



## Full wwPDB EM Validation Report ⓘ

Apr 22, 2026 – 02:55 PM EDT

PDB ID : 9ZKS / pdb\_00009zks  
EMDB ID : EMD-74383  
Title : The LBD-TMD structure of native mouse AMPAR with 2 TARPs 2 CNIHs and PRRT1/SynDIG4  
Authors : Fang, C.L.; Gouaux, E.  
Deposited on : 2025-12-07  
Resolution : 3.36 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](mailto: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>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

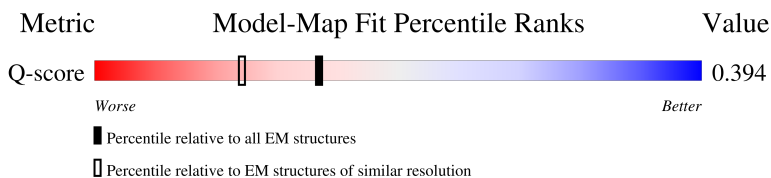
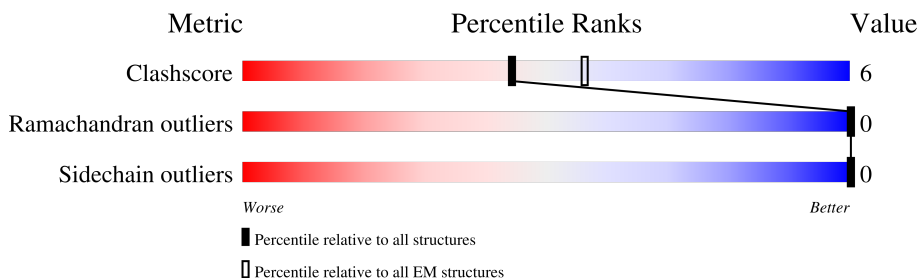
EMDB validation analysis : 0.0.1.dev132  
Mogul : 2022.3.0, CSD as543be (2022)  
MolProbity : 4-5-2 with Phenix2.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
EM percentile statistics : 202505.v01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




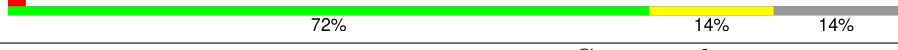
The reported resolution of this entry is 3.36 Å.

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	14332 ( 2.86 - 3.86 )

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  $\geq 3$ , 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	500	
1	C	500	
2	B	467	
2	D	467	

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Mol	Chain	Length	Quality of chain
3	E	160	<div><div></div><div>14%</div><div>60%</div><div>18%</div><div>22%</div></div>
3	F	160	<div><div></div><div>77%</div><div>14%</div><div>9%</div></div>
4	G	423	<div><div></div><div>33%</div><div>5%</div><div>62%</div></div>
4	H	423	<div><div></div><div>32%</div><div>6%</div><div>62%</div></div>
5	M	306	<div><div></div><div>9%</div><div>87%</div></div>

## 2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 16983 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 Glutamate receptor 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	403	Total	C	N	O	S	0	0
			2910	1902	462	530	16		
1	C	421	Total	C	N	O	S	0	0
			3063	1999	492	554	18		

- Molecule 2 is a protein called Glutamate receptor 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	407	Total	C	N	O	S	0	0
			2950	1915	480	538	17		
2	D	403	Total	C	N	O	S	0	0
			2902	1885	474	525	18		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	743	ARG	GLY	conflict	UNP C9K0Z0
D	743	ARG	GLY	conflict	UNP C9K0Z0

- Molecule 3 is a protein called Protein cornichon homolog 2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	124	Total	C	N	O	S	0	0
			969	665	142	151	11		
3	F	146	Total	C	N	O	S	0	0
			1177	794	183	188	12		

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	40	ASP	ASN	conflict	UNP O35089
F	40	ASP	ASN	conflict	UNP O35089

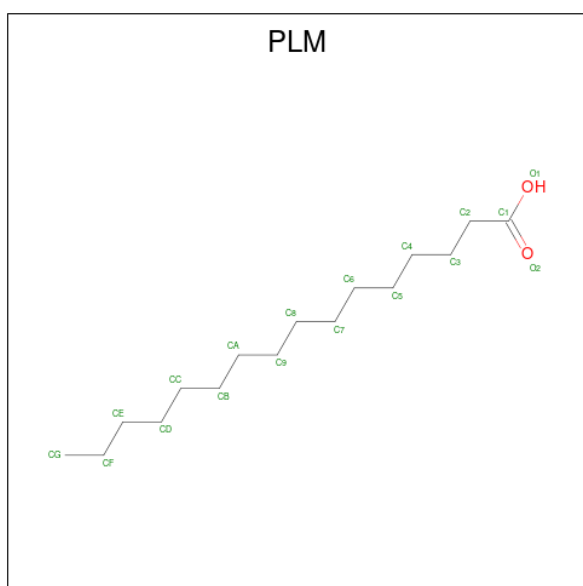
- Molecule 4 is a protein called Voltage-dependent calcium channel gamma-8 subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	162	Total	C	N	O	S	0	0
			1163	770	188	200	5		
4	H	162	Total	C	N	O	S	0	0
			1156	764	188	199	5		

- Molecule 5 is a protein called Proline-rich transmembrane protein 1.

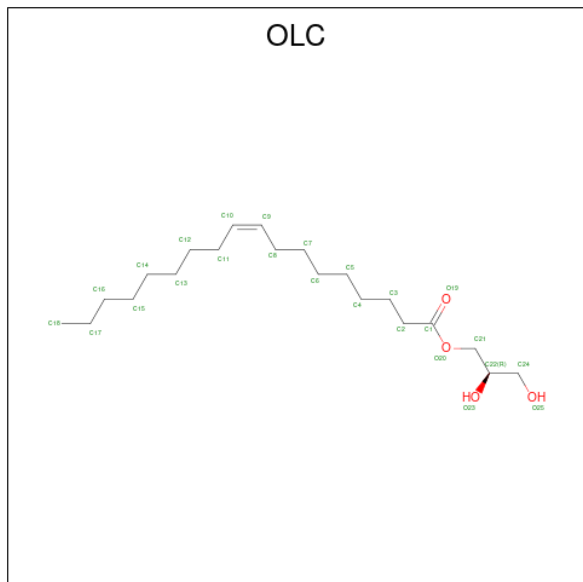
Mol	Chain	Residues	Atoms					AltConf	Trace
5	M	39	Total	C	N	O	S	0	0
			283	183	47	51	2		

- Molecule 6 is PALMITIC ACID (CCD ID: PLM) (formula:  $C_{16}H_{32}O_2$ ).



Mol	Chain	Residues	Atoms			AltConf
6	A	1	Total	C	O	0
			18	16	2	
6	A	1	Total	C	O	0
			13	11	2	
6	B	1	Total	C	O	0
			18	16	2	
6	C	1	Total	C	O	0
			18	16	2	
6	C	1	Total	C	O	0
			13	11	2	
6	D	1	Total	C	O	0
			18	16	2	

- Molecule 7 is (2R)-2,3-dihydroxypropyl (9Z)-octadec-9-enoate (CCD ID: OLC) (formula:  $C_{21}H_{40}O_4$ ).



Mol	Chain	Residues	Atoms			AltConf
7	A	1	Total	C	O	0
			25	21	4	
7	A	1	Total	C	O	0
			16	14	2	
7	B	1	Total	C	O	0
			25	21	4	
7	B	1	Total	C	O	0
			14	12	2	
7	C	1	Total	C	O	0
			25	21	4	
7	C	1	Total	C	O	0
			16	14	2	
7	D	1	Total	C	O	0
			25	21	4	
7	D	1	Total	C	O	0
			15	13	2	

- Molecule 8 is (2S)-3-(hexadecanoyloxy)-2-[(9Z)-octadec-9-enoyloxy]propyl 2-(trimethylamm onio)ethyl phosphate (CCD ID: POV) (formula:  $C_{42}H_{82}NO_8P$ ).

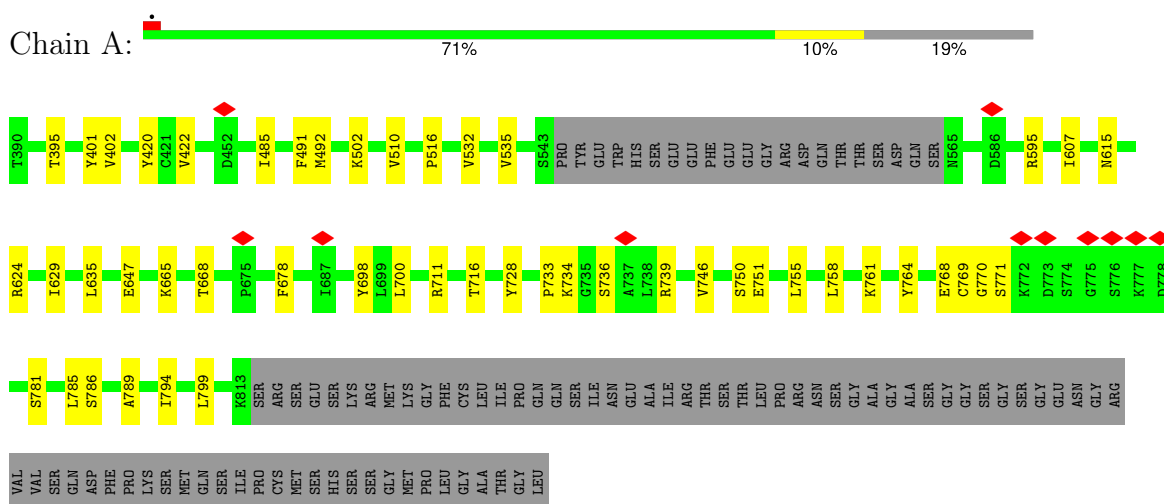


Mol	Chain	Residues	Atoms					AltConf
8	A	1	Total	C	N	O	P	0
			51	41	1	8	1	
8	D	1	Total	C	N	O	P	0
			50	40	1	8	1	
8	G	1	Total	C	N	O	P	0
			50	40	1	8	1	

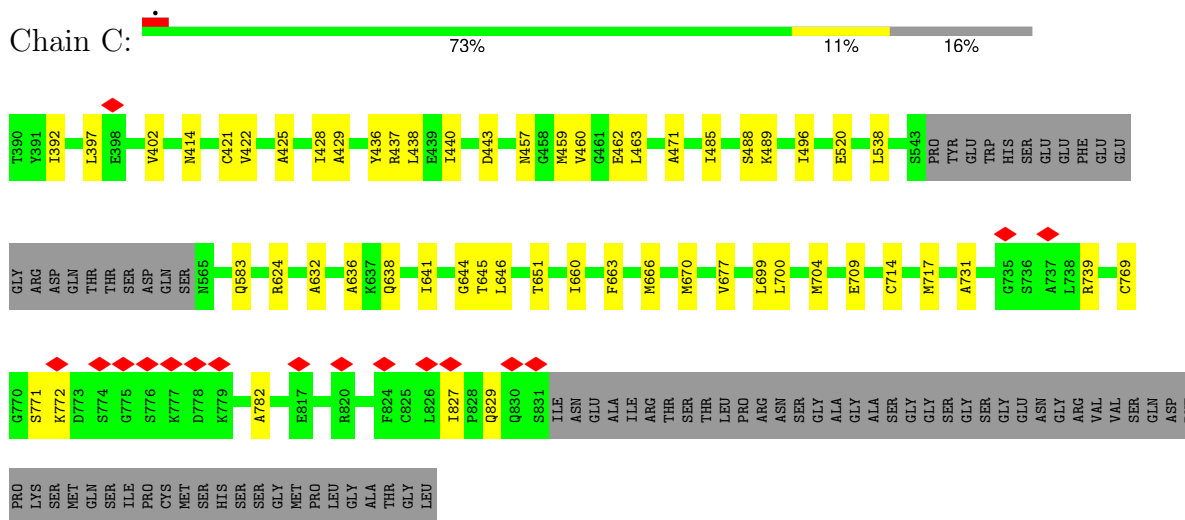
### 3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion in map density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red diamond above a residue indicates a poor fit to the EM map for this residue (all-atom inclusion < 40%). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

- Molecule 1: Glutamate receptor 1



- Molecule 1: Glutamate receptor 1



- Molecule 2: Glutamate receptor 2









ILE	VAL	PRO	ALA	MET
ILE	TYR	LEU	GLY	SER
ILE	PRO	PRO	GLY	SER
PHE	VAL	PRO	LEU	GLU
LYS	GLY	PRO	LEU	LYS
ALA	THR	PRO	SER	SER
VAL	PRO	PRO	ALA	GLY
GLN	TYR	ALA	ALA	LEU
VAL	ALA	ALA	ALA	ASP
ARG	GLY	ALA	SER	PRO
THR	GLY	ALA	ALA	SER
ALA	THR	PRO	GLN	VAL
LEU	PRO	PRO	ARG	PRO
ALA	GLY	PRO	GLY	HIS
ALA	GLY	PRO	PRO	THR
ARG	GLY	PRO	SER	SER
GLY	GLY	PRO	SER	PRO
ASP	GLY	ALA	SER	PRO
THR	VAL	ALA	SER	PRO
ALA	THR	PRO	ALA	PRO
LEU	THR	PRO	THR	TYR
S260	SER	ALA	ASN	ALA
1263	THR	LEU	LEU	PRO
	PRO	GLN	ARG	PRO
R266	PRO	PRO	PRO	GLN
	PRO	GLY	PRO	PRO
N270	PRO	PRO	HIS	PRO
	GLN	VAL	HIS	ALA
F273	GLY	VAL	ALA	GLU
1274	PRO	PRO	PRO	PRO
S275	GLY	THR	PRO	PRO
L276	LEU	HIS	GLY	ILE
	ALA	ALA	PRO	PRO
	LEU	GLY	ALA	PRO
C286	LEU	VAL	GLY	GLN
	GLU	VAL	ALA	THR
L289	PRO	GLY	ALA	ALA
T280	ARG	THR	PRO	ALA
	ARG	LEU	PRO	PRO
I283	PRO	PRO	PRO	SER
	PRO	LEU	GLY	SER
A287	HIS	GLY	CYS	HIS
Q288	ASP	GLY	ALA	HIS
	TYR	TYR	THR	HIS
HIS	MET	VAL	LEU	HIS
HIS	PRO	ALA	LEU	LEU
ASN	PRO	ARG	PRO	HIS
TYR	ILE	PRO	ARG	HIS
TRP	ALA	GLY	MET	TYR
ASP	VAL	TYR	PRO	THR
PRO	LEU	PRO	PRO	HIS
	THR	LEU	ASP	SER
	THR	GLN	PRO	SER
	ILE	LEU	TYR	GLY
	CYS	GLN	LEU	THR
	CYS	PRO	GLN	ALA
	PHE	CYS	GLU	THR
	TRP	THR	THR	PRO
	PRO	ALA	ARG	ARG
	THR	TYR	PHE	LEU
	GLY	VAL	GLY	GLY

## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	38978	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)	1200	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOCONTINUUM (6k x 4k)	Depositor
Maximum map value	0.707	Depositor
Minimum map value	-0.492	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.017	Depositor
Recommended contour level	0.082	Depositor
Map size (Å)	323.4, 323.4, 323.4	wwPDB
Map dimensions	392, 392, 392	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.825, 0.825, 0.825	Depositor

## 5 Model quality

### 5.1 Standard geometry

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

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	0.17	0/2976	0.42	0/4059
1	C	0.18	0/3132	0.44	0/4264
2	B	0.17	0/3015	0.43	0/4109
2	D	0.18	0/2964	0.48	0/4038
3	E	0.24	0/999	0.56	0/1364
3	F	0.25	0/1212	0.63	2/1648 (0.1%)
4	G	0.21	0/1185	0.53	0/1614
4	H	0.21	0/1177	0.51	1/1603 (0.1%)
5	M	0.30	0/284	0.84	0/385
All	All	0.19	0/16944	0.49	3/23084 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	125	ALA	CA-C-N	5.38	131.81	121.54
3	F	125	ALA	C-N-CA	5.38	131.81	121.54
4	H	139	LEU	CA-CB-CG	5.12	134.23	116.30

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	2910	0	2748	34	0
1	C	3063	0	2918	37	0
2	B	2950	0	2784	42	0
2	D	2902	0	2739	43	0
3	E	969	0	916	21	0
3	F	1177	0	1159	13	0
4	G	1163	0	1134	12	0
4	H	1156	0	1127	19	0
5	M	283	0	306	7	0
6	A	31	0	49	1	0
6	B	18	0	31	1	0
6	C	31	0	49	1	0
6	D	18	0	31	4	0
7	A	41	0	62	2	0
7	B	39	0	58	2	0
7	C	41	0	62	2	0
7	D	40	0	60	1	0
8	A	51	0	77	2	0
8	D	50	0	75	3	0
8	G	50	0	75	1	0
All	All	16983	0	16460	214	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 6.

All (214) 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)
2:D:713:GLU:HG2	2:D:721:MET:HA	1.69	0.74
3:E:13:THR:HG21	3:E:82:PHE:HE2	1.57	0.68
6:C:1001:PLM:H32	7:C:1002:OLC:H21	1.76	0.67
2:B:449:LYS:HD2	2:B:461:ASN:HB2	1.76	0.67
1:A:485:ILE:HD11	1:A:733:PRO:HA	1.77	0.66
4:H:88:ARG:HB3	4:H:102:LYS:HD3	1.79	0.64
1:C:638:GLN:HE22	1:C:641:ILE:HG13	1.62	0.64
2:B:520:PRO:O	2:B:619:ASN:ND2	2.31	0.64
1:A:510:VAL:HG21	8:A:1003:POV:H2	1.81	0.63
3:E:104:ARG:O	3:E:104:ARG:NH1	2.27	0.63
3:F:90:THR:O	3:F:94:ASN:ND2	2.31	0.63
1:C:644:GLY:HA2	1:C:670:MET:HE1	1.79	0.63
3:F:11:MET:HB3	3:F:153:MET:HE2	1.83	0.61
1:A:769:CYS:SG	1:A:770:GLY:N	2.73	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:475:ALA:HB3	2:D:735:ALA:HB3	1.82	0.60
2:B:488:VAL:HG13	2:B:489:ILE:HG12	1.83	0.60
4:H:129:ILE:HD12	4:H:132:ILE:HD11	1.84	0.60
2:D:405:TYR:HD2	2:D:478:PRO:HG3	1.67	0.60
4:H:154:LYS:O	4:H:229:ARG:NH2	2.36	0.58
2:D:433:ALA:HA	2:D:436:CYS:HB3	1.84	0.58
2:D:498:LEU:HD12	2:D:705:GLU:HB2	1.85	0.58
2:B:586:ARG:NH1	1:C:583:GLN:OE1	2.37	0.57
4:G:137:LEU:HD22	4:G:164:LEU:HD22	1.86	0.57
1:A:624:ARG:NH1	2:D:626:VAL:O	2.37	0.57
2:D:500:ILE:HB	2:D:727:LEU:HB2	1.87	0.56
2:B:619:ASN:HD21	1:C:782:ALA:HA	1.69	0.56
3:F:61:CYS:SG	3:F:65:ARG:NH1	2.78	0.56
1:C:402:VAL:HG22	1:C:422:VAL:HG23	1.87	0.56
1:A:595:ARG:HD2	2:B:578:TRP:CD2	2.41	0.56
2:D:520:PRO:O	2:D:619:ASN:ND2	2.38	0.55
2:B:481:ILE:O	2:B:730:LYS:NZ	2.38	0.55
3:E:87:GLU:OE1	3:E:90:THR:OG1	2.20	0.55
2:D:797:TYR:HB3	8:D:1003:POV:H28A	1.88	0.55
1:C:520:GLU:OE2	4:H:199:TYR:OH	2.26	0.54
2:D:424:TYR:O	2:D:428:LEU:HB2	2.07	0.54
1:A:516:PRO:O	1:A:615:ASN:ND2	2.39	0.54
2:B:626:VAL:O	1:C:624:ARG:NH1	2.40	0.54
4:G:214:LEU:HA	4:G:217:VAL:HG22	1.89	0.54
1:A:733:PRO:HG2	1:A:736:SER:HB2	1.90	0.54
2:D:531:PHE:HA	2:D:534:ILE:HG22	1.90	0.54
2:D:649:THR:HG22	2:D:703:LEU:HB2	1.90	0.54
3:F:21:ILE:O	3:F:25:ILE:HG12	2.08	0.54
2:B:646:ALA:HB3	2:B:700:TYR:HA	1.88	0.53
2:B:526:TRP:HB3	6:B:1001:PLM:H82	1.91	0.53
3:F:26:TRP:HD1	3:F:67:LEU:HD11	1.73	0.53
5:M:290:THR:HA	5:M:293:ILE:HG22	1.90	0.53
1:A:532:VAL:HG22	2:B:803:LEU:HD11	1.91	0.53
6:D:1001:PLM:H81	7:D:1002:OLC:H2	1.89	0.53
1:C:443:ASP:OD2	1:C:457:ASN:ND2	2.42	0.53
2:B:743:ARG:HH22	1:C:739:ARG:HH22	1.56	0.52
2:B:764:ASN:HA	2:B:768:TYR:HD2	1.74	0.52
4:G:89:ILE:HG23	4:G:103:ILE:HG12	1.92	0.52
4:H:44:TRP:HA	4:H:204:SER:HB3	1.90	0.52
5:M:270:ASN:HA	5:M:273:PHE:CE1	2.45	0.52
3:E:155:TYR:O	3:E:159:SER:CB	2.58	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:155:TYR:O	3:E:159:SER:HB2	2.10	0.51
2:B:480:THR:O	2:B:485:ARG:NH1	2.44	0.51
1:A:491:PHE:HB2	1:A:755:LEU:HD11	1.93	0.51
1:C:827:ILE:HG12	1:C:829:GLN:H	1.75	0.51
3:E:74:ILE:HA	3:E:77:LEU:HD12	1.93	0.51
1:C:471:ALA:HB3	1:C:731:ALA:HB3	1.92	0.51
2:D:521:LEU:HD22	2:D:525:ILE:HG21	1.93	0.50
2:D:467:LEU:HD11	2:D:475:ALA:HB2	1.93	0.50
4:H:136:ILE:HA	4:H:139:LEU:HD13	1.93	0.50
2:B:467:LEU:HD11	2:B:737:PRO:HD3	1.94	0.50
1:C:402:VAL:HG23	1:C:421:CYS:HB2	1.94	0.50
2:B:794:GLY:HA2	2:B:797:TYR:HD2	1.76	0.50
2:D:490:ASP:OD1	2:D:736:THR:OG1	2.30	0.49
4:G:170:LEU:O	4:G:174:ILE:HG12	2.12	0.49
2:D:445:VAL:HG22	2:D:447:ASP:H	1.78	0.49
2:D:526:TRP:HB3	6:D:1001:PLM:H92	1.94	0.49
2:D:453:ARG:HA	2:D:460:TRP:HD1	1.78	0.49
4:H:90:CYS:HA	4:H:100:CYS:HA	1.95	0.49
2:D:671:TRP:O	2:D:675:ARG:N	2.46	0.48
1:A:647:GLU:HA	1:A:678:PHE:HD2	1.78	0.48
7:B:1002:OLC:H12	8:G:501:POV:H31D	1.94	0.48
3:F:120:VAL:HB	3:F:122:ILE:HG12	1.94	0.48
2:B:714:GLN:HB3	2:B:771:GLY:HA3	1.95	0.48
6:A:1001:PLM:H72	7:A:1002:OLC:H4	1.95	0.48
1:C:644:GLY:HA3	1:C:677:VAL:HB	1.96	0.48
2:B:713:GLU:HG2	2:B:721:MET:HA	1.95	0.48
2:D:404:PRO:O	2:D:766:TRP:NE1	2.43	0.48
1:C:414:ASN:HA	1:C:437:ARG:HH22	1.79	0.48
3:E:13:THR:HG21	3:E:82:PHE:CE2	2.45	0.48
4:H:214:LEU:HA	4:H:217:VAL:HG12	1.94	0.48
2:D:481:ILE:HD11	2:D:494:PRO:HG3	1.95	0.47
1:A:751:GLU:OE2	2:D:482:THR:OG1	2.32	0.47
1:A:401:TYR:O	1:A:420:TYR:N	2.47	0.47
1:A:485:ILE:HA	1:A:734:LYS:HE3	1.96	0.47
2:D:801:GLY:HA2	8:D:1003:POV:H213	1.96	0.47
2:B:432:ILE:O	2:B:435:HIS:ND1	2.46	0.47
2:D:423:GLY:H	2:D:426:VAL:HB	1.79	0.47
5:M:266:ARG:O	5:M:270:ASN:OD1	2.32	0.47
4:G:145:VAL:O	4:G:148:SER:OG	2.31	0.47
4:H:129:ILE:O	4:H:133:LEU:HB2	2.14	0.47
2:B:761:LYS:HA	1:C:660:ILE:HD11	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:786:SER:HA	1:A:789:ALA:HB2	1.96	0.47
2:D:542:LEU:HD22	3:E:23:PHE:HZ	1.80	0.46
3:E:13:THR:O	3:E:17:CYS:HB2	2.14	0.46
3:F:90:THR:HG22	3:F:94:ASN:HD21	1.80	0.46
3:F:91:LEU:O	3:F:95:ILE:HG12	2.16	0.46
2:B:447:ASP:HB3	2:B:449:LYS:HG3	1.95	0.46
2:B:521:LEU:HD13	2:B:616:TYR:HD1	1.79	0.46
1:C:714:CYS:N	1:C:769:CYS:SG	2.89	0.46
2:B:475:ALA:HB3	2:B:735:ALA:HB3	1.98	0.46
1:C:425:ALA:HA	1:C:428:ILE:HG22	1.98	0.46
1:A:502:LYS:H	1:A:716:THR:HA	1.81	0.46
3:E:10:TYR:HB3	3:E:153:MET:HB3	1.97	0.46
1:A:629:ILE:HG21	1:A:635:LEU:HD13	1.97	0.46
2:B:800:VAL:HG11	4:G:174:ILE:HD11	1.98	0.46
1:C:496:ILE:HD12	1:C:651:THR:HG22	1.96	0.46
1:C:709:GLU:HG3	1:C:717:MET:HA	1.98	0.46
1:C:460:VAL:HA	1:C:463:LEU:HD12	1.97	0.45
1:C:488:SER:OG	1:C:489:LYS:N	2.49	0.45
3:E:91:LEU:O	3:E:95:ILE:HG12	2.16	0.45
4:H:45:LEU:HA	4:H:202:GLY:H	1.80	0.45
1:C:700:LEU:HD13	1:C:704:MET:HB3	1.97	0.45
1:A:785:LEU:HD22	2:D:524:GLU:HG2	1.98	0.45
1:A:799:LEU:HD22	2:D:539:VAL:HG21	1.97	0.45
2:D:458:LYS:O	2:D:469:TYR:OH	2.33	0.45
5:M:270:ASN:O	5:M:274:ILE:HG12	2.17	0.45
4:H:83:HIS:CE1	4:H:90:CYS:HB3	2.52	0.45
1:A:746:VAL:O	1:A:750:SER:N	2.50	0.45
1:A:711:ARG:HA	1:A:768:GLU:H	1.82	0.45
4:H:103:ILE:HG22	4:H:105:HIS:HD2	1.80	0.45
1:A:402:VAL:HG22	1:A:422:VAL:HG23	1.98	0.45
2:B:445:VAL:HG22	2:B:447:ASP:H	1.82	0.45
2:D:542:LEU:HD21	3:E:74:ILE:HD11	1.99	0.45
3:E:157:LEU:HD23	3:E:157:LEU:HA	1.84	0.45
4:H:140:LEU:HB3	4:H:164:LEU:HD21	2.00	0.44
4:G:42:ASP:HA	4:G:85:GLY:HA3	1.98	0.44
1:C:459:MET:SD	1:C:459:MET:N	2.89	0.44
1:C:645:THR:OG1	1:C:646:LEU:N	2.51	0.44
2:B:685:THR:OG1	2:B:688:GLU:OE1	2.29	0.44
2:B:648:GLY:HA2	2:B:681:VAL:HG22	2.00	0.44
1:A:607:ILE:HG21	2:B:795:VAL:HG21	2.00	0.43
1:A:624:ARG:NE	1:A:781:SER:OG	2.42	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:E:25:ILE:HA	3:E:28:ILE:HG12	1.99	0.43
4:G:210:LEU:HA	4:G:213:ILE:HG22	1.98	0.43
2:B:606:TRP:CD1	1:C:583:GLN:HG3	2.54	0.43
1:C:632:ALA:O	1:C:636:ALA:N	2.45	0.43
3:E:155:TYR:O	3:E:159:SER:OG	2.36	0.43
4:G:217:VAL:HA	4:G:220:VAL:HG12	2.01	0.43
2:B:600:ILE:HA	7:C:1002:OLC:H14	2.01	0.43
2:D:758:LEU:HD13	2:D:761:LYS:HD2	2.00	0.43
1:A:665:LYS:O	1:A:668:THR:OG1	2.23	0.43
1:A:734:LYS:HA	1:A:739:ARG:HH22	1.83	0.43
1:A:758:LEU:HA	1:A:761:LYS:HB3	2.01	0.43
1:C:429:ALA:HB2	1:C:436:TYR:HE2	1.83	0.43
2:D:460:TRP:HE3	2:D:465:GLY:HA2	1.83	0.43
1:A:764:TYR:O	1:A:771:SER:OG	2.25	0.43
1:A:794:ILE:HA	8:A:1003:POV:H38	2.01	0.43
2:D:400:ILE:HG23	2:D:402:GLU:H	1.84	0.43
2:D:664:ILE:HG23	2:D:666:VAL:H	1.83	0.43
4:G:83:HIS:NE2	4:G:90:CYS:HB3	2.34	0.43
1:A:535:VAL:HG21	2:B:803:LEU:HD23	2.00	0.43
2:B:702:TYR:CE2	2:B:704:LEU:HD23	2.54	0.43
2:B:707:THR:O	2:B:710:GLU:HG3	2.18	0.43
2:D:523:TYR:CD1	6:D:1001:PLM:H31	2.53	0.42
3:F:64:LEU:O	3:F:68:VAL:HG23	2.19	0.42
4:G:34:LEU:HB3	4:G:211:SER:HB2	2.01	0.42
2:D:423:GLY:HA2	2:D:766:TRP:CZ2	2.55	0.42
4:G:119:TYR:HA	4:G:122:ARG:HE	1.84	0.42
1:C:392:ILE:HG22	1:C:437:ARG:HB3	2.00	0.42
5:M:273:PHE:HA	5:M:276:LEU:HG	2.00	0.42
2:B:760:ASP:OD1	2:B:761:LYS:N	2.52	0.42
1:C:645:THR:HB	1:C:699:LEU:HB2	2.01	0.42
1:A:698:TYR:HD2	1:A:700:LEU:HD22	1.85	0.42
2:B:428:LEU:O	2:B:432:ILE:HG12	2.20	0.42
3:E:134:GLU:O	3:E:138:LYS:HG2	2.20	0.42
5:M:263:ILE:HG22	5:M:266:ARG:HH21	1.84	0.42
1:C:660:ILE:HB	1:C:663:PHE:HD1	1.85	0.42
2:D:397:VAL:HG12	2:D:474:ILE:HG22	2.02	0.42
2:D:523:TYR:HD1	6:D:1001:PLM:H31	1.85	0.42
2:D:721:MET:HE3	2:D:721:MET:HB2	1.83	0.42
2:B:713:GLU:OE2	2:B:722:LYS:N	2.50	0.42
2:D:538:VAL:O	2:D:542:LEU:HG	2.20	0.42
5:M:286:CYS:HA	5:M:289:LEU:HG	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:459:MET:HA	1:C:462:GLU:HG2	2.02	0.41
2:D:496:MET:HE2	2:D:496:MET:HB3	1.93	0.41
3:E:24:VAL:HG12	3:E:139:LEU:HD12	2.01	0.41
3:E:90:THR:O	3:E:94:ASN:ND2	2.52	0.41
3:E:126:ASP:N	3:E:126:ASP:OD1	2.50	0.41
3:F:7:ALA:O	3:F:11:MET:HG2	2.21	0.41
1:C:397:LEU:HB2	1:C:440:ILE:HD13	2.02	0.41
1:C:771:SER:OG	1:C:772:LYS:N	2.54	0.41
3:F:11:MET:HA	3:F:14:LEU:HG	2.03	0.41
2:B:442:LEU:HD23	2:B:442:LEU:HA	1.92	0.41
1:C:437:ARG:HH21	1:C:438:LEU:HB2	1.85	0.41
1:C:485:ILE:HG12	1:C:731:ALA:HB1	2.03	0.41
2:D:808:LEU:O	2:D:812:ILE:HG12	2.20	0.41
8:D:1003:POV:H22A	4:H:121:LEU:HA	2.02	0.41
3:E:18:ALA:HB2	3:E:146:PHE:HE1	1.86	0.41
3:F:17:CYS:HA	3:F:20:LEU:HD12	2.02	0.41
1:A:502:LYS:HE3	1:A:502:LYS:HB3	1.89	0.41
1:A:595:ARG:HA	1:A:595:ARG:HD3	1.89	0.41
2:B:401:LEU:HD23	2:B:406:VAL:HG23	2.03	0.41
3:F:75:HIS:CE1	3:F:97:LEU:HD23	2.55	0.41
1:A:395:THR:HG21	1:A:402:VAL:HG21	2.03	0.41
2:B:651:ASP:HA	2:B:682:PHE:HD2	1.86	0.41
2:D:607:PHE:O	2:D:611:ILE:HG12	2.21	0.41
4:H:140:LEU:HA	4:H:143:VAL:HG22	2.03	0.41
1:A:492:MET:HB3	1:A:728:TYR:HB2	2.03	0.41
1:C:538:LEU:HD21	4:H:163:ILE:HG12	2.03	0.41
1:C:666:MET:O	1:C:670:MET:HG2	2.21	0.41
3:E:21:ILE:HG12	3:E:143:LEU:HD12	2.02	0.41
2:B:518:LEU:HD11	7:B:1002:OLC:H2	2.01	0.40
7:A:1002:OLC:H14A	7:A:1004:OLC:H13	2.03	0.40
2:B:405:TYR:O	2:B:424:TYR:N	2.54	0.40
2:B:495:PHE:CD2	2:B:496:MET:HG3	2.56	0.40
4:H:160:GLY:O	4:H:164:LEU:HB2	2.21	0.40
2:D:792:VAL:O	2:D:795:VAL:HG12	2.22	0.40
4:H:45:LEU:HD22	4:H:179:TYR:OH	2.21	0.40
4:H:88:ARG:HB3	4:H:102:LYS:HA	2.03	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles

### 5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	399/500 (80%)	376 (94%)	23 (6%)	0	100	100
1	C	417/500 (83%)	398 (95%)	19 (5%)	0	100	100
2	B	403/467 (86%)	375 (93%)	28 (7%)	0	100	100
2	D	399/467 (85%)	374 (94%)	25 (6%)	0	100	100
3	E	118/160 (74%)	118 (100%)	0	0	100	100
3	F	140/160 (88%)	131 (94%)	9 (6%)	0	100	100
4	G	154/423 (36%)	149 (97%)	5 (3%)	0	100	100
4	H	154/423 (36%)	146 (95%)	8 (5%)	0	100	100
5	M	37/306 (12%)	37 (100%)	0	0	100	100
All	All	2221/3406 (65%)	2104 (95%)	117 (5%)	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	277/420 (66%)	277 (100%)	0	100	100
1	C	295/420 (70%)	295 (100%)	0	100	100
2	B	281/394 (71%)	281 (100%)	0	100	100
2	D	274/394 (70%)	274 (100%)	0	100	100
3	E	94/143 (66%)	94 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
3	F	122/143 (85%)	122 (100%)	0	100	100
4	G	110/310 (36%)	110 (100%)	0	100	100
4	H	109/310 (35%)	109 (100%)	0	100	100
5	M	30/236 (13%)	30 (100%)	0	100	100
All	All	1592/2770 (58%)	1592 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	619	ASN
2	B	756	GLN
2	B	764	ASN
1	C	615	ASN
1	C	638	GLN
1	C	705	ASN
2	D	435	HIS
2	D	726	ASN
3	F	75	HIS
3	F	94	ASN
3	F	107	HIS
4	G	105	HIS
4	H	224	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

17 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	$\# Z  > 2$
7	OLC	B	1002	-	24,24,24	0.24	0	25,25,25	0.23	0
8	POV	D	1003	-	49,49,51	0.31	0	55,57,59	0.31	0
7	OLC	D	1002	-	24,24,24	0.24	0	25,25,25	0.27	0
7	OLC	C	1002	-	24,24,24	0.25	0	25,25,25	0.33	0
7	OLC	C	1003	-	15,15,24	0.33	0	15,15,25	0.31	0
6	PLM	C	1004	-	12,12,17	0.63	0	12,12,17	0.59	0
6	PLM	A	1001	-	17,17,17	0.54	0	17,17,17	0.51	0
8	POV	G	501	-	49,49,51	0.29	0	55,57,59	0.47	0
8	POV	A	1003	-	50,50,51	0.30	0	56,58,59	0.34	0
7	OLC	A	1002	-	24,24,24	0.24	0	25,25,25	0.34	0
6	PLM	B	1001	-	17,17,17	0.54	0	17,17,17	0.50	0
7	OLC	D	1004	-	14,14,24	0.34	0	14,14,25	0.32	0
6	PLM	C	1001	-	17,17,17	0.54	0	17,17,17	0.50	0
7	OLC	A	1004	-	15,15,24	0.34	0	15,15,25	0.31	0
6	PLM	D	1001	-	17,17,17	0.53	0	17,17,17	0.51	0
7	OLC	B	1003	-	13,13,24	0.34	0	13,13,25	0.32	0
6	PLM	A	1005	-	12,12,17	0.63	0	12,12,17	0.68	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
7	OLC	B	1002	-	-	0/24/24/24	-
8	POV	D	1003	-	-	3/53/53/55	-
7	OLC	D	1002	-	-	2/24/24/24	-
7	OLC	C	1002	-	-	2/24/24/24	-
7	OLC	C	1003	-	-	3/13/13/24	-
6	PLM	C	1004	-	-	2/10/10/15	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	PLM	A	1001	-	-	3/15/15/15	-
8	POV	G	501	-	-	11/53/53/55	-
8	POV	A	1003	-	-	6/54/54/55	-
7	OLC	A	1002	-	-	2/24/24/24	-
6	PLM	B	1001	-	-	1/15/15/15	-
7	OLC	D	1004	-	-	1/12/12/24	-
6	PLM	C	1001	-	-	4/15/15/15	-
7	OLC	A	1004	-	-	0/13/13/24	-
6	PLM	D	1001	-	-	4/15/15/15	-
7	OLC	B	1003	-	-	1/11/11/24	-
6	PLM	A	1005	-	-	2/10/10/15	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (47) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	1003	POV	C11-O12-P-O14
8	A	1003	POV	C32-C31-O31-C3
8	A	1003	POV	O32-C31-O31-C3
8	G	501	POV	C1-O11-P-O13
8	G	501	POV	C32-C31-O31-C3
8	G	501	POV	O32-C31-O31-C3
8	G	501	POV	C2-C1-O11-P
6	A	1001	PLM	CB-CC-CD-CE
8	G	501	POV	O11-C1-C2-C3
6	C	1001	PLM	C6-C7-C8-C9
8	A	1003	POV	C24-C25-C26-C27
8	G	501	POV	C24-C25-C26-C27
8	G	501	POV	O11-C1-C2-O21
8	D	1003	POV	C12-C11-O12-P
8	A	1003	POV	O12-C11-C12-N
8	D	1003	POV	C11-O12-P-O14
8	G	501	POV	C11-O12-P-O14
8	G	501	POV	C29-C210-C211-C212
7	C	1003	OLC	C7-C8-C9-C10
8	D	1003	POV	C29-C210-C211-C212
6	C	1001	PLM	C5-C6-C7-C8

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Mol	Chain	Res	Type	Atoms
7	C	1002	OLC	C9-C10-C11-C12
7	A	1002	OLC	C7-C8-C9-C10
7	B	1003	OLC	C7-C8-C9-C10
7	D	1002	OLC	C7-C8-C9-C10
6	A	1005	PLM	O2-C1-C2-C3
6	A	1005	PLM	O1-C1-C2-C3
7	C	1002	OLC	C7-C8-C9-C10
6	C	1004	PLM	O2-C1-C2-C3
6	D	1001	PLM	C2-C3-C4-C5
6	C	1004	PLM	O1-C1-C2-C3
8	G	501	POV	C27-C28-C29-C210
8	A	1003	POV	C39-C310-C311-C312
6	C	1001	PLM	O1-C1-C2-C3
6	D	1001	PLM	O1-C1-C2-C3
7	D	1004	OLC	C7-C8-C9-C10
7	C	1003	OLC	O20-C1-C2-C3
6	D	1001	PLM	O2-C1-C2-C3
7	C	1003	OLC	O19-C1-C2-C3
6	C	1001	PLM	O2-C1-C2-C3
6	A	1001	PLM	C6-C7-C8-C9
7	D	1002	OLC	C9-C10-C11-C12
7	A	1002	OLC	C5-C6-C7-C8
6	A	1001	PLM	C5-C6-C7-C8
6	B	1001	PLM	C5-C6-C7-C8
8	G	501	POV	O31-C31-C32-C33
6	D	1001	PLM	C4-C5-C6-C7

There are no ring outliers.

12 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	1002	OLC	2	0
8	D	1003	POV	3	0
7	D	1002	OLC	1	0
7	C	1002	OLC	2	0
6	A	1001	PLM	1	0
8	G	501	POV	1	0
8	A	1003	POV	2	0
7	A	1002	OLC	2	0
6	B	1001	PLM	1	0
6	C	1001	PLM	1	0
7	A	1004	OLC	1	0

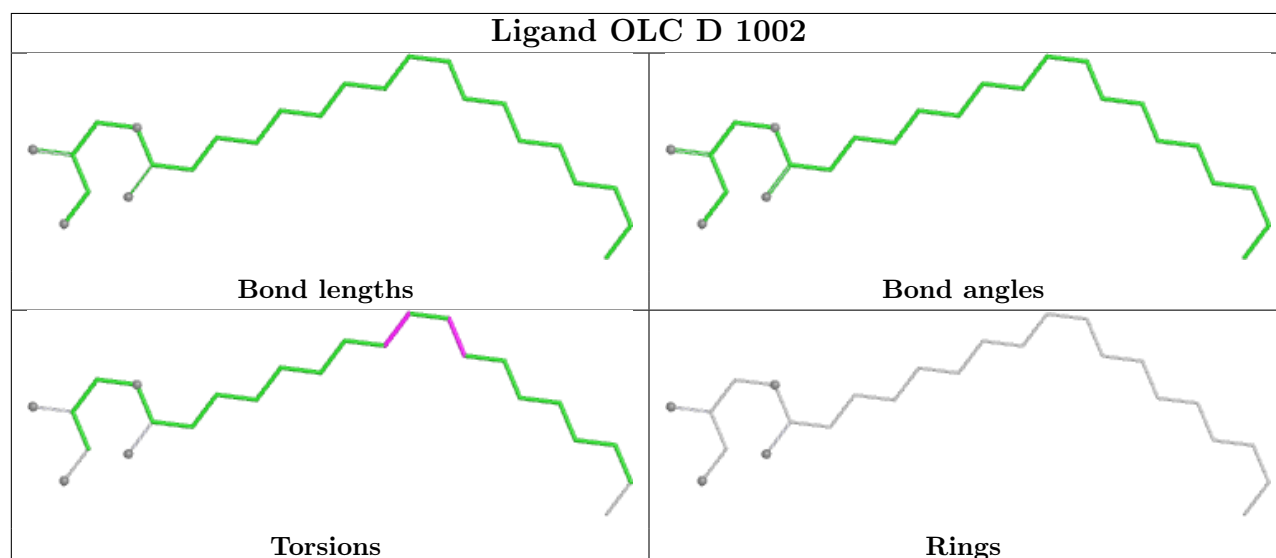
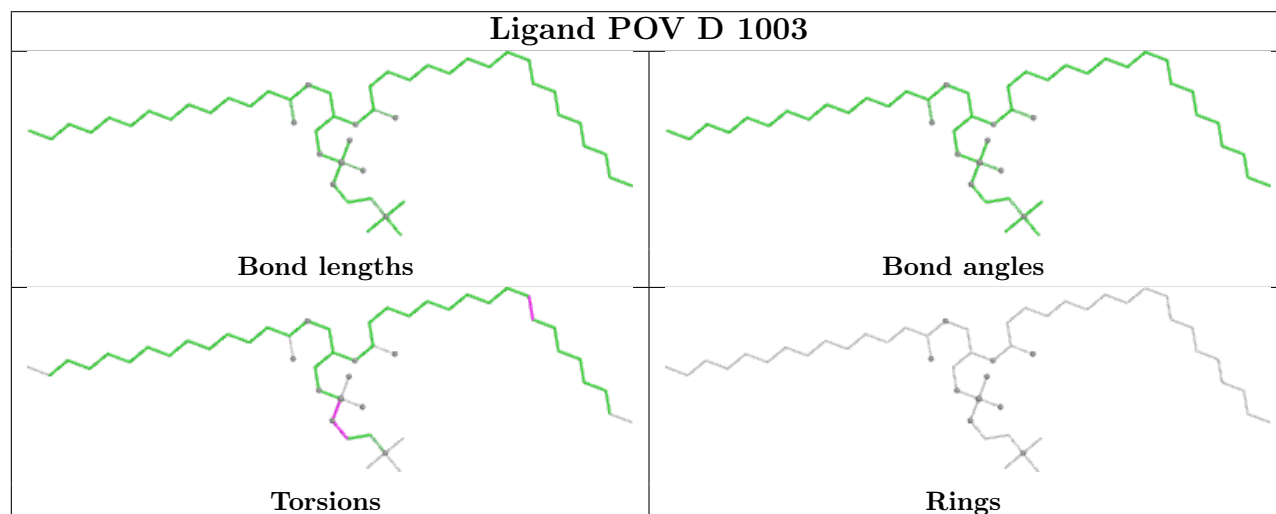
*Continued on next page...*

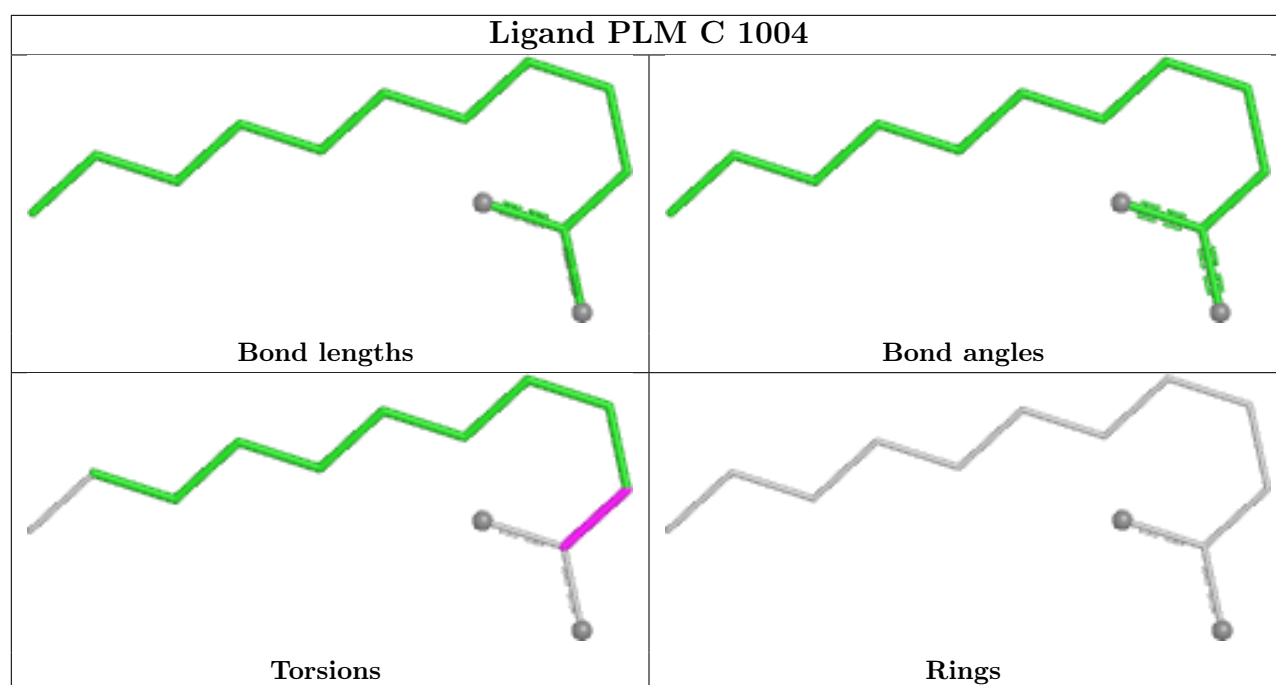
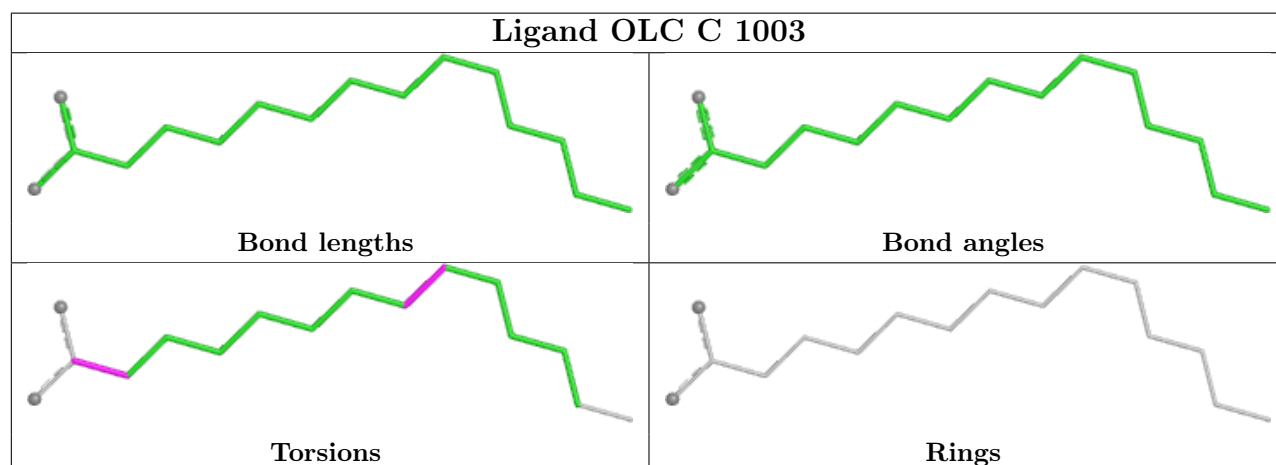
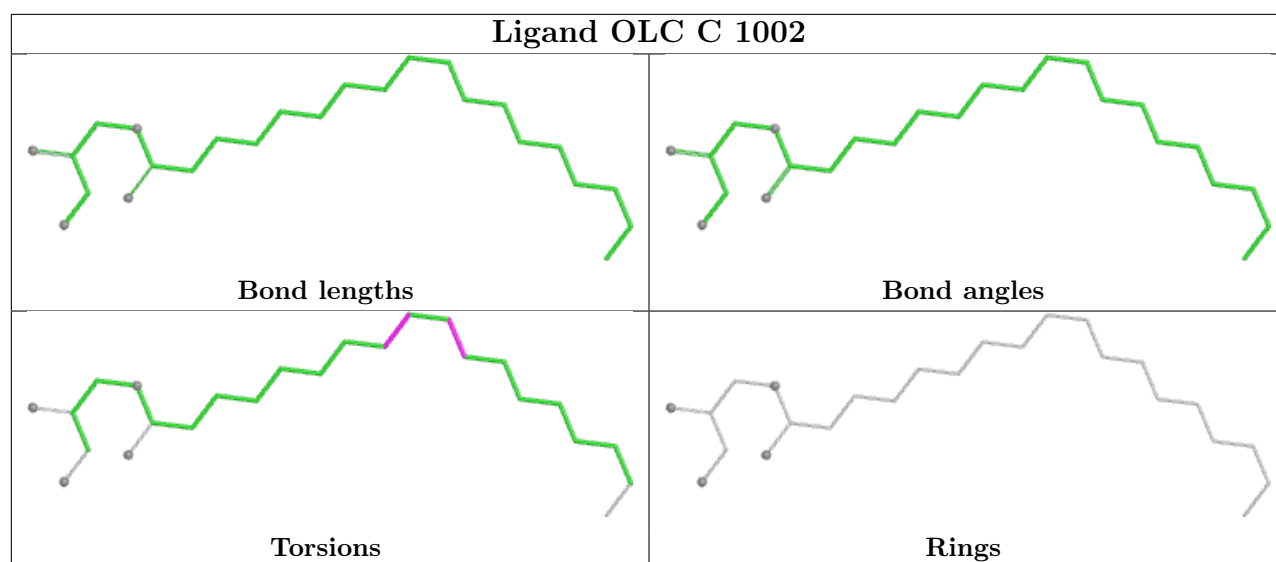


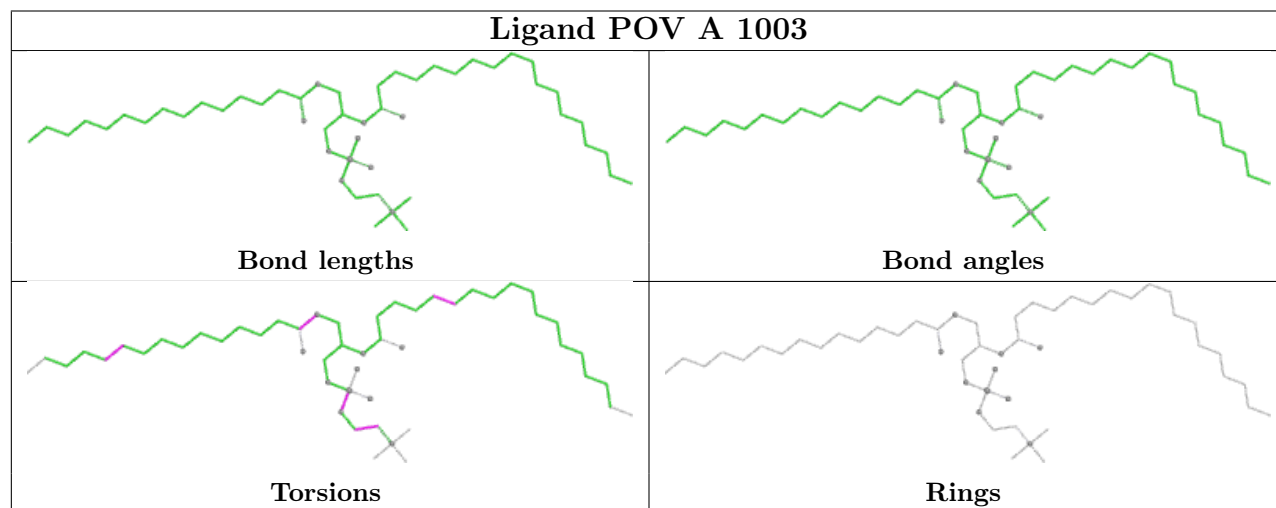
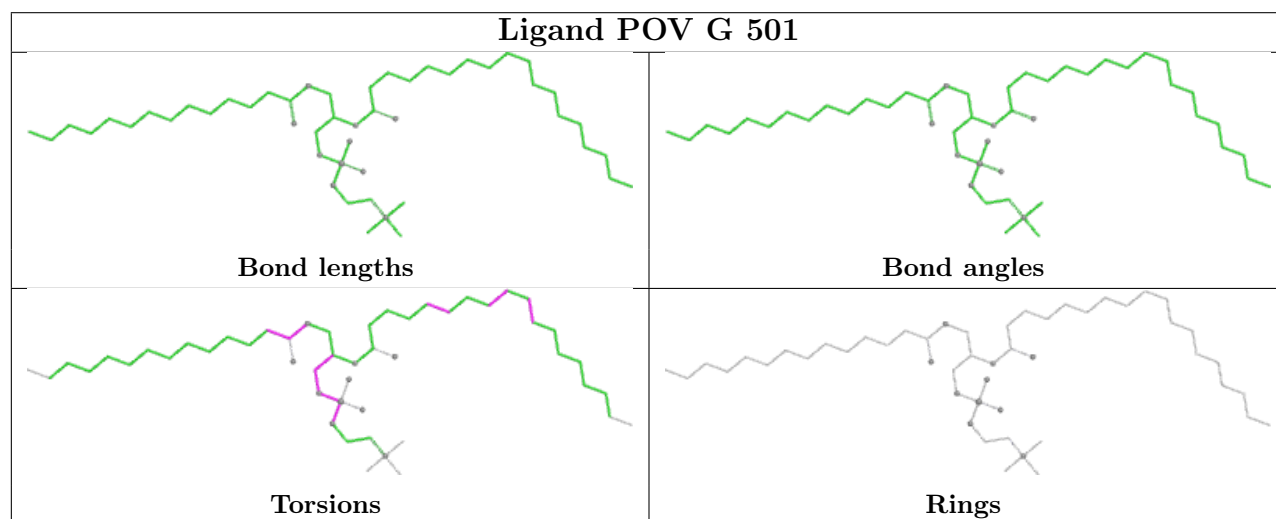
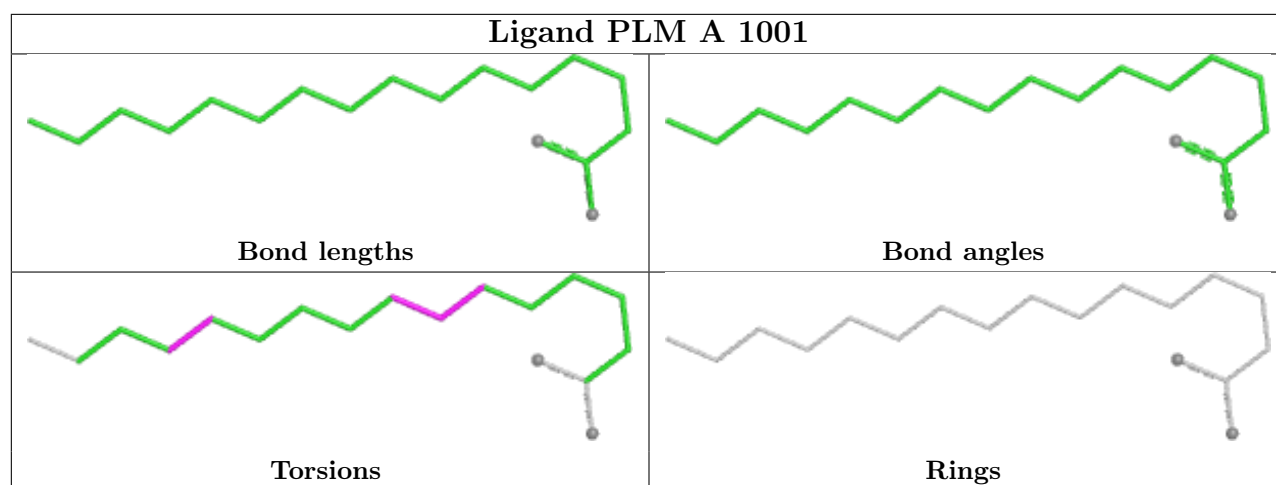
*Continued from previous page...*

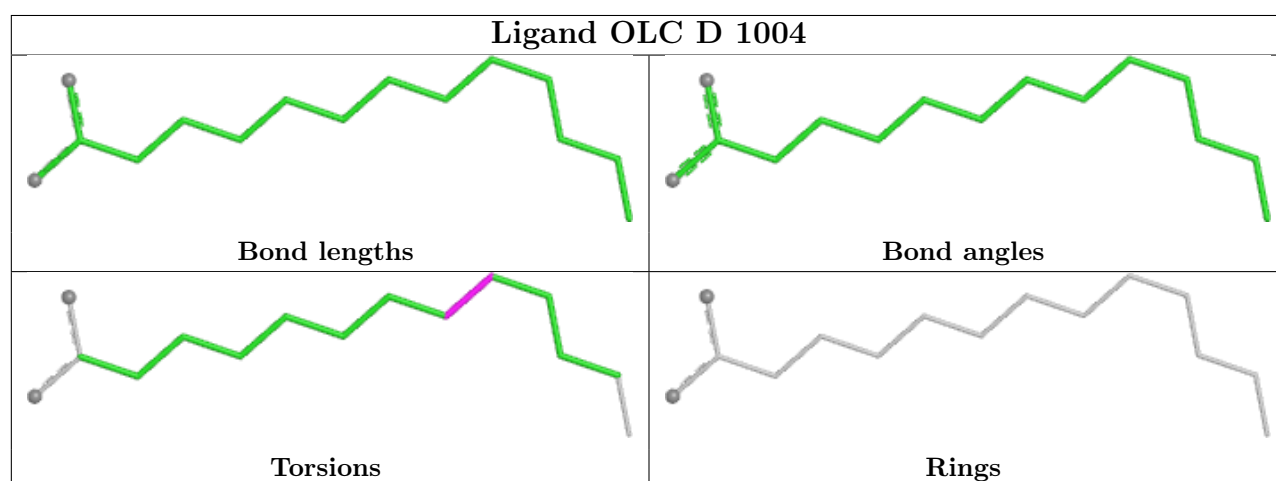
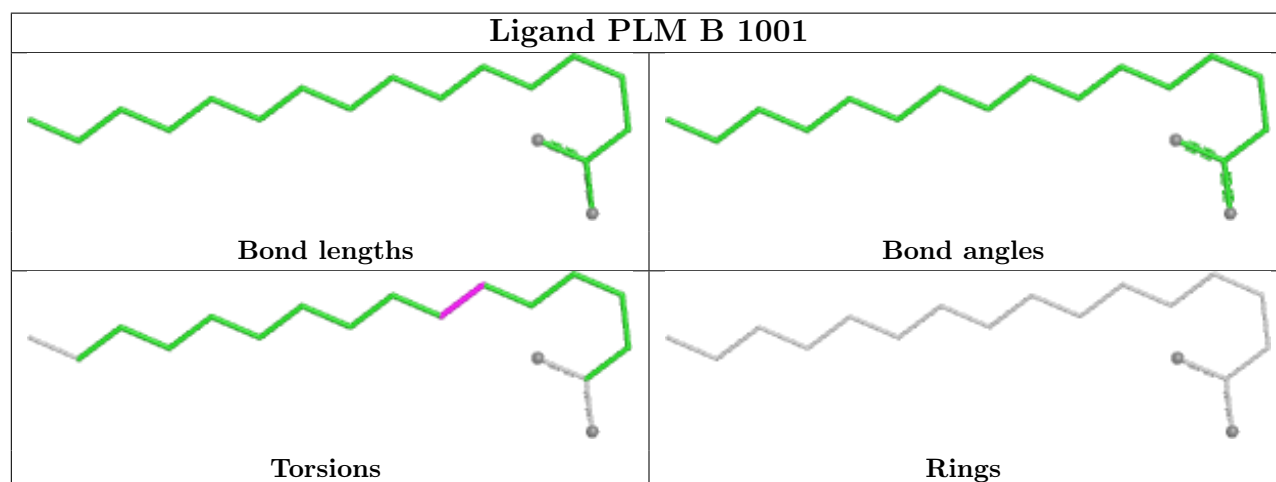
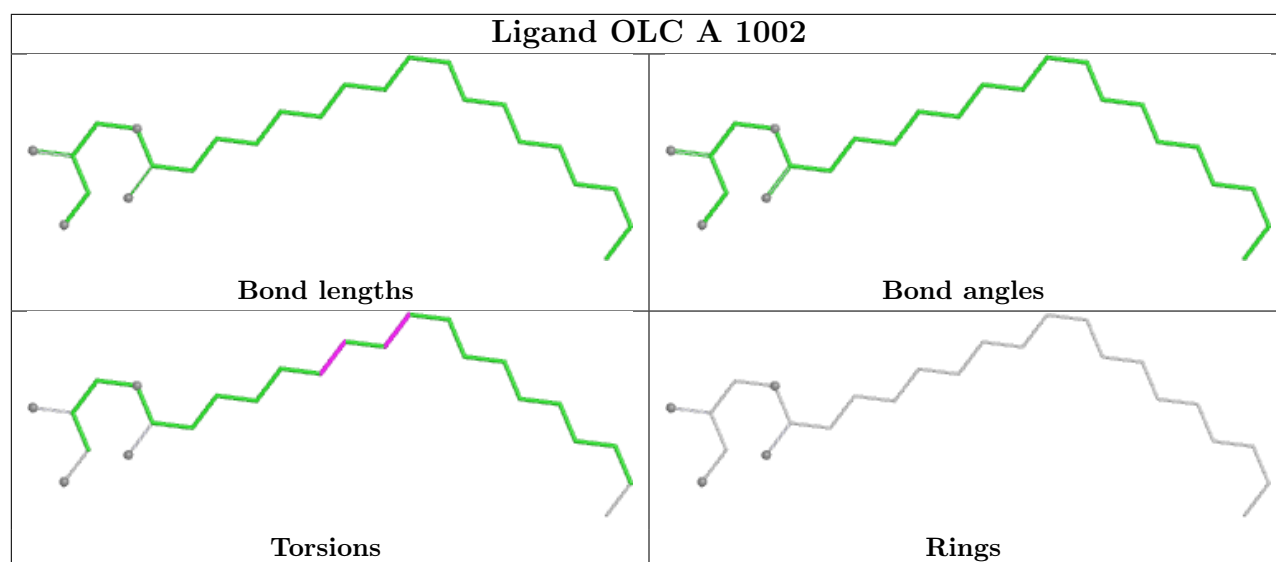
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	D	1001	PLM	4	0

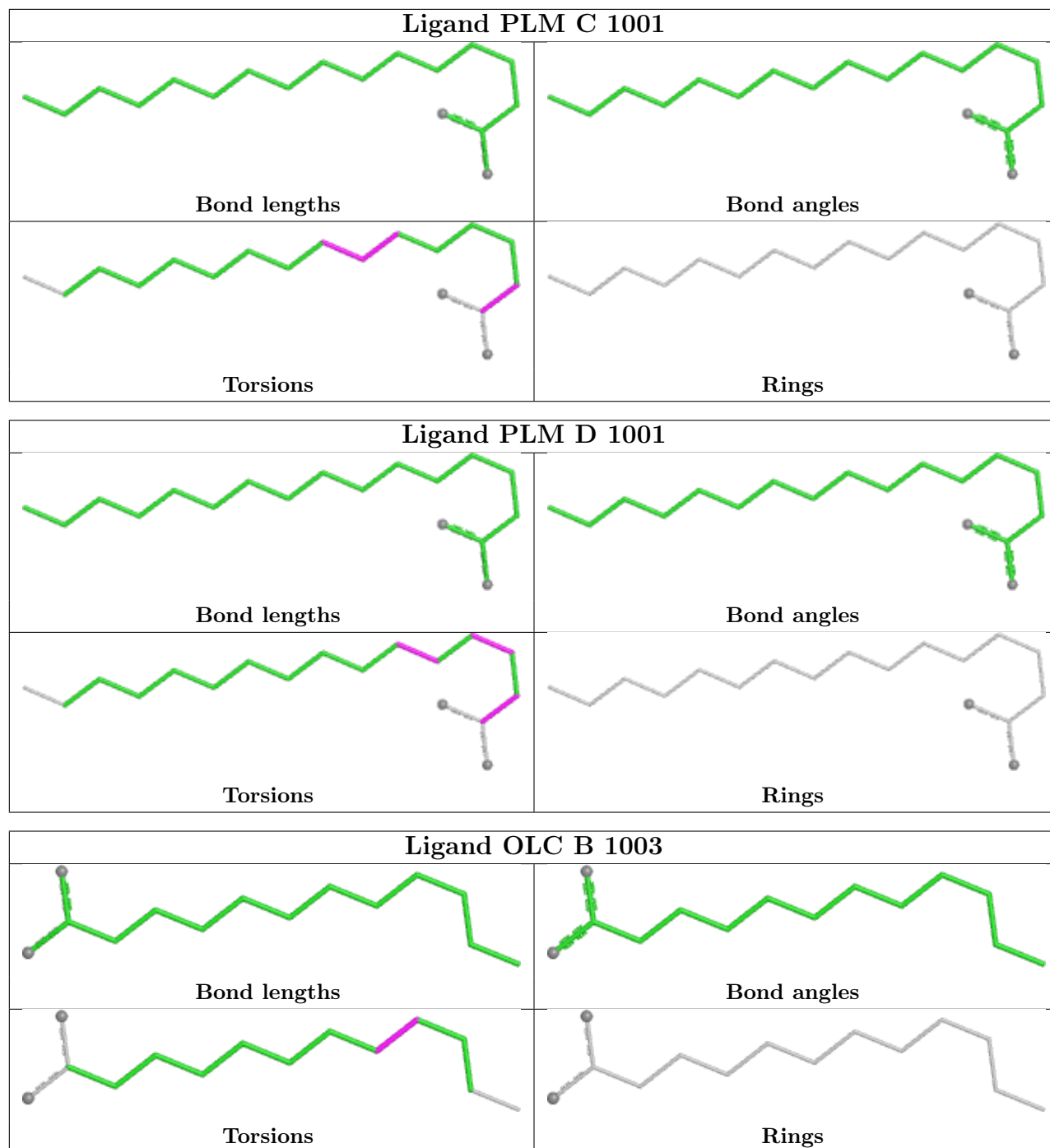
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

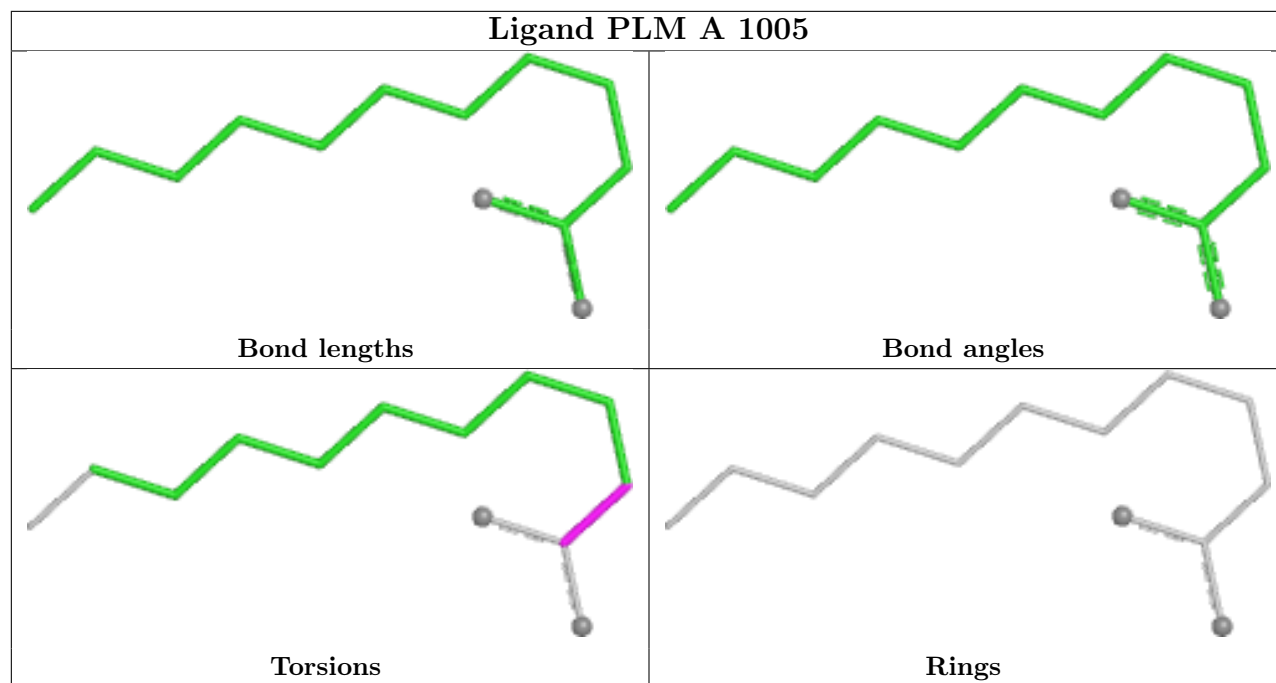












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

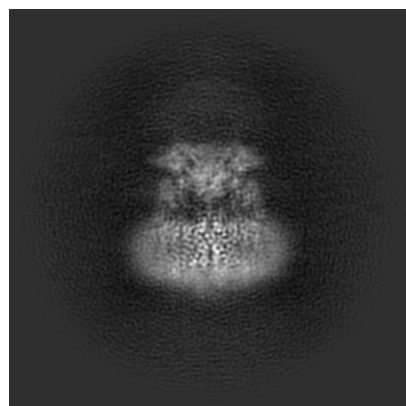
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-74383. These allow visual inspection of the internal detail of the map and identification of artifacts.

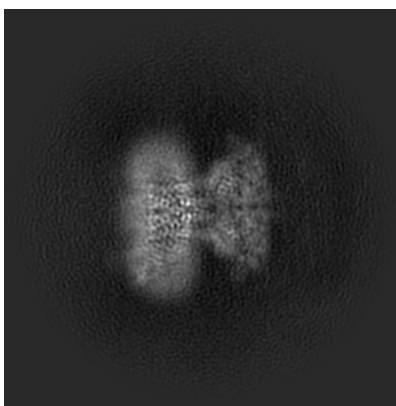
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

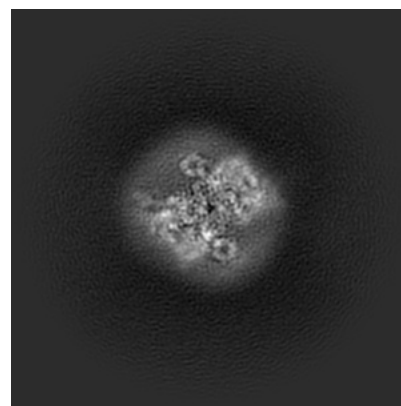
#### 6.1.1 Primary map



X

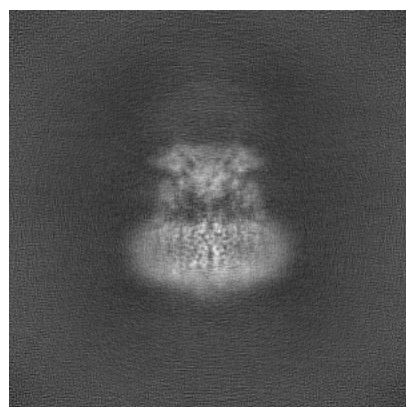


Y

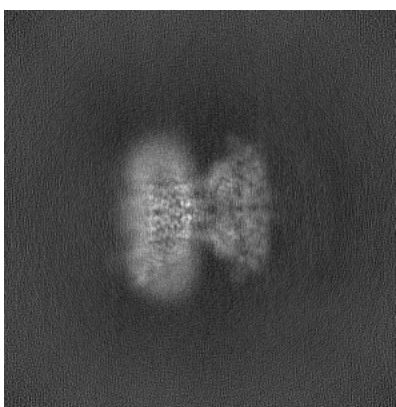


Z

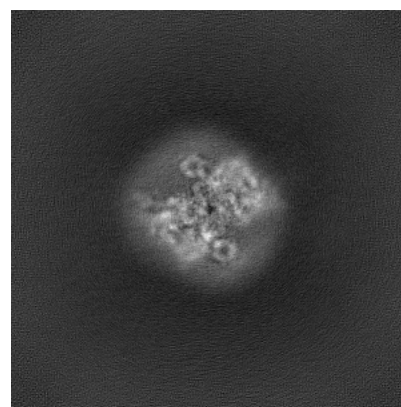
#### 6.1.2 Raw map



X



Y

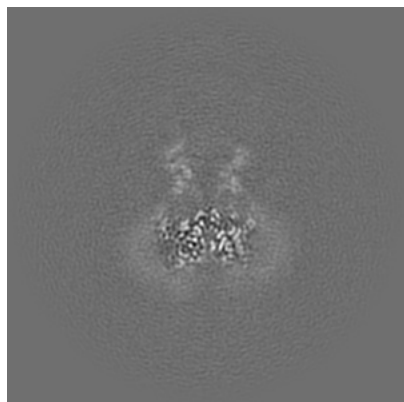


Z

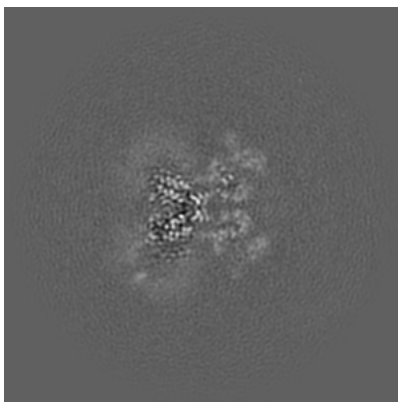
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

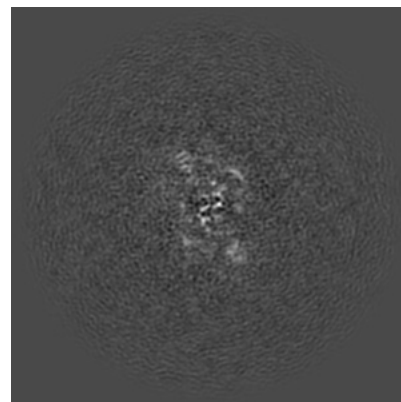
### 6.2.1 Primary map



X Index: 196

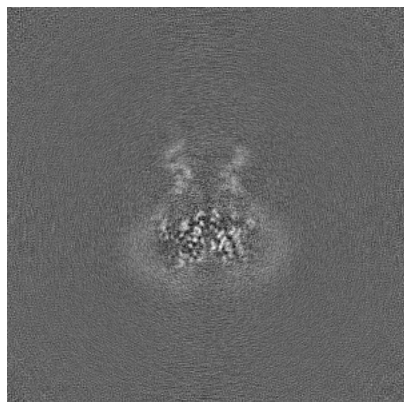


Y Index: 196

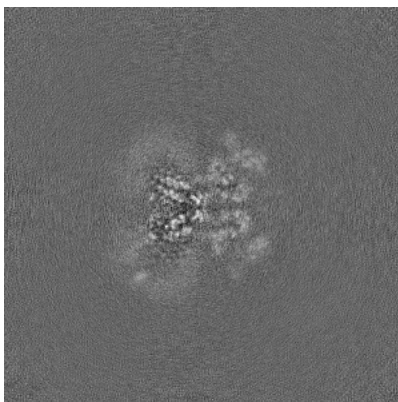


Z Index: 196

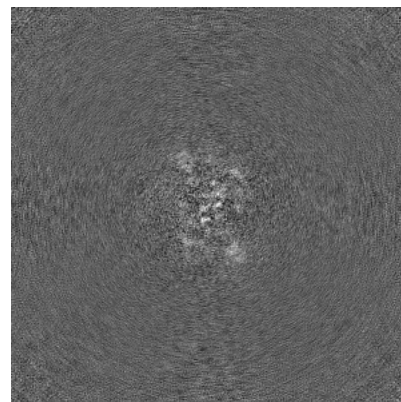
### 6.2.2 Raw map



X Index: 196



Y Index: 196



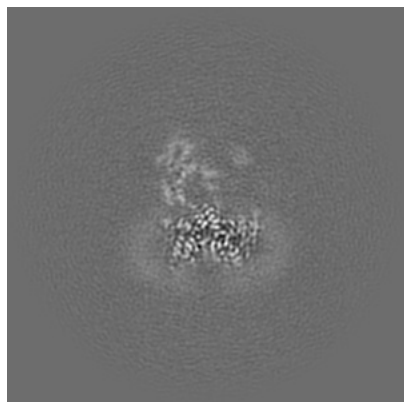
Z Index: 196

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

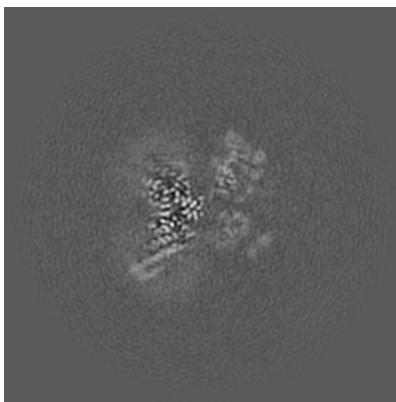


## 6.3 Largest variance slices [i](#)

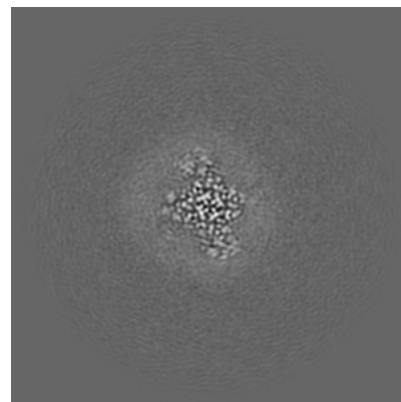
### 6.3.1 Primary map



X Index: 189

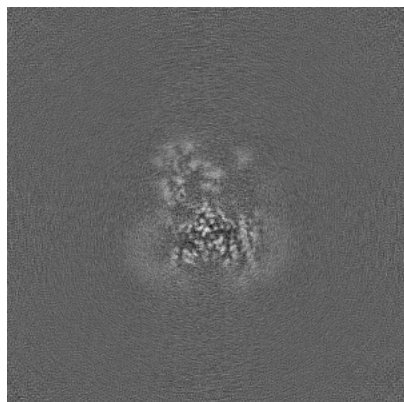


Y Index: 201

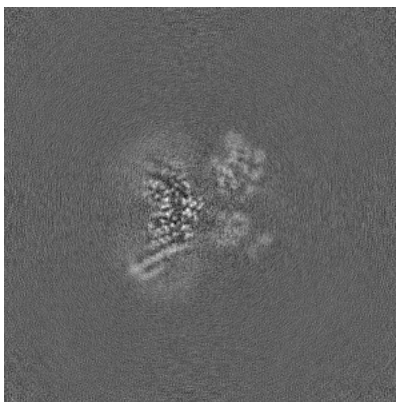


Z Index: 175

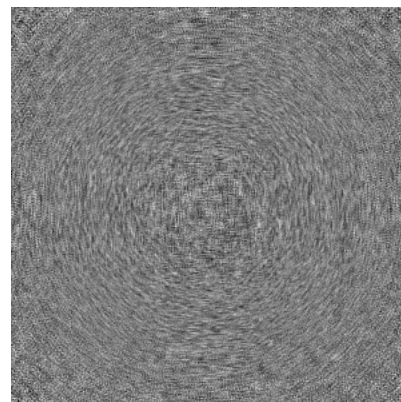
### 6.3.2 Raw map



X Index: 186



Y Index: 202

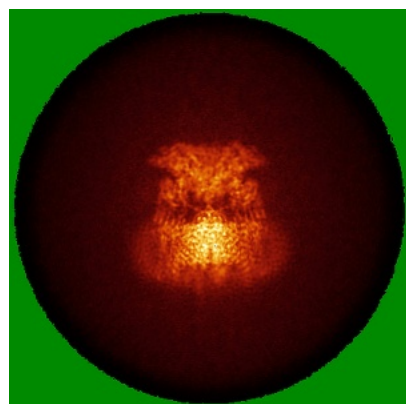


Z Index: 0

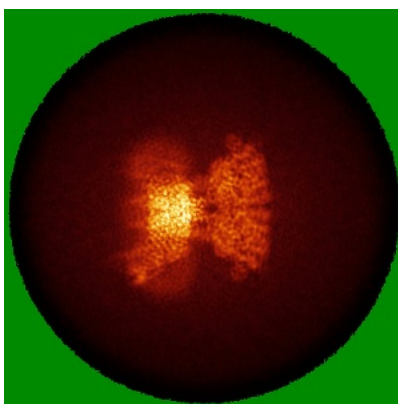
The images above show the largest variance slices of the map in three orthogonal directions.

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

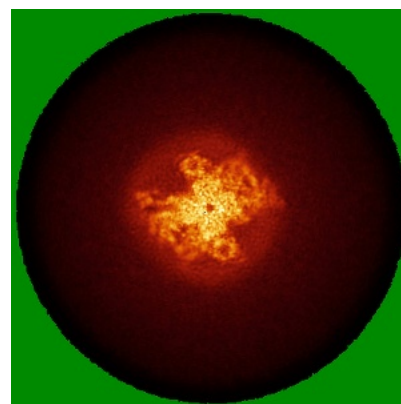
### 6.4.1 Primary map



X

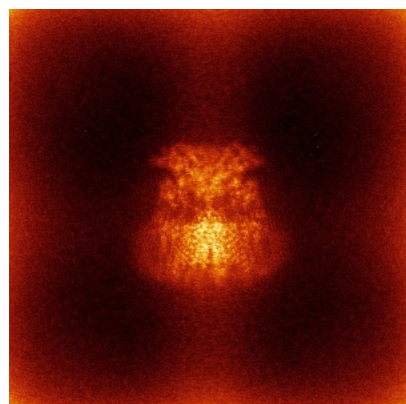


Y

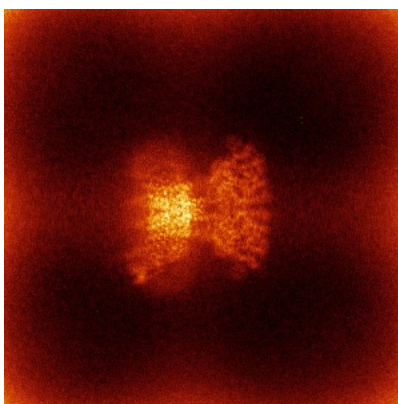


Z

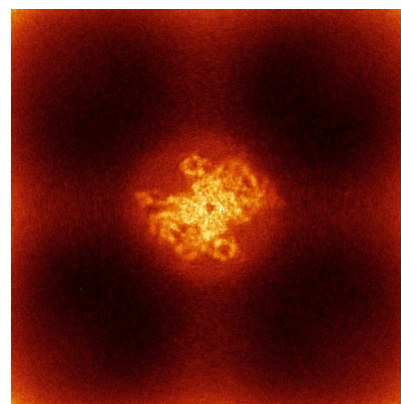
### 6.4.2 Raw map



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



X



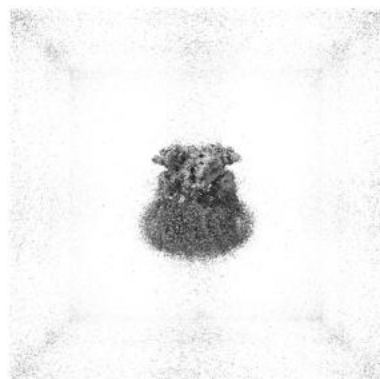
Y



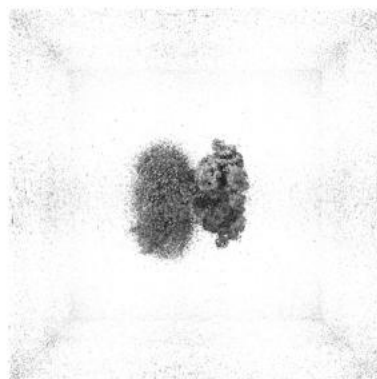
Z

The images above show the 3D surface view of the map at the recommended contour level 0.082. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

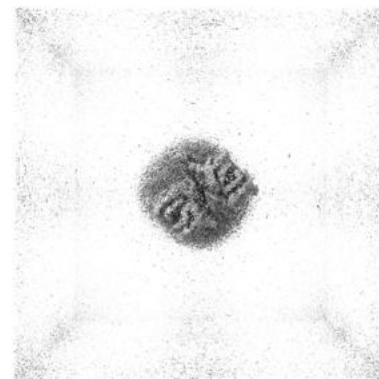
### 6.5.2 Raw map



X



Y



Z

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.

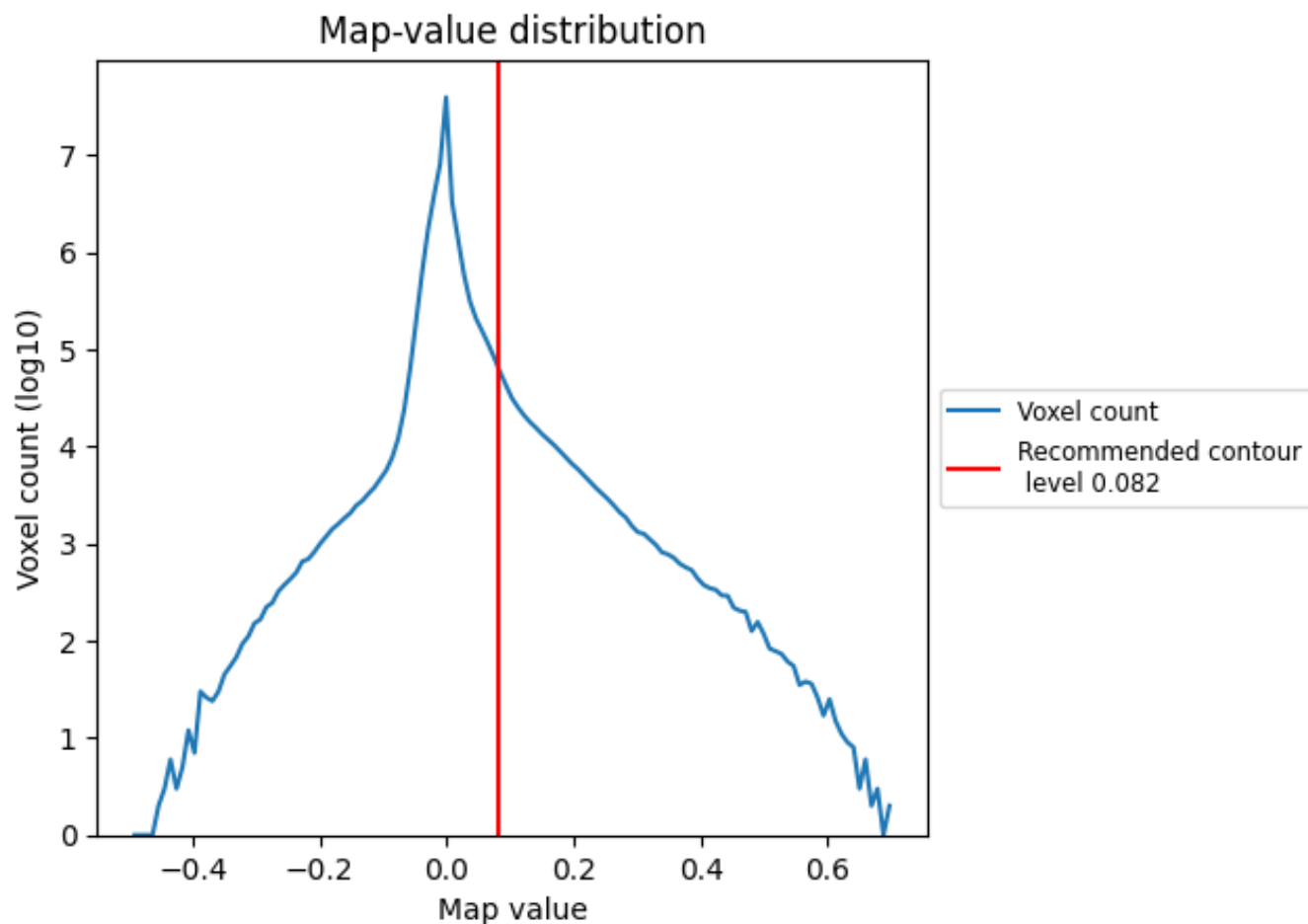
## 6.6 Mask visualisation [i](#)

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

## 7 Map analysis [i](#)

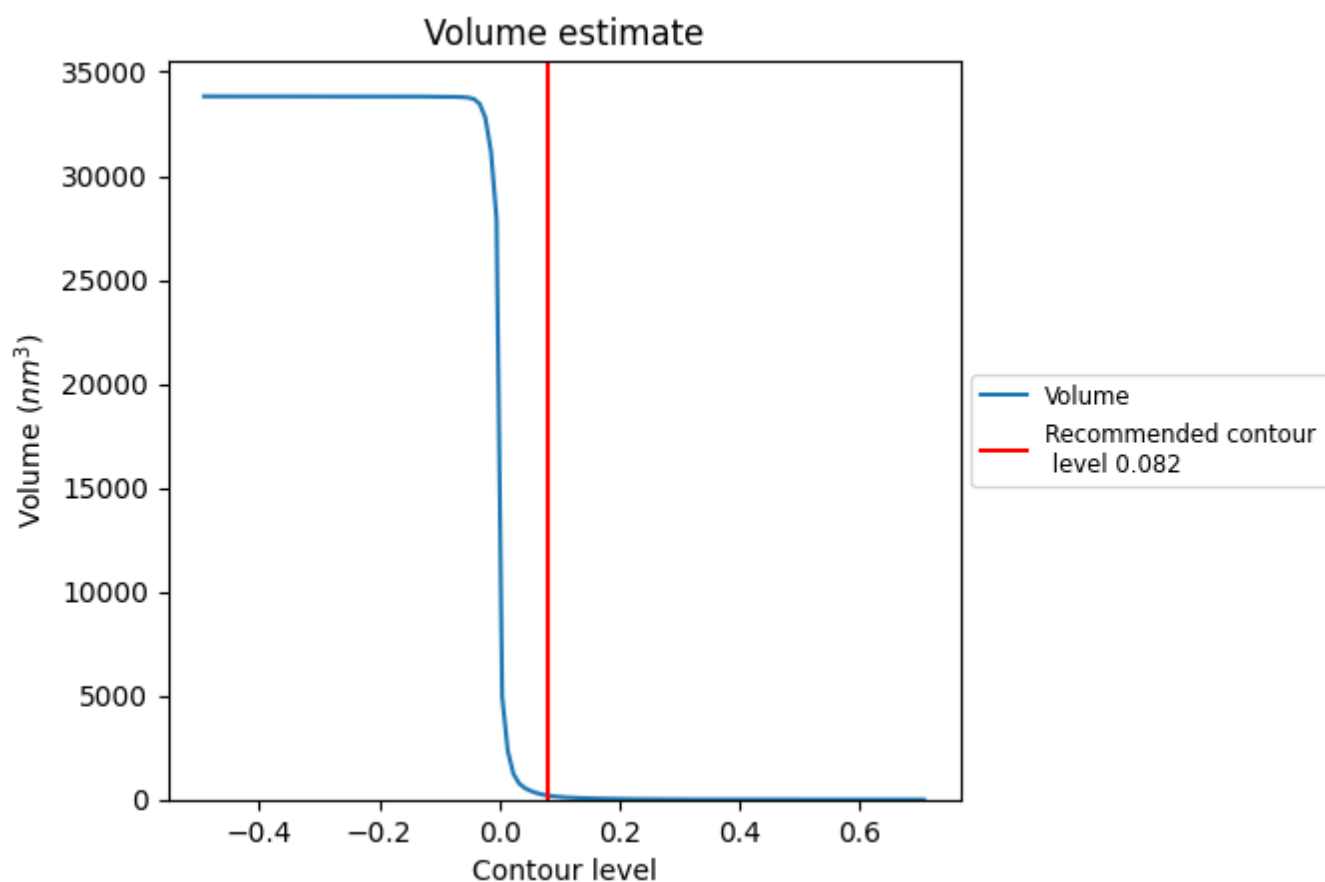
This section contains the results of statistical analysis of the map.

### 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.

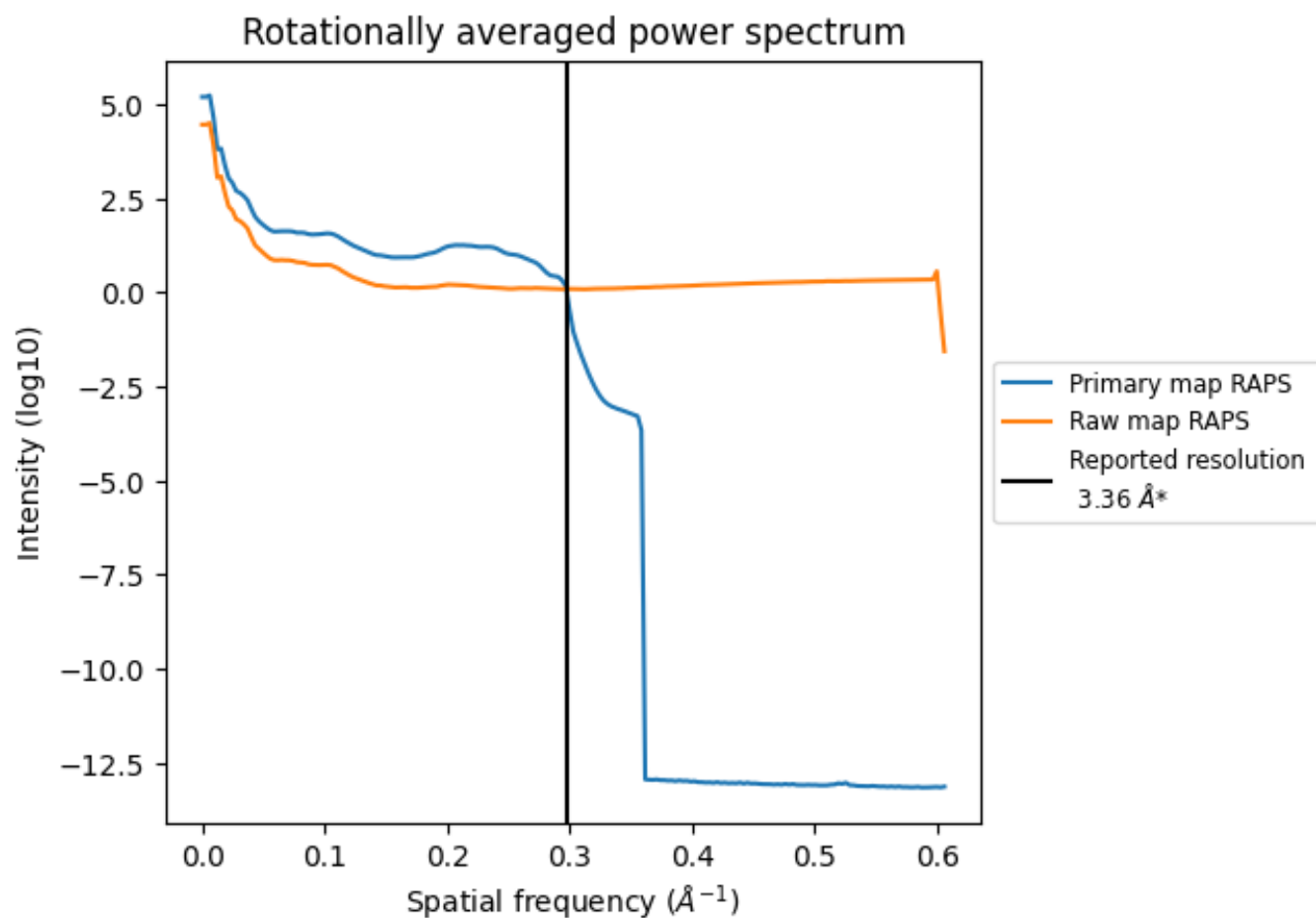
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 187 nm<sup>3</sup>; this corresponds to an approximate mass of 169 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 ⓘ

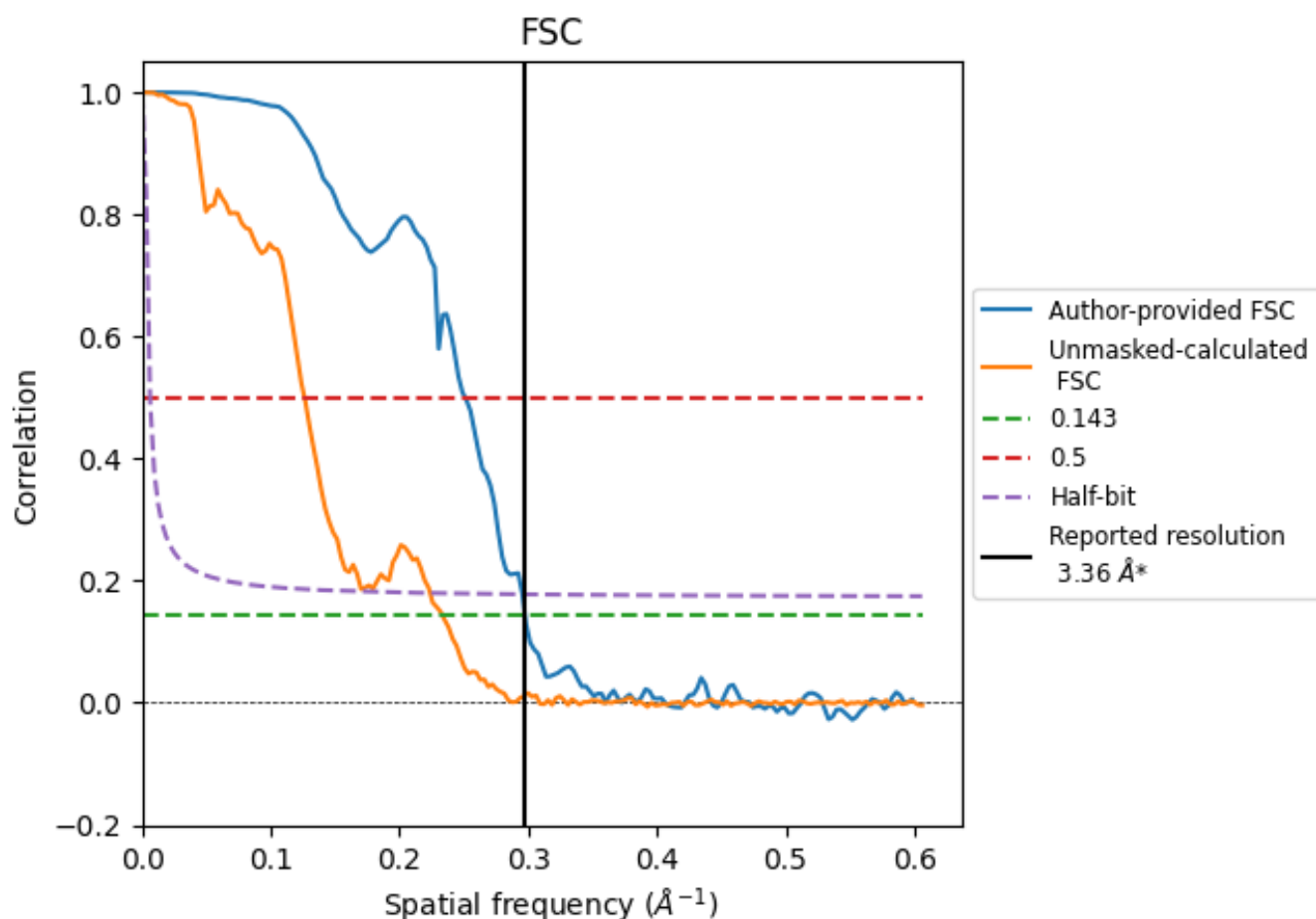


\*Reported resolution corresponds to spatial frequency of 0.298 Å<sup>-1</sup>

## 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 0.298  $\text{\AA}^{-1}$



## 8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.36	-	-
Author-provided FSC curve	3.36	3.98	3.39
Unmasked-calculated*	4.29	7.93	4.47

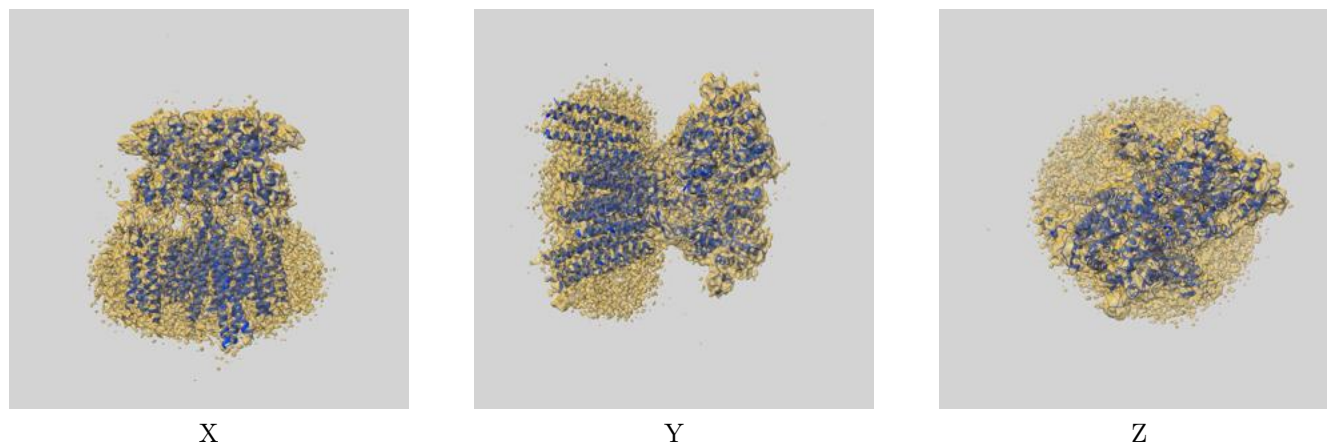
\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 4.29 differs from the reported value 3.36 by more than 10 %



## 9 Map-model fit [i](#)

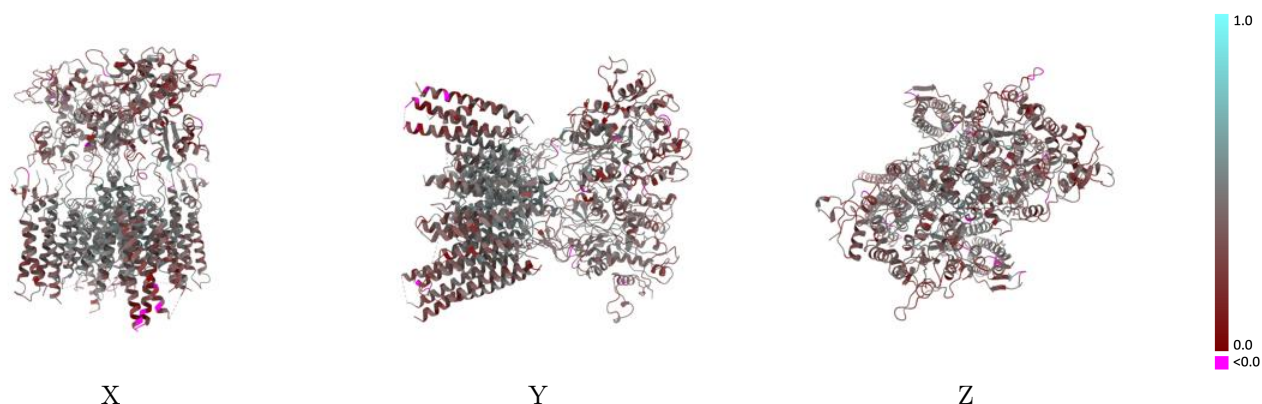
This section contains information regarding the fit between EMDB map EMD-74383 and PDB model 9ZKS. Per-residue inclusion information can be found in section [3](#) on page [8](#).

### 9.1 Map-model overlay [i](#)



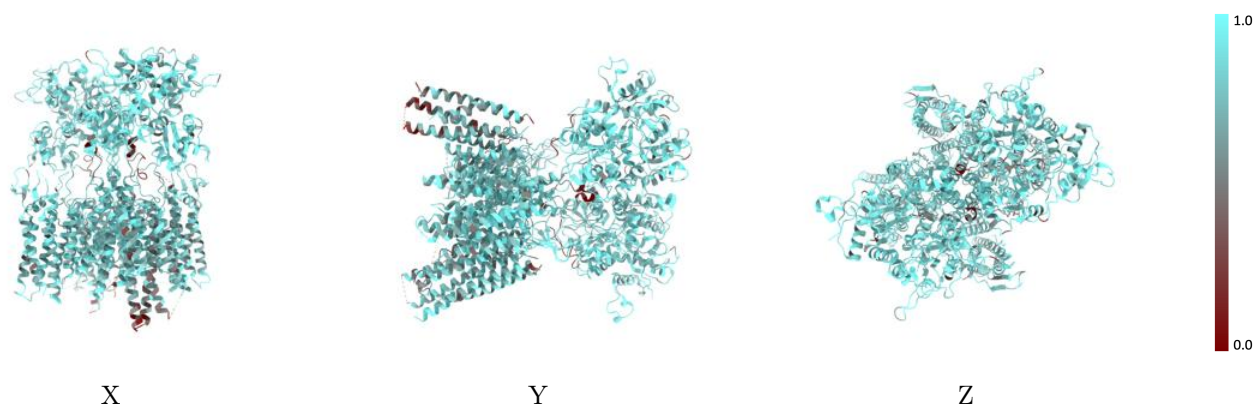
The images above show the 3D surface view of the map at the recommended contour level 0.082 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



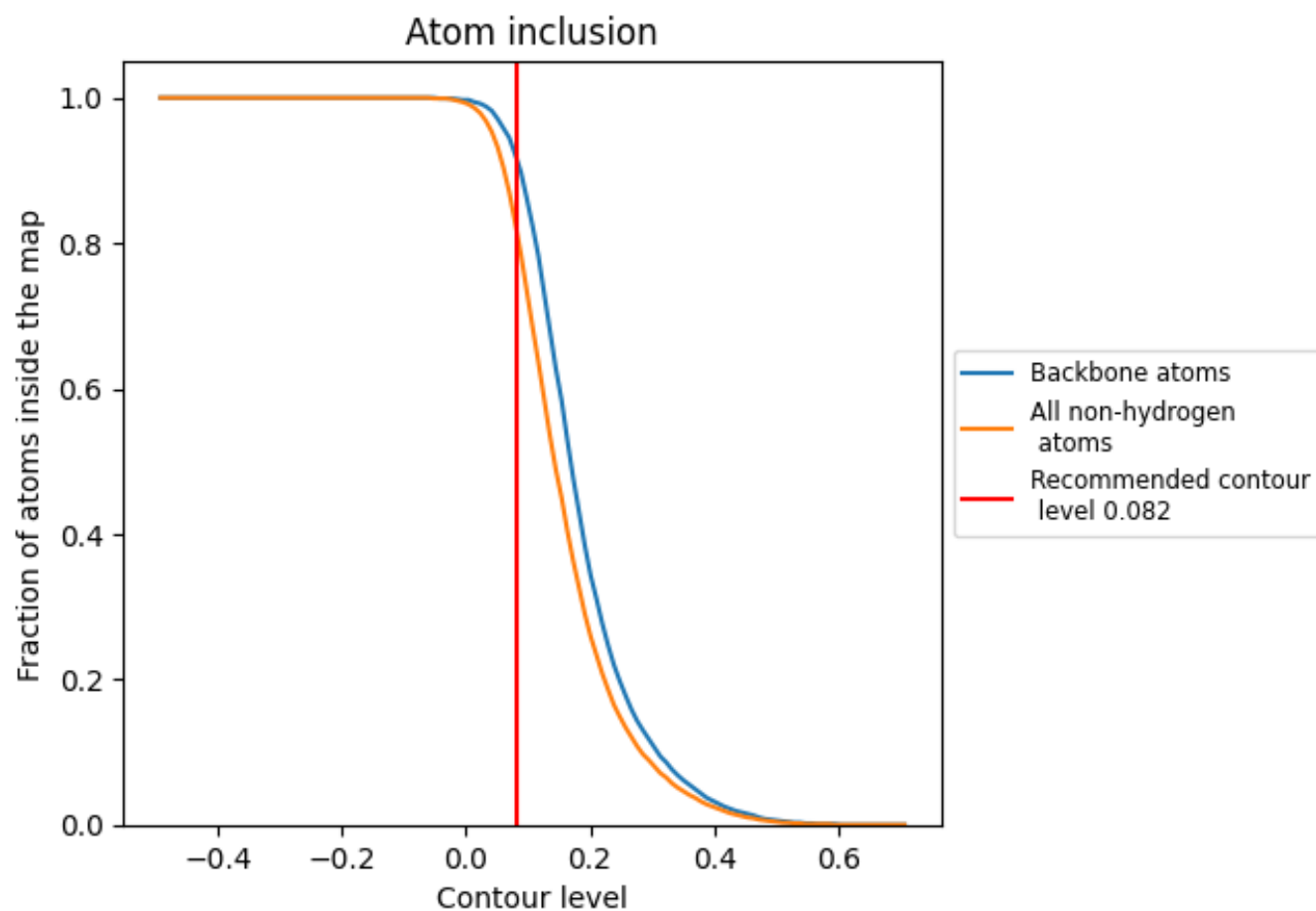
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.082).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 82% 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 (0.082) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.8170	<div><div></div></div> 0.3940
A	<div><div></div></div> 0.8410	<div><div></div></div> 0.4100
B	<div><div></div></div> 0.8400	<div><div></div></div> 0.4020
C	<div><div></div></div> 0.8390	<div><div></div></div> 0.4110
D	<div><div></div></div> 0.8430	<div><div></div></div> 0.4010
E	<div><div></div></div> 0.6210	<div><div></div></div> 0.3100
F	<div><div></div></div> 0.7810	<div><div></div></div> 0.3660
G	<div><div></div></div> 0.8060	<div><div></div></div> 0.3980
H	<div><div></div></div> 0.7930	<div><div></div></div> 0.3870
M	<div><div></div></div> 0.7670	<div><div></div></div> 0.2970

1.0

0.0

<0.0