



Full wwPDB EM Validation Report ⓘ

Jun 8, 2026 – 01:23 pm BST

PDB ID : 9RGE / pdb_00009rge
EMDB ID : EMD-53949
Title : CryoEM structure of human alpha1beta3gamma2 GABA(A)R in complex with GARLH4 and the Neuroligin2 transmembrane helix, in CL47a
Authors : Kasaragod, V.B.; Aricescu, A.R.
Deposited on : 2025-06-06
Resolution : 3.10 Å(reported)
Based on initial model : 9FAT

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:

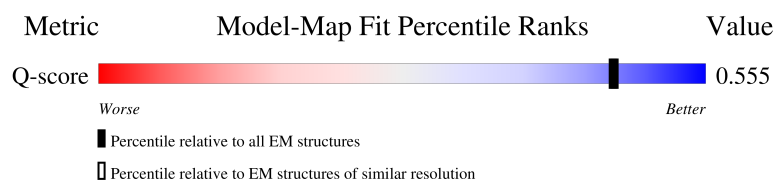
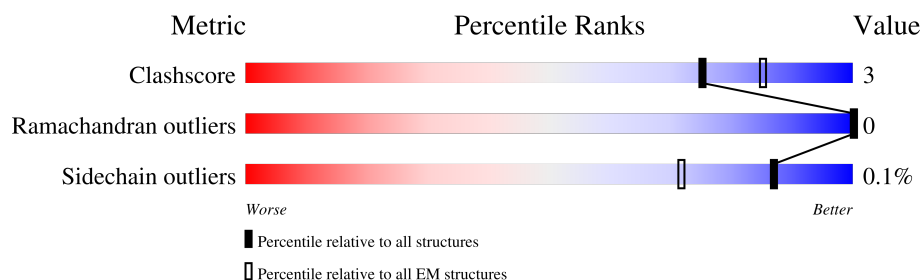
EMDB validation analysis : 0.0.1.dev132
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data 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	405	 81% 15%
1	D	405	 78% 8% 14%
2	B	439	 71% 5% 24%
2	E	439	 70% 5% 24%

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Mol	Chain	Length	Quality of chain
3	C	411	
4	H	29	
5	L	185	
6	F	10	
7	G	6	
7	K	6	
8	I	2	
8	M	2	
9	J	4	

2 Entry composition

There are 18 unique types of molecules in this entry. The entry contains 16281 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 Gamma-aminobutyric acid receptor subunit alpha-1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	345	Total	C	N	O	S	0	0
			2788	1807	466	499	16		
1	D	349	Total	C	N	O	S	0	0
			2822	1829	472	505	16		

- Molecule 2 is a protein called Gamma-aminobutyric acid receptor subunit beta-3.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	335	Total	C	N	O	S	0	0
			2752	1800	452	484	16		
2	E	332	Total	C	N	O	S	1	0
			2739	1796	449	478	16		

- Molecule 3 is a protein called Gamma-aminobutyric acid receptor subunit gamma-2.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	345	Total	C	N	O	S	0	0
			2833	1850	463	503	17		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	429	GLY	-	expression tag	UNP P18507

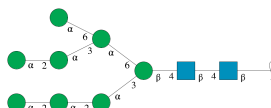
- Molecule 4 is a protein called Neuroligin-2.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	H	29	Total	C	N	O	0	0
			214	138	33	43		

- Molecule 5 is a protein called LHFPL tetraspan subfamily member 4 protein.

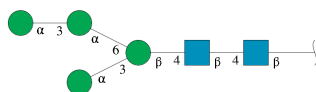
Mol	Chain	Residues	Atoms					AltConf	Trace
5	L	185	Total	C	N	O	S	0	0
			1421	943	222	241	15		

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



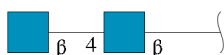
Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	10	Total	C	N	O		0	0
			116	64	2	50			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	6	Total	C	N	O		0	0
			72	40	2	30			
7	K	6	Total	C	N	O		0	0
			72	40	2	30			

- Molecule 8 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms					AltConf	Trace
8	I	2	Total	C	N	O		0	0
			28	16	2	10			

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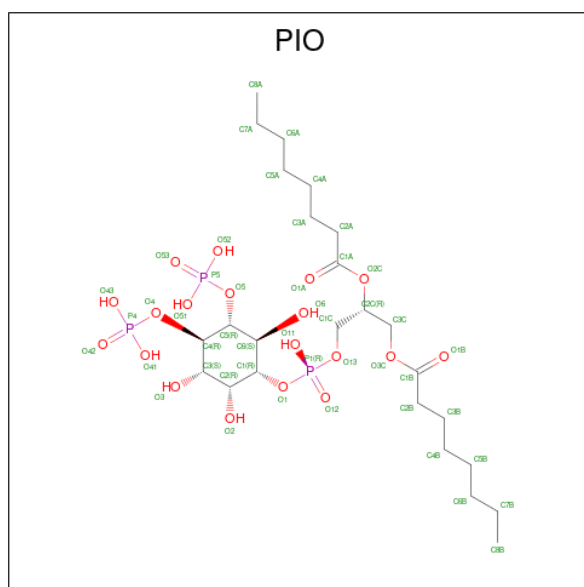
Mol	Chain	Residues	Atoms				AltConf	Trace
8	M	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



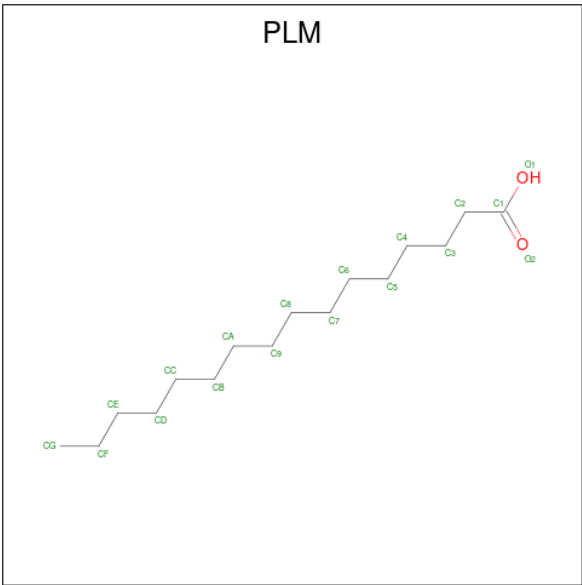
Mol	Chain	Residues	Atoms				AltConf	Trace
9	J	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 10 is [(2R)-2-octanoyloxy-3-[oxidanyl-[(1R,2R,3S,4R,5R,6S)-2,3,6-tris(oxidanyl)-4,5-diphosphonoxy-cyclohexyl]oxy-phosphoryl]oxy-propyl] octanoate (CCD ID: PIO) (formula: C₂₅H₄₉O₁₉P₃).



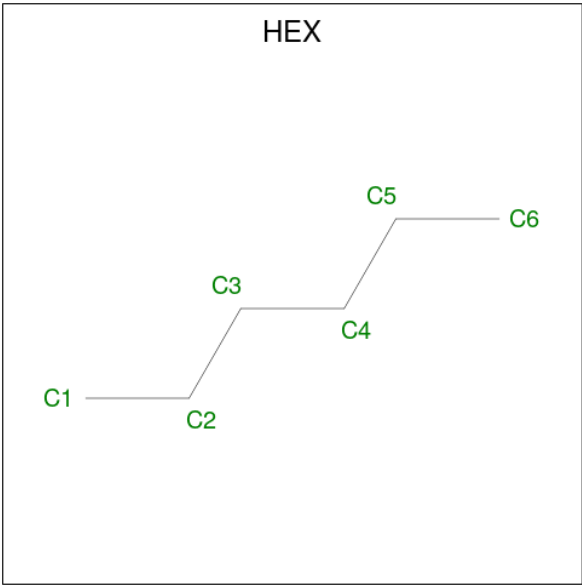
Mol	Chain	Residues	Atoms				AltConf
10	A	1	Total	C	O	P	0
			47	25	19	3	
10	D	1	Total	C	O	P	0
			47	25	19	3	

- Molecule 11 is PALMITIC ACID (CCD ID: PLM) (formula: C₁₆H₃₂O₂).



Mol	Chain	Residues	Atoms			AltConf
11	A	1	Total	C	O	0
			18	16	2	

- Molecule 12 is HEXANE (CCD ID: HEX) (formula: C₆H₁₄).



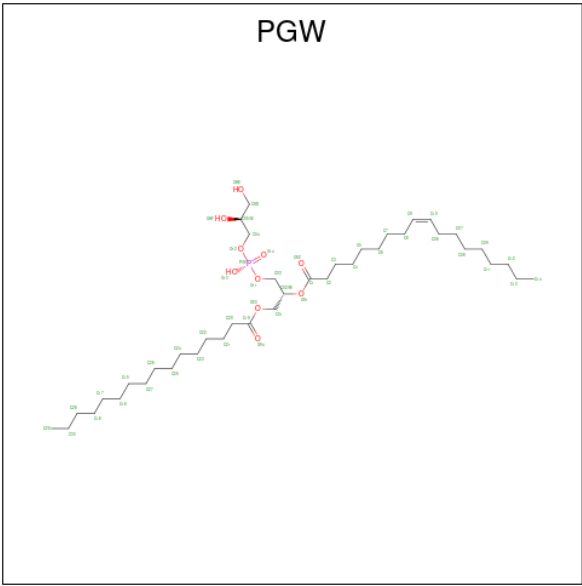
Mol	Chain	Residues	Atoms		AltConf
12	A	1	Total	C	0
			6	6	
12	A	1	Total	C	0
			6	6	
12	A	1	Total	C	0
			6	6	

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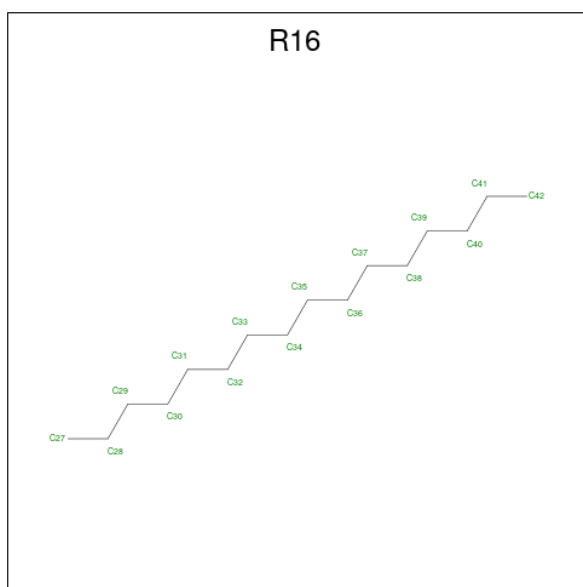
Mol	Chain	Residues	Atoms		AltConf
12	B	1	Total	C	0
			6	6	
12	D	1	Total	C	0
			6	6	

- Molecule 13 is (1R)-2-{[(S)-{[(2S)-2,3-dihydroxypropyl]oxy}(hydroxy)phosphoryl]oxy}-1-[(hexadecanoyloxy)methyl]ethyl (9Z)-octadec-9-enoate (CCD ID: PGW) (formula: C₄₀H₇₇O₁₀P).



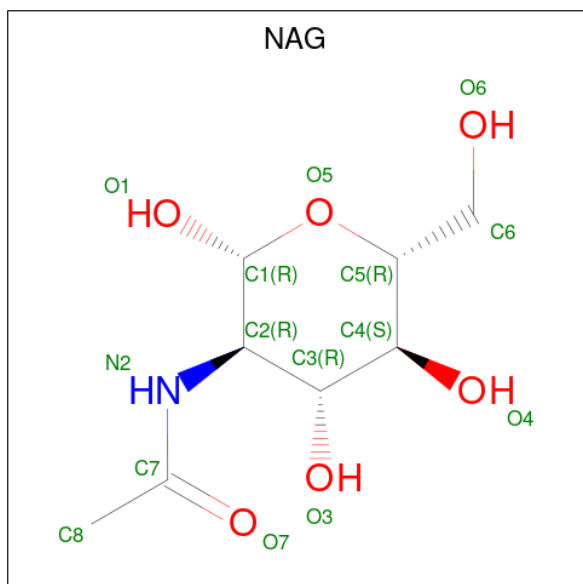
Mol	Chain	Residues	Atoms				AltConf
13	B	1	Total	C	O	P	0
			51	40	10	1	

- Molecule 14 is HEXADECANE (CCD ID: R16) (formula: C₁₆H₃₄).



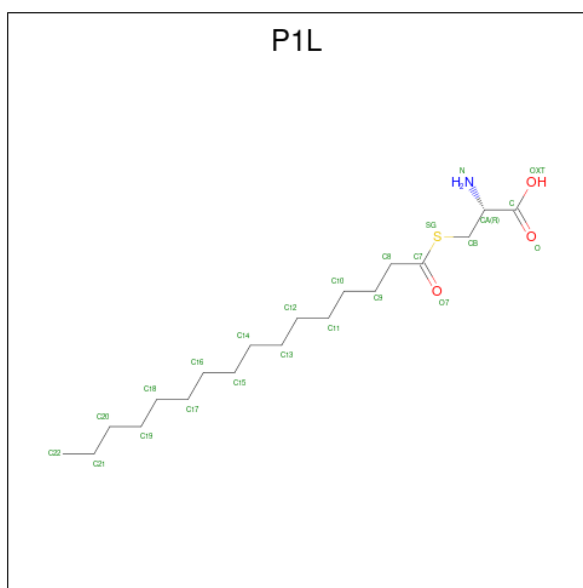
Mol	Chain	Residues	Atoms		AltConf
14	B	1	Total	C	0
			16	16	
14	D	1	Total	C	0
			16	16	

- Molecule 15 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



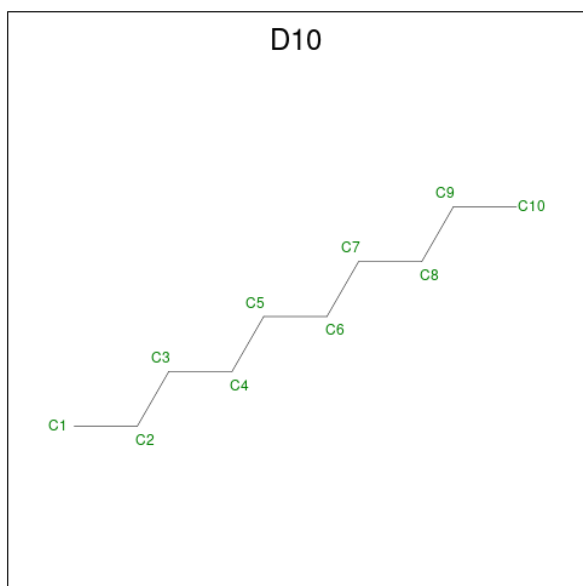
Mol	Chain	Residues	Atoms				AltConf
15	C	1	Total	C	N	O	0
			14	8	1	5	

- Molecule 16 is S-PALMITOYL-L-CYSTEINE (CCD ID: P1L) (formula: $C_{19}H_{37}NO_3S$).



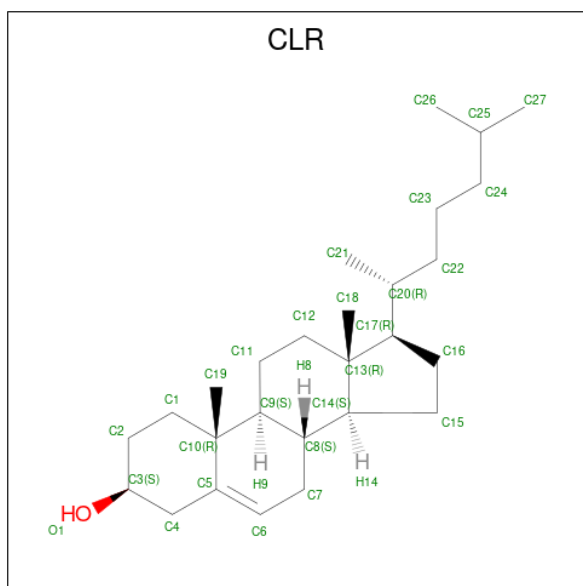
Mol	Chain	Residues	Atoms					AltConf
16	C	1	Total	C	N	O	S	0
			23	19	1	2	1	
16	C	1	Total	C	N	O	S	0
			23	19	1	2	1	
16	C	1	Total	C	N	O	S	0
			23	19	1	2	1	

- Molecule 17 is DECANE (CCD ID: D10) (formula: $C_{10}H_{22}$).



Mol	Chain	Residues	Atoms		AltConf
17	D	1	Total	C	0
			10	10	

- Molecule 18 is CHOLESTEROL (CCD ID: CLR) (formula: $C_{27}H_{46}O$) (labeled as "Ligand of Interest" by depositor).




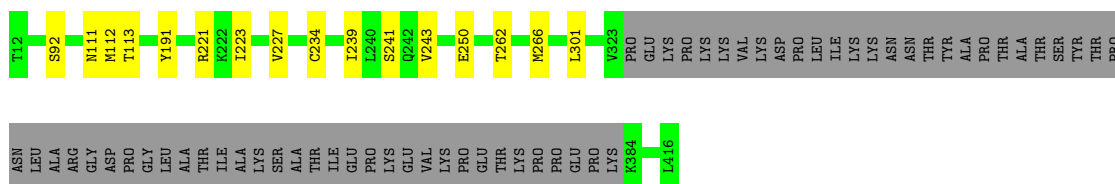
Mol	Chain	Residues	Atoms			AltConf
18	L	1	Total	C	O	0
			28	27	1	

3 Residue-property plots [i](#)


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.

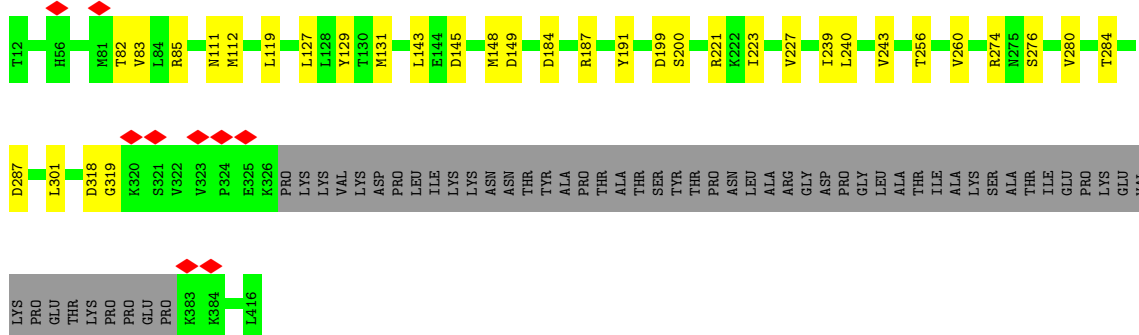
- Molecule 1: Gamma-aminobutyric acid receptor subunit alpha-1

Chain A: 



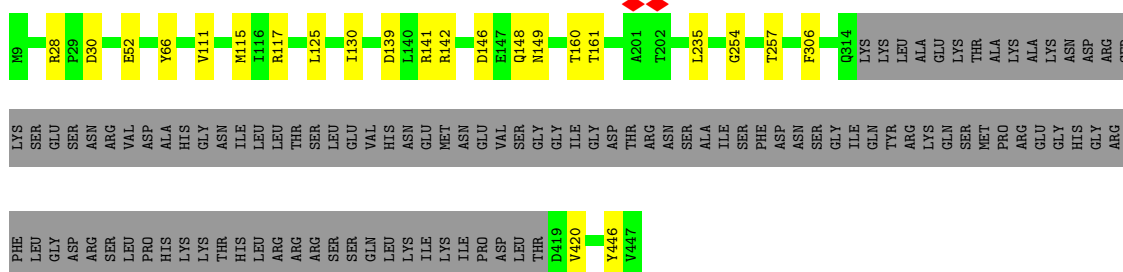
- Molecule 1: Gamma-aminobutyric acid receptor subunit alpha-1

Chain D: 

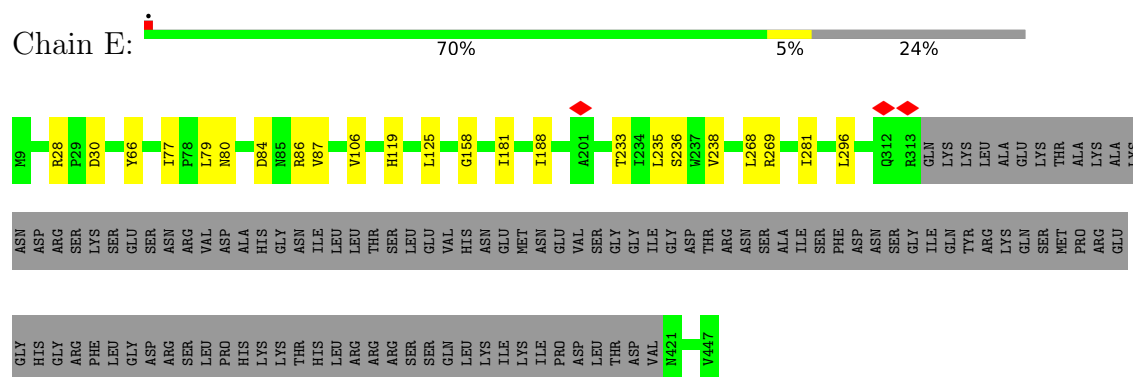


- Molecule 2: Gamma-aminobutyric acid receptor subunit beta-3

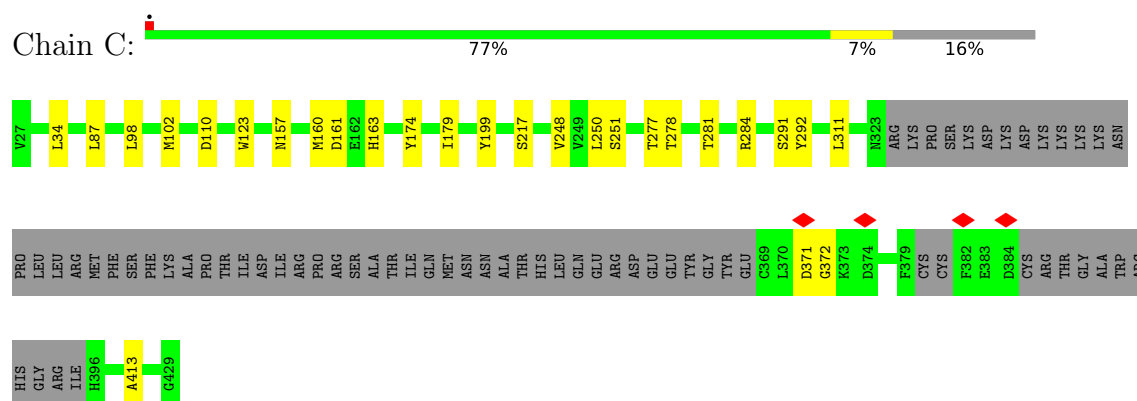
Chain B: 



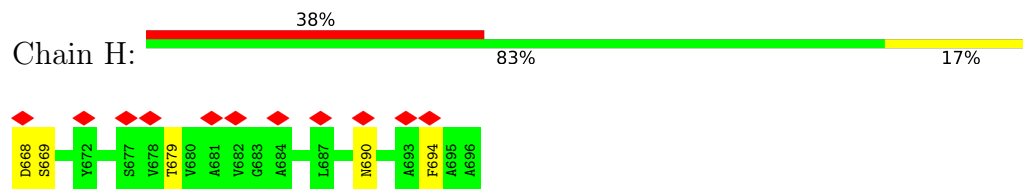
- Molecule 2: Gamma-aminobutyric acid receptor subunit beta-3



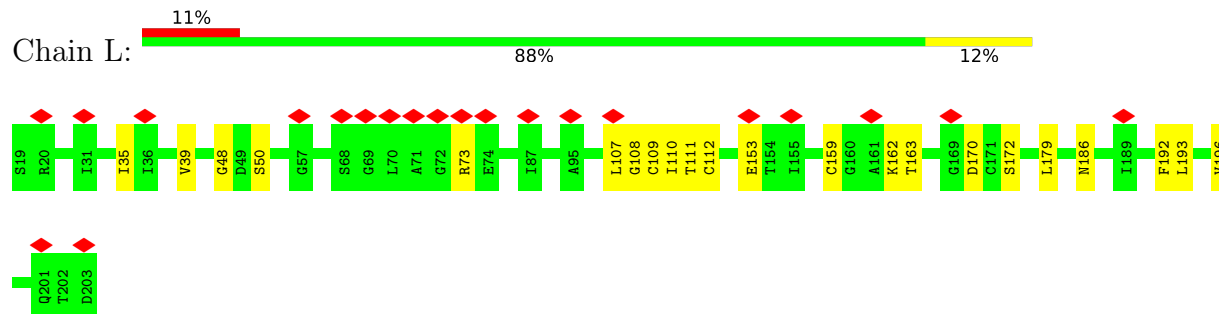
- Molecule 3: Gamma-aminobutyric acid receptor subunit gamma-2



- Molecule 4: Neuroligin-2



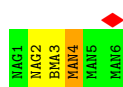
- Molecule 5: LHFPL tetraspan subfamily member 4 protein



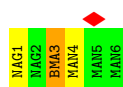
- Molecule 6: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 7: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 7: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



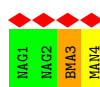
- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 8: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 9: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	22879	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 ($e^-/\text{\AA}^2$)	53.2	Depositor
Minimum defocus (nm)	600	Depositor
Maximum defocus (nm)	1600	Depositor
Magnification	165000	Depositor
Image detector	TFS FALCON 4i (4k x 4k)	Depositor
Maximum map value	0.018	Depositor
Minimum map value	-0.008	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.001	Depositor
Recommended contour level	0.0034	Depositor
Map size (Å)	262.44, 262.44, 262.44	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.72900003, 0.72900003, 0.72900003	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: P1L, PIO, PGW, CLR, NAG, HEX, D10, BMA, R16, MAN, PLM

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	0.16	0/2859	0.32	0/3884
1	D	0.15	0/2894	0.31	0/3930
2	B	0.17	0/2825	0.36	0/3842
2	E	0.15	0/2817	0.34	0/3832
3	C	0.16	0/2909	0.34	0/3955
4	H	0.11	0/216	0.25	0/294
5	L	0.15	0/1459	0.30	0/1983
All	All	0.16	0/15979	0.33	0/21720

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 [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2788	0	2783	13	0
1	D	2822	0	2823	24	0
2	B	2752	0	2746	16	0
2	E	2739	0	2735	17	0
3	C	2833	0	2805	20	0
4	H	214	0	214	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	L	1421	0	1429	15	0
6	F	116	0	97	6	0
7	G	72	0	61	1	0
7	K	72	0	61	2	0
8	I	28	0	25	0	0
8	M	28	0	25	0	0
9	J	50	0	43	1	0
10	A	47	0	44	0	0
10	D	47	0	44	0	0
11	A	18	0	31	3	0
12	A	18	0	42	0	0
12	B	6	0	14	0	0
12	D	6	0	14	0	0
13	B	51	0	76	2	0
14	B	16	0	34	0	0
14	D	16	0	34	0	0
15	C	14	0	13	0	0
16	C	69	0	105	1	0
17	D	10	0	22	0	0
18	L	28	0	46	1	0
All	All	16281	0	16366	101	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 3.

All (101) 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)
7:K:3:BMA:H4	7:K:4:MAN:H3	1.68	0.73
3:C:278:THR:O	3:C:281:THR:OG1	2.06	0.73
2:B:142:ARG:NH2	2:B:446:TYR:O	2.23	0.71
1:D:274:ARG:NH2	1:D:280:VAL:O	2.24	0.71
3:C:413:ALA:HB1	18:L:601:CLR:H271	1.74	0.70
1:D:85:ARG:NH1	2:E:158:GLY:O	2.25	0.69
1:A:301:LEU:HD23	2:E:235:LEU:HD22	1.74	0.68
16:C:504:P1L:SG	16:C:504:P1L:O	2.52	0.68
1:D:274:ARG:NH1	1:D:287:ASP:OD2	2.27	0.67
3:C:248:VAL:O	3:C:251:SER:OG	2.14	0.66
3:C:123:TRP:CZ2	6:F:3:BMA:H61	2.33	0.64
13:B:501:PGW:O14	3:C:291:SER:OG	2.16	0.62
4:H:690:ASN:ND2	5:L:186:ASN:OD1	2.33	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:82:THR:HG23	1:D:83:VAL:HG23	1.82	0.61
4:H:668:ASP:OD2	5:L:73:ARG:NH2	2.36	0.59
3:C:250:LEU:HD22	1:D:301:LEU:HD23	1.85	0.58
1:A:92:SER:O	2:E:86:ARG:NH1	2.37	0.58
3:C:123:TRP:CH2	6:F:3:BMA:H61	2.39	0.56
2:B:235:LEU:HD22	3:C:311:LEU:HD23	1.88	0.56
1:D:112:MET:HE1	2:E:106:VAL:HG23	1.88	0.55
1:A:191:TYR:OH	1:A:221:ARG:NH2	2.40	0.55
5:L:109:CYS:HA	5:L:112:CYS:SG	2.47	0.54
1:D:129:TYR:CE2	1:D:131:MET:HE2	2.43	0.54
2:B:146:ASP:OD2	2:B:148:GLN:NE2	2.40	0.53
7:G:3:BMA:H4	7:G:4:MAN:H3	1.91	0.53
2:E:80:ASN:OD1	2:E:119:HIS:ND1	2.41	0.53
2:E:233:THR:O	2:E:236:SER:OG	2.22	0.52
1:D:199:ASP:OD1	1:D:200:SER:N	2.43	0.51
3:C:277:THR:O	3:C:281:THR:HG23	2.10	0.51
3:C:157:ASN:O	3:C:160:MET:N	2.43	0.50
11:A:3902:PLM:HE1	2:E:238:VAL:HG22	1.94	0.50
5:L:192:PHE:O	5:L:196:VAL:HG23	2.12	0.49
7:K:3:BMA:C4	7:K:4:MAN:H3	2.39	0.49
1:A:234:CYS:SG	1:A:266:MET:HE1	2.52	0.49
5:L:50:SER:OG	5:L:170:ASP:O	2.15	0.49
2:B:52:GLU:OE1	2:B:52:GLU:N	2.46	0.48
3:C:34:LEU:HD22	3:C:87:LEU:HA	1.94	0.48
1:D:318:ASP:OD1	1:D:319:GLY:N	2.46	0.48
1:A:111:ASN:HB3	6:F:1:NAG:HN2	1.78	0.47
1:A:241:SER:HB3	1:A:262:THR:HG21	1.96	0.47
1:A:112:MET:HE3	1:A:113:THR:HA	1.96	0.47
2:B:28:ARG:NH1	2:B:30:ASP:O	2.48	0.47
2:B:111:VAL:HG21	3:C:110:ASP:HA	1.97	0.47
4:H:694:PHE:HE1	5:L:196:VAL:HG21	1.80	0.47
11:A:3902:PLM:H51	11:A:3902:PLM:H22	1.54	0.46
1:A:112:MET:HE2	2:B:130:ILE:HD11	1.96	0.46
3:C:371:ASP:OD1	3:C:372:GLY:N	2.48	0.46
3:C:199:TYR:CD1	1:D:280:VAL:HG12	2.51	0.46
2:E:66:TYR:CZ	2:E:125:LEU:HD13	2.50	0.46
11:A:3902:PLM:H51	11:A:3902:PLM:H82	1.77	0.46
3:C:284:ARG:NH2	3:C:292:TYR:O	2.49	0.46
1:D:184:ASP:O	1:D:187:ARG:NH1	2.47	0.46
5:L:48:GLY:N	5:L:172:SER:O	2.48	0.46
5:L:153:GLU:N	5:L:153:GLU:OE1	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:ASP:O	1:D:148:MET:N	2.49	0.46
2:B:117:ARG:NH1	3:C:217:SER:O	2.48	0.45
9:J:3:BMA:O2	9:J:4:MAN:H61	2.15	0.45
2:B:66:TYR:CZ	2:B:125:LEU:HD13	2.52	0.45
5:L:107:LEU:HD23	5:L:110:ILE:HD12	1.99	0.45
3:C:174:TYR:CG	3:C:179:ILE:HD12	2.52	0.45
1:D:239:ILE:O	1:D:243:VAL:HG23	2.16	0.45
1:D:129:TYR:HE2	1:D:131:MET:HE2	1.82	0.44
4:H:679:THR:HG23	5:L:179:LEU:HD12	1.98	0.44
3:C:161:ASP:OD2	3:C:163:HIS:NE2	2.49	0.44
1:D:143:LEU:O	1:D:284:THR:HG22	2.18	0.44
5:L:159:CYS:HB3	5:L:163:THR:HG21	2.00	0.43
2:B:306:PHE:HB2	2:B:420:VAL:HG11	2.01	0.43
1:D:111:ASN:HD22	6:F:9:MAN:H61	1.83	0.43
2:E:28:ARG:NH1	2:E:30:ASP:O	2.51	0.43
5:L:35:ILE:O	5:L:39:VAL:HG23	2.19	0.43
1:A:239:ILE:O	1:A:243:VAL:HG23	2.17	0.43
6:F:2:NAG:O3	6:F:3:BMA:O5	2.35	0.43
2:B:139:ASP:OD2	2:B:141:ARG:NH2	2.51	0.43
1:A:112:MET:HE2	2:B:130:ILE:CD1	2.49	0.42
2:E:268:LEU:HD11	2:E:281:ILE:HD11	2.01	0.42
1:D:256:THR:O	1:D:260:VAL:HG23	2.19	0.42
1:D:119:LEU:HD11	1:D:127:LEU:HD23	2.00	0.42
3:C:174:TYR:CD1	3:C:179:ILE:HD12	2.54	0.42
1:D:240:LEU:HD22	2:E:296:LEU:HD23	2.02	0.42
2:B:160:THR:OG1	2:B:161:THR:N	2.53	0.42
4:H:668:ASP:OD1	4:H:669:SER:N	2.53	0.42
5:L:162:LYS:NZ	5:L:170:ASP:OD1	2.30	0.42
1:A:301:LEU:CD2	2:E:235:LEU:HD22	2.48	0.41
2:B:66:TYR:CE2	2:B:125:LEU:HD13	2.54	0.41
1:A:223:ILE:O	1:A:227:VAL:HG23	2.20	0.41
2:B:254:GLY:O	2:B:257:THR:OG1	2.32	0.41
3:C:98:LEU:HD13	3:C:102:MET:HG3	2.02	0.41
1:D:149:ASP:OD1	1:D:149:ASP:N	2.51	0.41
1:D:276:SER:OG	2:E:269:ARG:NH2	2.53	0.41
5:L:108:GLY:O	5:L:111:THR:OG1	2.31	0.41
1:A:250:GLU:OE1	1:A:250:GLU:N	2.53	0.41
13:B:501:PGW:H26A	13:B:501:PGW:H8A	2.02	0.41
2:B:115:MET:HE3	2:B:125:LEU:HG	2.03	0.41
1:D:112:MET:HE1	2:E:106:VAL:CG2	2.51	0.41
1:D:191:TYR:OH	1:D:221:ARG:NH2	2.54	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:223:ILE:HG22	1:D:227:VAL:HG23	2.03	0.41
2:E:181:ILE:HD11	2:E:188:ILE:HD12	2.03	0.41
4:H:694:PHE:HA	5:L:193:LEU:HD13	2.03	0.40
6:F:7:MAN:H3	6:F:8:MAN:H5	2.03	0.40
2:E:77:ILE:HG22	2:E:79:LEU:H	1.86	0.40
2:E:84:ASP:O	2:E:87:VAL:HG12	2.21	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	341/405 (84%)	331 (97%)	10 (3%)	0	100	100
1	D	345/405 (85%)	333 (96%)	12 (4%)	0	100	100
2	B	331/439 (75%)	325 (98%)	6 (2%)	0	100	100
2	E	329/439 (75%)	320 (97%)	9 (3%)	0	100	100
3	C	337/411 (82%)	322 (96%)	15 (4%)	0	100	100
4	H	27/29 (93%)	26 (96%)	1 (4%)	0	100	100
5	L	183/185 (99%)	175 (96%)	8 (4%)	0	100	100
All	All	1893/2313 (82%)	1832 (97%)	61 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	307/359 (86%)	307 (100%)	0	100	100
1	D	311/359 (87%)	311 (100%)	0	100	100
2	B	301/392 (77%)	300 (100%)	1 (0%)	86	87
2	E	299/392 (76%)	299 (100%)	0	100	100
3	C	317/376 (84%)	317 (100%)	0	100	100
4	H	23/23 (100%)	23 (100%)	0	100	100
5	L	151/151 (100%)	151 (100%)	0	100	100
All	All	1709/2052 (83%)	1708 (100%)	1 (0%)	87	90

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

Mol	Chain	Res	Type
2	B	149	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	149	ASN
2	B	191	HIS
2	B	439	ASN
3	C	80	GLN
1	D	111	ASN
2	E	65	GLN
2	E	85	ASN
2	E	267	HIS
4	H	690	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 ⓘ

30 monosaccharides 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	$\# Z > 2$
6	NAG	F	1	6,1	14,14,15	0.39	0	17,19,21	0.66	0
6	MAN	F	10	6	11,11,12	0.32	0	15,15,17	0.94	1 (6%)
6	NAG	F	2	6	14,14,15	0.40	0	17,19,21	0.71	0
6	BMA	F	3	6	11,11,12	0.63	0	15,15,17	0.65	0
6	MAN	F	4	6	11,11,12	0.33	0	15,15,17	0.80	1 (6%)
6	MAN	F	5	6	11,11,12	0.40	0	15,15,17	0.80	1 (6%)
6	MAN	F	6	6	11,11,12	0.35	0	15,15,17	0.51	0
6	MAN	F	7	6	11,11,12	0.50	0	15,15,17	0.57	0
6	MAN	F	8	6	11,11,12	0.31	0	15,15,17	0.74	0
6	MAN	F	9	6	11,11,12	0.34	0	15,15,17	0.54	0
7	NAG	G	1	7	14,14,15	0.40	0	17,19,21	0.75	0
7	NAG	G	2	7	14,14,15	0.39	0	17,19,21	0.93	1 (5%)
7	BMA	G	3	7	11,11,12	0.35	0	15,15,17	0.60	0
7	MAN	G	4	7	11,11,12	0.42	0	15,15,17	0.90	1 (6%)
7	MAN	G	5	7	11,11,12	0.36	0	15,15,17	0.58	0
7	MAN	G	6	7	11,11,12	0.29	0	15,15,17	0.49	0
8	NAG	I	1	8	14,14,15	0.41	0	17,19,21	0.50	0
8	NAG	I	2	8	14,14,15	0.38	0	17,19,21	0.51	0
9	NAG	J	1	9	14,14,15	0.39	0	17,19,21	0.74	0
9	NAG	J	2	9	14,14,15	0.38	0	17,19,21	0.70	0
9	BMA	J	3	9	11,11,12	0.28	0	15,15,17	0.86	1 (6%)
9	MAN	J	4	9	11,11,12	0.33	0	15,15,17	0.56	0
7	NAG	K	1	7	14,14,15	0.43	0	17,19,21	0.91	1 (5%)
7	NAG	K	2	7	14,14,15	0.39	0	17,19,21	0.68	0
7	BMA	K	3	7	11,11,12	0.65	0	15,15,17	0.83	1 (6%)
7	MAN	K	4	7	11,11,12	0.32	0	15,15,17	0.72	0
7	MAN	K	5	7	11,11,12	0.36	0	15,15,17	0.61	0
7	MAN	K	6	7	11,11,12	0.33	0	15,15,17	0.58	0
8	NAG	M	1	8	14,14,15	0.39	0	17,19,21	0.78	0
8	NAG	M	2	8	14,14,15	0.39	0	17,19,21	0.61	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
6	NAG	F	1	6,1	-	2/6/23/26	0/1/1/1
6	MAN	F	10	6	-	1/2/19/22	0/1/1/1
6	NAG	F	2	6	-	1/6/23/26	0/1/1/1
6	BMA	F	3	6	-	0/2/19/22	0/1/1/1
6	MAN	F	4	6	-	0/2/19/22	0/1/1/1
6	MAN	F	5	6	-	0/2/19/22	0/1/1/1
6	MAN	F	6	6	-	0/2/19/22	0/1/1/1
6	MAN	F	7	6	-	2/2/19/22	0/1/1/1
6	MAN	F	8	6	-	0/2/19/22	0/1/1/1
6	MAN	F	9	6	-	1/2/19/22	0/1/1/1
7	NAG	G	1	7	-	0/6/23/26	0/1/1/1
7	NAG	G	2	7	-	0/6/23/26	0/1/1/1
7	BMA	G	3	7	-	2/2/19/22	0/1/1/1
7	MAN	G	4	7	-	0/2/19/22	0/1/1/1
7	MAN	G	5	7	-	0/2/19/22	0/1/1/1
7	MAN	G	6	7	-	1/2/19/22	0/1/1/1
8	NAG	I	1	8	-	0/6/23/26	0/1/1/1
8	NAG	I	2	8	-	1/6/23/26	0/1/1/1
9	NAG	J	1	9	-	4/6/23/26	0/1/1/1
9	NAG	J	2	9	-	0/6/23/26	0/1/1/1
9	BMA	J	3	9	-	2/2/19/22	0/1/1/1
9	MAN	J	4	9	-	2/2/19/22	0/1/1/1
7	NAG	K	1	7	-	1/6/23/26	0/1/1/1
7	NAG	K	2	7	-	2/6/23/26	0/1/1/1
7	BMA	K	3	7	-	2/2/19/22	0/1/1/1
7	MAN	K	4	7	-	1/2/19/22	0/1/1/1
7	MAN	K	5	7	-	0/2/19/22	0/1/1/1
7	MAN	K	6	7	-	0/2/19/22	0/1/1/1
8	NAG	M	1	8	-	4/6/23/26	0/1/1/1
8	NAG	M	2	8	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	K	1	NAG	C1-C2-N2	-2.87	105.59	110.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	J	3	BMA	C1-O5-C5	2.47	115.54	112.19
6	F	10	MAN	C1-O5-C5	2.45	115.52	112.19
7	G	2	NAG	O4-C4-C3	-2.34	104.94	110.35
6	F	4	MAN	C1-O5-C5	2.28	115.28	112.19
6	F	5	MAN	C1-O5-C5	2.20	115.17	112.19
7	G	4	MAN	O3-C3-C2	2.16	114.12	109.99
7	K	3	BMA	C1-O5-C5	2.04	114.96	112.19

There are no chirality outliers.

All (31) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	M	1	NAG	C3-C2-N2-C7
8	M	1	NAG	C8-C7-N2-C2
8	M	1	NAG	O7-C7-N2-C2
9	J	1	NAG	C3-C2-N2-C7
9	J	1	NAG	C8-C7-N2-C2
9	J	1	NAG	O7-C7-N2-C2
8	M	2	NAG	C8-C7-N2-C2
7	K	3	BMA	O5-C5-C6-O6
8	M	2	NAG	O7-C7-N2-C2
9	J	3	BMA	O5-C5-C6-O6
9	J	3	BMA	C4-C5-C6-O6
6	F	7	MAN	O5-C5-C6-O6
7	G	3	BMA	O5-C5-C6-O6
9	J	4	MAN	O5-C5-C6-O6
7	K	1	NAG	O5-C5-C6-O6
6	F	7	MAN	C4-C5-C6-O6
7	G	3	BMA	C4-C5-C6-O6
7	K	2	NAG	C8-C7-N2-C2
7	G	6	MAN	O5-C5-C6-O6
6	F	1	NAG	C1-C2-N2-C7
6	F	10	MAN	O5-C5-C6-O6
6	F	2	NAG	O5-C5-C6-O6
7	K	4	MAN	O5-C5-C6-O6
7	K	2	NAG	O7-C7-N2-C2
6	F	9	MAN	O5-C5-C6-O6
7	K	3	BMA	C4-C5-C6-O6
9	J	1	NAG	C1-C2-N2-C7
8	M	1	NAG	C1-C2-N2-C7
8	I	2	NAG	C1-C2-N2-C7
9	J	4	MAN	C4-C5-C6-O6

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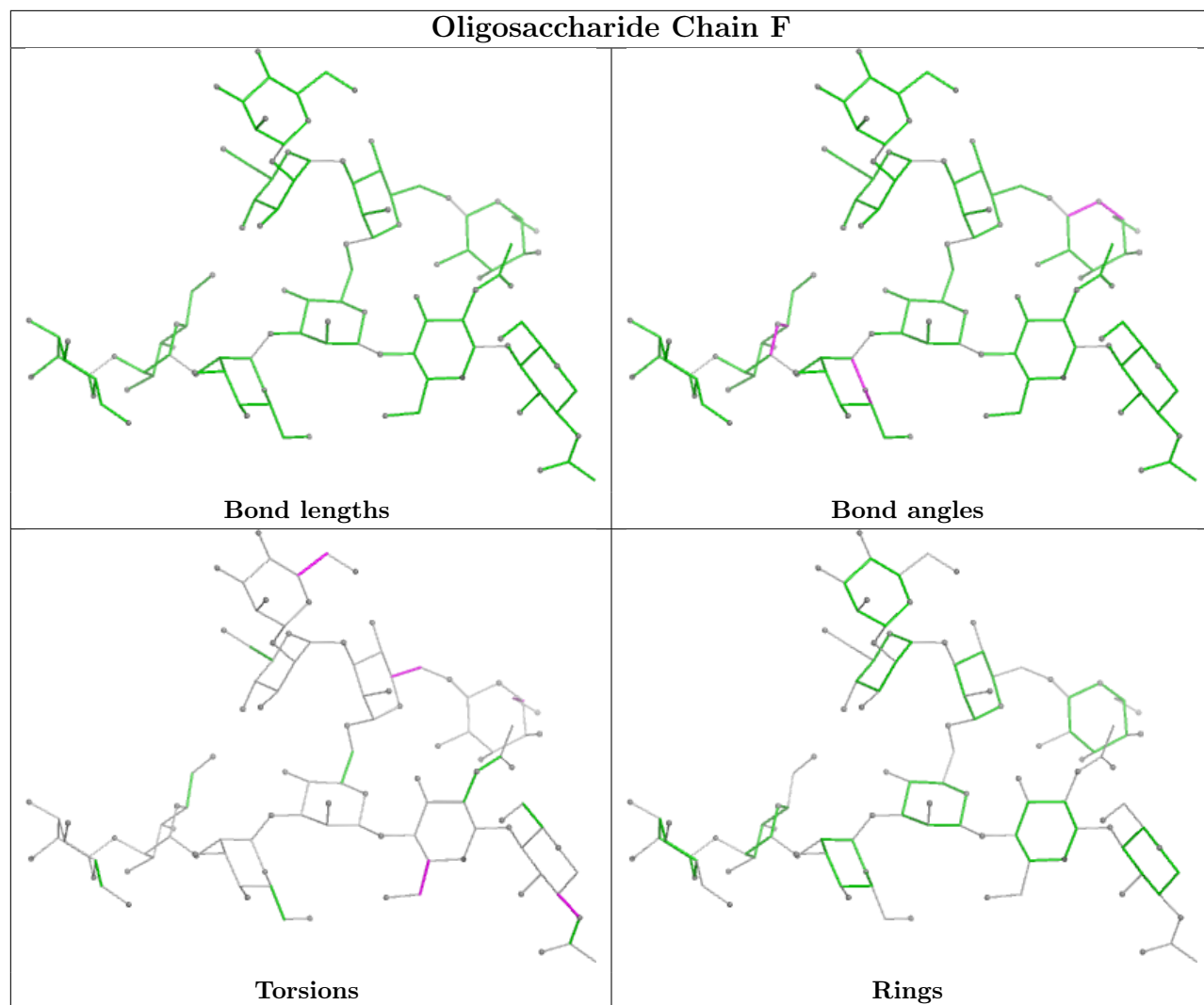
Mol	Chain	Res	Type	Atoms
6	F	1	NAG	C3-C2-N2-C7

There are no ring outliers.

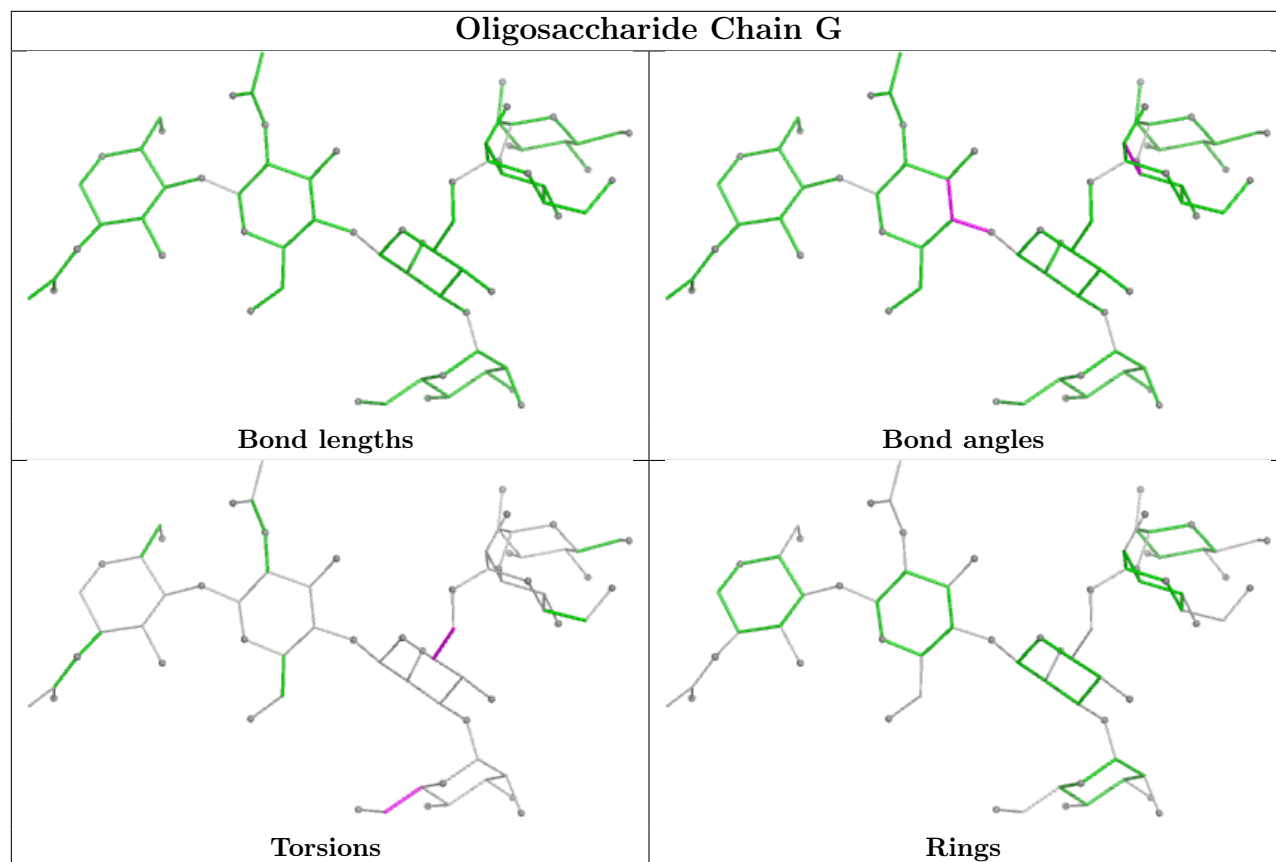
12 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	F	7	MAN	1	0
7	K	3	BMA	2	0
6	F	8	MAN	1	0
9	J	3	BMA	1	0
6	F	3	BMA	3	0
6	F	9	MAN	1	0
9	J	4	MAN	1	0
7	G	4	MAN	1	0
6	F	1	NAG	1	0
7	G	3	BMA	1	0
6	F	2	NAG	1	0
7	K	4	MAN	2	0

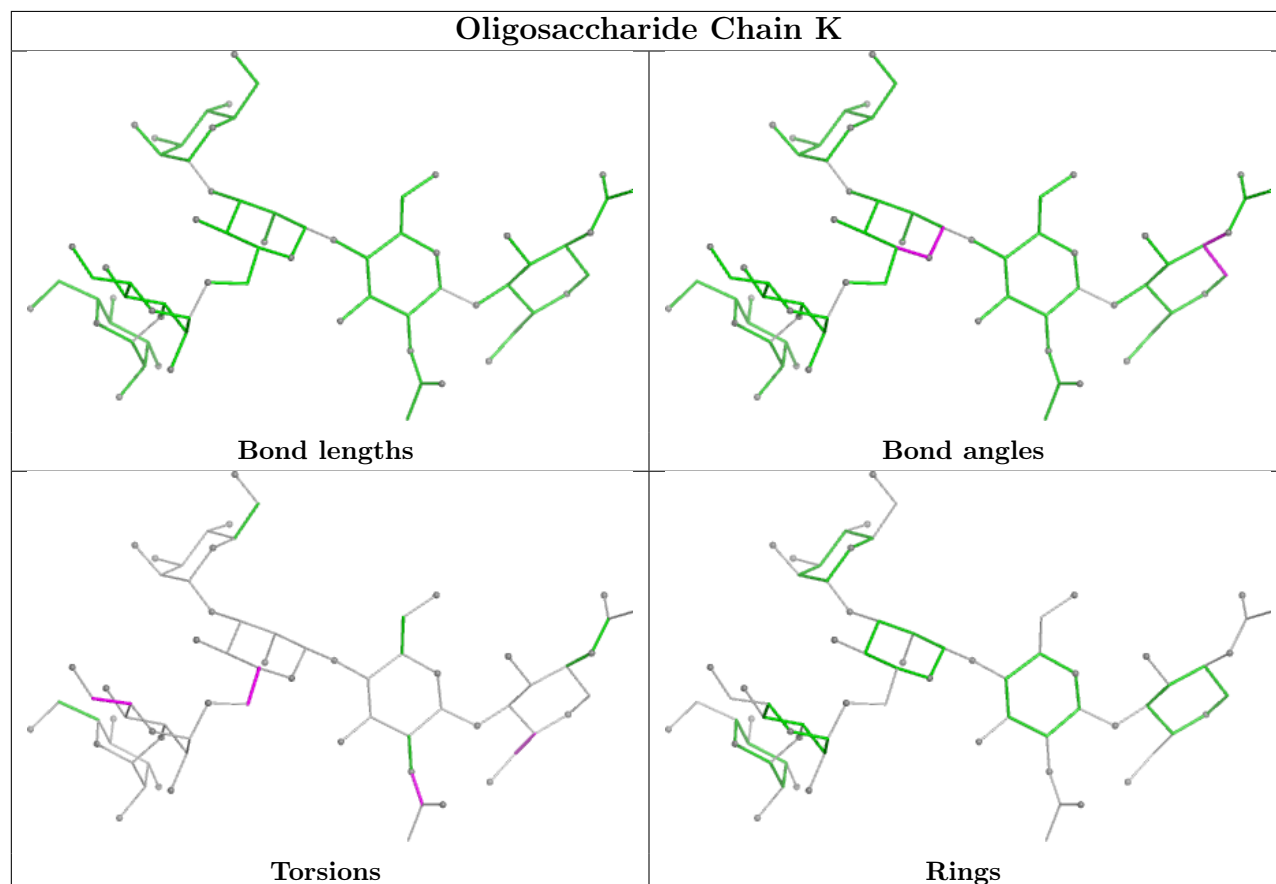
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

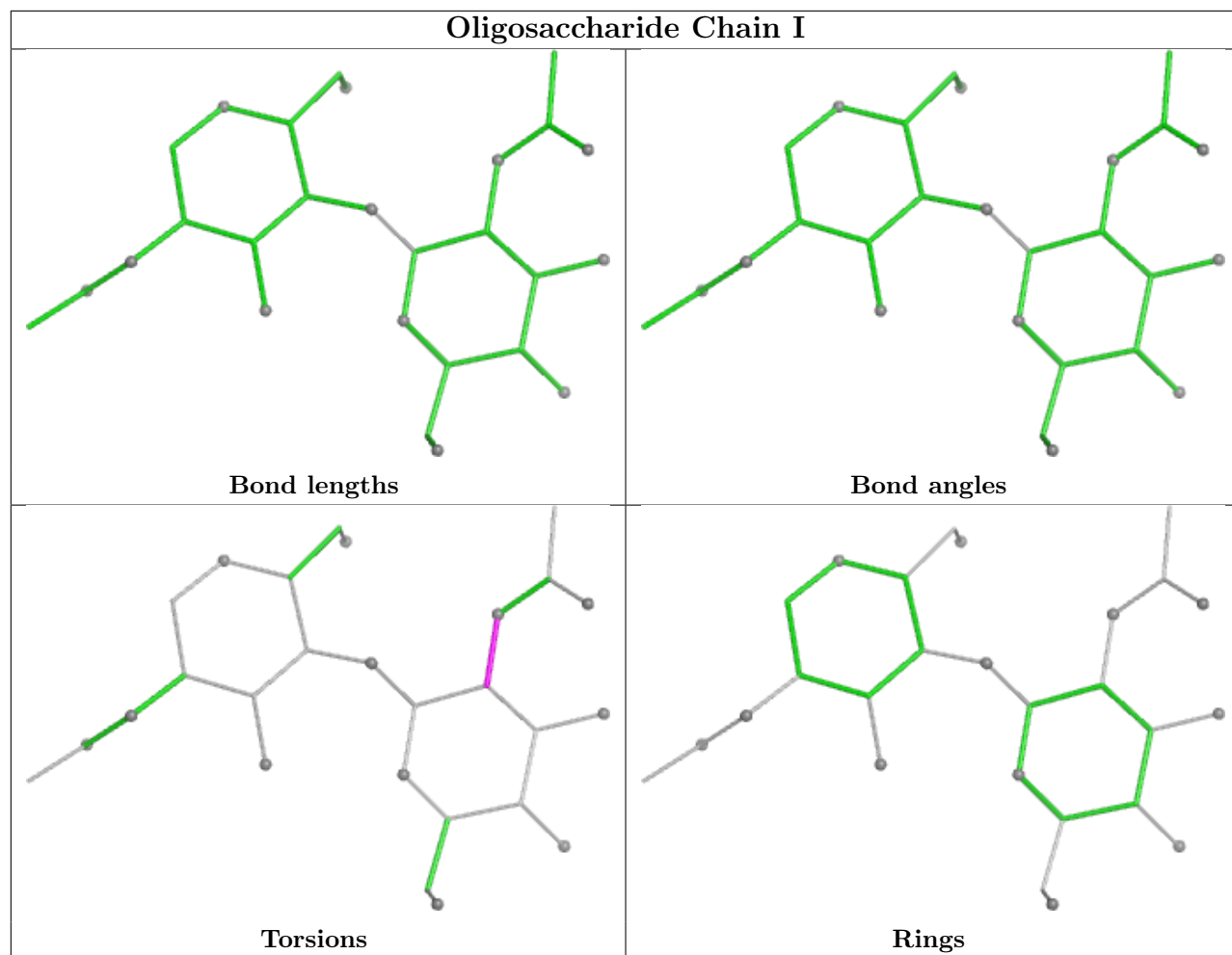


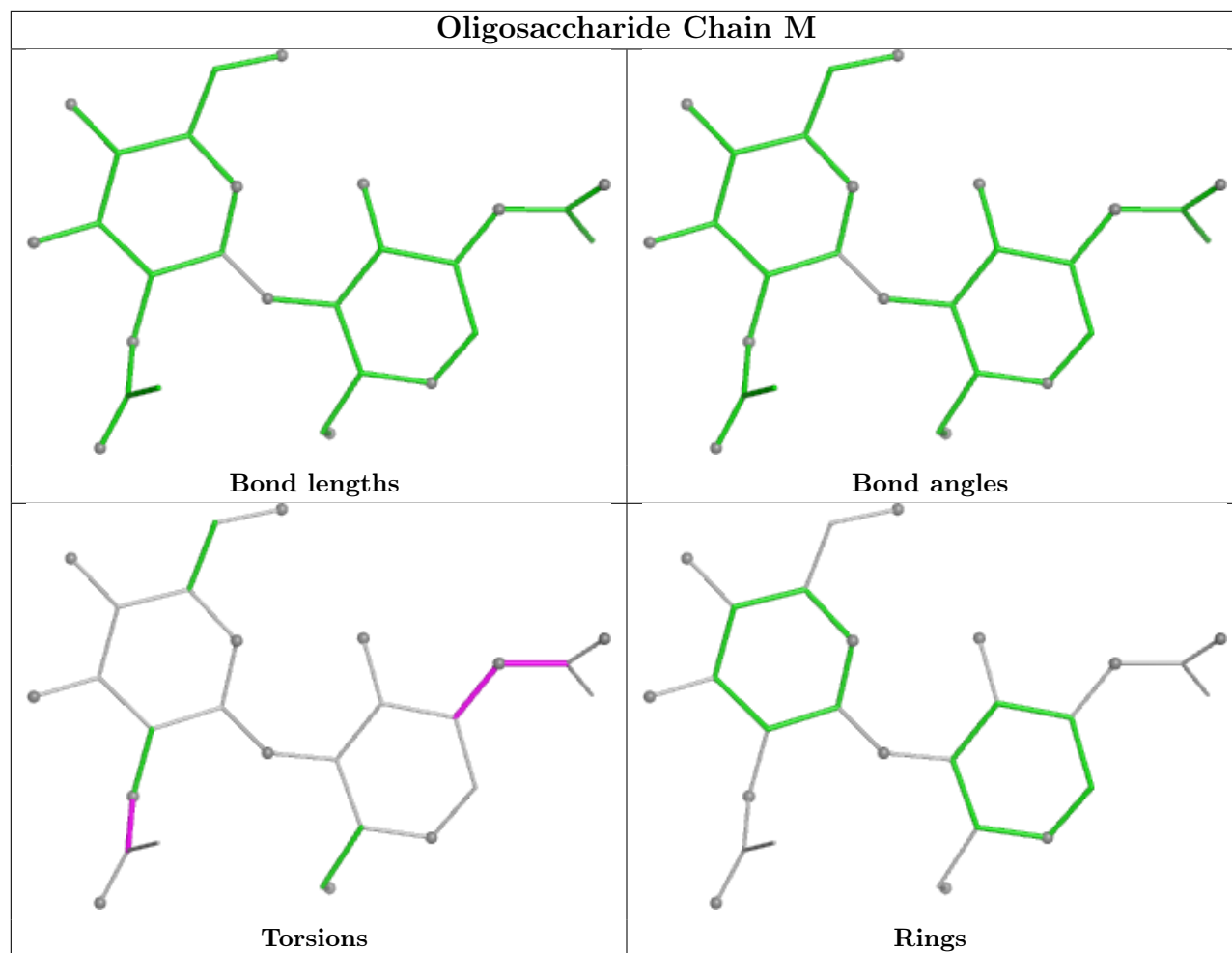
Oligosaccharide Chain G

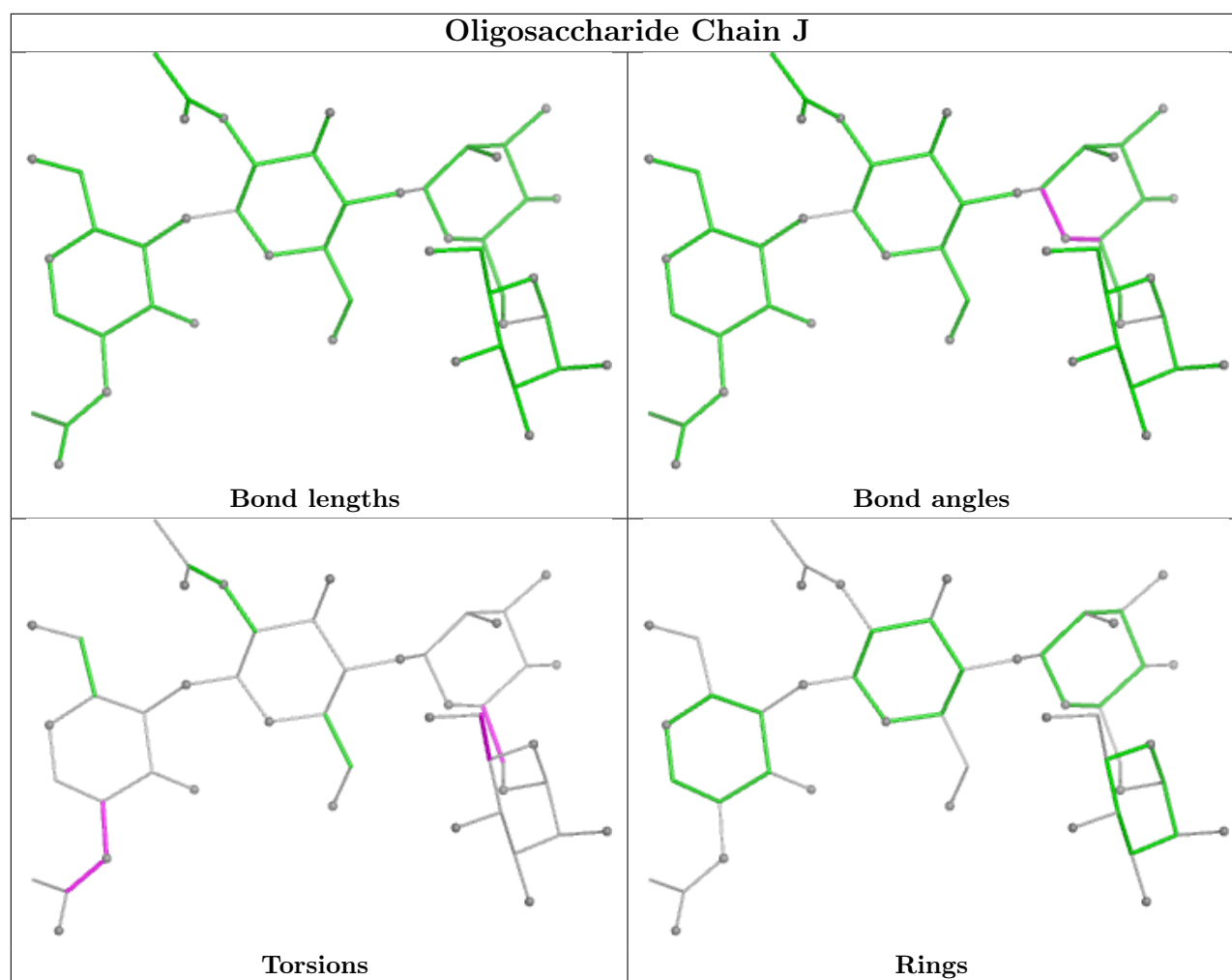


Oligosaccharide Chain K









5.6 Ligand geometry [i](#)

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	$\# Z > 2$
17	D10	D	502	-	9,9,9	0.30	0	8,8,8	0.73	0
12	HEX	A	3904	-	5,5,5	0.30	0	4,4,4	0.59	0
12	HEX	D	503	-	5,5,5	0.32	0	4,4,4	0.49	0
14	R16	B	502	-	15,15,15	0.31	0	14,14,14	0.77	0
16	P1L	C	504	-	21,22,23	1.09	1 (4%)	18,23,25	1.62	3 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
13	PGW	B	501	-	50,50,50	1.06	3 (6%)	53,56,56	0.96	2 (3%)
10	PIO	D	504	-	47,47,47	1.15	6 (12%)	61,65,65	0.85	2 (3%)
12	HEX	B	503	-	5,5,5	0.31	0	4,4,4	0.56	0
16	P1L	C	502	16	21,22,23	1.08	1 (4%)	18,23,25	1.18	3 (16%)
16	P1L	C	503	16	21,22,23	0.66	0	18,23,25	1.53	3 (16%)
11	PLM	A	3902	-	17,17,17	0.56	0	17,17,17	1.15	0
10	PIO	A	3901	-	47,47,47	1.16	6 (12%)	61,65,65	0.96	2 (3%)
12	HEX	A	3905	-	5,5,5	0.32	0	4,4,4	0.51	0
18	CLR	L	601	-	31,31,31	0.16	0	48,48,48	0.28	0
15	NAG	C	501	-	14,14,15	0.76	0	17,19,21	1.46	1 (5%)
12	HEX	A	3903	-	5,5,5	0.30	0	4,4,4	0.57	0
14	R16	D	501	-	15,15,15	0.30	0	14,14,14	0.78	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
17	D10	D	502	-	-	3/7/7/7	-
12	HEX	A	3904	-	-	0/3/3/3	-
12	HEX	D	503	-	-	1/3/3/3	-
14	R16	B	502	-	-	3/13/13/13	-
16	P1L	C	504	-	-	8/20/22/24	-
13	PGW	B	501	-	-	30/55/55/55	-
10	PIO	D	504	-	-	21/44/68/68	0/1/1/1
12	HEX	B	503	-	-	0/3/3/3	-
16	P1L	C	502	16	-	7/20/22/24	-
16	P1L	C	503	16	-	4/20/22/24	-
11	PLM	A	3902	-	-	8/15/15/15	-
10	PIO	A	3901	-	-	10/44/68/68	0/1/1/1
12	HEX	A	3905	-	-	0/3/3/3	-
18	CLR	L	601	-	-	0/10/68/68	0/4/4/4
15	NAG	C	501	-	-	2/6/23/26	0/1/1/1
12	HEX	A	3903	-	-	0/3/3/3	-
14	R16	D	501	-	-	5/13/13/13	-

All (17) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	C	504	P1L	O-C	4.14	1.36	1.19
16	C	502	P1L	O-C	4.13	1.36	1.19
13	B	501	PGW	O03-C19	3.13	1.42	1.33
10	A	3901	PIO	P5-O5	3.05	1.65	1.59
10	A	3901	PIO	P4-O4	3.00	1.65	1.59
10	D	504	PIO	P5-O5	2.89	1.64	1.59
10	D	504	PIO	P4-O4	2.87	1.64	1.59
10	D	504	PIO	O2C-C2C	-2.64	1.40	1.46
13	B	501	PGW	O01-C02	-2.64	1.40	1.46
13	B	501	PGW	O01-C1	2.42	1.41	1.34
10	A	3901	PIO	O3C-C1B	2.37	1.40	1.33
10	A	3901	PIO	O2C-C1A	2.35	1.40	1.34
10	A	3901	PIO	O2C-C2C	-2.32	1.40	1.46
10	D	504	PIO	O3C-C1B	2.30	1.40	1.33
10	D	504	PIO	O3C-C3C	-2.19	1.40	1.45
10	A	3901	PIO	O3C-C3C	-2.13	1.40	1.45
10	D	504	PIO	O2C-C1A	2.11	1.40	1.34

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	C	504	P1L	C8-C7-SG	-4.97	107.67	113.46
15	C	501	NAG	C2-N2-C7	4.66	129.54	122.90
10	A	3901	PIO	O2C-C1A-C2A	4.55	121.32	111.50
16	C	503	P1L	C8-C7-SG	-4.18	108.59	113.46
13	B	501	PGW	O01-C1-C2	3.96	120.03	111.50
16	C	503	P1L	CB-SG-C7	3.80	106.15	100.84
16	C	504	P1L	CB-SG-C7	3.71	106.02	100.84
10	D	504	PIO	O2C-C1A-C2A	3.60	119.25	111.50
16	C	502	P1L	C8-C7-SG	-3.16	109.78	113.46
16	C	502	P1L	CB-SG-C7	2.77	104.72	100.84
10	D	504	PIO	O3C-C1B-C2B	2.67	120.30	111.91
16	C	503	P1L	O7-C7-SG	2.59	125.98	122.61
13	B	501	PGW	O03-C19-C20	2.49	119.72	111.91
10	A	3901	PIO	O3C-C1B-C2B	2.48	119.71	111.91
16	C	504	P1L	O7-C7-SG	2.35	125.67	122.61
16	C	502	P1L	O7-C7-SG	2.18	125.44	122.61

There are no chirality outliers.

All (102) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	A	3901	PIO	C5-O5-P5-O53

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Mol	Chain	Res	Type	Atoms
10	A	3901	PIO	O1A-C1A-O2C-C2C
10	D	504	PIO	C1-O1-P1-O11
10	D	504	PIO	C5-O5-P5-O53
13	B	501	PGW	C03-O11-P-O12
13	B	501	PGW	C03-O11-P-O13
13	B	501	PGW	C03-O11-P-O14
16	C	502	P1L	N-CA-CB-SG
16	C	502	P1L	C-CA-CB-SG
16	C	503	P1L	C8-C7-SG-CB
16	C	504	P1L	N-CA-CB-SG
16	C	504	P1L	C-CA-CB-SG
16	C	504	P1L	C8-C7-SG-CB
10	A	3901	PIO	C2B-C1B-O3C-C3C
10	A	3901	PIO	O1B-C1B-O3C-C3C
10	A	3901	PIO	C2A-C1A-O2C-C2C
13	B	501	PGW	C1-C2-C3-C4
15	C	501	NAG	C8-C7-N2-C2
15	C	501	NAG	O7-C7-N2-C2
13	B	501	PGW	C20-C19-O03-C01
13	B	501	PGW	O04-C19-O03-C01
13	B	501	PGW	C3-C4-C5-C6
10	D	504	PIO	C1C-O13-P1-O1
12	D	503	HEX	C2-C3-C4-C5
14	B	502	R16	C29-C30-C31-C32
16	C	504	P1L	C12-C13-C14-C15
13	B	501	PGW	C04-C05-CAD-OAE
16	C	502	P1L	C10-C11-C12-C13
11	A	3902	PLM	C2-C3-C4-C5
13	B	501	PGW	C2-C1-O01-C02
13	B	501	PGW	O02-C1-O01-C02
11	A	3902	PLM	CC-CD-CE-CF
13	B	501	PGW	C6-C7-C8-C9
10	D	504	PIO	O1A-C1A-O2C-C2C
10	D	504	PIO	C2A-C1A-O2C-C2C
16	C	503	P1L	C11-C10-C9-C8
13	B	501	PGW	C07-C08-C09-C11
13	B	501	PGW	C16-C15-C27-C26
11	A	3902	PLM	CB-CC-CD-CE
16	C	504	P1L	C9-C10-C11-C12
13	B	501	PGW	C23-C24-C25-C26
13	B	501	PGW	C01-C02-C03-O11
10	A	3901	PIO	C1B-C2B-C3B-C4B

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Mol	Chain	Res	Type	Atoms
10	D	504	PIO	C1-O1-P1-O13
10	D	504	PIO	C3B-C4B-C5B-C6B
13	B	501	PGW	O03-C01-C02-C03
10	D	504	PIO	C1-O1-P1-O12
16	C	502	P1L	C12-C13-C14-C15
17	D	502	D10	C6-C7-C8-C9
10	D	504	PIO	O13-C1C-C2C-O2C
10	D	504	PIO	C2B-C1B-O3C-C3C
10	D	504	PIO	O2C-C2C-C3C-O3C
13	B	501	PGW	O03-C01-C02-O01
13	B	501	PGW	C05-C04-O12-P
16	C	502	P1L	O7-C7-SG-CB
16	C	503	P1L	O7-C7-SG-CB
10	D	504	PIO	O13-C1C-C2C-C3C
14	B	502	R16	C38-C39-C40-C41
13	B	501	PGW	C19-C20-C21-C22
14	D	501	R16	C36-C37-C38-C39
14	D	501	R16	C33-C34-C35-C36
13	B	501	PGW	C09-C11-C12-C13
16	C	502	P1L	C15-C16-C17-C18
10	D	504	PIO	O1B-C1B-O3C-C3C
14	D	501	R16	C35-C36-C37-C38
14	B	502	R16	C37-C38-C39-C40
10	A	3901	PIO	C3B-C4B-C5B-C6B
10	D	504	PIO	C1C-O13-P1-O11
10	D	504	PIO	C1C-O13-P1-O12
13	B	501	PGW	C20-C21-C22-C23
13	B	501	PGW	O01-C02-C03-O11
17	D	502	D10	C4-C5-C6-C7
17	D	502	D10	C5-C6-C7-C8
10	D	504	PIO	C3A-C4A-C5A-C6A
11	A	3902	PLM	C5-C6-C7-C8
14	D	501	R16	C38-C39-C40-C41
13	B	501	PGW	C04-O12-P-O11
13	B	501	PGW	C15-C16-C17-C18
13	B	501	PGW	OAF-C05-CAD-OAE
14	D	501	R16	C34-C35-C36-C37
13	B	501	PGW	C16-C17-C18-C28
10	A	3901	PIO	C3A-C4A-C5A-C6A
16	C	503	P1L	C10-C11-C12-C13
13	B	501	PGW	C07-C06-C10-C9
16	C	502	P1L	C11-C12-C13-C14

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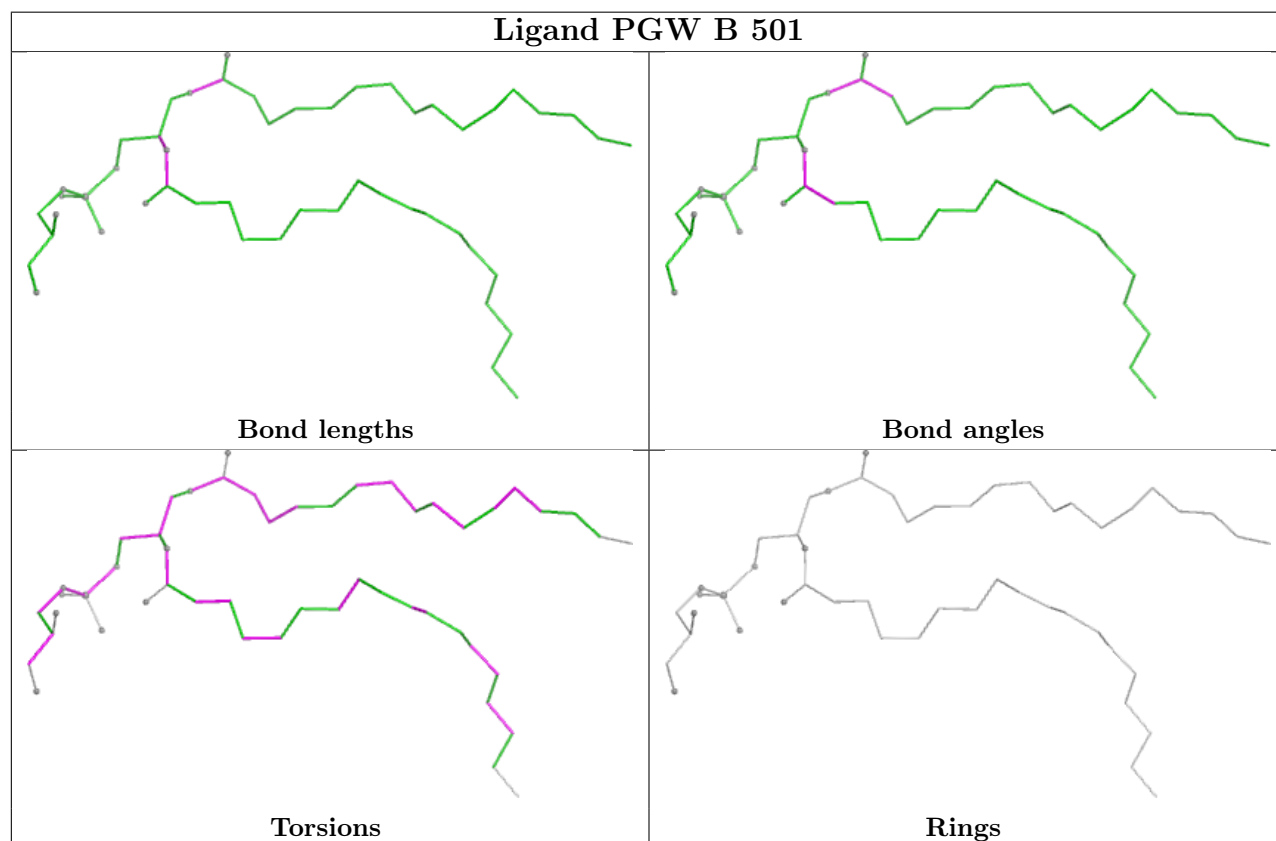
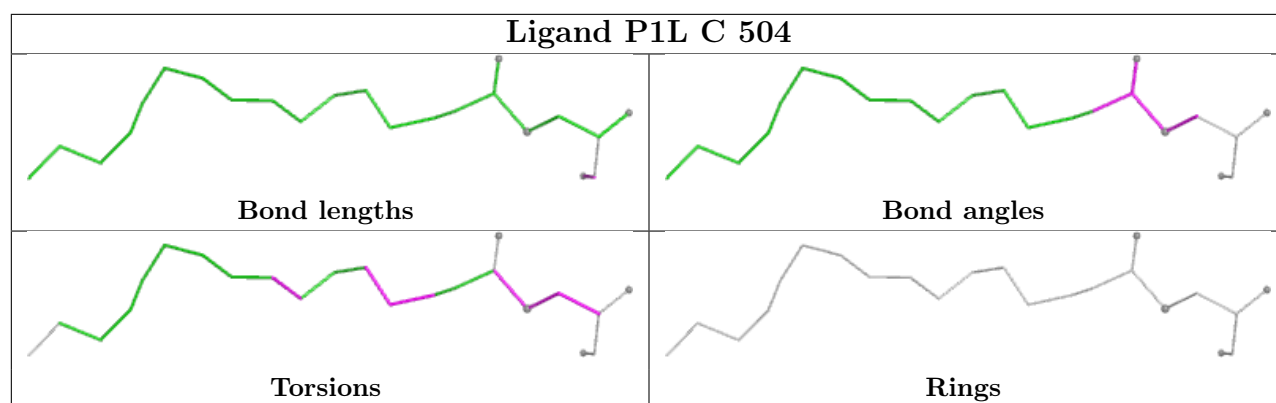
Mol	Chain	Res	Type	Atoms
11	A	3902	PLM	C7-C8-C9-CA
10	D	504	PIO	C5B-C6B-C7B-C8B
11	A	3902	PLM	O1-C1-C2-C3
11	A	3902	PLM	O2-C1-C2-C3
16	C	504	P1L	O7-C7-SG-CB
16	C	504	P1L	C11-C10-C9-C8
10	D	504	PIO	C1B-C2B-C3B-C4B
11	A	3902	PLM	CA-CB-CC-CD
13	B	501	PGW	O03-C19-C20-C21
10	D	504	PIO	C1C-C2C-C3C-O3C
10	D	504	PIO	O2C-C1A-C2A-C3A
13	B	501	PGW	C24-C25-C26-C27
16	C	504	P1L	CA-CB-SG-C7
13	B	501	PGW	O04-C19-C20-C21
10	A	3901	PIO	C1C-O13-P1-O12
10	D	504	PIO	O1A-C1A-C2A-C3A
10	A	3901	PIO	C2B-C3B-C4B-C5B

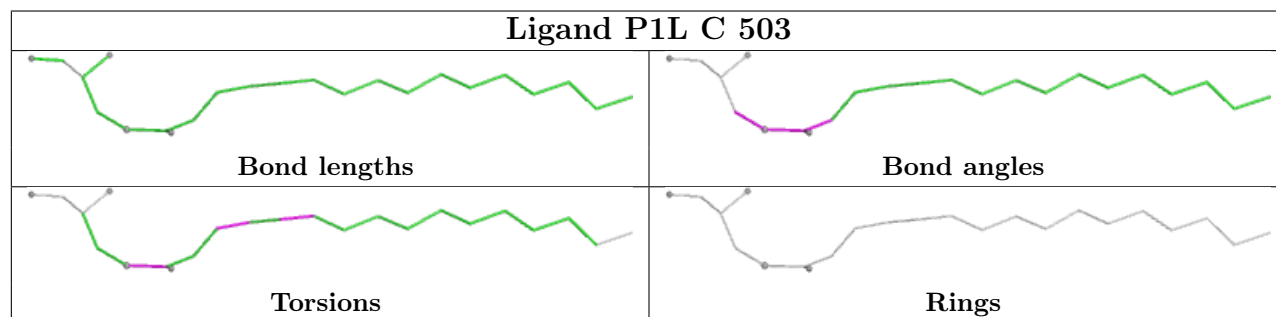
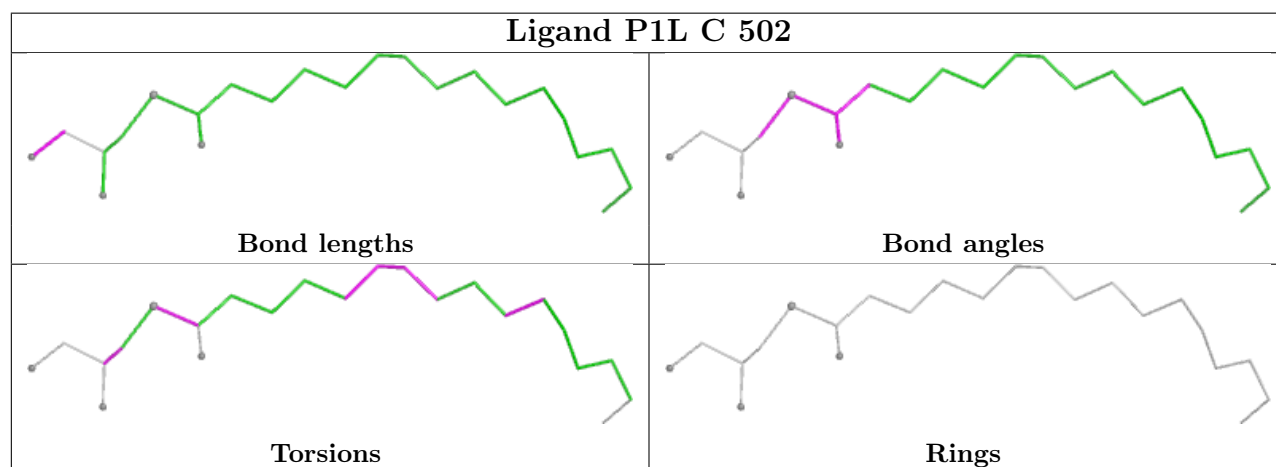
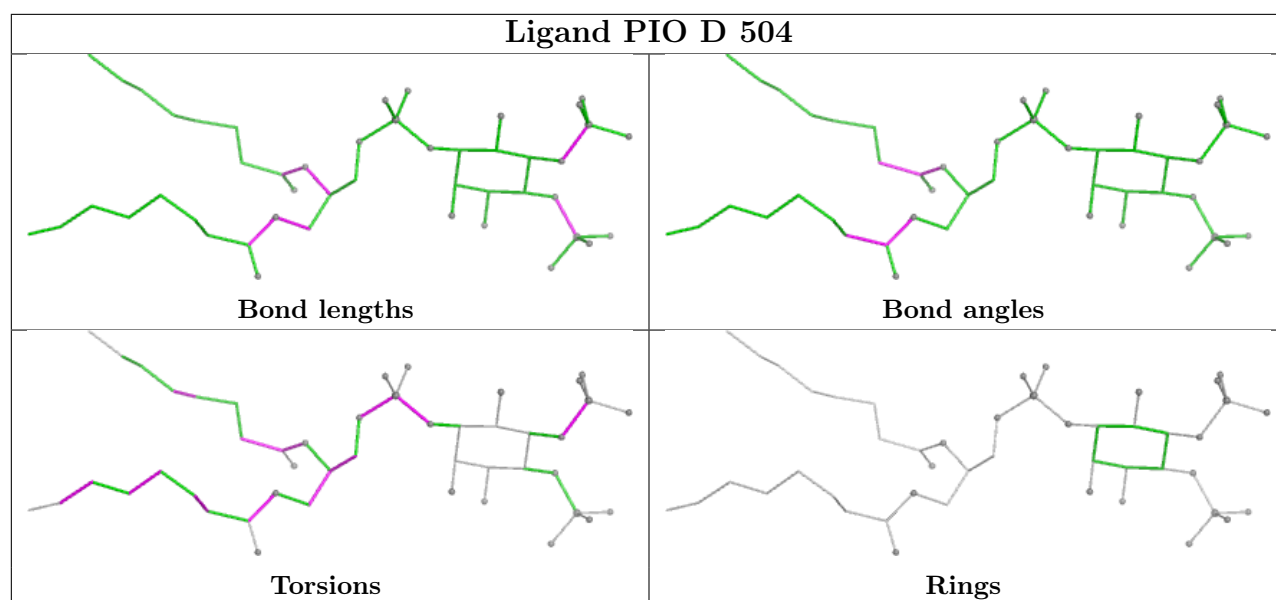
There are no ring outliers.

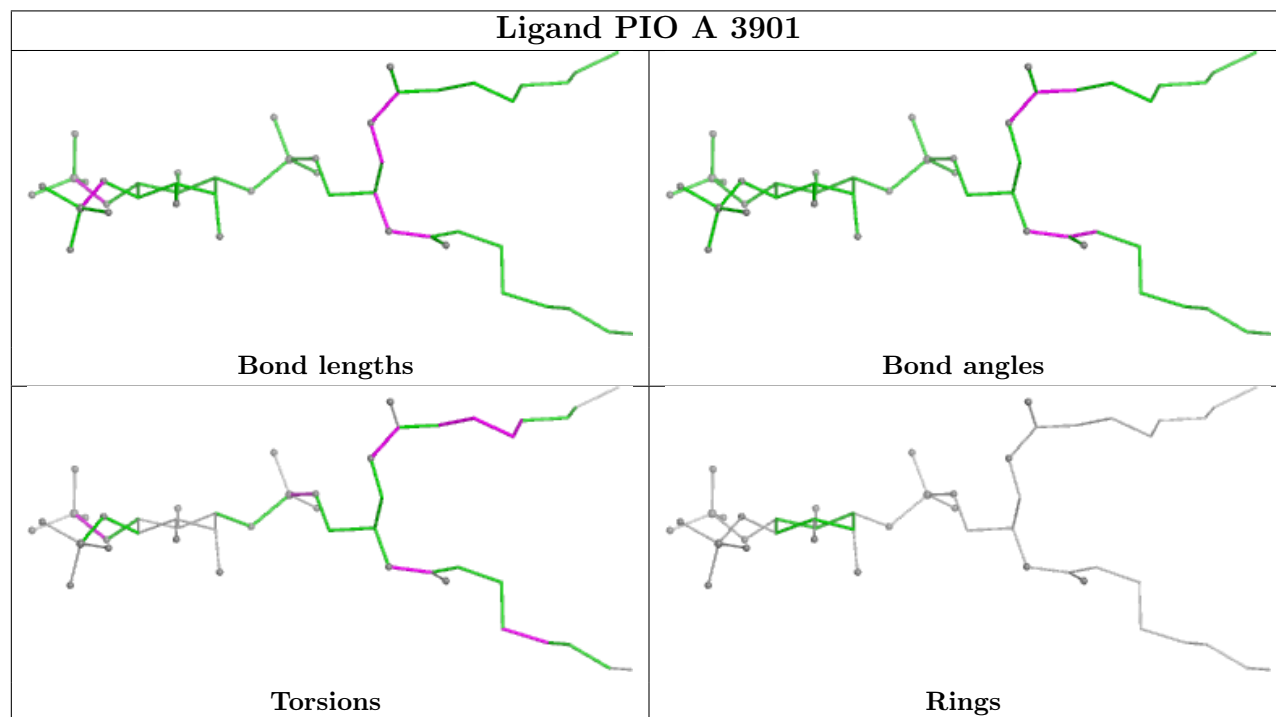
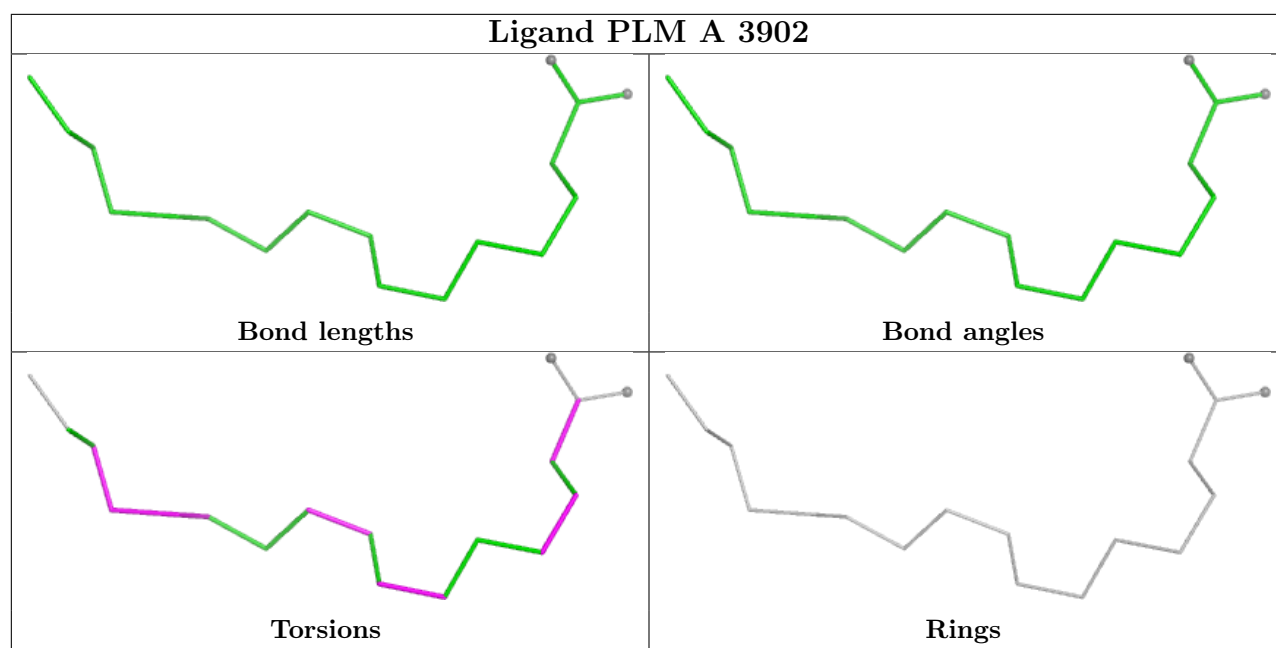
4 monomers are involved in 7 short contacts:

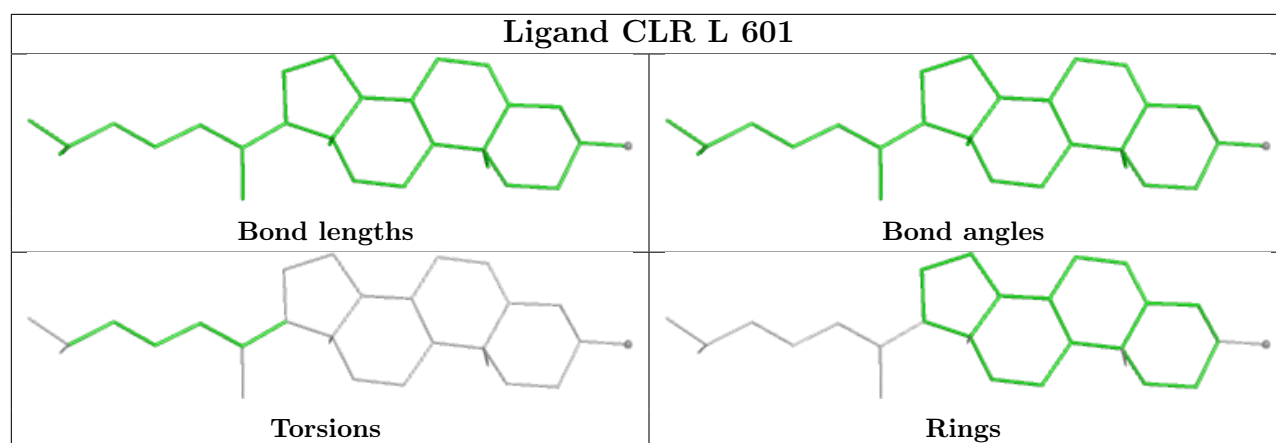
Mol	Chain	Res	Type	Clashes	Symm-Clashes
16	C	504	P1L	1	0
13	B	501	PGW	2	0
11	A	3902	PLM	3	0
18	L	601	CLR	1	0

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









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.

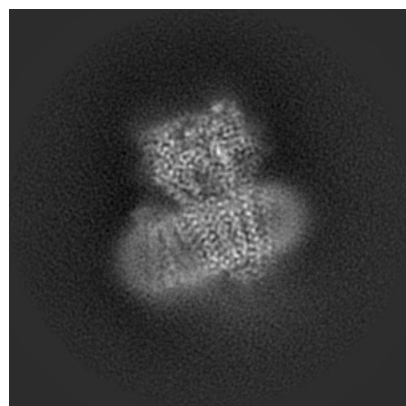
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-53949. These allow visual inspection of the internal detail of the map and identification of artifacts.

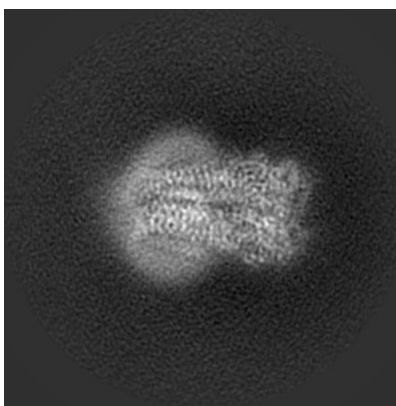
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

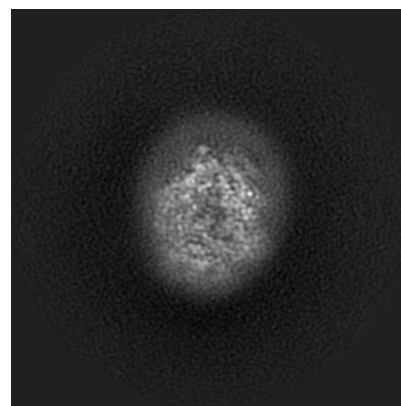
6.1.1 Primary map



X

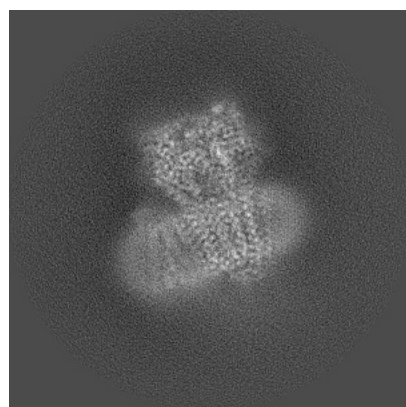


Y

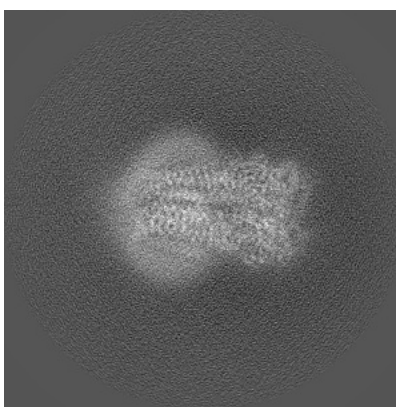


Z

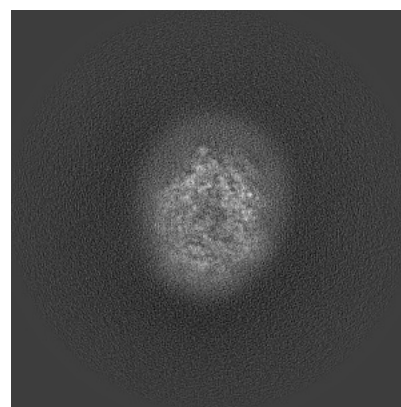
6.1.2 Raw map



X



Y

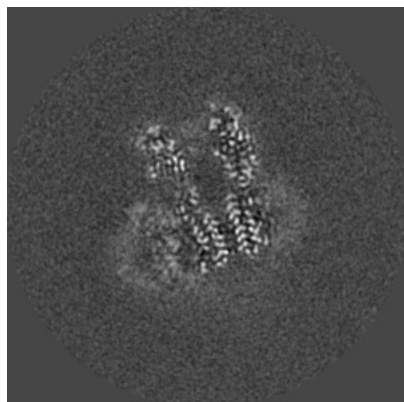


Z

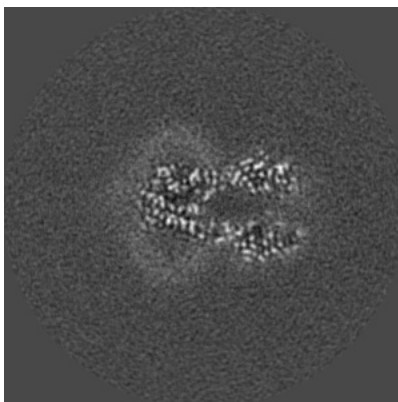
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

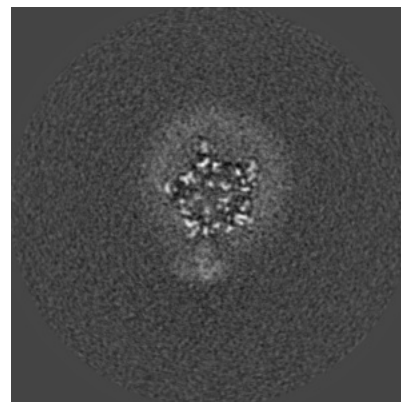
6.2.1 Primary map



X Index: 180

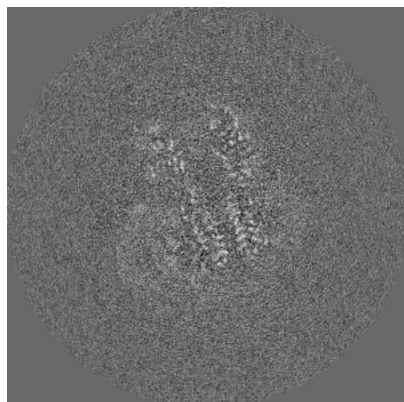


Y Index: 180

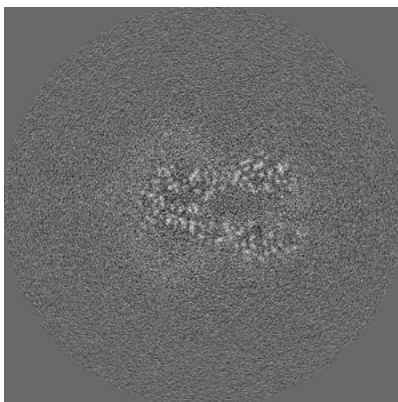


Z Index: 180

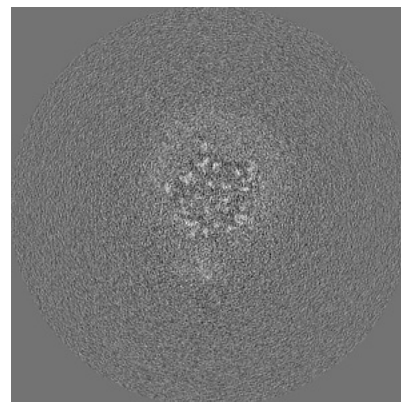
6.2.2 Raw map



X Index: 180



Y Index: 180

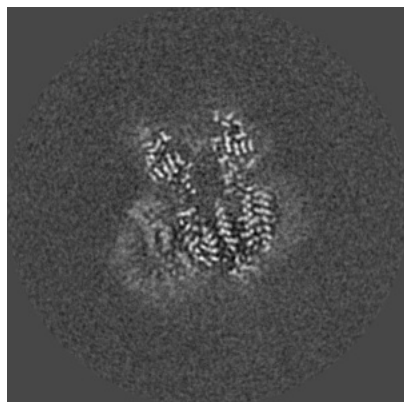


Z Index: 180

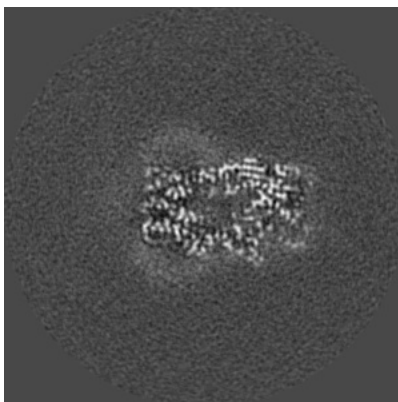
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

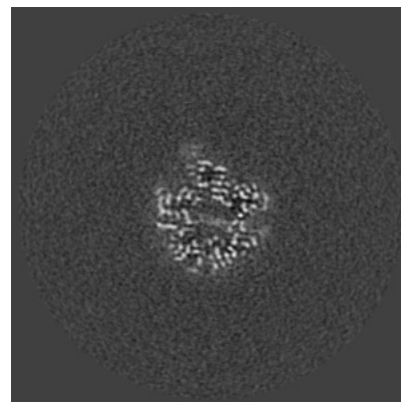
6.3.1 Primary map



X Index: 170

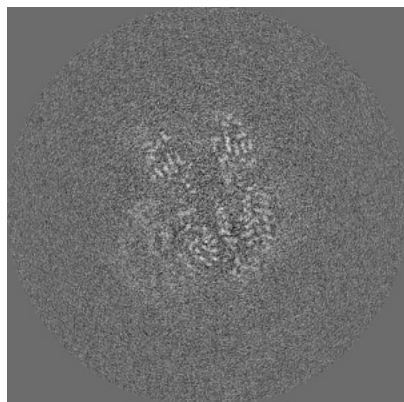


Y Index: 194

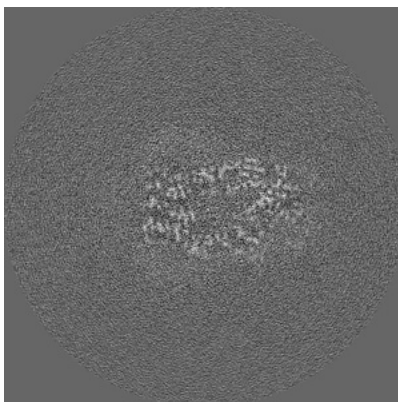


Z Index: 231

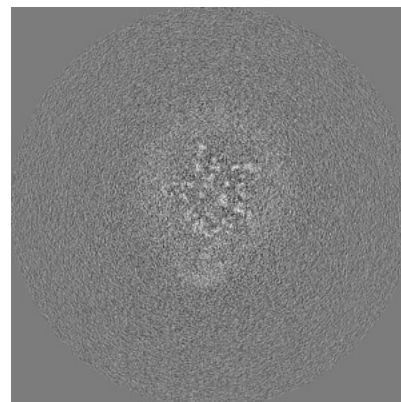
6.3.2 Raw map



X Index: 170



Y Index: 195

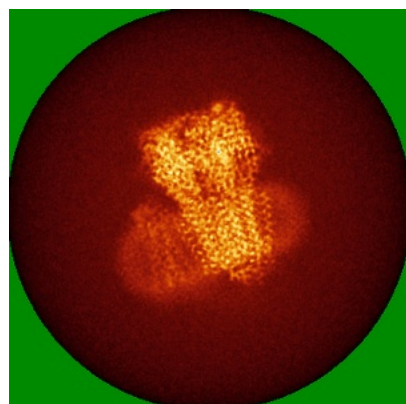


Z Index: 175

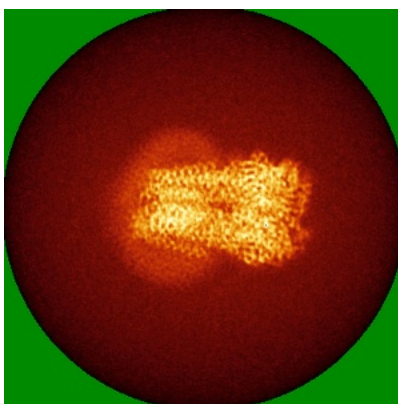
The images above show the largest variance slices of the map in three orthogonal directions.

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

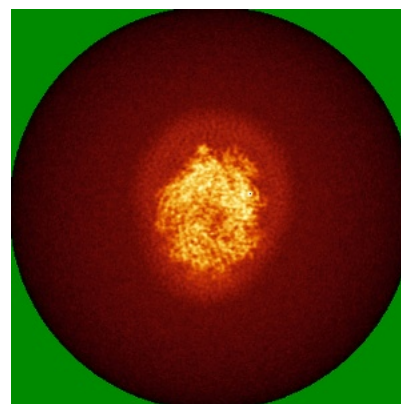
6.4.1 Primary map



X

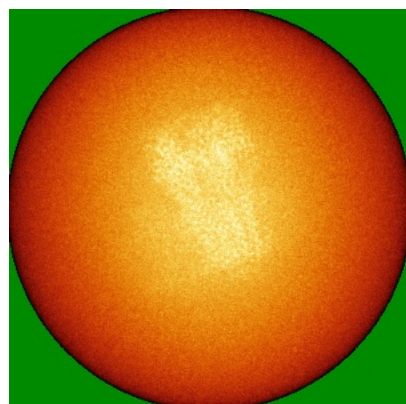


Y

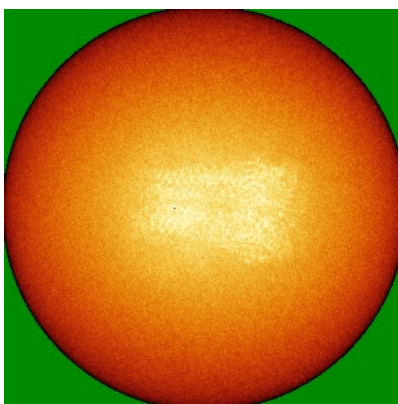


Z

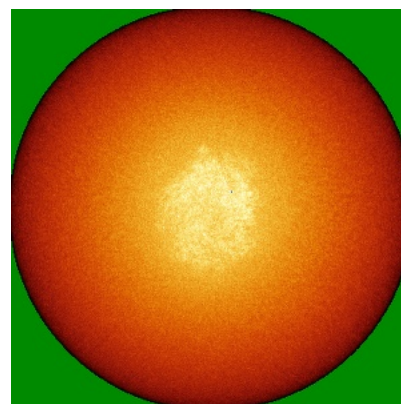
6.4.2 Raw map



X



Y

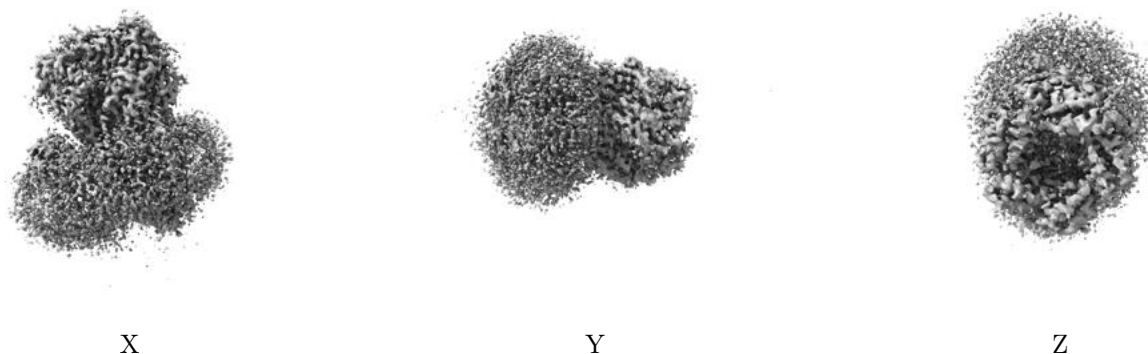


Z

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.

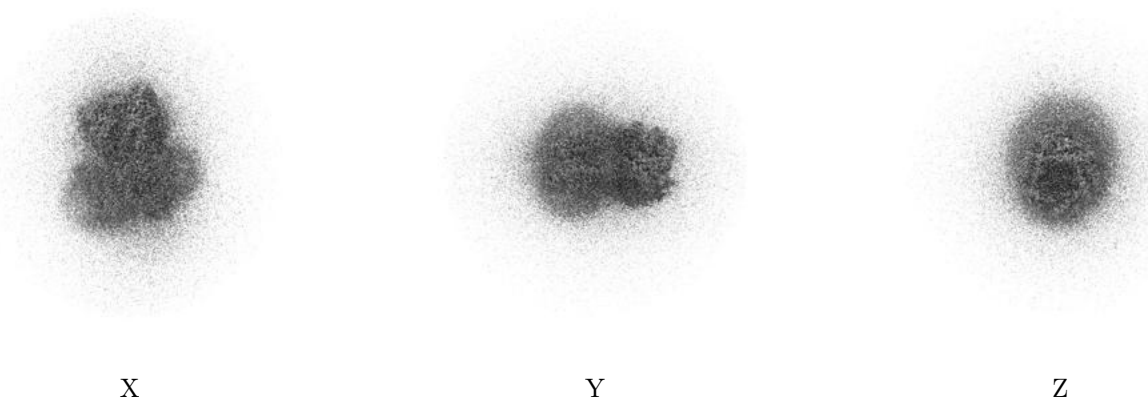
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0034. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

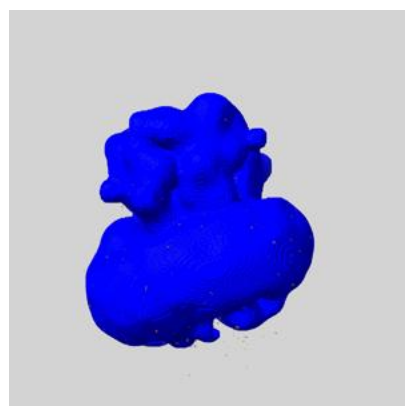
6.6 Mask visualisation [i](#)

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

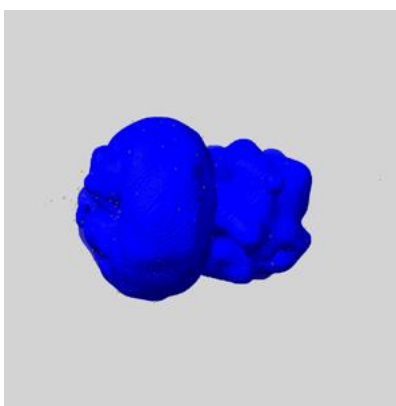
A mask typically either:

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

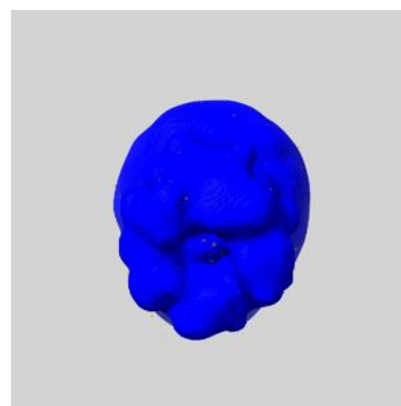
6.6.1 emd_53949_msk_1.map [i](#)



X



Y

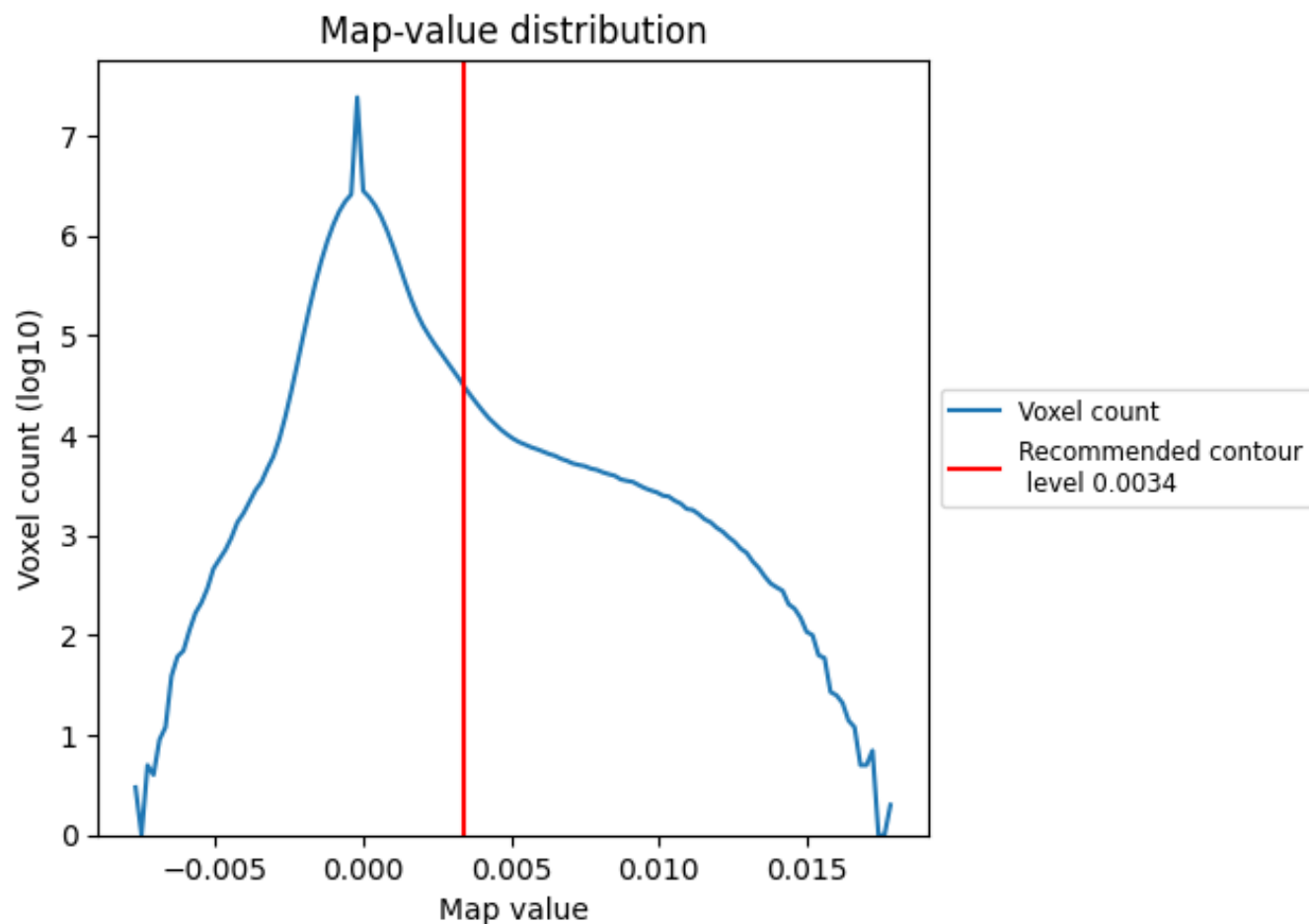


Z

7 Map analysis [i](#)

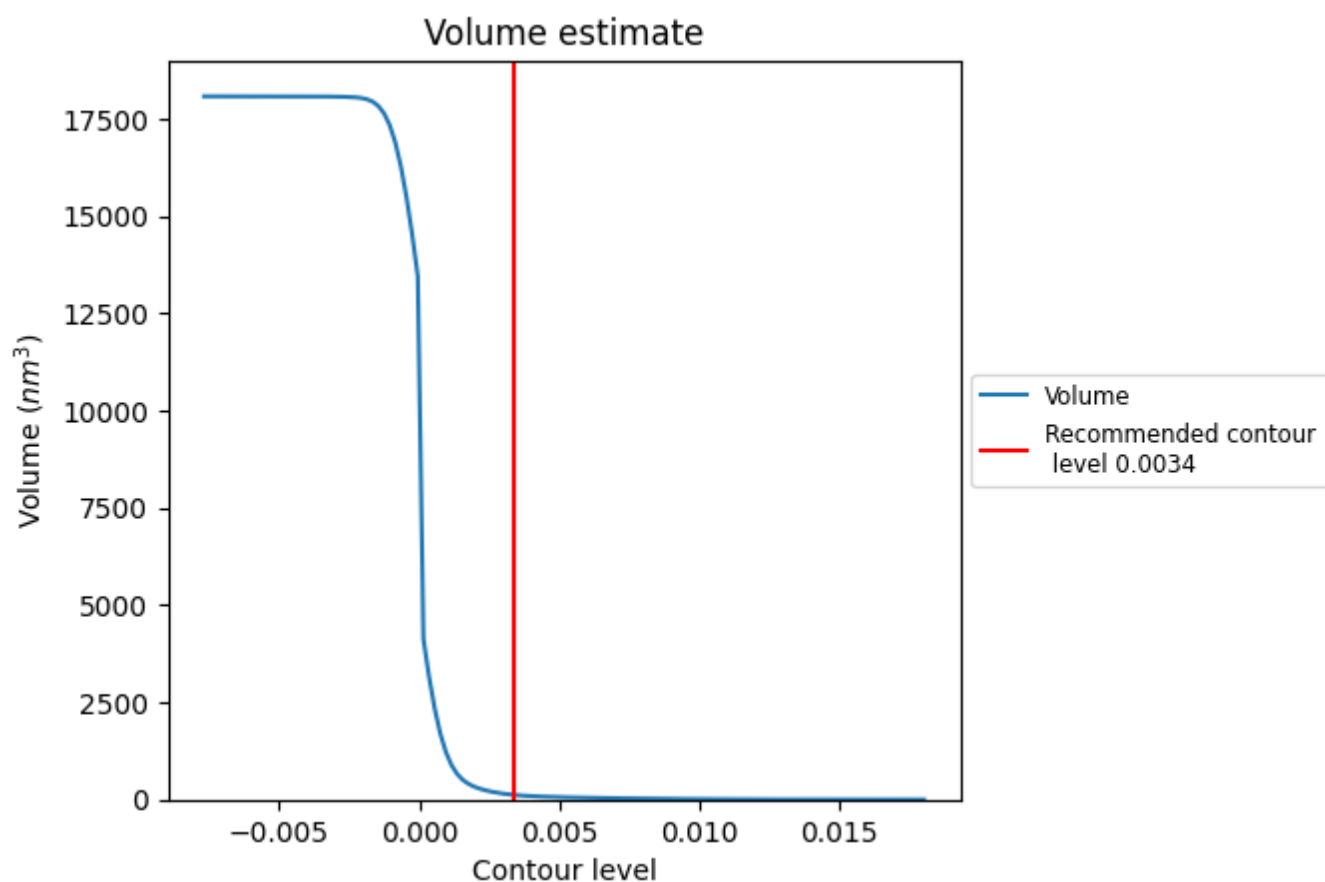
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

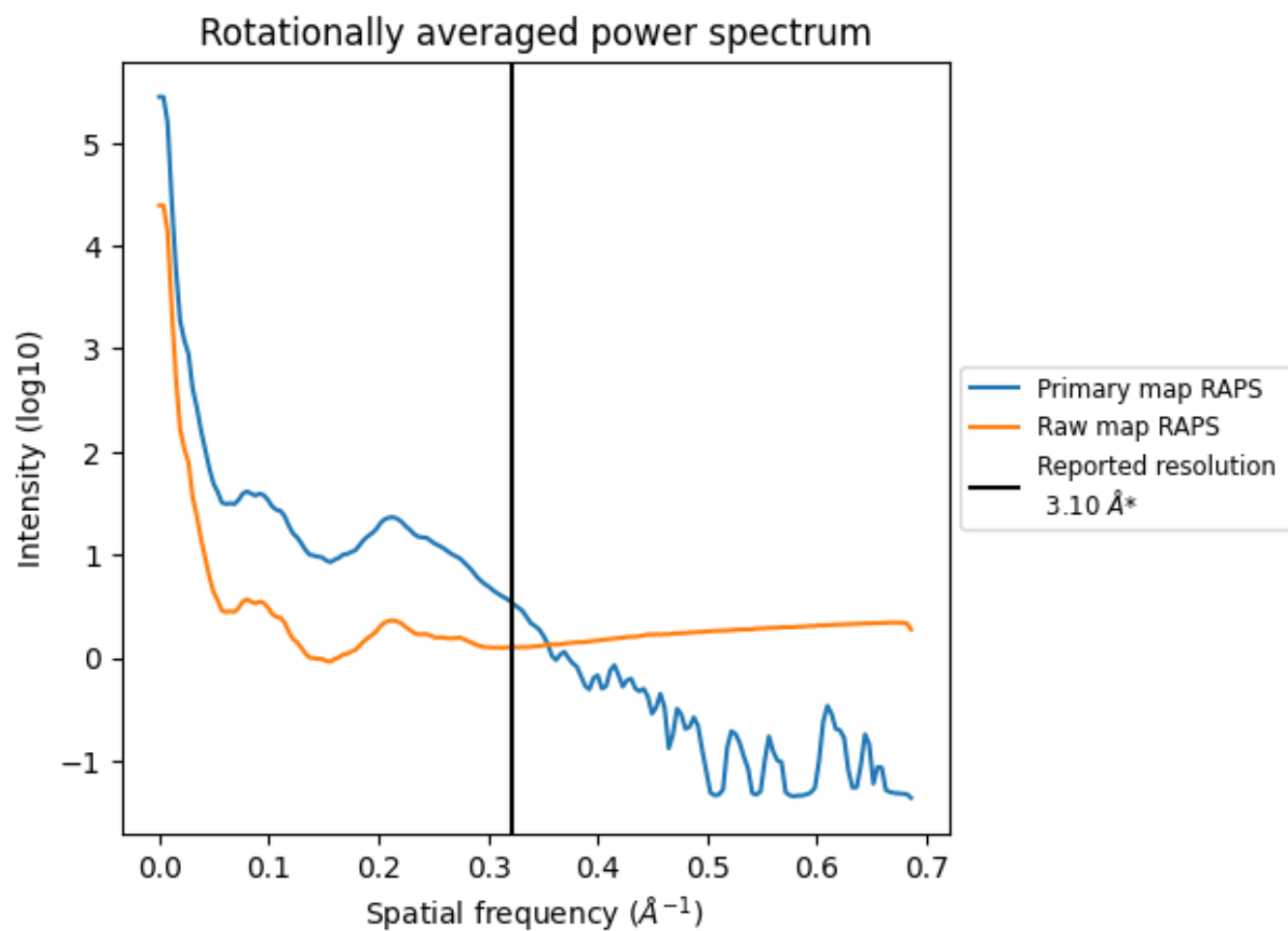
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 119 nm³; this corresponds to an approximate mass of 108 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 ⓘ

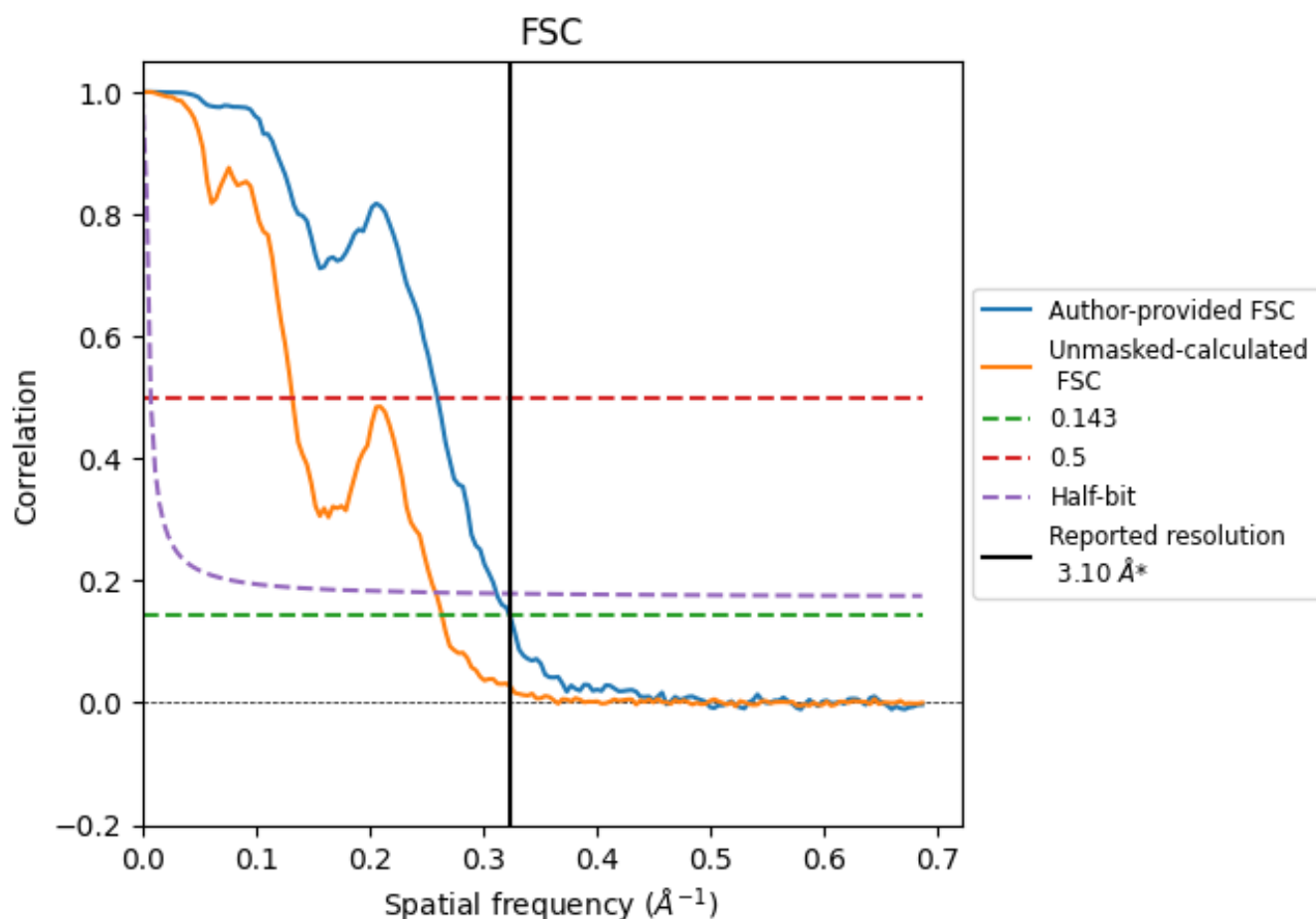


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

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 \AA^{-1}

8.2 Resolution estimates [i](#)

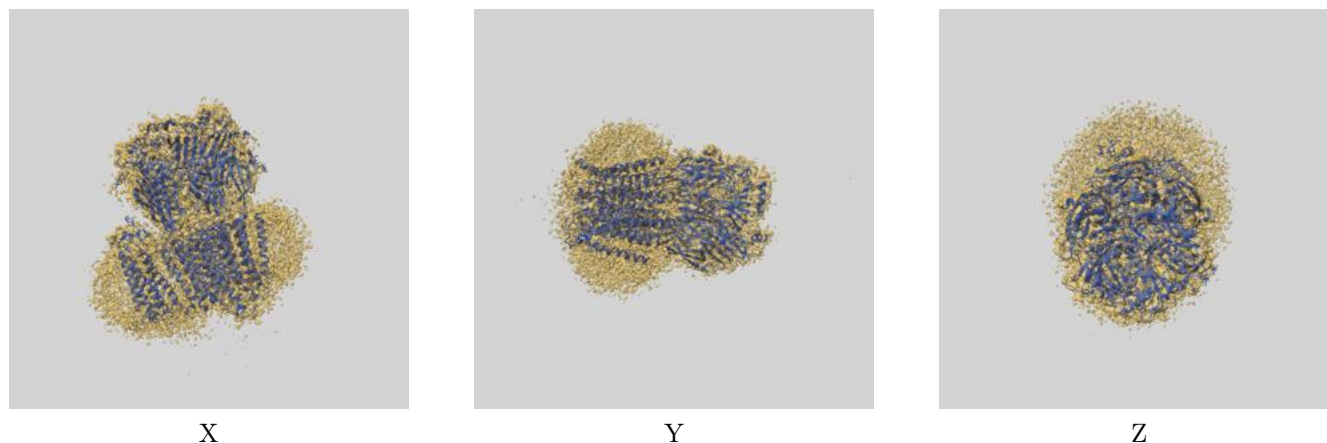
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.10	3.85	3.21
Unmasked-calculated*	3.79	7.58	3.88

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.79 differs from the reported value 3.1 by more than 10 %

9 Map-model fit [i](#)

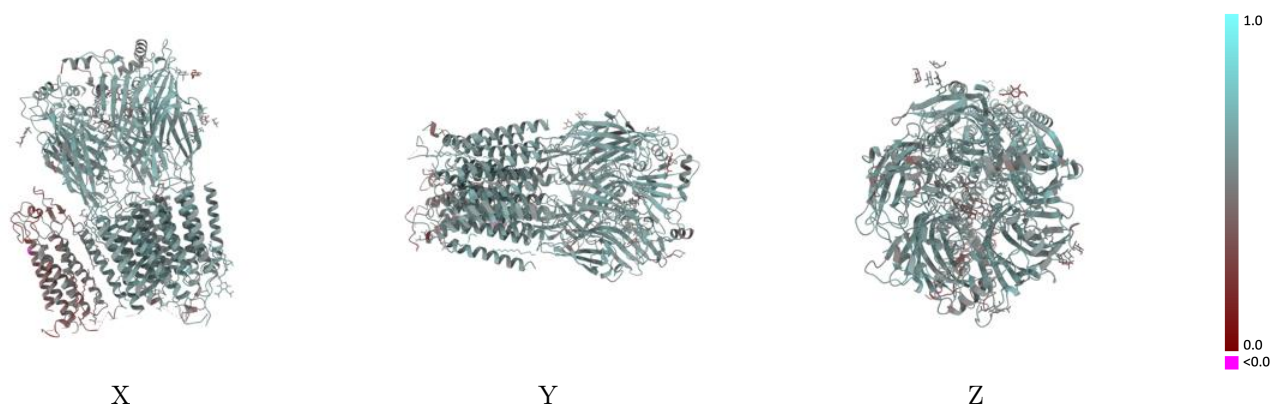
This section contains information regarding the fit between EMDB map EMD-53949 and PDB model 9RGE. Per-residue inclusion information can be found in section 3 on page 12.

9.1 Map-model overlay [i](#)



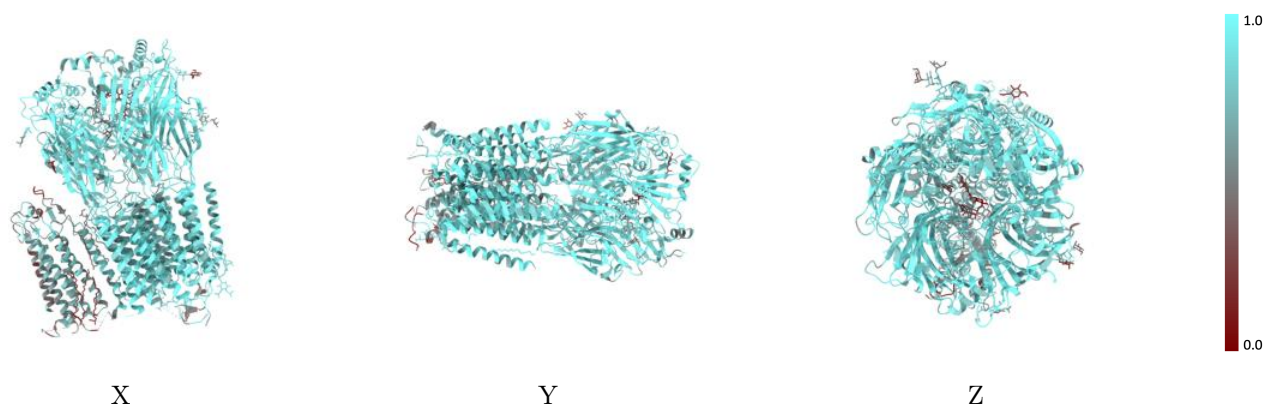
The images above show the 3D surface view of the map at the recommended contour level 0.0034 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



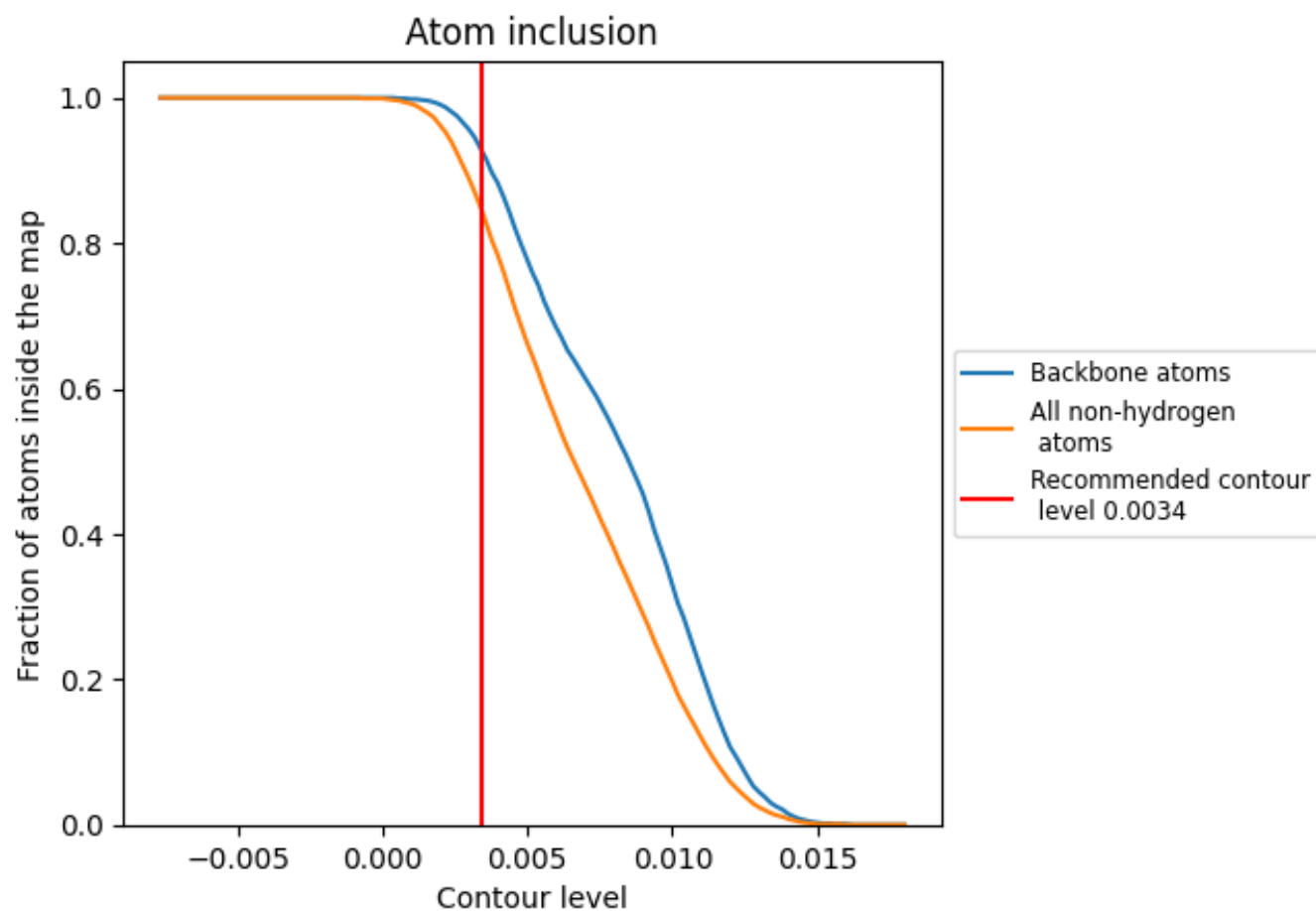
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0034).





























9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 85% 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 (0.0034) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8470	 0.5550
A	 0.8920	 0.5790
B	 0.8910	 0.5830
C	 0.8670	 0.5690
D	 0.8640	 0.5650
E	 0.8990	 0.5780
F	 0.5340	 0.4800
G	 0.6810	 0.4860
H	 0.4790	 0.3380
I	 0.6790	 0.4450
J	 0.2000	 0.3900
K	 0.7080	 0.4870
L	 0.6320	 0.4190
M	 0.5000	 0.3650

