



wwPDB EM Validation Summary Report ⓘ

Nov 22, 2022 – 12:48 AM EST

PDB ID : 7TFO
EMDB ID : EMD-25878
Title : Cryo-EM structure of HIV-1 Env trimer BG505 SOSIP.664 in complex with CD4bs antibody Ab1573
Authors : Yang, Z.; Bjorkman, P.J.
Deposited on : 2022-01-06
Resolution : 4.10 Å(reported)

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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.dev43
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.9
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.31.3

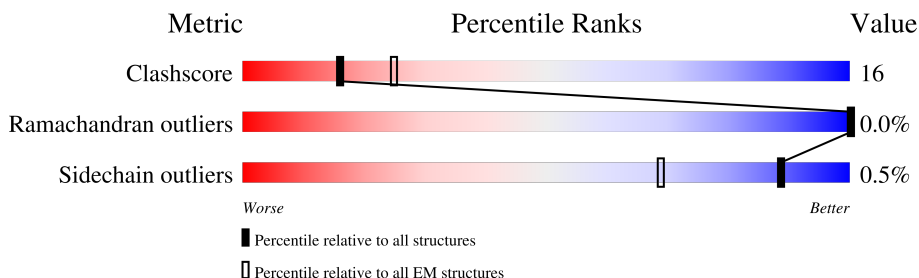
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.10 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)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

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	481	
1	B	481	
1	C	481	
2	H	228	
2	I	228	
2	P	228	
3	J	216	
3	L	216	

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Mol	Chain	Length	Quality of chain
3	Q	216	
4	X	153	
4	Y	153	
4	Z	153	
5	D	2	
6	E	3	
6	F	3	
6	K	3	
7	G	5	

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 17774 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 Envelope glycoprotein BG505 SOSIP.664 - gp120.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	421	Total	C	N	O	S	0	0
			3245	2058	568	593	26		
1	B	411	Total	C	N	O	S	0	0
			3175	2007	552	593	23		
1	C	404	Total	C	N	O	S	0	0
			3110	1964	543	576	27		

There are 18 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	64	LYS	GLU	conflict	UNP A0A6H1VH54
A	375	SER	TYR	conflict	UNP A0A6H1VH54
A	501	CYS	ALA	conflict	UNP A0A6H1VH54
A	509	ARG	GLU	conflict	UNP A0A6H1VH54
A	512	ARG	-	expression tag	UNP A0A6H1VH54
A	513	ARG	-	expression tag	UNP A0A6H1VH54
B	64	LYS	GLU	conflict	UNP A0A6H1VH54
B	375	SER	TYR	conflict	UNP A0A6H1VH54
B	501	CYS	ALA	conflict	UNP A0A6H1VH54
B	509	ARG	GLU	conflict	UNP A0A6H1VH54
B	512	ARG	-	expression tag	UNP A0A6H1VH54
B	513	ARG	-	expression tag	UNP A0A6H1VH54
C	64	LYS	GLU	conflict	UNP A0A6H1VH54
C	375	SER	TYR	conflict	UNP A0A6H1VH54
C	501	CYS	ALA	conflict	UNP A0A6H1VH54
C	509	ARG	GLU	conflict	UNP A0A6H1VH54
C	512	ARG	-	expression tag	UNP A0A6H1VH54
C	513	ARG	-	expression tag	UNP A0A6H1VH54

- Molecule 2 is a protein called CD4 binding site antibody Ab1573 - Fab heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	122	Total	C	N	O	S	0	0
			914	578	159	173	4		
2	I	122	Total	C	N	O	S	0	0
			905	573	156	172	4		
2	P	117	Total	C	N	O	S	0	0
			877	559	153	161	4		

- Molecule 3 is a protein called CD4 binding site antibody Ab1573 - Fab light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	J	110	Total	C	N	O	S	0	0
			828	520	136	169	3		
3	L	110	Total	C	N	O	S	0	0
			834	525	136	170	3		
3	Q	108	Total	C	N	O	S	0	0
			816	510	134	169	3		

- Molecule 4 is a protein called Envelope glycoprotein BG505 SOSIP.664 - gp41.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	X	124	Total	C	N	O	S	0	0
			947	602	164	175	6		
4	Y	117	Total	C	N	O	S	0	0
			917	579	161	172	5		
4	Z	117	Total	C	N	O	S	0	0
			930	588	163	174	5		

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
X	559	PRO	ILE	conflict	UNP Q2N0S6
X	605	CYS	THR	conflict	UNP Q2N0S6
Y	559	PRO	ILE	conflict	UNP Q2N0S6
Y	605	CYS	THR	conflict	UNP Q2N0S6
Z	559	PRO	ILE	conflict	UNP Q2N0S6
Z	605	CYS	THR	conflict	UNP Q2N0S6

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



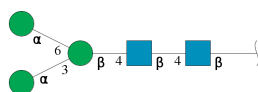
Mol	Chain	Residues	Atoms				AltConf	Trace
5	D	2	Total	C	N	O	0	0
			28	16	2	10		

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



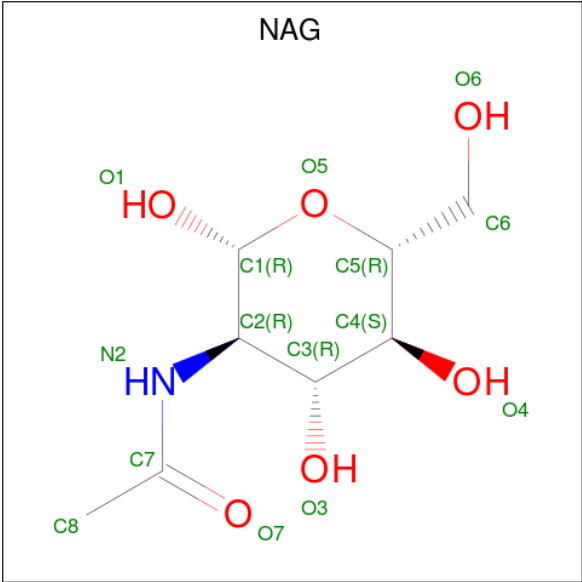
Mol	Chain	Residues	Atoms				AltConf	Trace
6	E	3	Total	C	N	O	0	0
			39	22	2	15		
6	F	3	Total	C	N	O	0	0
			39	22	2	15		
6	K	3	Total	C	N	O	0	0
			39	22	2	15		

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[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.



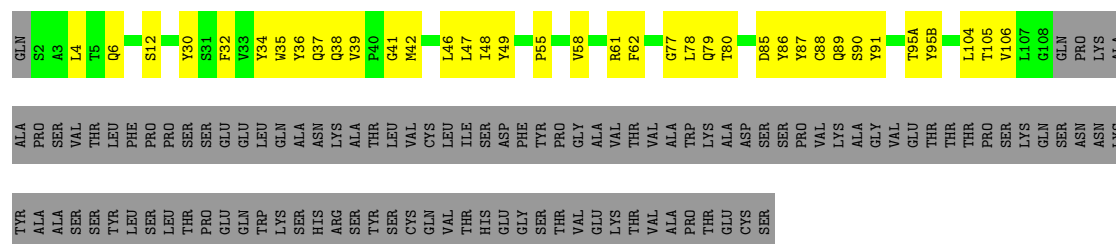
Mol	Chain	Residues	Atoms				AltConf	Trace
7	G	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



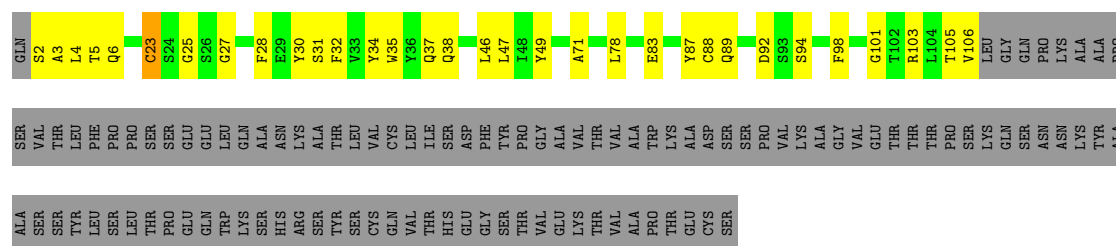
Mol	Chain	Residues	Atoms				AltConf
8	A	1	Total	C	N	O	0
			14	8	1	5	
8	B	1	Total	C	N	O	0
			28	16	2	10	
8	B	1	Total	C	N	O	0
			28	16	2	10	
8	C	1	Total	C	N	O	0
			28	16	2	10	
8	C	1	Total	C	N	O	0
			28	16	2	10	

34% 17% 49%



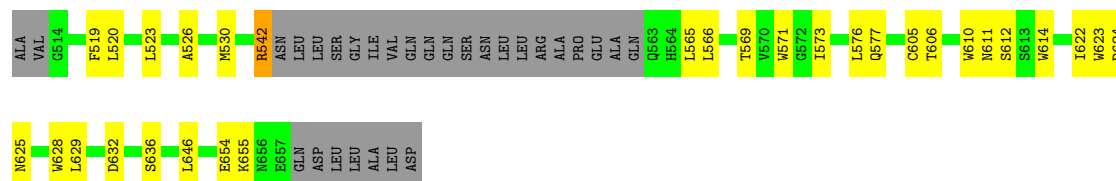
• Molecule 3: CD4 binding site antibody Ab1573 - Fab light chain

Chain Q: 35% 14% 50%



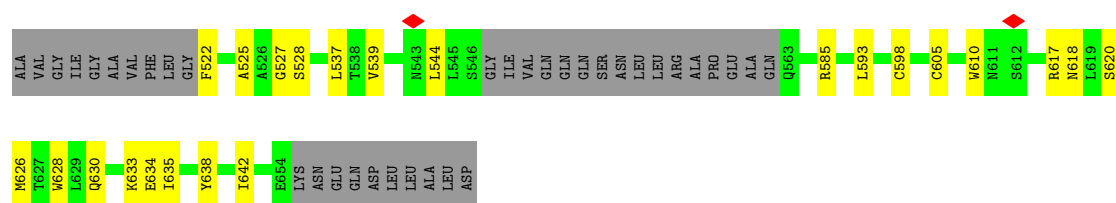
• Molecule 4: Envelope glycoprotein BG505 SOSIP.664 - gp41

Chain X: 61% 19% 19%



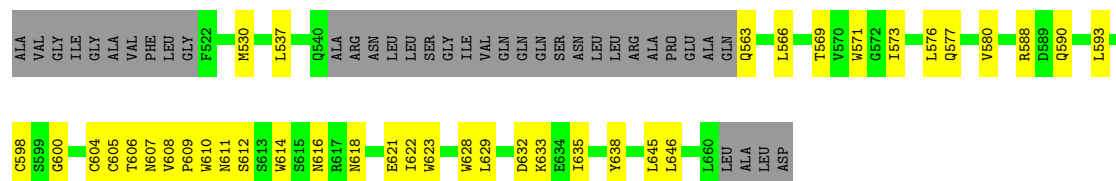
• Molecule 4: Envelope glycoprotein BG505 SOSIP.664 - gp41

Chain Y: 61% 15% 24%



• Molecule 4: Envelope glycoprotein BG505 SOSIP.664 - gp41

Chain Z: 52% 25% 24%



- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:  50% 50%

MAG1
MAG2

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

Chain E:  67% 33%

MAG1
MAG2
BMA3

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

Chain F:  33% 33% 33%

MAG1
MAG2
BMA3

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

Chain K:  100%

MAG1
MAG2
BMA3

- Molecule 7: alpha-D-mannopyranose-(1-3)-[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

Chain G:  40% 60%

MAG1
MAG2
BMA3
MAN4
MAN5

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	443817	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	1400	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.056	Depositor
Minimum map value	-0.024	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.002	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	334.40002, 334.40002, 334.40002	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8360001, 0.8360001, 0.8360001	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: MAN, BMA, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.33	0/3309	0.51	0/4493
1	B	0.29	0/3237	0.50	0/4393
1	C	0.38	0/3171	0.56	0/4310
2	H	0.36	0/932	0.56	0/1264
2	I	0.28	0/923	0.52	0/1254
2	P	0.42	0/894	0.57	0/1211
3	J	0.30	0/849	0.51	0/1155
3	L	0.43	0/856	0.54	0/1165
3	Q	0.47	0/837	0.57	0/1139
4	X	0.29	0/965	0.48	0/1311
4	Y	0.27	0/933	0.43	0/1267
4	Z	0.32	0/948	0.48	0/1287
All	All	0.34	0/17854	0.52	0/24249

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	3245	0	3137	114	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3175	0	3066	97	0
1	C	3110	0	3004	126	0
2	H	914	0	907	33	0
2	I	905	0	889	38	0
2	P	877	0	871	29	0
3	J	828	0	774	37	0
3	L	834	0	779	27	0
3	Q	816	0	753	36	0
4	X	947	0	896	27	0
4	Y	917	0	889	23	0
4	Z	930	0	899	31	0
5	D	28	0	25	1	0
6	E	39	0	34	0	0
6	F	39	0	34	1	0
6	K	39	0	34	0	0
7	G	61	0	52	6	0
8	A	14	0	13	0	0
8	B	28	0	26	0	0
8	C	28	0	26	0	0
All	All	17774	0	17108	558	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 16.

The worst 5 of 558 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:262:ASN:ND2	7:G:1:NAG:C1	1.76	1.46
1:A:128:THR:HG22	1:A:192:ARG:HH21	1.16	1.06
1:A:262:ASN:HD22	7:G:1:NAG:C1	1.50	1.01
1:A:128:THR:CG2	1:A:192:ARG:HH21	1.78	0.96
3:Q:35:TRP:CZ3	3:Q:88:CYS:HB3	2.05	0.90

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	399/481 (83%)	368 (92%)	31 (8%)	0	100	100
1	B	389/481 (81%)	355 (91%)	33 (8%)	1 (0%)	41	75
1	C	386/481 (80%)	326 (84%)	60 (16%)	0	100	100
2	H	120/228 (53%)	116 (97%)	4 (3%)	0	100	100
2	I	120/228 (53%)	119 (99%)	1 (1%)	0	100	100
2	P	113/228 (50%)	102 (90%)	11 (10%)	0	100	100
3	J	108/216 (50%)	101 (94%)	7 (6%)	0	100	100
3	L	108/216 (50%)	101 (94%)	7 (6%)	0	100	100
3	Q	106/216 (49%)	94 (89%)	12 (11%)	0	100	100
4	X	120/153 (78%)	108 (90%)	12 (10%)	0	100	100
4	Y	113/153 (74%)	103 (91%)	10 (9%)	0	100	100
4	Z	113/153 (74%)	106 (94%)	7 (6%)	0	100	100
All	All	2195/3234 (68%)	1999 (91%)	195 (9%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	365	SER

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	353/428 (82%)	353 (100%)	0	100	100
1	B	346/428 (81%)	345 (100%)	1 (0%)	92	95
1	C	341/428 (80%)	337 (99%)	4 (1%)	71	83
2	H	97/190 (51%)	95 (98%)	2 (2%)	53	72
2	I	95/190 (50%)	95 (100%)	0	100	100
2	P	91/190 (48%)	90 (99%)	1 (1%)	73	84
3	J	91/185 (49%)	91 (100%)	0	100	100
3	L	92/185 (50%)	92 (100%)	0	100	100
3	Q	90/185 (49%)	89 (99%)	1 (1%)	73	84
4	X	94/129 (73%)	93 (99%)	1 (1%)	73	84
4	Y	97/129 (75%)	97 (100%)	0	100	100
4	Z	99/129 (77%)	99 (100%)	0	100	100
All	All	1886/2796 (68%)	1876 (100%)	10 (0%)	89	93

5 of 10 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	P	62	LYS
3	Q	23	CYS
4	X	542	ARG
1	C	368	ASP
1	C	369	LEU

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

Mol	Chain	Res	Type
1	C	289	ASN
3	J	89	GLN
4	Z	607	ASN
3	Q	38	GLN
1	B	478	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 ⓘ

16 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$
5	NAG	D	1	5,1	14,14,15	0.75	1 (7%)	17,19,21	0.55	0
5	NAG	D	2	5	14,14,15	0.41	0	17,19,21	0.55	0
6	NAG	E	1	6,1	14,14,15	0.62	1 (7%)	17,19,21	0.72	0
6	NAG	E	2	6	14,14,15	0.30	0	17,19,21	0.42	0
6	BMA	E	3	6	11,11,12	0.52	0	15,15,17	0.73	0
6	NAG	F	1	6,1	14,14,15	0.29	0	17,19,21	0.79	1 (5%)
6	NAG	F	2	6	14,14,15	0.29	0	17,19,21	0.61	0
6	BMA	F	3	6	11,11,12	0.22	0	15,15,17	0.61	0
7	NAG	G	1	7	14,14,15	0.49	0	17,19,21	0.59	0
7	NAG	G	2	7	14,14,15	0.22	0	17,19,21	0.59	0
7	BMA	G	3	7	11,11,12	0.62	0	15,15,17	0.79	0
7	MAN	G	4	7	11,11,12	0.68	0	15,15,17	1.08	2 (13%)
7	MAN	G	5	7	11,11,12	0.62	0	15,15,17	1.12	2 (13%)
6	NAG	K	1	6,1	14,14,15	0.39	0	17,19,21	0.59	0
6	NAG	K	2	6	14,14,15	0.19	0	17,19,21	0.47	0
6	BMA	K	3	6	11,11,12	0.25	0	15,15,17	0.59	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
5	NAG	D	1	5,1	-	3/6/23/26	0/1/1/1
5	NAG	D	2	5	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	E	1	6,1	-	3/6/23/26	0/1/1/1
6	NAG	E	2	6	-	2/6/23/26	0/1/1/1
6	BMA	E	3	6	-	1/2/19/22	0/1/1/1
6	NAG	F	1	6,1	-	2/6/23/26	0/1/1/1
6	NAG	F	2	6	-	0/6/23/26	0/1/1/1
6	BMA	F	3	6	-	0/2/19/22	0/1/1/1
7	NAG	G	1	7	-	4/6/23/26	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
6	NAG	K	1	6,1	-	2/6/23/26	0/1/1/1
6	NAG	K	2	6	-	0/6/23/26	0/1/1/1
6	BMA	K	3	6	-	1/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	1	NAG	O5-C1	-2.68	1.39	1.43
6	E	1	NAG	O5-C1	-2.25	1.40	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	G	5	MAN	C1-O5-C5	2.79	115.97	112.19
7	G	4	MAN	C1-O5-C5	2.63	115.75	112.19
7	G	5	MAN	O2-C2-C3	-2.44	105.25	110.14
7	G	4	MAN	O2-C2-C3	-2.17	105.79	110.14
6	F	1	NAG	C1-O5-C5	2.04	114.96	112.19

There are no chirality outliers.

5 of 22 torsion outliers are listed below:

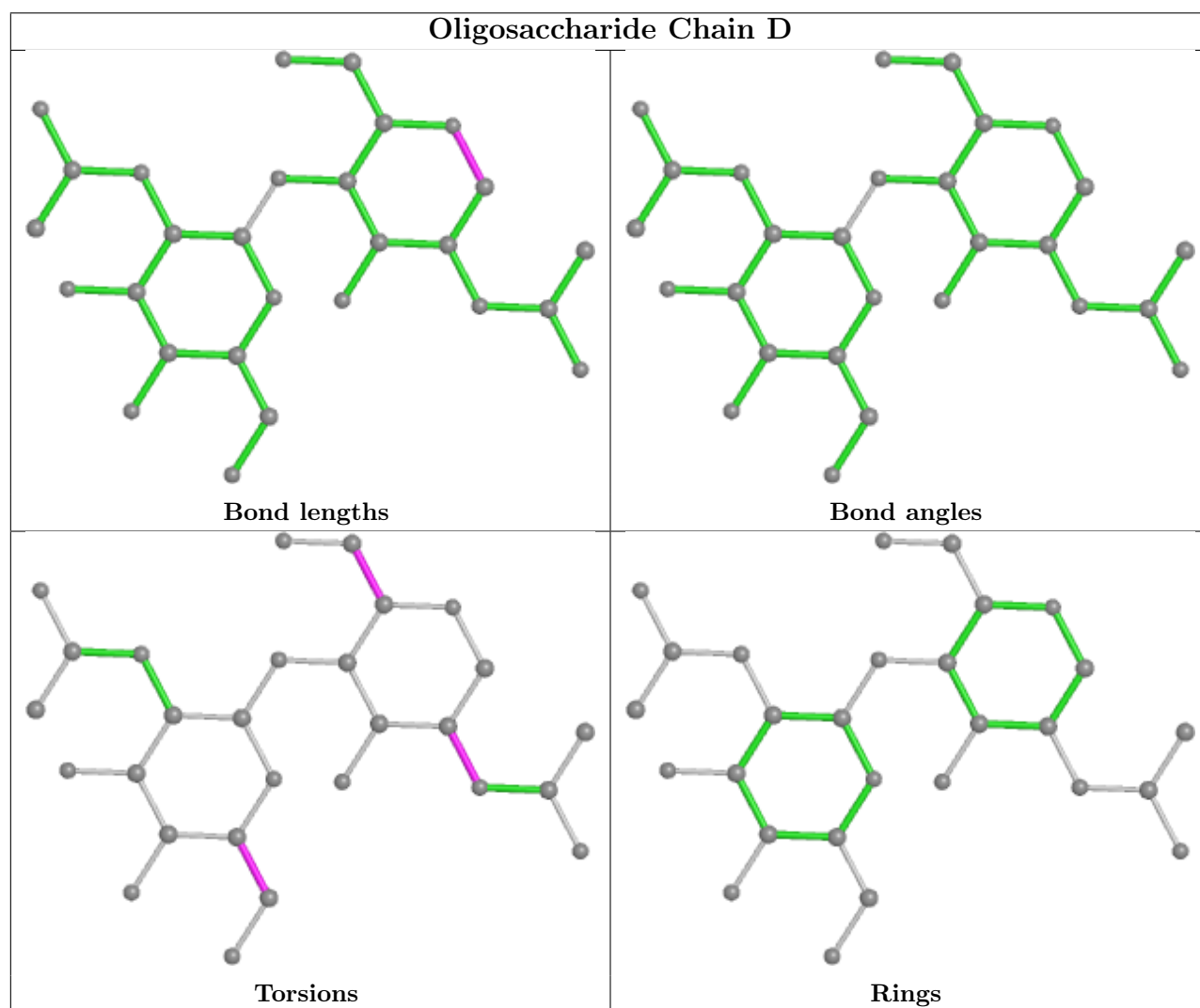
Mol	Chain	Res	Type	Atoms
6	F	1	NAG	C8-C7-N2-C2
6	F	1	NAG	O7-C7-N2-C2
5	D	1	NAG	O5-C5-C6-O6
6	E	1	NAG	O5-C5-C6-O6
5	D	2	NAG	C4-C5-C6-O6

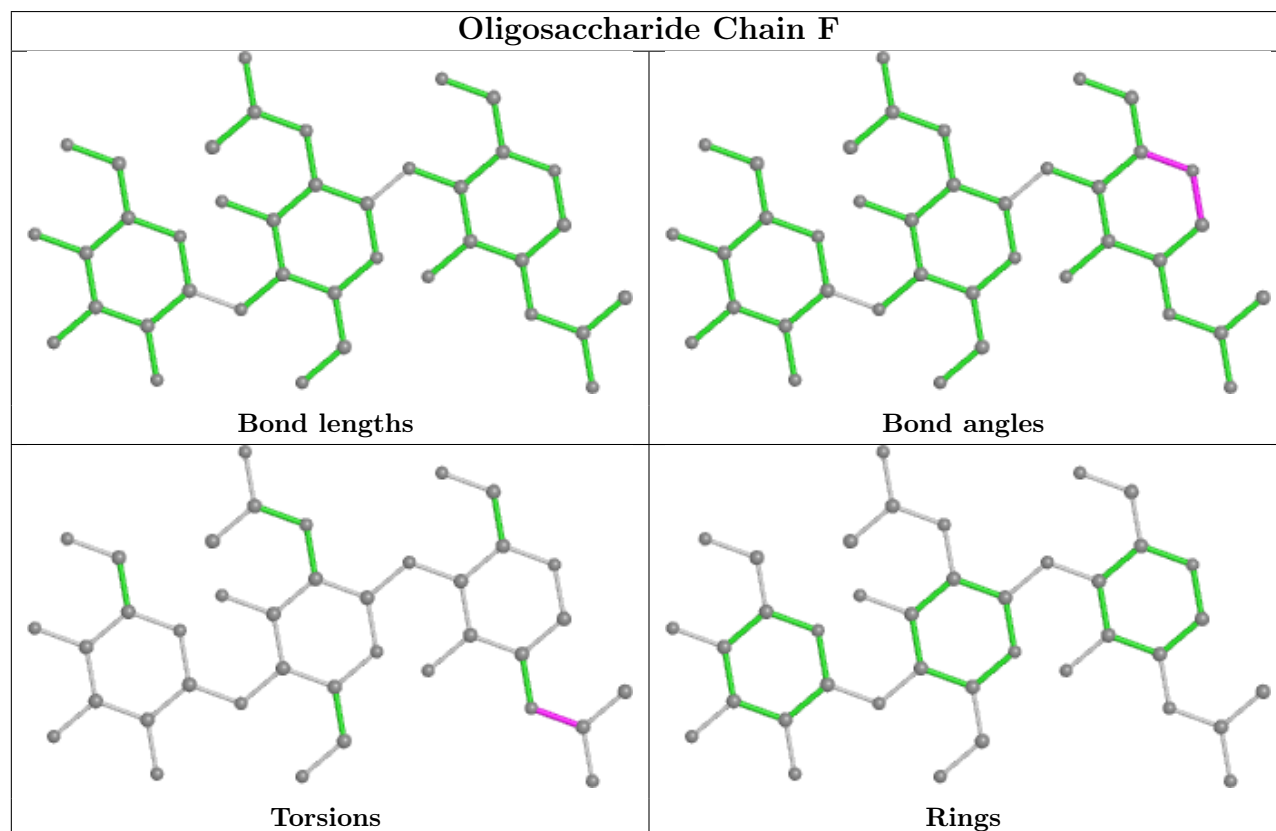
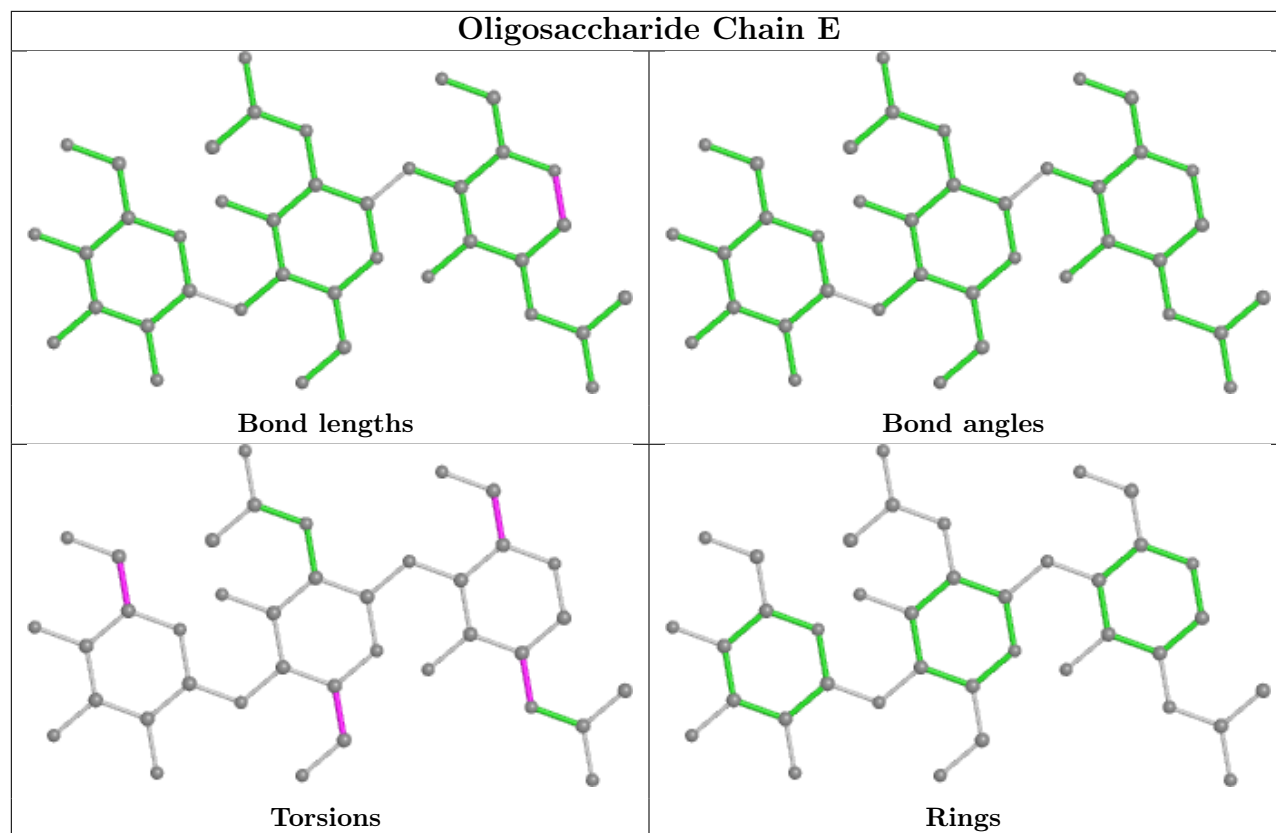
There are no ring outliers.

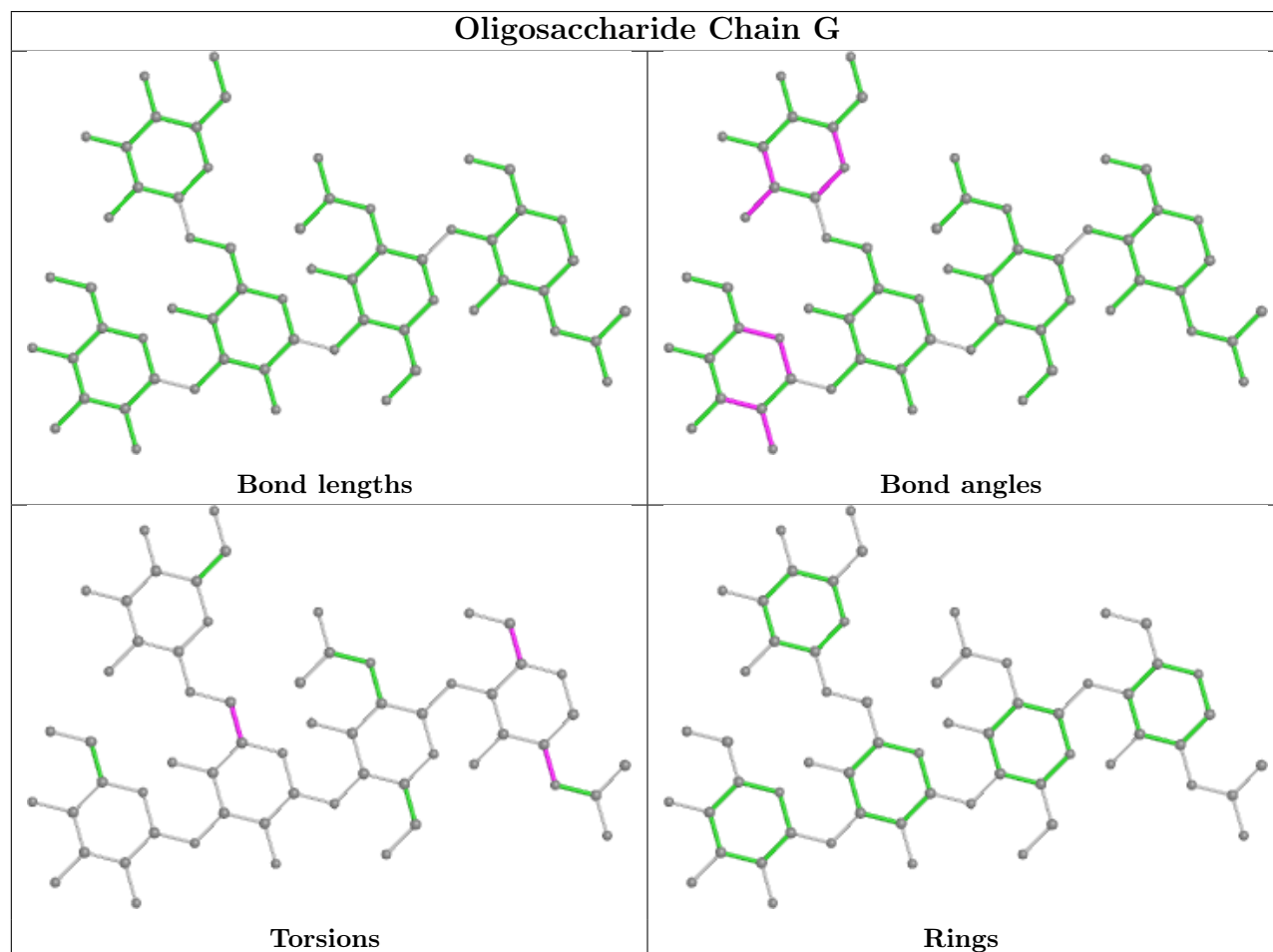
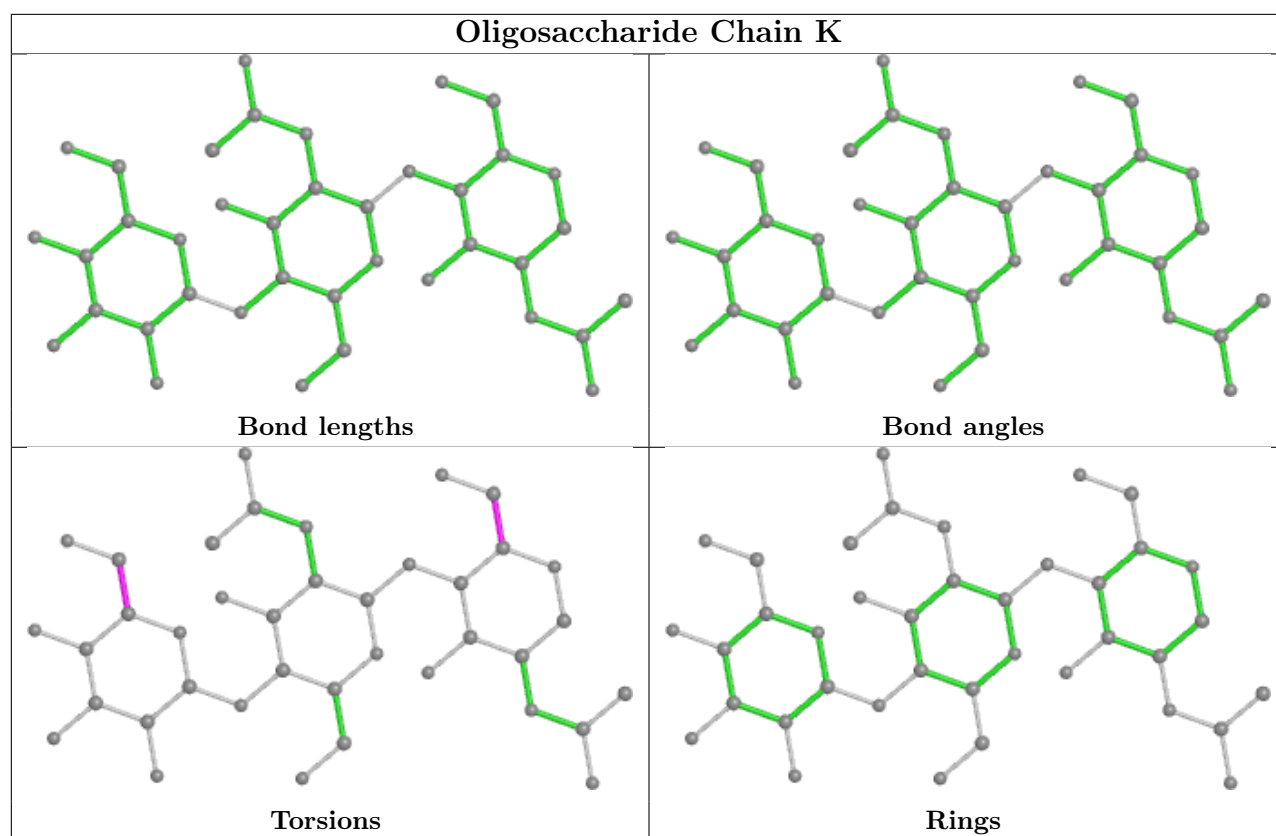
5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	G	1	NAG	6	0
5	D	2	NAG	1	0
6	F	2	NAG	1	0
6	F	1	NAG	1	0
5	D	1	NAG	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 oligosaccharide.







5.6 Ligand geometry

5 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	$\# Z > 2$
8	NAG	C	601	1	14,14,15	0.40	0	17,19,21	0.82	0
8	NAG	A	601	1	14,14,15	0.25	0	17,19,21	0.59	0
8	NAG	B	601	1	14,14,15	0.34	0	17,19,21	1.04	2 (11%)
8	NAG	B	602	1	14,14,15	0.30	0	17,19,21	0.70	0
8	NAG	C	602	1	14,14,15	0.27	0	17,19,21	0.85	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
8	NAG	C	601	1	-	0/6/23/26	0/1/1/1
8	NAG	A	601	1	-	3/6/23/26	0/1/1/1
8	NAG	B	601	1	-	3/6/23/26	0/1/1/1
8	NAG	B	602	1	-	2/6/23/26	0/1/1/1
8	NAG	C	602	1	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	601	NAG	C4-C3-C2	-2.35	107.57	111.02
8	B	601	NAG	O5-C5-C6	2.12	110.53	107.20

There are no chirality outliers.

5 of 9 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	601	NAG	C8-C7-N2-C2
8	A	601	NAG	O7-C7-N2-C2
8	B	601	NAG	C1-C2-N2-C7
8	B	601	NAG	C8-C7-N2-C2
8	B	601	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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.

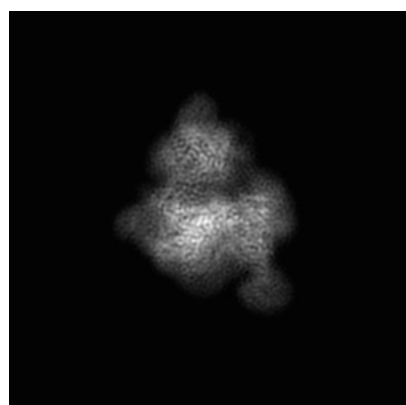
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-25878. These allow visual inspection of the internal detail of the map and identification of artifacts.

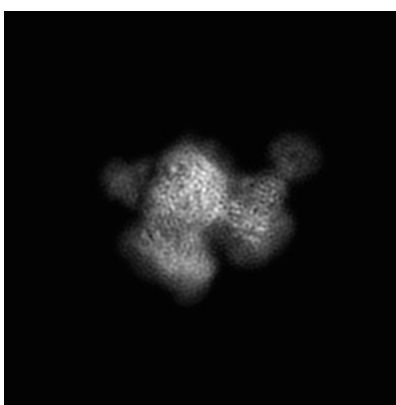
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

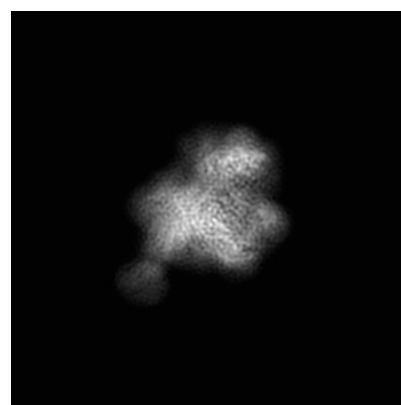
6.1.1 Primary map



X



Y

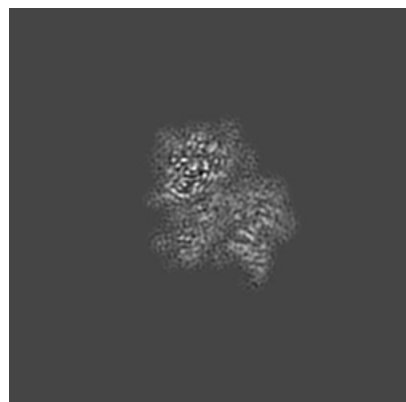


Z

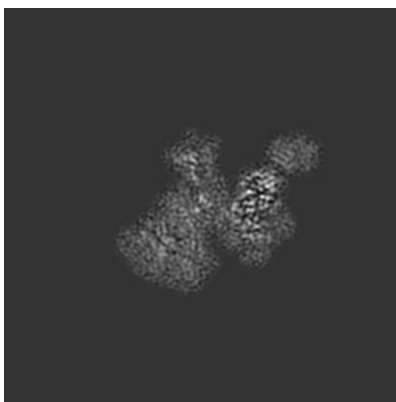
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

6.2.1 Primary map



X Index: 200



Y Index: 200

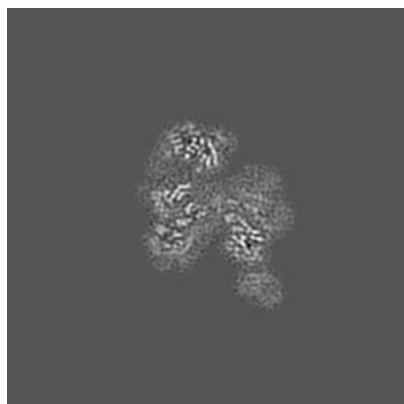


Z Index: 200

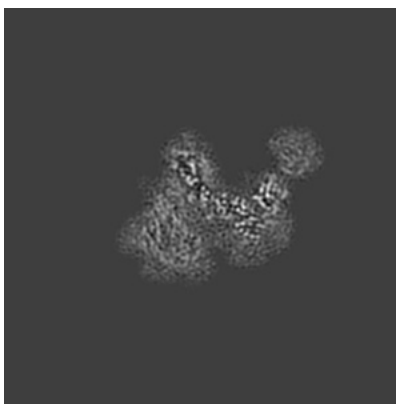
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

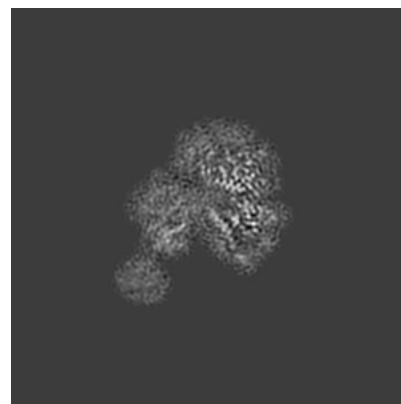
6.3.1 Primary map



X Index: 215



Y Index: 184



Z Index: 188

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

6.4 Orthogonal surface views [i](#)

6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.01. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

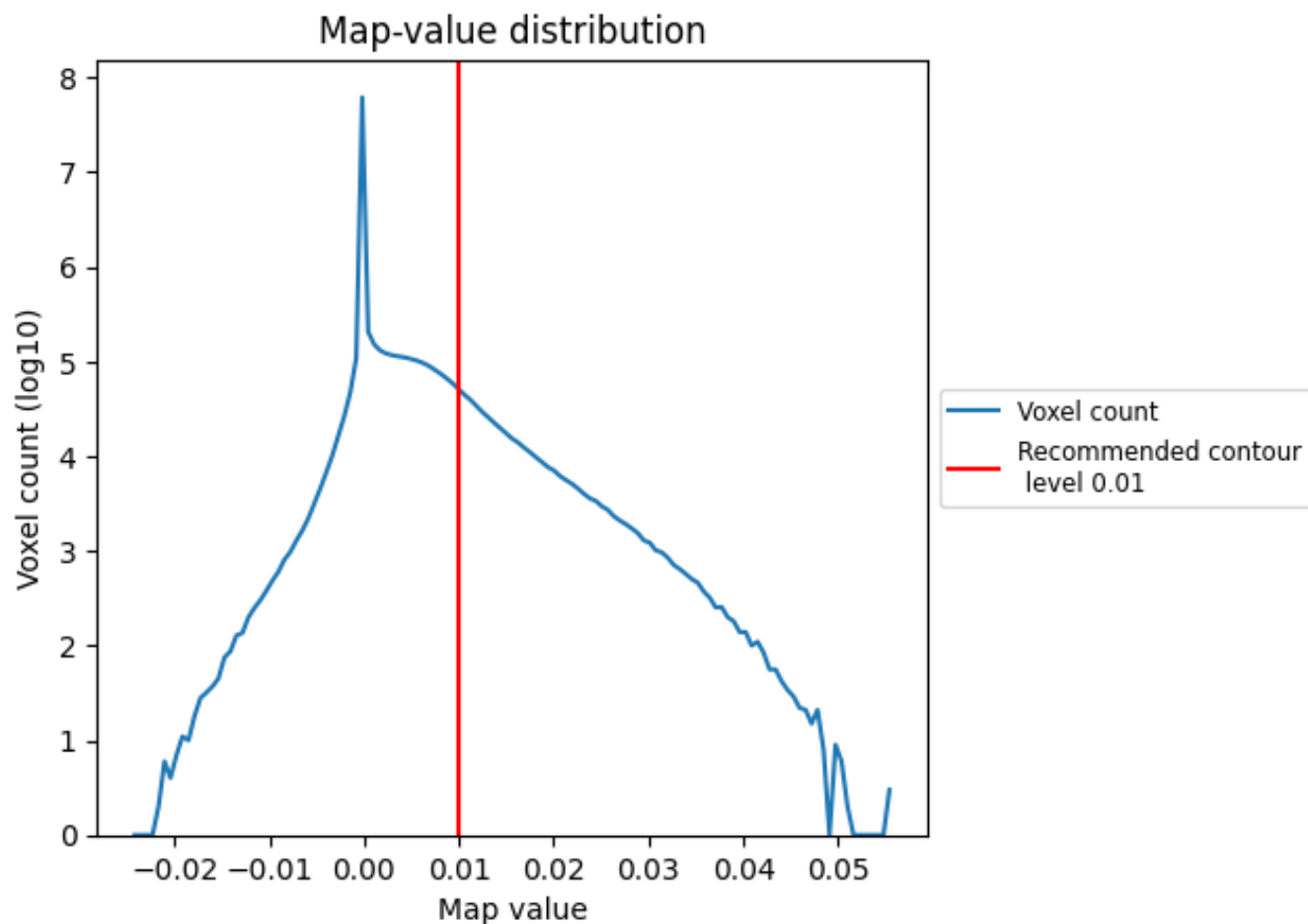
6.5 Mask visualisation

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

7 Map analysis [i](#)

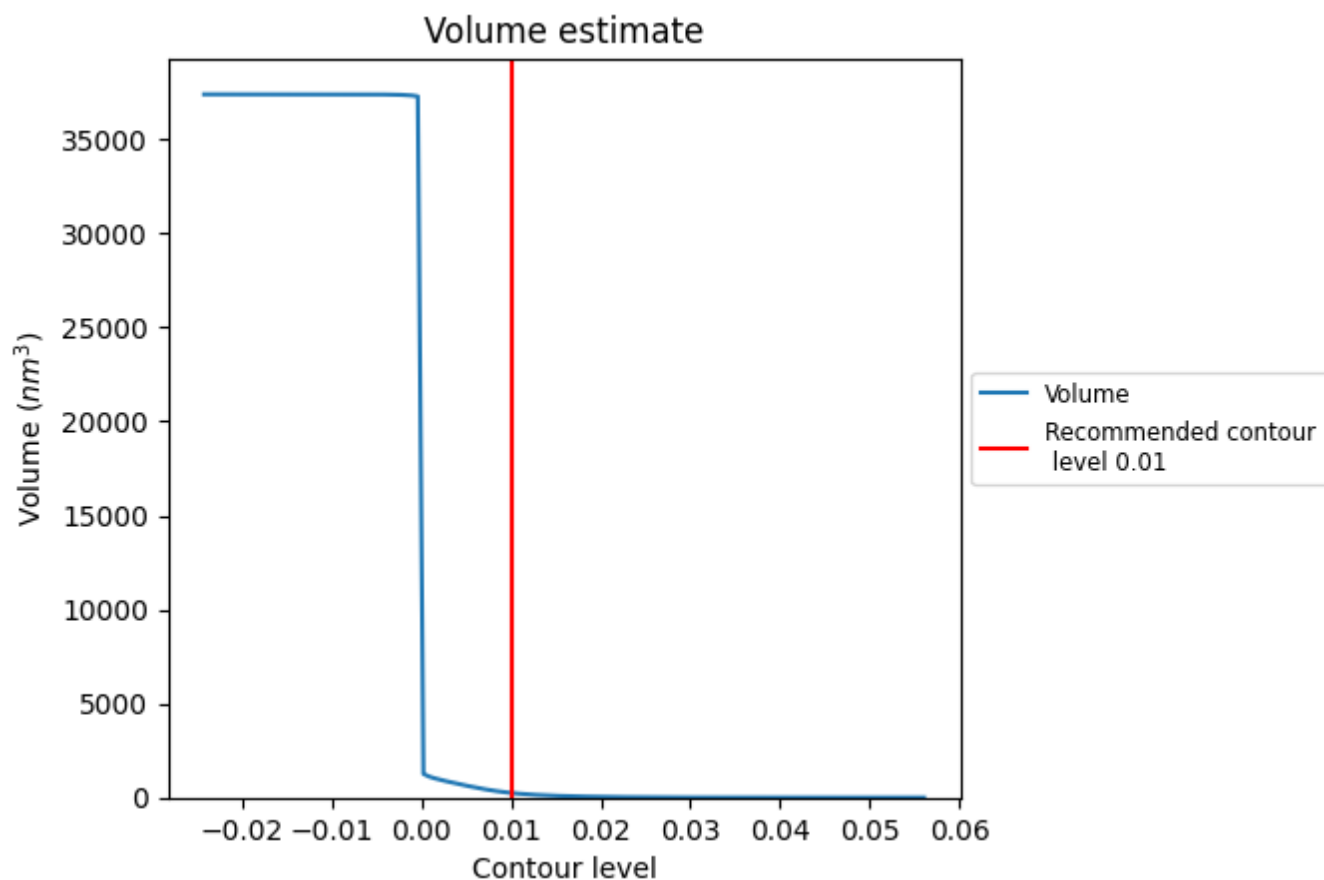
This section contains the results of statistical analysis of the map.

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.

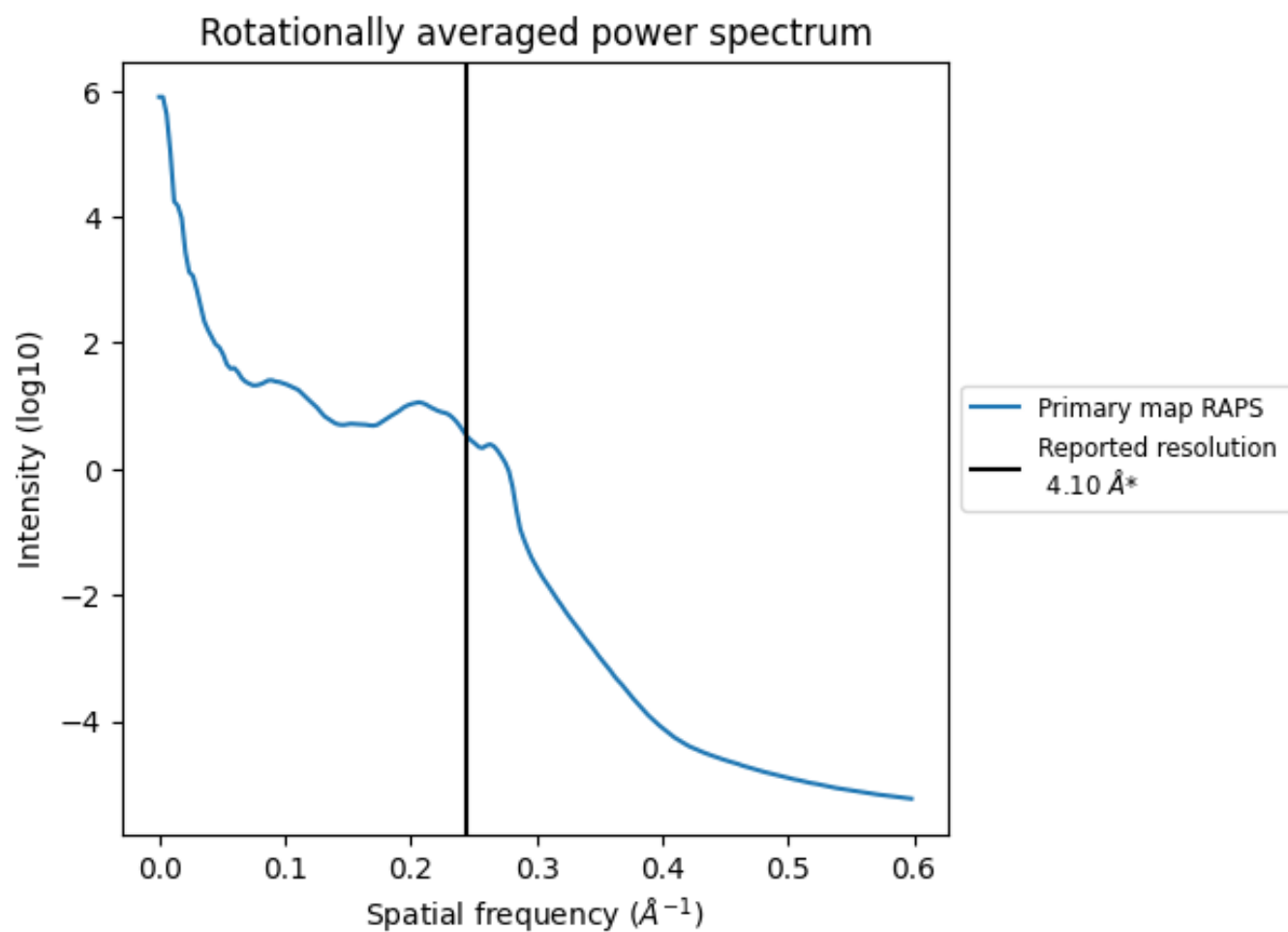
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 247 nm³; this corresponds to an approximate mass of 223 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 0.244 Å⁻¹

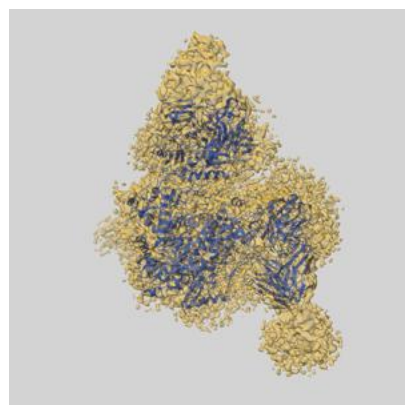
8 Fourier-Shell correlation

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

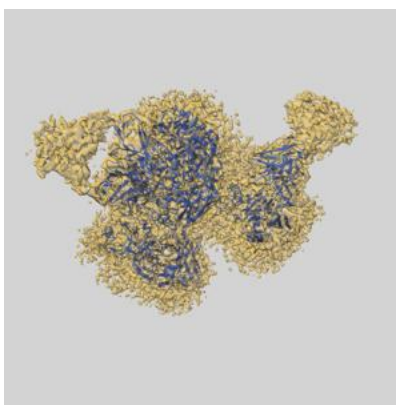
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-25878 and PDB model 7TFO. Per-residue inclusion information can be found in section [3](#) on page [8](#).

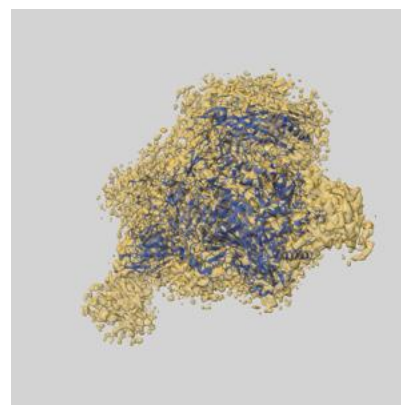
9.1 Map-model overlay [i](#)



X



Y



Z

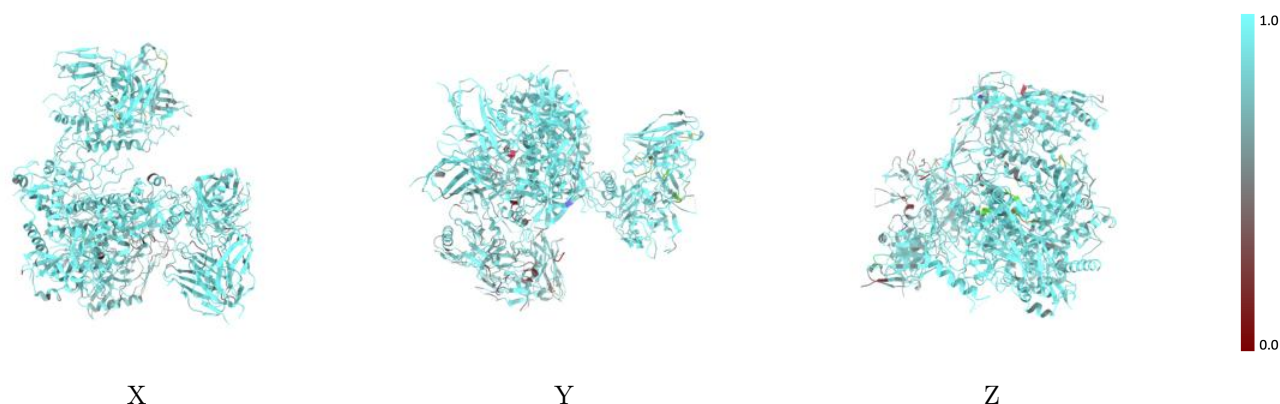
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



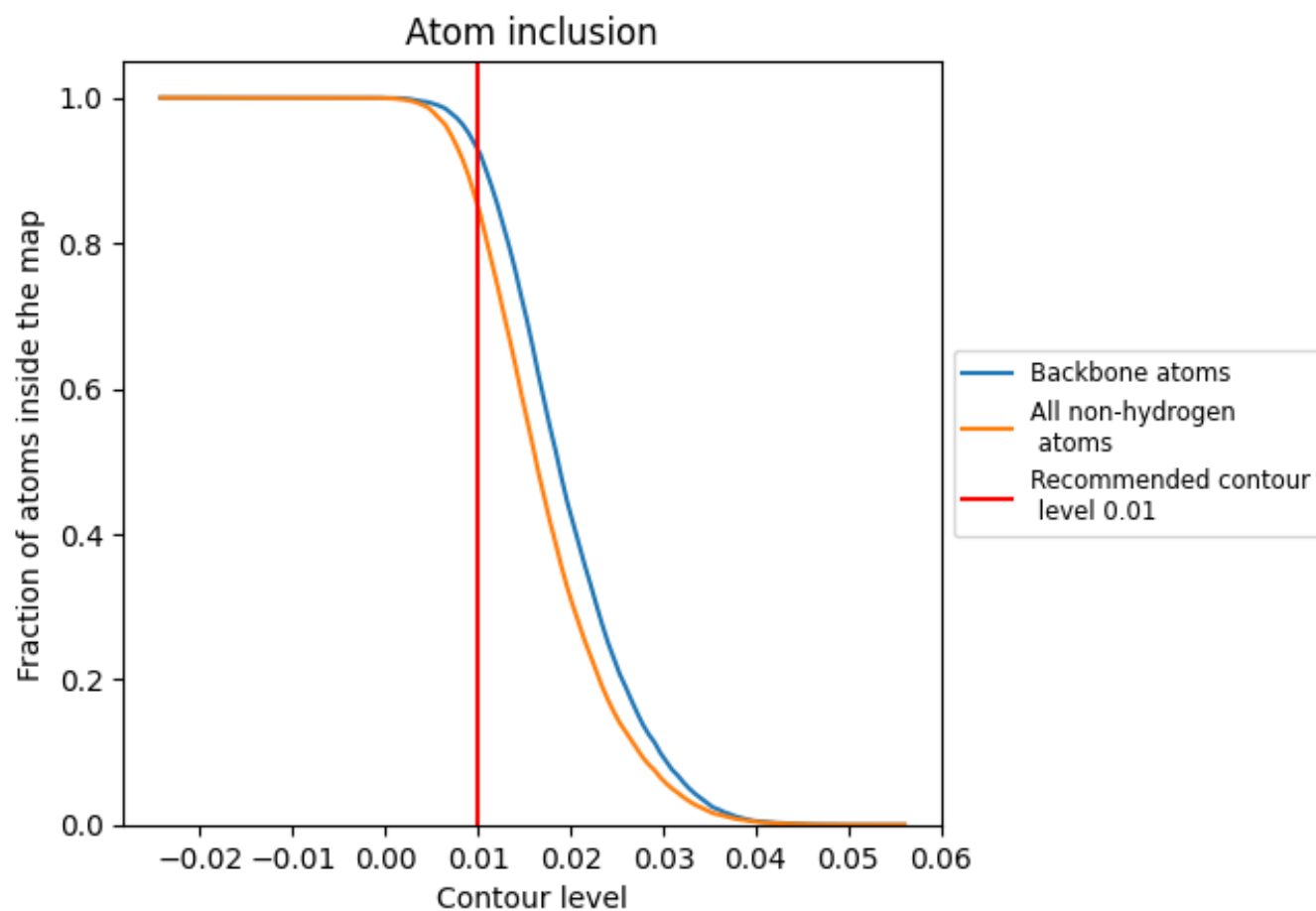
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.01).





































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 (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8521	 0.3590
A	 0.8714	 0.3750
B	 0.7570	 0.2520
C	 0.8847	 0.4010
D	 0.8571	 0.3460
E	 0.8718	 0.3650
F	 0.7436	 0.3480
G	 0.7541	 0.2660
H	 0.8845	 0.4110
I	 0.7864	 0.2850
J	 0.8403	 0.3280
K	 0.7179	 0.2830
L	 0.9041	 0.4420
P	 0.8936	 0.4260
Q	 0.8995	 0.4460
X	 0.8891	 0.3650
Y	 0.8641	 0.3410
Z	 0.8849	 0.3770

