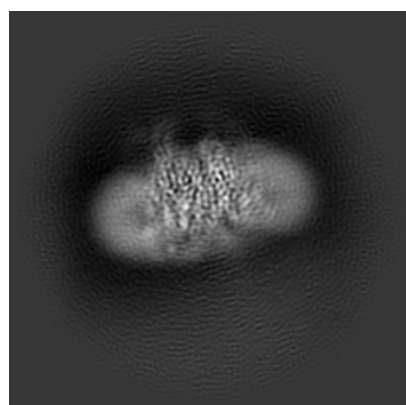
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-32584. These allow visual inspection of the internal detail of the map and identification of artifacts.

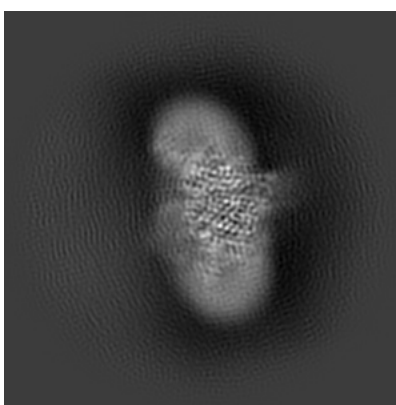
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

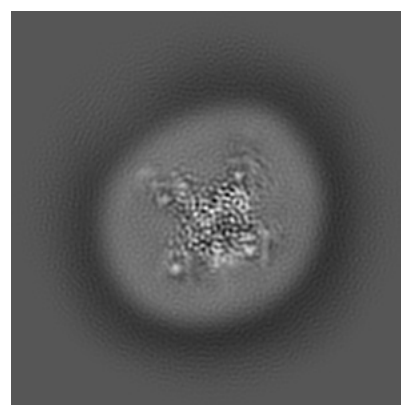
6.1.1 Primary map



X



Y

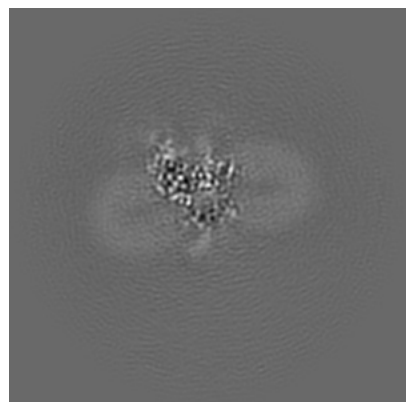


Z

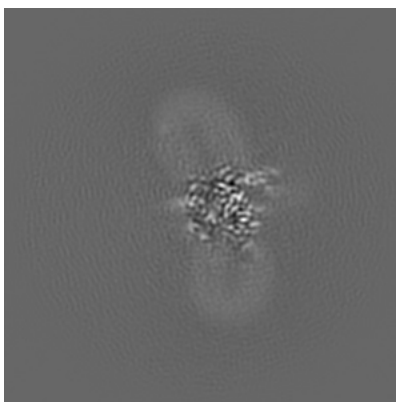
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

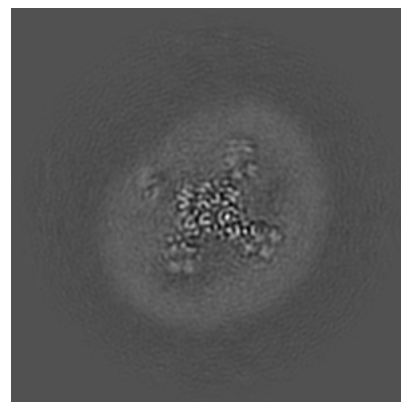
6.2.1 Primary map



X Index: 128



Y Index: 128

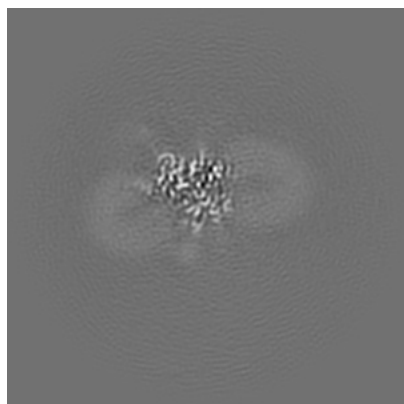


Z Index: 128

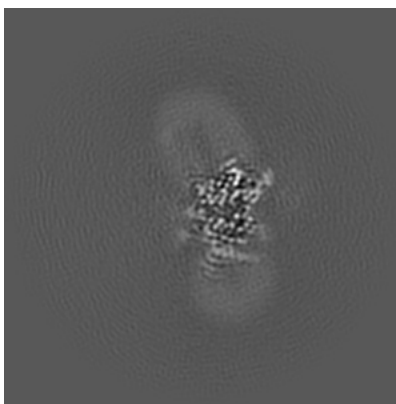
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

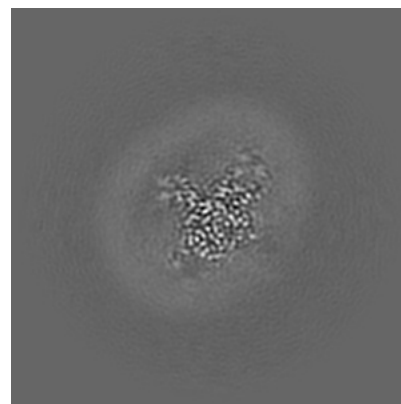
6.3.1 Primary map



X Index: 122



Y Index: 135

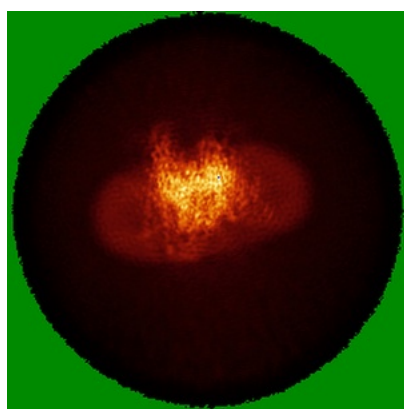


Z Index: 146

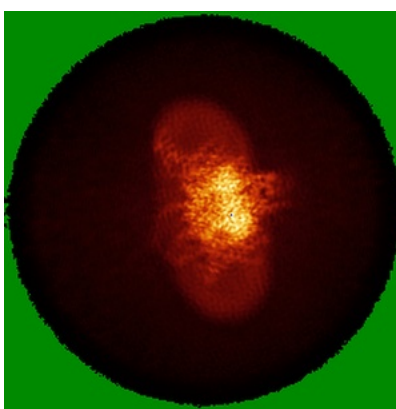
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) [i](#)

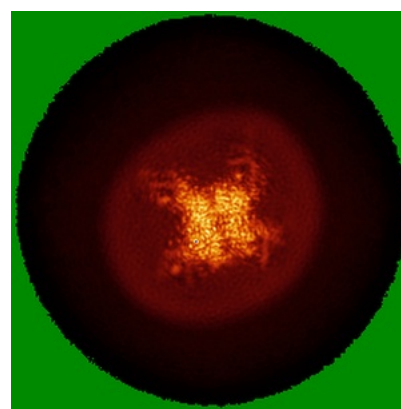
6.4.1 Primary map



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 0.35. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

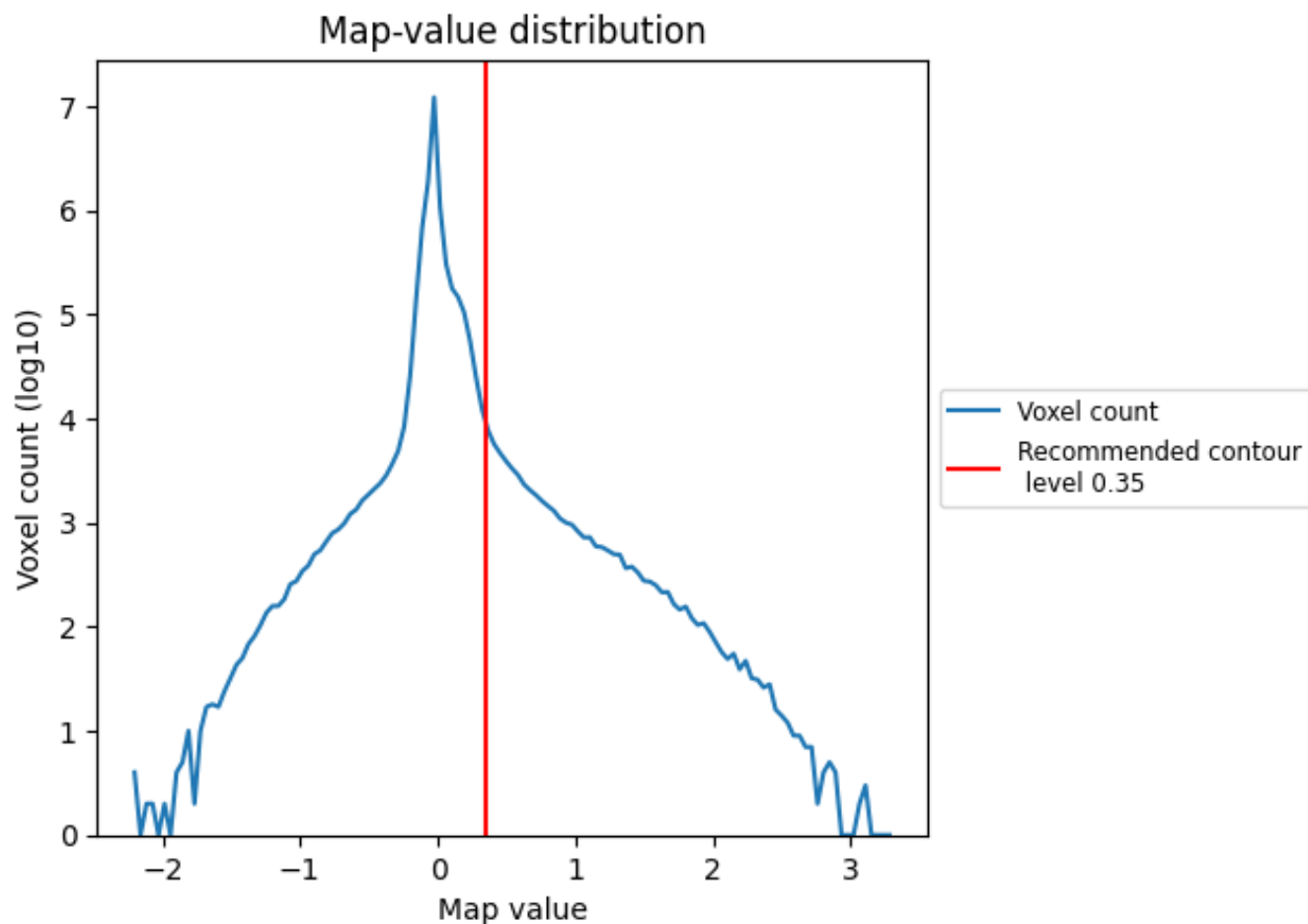
6.6 Mask visualisation [i](#)

This section was not generated. No masks/segmentation were deposited.

7 Map analysis [i](#)

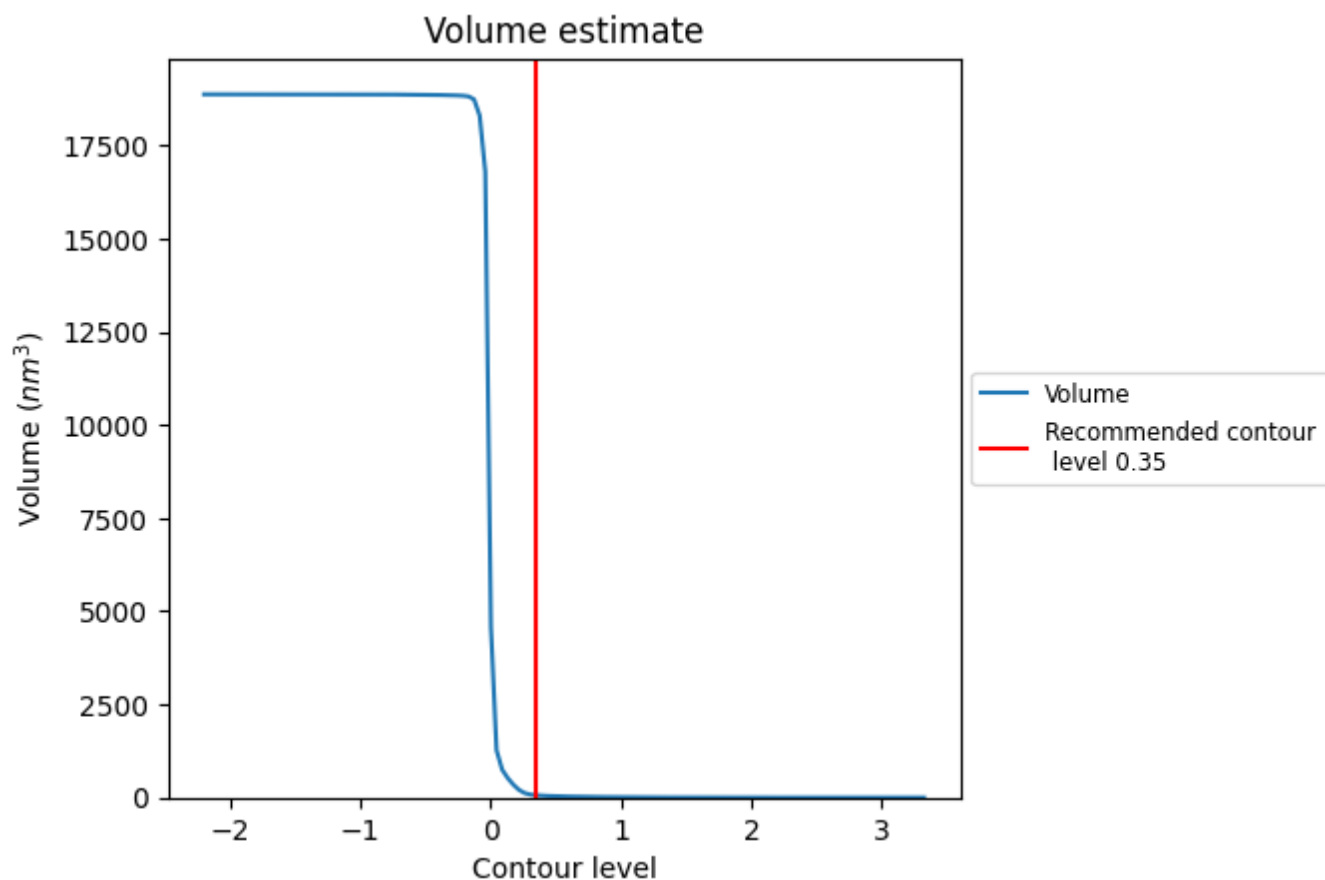
This section contains the results of statistical analysis of the map.

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.

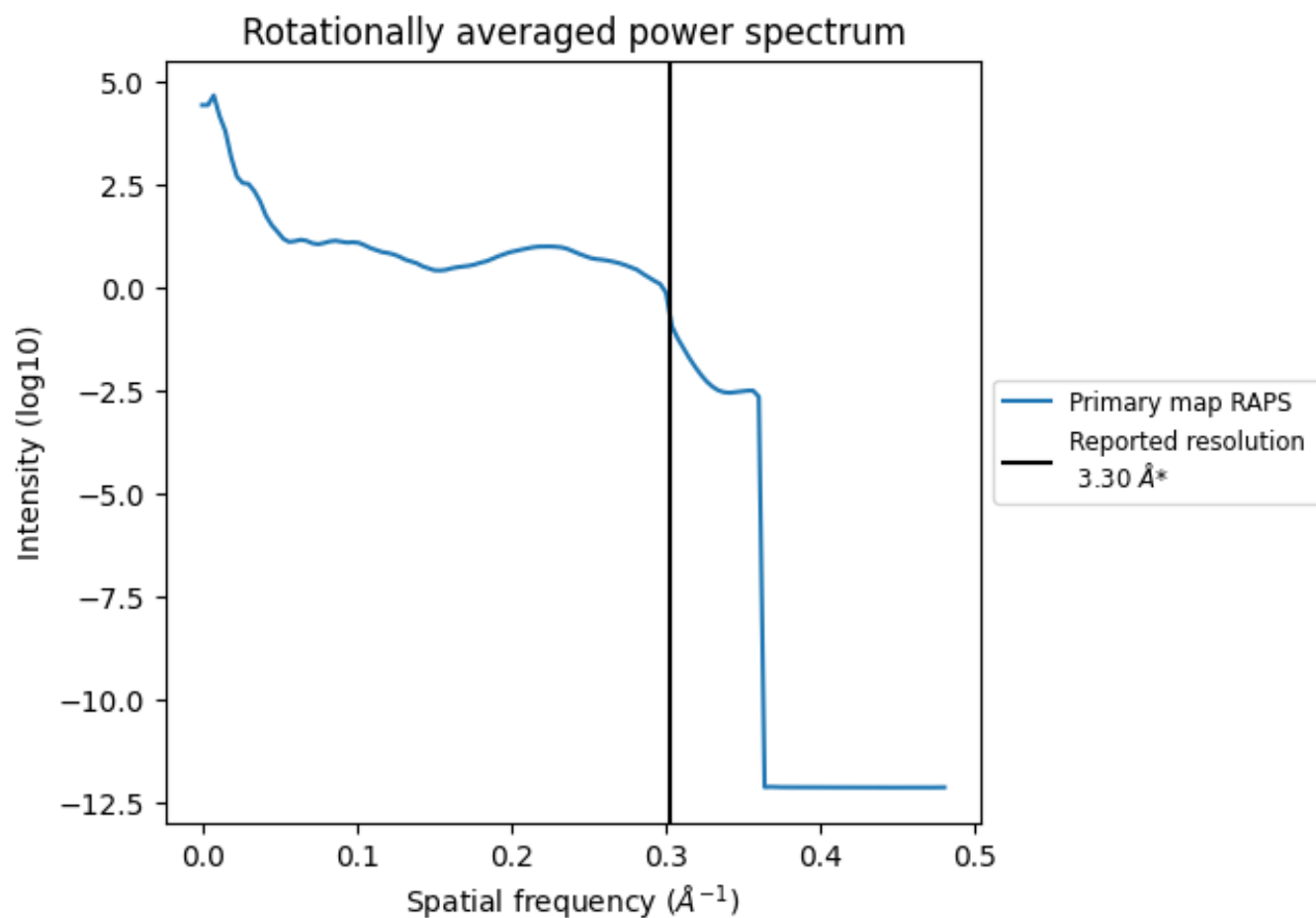
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 61 nm³; this corresponds to an approximate mass of 55 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 0.303 Å⁻¹

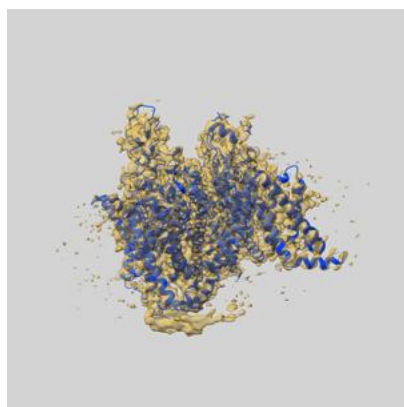
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

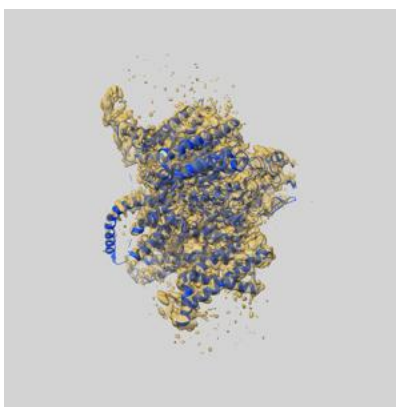
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-32584 and PDB model 7WLI. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

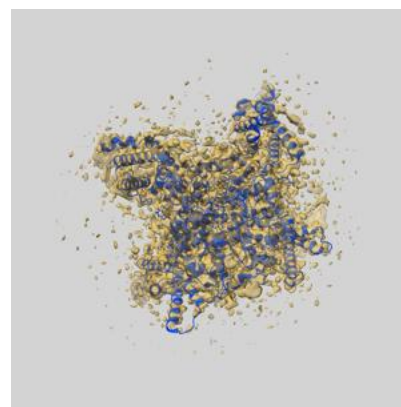
9.1 Map-model overlay [i](#)



X



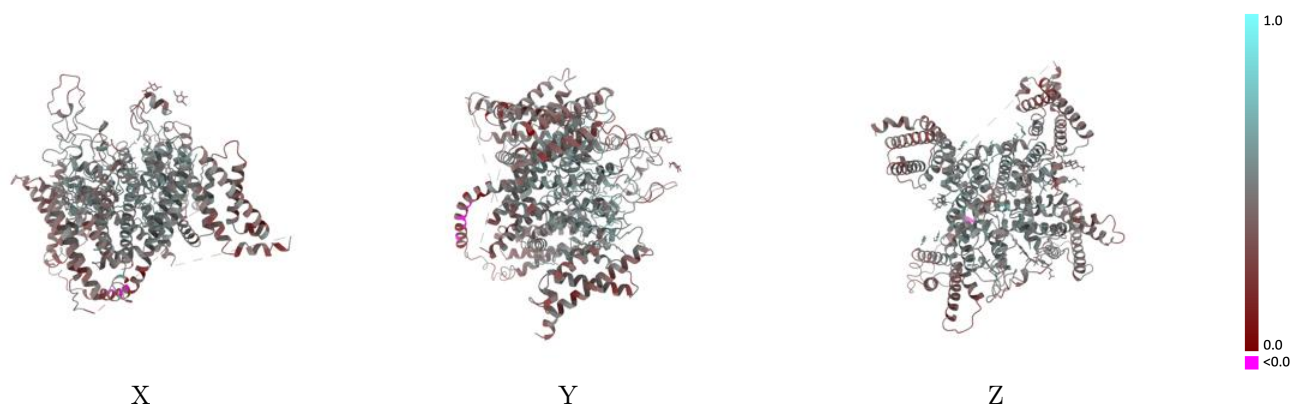
Y



Z

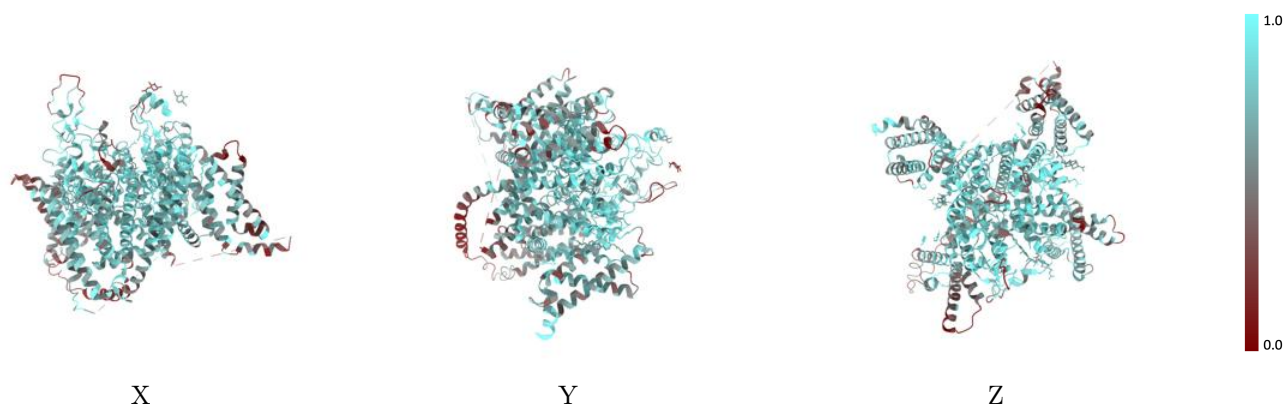
The images above show the 3D surface view of the map at the recommended contour level 0.35 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



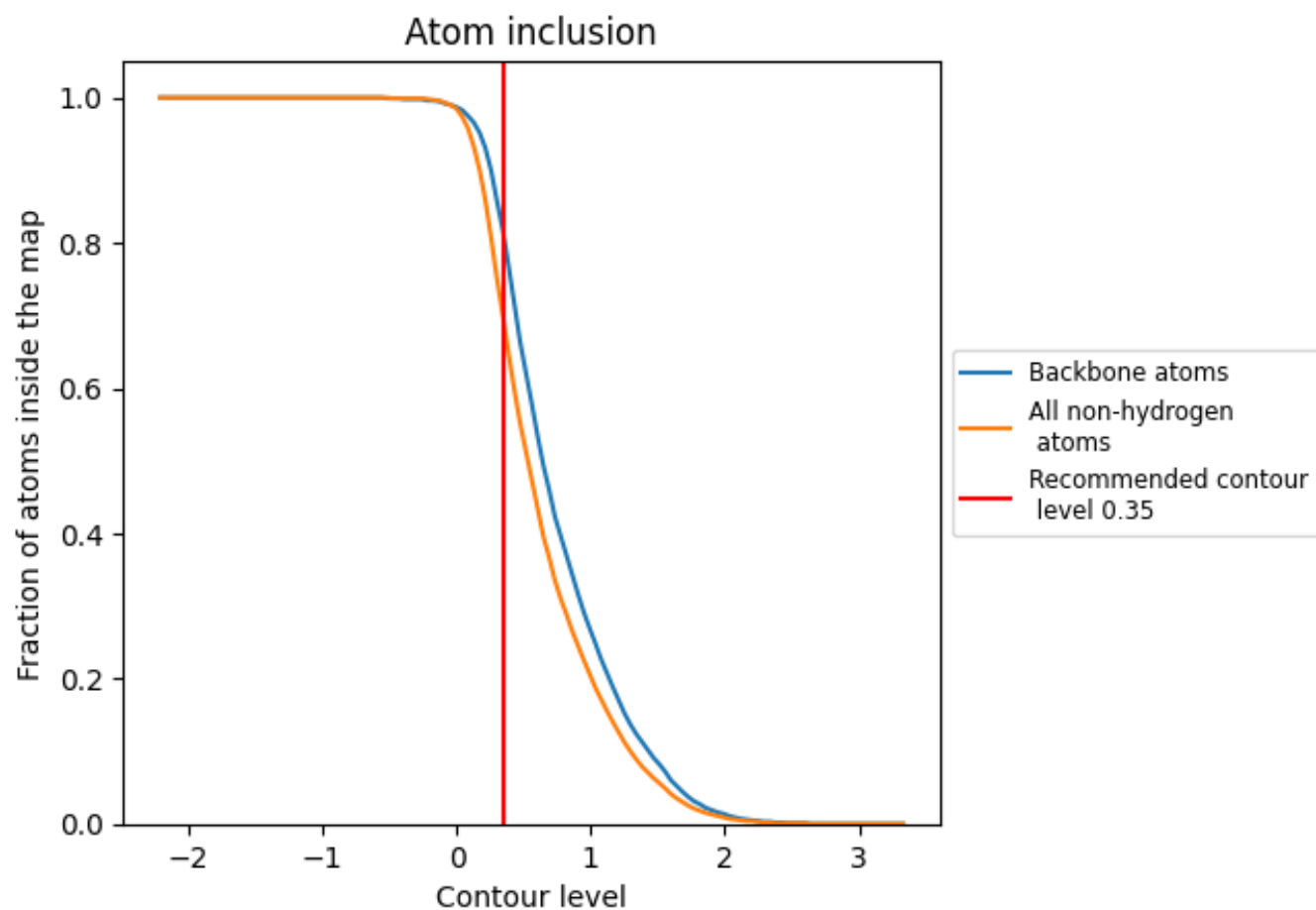
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.35).

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 70% 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 (0.35) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6970	0.4430
A	0.6970	0.4430

