



Full wwPDB EM Validation Report ⓘ

Mar 8, 2026 – 04:42 PM UTC

PDB ID : 6PWN / pdb_00006pwn
EMDB ID : EMD-20508
Title : MscS Nanodisc with N-terminal His-Tag
Authors : Reddy, B.G.; Perozo, E.
Deposited on : 2019-07-23
Resolution : 3.10 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:

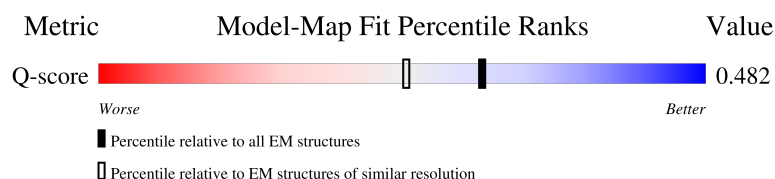
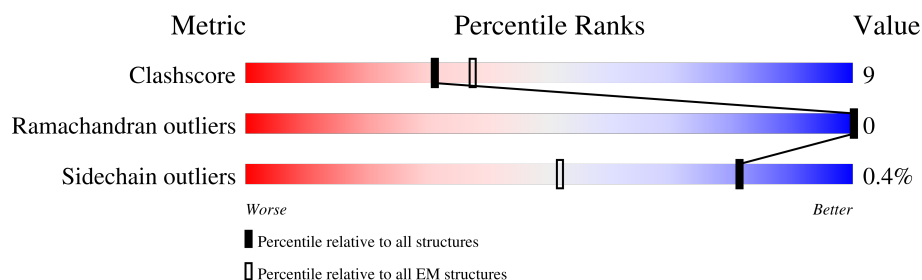
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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

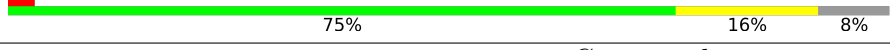
The reported resolution of this entry is 3.10 Å.

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	14724 (2.60 - 3.60)

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	306	
1	B	306	
1	C	306	
1	D	306	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	E	306	75%16%8%
1	F	306	75%16%8%
1	G	306	74%17%8%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15379 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 Small-conductance mechanosensitive channel.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	280	Total	C	N	O	S	0	0
			2129	1364	370	388	7		
1	B	280	Total	C	N	O	S	0	0
			2129	1364	370	388	7		
1	C	280	Total	C	N	O	S	0	0
			2129	1364	370	388	7		
1	D	280	Total	C	N	O	S	0	0
			2129	1364	370	388	7		
1	E	280	Total	C	N	O	S	0	0
			2129	1364	370	388	7		
1	F	280	Total	C	N	O	S	0	0
			2129	1364	370	388	7		
1	G	280	Total	C	N	O	S	0	0
			2129	1364	370	388	7		

There are 140 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P0C0S1
A	-18	GLY	-	expression tag	UNP P0C0S1
A	-17	SER	-	expression tag	UNP P0C0S1
A	-16	SER	-	expression tag	UNP P0C0S1
A	-15	HIS	-	expression tag	UNP P0C0S1
A	-14	HIS	-	expression tag	UNP P0C0S1
A	-13	HIS	-	expression tag	UNP P0C0S1
A	-12	HIS	-	expression tag	UNP P0C0S1
A	-11	HIS	-	expression tag	UNP P0C0S1
A	-10	HIS	-	expression tag	UNP P0C0S1
A	-9	SER	-	expression tag	UNP P0C0S1
A	-8	SER	-	expression tag	UNP P0C0S1
A	-7	GLY	-	expression tag	UNP P0C0S1
A	-6	LEU	-	expression tag	UNP P0C0S1
A	-5	VAL	-	expression tag	UNP P0C0S1
A	-4	PRO	-	expression tag	UNP P0C0S1

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	ARG	-	expression tag	UNP P0C0S1
A	-2	GLY	-	expression tag	UNP P0C0S1
A	-1	SER	-	expression tag	UNP P0C0S1
A	0	HIS	-	expression tag	UNP P0C0S1
B	-19	MET	-	initiating methionine	UNP P0C0S1
B	-18	GLY	-	expression tag	UNP P0C0S1
B	-17	SER	-	expression tag	UNP P0C0S1
B	-16	SER	-	expression tag	UNP P0C0S1
B	-15	HIS	-	expression tag	UNP P0C0S1
B	-14	HIS	-	expression tag	UNP P0C0S1
B	-13	HIS	-	expression tag	UNP P0C0S1
B	-12	HIS	-	expression tag	UNP P0C0S1
B	-11	HIS	-	expression tag	UNP P0C0S1
B	-10	HIS	-	expression tag	UNP P0C0S1
B	-9	SER	-	expression tag	UNP P0C0S1
B	-8	SER	-	expression tag	UNP P0C0S1
B	-7	GLY	-	expression tag	UNP P0C0S1
B	-6	LEU	-	expression tag	UNP P0C0S1
B	-5	VAL	-	expression tag	UNP P0C0S1
B	-4	PRO	-	expression tag	UNP P0C0S1
B	-3	ARG	-	expression tag	UNP P0C0S1
B	-2	GLY	-	expression tag	UNP P0C0S1
B	-1	SER	-	expression tag	UNP P0C0S1
B	0	HIS	-	expression tag	UNP P0C0S1
C	-19	MET	-	initiating methionine	UNP P0C0S1
C	-18	GLY	-	expression tag	UNP P0C0S1
C	-17	SER	-	expression tag	UNP P0C0S1
C	-16	SER	-	expression tag	UNP P0C0S1
C	-15	HIS	-	expression tag	UNP P0C0S1
C	-14	HIS	-	expression tag	UNP P0C0S1
C	-13	HIS	-	expression tag	UNP P0C0S1
C	-12	HIS	-	expression tag	UNP P0C0S1
C	-11	HIS	-	expression tag	UNP P0C0S1
C	-10	HIS	-	expression tag	UNP P0C0S1
C	-9	SER	-	expression tag	UNP P0C0S1
C	-8	SER	-	expression tag	UNP P0C0S1
C	-7	GLY	-	expression tag	UNP P0C0S1
C	-6	LEU	-	expression tag	UNP P0C0S1
C	-5	VAL	-	expression tag	UNP P0C0S1
C	-4	PRO	-	expression tag	UNP P0C0S1
C	-3	ARG	-	expression tag	UNP P0C0S1
C	-2	GLY	-	expression tag	UNP P0C0S1

Continued on next page...

Continued from previous page...

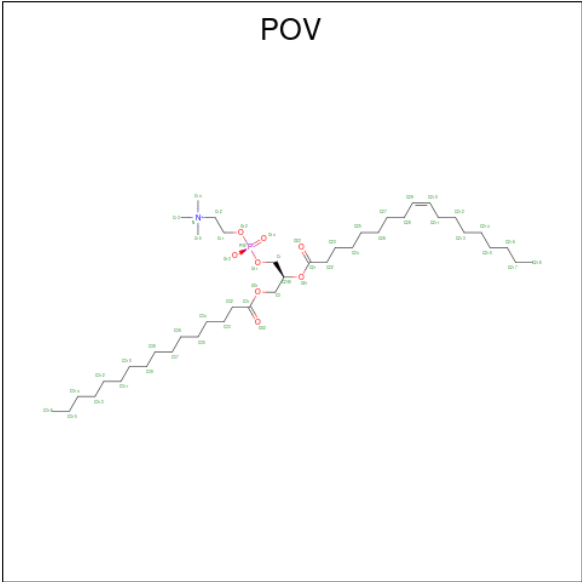
Chain	Residue	Modelled	Actual	Comment	Reference
C	-1	SER	-	expression tag	UNP P0C0S1
C	0	HIS	-	expression tag	UNP P0C0S1
D	-19	MET	-	initiating methionine	UNP P0C0S1
D	-18	GLY	-	expression tag	UNP P0C0S1
D	-17	SER	-	expression tag	UNP P0C0S1
D	-16	SER	-	expression tag	UNP P0C0S1
D	-15	HIS	-	expression tag	UNP P0C0S1
D	-14	HIS	-	expression tag	UNP P0C0S1
D	-13	HIS	-	expression tag	UNP P0C0S1
D	-12	HIS	-	expression tag	UNP P0C0S1
D	-11	HIS	-	expression tag	UNP P0C0S1
D	-10	HIS	-	expression tag	UNP P0C0S1
D	-9	SER	-	expression tag	UNP P0C0S1
D	-8	SER	-	expression tag	UNP P0C0S1
D	-7	GLY	-	expression tag	UNP P0C0S1
D	-6	LEU	-	expression tag	UNP P0C0S1
D	-5	VAL	-	expression tag	UNP P0C0S1
D	-4	PRO	-	expression tag	UNP P0C0S1
D	-3	ARG	-	expression tag	UNP P0C0S1
D	-2	GLY	-	expression tag	UNP P0C0S1
D	-1	SER	-	expression tag	UNP P0C0S1
D	0	HIS	-	expression tag	UNP P0C0S1
E	-19	MET	-	initiating methionine	UNP P0C0S1
E	-18	GLY	-	expression tag	UNP P0C0S1
E	-17	SER	-	expression tag	UNP P0C0S1
E	-16	SER	-	expression tag	UNP P0C0S1
E	-15	HIS	-	expression tag	UNP P0C0S1
E	-14	HIS	-	expression tag	UNP P0C0S1
E	-13	HIS	-	expression tag	UNP P0C0S1
E	-12	HIS	-	expression tag	UNP P0C0S1
E	-11	HIS	-	expression tag	UNP P0C0S1
E	-10	HIS	-	expression tag	UNP P0C0S1
E	-9	SER	-	expression tag	UNP P0C0S1
E	-8	SER	-	expression tag	UNP P0C0S1
E	-7	GLY	-	expression tag	UNP P0C0S1
E	-6	LEU	-	expression tag	UNP P0C0S1
E	-5	VAL	-	expression tag	UNP P0C0S1
E	-4	PRO	-	expression tag	UNP P0C0S1
E	-3	ARG	-	expression tag	UNP P0C0S1
E	-2	GLY	-	expression tag	UNP P0C0S1
E	-1	SER	-	expression tag	UNP P0C0S1
E	0	HIS	-	expression tag	UNP P0C0S1

Continued on next page...

Continued from previous page...

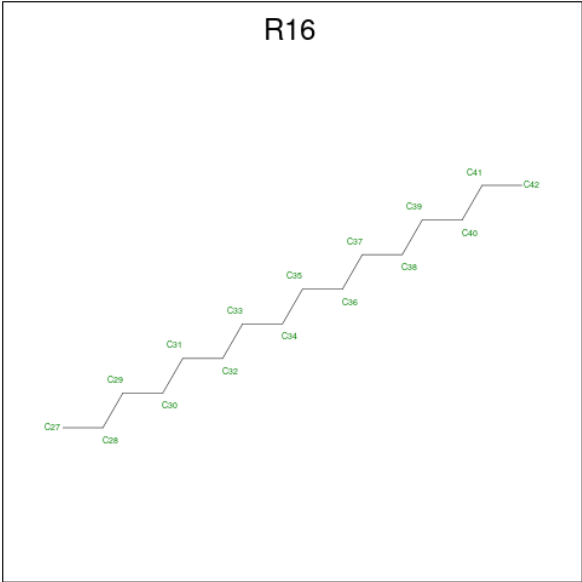
Chain	Residue	Modelled	Actual	Comment	Reference
F	-19	MET	-	initiating methionine	UNP P0C0S1
F	-18	GLY	-	expression tag	UNP P0C0S1
F	-17	SER	-	expression tag	UNP P0C0S1
F	-16	SER	-	expression tag	UNP P0C0S1
F	-15	HIS	-	expression tag	UNP P0C0S1
F	-14	HIS	-	expression tag	UNP P0C0S1
F	-13	HIS	-	expression tag	UNP P0C0S1
F	-12	HIS	-	expression tag	UNP P0C0S1
F	-11	HIS	-	expression tag	UNP P0C0S1
F	-10	HIS	-	expression tag	UNP P0C0S1
F	-9	SER	-	expression tag	UNP P0C0S1
F	-8	SER	-	expression tag	UNP P0C0S1
F	-7	GLY	-	expression tag	UNP P0C0S1
F	-6	LEU	-	expression tag	UNP P0C0S1
F	-5	VAL	-	expression tag	UNP P0C0S1
F	-4	PRO	-	expression tag	UNP P0C0S1
F	-3	ARG	-	expression tag	UNP P0C0S1
F	-2	GLY	-	expression tag	UNP P0C0S1
F	-1	SER	-	expression tag	UNP P0C0S1
F	0	HIS	-	expression tag	UNP P0C0S1
G	-19	MET	-	initiating methionine	UNP P0C0S1
G	-18	GLY	-	expression tag	UNP P0C0S1
G	-17	SER	-	expression tag	UNP P0C0S1
G	-16	SER	-	expression tag	UNP P0C0S1
G	-15	HIS	-	expression tag	UNP P0C0S1
G	-14	HIS	-	expression tag	UNP P0C0S1
G	-13	HIS	-	expression tag	UNP P0C0S1
G	-12	HIS	-	expression tag	UNP P0C0S1
G	-11	HIS	-	expression tag	UNP P0C0S1
G	-10	HIS	-	expression tag	UNP P0C0S1
G	-9	SER	-	expression tag	UNP P0C0S1
G	-8	SER	-	expression tag	UNP P0C0S1
G	-7	GLY	-	expression tag	UNP P0C0S1
G	-6	LEU	-	expression tag	UNP P0C0S1
G	-5	VAL	-	expression tag	UNP P0C0S1
G	-4	PRO	-	expression tag	UNP P0C0S1
G	-3	ARG	-	expression tag	UNP P0C0S1
G	-2	GLY	-	expression tag	UNP P0C0S1
G	-1	SER	-	expression tag	UNP P0C0S1
G	0	HIS	-	expression tag	UNP P0C0S1

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



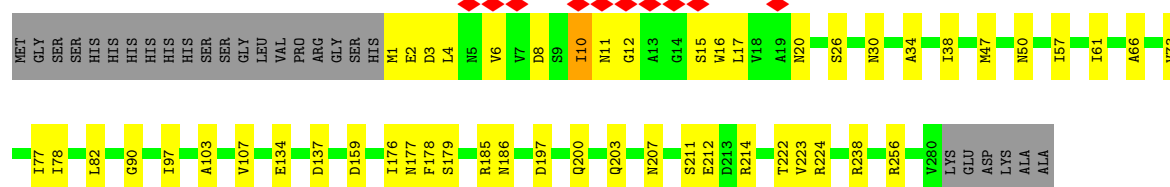
Mol	Chain	Residues	Atoms					AltConf
2	A	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	B	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	C	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	D	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	E	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	F	1	Total	C	N	O	P	0
			52	42	1	8	1	
2	G	1	Total	C	N	O	P	0
			52	42	1	8	1	

- Molecule 3 is HEXADECANE (CCD ID: R16) (formula: C₁₆H₃₄).



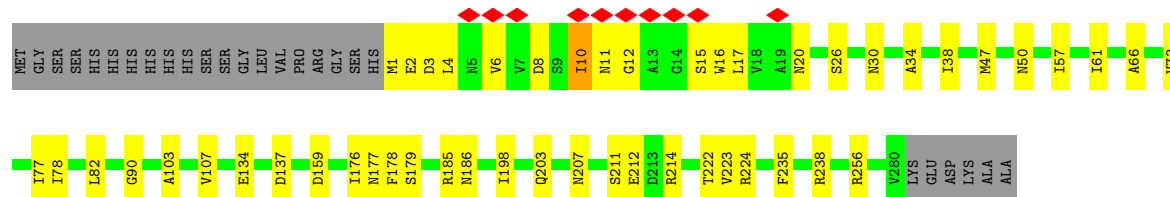
Mol	Chain	Residues	Atoms		AltConf
3	A	1	Total	C	0
			16	16	
3	B	1	Total	C	0
			16	16	
3	C	1	Total	C	0
			16	16	
3	C	1	Total	C	0
			16	16	
3	E	1	Total	C	0
			16	16	
3	F	1	Total	C	0
			16	16	
3	F	1	Total	C	0
			16	16	

Chain D: 



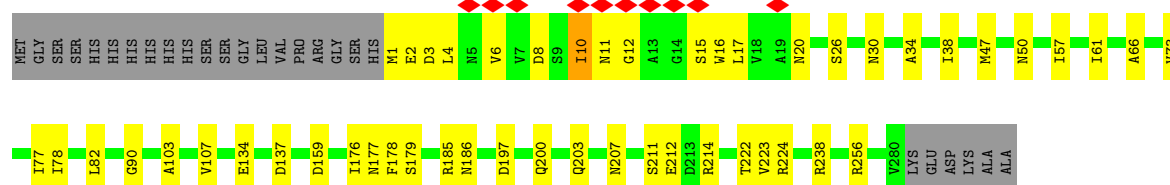
- Molecule 1: Small-conductance mechanosensitive channel

Chain E: 



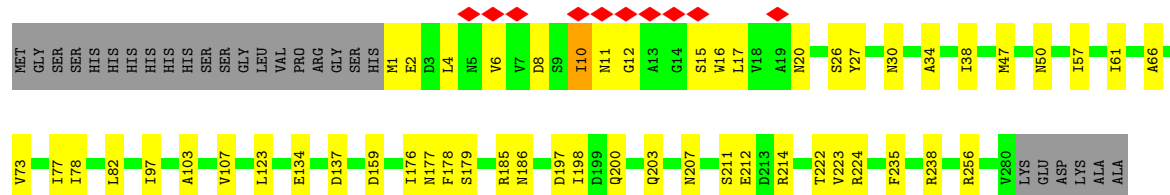
- Molecule 1: Small-conductance mechanosensitive channel

Chain F: 



- Molecule 1: Small-conductance mechanosensitive channel

Chain G: 



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C7	Depositor
Number of particles used	372870	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	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	0.073	Depositor
Minimum map value	-0.027	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.0092	Depositor
Map size ($\text{\AA}$)	272.384, 272.384, 272.384	wwPDB
Map dimensions	256, 256, 256	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.064, 1.064, 1.064	Depositor

5 Model quality

5.1 Standard geometry

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

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	0.26	0/2156	0.37	0/2926
1	B	0.26	0/2156	0.36	0/2926
1	C	0.26	0/2156	0.36	0/2926
1	D	0.26	0/2156	0.36	0/2926
1	E	0.26	0/2156	0.36	0/2926
1	F	0.26	0/2156	0.36	0/2926
1	G	0.26	0/2156	0.36	0/2926
All	All	0.26	0/15092	0.36	0/20482

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	2129	0	2223	44	0
1	B	2129	0	2223	45	0
1	C	2129	0	2223	40	0
1	D	2129	0	2223	40	0
1	E	2129	0	2223	38	0
1	F	2129	0	2223	39	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	2129	0	2223	40	0
2	A	52	0	82	1	0
2	B	52	0	82	2	0
2	C	52	0	82	1	0
2	D	52	0	82	1	0
2	E	52	0	82	2	0
2	F	52	0	82	2	0
2	G	52	0	82	3	0
3	A	16	0	34	6	0
3	B	16	0	34	5	0
3	C	32	0	68	4	0
3	E	16	0	34	1	0
3	F	32	0	68	4	0
All	All	15379	0	16373	278	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 9.

All (278) 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:78:ILE:O	1:A:82:LEU:HD13	1.69	0.93
3:C:302:R16:H331	1:D:97:ILE:HD11	1.57	0.87
3:A:302:R16:H312	3:B:302:R16:H291	1.60	0.81
1:F:16:TRP:O	1:F:20:ASN:HB2	1.82	0.80
1:G:16:TRP:O	1:G:20:ASN:HB2	1.82	0.80
1:C:16:TRP:O	1:C:20:ASN:HB2	1.82	0.79
1:D:16:TRP:O	1:D:20:ASN:HB2	1.82	0.79
1:A:16:TRP:O	1:A:20:ASN:HB2	1.82	0.79
3:A:302:R16:H291	3:F:302:R16:H311	1.63	0.79
1:E:16:TRP:O	1:E:20:ASN:HB2	1.82	0.79
1:B:16:TRP:O	1:B:20:ASN:HB2	1.82	0.78
1:F:1:MET:HE1	1:F:17:LEU:H	1.49	0.78
1:C:1:MET:HE1	1:C:17:LEU:H	1.49	0.78
1:G:1:MET:HE1	1:G:17:LEU:H	1.49	0.78
1:A:99:VAL:CG2	1:B:100:LEU:HD21	2.14	0.77
1:A:10:ILE:HG22	1:A:12:GLY:H	1.49	0.77
1:A:1:MET:HE1	1:A:17:LEU:H	1.49	0.77
1:G:10:ILE:HG22	1:G:12:GLY:H	1.50	0.77
1:B:10:ILE:HG22	1:B:12:GLY:H	1.50	0.77
1:D:1:MET:HE1	1:D:17:LEU:H	1.49	0.76

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:1:MET:HE1	1:B:17:LEU:H	1.49	0.76
1:E:1:MET:HE1	1:E:17:LEU:H	1.49	0.76
1:E:10:ILE:HG22	1:E:12:GLY:H	1.49	0.76
1:F:10:ILE:HG22	1:F:12:GLY:H	1.49	0.76
1:C:10:ILE:HG22	1:C:12:GLY:H	1.49	0.76
1:E:78:ILE:O	1:E:82:LEU:HD13	1.86	0.75
1:C:78:ILE:O	1:C:82:LEU:HD13	1.86	0.75
1:D:78:ILE:O	1:D:82:LEU:HD13	1.87	0.75
1:B:78:ILE:O	1:B:82:LEU:HD13	1.86	0.75
1:B:57:ILE:HD12	1:B:61:ILE:HD11	1.69	0.74
1:F:78:ILE:O	1:F:82:LEU:HD13	1.86	0.74
1:F:57:ILE:HD12	1:F:61:ILE:HD11	1.70	0.74
1:C:57:ILE:HD12	1:C:61:ILE:HD11	1.70	0.74
1:A:57:ILE:HD12	1:A:61:ILE:HD11	1.69	0.74
1:G:57:ILE:HD12	1:G:61:ILE:HD11	1.70	0.74
1:D:10:ILE:HG22	1:D:12:GLY:H	1.50	0.74
1:E:57:ILE:HD12	1:E:61:ILE:HD11	1.70	0.74
1:D:57:ILE:HD12	1:D:61:ILE:HD11	1.69	0.74
1:G:78:ILE:O	1:G:82:LEU:HD13	1.87	0.74
1:G:134:GLU:OE2	1:G:177:ASN:ND2	2.28	0.67
1:D:134:GLU:OE2	1:D:177:ASN:ND2	2.28	0.67
1:A:78:ILE:O	1:A:82:LEU:CD1	2.44	0.65
1:A:134:GLU:OE2	1:A:177:ASN:ND2	2.28	0.65
1:F:159:ASP:OD1	1:G:185:ARG:NH2	2.30	0.65
1:C:159:ASP:OD1	1:D:185:ARG:NH2	2.30	0.64
1:G:1:MET:HE3	1:G:4:LEU:HB2	1.80	0.64
1:E:134:GLU:OE2	1:E:177:ASN:ND2	2.28	0.64
1:F:1:MET:HE3	1:F:4:LEU:HB2	1.80	0.64
1:A:1:MET:HE3	1:A:4:LEU:HB2	1.80	0.64
1:C:16:TRP:O	1:C:20:ASN:CB	2.46	0.64
1:E:1:MET:HE3	1:E:4:LEU:HB2	1.80	0.64
1:D:1:MET:HE3	1:D:4:LEU:HB2	1.80	0.64
1:B:134:GLU:OE2	1:B:177:ASN:ND2	2.28	0.64
1:D:159:ASP:OD1	1:E:185:ARG:NH2	2.31	0.64
3:A:302:R16:H281	3:F:302:R16:H292	1.80	0.63
1:F:16:TRP:O	1:F:20:ASN:CB	2.46	0.63
1:G:16:TRP:O	1:G:20:ASN:CB	2.46	0.63
1:A:16:TRP:O	1:A:20:ASN:CB	2.46	0.63
1:C:1:MET:HE3	1:C:4:LEU:HB2	1.80	0.63
1:E:16:TRP:O	1:E:20:ASN:CB	2.46	0.63
1:B:1:MET:HE3	1:B:4:LEU:HB2	1.80	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:159:ASP:OD1	1:C:185:ARG:NH2	2.30	0.63
1:A:159:ASP:OD1	1:B:185:ARG:NH2	2.31	0.63
1:F:134:GLU:OE2	1:F:177:ASN:ND2	2.28	0.63
3:A:302:R16:H301	3:B:302:R16:H272	1.81	0.63
1:D:16:TRP:O	1:D:20:ASN:CB	2.46	0.63
1:B:16:TRP:O	1:B:20:ASN:CB	2.46	0.63
1:A:185:ARG:NH2	1:G:159:ASP:OD1	2.30	0.62
1:E:159:ASP:OD1	1:F:185:ARG:NH2	2.30	0.62
1:C:134:GLU:OE2	1:C:177:ASN:ND2	2.28	0.60
3:F:302:R16:H341	1:G:97:ILE:HD11	1.83	0.59
1:E:222:THR:HG23	1:E:224:ARG:HG2	1.84	0.58
1:A:222:THR:HG23	1:A:224:ARG:HG2	1.85	0.58
1:C:90:GLY:HA2	2:C:301:POV:H12	1.85	0.58
1:D:222:THR:HG23	1:D:224:ARG:HG2	1.84	0.58
1:F:90:GLY:HA2	2:F:301:POV:H12	1.85	0.58
1:B:222:THR:HG23	1:B:224:ARG:HG2	1.84	0.58
1:D:90:GLY:HA2	2:D:301:POV:H12	1.84	0.58
1:B:61:ILE:HD12	1:B:66:ALA:HB2	1.86	0.58
1:C:222:THR:HG23	1:C:224:ARG:HG2	1.84	0.58
1:F:222:THR:HG23	1:F:224:ARG:HG2	1.85	0.58
1:G:222:THR:HG23	1:G:224:ARG:HG2	1.84	0.58
1:C:61:ILE:HD12	1:C:66:ALA:HB2	1.86	0.58
1:A:11:ASN:HB3	1:B:17:LEU:HD22	1.87	0.57
1:A:61:ILE:HD12	1:A:66:ALA:HB2	1.86	0.57
1:F:61:ILE:HD12	1:F:66:ALA:HB2	1.86	0.57
1:D:61:ILE:HD12	1:D:66:ALA:HB2	1.86	0.57
1:E:61:ILE:HD12	1:E:66:ALA:HB2	1.86	0.57
1:A:99:VAL:HG23	1:B:100:LEU:HD21	1.87	0.57
1:G:61:ILE:HD12	1:G:66:ALA:HB2	1.86	0.56
1:E:11:ASN:HB3	1:F:17:LEU:HD22	1.87	0.55
1:D:11:ASN:HB3	1:E:17:LEU:HD22	1.88	0.54
1:F:11:ASN:HB3	1:G:17:LEU:HD22	1.89	0.54
1:A:90:GLY:HA2	2:A:301:POV:H12	1.89	0.54
1:F:212:GLU:OE2	1:F:214:ARG:NH2	2.38	0.53
1:B:97:ILE:HD11	3:B:302:R16:H331	1.91	0.53
1:B:185:ARG:HD2	1:B:238:ARG:HB3	1.91	0.53
1:G:27:TYR:CE1	2:G:301:POV:H3A	2.44	0.53
1:A:185:ARG:HD2	1:A:238:ARG:HB3	1.91	0.53
1:C:185:ARG:HD2	1:C:238:ARG:HB3	1.91	0.53
1:G:185:ARG:HD2	1:G:238:ARG:HB3	1.91	0.53
1:A:17:LEU:HD22	1:G:11:ASN:HB3	1.91	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:212:GLU:OE2	1:A:214:ARG:NH2	2.38	0.53
1:C:212:GLU:OE2	1:C:214:ARG:NH2	2.38	0.53
1:F:185:ARG:HD2	1:F:238:ARG:HB3	1.91	0.52
1:A:137:ASP:HB3	1:A:176:ILE:HB	1.92	0.52
1:D:185:ARG:HD2	1:D:238:ARG:HB3	1.91	0.52
1:G:137:ASP:HB3	1:G:176:ILE:HB	1.92	0.52
1:B:11:ASN:HB3	1:C:17:LEU:HD22	1.90	0.52
1:B:137:ASP:HB3	1:B:176:ILE:HB	1.92	0.52
1:C:97:ILE:HD11	3:C:303:R16:H331	1.91	0.52
1:F:137:ASP:HB3	1:F:176:ILE:HB	1.92	0.52
1:E:90:GLY:HA2	2:E:301:POV:H12	1.90	0.52
1:E:1:MET:HB3	1:E:4:LEU:HB3	1.92	0.52
1:A:1:MET:HB3	1:A:4:LEU:HB3	1.93	0.51
1:C:26:SER:O	1:C:30:ASN:ND2	2.43	0.51
1:D:197:ASP:OD2	1:D:200:GLN:N	2.33	0.51
1:G:1:MET:HB3	1:G:4:LEU:HB3	1.93	0.51
1:D:1:MET:HB3	1:D:4:LEU:HB3	1.92	0.51
1:E:137:ASP:HB3	1:E:176:ILE:HB	1.92	0.51
1:C:137:ASP:HB3	1:C:176:ILE:HB	1.92	0.51
1:F:1:MET:HB3	1:F:4:LEU:HB3	1.93	0.51
1:B:1:MET:HB3	1:B:4:LEU:HB3	1.93	0.51
1:D:26:SER:O	1:D:30:ASN:ND2	2.43	0.51
1:E:26:SER:O	1:E:30:ASN:ND2	2.43	0.51
1:C:1:MET:HB3	1:C:4:LEU:HB3	1.92	0.51
1:A:26:SER:O	1:A:30:ASN:ND2	2.43	0.51
1:F:26:SER:O	1:F:30:ASN:ND2	2.43	0.51
1:G:26:SER:O	1:G:30:ASN:ND2	2.44	0.51
1:D:137:ASP:HB3	1:D:176:ILE:HB	1.92	0.51
1:C:11:ASN:HB3	1:D:17:LEU:HD22	1.92	0.51
1:B:90:GLY:HA2	2:B:301:POV:H12	1.92	0.50
1:G:212:GLU:OE2	1:G:214:ARG:NH2	2.38	0.50
1:E:185:ARG:HD2	1:E:238:ARG:HB3	1.91	0.50
1:B:26:SER:O	1:B:30:ASN:ND2	2.44	0.50
1:G:203:GLN:HE21	1:G:207:ASN:HD21	1.60	0.50
1:E:212:GLU:OE2	1:E:214:ARG:NH2	2.38	0.49
1:B:203:GLN:HE21	1:B:207:ASN:HD21	1.60	0.49
1:E:34:ALA:O	1:E:38:ILE:HG13	2.12	0.49
1:D:34:ALA:O	1:D:38:ILE:HG13	2.12	0.49
1:G:34:ALA:O	1:G:38:ILE:HG13	2.12	0.49
1:A:203:GLN:HE21	1:A:207:ASN:HD21	1.60	0.49
1:A:34:ALA:O	1:A:38:ILE:HG13	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:134:GLU:HG3	1:F:179:SER:OG	2.13	0.49
1:E:134:GLU:HG3	1:E:179:SER:OG	2.13	0.49
1:E:203:GLN:HE21	1:E:207:ASN:HD21	1.60	0.48
1:C:34:ALA:O	1:C:38:ILE:HG13	2.13	0.48
1:C:197:ASP:OD2	1:C:200:GLN:N	2.33	0.48
1:C:203:GLN:HE21	1:C:207:ASN:HD21	1.60	0.48
1:C:134:GLU:HG3	1:C:179:SER:OG	2.13	0.48
1:G:134:GLU:HG3	1:G:179:SER:OG	2.13	0.48
1:B:34:ALA:O	1:B:38:ILE:HG13	2.12	0.48
1:B:212:GLU:OE2	1:B:214:ARG:NH2	2.38	0.48
1:D:134:GLU:HG3	1:D:179:SER:OG	2.13	0.48
1:F:203:GLN:HE21	1:F:207:ASN:HD21	1.60	0.48
1:F:34:ALA:O	1:F:38:ILE:HG13	2.13	0.48
1:B:134:GLU:HG3	1:B:179:SER:OG	2.13	0.48
1:A:197:ASP:OD2	1:A:200:GLN:N	2.33	0.48
1:A:134:GLU:HG3	1:A:179:SER:OG	2.13	0.47
1:G:197:ASP:OD2	1:G:200:GLN:N	2.33	0.47
1:B:197:ASP:OD2	1:B:200:GLN:N	2.33	0.47
1:D:203:GLN:HE21	1:D:207:ASN:HD21	1.60	0.47
1:D:212:GLU:OE2	1:D:214:ARG:NH2	2.38	0.46
1:D:78:ILE:O	1:D:82:LEU:CD1	2.61	0.46
1:E:103:ALA:O	1:E:107:VAL:HG23	2.16	0.46
1:G:103:ALA:O	1:G:107:VAL:HG23	2.16	0.46
1:B:103:ALA:O	1:B:107:VAL:HG23	2.16	0.46
1:E:177:ASN:OD1	1:E:178:PHE:N	2.49	0.46
1:A:99:VAL:CG2	1:B:100:LEU:CD2	2.89	0.45
1:B:2:GLU:O	1:B:6:VAL:HG23	2.17	0.45
1:G:2:GLU:O	1:G:6:VAL:HG23	2.17	0.45
1:G:47:MET:HA	1:G:50:ASN:HD21	1.81	0.45
1:C:103:ALA:O	1:C:107:VAL:HG23	2.16	0.45
1:D:103:ALA:O	1:D:107:VAL:HG23	2.16	0.45
1:F:177:ASN:OD1	1:F:178:PHE:N	2.49	0.45
1:A:103:ALA:O	1:A:107:VAL:HG23	2.16	0.45
1:A:177:ASN:OD1	1:A:178:PHE:N	2.49	0.45
3:A:302:R16:C30	3:B:302:R16:H272	2.46	0.45
3:C:302:R16:H301	3:E:302:R16:H272	1.98	0.45
1:F:197:ASP:OD2	1:F:200:GLN:N	2.33	0.45
1:A:2:GLU:O	1:A:6:VAL:HG23	2.17	0.45
1:E:2:GLU:O	1:E:6:VAL:HG23	2.17	0.45
1:G:78:ILE:O	1:G:82:LEU:CD1	2.61	0.45
1:G:177:ASN:OD1	1:G:178:PHE:N	2.49	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:78:ILE:O	1:F:82:LEU:CD1	2.61	0.45
1:B:47:MET:HA	1:B:50:ASN:HD21	1.82	0.45
1:B:177:ASN:OD1	1:B:178:PHE:N	2.49	0.45
1:C:177:ASN:OD1	1:C:178:PHE:N	2.49	0.45
1:D:177:ASN:OD1	1:D:178:PHE:N	2.49	0.45
1:F:103:ALA:O	1:F:107:VAL:HG23	2.16	0.45
1:A:47:MET:HA	1:A:50:ASN:HD21	1.82	0.45
1:C:47:MET:HA	1:C:50:ASN:HD21	1.82	0.45
1:D:2:GLU:O	1:D:6:VAL:HG23	2.17	0.44
1:D:47:MET:HA	1:D:50:ASN:HD21	1.82	0.44
1:E:47:MET:HA	1:E:50:ASN:HD21	1.82	0.44
1:F:47:MET:HA	1:F:50:ASN:HD21	1.81	0.44
1:F:2:GLU:O	1:F:6:VAL:HG23	2.17	0.44
1:A:8:ASP:OD2	1:A:16:TRP:NE1	2.51	0.44
1:C:2:GLU:O	1:C:6:VAL:HG23	2.17	0.44
1:C:78:ILE:O	1:C:82:LEU:CD1	2.61	0.44
1:E:8:ASP:OD2	1:E:16:TRP:NE1	2.51	0.44
1:E:211:SER:HB2	1:E:256:ARG:HH12	1.83	0.44
1:F:211:SER:HB2	1:F:256:ARG:HH12	1.83	0.44
1:G:8:ASP:OD2	1:G:16:TRP:NE1	2.51	0.44
2:E:301:POV:H212	2:E:301:POV:H21E	1.90	0.44
1:A:99:VAL:HG23	1:B:100:LEU:CD2	2.48	0.44
1:D:211:SER:HB2	1:D:256:ARG:HH12	1.83	0.44
1:F:8:ASP:OD2	1:F:16:TRP:NE1	2.51	0.44
1:B:211:SER:HB2	1:B:256:ARG:HH12	1.83	0.43
3:F:302:R16:C34	1:G:97:ILE:HD11	2.47	0.43
1:G:211:SER:HB2	1:G:256:ARG:HH12	1.83	0.43
1:A:211:SER:HB2	1:A:256:ARG:HH12	1.83	0.43
1:C:211:SER:HB2	1:C:256:ARG:HH12	1.83	0.43
1:D:8:ASP:OD2	1:D:16:TRP:NE1	2.51	0.43
1:B:8:ASP:OD2	1:B:16:TRP:NE1	2.51	0.43
2:B:301:POV:H212	2:B:301:POV:H21E	1.90	0.43
1:E:73:VAL:O	1:E:77:ILE:HG12	2.19	0.43
1:F:73:VAL:O	1:F:77:ILE:HG12	2.19	0.43
1:E:1:MET:HE2	1:E:15:SER:HA	2.00	0.43
2:G:301:POV:H212	2:G:301:POV:H21E	1.90	0.43
1:C:8:ASP:OD2	1:C:16:TRP:NE1	2.51	0.43
1:F:1:MET:HE2	1:F:15:SER:HA	2.00	0.42
1:G:73:VAL:O	1:G:77:ILE:HG12	2.19	0.42
1:D:73:VAL:O	1:D:77:ILE:HG12	2.19	0.42
1:B:1:MET:HE2	1:B:15:SER:HA	2.00	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1:MET:HE2	1:A:15:SER:HA	2.00	0.42
1:C:73:VAL:O	1:C:77:ILE:HG12	2.19	0.42
1:D:1:MET:HE2	1:D:15:SER:HA	2.00	0.42
1:D:223:VAL:O	1:D:224:ARG:HD2	2.19	0.42
1:A:223:VAL:O	1:A:224:ARG:HD2	2.20	0.42
1:B:73:VAL:O	1:B:77:ILE:HG12	2.19	0.42
3:A:302:R16:C31	3:B:302:R16:H291	2.42	0.42
1:B:223:VAL:O	1:B:224:ARG:HD2	2.19	0.42
1:G:123:LEU:HD23	1:G:123:LEU:HA	1.89	0.42
1:A:73:VAL:O	1:A:77:ILE:HG12	2.19	0.42
1:B:78:ILE:O	1:B:82:LEU:CD1	2.61	0.42
1:G:223:VAL:O	1:G:224:ARG:HD2	2.19	0.42
1:E:78:ILE:O	1:E:82:LEU:CD1	2.61	0.42
1:G:1:MET:HE2	1:G:15:SER:HA	2.00	0.42
1:B:246:LEU:HD12	1:B:246:LEU:HA	1.85	0.42
1:C:223:VAL:O	1:C:224:ARG:HD2	2.19	0.42
1:A:11:ASN:ND2	1:B:1:MET:SD	2.93	0.41
3:C:302:R16:H272	3:C:303:R16:H273	2.02	0.41
1:E:11:ASN:ND2	1:F:3:ASP:OD1	2.51	0.41
1:E:223:VAL:O	1:E:224:ARG:HD2	2.19	0.41
1:B:134:GLU:CD	1:B:177:ASN:HD21	2.25	0.41
1:C:1:MET:HE2	1:C:15:SER:HA	2.00	0.41
1:D:4:LEU:HD12	1:D:15:SER:HA	2.02	0.41
1:B:4:LEU:HD12	1:B:15:SER:HA	2.02	0.41
1:F:223:VAL:O	1:F:224:ARG:HD2	2.20	0.41
1:C:4:LEU:HD12	1:C:15:SER:HA	2.02	0.41
1:C:123:LEU:HD23	1:C:123:LEU:HA	1.89	0.41
1:E:4:LEU:HD12	1:E:15:SER:HA	2.02	0.41
1:B:185:ARG:HG2	1:B:186:ASN:O	2.21	0.41
1:F:3:ASP:OD2	1:F:17:LEU:HB2	2.21	0.41
1:F:185:ARG:HG2	1:F:186:ASN:O	2.21	0.41
1:G:185:ARG:HG2	1:G:186:ASN:O	2.21	0.41
1:A:3:ASP:OD2	1:A:17:LEU:HB2	2.21	0.41
1:E:185:ARG:HG2	1:E:186:ASN:O	2.21	0.41
1:F:134:GLU:CD	1:F:177:ASN:HD21	2.25	0.41
1:A:4:LEU:HD12	1:A:15:SER:HA	2.02	0.41
1:A:198:ILE:HG23	1:A:235:PHE:HZ	1.85	0.41
1:G:198:ILE:HG23	1:G:235:PHE:HZ	1.86	0.41
1:A:78:ILE:HD13	1:A:78:ILE:HA	1.91	0.41
1:D:3:ASP:OD2	1:D:17:LEU:HB2	2.21	0.40
1:D:11:ASN:ND2	1:E:3:ASP:OD1	2.52	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:ARG:HG2	1:D:186:ASN:O	2.21	0.40
1:C:3:ASP:OD2	1:C:17:LEU:HB2	2.21	0.40
1:F:4:LEU:HD12	1:F:15:SER:HA	2.02	0.40
1:B:198:ILE:HG23	1:B:235:PHE:HZ	1.86	0.40
2:F:301:POV:H212	2:F:301:POV:H21E	1.90	0.40
2:G:301:POV:H11	2:G:301:POV:H13A	1.82	0.40
1:A:185:ARG:HG2	1:A:186:ASN:O	2.21	0.40
1:C:185:ARG:HG2	1:C:186:ASN:O	2.21	0.40
1:C:198:ILE:HG23	1:C:235:PHE:HZ	1.86	0.40
1:D:134:GLU:CD	1:D:177:ASN:HD21	2.25	0.40
1:E:198:ILE:HG23	1:E:235:PHE:HZ	1.86	0.40
1:G:4:LEU:HD12	1:G:15:SER:HA	2.02	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	278/306 (91%)	268 (96%)	10 (4%)	0	100	100
1	B	278/306 (91%)	268 (96%)	10 (4%)	0	100	100
1	C	278/306 (91%)	268 (96%)	10 (4%)	0	100	100
1	D	278/306 (91%)	268 (96%)	10 (4%)	0	100	100
1	E	278/306 (91%)	268 (96%)	10 (4%)	0	100	100
1	F	278/306 (91%)	269 (97%)	9 (3%)	0	100	100
1	G	278/306 (91%)	268 (96%)	10 (4%)	0	100	100
All	All	1946/2142 (91%)	1877 (96%)	69 (4%)	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	227/248 (92%)	226 (100%)	1 (0%)	84	86
1	B	227/248 (92%)	226 (100%)	1 (0%)	84	86
1	C	227/248 (92%)	226 (100%)	1 (0%)	84	86
1	D	227/248 (92%)	226 (100%)	1 (0%)	84	86
1	E	227/248 (92%)	226 (100%)	1 (0%)	84	86
1	F	227/248 (92%)	226 (100%)	1 (0%)	84	86
1	G	227/248 (92%)	226 (100%)	1 (0%)	84	86
All	All	1589/1736 (92%)	1582 (100%)	7 (0%)	81	86

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	10	ILE
1	B	10	ILE
1	C	10	ILE
1	D	10	ILE
1	E	10	ILE
1	F	10	ILE
1	G	10	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	GLN
1	A	207	ASN
1	B	92	GLN
1	B	207	ASN
1	C	92	GLN
1	C	207	ASN
1	D	92	GLN
1	D	207	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	E	92	GLN
1	E	207	ASN
1	F	92	GLN
1	F	207	ASN
1	G	92	GLN
1	G	207	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

14 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	$\# Z > 2$
3	R16	C	303	-	15,15,15	0.12	0	14,14,14	0.18	0
3	R16	A	302	-	15,15,15	0.12	0	14,14,14	0.18	0
2	POV	E	301	-	51,51,51	1.38	4 (7%)	57,59,59	1.11	3 (5%)
3	R16	C	302	-	15,15,15	0.12	0	14,14,14	0.18	0
2	POV	D	301	-	51,51,51	1.38	4 (7%)	57,59,59	1.11	3 (5%)
3	R16	B	302	-	15,15,15	0.12	0	14,14,14	0.18	0
2	POV	G	301	-	51,51,51	1.38	4 (7%)	57,59,59	1.11	3 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	POV	C	301	-	51,51,51	1.38	4 (7%)	57,59,59	1.11	3 (5%)
3	R16	F	303	-	15,15,15	0.12	0	14,14,14	0.18	0
3	R16	E	302	-	15,15,15	0.12	0	14,14,14	0.18	0
2	POV	A	301	-	51,51,51	1.38	4 (7%)	57,59,59	1.11	3 (5%)
3	R16	F	302	-	15,15,15	0.12	0	14,14,14	0.18	0
2	POV	B	301	-	51,51,51	1.38	4 (7%)	57,59,59	1.11	3 (5%)
2	POV	F	301	-	51,51,51	1.38	4 (7%)	57,59,59	1.11	3 (5%)

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
3	R16	C	303	-	-	1/13/13/13	-
3	R16	A	302	-	-	1/13/13/13	-
2	POV	E	301	-	-	22/55/55/55	-
3	R16	C	302	-	-	1/13/13/13	-
2	POV	D	301	-	-	22/55/55/55	-
3	R16	B	302	-	-	1/13/13/13	-
2	POV	G	301	-	-	22/55/55/55	-
2	POV	C	301	-	-	22/55/55/55	-
3	R16	F	303	-	-	1/13/13/13	-
3	R16	E	302	-	-	1/13/13/13	-
2	POV	A	301	-	-	22/55/55/55	-
3	R16	F	302	-	-	1/13/13/13	-
2	POV	B	301	-	-	22/55/55/55	-
2	POV	F	301	-	-	22/55/55/55	-

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	301	POV	O31-C31	3.48	1.43	1.33
2	C	301	POV	O31-C31	3.47	1.43	1.33
2	A	301	POV	O31-C31	3.47	1.43	1.33
2	E	301	POV	O31-C31	3.46	1.43	1.33
2	B	301	POV	O31-C31	3.46	1.43	1.33
2	D	301	POV	O31-C31	3.46	1.43	1.33
2	G	301	POV	O31-C31	3.45	1.43	1.33
2	F	301	POV	O21-C2	-3.13	1.39	1.46

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	301	POV	O21-C2	-3.11	1.39	1.46
2	B	301	POV	O21-C2	-3.09	1.39	1.46
2	C	301	POV	O21-C2	-3.08	1.39	1.46
2	D	301	POV	O21-C2	-3.08	1.39	1.46
2	A	301	POV	O21-C2	-3.07	1.39	1.46
2	E	301	POV	O21-C2	-3.07	1.39	1.46
2	G	301	POV	O21-C21	2.94	1.42	1.34
2	D	301	POV	O21-C21	2.94	1.42	1.34
2	F	301	POV	O21-C21	2.93	1.42	1.34
2	B	301	POV	O21-C21	2.93	1.42	1.34
2	E	301	POV	O21-C21	2.92	1.42	1.34
2	A	301	POV	O21-C21	2.92	1.42	1.34
2	C	301	POV	O21-C21	2.91	1.42	1.34
2	C	301	POV	P-O11	2.68	1.69	1.59
2	E	301	POV	P-O11	2.66	1.69	1.59
2	A	301	POV	P-O11	2.66	1.69	1.59
2	D	301	POV	P-O11	2.65	1.69	1.59
2	F	301	POV	P-O11	2.65	1.69	1.59
2	B	301	POV	P-O11	2.65	1.69	1.59
2	G	301	POV	P-O11	2.63	1.69	1.59

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	301	POV	O21-C21-C22	4.38	120.97	111.48
2	F	301	POV	O21-C21-C22	4.38	120.95	111.48
2	A	301	POV	O21-C21-C22	4.37	120.94	111.48
2	C	301	POV	O21-C21-C22	4.37	120.93	111.48
2	E	301	POV	O21-C21-C22	4.36	120.92	111.48
2	B	301	POV	O21-C21-C22	4.36	120.91	111.48
2	D	301	POV	O21-C21-C22	4.34	120.88	111.48
2	C	301	POV	O31-C31-C32	2.57	119.66	111.83
2	B	301	POV	O31-C31-C32	2.56	119.64	111.83
2	E	301	POV	O31-C31-C32	2.56	119.64	111.83
2	D	301	POV	O31-C31-C32	2.56	119.63	111.83
2	A	301	POV	O31-C31-C32	2.55	119.62	111.83
2	G	301	POV	O31-C31-C32	2.55	119.62	111.83
2	F	301	POV	O31-C31-C32	2.54	119.59	111.83
2	D	301	POV	C2-O21-C21	-2.13	112.70	117.80
2	B	301	POV	C2-O21-C21	-2.13	112.71	117.80
2	A	301	POV	C2-O21-C21	-2.12	112.72	117.80
2	G	301	POV	C2-O21-C21	-2.12	112.73	117.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	301	POV	C2-O21-C21	-2.12	112.73	117.80
2	C	301	POV	C2-O21-C21	-2.11	112.76	117.80
2	F	301	POV	C2-O21-C21	-2.10	112.76	117.80

There are no chirality outliers.

All (161) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	POV	C1-O11-P-O14
2	B	301	POV	C1-O11-P-O14
2	C	301	POV	C1-O11-P-O14
2	D	301	POV	C1-O11-P-O14
2	E	301	POV	C1-O11-P-O14
2	F	301	POV	C1-O11-P-O14
2	G	301	POV	C1-O11-P-O14
2	A	301	POV	C21-C22-C23-C24
2	B	301	POV	C21-C22-C23-C24
2	C	301	POV	C21-C22-C23-C24
2	D	301	POV	C21-C22-C23-C24
2	E	301	POV	C21-C22-C23-C24
2	F	301	POV	C21-C22-C23-C24
2	G	301	POV	C21-C22-C23-C24
2	A	301	POV	C23-C24-C25-C26
2	B	301	POV	C23-C24-C25-C26
2	C	301	POV	C23-C24-C25-C26
2	D	301	POV	C23-C24-C25-C26
2	E	301	POV	C23-C24-C25-C26
2	F	301	POV	C23-C24-C25-C26
2	G	301	POV	C23-C24-C25-C26
2	B	301	POV	C25-C26-C27-C28
2	C	301	POV	C25-C26-C27-C28
2	E	301	POV	C25-C26-C27-C28
2	A	301	POV	C25-C26-C27-C28
2	D	301	POV	C25-C26-C27-C28
2	F	301	POV	C25-C26-C27-C28
2	G	301	POV	C25-C26-C27-C28
2	A	301	POV	C31-C32-C33-C34
2	B	301	POV	C31-C32-C33-C34
2	C	301	POV	C31-C32-C33-C34
2	D	301	POV	C31-C32-C33-C34
2	E	301	POV	C31-C32-C33-C34
2	F	301	POV	C31-C32-C33-C34

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	G	301	POV	C31-C32-C33-C34
2	C	301	POV	C212-C213-C214-C215
2	E	301	POV	C212-C213-C214-C215
2	F	301	POV	C212-C213-C214-C215
2	G	301	POV	C212-C213-C214-C215
2	A	301	POV	C212-C213-C214-C215
2	B	301	POV	C212-C213-C214-C215
2	D	301	POV	C212-C213-C214-C215
2	C	301	POV	C213-C214-C215-C216
2	E	301	POV	C213-C214-C215-C216
2	F	301	POV	C213-C214-C215-C216
2	G	301	POV	C213-C214-C215-C216
2	A	301	POV	C213-C214-C215-C216
2	B	301	POV	C213-C214-C215-C216
2	D	301	POV	C213-C214-C215-C216
2	A	301	POV	C210-C211-C212-C213
2	B	301	POV	C210-C211-C212-C213
2	C	301	POV	C210-C211-C212-C213
2	D	301	POV	C210-C211-C212-C213
2	E	301	POV	C210-C211-C212-C213
2	F	301	POV	C210-C211-C212-C213
2	G	301	POV	C210-C211-C212-C213
2	A	301	POV	C1-C2-C3-O31
2	B	301	POV	C1-C2-C3-O31
2	C	301	POV	C1-C2-C3-O31
2	D	301	POV	C1-C2-C3-O31
2	E	301	POV	C1-C2-C3-O31
2	F	301	POV	C1-C2-C3-O31
2	G	301	POV	C1-C2-C3-O31
2	A	301	POV	O11-C1-C2-O21
2	B	301	POV	O11-C1-C2-O21
2	C	301	POV	O11-C1-C2-O21
2	D	301	POV	O11-C1-C2-O21
2	E	301	POV	O11-C1-C2-O21
2	F	301	POV	O11-C1-C2-O21
2	G	301	POV	O11-C1-C2-O21
2	A	301	POV	O21-C2-C3-O31
2	B	301	POV	O21-C2-C3-O31
2	C	301	POV	O21-C2-C3-O31
2	D	301	POV	O21-C2-C3-O31
2	E	301	POV	O21-C2-C3-O31
2	F	301	POV	O21-C2-C3-O31

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	G	301	POV	O21-C2-C3-O31
2	A	301	POV	C34-C35-C36-C37
2	B	301	POV	C34-C35-C36-C37
2	D	301	POV	C34-C35-C36-C37
2	E	301	POV	C34-C35-C36-C37
2	C	301	POV	C34-C35-C36-C37
2	F	301	POV	C34-C35-C36-C37
2	G	301	POV	C34-C35-C36-C37
2	A	301	POV	C2-C1-O11-P
2	B	301	POV	C2-C1-O11-P
2	C	301	POV	C2-C1-O11-P
2	D	301	POV	C2-C1-O11-P
2	E	301	POV	C2-C1-O11-P
2	F	301	POV	C2-C1-O11-P
2	G	301	POV	C2-C1-O11-P
2	B	301	POV	C310-C311-C312-C313
2	A	301	POV	C310-C311-C312-C313
2	G	301	POV	C310-C311-C312-C313
2	C	301	POV	C310-C311-C312-C313
2	E	301	POV	C310-C311-C312-C313
2	F	301	POV	C310-C311-C312-C313
2	D	301	POV	C310-C311-C312-C313
2	A	301	POV	O11-C1-C2-C3
2	B	301	POV	O11-C1-C2-C3
2	C	301	POV	O11-C1-C2-C3
2	D	301	POV	O11-C1-C2-C3
2	E	301	POV	O11-C1-C2-C3
2	F	301	POV	O11-C1-C2-C3
2	G	301	POV	O11-C1-C2-C3
2	A	301	POV	O12-C11-C12-N
2	B	301	POV	O12-C11-C12-N
2	C	301	POV	O12-C11-C12-N
2	D	301	POV	O12-C11-C12-N
2	E	301	POV	O12-C11-C12-N
2	F	301	POV	O12-C11-C12-N
2	G	301	POV	O12-C11-C12-N
2	B	301	POV	C24-C25-C26-C27
2	A	301	POV	C24-C25-C26-C27
2	D	301	POV	C24-C25-C26-C27
2	C	301	POV	C24-C25-C26-C27
2	F	301	POV	C24-C25-C26-C27
2	E	301	POV	C24-C25-C26-C27

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms
2	G	301	POV	C24-C25-C26-C27
2	B	301	POV	C22-C23-C24-C25
2	G	301	POV	C22-C23-C24-C25
2	C	301	POV	C22-C23-C24-C25
2	E	301	POV	C22-C23-C24-C25
2	A	301	POV	C22-C23-C24-C25
2	D	301	POV	C22-C23-C24-C25
2	F	301	POV	C22-C23-C24-C25
2	A	301	POV	C27-C28-C29-C210
2	E	301	POV	C27-C28-C29-C210
2	F	301	POV	C27-C28-C29-C210
2	G	301	POV	C27-C28-C29-C210
2	A	301	POV	C29-C210-C211-C212
2	B	301	POV	C29-C210-C211-C212
2	B	301	POV	C27-C28-C29-C210
2	C	301	POV	C29-C210-C211-C212
2	C	301	POV	C27-C28-C29-C210
2	D	301	POV	C29-C210-C211-C212
2	D	301	POV	C27-C28-C29-C210
2	E	301	POV	C29-C210-C211-C212
2	F	301	POV	C29-C210-C211-C212
2	G	301	POV	C29-C210-C211-C212
2	C	301	POV	C311-C310-C39-C38
2	D	301	POV	C311-C310-C39-C38
2	E	301	POV	C311-C310-C39-C38
2	A	301	POV	C311-C310-C39-C38
2	F	301	POV	C311-C310-C39-C38
2	G	301	POV	C311-C310-C39-C38
2	B	301	POV	C311-C310-C39-C38
2	A	301	POV	O31-C31-C32-C33
2	B	301	POV	O31-C31-C32-C33
2	C	301	POV	O31-C31-C32-C33
2	E	301	POV	O31-C31-C32-C33
2	F	301	POV	O31-C31-C32-C33
2	G	301	POV	O31-C31-C32-C33
2	D	301	POV	O31-C31-C32-C33
3	A	302	R16	C27-C28-C29-C30
3	F	302	R16	C27-C28-C29-C30
3	B	302	R16	C27-C28-C29-C30
3	C	303	R16	C27-C28-C29-C30
3	F	303	R16	C27-C28-C29-C30
3	C	302	R16	C27-C28-C29-C30

Continued on next page...

Continued from previous page...

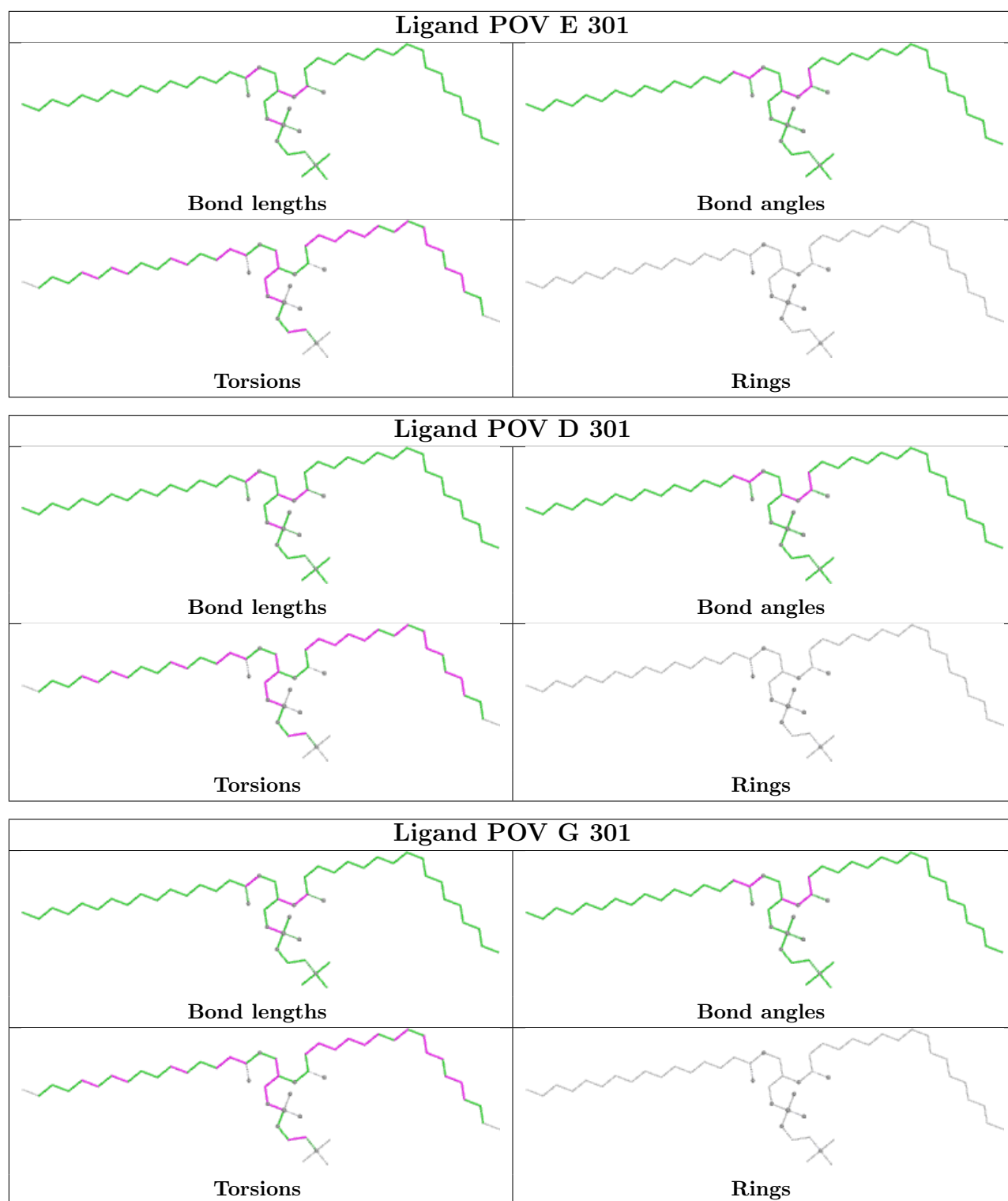
Mol	Chain	Res	Type	Atoms
3	E	302	R16	C27-C28-C29-C30

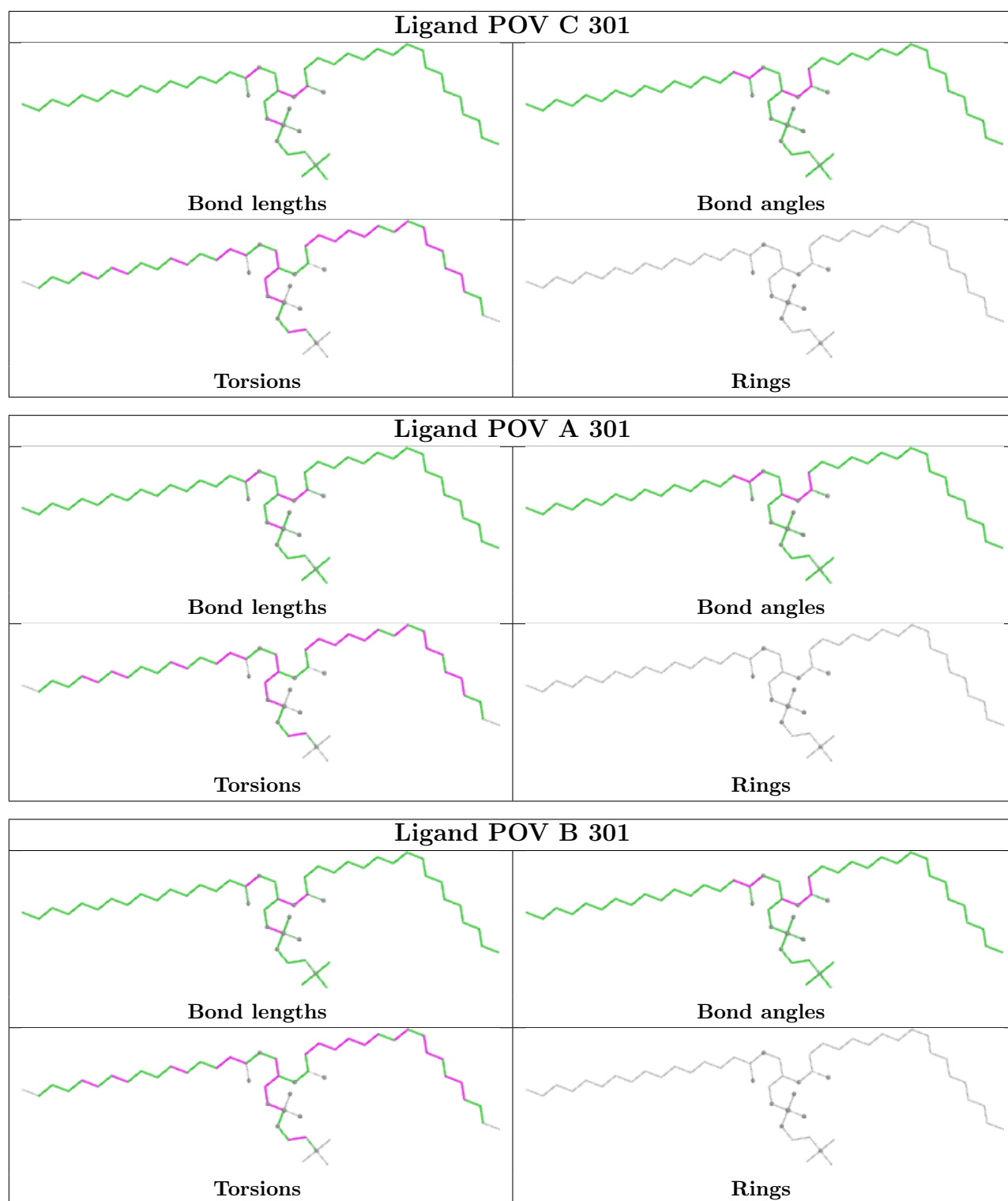
There are no ring outliers.

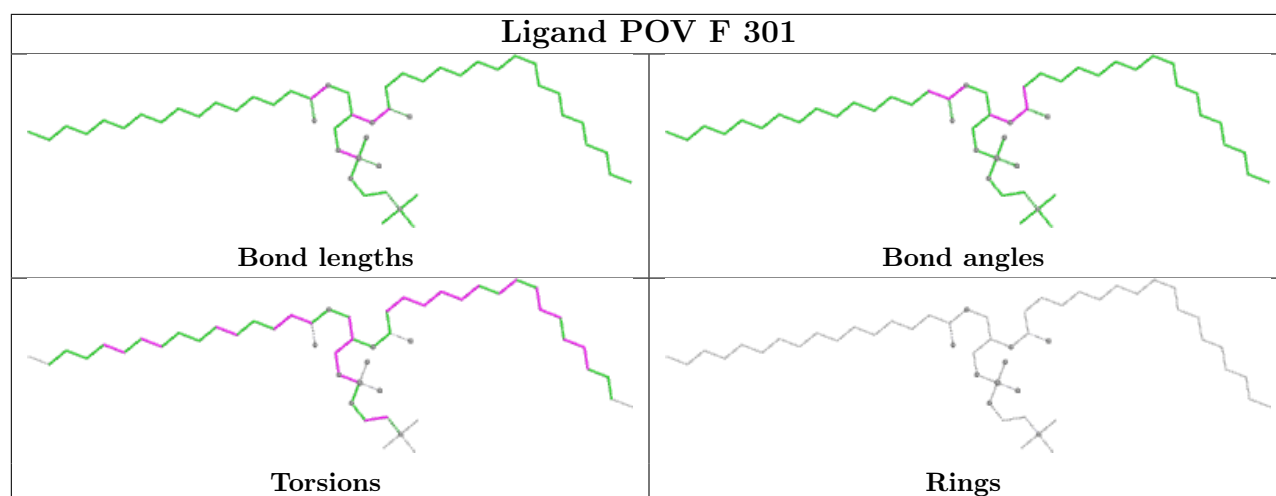
13 monomers are involved in 25 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	303	R16	2	0
3	A	302	R16	6	0
2	E	301	POV	2	0
3	C	302	R16	3	0
2	D	301	POV	1	0
3	B	302	R16	5	0
2	G	301	POV	3	0
2	C	301	POV	1	0
3	E	302	R16	1	0
2	A	301	POV	1	0
3	F	302	R16	4	0
2	B	301	POV	2	0
2	F	301	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 [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-20508. These allow visual inspection of the internal detail of the map and identification of artifacts.

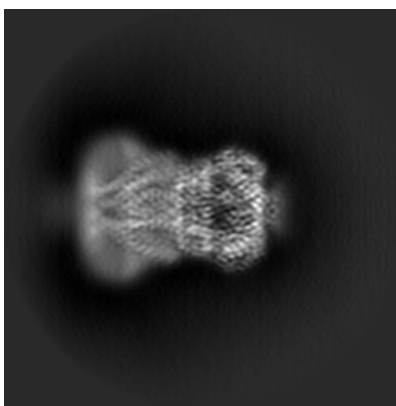
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

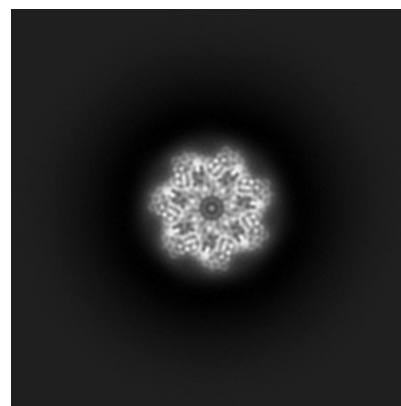
6.1.1 Primary map



X

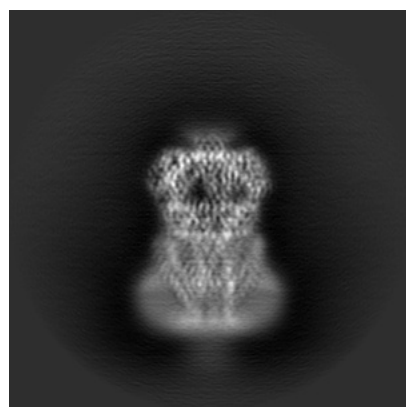


Y

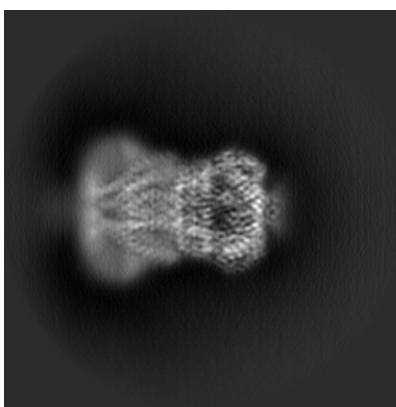


Z

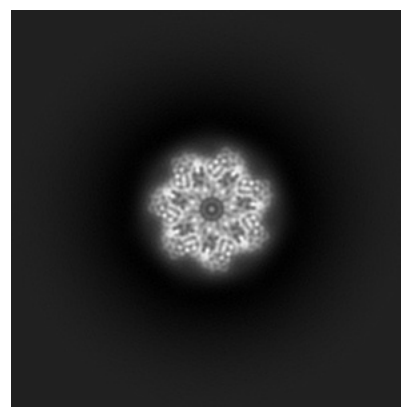
6.1.2 Raw map



X



Y



Z

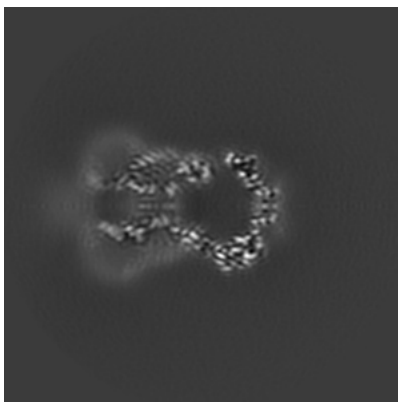
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

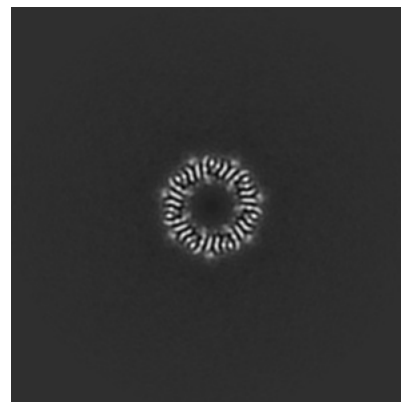
6.2.1 Primary map



X Index: 128

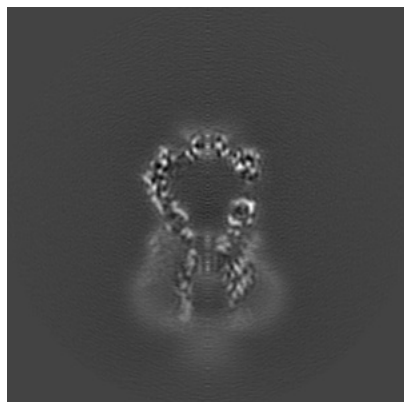


Y Index: 128

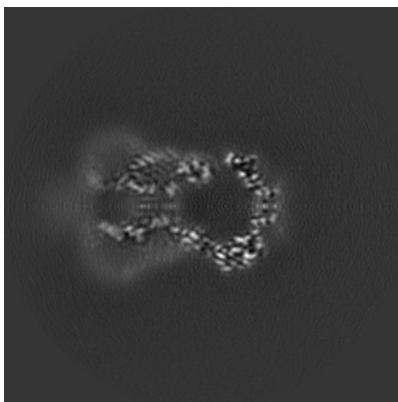


Z Index: 128

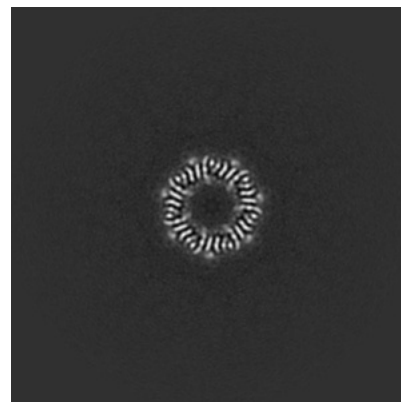
6.2.2 Raw map



X Index: 128



Y Index: 128

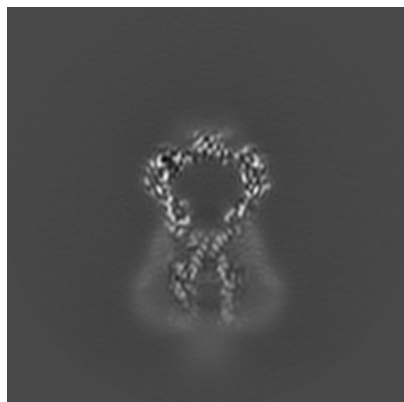


Z Index: 128

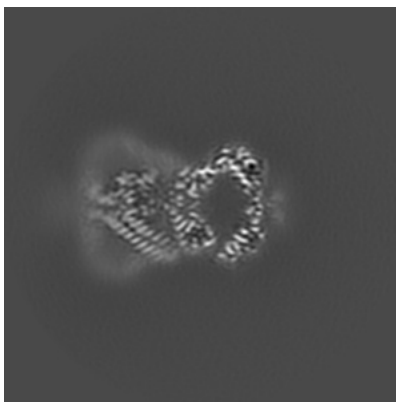
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

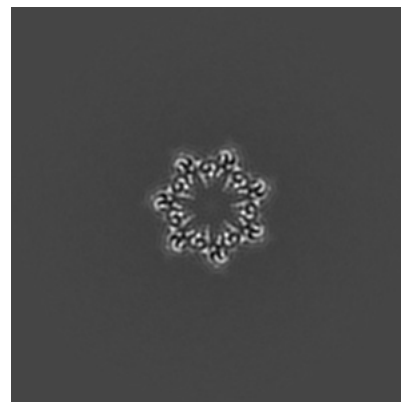
6.3.1 Primary map



X Index: 135

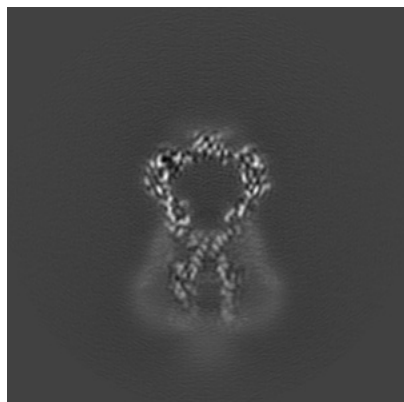


Y Index: 141

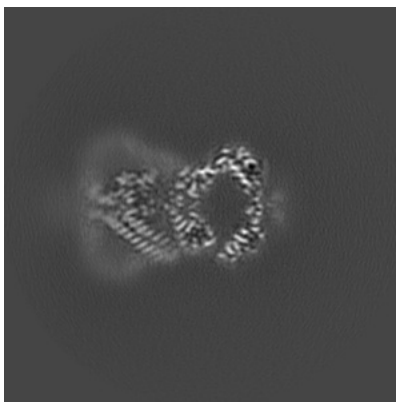


Z Index: 154

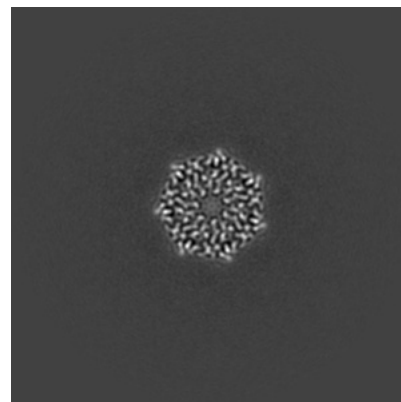
6.3.2 Raw map



X Index: 135



Y Index: 141



Z Index: 159

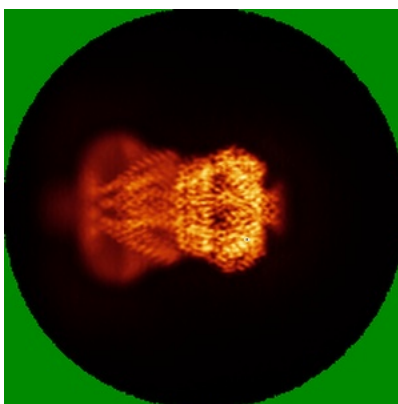
The images above show the largest variance slices of the map in three orthogonal directions.

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

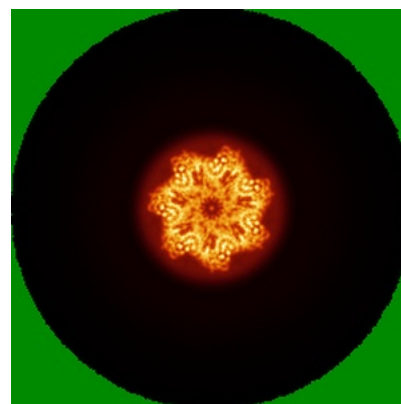
6.4.1 Primary map



X

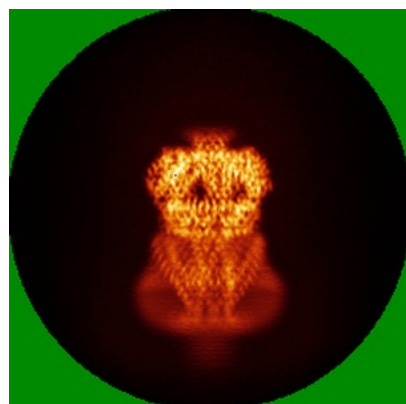


Y

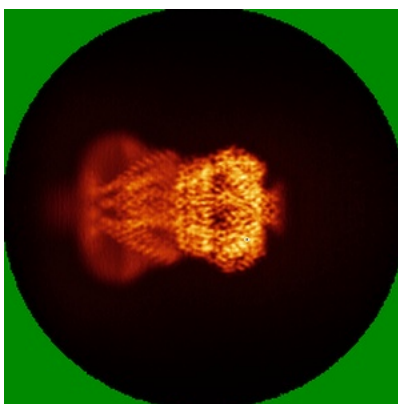


Z

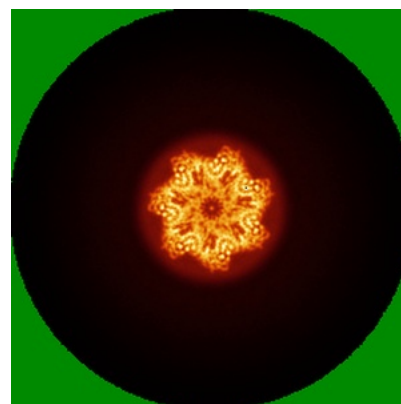
6.4.2 Raw map



X



Y

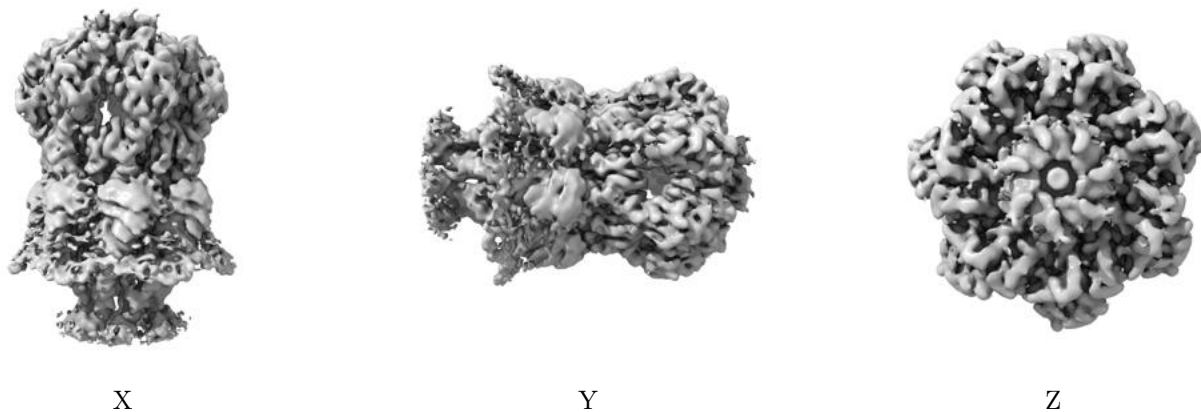


Z

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.

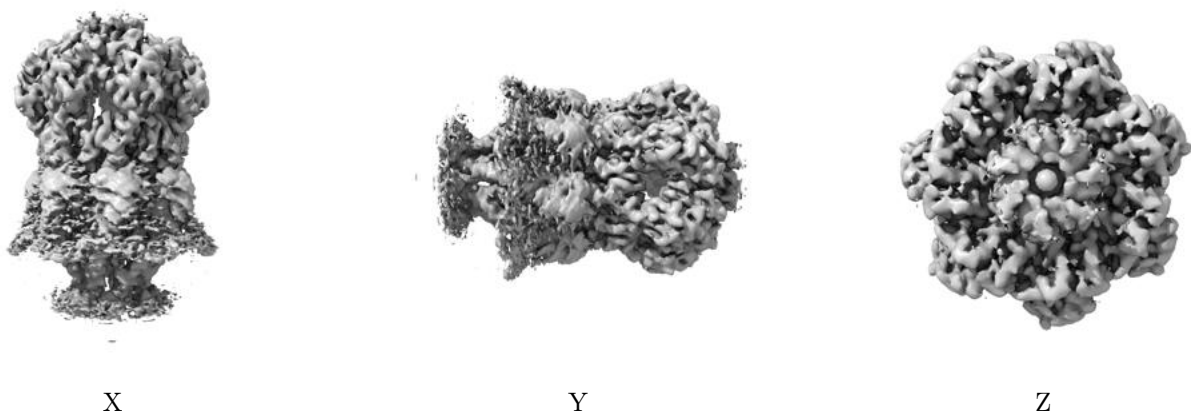
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0092. 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.

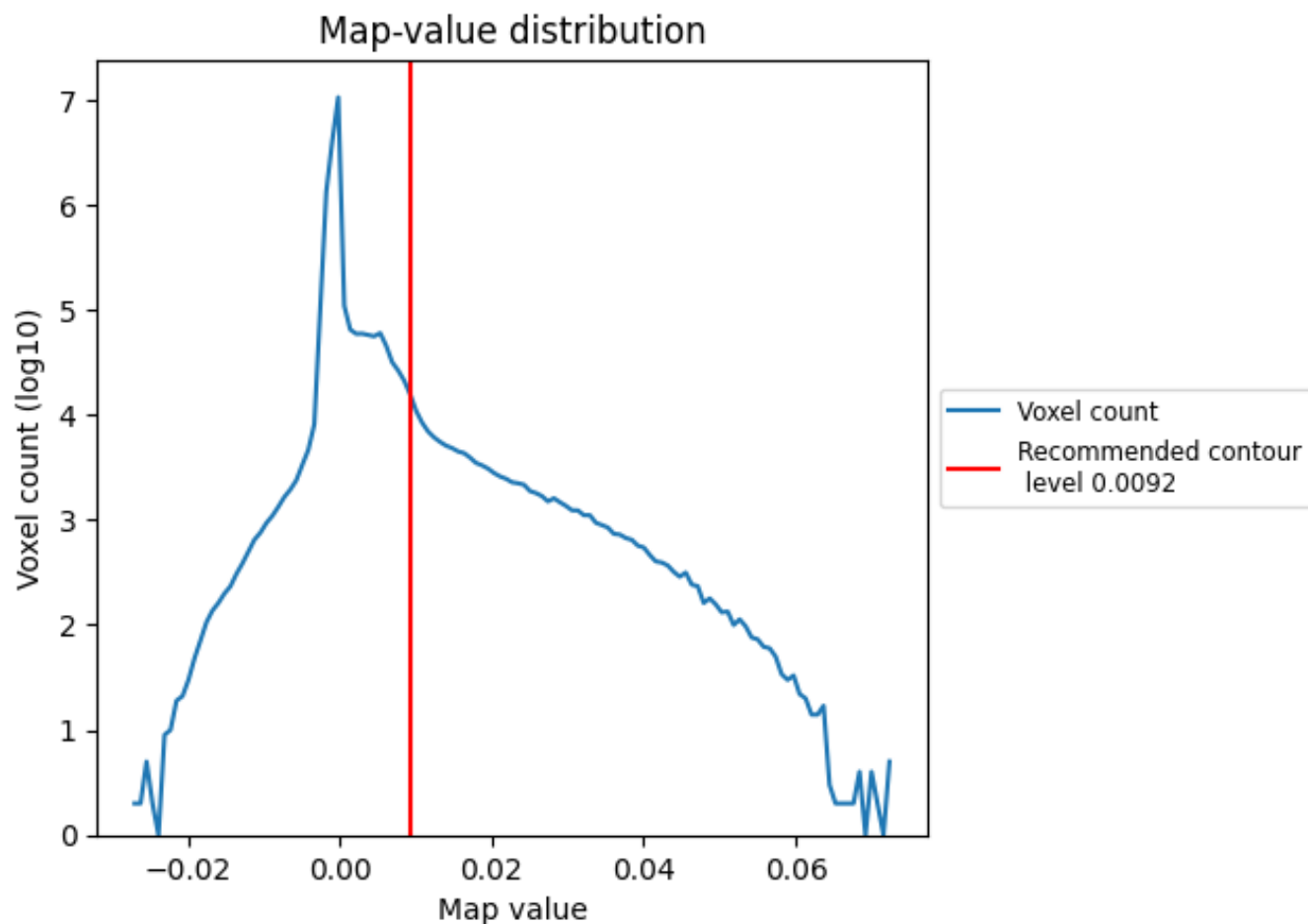
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

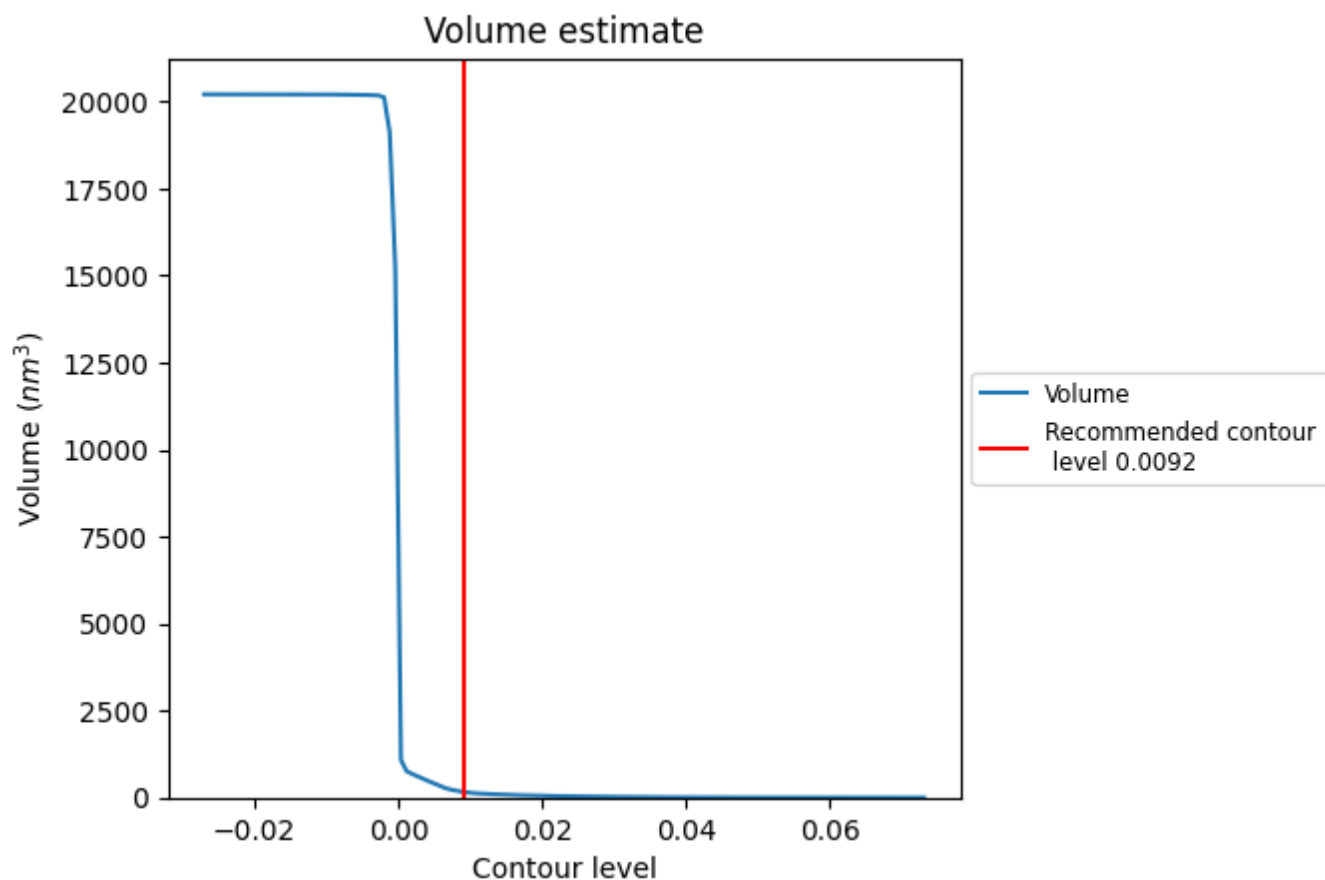
This section contains the results of statistical analysis of the map.

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.

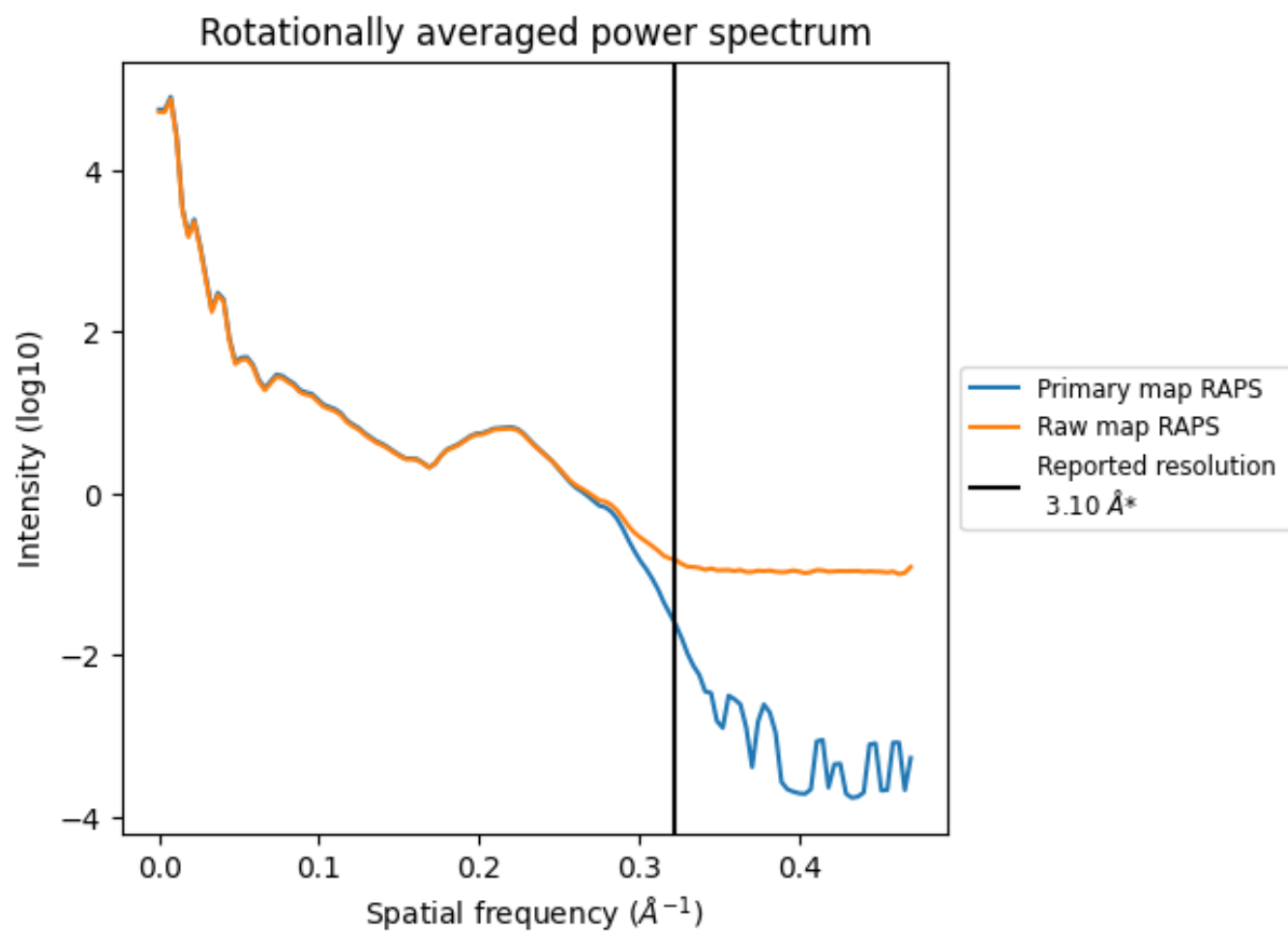
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 156 nm^3 ; this corresponds to an approximate mass of 141 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 ⓘ

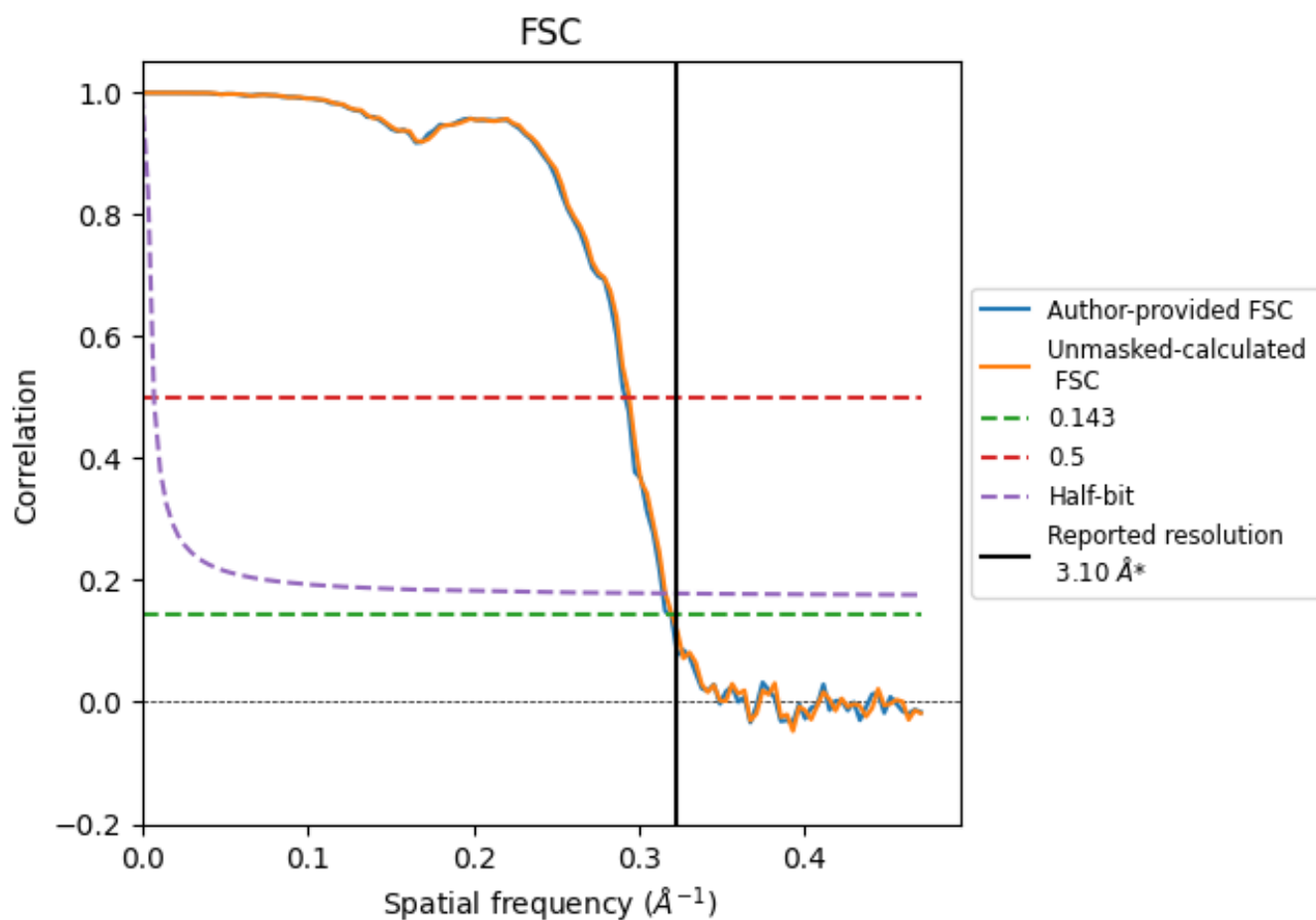


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

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 0.323 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

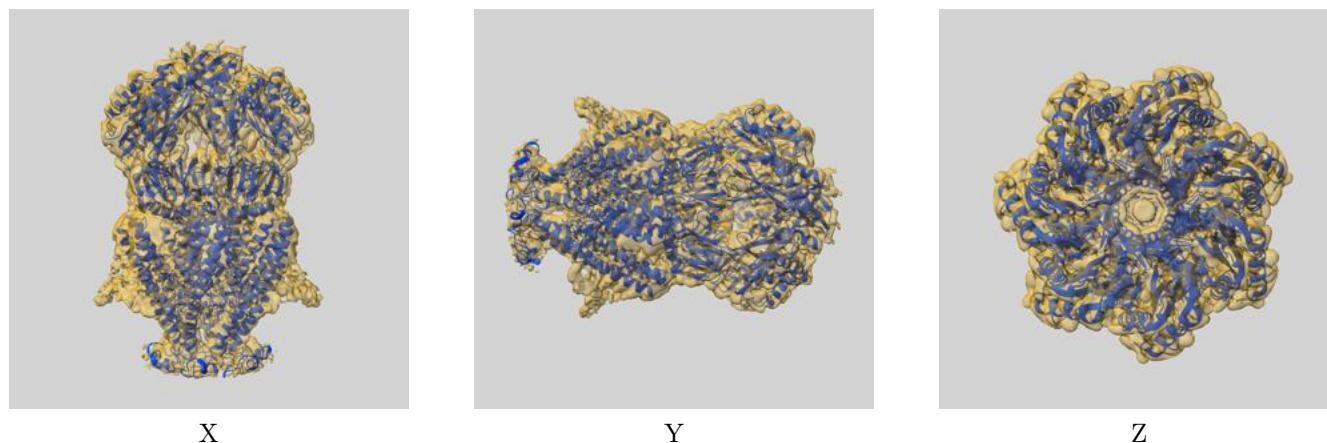
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.14	3.43	3.18
Unmasked-calculated*	3.13	3.40	3.17

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

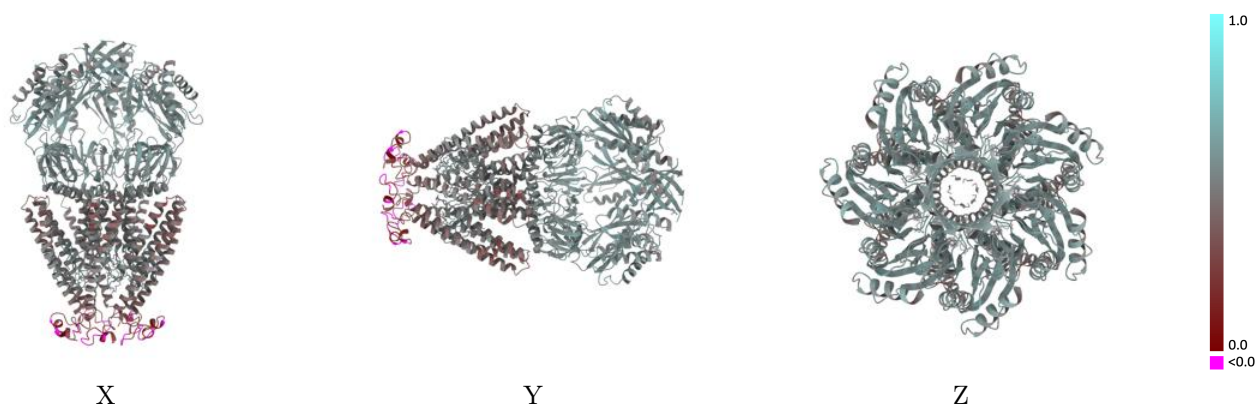
This section contains information regarding the fit between EMDB map EMD-20508 and PDB model 6PWN. Per-residue inclusion information can be found in section [3](#) on page [10](#).

9.1 Map-model overlay [i](#)



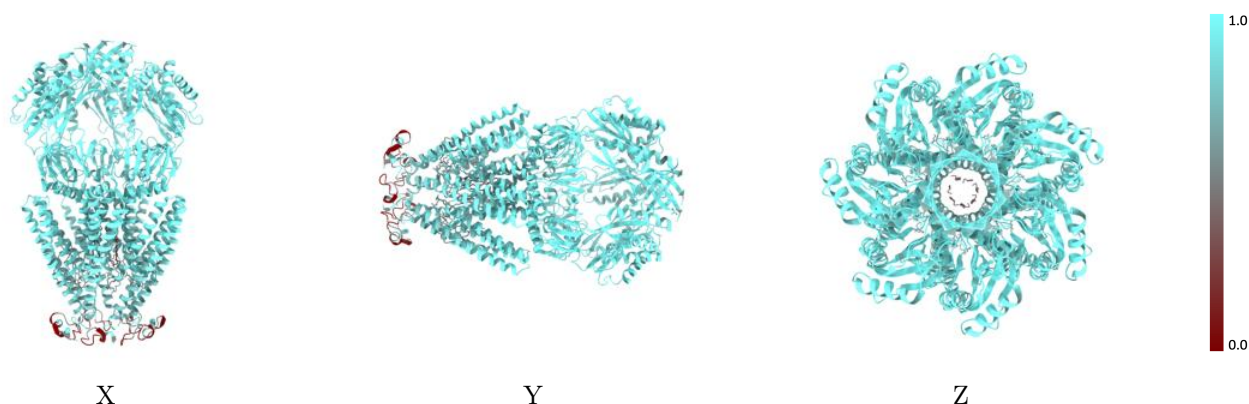
The images above show the 3D surface view of the map at the recommended contour level 0.0092 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



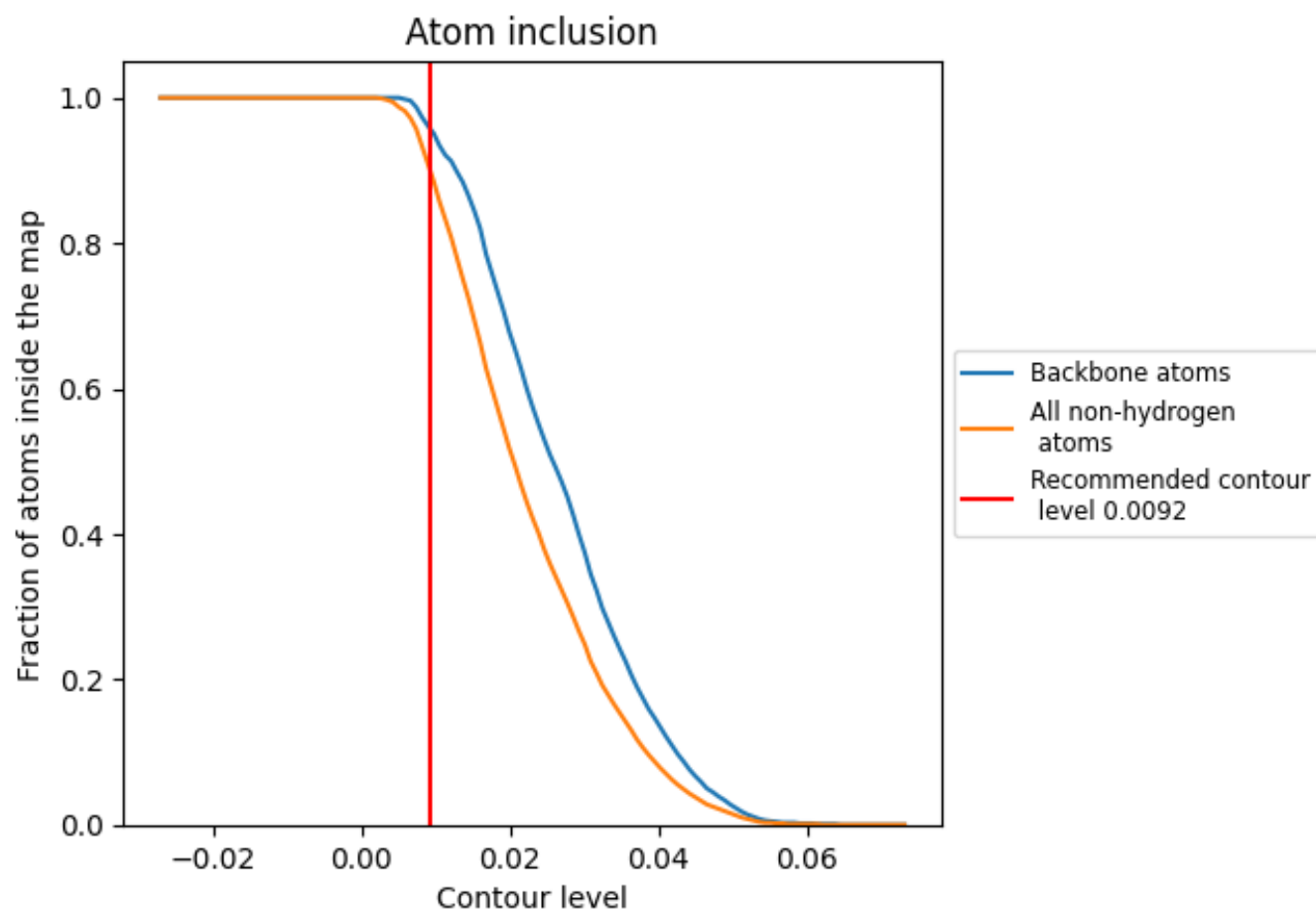
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0092).

9.4 Atom inclusion [i](#)



At the recommended contour level, 96% of all backbone atoms, 90% 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 (0.0092) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9020	0.4820
A	0.9020	0.4810
B	0.9010	0.4830
C	0.8980	0.4830
D	0.9070	0.4840
E	0.9020	0.4820
F	0.8980	0.4810
G	0.9090	0.4810

1.0

0.0

<0.0