



Full wwPDB EM Validation Report ⓘ

Jun 24, 2026 – 01:48 PM EDT

PDB ID : 13LR / pdb_000013lr
EMDB ID : EMD-77142
Title : Structure of human TRPV3-G568D Olmsted syndrome mutant in the 4-fold symmetrical inactivated state
Authors : Nadezhdin, K.D.; Khau, J.; Purohit, R.; Sobolevsky, A.I.
Deposited on : 2026-05-13
Resolution : 3.33 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.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

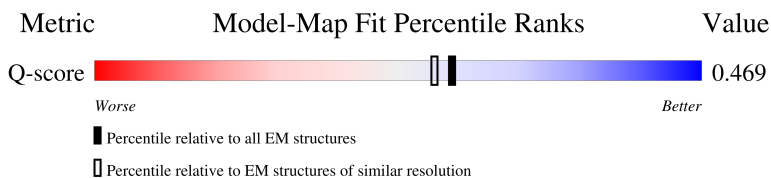
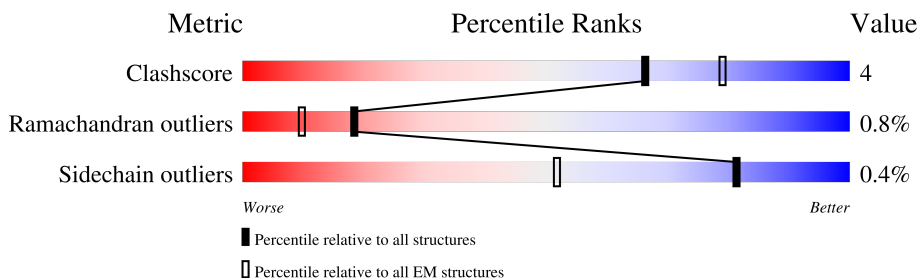
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.33 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	14484 (2.83 - 3.83)

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	1052	
1	B	1052	
1	C	1052	
1	D	1052	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 20575 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 Isoform 2 of Transient receptor potential cation channel sub-family V member 3, Green fluorescent protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	612	Total	C	N	O	S	0	0
			4987	3249	821	886	31		
1	B	612	Total	C	N	O	S	0	0
			4987	3249	821	886	31		
1	C	612	Total	C	N	O	S	0	0
			4987	3249	821	886	31		
1	D	612	Total	C	N	O	S	0	0
			4987	3249	821	886	31		

There are 116 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	568	ASP	GLY	engineered mutation	UNP Q8NET8
A	792	LEU	-	linker	UNP Q8NET8
A	793	VAL	-	linker	UNP Q8NET8
A	794	PRO	-	linker	UNP Q8NET8
A	795	ARG	-	linker	UNP Q8NET8
A	796	GLY	-	linker	UNP Q8NET8
A	797	SER	-	linker	UNP Q8NET8
A	798	ALA	-	linker	UNP Q8NET8
A	799	ALA	-	linker	UNP Q8NET8
A	800	ALA	-	linker	UNP Q8NET8
A	801	ALA	-	linker	UNP Q8NET8
A	802	VAL	-	linker	UNP Q8NET8
A	865	LEU	PHE	conflict	UNP P42212
A	866	THR	SER	conflict	UNP P42212
A	1007	LYS	ALA	conflict	UNP P42212
A	1032	LEU	HIS	conflict	UNP P42212
A	1040	SER	-	expression tag	UNP P42212
A	1041	GLY	-	expression tag	UNP P42212
A	1042	LEU	-	expression tag	UNP P42212
A	1043	ARG	-	expression tag	UNP P42212
A	1044	SER	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1045	TRP	-	expression tag	UNP P42212
A	1046	SER	-	expression tag	UNP P42212
A	1047	HIS	-	expression tag	UNP P42212
A	1048	PRO	-	expression tag	UNP P42212
A	1049	GLN	-	expression tag	UNP P42212
A	1050	PHE	-	expression tag	UNP P42212
A	1051	GLU	-	expression tag	UNP P42212
A	1052	LYS	-	expression tag	UNP P42212
B	568	ASP	GLY	engineered mutation	UNP Q8NET8
B	792	LEU	-	linker	UNP Q8NET8
B	793	VAL	-	linker	UNP Q8NET8
B	794	PRO	-	linker	UNP Q8NET8
B	795	ARG	-	linker	UNP Q8NET8
B	796	GLY	-	linker	UNP Q8NET8
B	797	SER	-	linker	UNP Q8NET8
B	798	ALA	-	linker	UNP Q8NET8
B	799	ALA	-	linker	UNP Q8NET8
B	800	ALA	-	linker	UNP Q8NET8
B	801	ALA	-	linker	UNP Q8NET8
B	802	VAL	-	linker	UNP Q8NET8
B	865	LEU	PHE	conflict	UNP P42212
B	866	THR	SER	conflict	UNP P42212
B	1007	LYS	ALA	conflict	UNP P42212
B	1032	LEU	HIS	conflict	UNP P42212
B	1040	SER	-	expression tag	UNP P42212
B	1041	GLY	-	expression tag	UNP P42212
B	1042	LEU	-	expression tag	UNP P42212
B	1043	ARG	-	expression tag	UNP P42212
B	1044	SER	-	expression tag	UNP P42212
B	1045	TRP	-	expression tag	UNP P42212
B	1046	SER	-	expression tag	UNP P42212
B	1047	HIS	-	expression tag	UNP P42212
B	1048	PRO	-	expression tag	UNP P42212
B	1049	GLN	-	expression tag	UNP P42212
B	1050	PHE	-	expression tag	UNP P42212
B	1051	GLU	-	expression tag	UNP P42212
B	1052	LYS	-	expression tag	UNP P42212
C	568	ASP	GLY	engineered mutation	UNP Q8NET8
C	792	LEU	-	linker	UNP Q8NET8
C	793	VAL	-	linker	UNP Q8NET8
C	794	PRO	-	linker	UNP Q8NET8
C	795	ARG	-	linker	UNP Q8NET8

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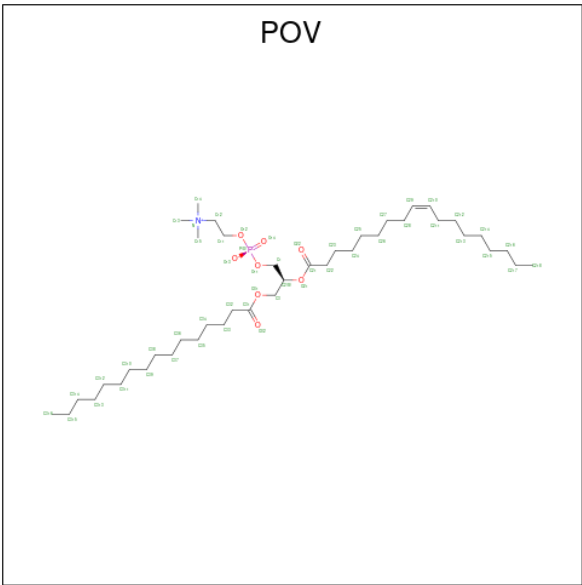
Chain	Residue	Modelled	Actual	Comment	Reference
C	796	GLY	-	linker	UNP Q8NET8
C	797	SER	-	linker	UNP Q8NET8
C	798	ALA	-	linker	UNP Q8NET8
C	799	ALA	-	linker	UNP Q8NET8
C	800	ALA	-	linker	UNP Q8NET8
C	801	ALA	-	linker	UNP Q8NET8
C	802	VAL	-	linker	UNP Q8NET8
C	865	LEU	PHE	conflict	UNP P42212
C	866	THR	SER	conflict	UNP P42212
C	1007	LYS	ALA	conflict	UNP P42212
C	1032	LEU	HIS	conflict	UNP P42212
C	1040	SER	-	expression tag	UNP P42212
C	1041	GLY	-	expression tag	UNP P42212
C	1042	LEU	-	expression tag	UNP P42212
C	1043	ARG	-	expression tag	UNP P42212
C	1044	SER	-	expression tag	UNP P42212
C	1045	TRP	-	expression tag	UNP P42212
C	1046	SER	-	expression tag	UNP P42212
C	1047	HIS	-	expression tag	UNP P42212
C	1048	PRO	-	expression tag	UNP P42212
C	1049	GLN	-	expression tag	UNP P42212
C	1050	PHE	-	expression tag	UNP P42212
C	1051	GLU	-	expression tag	UNP P42212
C	1052	LYS	-	expression tag	UNP P42212
D	568	ASP	GLY	engineered mutation	UNP Q8NET8
D	792	LEU	-	linker	UNP Q8NET8
D	793	VAL	-	linker	UNP Q8NET8
D	794	PRO	-	linker	UNP Q8NET8
D	795	ARG	-	linker	UNP Q8NET8
D	796	GLY	-	linker	UNP Q8NET8
D	797	SER	-	linker	UNP Q8NET8
D	798	ALA	-	linker	UNP Q8NET8
D	799	ALA	-	linker	UNP Q8NET8
D	800	ALA	-	linker	UNP Q8NET8
D	801	ALA	-	linker	UNP Q8NET8
D	802	VAL	-	linker	UNP Q8NET8
D	865	LEU	PHE	conflict	UNP P42212
D	866	THR	SER	conflict	UNP P42212
D	1007	LYS	ALA	conflict	UNP P42212
D	1032	LEU	HIS	conflict	UNP P42212
D	1040	SER	-	expression tag	UNP P42212
D	1041	GLY	-	expression tag	UNP P42212

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Chain	Residue	Modelled	Actual	Comment	Reference
D	1042	LEU	-	expression tag	UNP P42212
D	1043	ARG	-	expression tag	UNP P42212
D	1044	SER	-	expression tag	UNP P42212
D	1045	TRP	-	expression tag	UNP P42212
D	1046	SER	-	expression tag	UNP P42212
D	1047	HIS	-	expression tag	UNP P42212
D	1048	PRO	-	expression tag	UNP P42212
D	1049	GLN	-	expression tag	UNP P42212
D	1050	PHE	-	expression tag	UNP P42212
D	1051	GLU	-	expression tag	UNP P42212
D	1052	LYS	-	expression tag	UNP P42212

- Molecule 2 is (2S)-3-(hexadecanoyloxy)-2-[(9Z)-octadec-9-enoyloxy]propyl 2-(trimethylamm onio)ethyl phosphate (CCD ID: POV) (formula: C₄₂H₈₂NO₈P).



Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	A	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	A	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	A	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	B	1	Total	C	N	O	P	0
			52	42	1	8	1	

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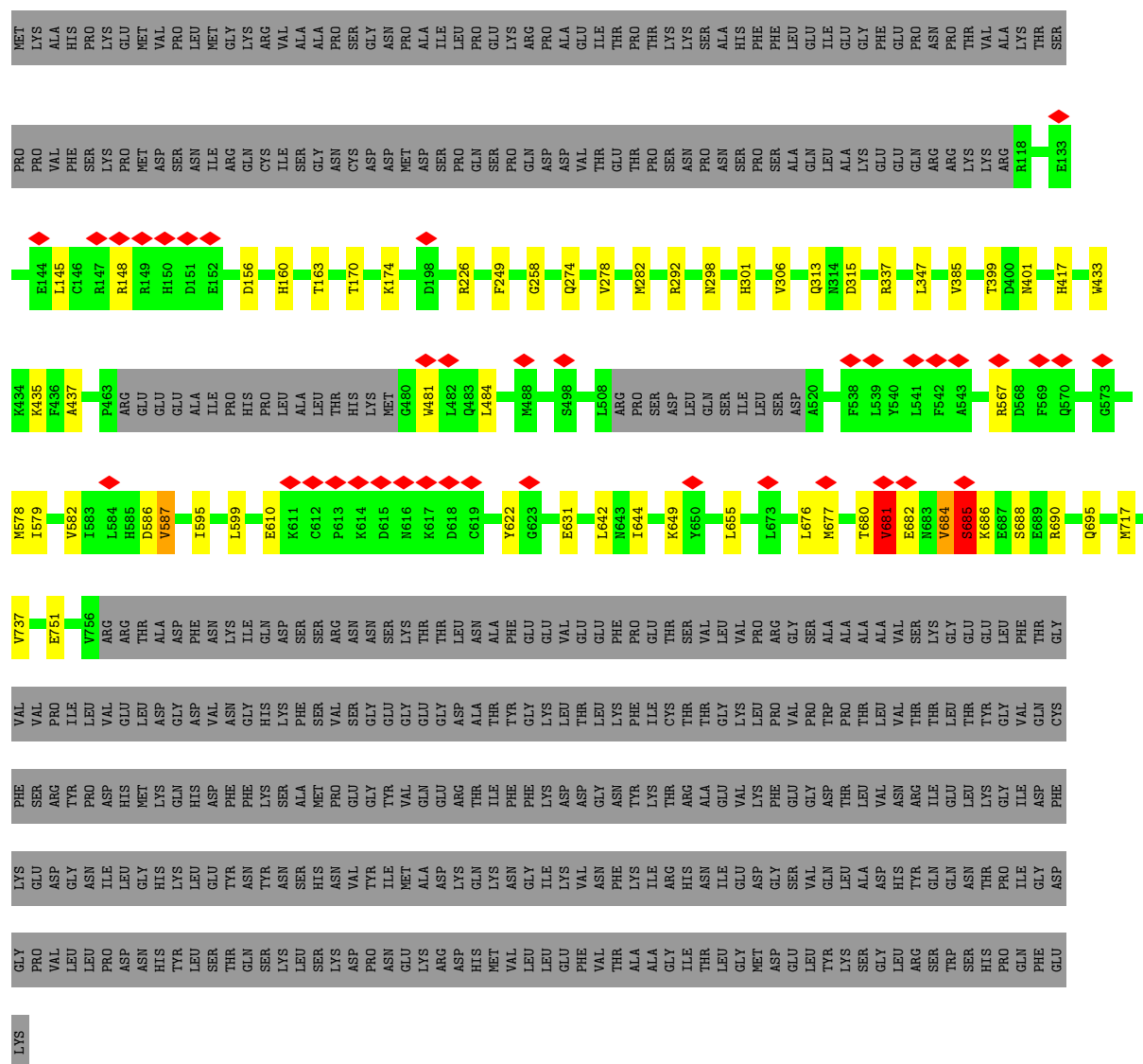
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Mol	Chain	Residues	Atoms					AltConf
2	B	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	C	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	C	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	C	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	D	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	D	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	D	1	Total	C	N	O	P	0
			52	42	1	8	1	

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

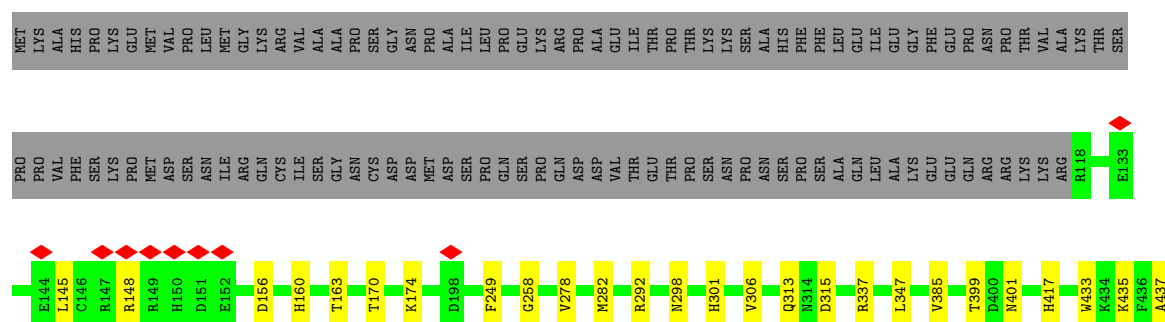
Mol	Chain	Residues	Atoms		AltConf
3	A	3	Total	Na	0
			3	3	

Chain B:



- Molecule 1: Isoform 2 of Transient receptor potential cation channel subfamily V member 3, Green fluorescent protein

Chain C:





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

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	45261	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.820	Depositor
Minimum map value	-0.400	Depositor
Average map value	0.010	Depositor
Map value standard deviation	0.027	Depositor
Recommended contour level	0.107	Depositor
Map size (Å)	251.136, 251.136, 251.136	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.981, 0.981, 0.981	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: POV, NA

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	0.34	0/5096	0.71	2/6893 (0.0%)
1	B	0.34	0/5096	0.71	2/6893 (0.0%)
1	C	0.35	0/5096	0.71	2/6893 (0.0%)
1	D	0.35	0/5096	0.70	2/6893 (0.0%)
All	All	0.34	0/20384	0.71	8/27572 (0.0%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	586	ASP	CA-C-N	5.91	132.60	121.97
1	B	586	ASP	C-N-CA	5.91	132.60	121.97
1	A	586	ASP	CA-C-N	5.89	132.57	121.97
1	A	586	ASP	C-N-CA	5.89	132.57	121.97
1	D	586	ASP	CA-C-N	5.89	132.57	121.97
1	D	586	ASP	C-N-CA	5.89	132.57	121.97
1	C	586	ASP	CA-C-N	5.88	132.54	121.97
1	C	586	ASP	C-N-CA	5.88	132.54	121.97

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	4987	0	5045	36	0
1	B	4987	0	5045	40	0
1	C	4987	0	5045	37	0
1	D	4987	0	5045	37	0
2	A	208	0	328	9	0
2	B	104	0	164	4	0
2	C	156	0	246	7	0
2	D	156	0	246	6	0
3	A	3	0	0	0	0
All	All	20575	0	21164	161	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 4.

All (161) 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:417:HIS:HB3	1:A:690:ARG:HG3	1.83	0.61
1:B:417:HIS:HB3	1:B:690:ARG:HG3	1.83	0.61
1:C:417:HIS:HB3	1:C:690:ARG:HG3	1.83	0.60
1:D:417:HIS:HB3	1:D:690:ARG:HG3	1.83	0.60
1:A:298:ASN:OD1	1:A:301:HIS:ND1	2.34	0.60
1:C:481:TRP:HB3	1:C:484:LEU:HB2	1.84	0.60
1:D:298:ASN:OD1	1:D:301:HIS:ND1	2.34	0.59
1:B:685:SER:O	1:B:688:SER:N	2.36	0.59
1:B:481:TRP:HB3	1:B:484:LEU:HB2	1.84	0.59
1:A:481:TRP:HB3	1:A:484:LEU:HB2	1.84	0.58
1:C:298:ASN:OD1	1:C:301:HIS:ND1	2.34	0.58
1:D:685:SER:O	1:D:688:SER:N	2.36	0.58
1:B:298:ASN:OD1	1:B:301:HIS:ND1	2.34	0.58
1:A:163:THR:HG22	1:A:170:THR:HG22	1.86	0.58
1:A:685:SER:O	1:A:688:SER:N	2.36	0.58
1:D:163:THR:HG22	1:D:170:THR:HG22	1.86	0.58
1:D:481:TRP:HB3	1:D:484:LEU:HB2	1.84	0.57
1:C:685:SER:O	1:C:688:SER:N	2.36	0.57
1:B:684:VAL:O	1:B:686:LYS:N	2.38	0.57
1:C:684:VAL:O	1:C:686:LYS:N	2.38	0.57
1:D:684:VAL:O	1:D:686:LYS:N	2.38	0.57
1:A:684:VAL:O	1:A:686:LYS:N	2.38	0.57
1:C:163:THR:HG22	1:C:170:THR:HG22	1.86	0.57
2:D:1101:POV:H36A	2:D:1101:POV:H27	1.87	0.56
1:B:163:THR:HG22	1:B:170:THR:HG22	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:684:VAL:O	1:D:685:SER:C	2.49	0.56
1:B:684:VAL:O	1:B:685:SER:C	2.49	0.56
2:C:1101:POV:H36A	2:C:1101:POV:H27	1.89	0.55
1:B:680:THR:O	1:B:681:VAL:C	2.50	0.55
2:A:1103:POV:H215	2:A:1103:POV:H31D	1.89	0.55
1:C:680:THR:O	1:C:681:VAL:C	2.50	0.55
1:B:751:GLU:OE2	1:C:174:LYS:NZ	2.36	0.55
1:C:684:VAL:O	1:C:685:SER:C	2.49	0.54
1:A:684:VAL:O	1:A:685:SER:C	2.50	0.54
1:D:680:THR:O	1:D:681:VAL:C	2.50	0.54
1:B:595:ILE:HG23	1:B:599:LEU:HD23	1.90	0.54
1:C:156:ASP:O	1:C:160:HIS:ND1	2.37	0.54
2:D:1103:POV:H31D	2:D:1103:POV:H215	1.90	0.54
1:A:174:LYS:NZ	1:D:751:GLU:OE2	2.38	0.54
1:A:680:THR:O	1:A:681:VAL:C	2.50	0.54
2:A:1104:POV:H36A	2:A:1104:POV:H27	1.90	0.54
1:A:156:ASP:O	1:A:160:HIS:ND1	2.37	0.53
1:C:595:ILE:HG23	1:C:599:LEU:HD23	1.90	0.53
1:A:595:ILE:HG23	1:A:599:LEU:HD23	1.90	0.53
1:C:680:THR:HG22	1:D:677:MET:HG2	1.91	0.53
1:D:595:ILE:HG23	1:D:599:LEU:HD23	1.90	0.53
1:A:680:THR:HG22	1:B:677:MET:HG2	1.91	0.53
1:C:567:ARG:HH12	1:C:695:GLN:HG3	1.74	0.53
1:D:156:ASP:O	1:D:160:HIS:ND1	2.37	0.53
1:B:567:ARG:HH12	1:B:695:GLN:HG3	1.74	0.53
1:B:680:THR:HG22	1:C:677:MET:HG2	1.91	0.52
1:A:751:GLU:OE2	1:B:174:LYS:NZ	2.39	0.52
2:A:1101:POV:H3	1:B:655:LEU:HD12	1.91	0.52
1:B:610:GLU:OE1	1:B:649:LYS:NZ	2.42	0.52
1:A:677:MET:HG2	1:D:680:THR:HG22	1.92	0.51
1:A:717:MET:SD	1:A:717:MET:N	2.84	0.51
1:B:717:MET:SD	1:B:717:MET:N	2.83	0.51
2:C:1103:POV:H31D	2:C:1103:POV:H215	1.91	0.51
1:D:717:MET:SD	1:D:717:MET:N	2.83	0.51
1:C:717:MET:SD	1:C:717:MET:N	2.83	0.51
1:D:567:ARG:HH12	1:D:695:GLN:HG3	1.74	0.51
1:B:433:TRP:HA	1:B:437:ALA:HB3	1.93	0.51
1:A:567:ARG:HH12	1:A:695:GLN:HG3	1.74	0.51
2:B:1102:POV:H31C	2:B:1102:POV:H215	1.93	0.51
1:C:655:LEU:HD12	2:C:1101:POV:H3	1.92	0.51
1:A:655:LEU:HD12	2:A:1104:POV:H3	1.92	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:655:LEU:HD12	2:D:1101:POV:H3	1.92	0.50
1:A:433:TRP:HA	1:A:437:ALA:HB3	1.93	0.50
1:C:433:TRP:HA	1:C:437:ALA:HB3	1.93	0.50
1:C:751:GLU:OE2	1:D:174:LYS:NZ	2.39	0.50
1:B:156:ASP:O	1:B:160:HIS:ND1	2.37	0.50
1:D:433:TRP:HA	1:D:437:ALA:HB3	1.93	0.49
1:A:610:GLU:OE1	1:A:649:LYS:NZ	2.42	0.49
1:C:676:LEU:O	1:C:680:THR:HG23	2.13	0.49
1:C:610:GLU:OE1	1:C:649:LYS:NZ	2.42	0.48
1:B:435:LYS:HD2	1:B:717:MET:HE3	1.96	0.48
1:B:676:LEU:O	1:B:680:THR:HG23	2.13	0.48
1:C:435:LYS:HD2	1:C:717:MET:HE3	1.96	0.48
1:D:676:LEU:O	1:D:680:THR:HG23	2.13	0.48
2:B:1102:POV:H3A	2:B:1102:POV:H32	1.49	0.48
1:A:676:LEU:O	1:A:680:THR:HG23	2.13	0.48
1:D:610:GLU:OE1	1:D:649:LYS:NZ	2.42	0.48
1:A:435:LYS:HD2	1:A:717:MET:HE3	1.96	0.47
1:D:435:LYS:HD2	1:D:717:MET:HE3	1.96	0.47
2:A:1103:POV:H3A	2:A:1103:POV:H32	1.49	0.47
1:D:579:ILE:HA	1:D:582:VAL:HG22	1.98	0.46
1:B:579:ILE:HA	1:B:582:VAL:HG22	1.98	0.46
1:B:578:MET:O	1:B:582:VAL:N	2.48	0.46
1:C:579:ILE:HA	1:C:582:VAL:HG22	1.98	0.46
1:A:591:LEU:HD12	2:A:1101:POV:H21C	1.97	0.46
1:D:578:MET:O	1:D:582:VAL:N	2.48	0.46
2:D:1103:POV:H3A	2:D:1103:POV:H32	1.49	0.46
1:A:578:MET:O	1:A:582:VAL:N	2.48	0.46
1:A:579:ILE:HA	1:A:582:VAL:HG22	1.98	0.46
1:C:278:VAL:HG12	1:C:282:MET:HE2	1.98	0.46
1:D:278:VAL:HG12	1:D:282:MET:HE2	1.98	0.46
1:B:278:VAL:HG12	1:B:282:MET:HE2	1.98	0.45
2:C:1103:POV:H38	2:C:1103:POV:H31B	1.84	0.45
1:C:399:THR:OG1	1:C:401:ASN:O	2.29	0.45
2:C:1103:POV:H22A	2:C:1103:POV:H2	1.60	0.45
1:D:306:VAL:O	1:D:313:GLN:NE2	2.47	0.45
1:A:278:VAL:HG12	1:A:282:MET:HE2	1.98	0.45
1:C:644:ILE:HG12	1:C:655:LEU:HD23	2.00	0.44
1:B:680:THR:O	1:B:682:GLU:N	2.51	0.44
1:C:680:THR:O	1:C:682:GLU:N	2.50	0.44
2:C:1101:POV:H29	2:C:1101:POV:H26A	1.85	0.44
2:B:1102:POV:H2	2:B:1102:POV:H22A	1.61	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:249:PHE:CE1	1:C:258:GLY:HA2	2.53	0.44
1:C:578:MET:O	1:C:582:VAL:N	2.49	0.44
1:D:249:PHE:CE1	1:D:258:GLY:HA2	2.53	0.44
1:D:680:THR:O	1:D:682:GLU:N	2.50	0.44
2:D:1103:POV:H38	2:D:1103:POV:H31B	1.84	0.44
1:A:249:PHE:CE1	1:A:258:GLY:HA2	2.53	0.44
1:C:306:VAL:O	1:C:313:GLN:NE2	2.47	0.44
1:D:585:HIS:O	1:D:589:LYS:NZ	2.43	0.44
1:B:301:HIS:HD2	1:B:347:LEU:HD13	1.83	0.43
1:A:680:THR:O	1:A:682:GLU:N	2.51	0.43
2:C:1103:POV:H3A	2:C:1103:POV:H32	1.50	0.43
1:D:644:ILE:HG12	1:D:655:LEU:HD23	1.99	0.43
1:B:644:ILE:HG12	1:B:655:LEU:HD23	2.00	0.43
1:D:301:HIS:HD2	1:D:347:LEU:HD13	1.83	0.43
1:B:306:VAL:O	1:B:313:GLN:NE2	2.47	0.43
1:B:399:THR:OG1	1:B:401:ASN:O	2.29	0.43
1:B:292:ARG:NH1	1:B:337:ARG:O	2.52	0.43
1:B:249:PHE:CE1	1:B:258:GLY:HA2	2.53	0.43
1:A:292:ARG:NH1	1:A:337:ARG:O	2.52	0.43
1:C:292:ARG:NH1	1:C:337:ARG:O	2.52	0.43
1:C:301:HIS:HD2	1:C:347:LEU:HD13	1.83	0.43
1:A:644:ILE:HG12	1:A:655:LEU:HD23	1.99	0.42
2:D:1103:POV:H22A	2:D:1103:POV:H2	1.61	0.42
1:A:301:HIS:HD2	1:A:347:LEU:HD13	1.83	0.42
1:D:292:ARG:NH1	1:D:337:ARG:O	2.52	0.42
1:A:685:SER:O	1:A:686:LYS:C	2.63	0.42
1:B:599:LEU:HB3	2:B:1102:POV:H311	2.02	0.42
1:B:685:SER:O	1:B:686:LYS:C	2.62	0.41
1:C:685:SER:O	1:C:686:LYS:C	2.62	0.41
1:A:399:THR:OG1	1:A:401:ASN:O	2.29	0.41
1:C:674:ILE:O	1:C:678:GLY:N	2.54	0.41
1:A:306:VAL:O	1:A:313:GLN:NE2	2.47	0.41
1:C:622:TYR:HE1	1:C:631:GLU:HB2	1.86	0.41
1:B:622:TYR:HE1	1:B:631:GLU:HB2	1.86	0.41
1:D:681:VAL:O	1:D:682:GLU:C	2.63	0.41
1:D:385:VAL:HG22	1:D:737:VAL:HG12	2.02	0.41
1:B:681:VAL:O	1:B:682:GLU:C	2.63	0.41
1:C:385:VAL:HG22	1:C:737:VAL:HG12	2.02	0.41
1:D:399:THR:OG1	1:D:401:ASN:O	2.29	0.41
1:D:622:TYR:HE1	1:D:631:GLU:HB2	1.85	0.41
1:A:585:HIS:O	1:A:589:LYS:NZ	2.43	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:1103:POV:H22A	2:A:1103:POV:H2	1.61	0.41
1:B:587:VAL:H	1:B:587:VAL:HG23	1.55	0.41
1:C:145:LEU:HA	1:C:148:ARG:HH21	1.86	0.41
1:B:145:LEU:HA	1:B:148:ARG:HH21	1.86	0.41
1:A:639:LEU:HD13	1:B:642:LEU:HB2	2.03	0.40
1:B:385:VAL:HG22	1:B:737:VAL:HG12	2.02	0.40
1:C:681:VAL:O	1:C:682:GLU:C	2.63	0.40
1:A:681:VAL:O	1:A:682:GLU:C	2.63	0.40
2:A:1104:POV:H214	2:A:1104:POV:H38A	2.03	0.40
1:D:570:GLN:HB3	1:D:571:SER:H	1.78	0.40
1:D:674:ILE:O	1:D:678:GLY:N	2.54	0.40
2:A:1104:POV:H29	2:A:1104:POV:H26A	1.84	0.40
1:B:226:ARG:HD2	1:B:274:GLN:HE21	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	606/1052 (58%)	557 (92%)	44 (7%)	5 (1%)	16	46
1	B	606/1052 (58%)	557 (92%)	44 (7%)	5 (1%)	16	46
1	C	606/1052 (58%)	557 (92%)	44 (7%)	5 (1%)	16	46
1	D	606/1052 (58%)	557 (92%)	44 (7%)	5 (1%)	16	46
All	All	2424/4208 (58%)	2228 (92%)	176 (7%)	20 (1%)	18	46

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	315	ASP
1	A	681	VAL

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Mol	Chain	Res	Type
1	A	685	SER
1	B	315	ASP
1	B	681	VAL
1	B	685	SER
1	C	315	ASP
1	C	681	VAL
1	C	685	SER
1	D	315	ASP
1	D	681	VAL
1	D	685	SER
1	A	587	VAL
1	A	684	VAL
1	B	587	VAL
1	B	684	VAL
1	C	587	VAL
1	C	684	VAL
1	D	587	VAL
1	D	684	VAL

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	542/930 (58%)	540 (100%)	2 (0%)	84	83
1	B	542/930 (58%)	540 (100%)	2 (0%)	84	83
1	C	542/930 (58%)	540 (100%)	2 (0%)	84	83
1	D	542/930 (58%)	540 (100%)	2 (0%)	84	83
All	All	2168/3720 (58%)	2160 (100%)	8 (0%)	81	83

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	681	VAL
1	A	685	SER
1	B	681	VAL

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Mol	Chain	Res	Type
1	B	685	SER
1	C	681	VAL
1	C	685	SER
1	D	681	VAL
1	D	685	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	220	ASN
1	A	244	HIS
1	A	410	ASN
1	A	439	HIS
1	A	452	ASN
1	A	570	GLN
1	A	646	GLN
1	A	647	ASN
1	A	671	ASN
1	A	745	HIS
1	A	750	ASN
1	B	220	ASN
1	B	244	HIS
1	B	410	ASN
1	B	452	ASN
1	B	570	GLN
1	B	647	ASN
1	B	671	ASN
1	B	745	HIS
1	B	750	ASN
1	C	220	ASN
1	C	244	HIS
1	C	410	ASN
1	C	570	GLN
1	C	585	HIS
1	C	671	ASN
1	C	745	HIS
1	C	750	ASN
1	D	220	ASN
1	D	244	HIS
1	D	410	ASN
1	D	452	ASN
1	D	570	GLN

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Mol	Chain	Res	Type
1	D	585	HIS
1	D	647	ASN
1	D	671	ASN
1	D	745	HIS
1	D	750	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 15 ligands modelled in this entry, 3 are monoatomic - leaving 12 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	$\# Z > 2$
2	POV	A	1104	-	51,51,51	0.30	0	57,59,59	0.34	0
2	POV	B	1102	-	51,51,51	0.33	0	57,59,59	0.46	0
2	POV	A	1103	-	51,51,51	0.33	0	57,59,59	0.47	0
2	POV	C	1101	-	51,51,51	0.30	0	57,59,59	0.34	0
2	POV	D	1102	-	51,51,51	0.29	0	57,59,59	0.26	0
2	POV	D	1103	-	51,51,51	0.33	0	57,59,59	0.46	0
2	POV	C	1103	-	51,51,51	0.33	0	57,59,59	0.47	0
2	POV	D	1101	-	51,51,51	0.30	0	57,59,59	0.34	0
2	POV	A	1101	-	51,51,51	0.30	0	57,59,59	0.35	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	POV	A	1102	-	51,51,51	0.29	0	57,59,59	0.26	0
2	POV	B	1101	-	51,51,51	0.29	0	57,59,59	0.26	0
2	POV	C	1102	-	51,51,51	0.29	0	57,59,59	0.26	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	POV	A	1104	-	-	16/55/55/55	-
2	POV	B	1102	-	-	23/55/55/55	-
2	POV	A	1103	-	-	23/55/55/55	-
2	POV	C	1101	-	-	15/55/55/55	-
2	POV	D	1102	-	-	9/55/55/55	-
2	POV	D	1103	-	-	22/55/55/55	-
2	POV	C	1103	-	-	22/55/55/55	-
2	POV	D	1101	-	-	16/55/55/55	-
2	POV	A	1101	-	-	15/55/55/55	-
2	POV	A	1102	-	-	11/55/55/55	-
2	POV	B	1101	-	-	10/55/55/55	-
2	POV	C	1102	-	-	10/55/55/55	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (192) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1101	POV	C1-O11-P-O12
2	A	1101	POV	O12-C11-C12-N
2	A	1101	POV	C32-C31-O31-C3
2	A	1101	POV	O32-C31-O31-C3
2	A	1103	POV	C1-O11-P-O12
2	A	1103	POV	C1-O11-P-O13
2	A	1103	POV	C22-C21-O21-C2
2	A	1103	POV	O22-C21-O21-C2
2	A	1103	POV	C32-C31-O31-C3

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Mol	Chain	Res	Type	Atoms
2	A	1103	POV	O32-C31-O31-C3
2	A	1104	POV	C1-O11-P-O12
2	A	1104	POV	O12-C11-C12-N
2	A	1104	POV	C32-C31-O31-C3
2	A	1104	POV	O32-C31-O31-C3
2	B	1102	POV	C1-O11-P-O12
2	B	1102	POV	C1-O11-P-O13
2	B	1102	POV	C22-C21-O21-C2
2	B	1102	POV	O22-C21-O21-C2
2	B	1102	POV	C32-C31-O31-C3
2	B	1102	POV	O32-C31-O31-C3
2	C	1101	POV	C1-O11-P-O12
2	C	1101	POV	O12-C11-C12-N
2	C	1101	POV	C32-C31-O31-C3
2	C	1101	POV	O32-C31-O31-C3
2	C	1103	POV	C1-O11-P-O12
2	C	1103	POV	C1-O11-P-O13
2	C	1103	POV	C22-C21-O21-C2
2	C	1103	POV	O22-C21-O21-C2
2	C	1103	POV	C32-C31-O31-C3
2	C	1103	POV	O32-C31-O31-C3
2	D	1101	POV	C1-O11-P-O12
2	D	1101	POV	O12-C11-C12-N
2	D	1101	POV	C32-C31-O31-C3
2	D	1101	POV	O32-C31-O31-C3
2	D	1103	POV	C1-O11-P-O12
2	D	1103	POV	C1-O11-P-O13
2	D	1103	POV	C22-C21-O21-C2
2	D	1103	POV	O22-C21-O21-C2
2	D	1103	POV	C32-C31-O31-C3
2	D	1103	POV	O32-C31-O31-C3
2	D	1101	POV	C37-C38-C39-C310
2	B	1101	POV	C36-C37-C38-C39
2	C	1102	POV	C36-C37-C38-C39
2	D	1102	POV	C23-C24-C25-C26
2	D	1101	POV	C310-C311-C312-C313
2	D	1102	POV	C36-C37-C38-C39
2	C	1101	POV	C310-C311-C312-C313
2	A	1101	POV	C310-C311-C312-C313
2	A	1104	POV	C310-C311-C312-C313
2	B	1101	POV	C23-C24-C25-C26
2	C	1101	POV	C37-C38-C39-C310

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Mol	Chain	Res	Type	Atoms
2	C	1102	POV	C23-C24-C25-C26
2	A	1102	POV	C36-C37-C38-C39
2	A	1102	POV	C34-C35-C36-C37
2	A	1104	POV	C37-C38-C39-C310
2	D	1101	POV	C35-C36-C37-C38
2	D	1102	POV	C34-C35-C36-C37
2	C	1101	POV	C35-C36-C37-C38
2	A	1104	POV	O11-C1-C2-C3
2	A	1104	POV	C35-C36-C37-C38
2	B	1101	POV	C1-C2-C3-O31
2	C	1102	POV	C1-C2-C3-O31
2	B	1101	POV	C34-C35-C36-C37
2	C	1102	POV	C34-C35-C36-C37
2	B	1102	POV	C36-C37-C38-C39
2	A	1103	POV	C26-C27-C28-C29
2	C	1102	POV	C35-C36-C37-C38
2	A	1103	POV	C27-C28-C29-C210
2	C	1103	POV	C27-C28-C29-C210
2	D	1103	POV	C27-C28-C29-C210
2	A	1101	POV	O11-C1-C2-C3
2	B	1102	POV	C27-C28-C29-C210
2	B	1101	POV	C35-C36-C37-C38
2	A	1102	POV	C1-C2-C3-O31
2	D	1102	POV	C1-C2-C3-O31
2	A	1101	POV	O11-C1-C2-O21
2	A	1104	POV	O11-C1-C2-O21
2	C	1101	POV	O11-C1-C2-O21
2	D	1101	POV	O11-C1-C2-O21
2	A	1102	POV	O21-C2-C3-O31
2	D	1102	POV	O21-C2-C3-O31
2	D	1103	POV	C26-C27-C28-C29
2	A	1103	POV	C11-C12-N-C13
2	D	1103	POV	C11-C12-N-C13
2	B	1102	POV	C26-C27-C28-C29
2	C	1103	POV	C26-C27-C28-C29
2	C	1101	POV	O11-C1-C2-C3
2	D	1101	POV	O11-C1-C2-C3
2	D	1103	POV	C311-C312-C313-C314
2	C	1103	POV	C311-C312-C313-C314
2	B	1101	POV	O21-C2-C3-O31
2	C	1102	POV	O21-C2-C3-O31
2	A	1103	POV	C311-C312-C313-C314

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Mol	Chain	Res	Type	Atoms
2	A	1103	POV	C11-C12-N-C14
2	B	1102	POV	C11-C12-N-C13
2	B	1102	POV	C11-C12-N-C14
2	C	1103	POV	C11-C12-N-C13
2	C	1103	POV	C11-C12-N-C14
2	D	1103	POV	C11-C12-N-C14
2	A	1102	POV	O12-C11-C12-N
2	B	1101	POV	O12-C11-C12-N
2	C	1102	POV	O12-C11-C12-N
2	D	1102	POV	O12-C11-C12-N
2	B	1102	POV	C310-C311-C312-C313
2	B	1102	POV	C2-C1-O11-P
2	D	1102	POV	C35-C36-C37-C38
2	B	1102	POV	C34-C35-C36-C37
2	A	1101	POV	C1-O11-P-O14
2	A	1103	POV	C11-C12-N-C15
2	A	1103	POV	C11-O12-P-O14
2	A	1104	POV	C1-O11-P-O14
2	B	1102	POV	C11-C12-N-C15
2	C	1101	POV	C1-O11-P-O14
2	C	1103	POV	C11-C12-N-C15
2	D	1101	POV	C1-O11-P-O14
2	D	1103	POV	C11-C12-N-C15
2	A	1103	POV	C22-C23-C24-C25
2	A	1103	POV	C2-C1-O11-P
2	C	1103	POV	C2-C1-O11-P
2	D	1103	POV	C2-C1-O11-P
2	D	1101	POV	C39-C310-C311-C312
2	C	1102	POV	C29-C210-C211-C212
2	B	1101	POV	C29-C210-C211-C212
2	D	1102	POV	C29-C210-C211-C212
2	C	1103	POV	C36-C37-C38-C39
2	D	1103	POV	C36-C37-C38-C39
2	A	1102	POV	C21-C22-C23-C24
2	A	1102	POV	C29-C210-C211-C212
2	C	1103	POV	C29-C210-C211-C212
2	D	1103	POV	C22-C23-C24-C25
2	A	1102	POV	C23-C24-C25-C26
2	C	1101	POV	C39-C310-C311-C312
2	D	1101	POV	C2-C1-O11-P
2	D	1103	POV	C29-C210-C211-C212
2	A	1103	POV	C36-C37-C38-C39

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Mol	Chain	Res	Type	Atoms
2	A	1104	POV	C39-C310-C311-C312
2	B	1102	POV	C22-C23-C24-C25
2	C	1103	POV	C22-C23-C24-C25
2	A	1103	POV	C29-C210-C211-C212
2	C	1101	POV	C2-C1-O11-P
2	A	1102	POV	C35-C36-C37-C38
2	C	1101	POV	C33-C34-C35-C36
2	B	1102	POV	C29-C210-C211-C212
2	A	1104	POV	C2-C1-O11-P
2	B	1102	POV	C311-C312-C313-C314
2	A	1104	POV	C33-C34-C35-C36
2	A	1101	POV	C11-C12-N-C13
2	A	1101	POV	C33-C34-C35-C36
2	C	1103	POV	O21-C21-C22-C23
2	A	1103	POV	O21-C21-C22-C23
2	D	1103	POV	O21-C21-C22-C23
2	B	1102	POV	O21-C21-C22-C23
2	C	1102	POV	C25-C26-C27-C28
2	A	1104	POV	O21-C21-C22-C23
2	C	1101	POV	O21-C21-C22-C23
2	D	1101	POV	O21-C21-C22-C23
2	D	1103	POV	O31-C31-C32-C33
2	A	1101	POV	O21-C21-C22-C23
2	B	1102	POV	O31-C31-C32-C33
2	D	1101	POV	C33-C34-C35-C36
2	A	1103	POV	O31-C31-C32-C33
2	C	1103	POV	O31-C31-C32-C33
2	A	1101	POV	C11-C12-N-C14
2	B	1102	POV	C37-C38-C39-C310
2	A	1101	POV	C2-C1-O11-P
2	D	1103	POV	O22-C21-C22-C23
2	A	1104	POV	C25-C26-C27-C28
2	D	1103	POV	C34-C35-C36-C37
2	A	1103	POV	O22-C21-C22-C23
2	B	1101	POV	C25-C26-C27-C28
2	C	1103	POV	C34-C35-C36-C37
2	B	1102	POV	O22-C21-C22-C23
2	A	1101	POV	C11-C12-N-C15
2	A	1101	POV	O22-C21-C22-C23
2	C	1103	POV	C39-C310-C311-C312
2	A	1102	POV	O21-C21-C22-C23
2	A	1103	POV	C39-C310-C311-C312

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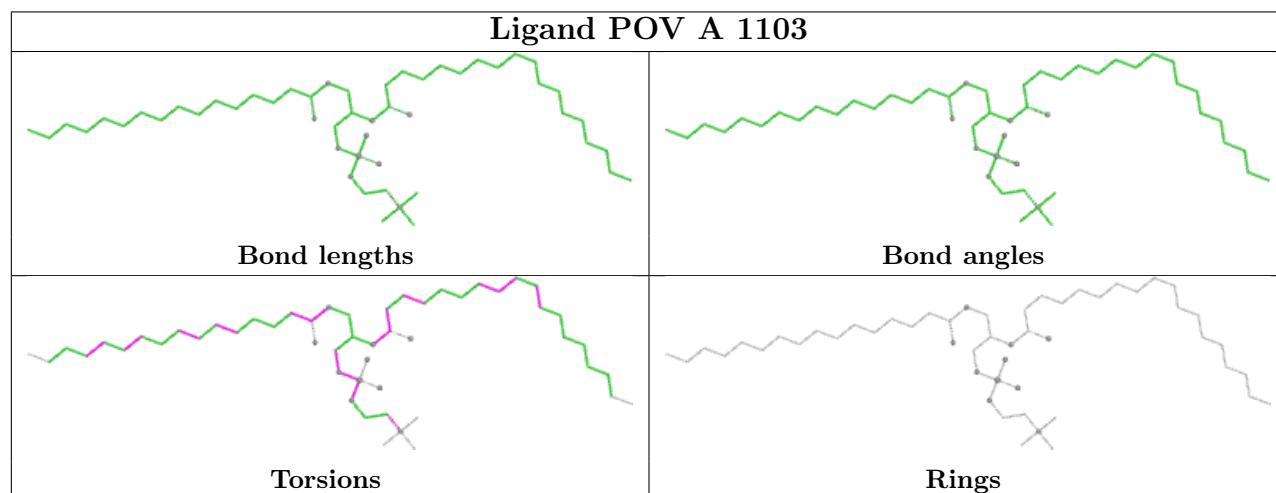
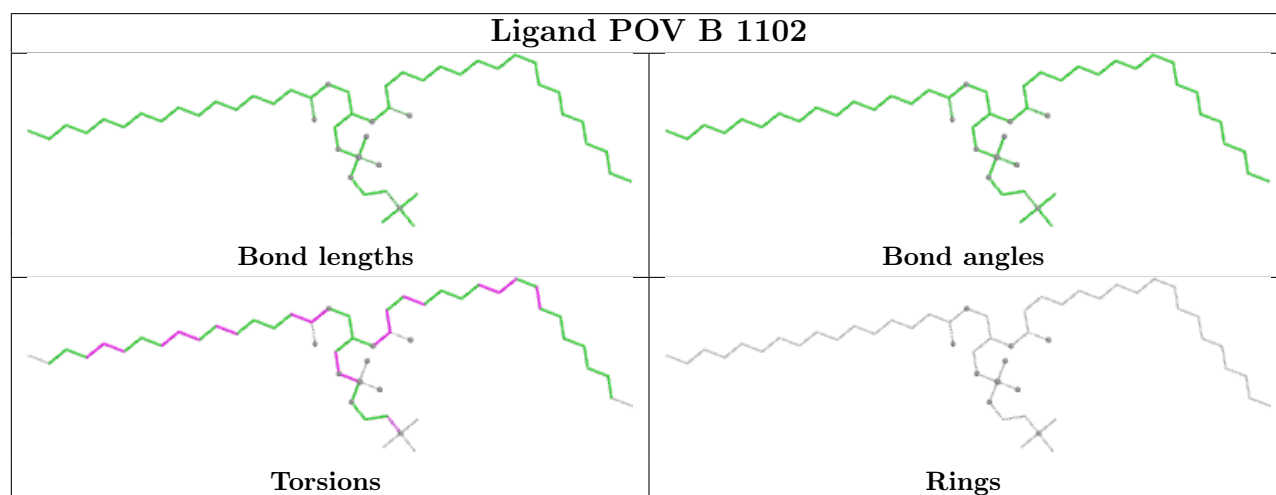
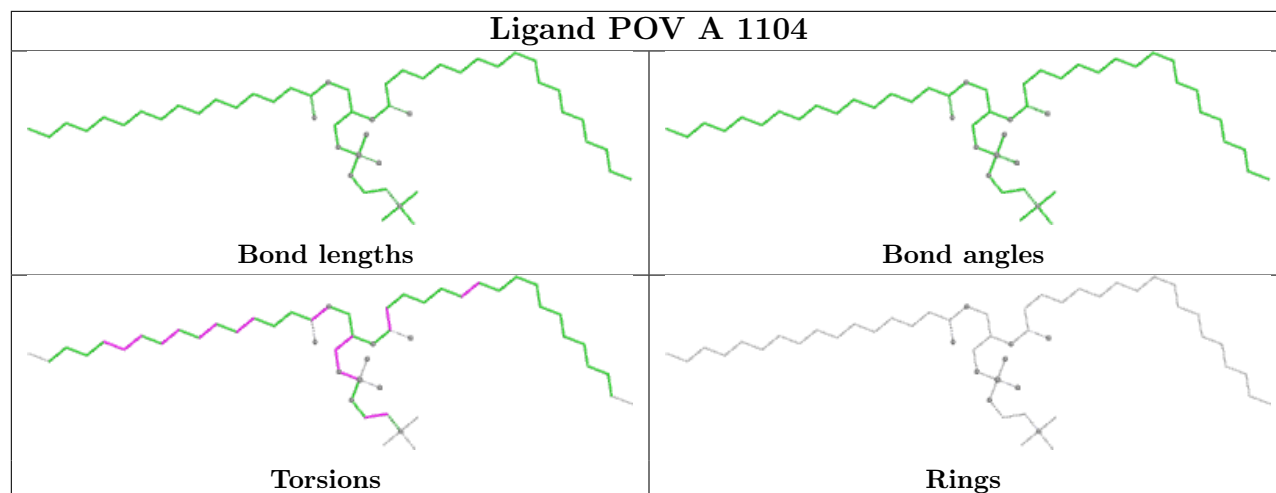
Mol	Chain	Res	Type	Atoms
2	A	1103	POV	C34-C35-C36-C37
2	A	1104	POV	O22-C21-C22-C23
2	C	1101	POV	O22-C21-C22-C23
2	D	1101	POV	O22-C21-C22-C23
2	B	1102	POV	O32-C31-C32-C33
2	A	1102	POV	C2-C1-O11-P
2	B	1101	POV	C2-C1-O11-P
2	D	1102	POV	C2-C1-O11-P
2	C	1103	POV	O22-C21-C22-C23
2	D	1101	POV	C25-C26-C27-C28
2	D	1103	POV	O32-C31-C32-C33
2	C	1103	POV	O32-C31-C32-C33
2	C	1102	POV	C39-C310-C311-C312
2	A	1103	POV	O32-C31-C32-C33
2	D	1103	POV	C39-C310-C311-C312

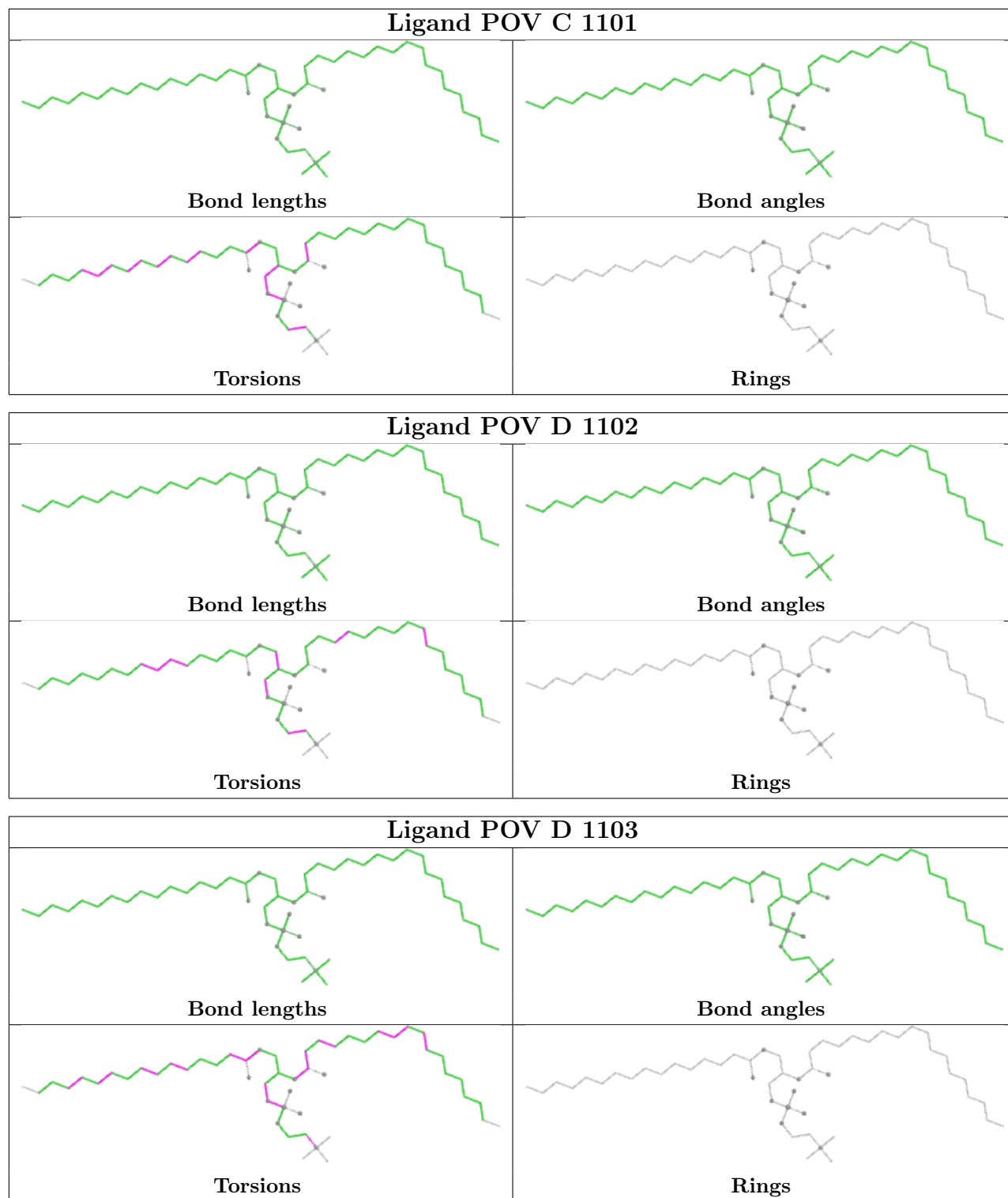
There are no ring outliers.

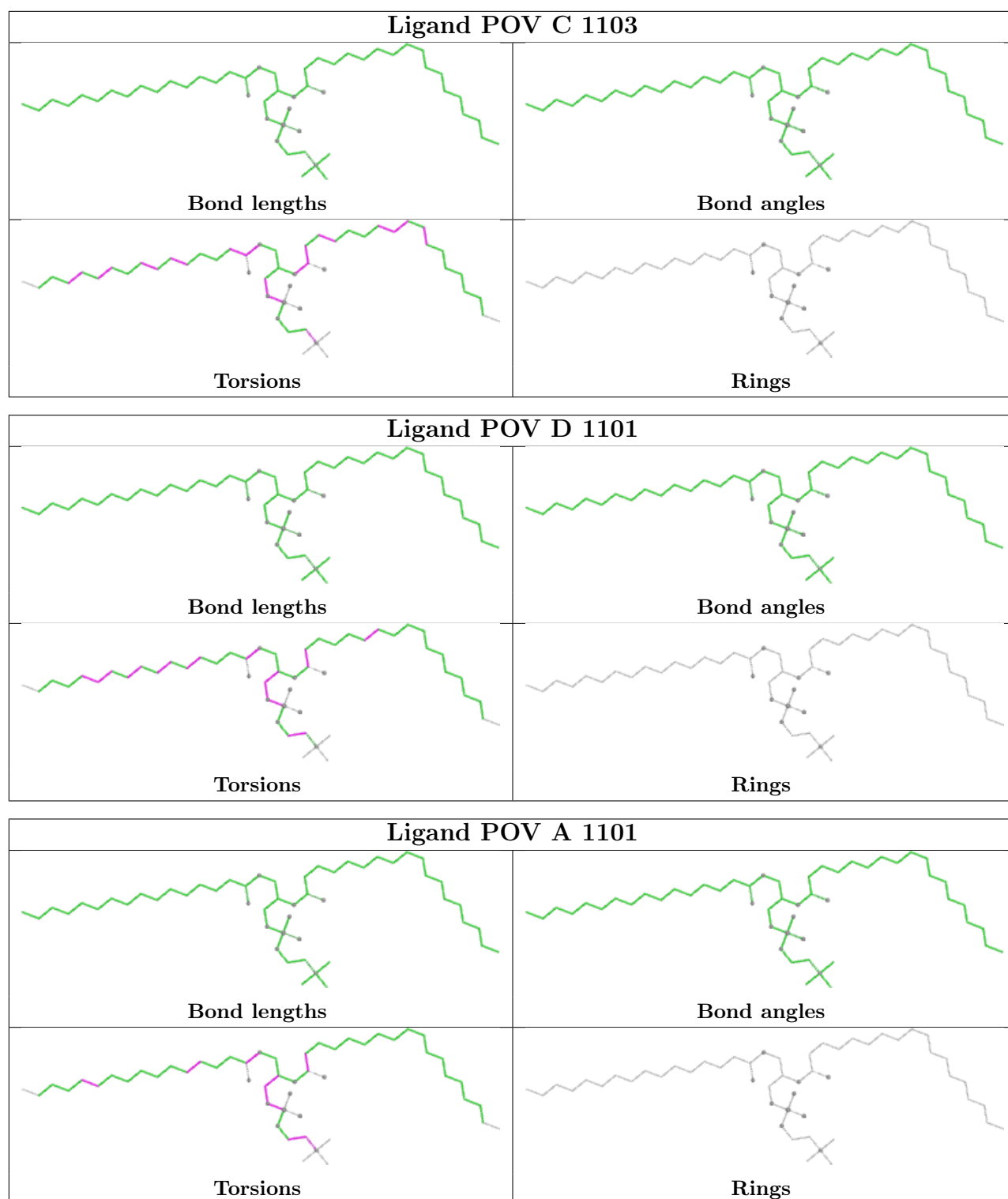
8 monomers are involved in 26 short contacts:

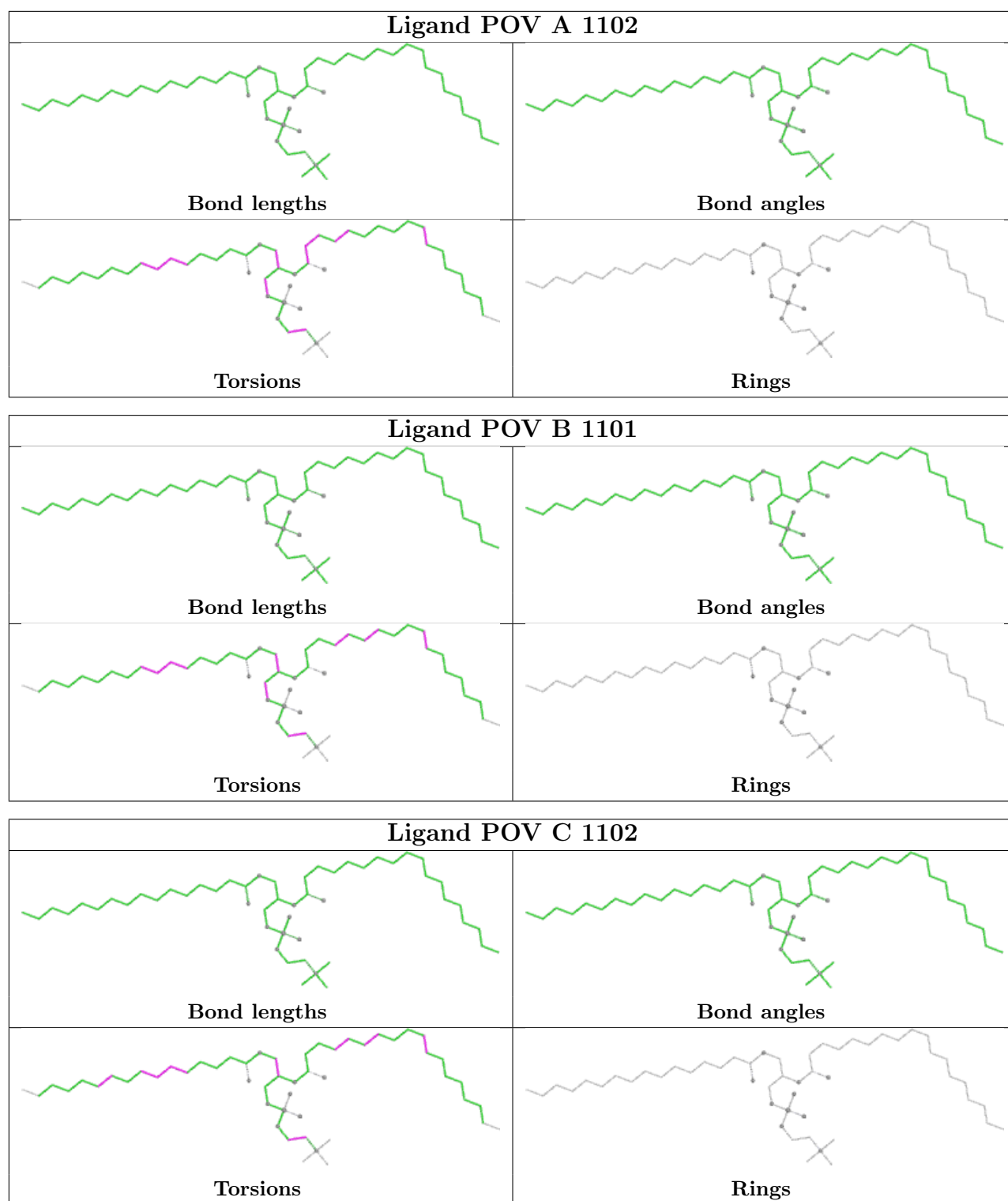
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1104	POV	4	0
2	B	1102	POV	4	0
2	A	1103	POV	3	0
2	C	1101	POV	3	0
2	D	1103	POV	4	0
2	C	1103	POV	4	0
2	D	1101	POV	2	0
2	A	1101	POV	2	0

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









5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

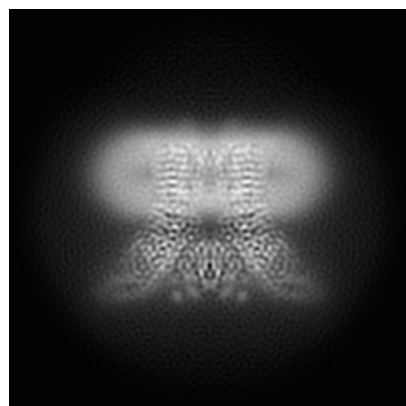
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-77142. These allow visual inspection of the internal detail of the map and identification of artifacts.

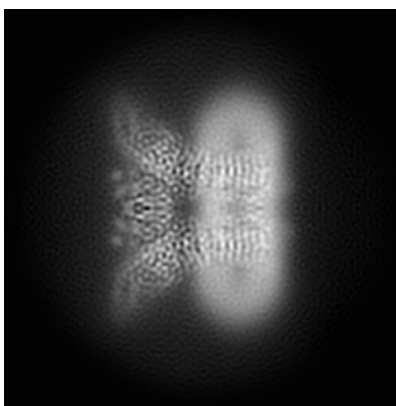
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

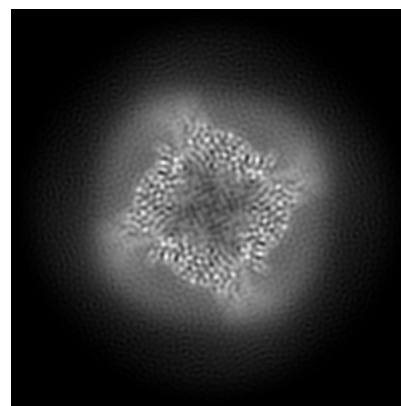
6.1.1 Primary map



X

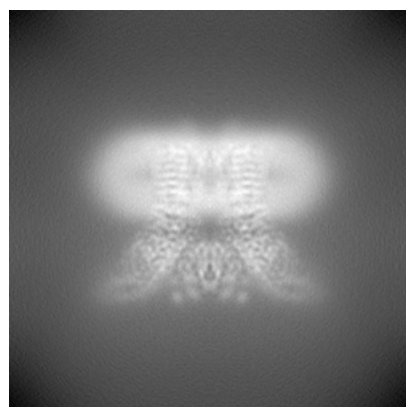


Y

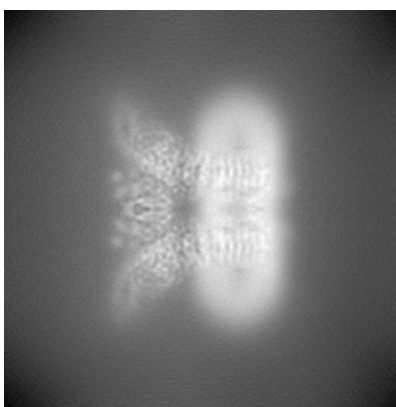


Z

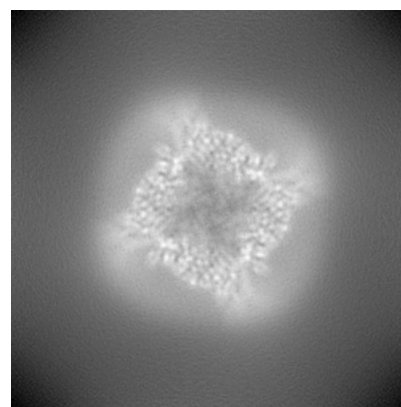
6.1.2 Raw map



X



Y

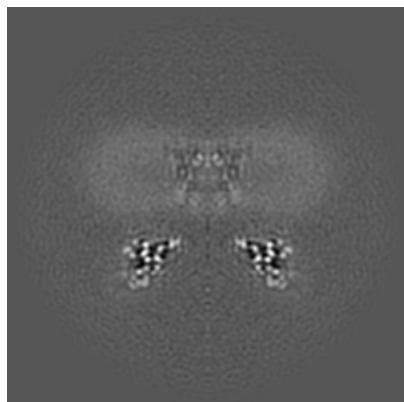


Z

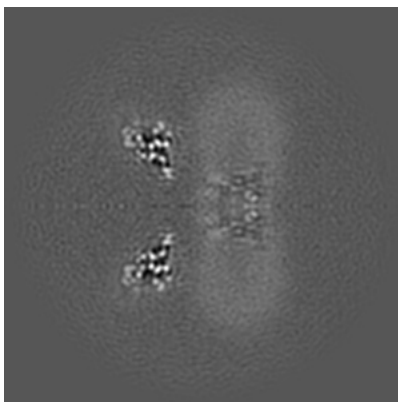
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

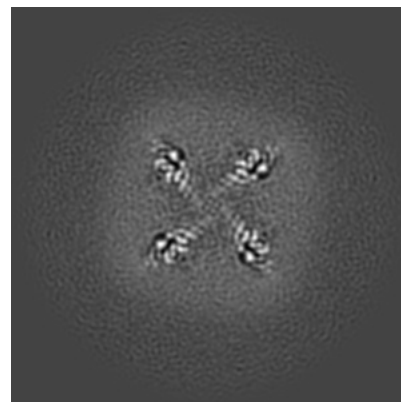
6.2.1 Primary map



X Index: 128

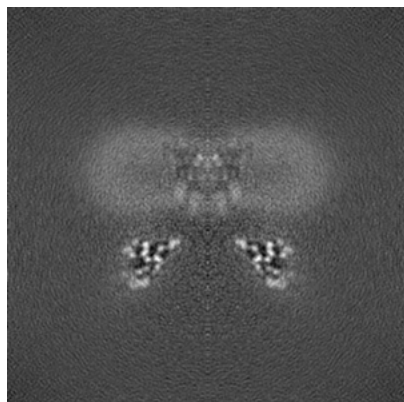


Y Index: 128

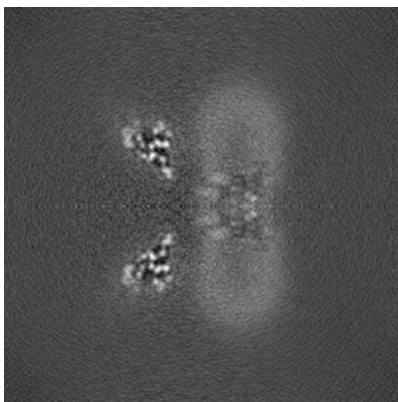


Z Index: 128

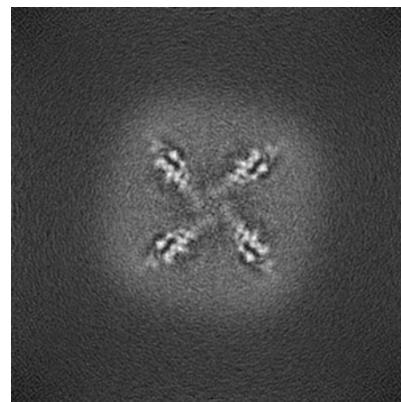
6.2.2 Raw map



X Index: 128



Y Index: 128

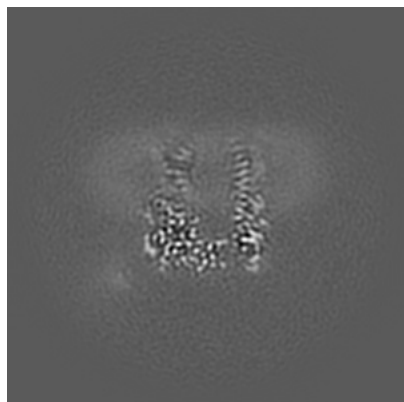


Z Index: 128

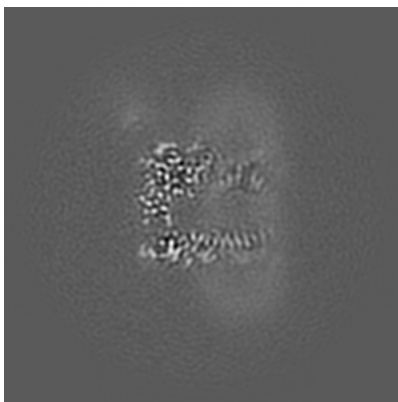
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

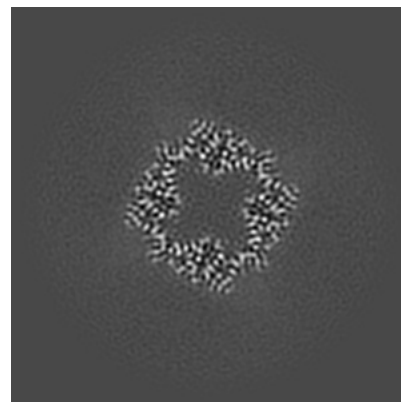
6.3.1 Primary map



X Index: 157

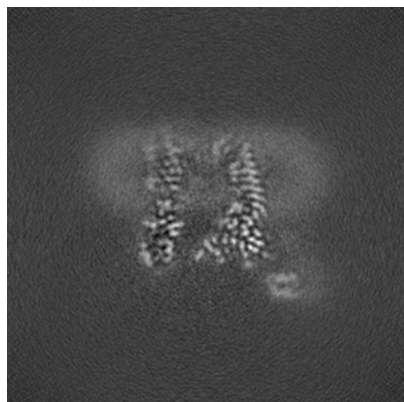


Y Index: 157

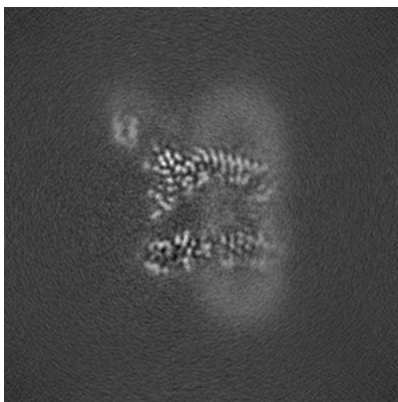


Z Index: 99

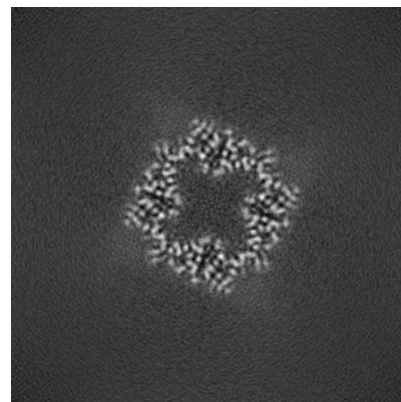
6.3.2 Raw map



X Index: 105



Y Index: 151

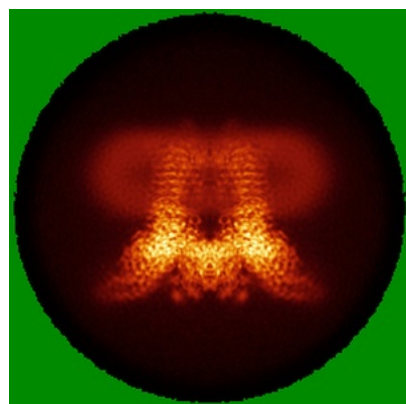


Z Index: 99

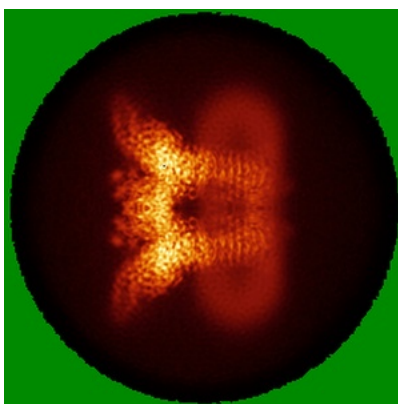
The images above show the largest variance slices of the map in three orthogonal directions.

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

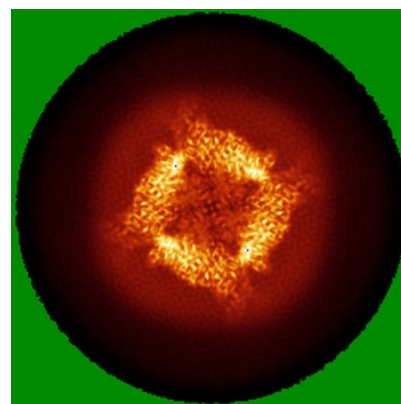
6.4.1 Primary map



X

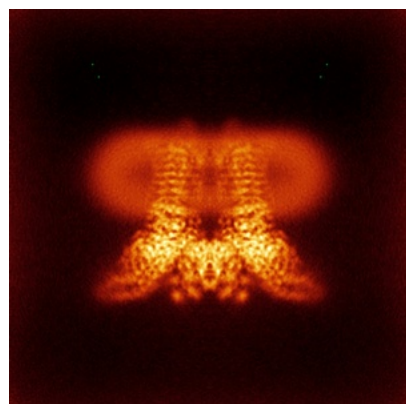


Y

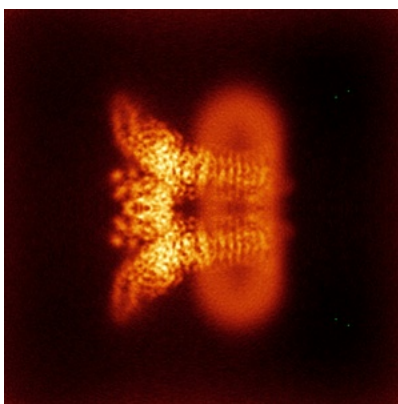


Z

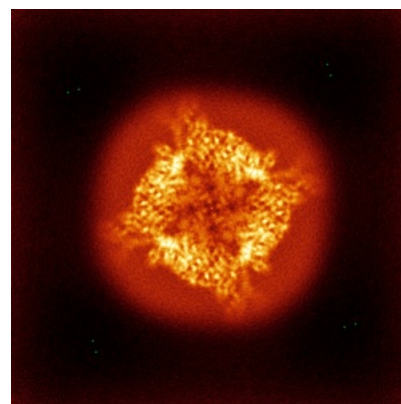
6.4.2 Raw map



X



Y

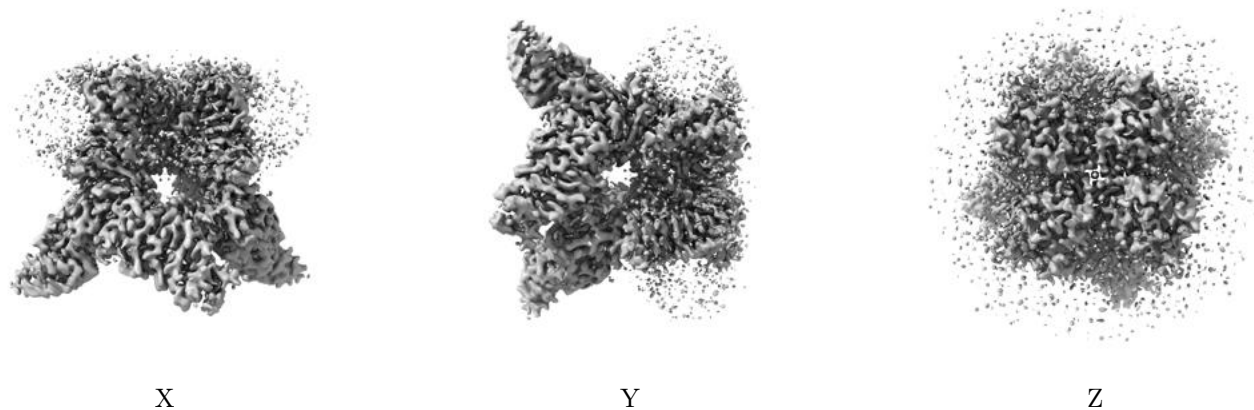


Z

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.

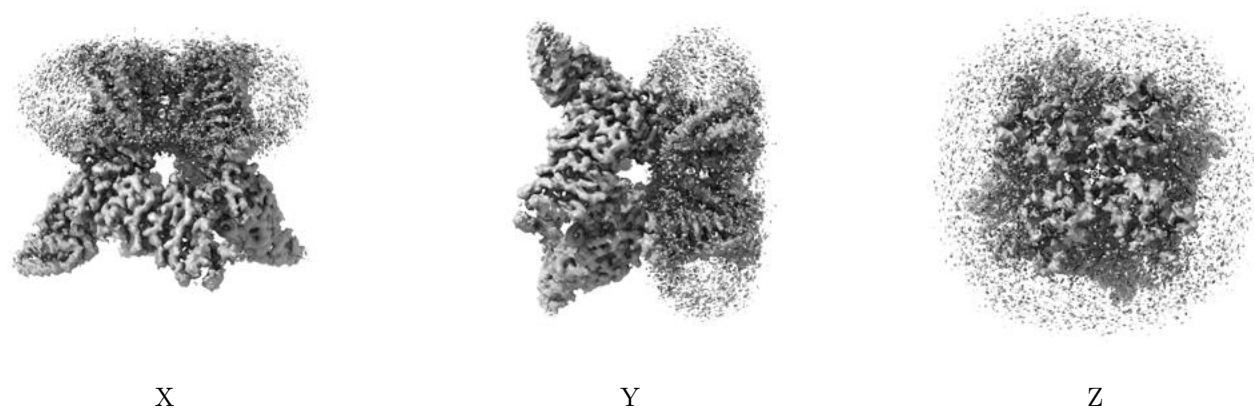
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.107. 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.

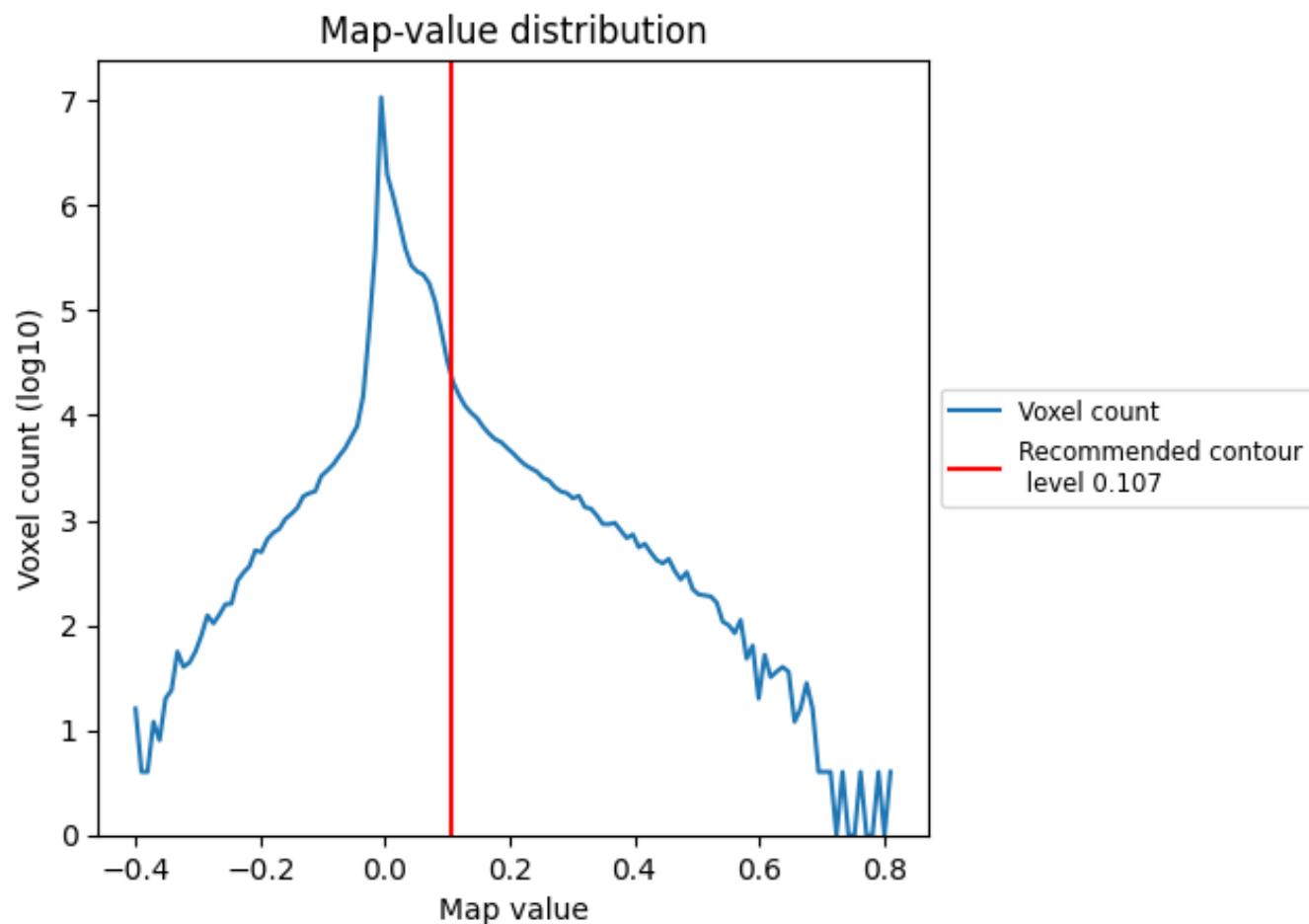
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

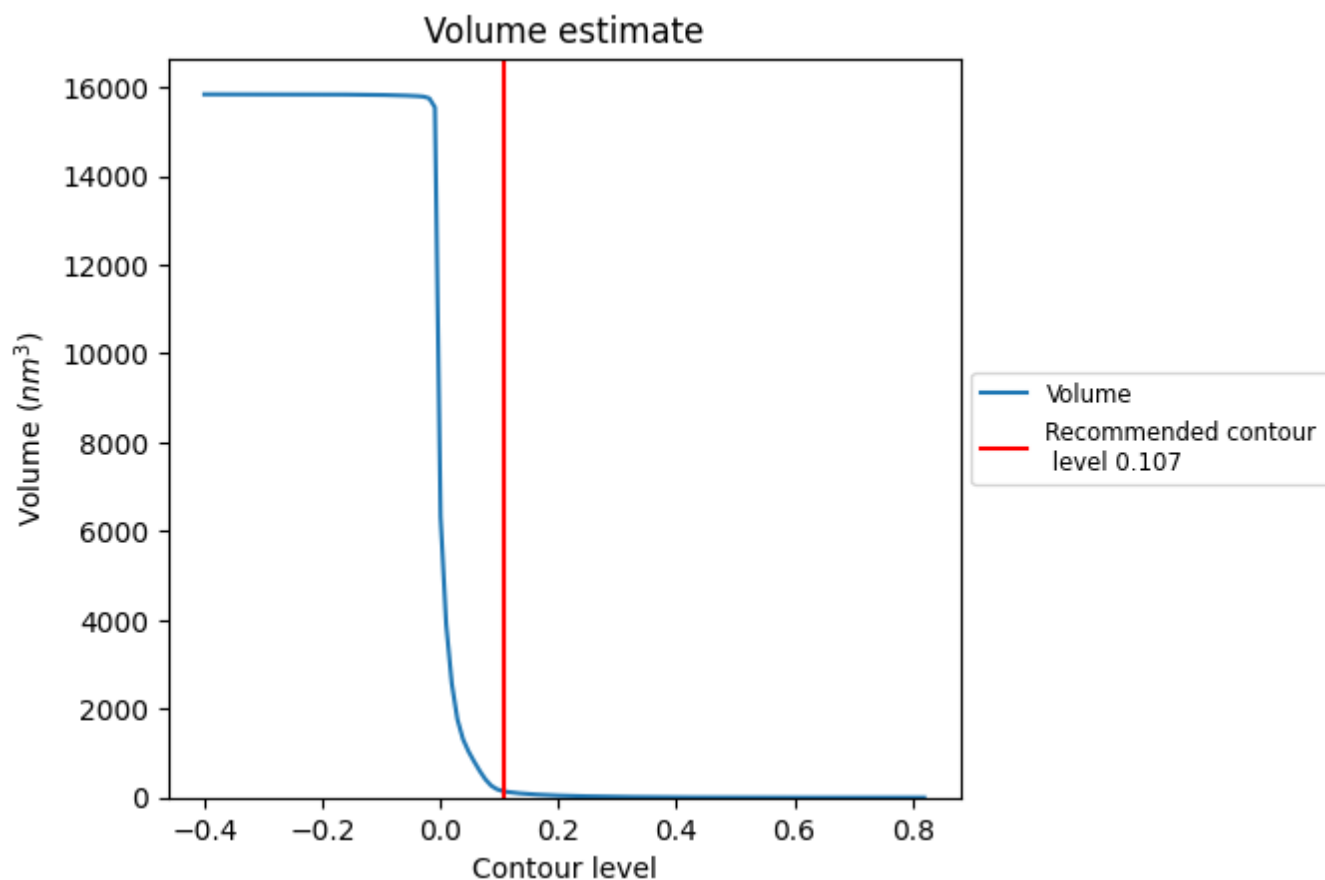
This section contains the results of statistical analysis of the map.

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.

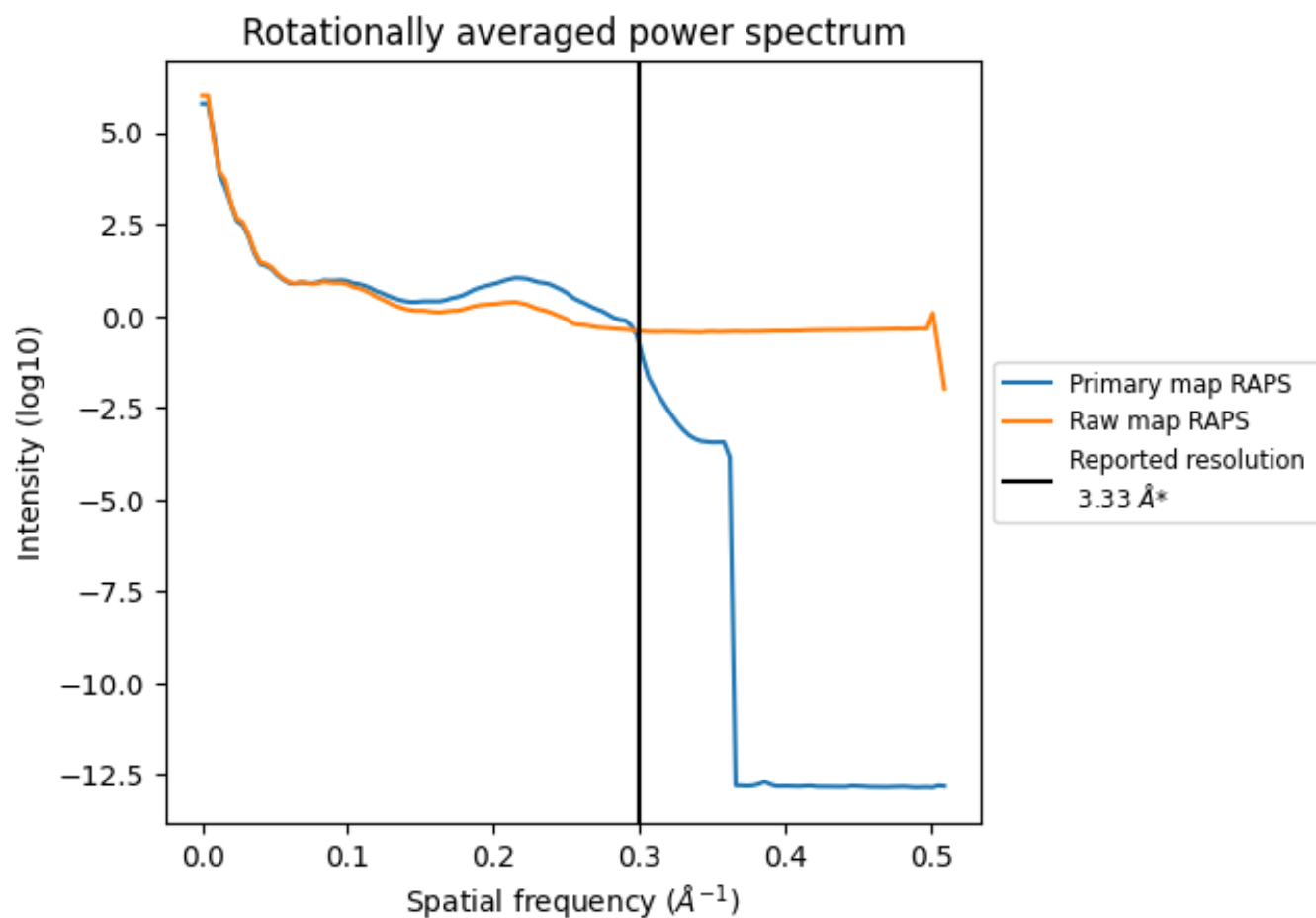
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 143 nm³; this corresponds to an approximate mass of 130 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 ⓘ

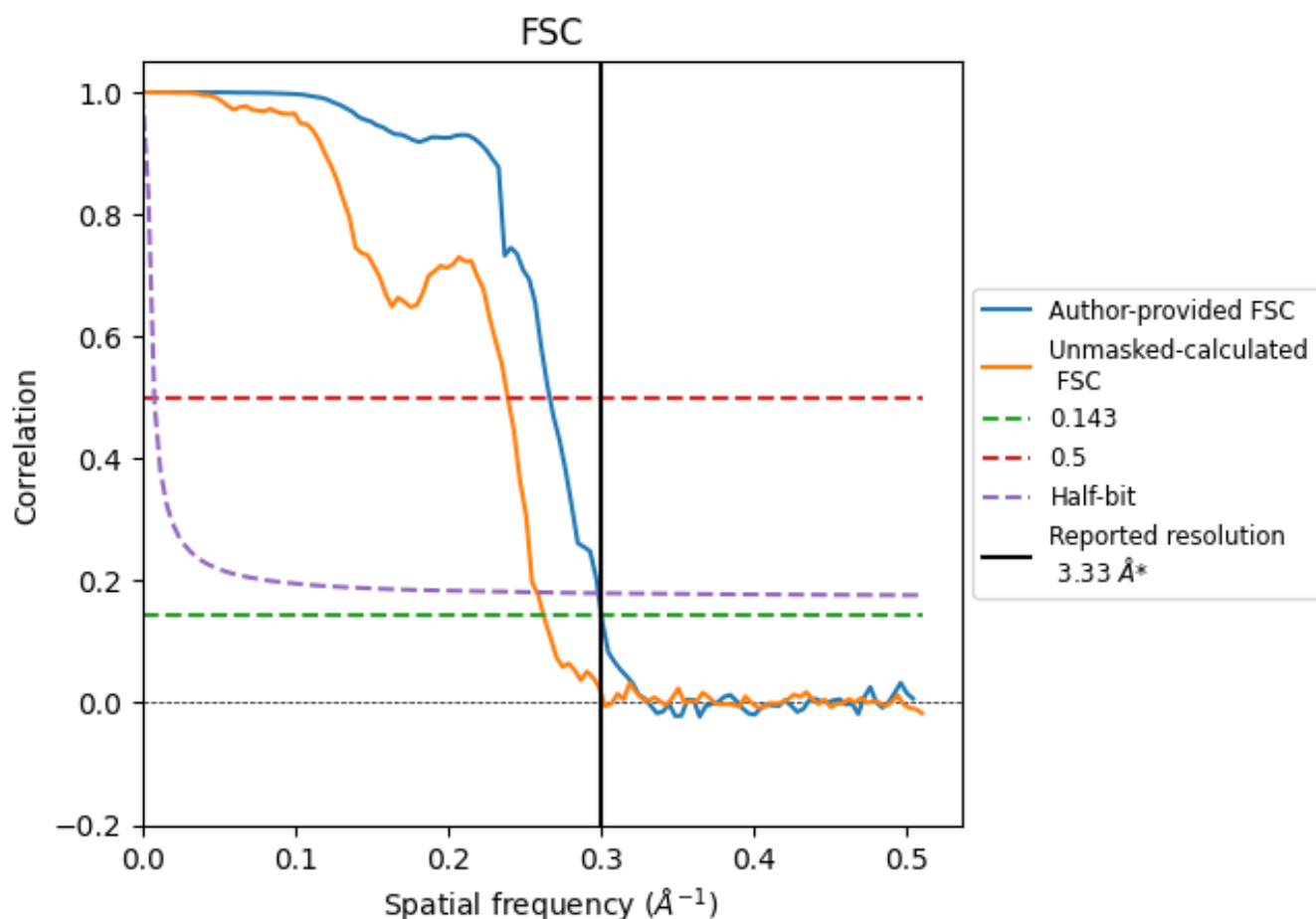


*Reported resolution corresponds to spatial frequency of 0.300 Å⁻¹

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 0.300 \AA^{-1}

8.2 Resolution estimates [i](#)

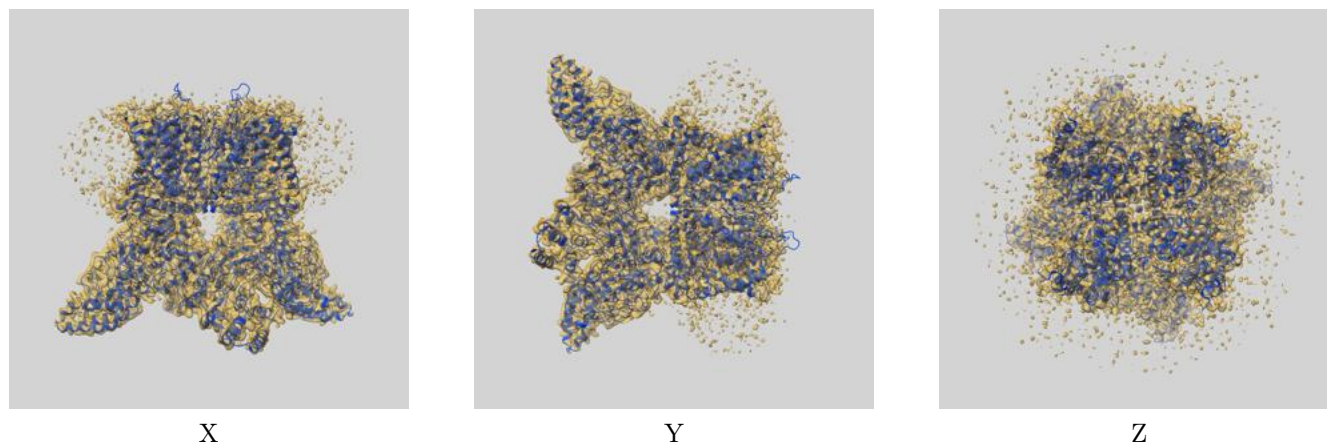
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.33	-	-
Author-provided FSC curve	3.33	3.76	3.35
Unmasked-calculated*	3.81	4.19	3.87

*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 3.81 differs from the reported value 3.33 by more than 10 %

9 Map-model fit [i](#)

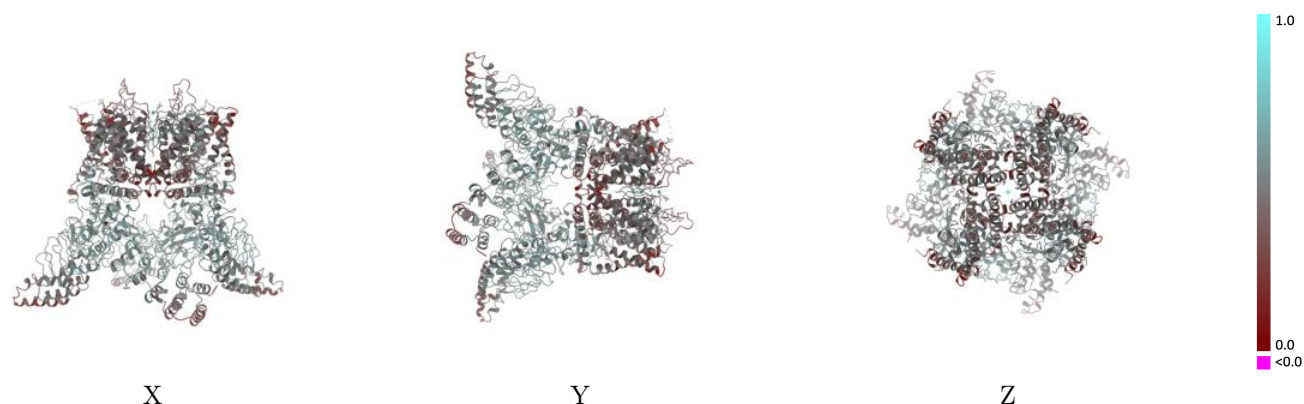
This section contains information regarding the fit between EMDB map EMD-77142 and PDB model 13LR. Per-residue inclusion information can be found in section [3](#) on page [8](#).

9.1 Map-model overlay [i](#)



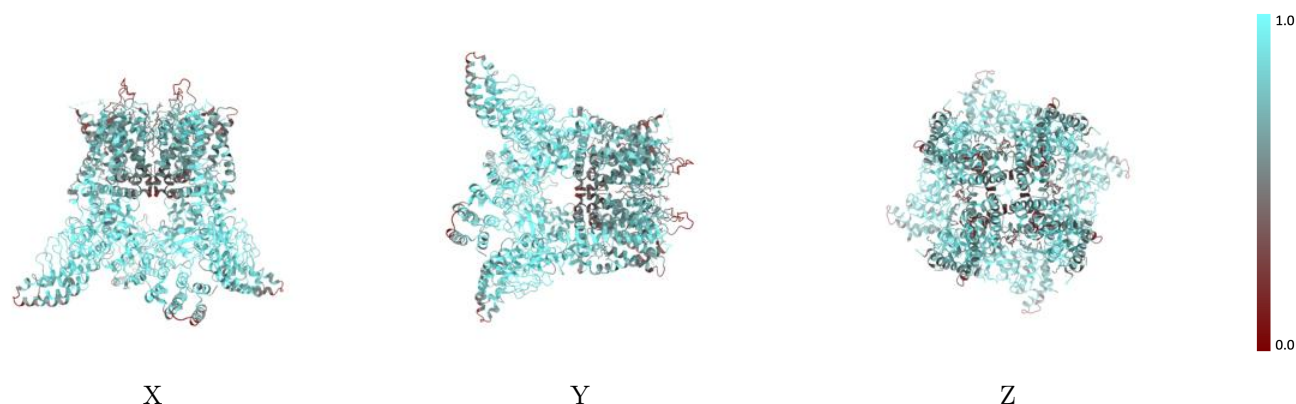
The images above show the 3D surface view of the map at the recommended contour level 0.107 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



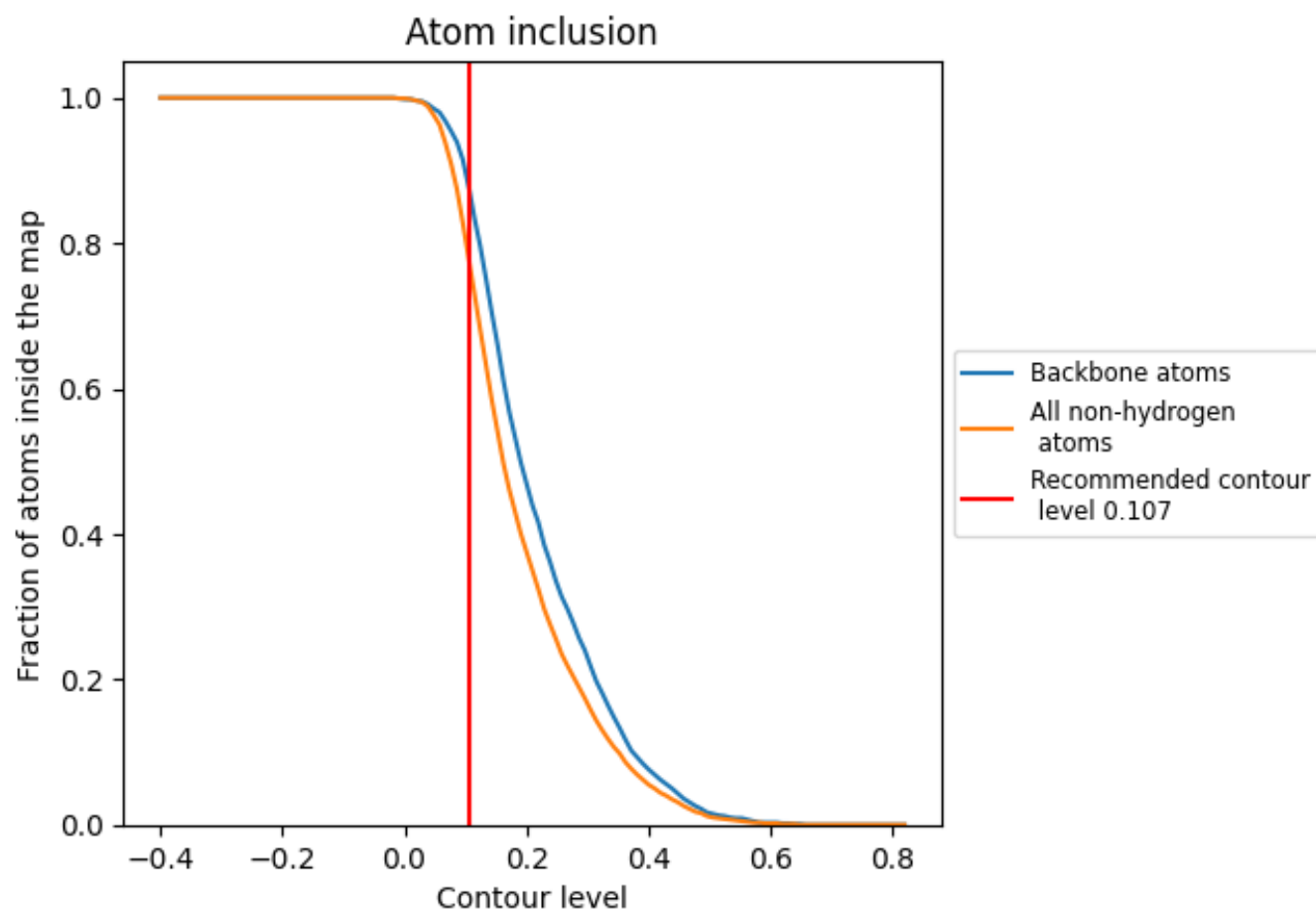
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.107).

9.4 Atom inclusion [i](#)



At the recommended contour level, 87% of all backbone atoms, 77% 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 (0.107) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div></div> 0.7710	<div></div> 0.4690
A	<div></div> 0.7670	<div></div> 0.4690
B	<div></div> 0.7760	<div></div> 0.4700
C	<div></div> 0.7700	<div></div> 0.4700
D	<div></div> 0.7710	<div></div> 0.4690

