



wwPDB EM Validation Summary Report ⓘ

Oct 15, 2024 – 12:01 AM JST

PDB ID : 8I4T
EMDB ID : EMD-35183
Title : Structure of the asymmetric unit of SFTSV virion
Authors : Du, S.; Peng, R.; Qi, J.; Li, C.
Deposited on : 2023-01-21
Resolution : 5.20 Å(reported)

This is a wwPDB EM Validation Summary 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.dev113
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.39

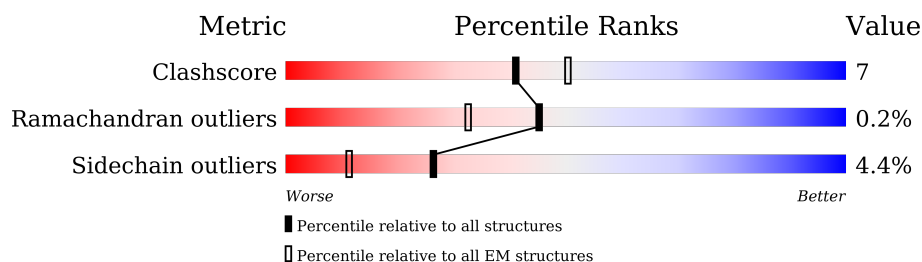
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 5.20 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	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	B	513	
1	G	513	
1	H	513	
1	I	513	
1	J	513	
1	K	513	
1	L	513	
1	O	513	

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Mol	Chain	Length	Quality of chain
1	P	513	
1	T	513	
1	V	513	
1	W	513	
2	A	560	
2	C	560	
2	D	560	
2	E	560	
2	F	560	
2	M	560	
2	N	560	
2	Q	560	
2	R	560	
2	S	560	
2	U	560	
2	X	560	
3	Y	2	
3	Z	2	
3	a	2	
3	b	2	
3	c	2	
3	d	2	

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 91976 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 Envelopment polypeptide.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	G	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	B	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	H	504	Total	C	N	O	S	0	0
			3778	2364	659	718	37		
1	I	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	J	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	K	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	L	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	T	505	Total	C	N	O	S	0	0
			3782	2366	660	719	37		
1	O	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	V	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	P	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		
1	W	509	Total	C	N	O	S	0	0
			3808	2381	665	725	37		

- Molecule 2 is a protein called Envelopment polypeptide.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	F	456	Total	C	N	O	S	0	0
			3478	2184	597	659	38		
2	A	499	Total	C	N	O	S	0	0
			3815	2404	657	712	42		
2	C	495	Total	C	N	O	S	0	0
			3795	2392	653	708	42		

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Mol	Chain	Residues	Atoms					AltConf	Trace
2	D	498	Total	C	N	O	S	0	0
			3817	2407	657	711	42		
2	E	498	Total	C	N	O	S	0	0
			3817	2407	657	711	42		
2	U	496	Total	C	N	O	S	0	0
			3803	2398	654	709	42		
2	X	498	Total	C	N	O	S	0	0
			3817	2407	657	711	42		
2	M	499	Total	C	N	O	S	0	0
			3820	2410	658	710	42		
2	Q	495	Total	C	N	O	S	0	0
			3790	2390	654	705	41		
2	R	501	Total	C	N	O	S	0	0
			3837	2421	660	714	42		
2	N	497	Total	C	N	O	S	0	0
			3812	2404	656	710	42		
2	S	498	Total	C	N	O	S	0	0
			3811	2402	656	711	42		

- Molecule 3 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
3	Y	2	Total	C	N	O	0	0
			28	16	2	10		
3	Z	2	Total	C	N	O	0	0
			28	16	2	10		
3	a	2	Total	C	N	O	0	0
			28	16	2	10		
3	b	2	Total	C	N	O	0	0
			28	16	2	10		
3	c	2	Total	C	N	O	0	0
			28	16	2	10		
3	d	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				AltConf
4	G	1	Total	C	N	O	0
			14	8	1	5	
4	G	1	Total	C	N	O	0
			14	8	1	5	
4	G	1	Total	C	N	O	0
			14	8	1	5	
4	F	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	A	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	B	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	C	1	Total	C	N	O	0
			14	8	1	5	
4	H	1	Total	C	N	O	0
			14	8	1	5	
4	H	1	Total	C	N	O	0
			14	8	1	5	
4	H	1	Total	C	N	O	0
			14	8	1	5	

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Mol	Chain	Residues	Atoms				AltConf
4	D	1	Total 14	C 8	N 1	O 5	0
4	D	1	Total 14	C 8	N 1	O 5	0
4	I	1	Total 14	C 8	N 1	O 5	0
4	I	1	Total 14	C 8	N 1	O 5	0
4	I	1	Total 14	C 8	N 1	O 5	0
4	E	1	Total 14	C 8	N 1	O 5	0
4	E	1	Total 14	C 8	N 1	O 5	0
4	J	1	Total 14	C 8	N 1	O 5	0
4	J	1	Total 14	C 8	N 1	O 5	0
4	J	1	Total 14	C 8	N 1	O 5	0
4	U	1	Total 14	C 8	N 1	O 5	0
4	U	1	Total 14	C 8	N 1	O 5	0
4	K	1	Total 14	C 8	N 1	O 5	0
4	K	1	Total 14	C 8	N 1	O 5	0
4	K	1	Total 14	C 8	N 1	O 5	0
4	X	1	Total 14	C 8	N 1	O 5	0
4	X	1	Total 14	C 8	N 1	O 5	0
4	L	1	Total 14	C 8	N 1	O 5	0
4	L	1	Total 14	C 8	N 1	O 5	0
4	L	1	Total 14	C 8	N 1	O 5	0
4	T	1	Total 14	C 8	N 1	O 5	0

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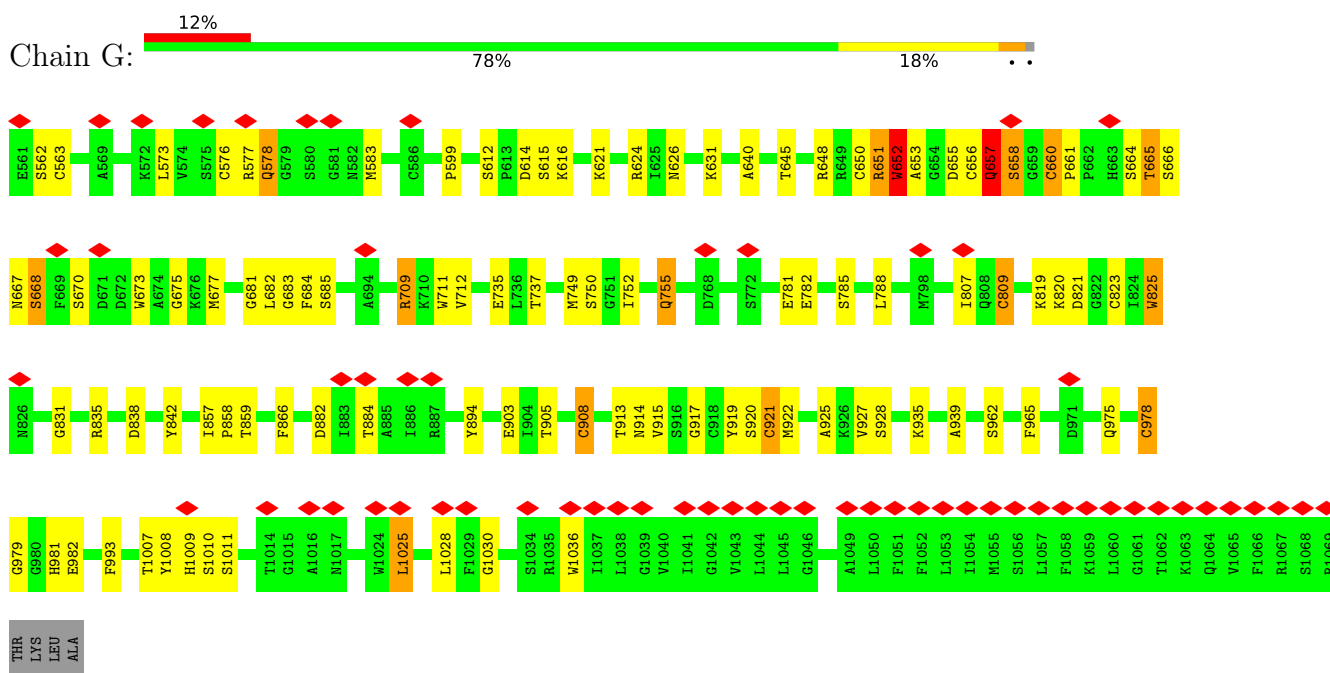
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Mol	Chain	Residues	Atoms				AltConf
4	T	1	Total 14	C 8	N 1	O 5	0
4	T	1	Total 14	C 8	N 1	O 5	0
4	M	1	Total 14	C 8	N 1	O 5	0
4	O	1	Total 14	C 8	N 1	O 5	0
4	O	1	Total 14	C 8	N 1	O 5	0
4	O	1	Total 14	C 8	N 1	O 5	0
4	Q	1	Total 14	C 8	N 1	O 5	0
4	V	1	Total 14	C 8	N 1	O 5	0
4	V	1	Total 14	C 8	N 1	O 5	0
4	V	1	Total 14	C 8	N 1	O 5	0
4	R	1	Total 14	C 8	N 1	O 5	0
4	N	1	Total 14	C 8	N 1	O 5	0
4	P	1	Total 14	C 8	N 1	O 5	0
4	P	1	Total 14	C 8	N 1	O 5	0
4	P	1	Total 14	C 8	N 1	O 5	0
4	S	1	Total 14	C 8	N 1	O 5	0
4	W	1	Total 14	C 8	N 1	O 5	0
4	W	1	Total 14	C 8	N 1	O 5	0
4	W	1	Total 14	C 8	N 1	O 5	0

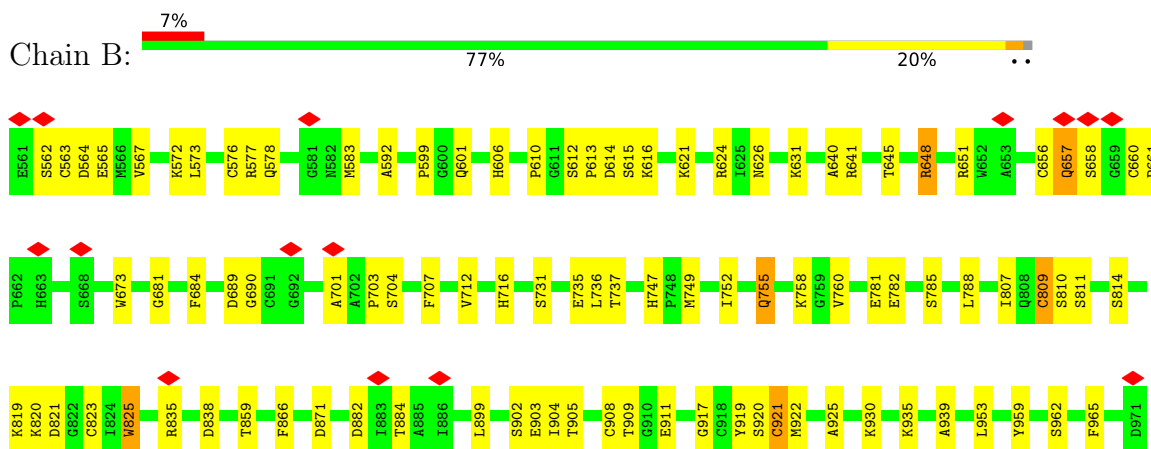
3 Residue-property plots

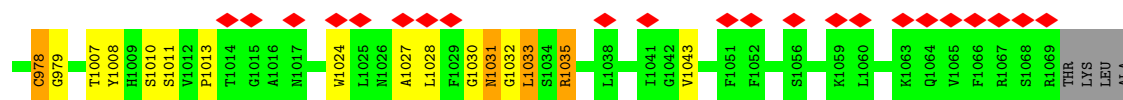
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Envelopment polypeptide

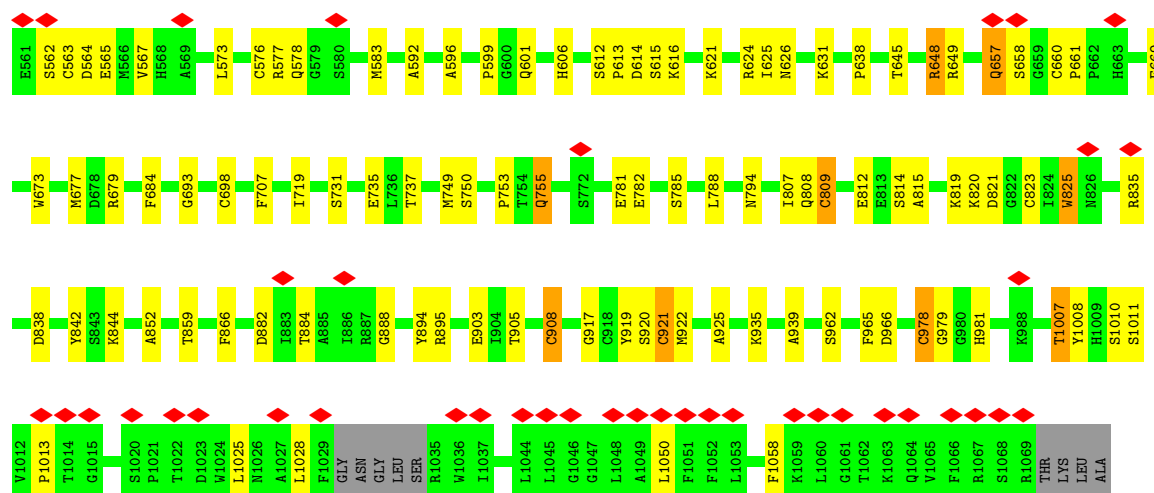
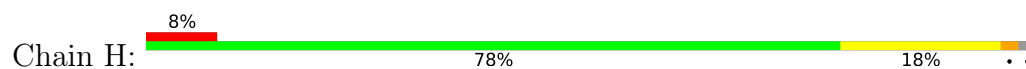


- Molecule 1: Envelopment polypeptide

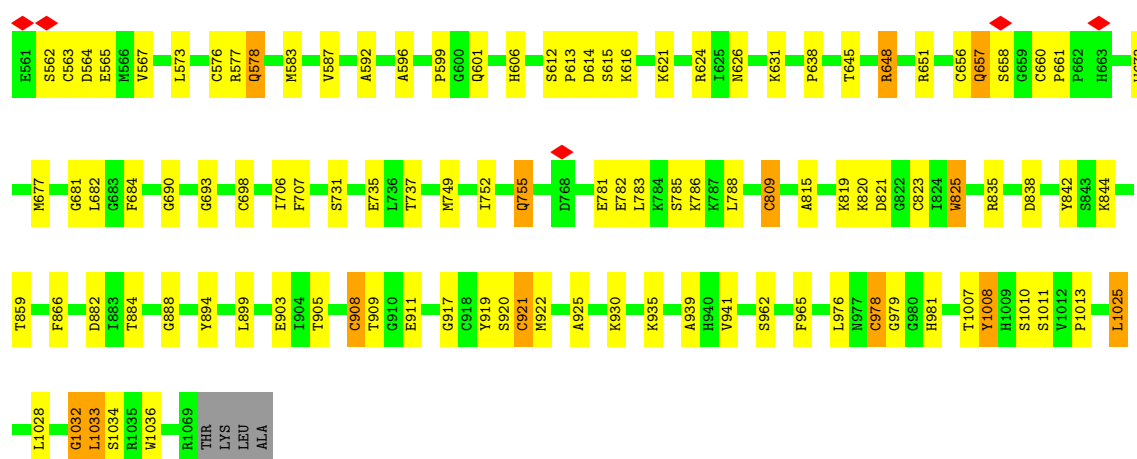
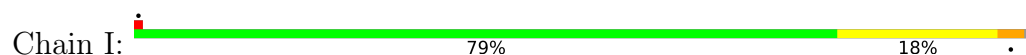




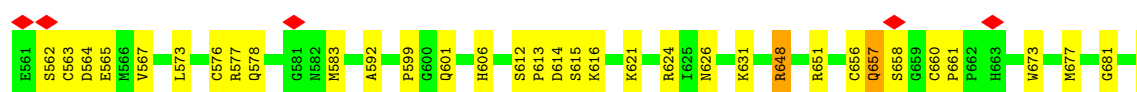
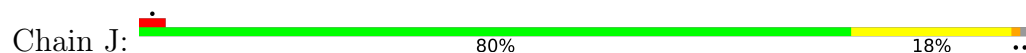
• Molecule 1: Envelopment polypeptide

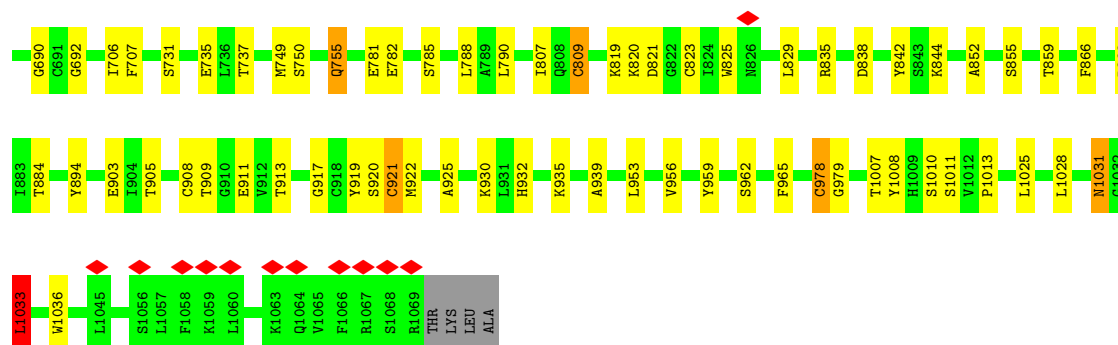


• Molecule 1: Envelopment polypeptide

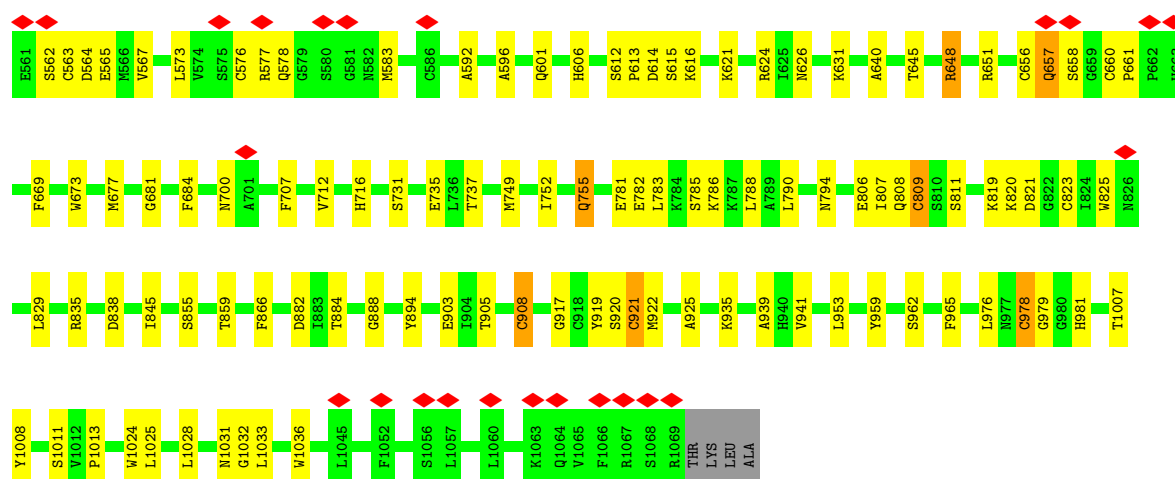
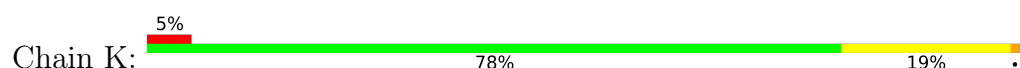


• Molecule 1: Envelopment polypeptide

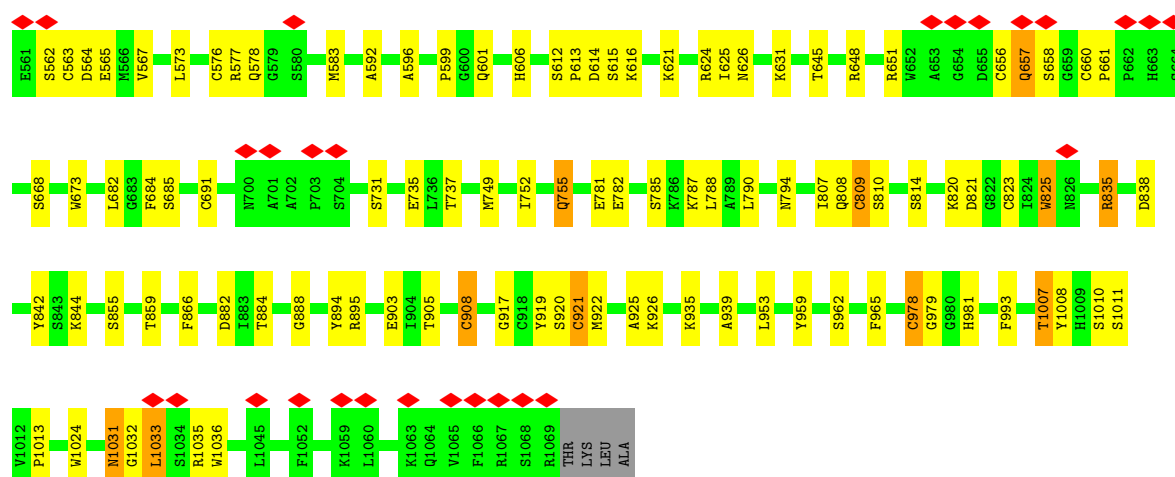
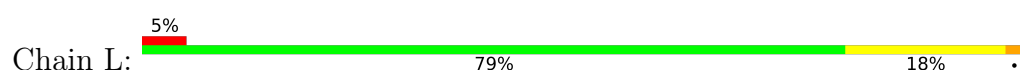




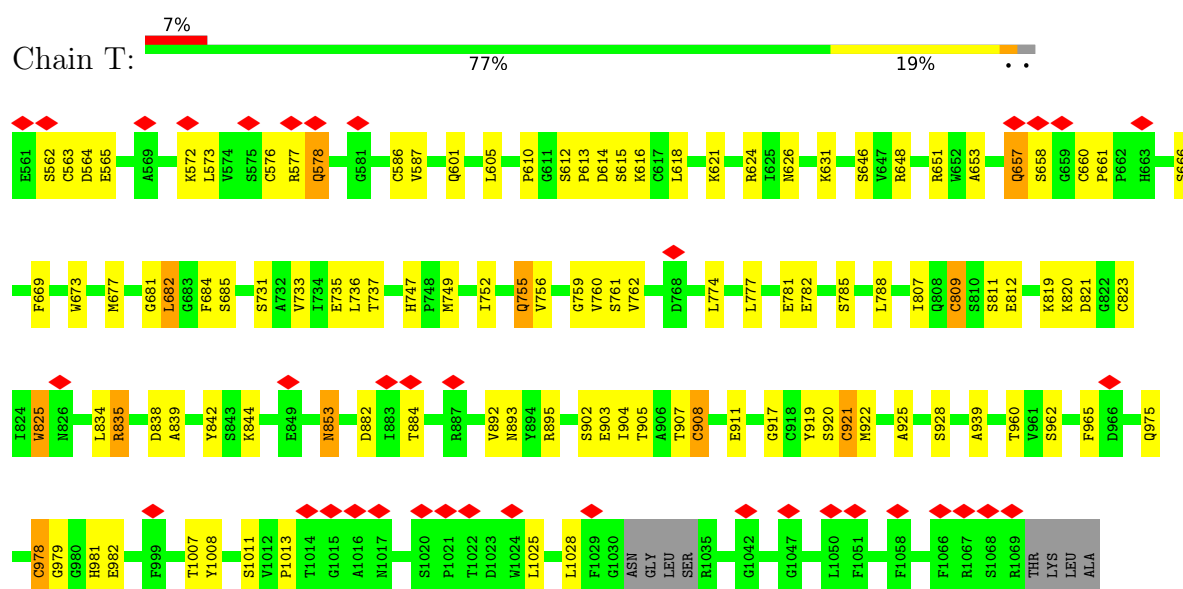
• Molecule 1: Envelopment polyprotein



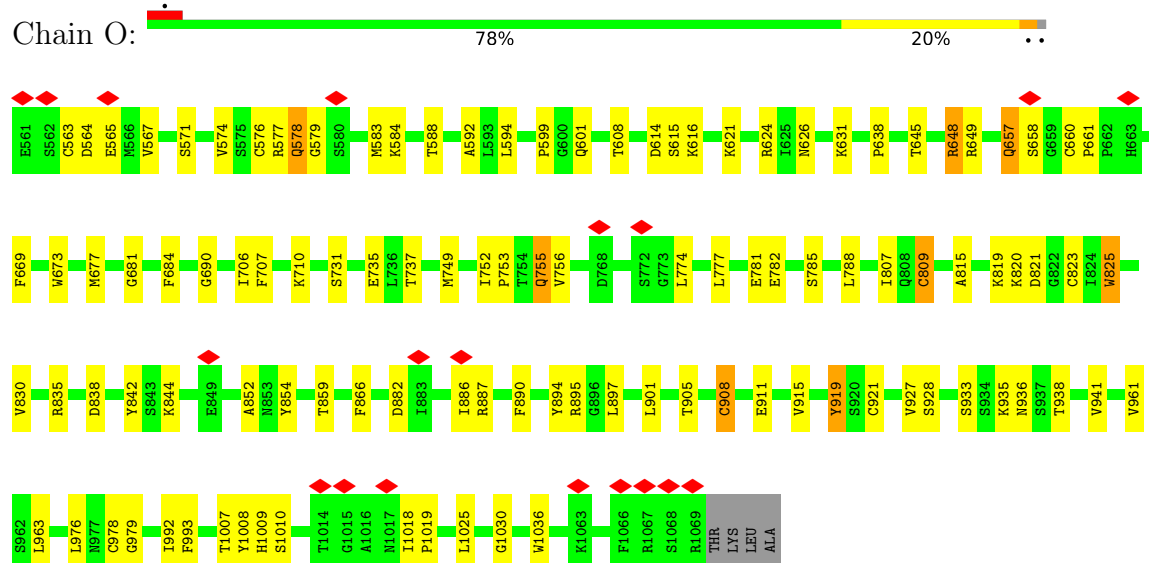
• Molecule 1: Envelopment polyprotein



• Molecule 1: Envelopment polyprotein

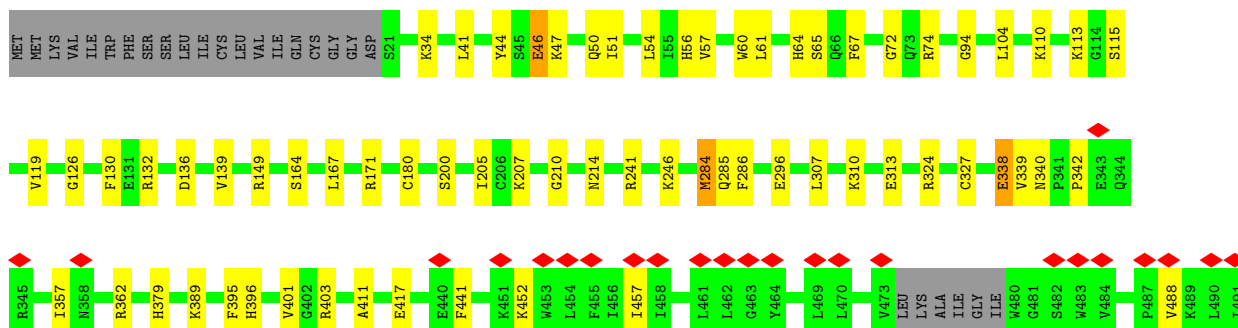


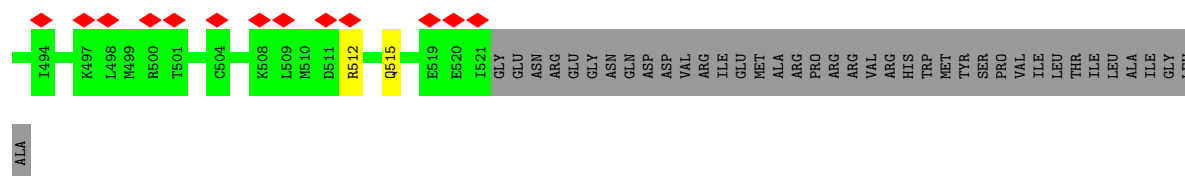
• Molecule 1: Envelopment polypeptide



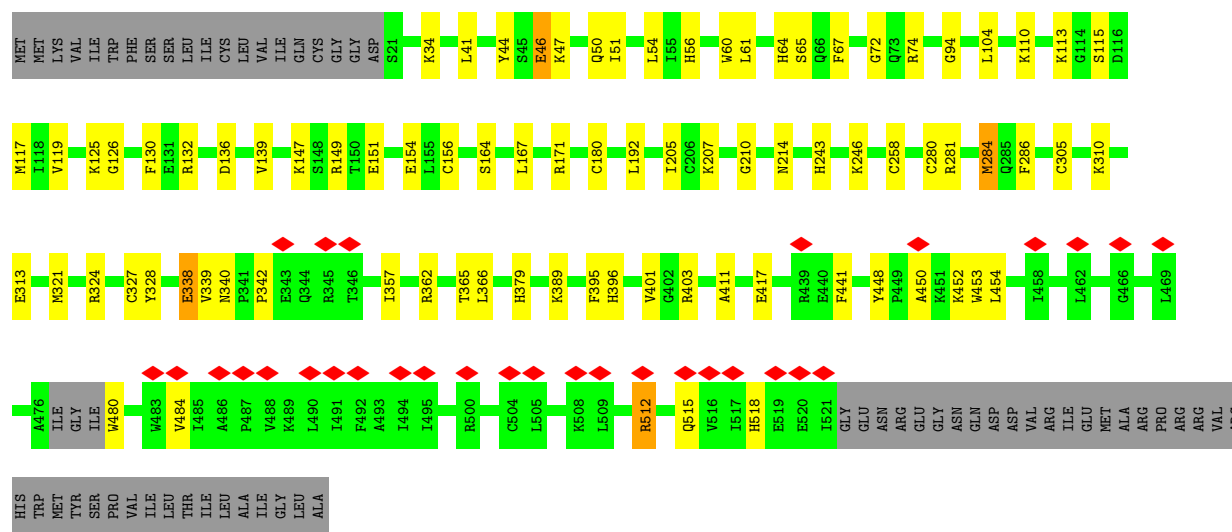
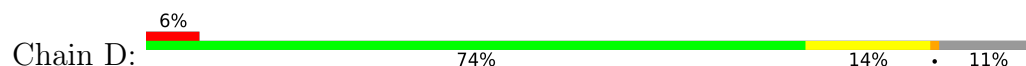
• Molecule 1: Envelopment polypeptide



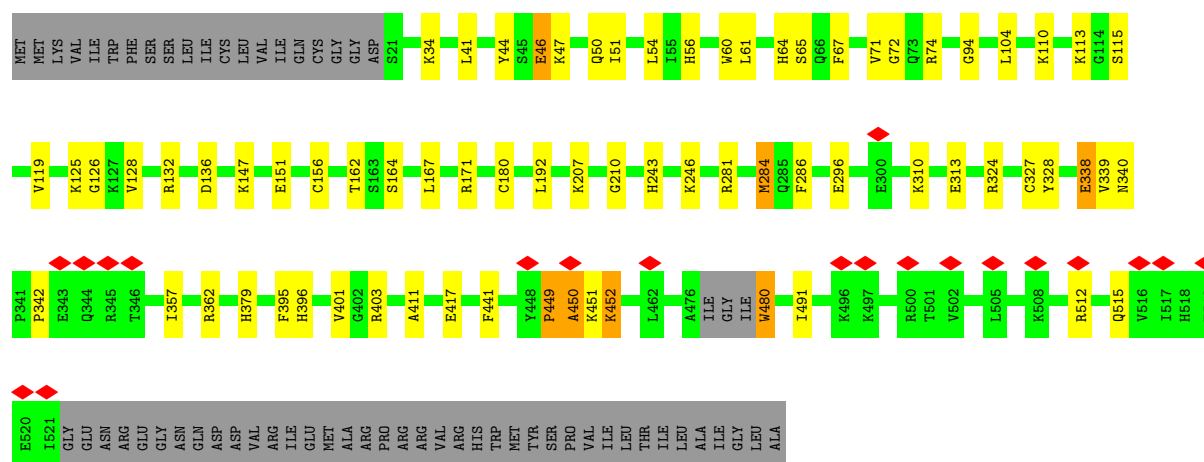




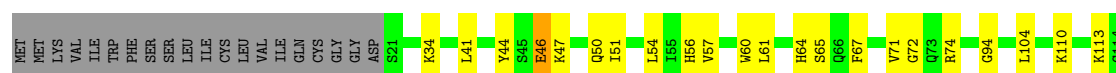
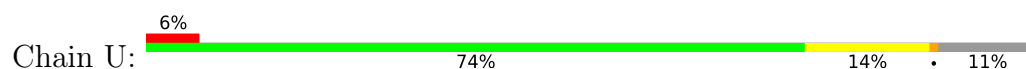
• Molecule 2: Envelopment polypeptide

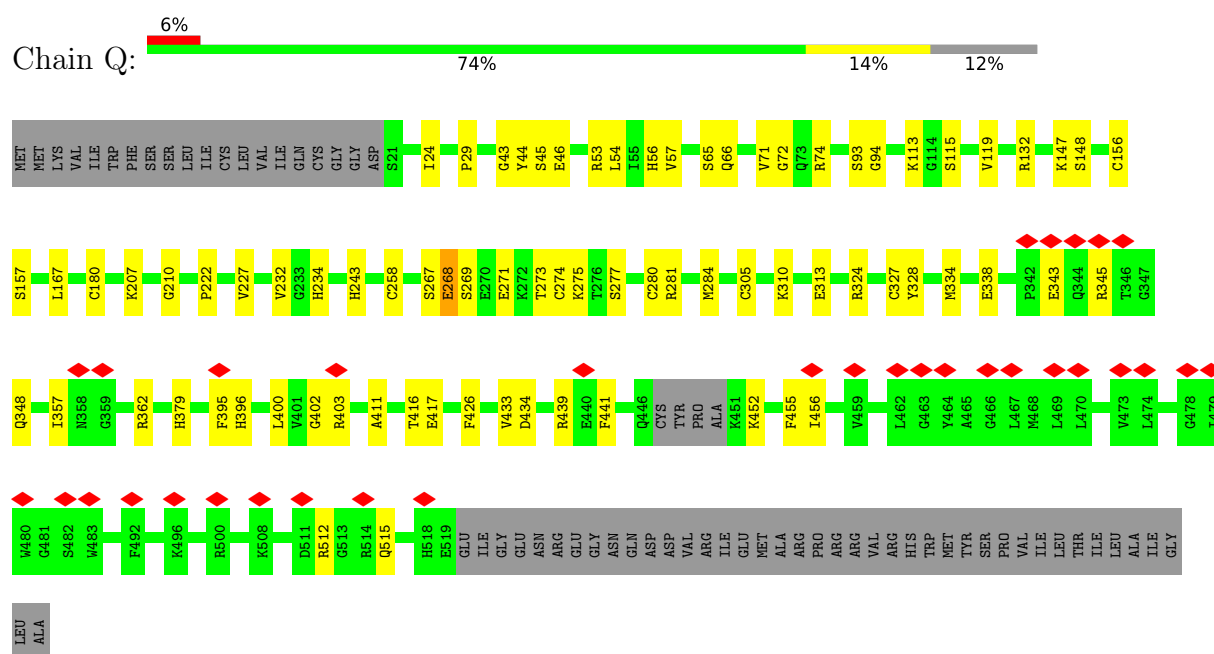


• Molecule 2: Envelopment polypeptide

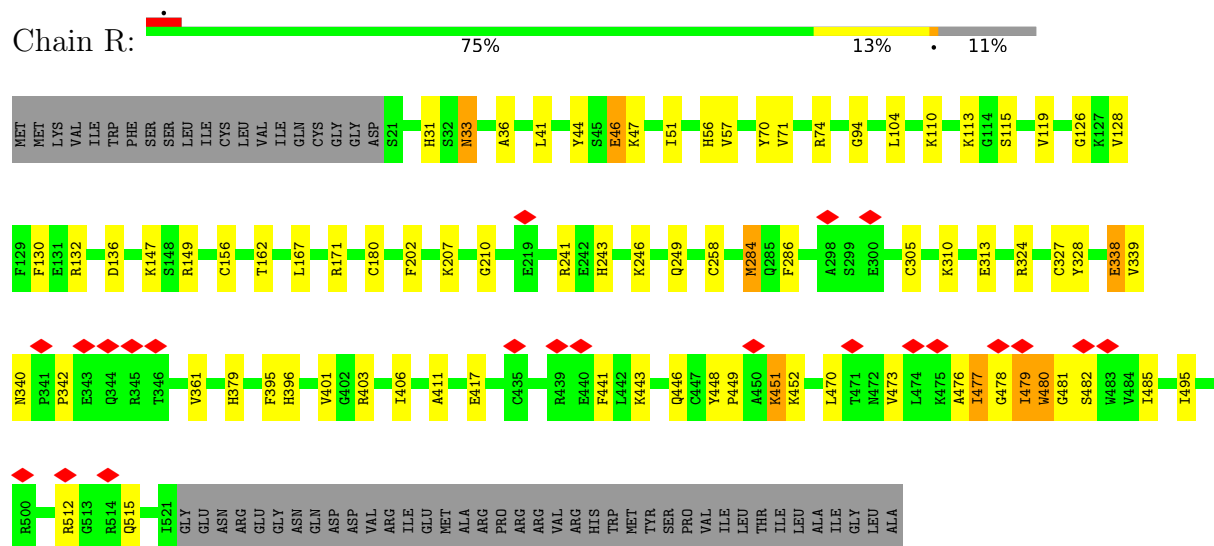


• Molecule 2: Envelopment polypeptide

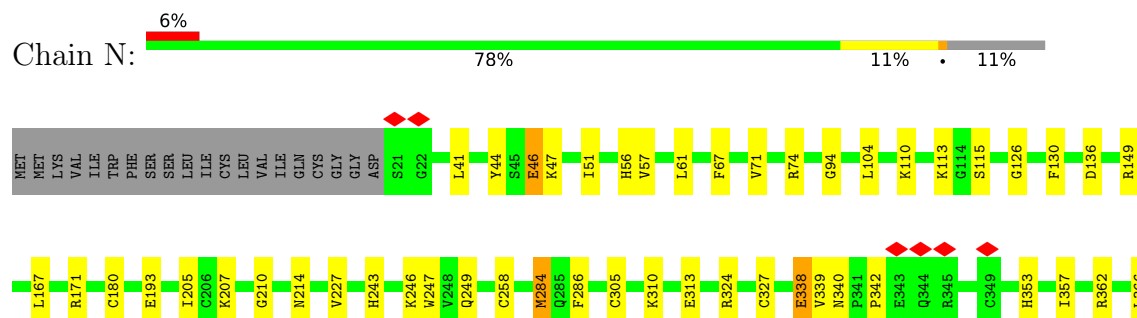


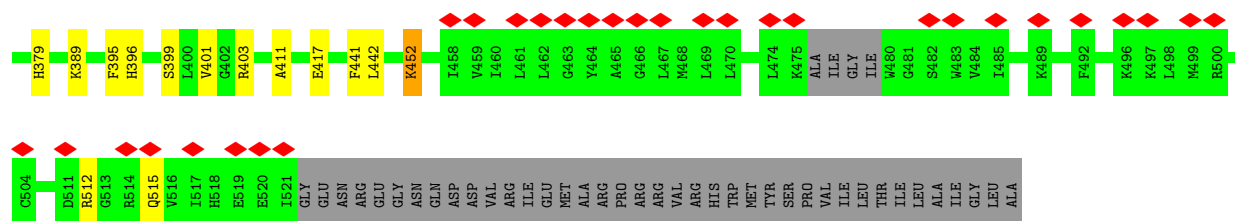


- Molecule 2: Envelopment polypeptide



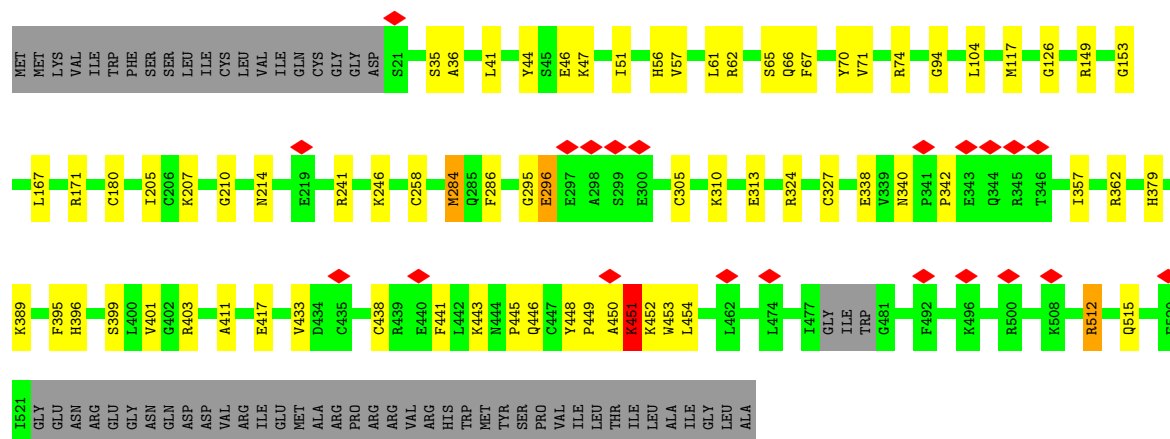
- Molecule 2: Envelopment polypeptide





• Molecule 2: Envelopment polyprotein

Chain S: 76% 12% 11%



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Y: 50% 50%



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Z: 50% 50%



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain a: 50% 50%



• Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain b:  50% 50%

MAG1
MAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain c:  50% 50%

MAG1
MAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain d:  100%

MAG1
MAG2

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	90188	Depositor
Resolution determination method	OTHER	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)	1000	Depositor
Maximum defocus (nm)	3500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	15.161	Depositor
Minimum map value	-5.531	Depositor
Average map value	0.039	Depositor
Map value standard deviation	0.437	Depositor
Recommended contour level	1.1	Depositor
Map size (Å)	503.99997, 503.99997, 503.99997	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.68, 1.68, 1.68	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: 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	B	0.25	0/3886	0.44	0/5258
1	G	0.32	0/3886	0.46	0/5258
1	H	0.24	0/3855	0.43	0/5215
1	I	0.25	0/3886	0.44	0/5258
1	J	0.25	0/3886	0.44	0/5258
1	K	0.25	0/3886	0.44	0/5258
1	L	0.25	0/3886	0.44	0/5258
1	O	0.24	0/3886	0.43	0/5258
1	P	0.24	0/3886	0.43	0/5258
1	T	0.24	0/3859	0.44	0/5220
1	V	0.25	0/3886	0.44	0/5258
1	W	0.25	0/3886	0.43	0/5258
2	A	0.25	0/3902	0.41	0/5263
2	C	0.24	0/3884	0.41	0/5241
2	D	0.24	0/3906	0.41	0/5270
2	E	0.24	0/3906	0.41	0/5270
2	F	0.25	0/3561	0.42	0/4808
2	M	0.24	0/3910	0.40	0/5277
2	N	0.24	0/3901	0.40	0/5263
2	Q	0.25	0/3877	0.41	0/5229
2	R	0.25	0/3927	0.41	0/5300
2	S	0.24	0/3898	0.41	0/5258
2	U	0.25	0/3892	0.41	0/5252
2	X	0.24	0/3906	0.41	0/5270
All	All	0.25	0/93044	0.42	0/125716

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	3808	0	3752	73	0
1	G	3808	0	3750	75	0
1	H	3778	0	3723	56	0
1	I	3808	0	3752	67	0
1	J	3808	0	3752	65	0
1	K	3808	0	3752	56	0
1	L	3808	0	3752	56	0
1	O	3808	0	3752	58	0
1	P	3808	0	3752	49	0
1	T	3782	0	3725	63	0
1	V	3808	0	3752	54	0
1	W	3808	0	3752	78	0
2	A	3815	0	3757	57	0
2	C	3795	0	3724	42	0
2	D	3817	0	3753	68	0
2	E	3817	0	3753	41	0
2	F	3478	0	3388	69	0
2	M	3820	0	3762	46	0
2	N	3812	0	3748	48	0
2	Q	3790	0	3741	41	0
2	R	3837	0	3779	52	0
2	S	3811	0	3754	61	0
2	U	3803	0	3735	48	0
2	X	3817	0	3753	44	0
3	Y	28	0	25	1	0
3	Z	28	0	25	0	0
3	a	28	0	25	0	0
3	b	28	0	25	0	0
3	c	28	0	25	0	0
3	d	28	0	25	0	0
4	A	28	0	26	0	0
4	B	42	0	39	0	0
4	C	28	0	26	0	0
4	D	28	0	26	0	0
4	E	28	0	26	0	0
4	F	14	0	13	0	0
4	G	42	0	39	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	H	42	0	39	1	0
4	I	42	0	39	0	0
4	J	42	0	39	1	0
4	K	42	0	39	0	0
4	L	42	0	39	0	0
4	M	14	0	13	0	0
4	N	14	0	13	0	0
4	O	42	0	39	1	0
4	P	42	0	39	1	0
4	Q	14	0	13	0	0
4	R	14	0	13	0	0
4	S	14	0	13	1	0
4	T	42	0	39	2	0
4	U	28	0	26	0	0
4	V	42	0	39	1	0
4	W	42	0	39	1	0
4	X	28	0	26	0	0
All	All	91976	0	90465	1223	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 7.

The worst 5 of 1223 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:F:61:LEU:HD21	2:F:129:PHE:CE1	1.38	1.56
1:G:653:ALA:HB2	2:F:309:HIS:NE2	1.55	1.21
1:I:911:GLU:OE2	2:M:389:LYS:HG3	1.40	1.18
2:A:448:TYR:H	2:A:449:PRO:CD	1.58	1.16
2:D:366:LEU:CD1	1:W:930:LYS:HD2	1.77	1.15

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	B	507/513 (99%)	449 (89%)	58 (11%)	0	100	100
1	G	507/513 (99%)	455 (90%)	48 (10%)	4 (1%)	16	53
1	H	500/513 (98%)	443 (89%)	57 (11%)	0	100	100
1	I	507/513 (99%)	445 (88%)	60 (12%)	2 (0%)	30	68
1	J	507/513 (99%)	450 (89%)	56 (11%)	1 (0%)	44	78
1	K	507/513 (99%)	448 (88%)	59 (12%)	0	100	100
1	L	507/513 (99%)	448 (88%)	59 (12%)	0	100	100
1	O	507/513 (99%)	447 (88%)	60 (12%)	0	100	100
1	P	507/513 (99%)	458 (90%)	49 (10%)	0	100	100
1	T	501/513 (98%)	443 (88%)	58 (12%)	0	100	100
1	V	507/513 (99%)	455 (90%)	52 (10%)	0	100	100
1	W	507/513 (99%)	456 (90%)	51 (10%)	0	100	100
2	A	495/560 (88%)	477 (96%)	16 (3%)	2 (0%)	30	68
2	C	491/560 (88%)	474 (96%)	17 (4%)	0	100	100
2	D	494/560 (88%)	474 (96%)	20 (4%)	0	100	100
2	E	494/560 (88%)	475 (96%)	16 (3%)	3 (1%)	22	60
2	F	454/560 (81%)	420 (92%)	32 (7%)	2 (0%)	30	68
2	M	497/560 (89%)	479 (96%)	18 (4%)	0	100	100
2	N	493/560 (88%)	475 (96%)	18 (4%)	0	100	100
2	Q	491/560 (88%)	469 (96%)	22 (4%)	0	100	100
2	R	499/560 (89%)	477 (96%)	20 (4%)	2 (0%)	30	68
2	S	494/560 (88%)	475 (96%)	17 (3%)	2 (0%)	30	68
2	U	492/560 (88%)	472 (96%)	18 (4%)	2 (0%)	30	68
2	X	494/560 (88%)	476 (96%)	18 (4%)	0	100	100
All	All	11959/12876 (93%)	11040 (92%)	899 (8%)	20 (0%)	45	78

5 of 20 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	651	ARG
1	G	657	GLN
1	G	660	CYS

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Mol	Chain	Res	Type
2	F	80	VAL
2	A	446	GLN

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	B	425/428 (99%)	398 (94%)	27 (6%)	14	36
1	G	425/428 (99%)	393 (92%)	32 (8%)	11	31
1	H	422/428 (99%)	396 (94%)	26 (6%)	15	37
1	I	425/428 (99%)	400 (94%)	25 (6%)	16	38
1	J	425/428 (99%)	399 (94%)	26 (6%)	15	37
1	K	425/428 (99%)	400 (94%)	25 (6%)	16	38
1	L	425/428 (99%)	399 (94%)	26 (6%)	15	37
1	O	425/428 (99%)	397 (93%)	28 (7%)	14	35
1	P	425/428 (99%)	402 (95%)	23 (5%)	18	40
1	T	422/428 (99%)	399 (94%)	23 (6%)	18	40
1	V	425/428 (99%)	403 (95%)	22 (5%)	19	41
1	W	425/428 (99%)	403 (95%)	22 (5%)	19	41
2	A	426/479 (89%)	412 (97%)	14 (3%)	33	53
2	C	424/479 (88%)	414 (98%)	10 (2%)	44	62
2	D	426/479 (89%)	413 (97%)	13 (3%)	35	55
2	E	426/479 (89%)	413 (97%)	13 (3%)	35	55
2	F	389/479 (81%)	374 (96%)	15 (4%)	27	48
2	M	426/479 (89%)	412 (97%)	14 (3%)	33	53
2	N	426/479 (89%)	416 (98%)	10 (2%)	45	64
2	Q	423/479 (88%)	415 (98%)	8 (2%)	52	69
2	R	428/479 (89%)	415 (97%)	13 (3%)	36	56
2	S	426/479 (89%)	414 (97%)	12 (3%)	38	58

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	U	425/479 (89%)	414 (97%)	11 (3%)	41	60
2	X	426/479 (89%)	412 (97%)	14 (3%)	33	53
All	All	10165/10884 (93%)	9713 (96%)	452 (4%)	26	46

5 of 452 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	K	935	LYS
1	W	835	ARG
1	T	737	THR
1	W	684	PHE
1	P	616	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 113 such sidechains are listed below:

Mol	Chain	Res	Type
2	X	66	GLN
1	W	755	GLN
2	M	56	HIS
1	W	626	ASN
2	N	446	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

12 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	$\# Z > 2$
3	NAG	Y	1	2,3	14,14,15	0.52	0	17,19,21	1.30	1 (5%)
3	NAG	Y	2	3	14,14,15	0.22	0	17,19,21	0.42	0
3	NAG	Z	1	2,3	14,14,15	0.20	0	17,19,21	0.41	0
3	NAG	Z	2	3	14,14,15	0.98	1 (7%)	17,19,21	1.36	1 (5%)
3	NAG	a	1	2,3	14,14,15	0.44	0	17,19,21	0.35	0
3	NAG	a	2	3	14,14,15	0.99	1 (7%)	17,19,21	1.27	1 (5%)
3	NAG	b	1	2,3	14,14,15	0.44	0	17,19,21	1.26	2 (11%)
3	NAG	b	2	3	14,14,15	0.38	0	17,19,21	0.35	0
3	NAG	c	1	2,3	14,14,15	0.33	0	17,19,21	0.63	0
3	NAG	c	2	3	14,14,15	0.51	0	17,19,21	1.27	2 (11%)
3	NAG	d	1	2,3	14,14,15	0.24	0	17,19,21	0.52	0
3	NAG	d	2	3	14,14,15	0.22	0	17,19,21	0.43	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
3	NAG	Y	1	2,3	-	5/6/23/26	0/1/1/1
3	NAG	Y	2	3	-	2/6/23/26	0/1/1/1
3	NAG	Z	1	2,3	-	0/6/23/26	0/1/1/1
3	NAG	Z	2	3	-	3/6/23/26	0/1/1/1
3	NAG	a	1	2,3	-	4/6/23/26	0/1/1/1
3	NAG	a	2	3	-	3/6/23/26	0/1/1/1
3	NAG	b	1	2,3	-	5/6/23/26	0/1/1/1
3	NAG	b	2	3	-	0/6/23/26	0/1/1/1
3	NAG	c	1	2,3	-	3/6/23/26	0/1/1/1
3	NAG	c	2	3	-	5/6/23/26	0/1/1/1
3	NAG	d	1	2,3	-	3/6/23/26	0/1/1/1
3	NAG	d	2	3	-	4/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	Z	2	NAG	O5-C1	3.54	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	a	2	NAG	O5-C1	3.50	1.49	1.43

The worst 5 of 7 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	Z	2	NAG	C1-O5-C5	5.16	119.18	112.19
3	a	2	NAG	C1-O5-C5	4.85	118.76	112.19
3	Y	1	NAG	C2-N2-C7	4.33	129.07	122.90
3	c	2	NAG	C2-N2-C7	4.30	129.02	122.90
3	b	1	NAG	C2-N2-C7	4.26	128.97	122.90

There are no chirality outliers.

5 of 37 torsion outliers are listed below:

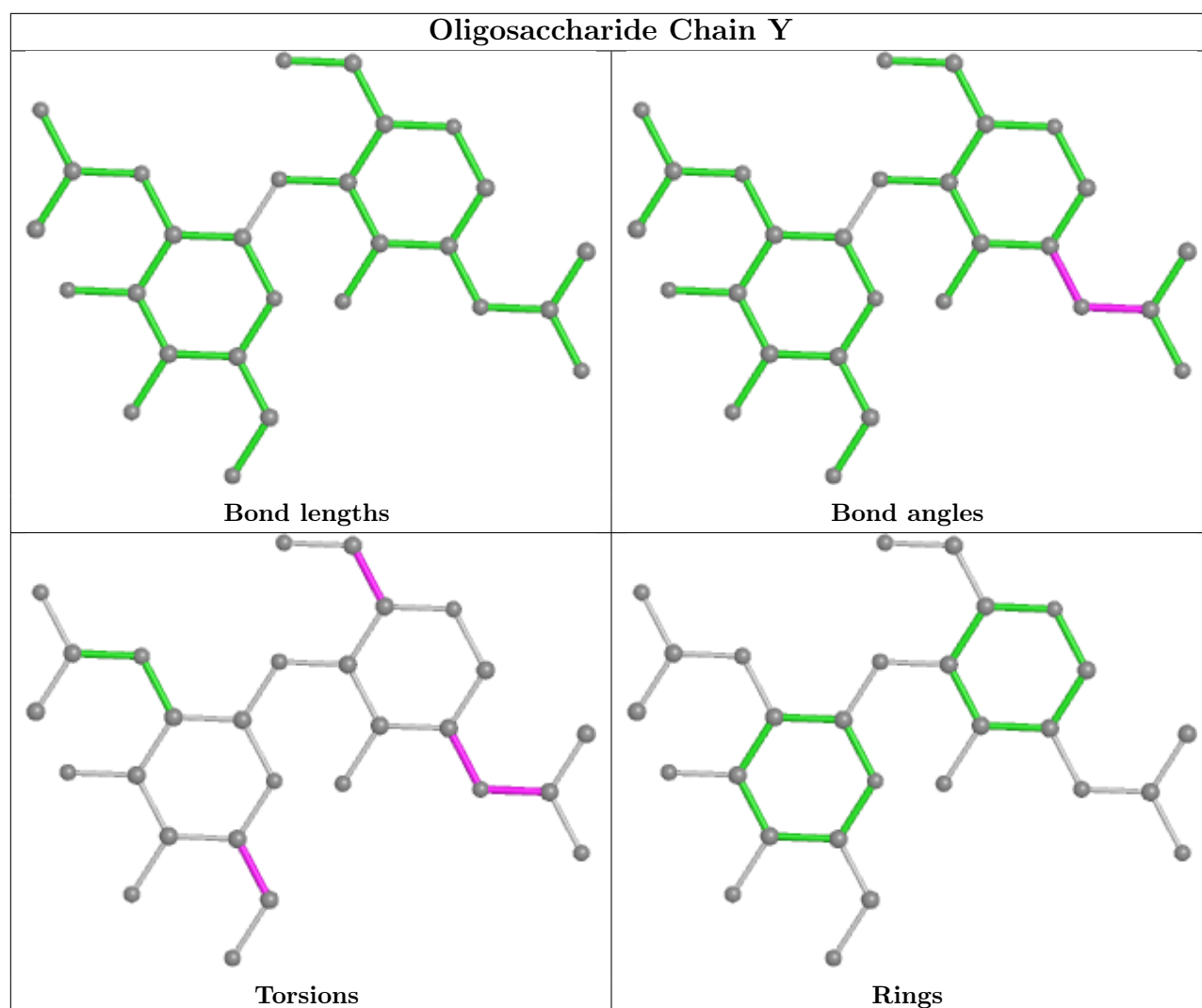
Mol	Chain	Res	Type	Atoms
3	Y	2	NAG	C4-C5-C6-O6
3	Y	1	NAG	C4-C5-C6-O6
3	Z	2	NAG	O5-C5-C6-O6
3	a	2	NAG	O5-C5-C6-O6
3	c	2	NAG	C4-C5-C6-O6

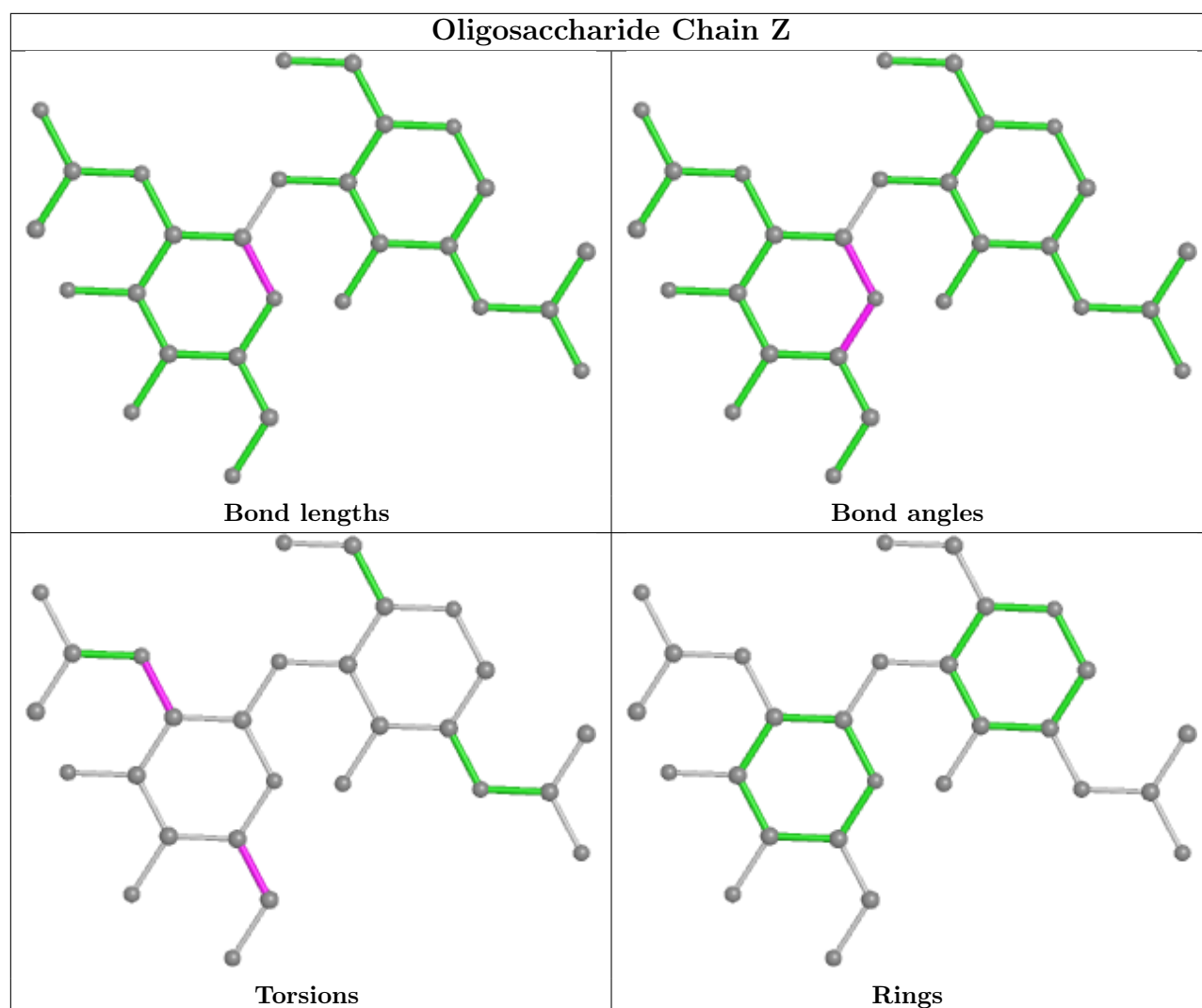
There are no ring outliers.

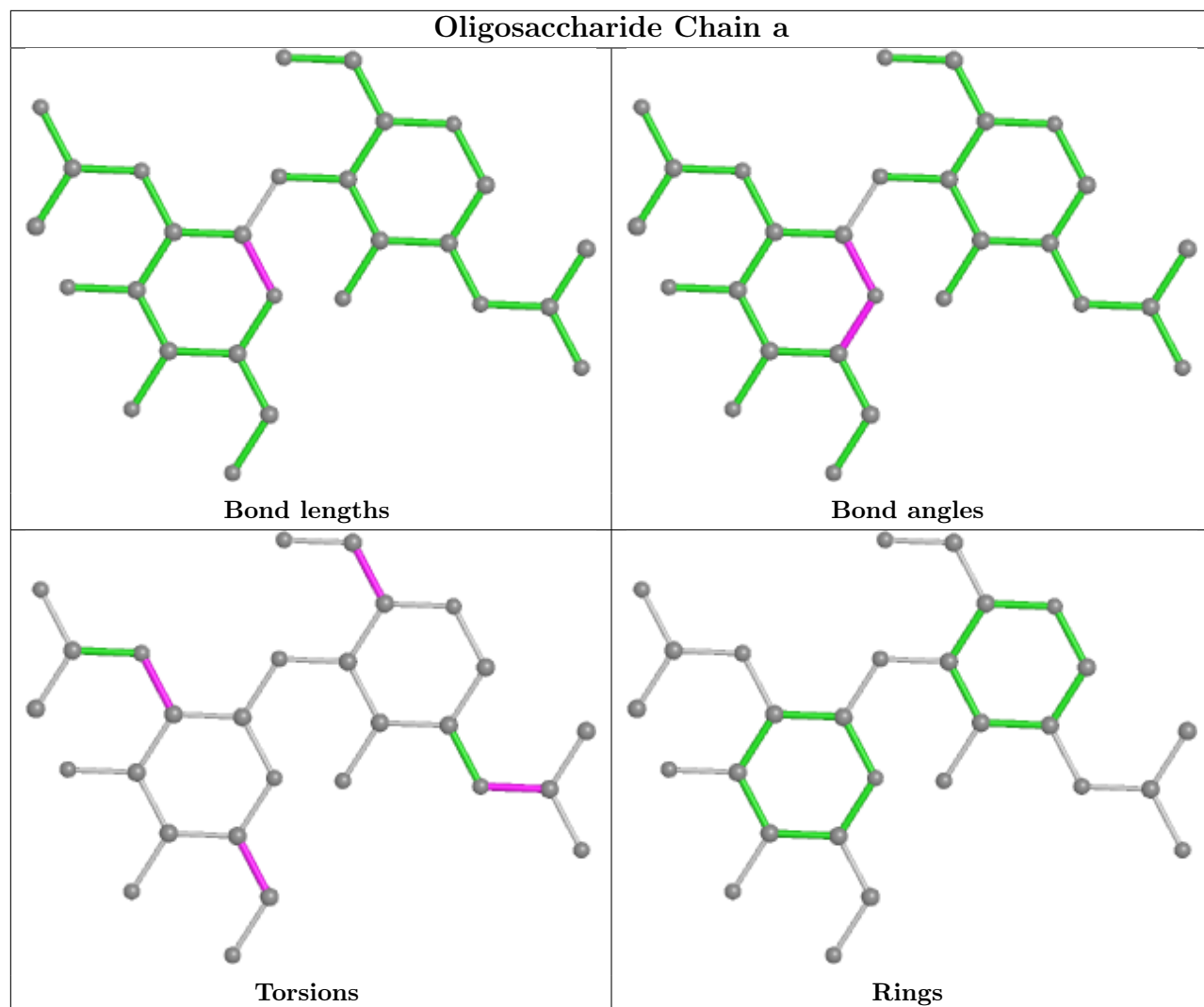
1 monomer is involved in 1 short contact:

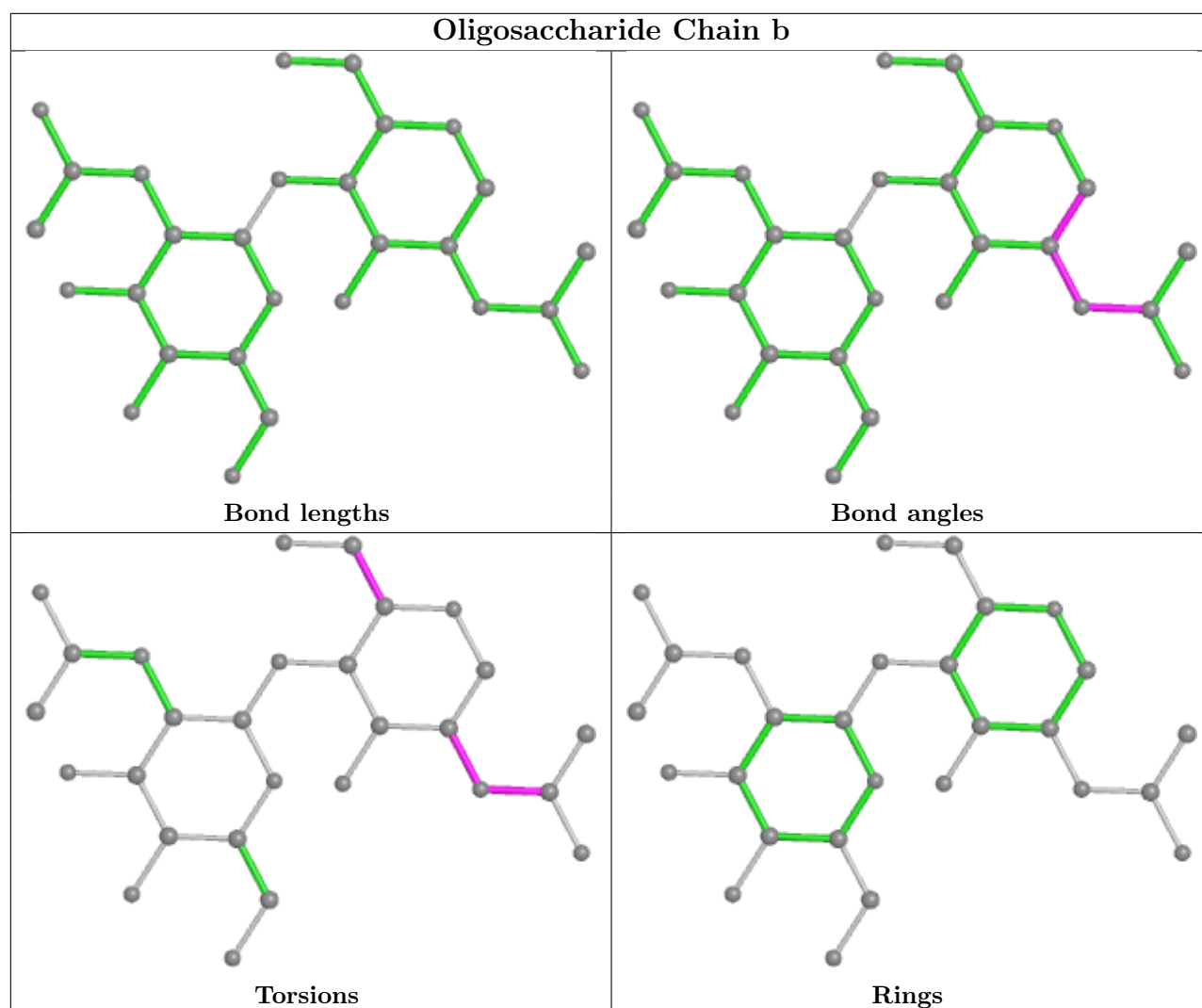
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	Y	1	NAG	1	0

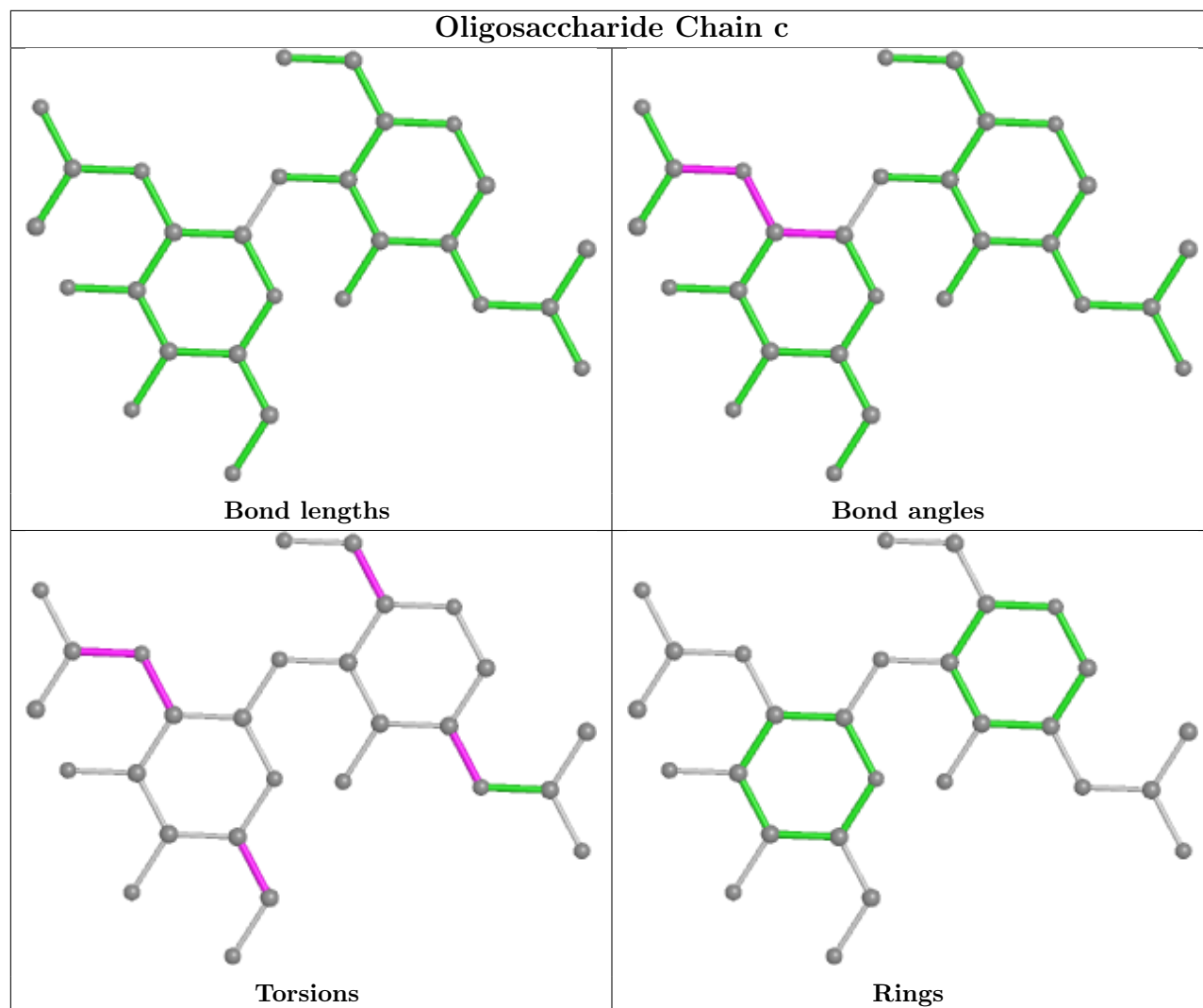
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

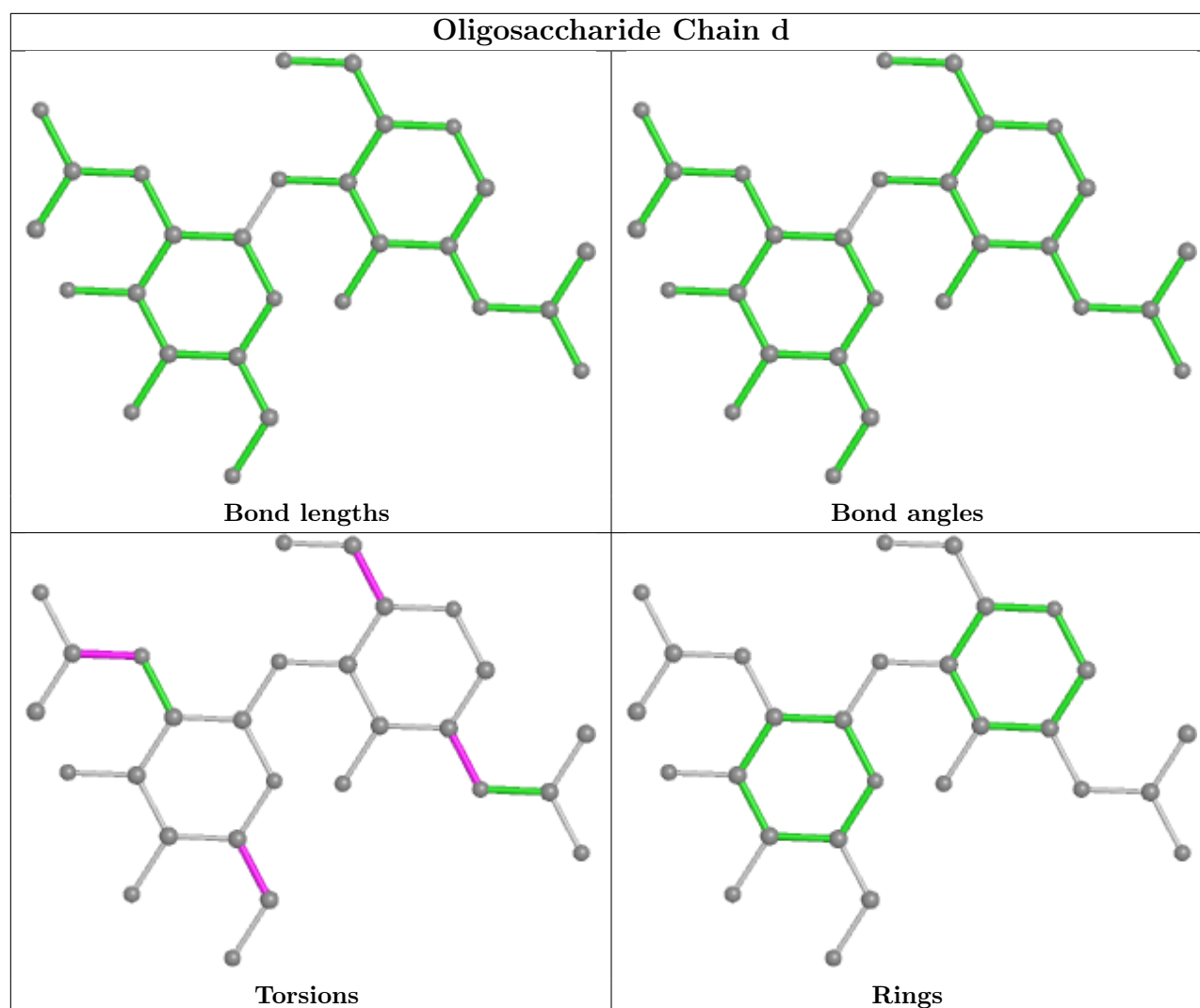












5.6 Ligand geometry [i](#)

54 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	$\# Z > 2$
4	NAG	V	1102	1	14,14,15	0.24	0	17,19,21	0.45	0
4	NAG	C	602	2	14,14,15	0.25	0	17,19,21	0.45	0
4	NAG	N	601	2	14,14,15	0.22	0	17,19,21	0.41	0
4	NAG	B	1102	1	14,14,15	0.22	0	17,19,21	0.42	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	K	1102	1	14,14,15	0.25	0	17,19,21	0.42	0
4	NAG	I	1102	1	14,14,15	0.20	0	17,19,21	0.43	0
4	NAG	T	1103	1	14,14,15	0.44	0	17,19,21	1.26	1 (5%)
4	NAG	Q	601	2	14,14,15	0.24	0	17,19,21	0.43	0
4	NAG	A	602	2	14,14,15	0.20	0	17,19,21	0.44	0
4	NAG	W	1101	1	14,14,15	0.20	0	17,19,21	0.42	0
4	NAG	B	1103	1	14,14,15	0.20	0	17,19,21	0.42	0
4	NAG	H	1103	1	14,14,15	0.24	0	17,19,21	0.44	0
4	NAG	D	602	2	14,14,15	0.22	0	17,19,21	0.45	0
4	NAG	I	1103	1	14,14,15	0.23	0	17,19,21	0.44	0
4	NAG	G	1102	1	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	J	1102	1	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	L	1103	1	14,14,15	0.23	0	17,19,21	0.43	0
4	NAG	T	1102	1	14,14,15	0.19	0	17,19,21	0.42	0
4	NAG	P	1102	1	14,14,15	0.22	0	17,19,21	0.39	0
4	NAG	W	1102	1	14,14,15	0.29	0	17,19,21	0.44	0
4	NAG	K	1103	1	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	F	601	2	14,14,15	0.26	0	17,19,21	0.53	0
4	NAG	G	1101	1	14,14,15	0.71	1 (7%)	17,19,21	1.52	3 (17%)
4	NAG	U	601	2	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	C	601	2	14,14,15	0.23	0	17,19,21	0.43	0
4	NAG	H	1101	1	14,14,15	0.89	1 (7%)	17,19,21	1.02	1 (5%)
4	NAG	O	1102	1	14,14,15	0.25	0	17,19,21	0.42	0
4	NAG	W	1103	1	14,14,15	0.27	0	17,19,21	0.53	0
4	NAG	J	1103	1	14,14,15	0.24	0	17,19,21	0.42	0
4	NAG	L	1101	1	14,14,15	0.88	1 (7%)	17,19,21	1.02	1 (5%)
4	NAG	V	1101	1	14,14,15	0.23	0	17,19,21	0.41	0
4	NAG	A	601	2	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	B	1101	1	14,14,15	0.91	1 (7%)	17,19,21	1.03	1 (5%)
4	NAG	K	1101	1	14,14,15	0.90	1 (7%)	17,19,21	1.02	1 (5%)
4	NAG	E	601	2	14,14,15	0.23	0	17,19,21	0.44	0
4	NAG	L	1102	1	14,14,15	0.23	0	17,19,21	0.42	0
4	NAG	O	1103	1	14,14,15	0.41	0	17,19,21	0.39	0
4	NAG	P	1101	1	14,14,15	0.66	1 (7%)	17,19,21	0.75	0
4	NAG	E	602	2	14,14,15	0.18	0	17,19,21	0.45	0
4	NAG	J	1101	1	14,14,15	0.89	1 (7%)	17,19,21	1.04	1 (5%)
4	NAG	G	1103	1	14,14,15	0.44	0	17,19,21	1.25	1 (5%)
4	NAG	X	602	2	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	S	601	2	14,14,15	0.40	0	17,19,21	0.37	0
4	NAG	T	1101	1	14,14,15	0.94	1 (7%)	17,19,21	1.62	3 (17%)
4	NAG	D	601	2	14,14,15	0.21	0	17,19,21	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	V	1103	1	14,14,15	0.43	0	17,19,21	0.36	0
4	NAG	O	1101	1	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	M	601	2	14,14,15	0.42	0	17,19,21	0.36	0
4	NAG	R	601	2	14,14,15	0.25	0	17,19,21	0.42	0
4	NAG	P	1103	1	14,14,15	0.45	0	17,19,21	1.26	1 (5%)
4	NAG	I	1101	1	14,14,15	0.90	1 (7%)	17,19,21	1.03	1 (5%)
4	NAG	H	1102	1	14,14,15	0.21	0	17,19,21	0.42	0
4	NAG	X	601	2	14,14,15	0.22	0	17,19,21	0.42	0
4	NAG	U	602	2	14,14,15	0.24	0	17,19,21	0.46	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
4	NAG	V	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	C	602	2	-	2/6/23/26	0/1/1/1
4	NAG	N	601	2	-	3/6/23/26	0/1/1/1
4	NAG	B	1102	1	-	0/6/23/26	0/1/1/1
4	NAG	K	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	I	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	T	1103	1	-	4/6/23/26	0/1/1/1
4	NAG	Q	601	2	-	0/6/23/26	0/1/1/1
4	NAG	A	602	2	-	2/6/23/26	0/1/1/1
4	NAG	W	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	B	1103	1	-	2/6/23/26	0/1/1/1
4	NAG	H	1103	1	-	2/6/23/26	0/1/1/1
4	NAG	D	602	2	-	2/6/23/26	0/1/1/1
4	NAG	I	1103	1	-	2/6/23/26	0/1/1/1
4	NAG	G	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	J	1102	1	-	0/6/23/26	0/1/1/1
4	NAG	L	1103	1	-	2/6/23/26	0/1/1/1
4	NAG	T	1102	1	-	0/6/23/26	0/1/1/1
4	NAG	P	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	W	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	K	1103	1	-	2/6/23/26	0/1/1/1
4	NAG	F	601	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	G	1101	1	-	3/6/23/26	0/1/1/1
4	NAG	U	601	2	-	2/6/23/26	0/1/1/1
4	NAG	C	601	2	-	3/6/23/26	0/1/1/1
4	NAG	H	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	O	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	W	1103	1	-	1/6/23/26	0/1/1/1
4	NAG	J	1103	1	-	2/6/23/26	0/1/1/1
4	NAG	L	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	V	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	A	601	2	-	4/6/23/26	0/1/1/1
4	NAG	B	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	K	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	E	601	2	-	4/6/23/26	0/1/1/1
4	NAG	L	1102	1	-	2/6/23/26	0/1/1/1
4	NAG	O	1103	1	-	2/6/23/26	0/1/1/1
4	NAG	P	1101	1	-	4/6/23/26	0/1/1/1
4	NAG	E	602	2	-	2/6/23/26	0/1/1/1
4	NAG	J	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	G	1103	1	-	5/6/23/26	0/1/1/1
4	NAG	X	602	2	-	2/6/23/26	0/1/1/1
4	NAG	S	601	2	-	4/6/23/26	0/1/1/1
4	NAG	T	1101	1	-	5/6/23/26	0/1/1/1
4	NAG	D	601	2	-	4/6/23/26	0/1/1/1
4	NAG	V	1103	1	-	0/6/23/26	0/1/1/1
4	NAG	O	1101	1	-	1/6/23/26	0/1/1/1
4	NAG	M	601	2	-	4/6/23/26	0/1/1/1
4	NAG	R	601	2	-	4/6/23/26	0/1/1/1
4	NAG	P	1103	1	-	3/6/23/26	0/1/1/1
4	NAG	I	1101	1	-	2/6/23/26	0/1/1/1
4	NAG	H	1102	1	-	0/6/23/26	0/1/1/1
4	NAG	X	601	2	-	4/6/23/26	0/1/1/1
4	NAG	U	602	2	-	2/6/23/26	0/1/1/1

The worst 5 of 9 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1101	NAG	O5-C1	3.00	1.48	1.43
4	I	1101	NAG	O5-C1	3.00	1.48	1.43
4	K	1101	NAG	O5-C1	2.97	1.48	1.43
4	J	1101	NAG	O5-C1	2.97	1.48	1.43
4	L	1101	NAG	O5-C1	2.96	1.48	1.43

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	1101	NAG	C1-O5-C5	4.71	118.57	112.19
4	T	1101	NAG	C1-O5-C5	4.30	118.02	112.19
4	T	1101	NAG	C2-N2-C7	4.28	129.00	122.90
4	G	1103	NAG	C2-N2-C7	4.28	129.00	122.90
4	T	1103	NAG	C2-N2-C7	4.28	129.00	122.90

There are no chirality outliers.

5 of 122 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	1101	NAG	C4-C5-C6-O6
4	R	601	NAG	O5-C5-C6-O6
4	K	1101	NAG	C4-C5-C6-O6
4	G	1102	NAG	O5-C5-C6-O6
4	J	1101	NAG	O5-C5-C6-O6

There are no ring outliers.

11 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	T	1103	NAG	1	0
4	W	1101	NAG	1	0
4	J	1102	NAG	1	0
4	G	1101	NAG	2	0
4	V	1101	NAG	1	0
4	G	1103	NAG	1	0
4	S	601	NAG	1	0
4	T	1101	NAG	1	0
4	O	1101	NAG	1	0
4	P	1103	NAG	1	0
4	H	1102	NAG	1	0

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.

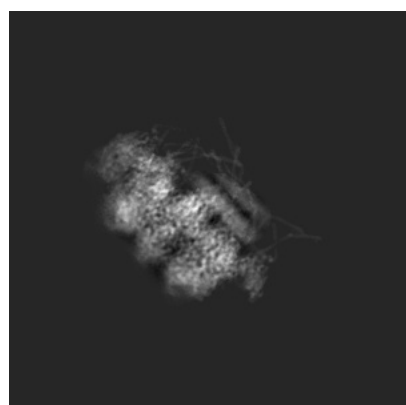
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-35183. These allow visual inspection of the internal detail of the map and identification of artifacts.

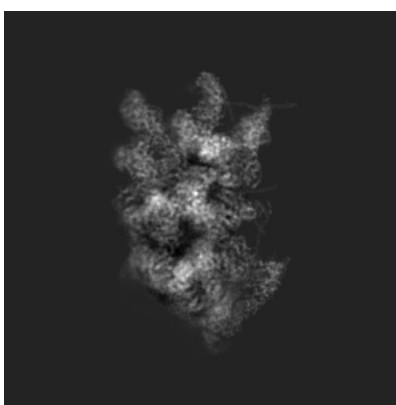
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

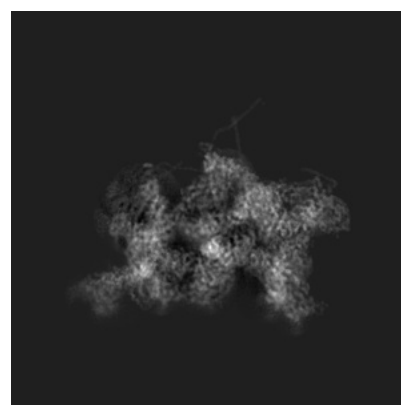
6.1.1 Primary map



X



Y

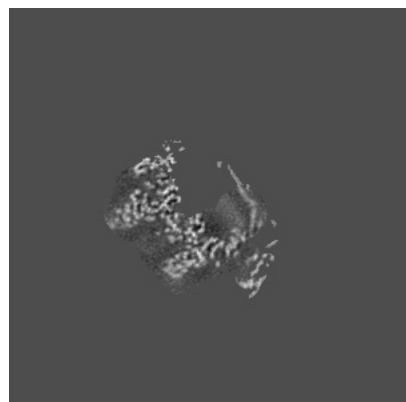


Z

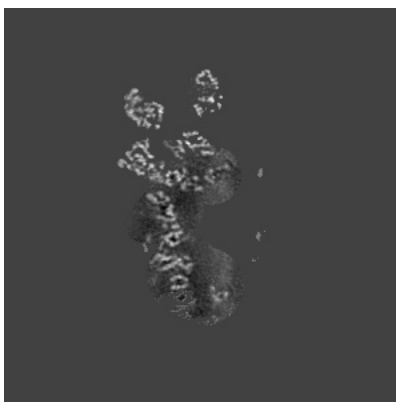
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

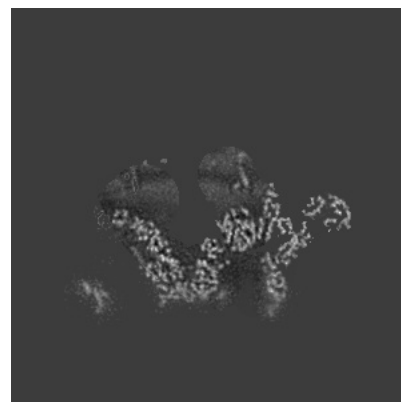
6.2.1 Primary map



X Index: 150



Y Index: 150

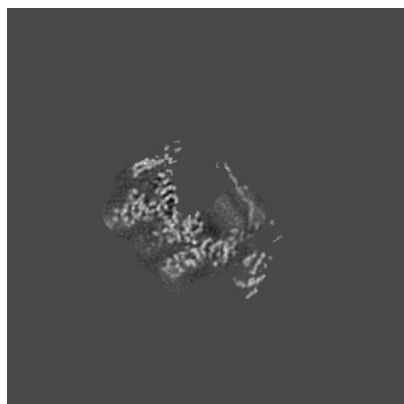


Z Index: 150

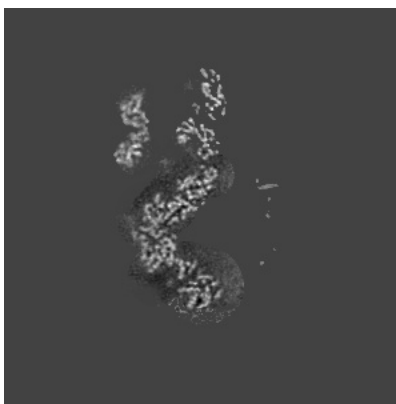
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

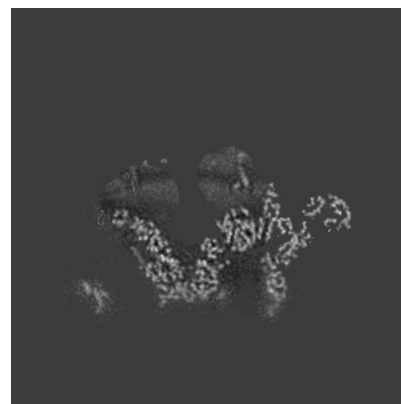
6.3.1 Primary map



X Index: 151



Y Index: 140

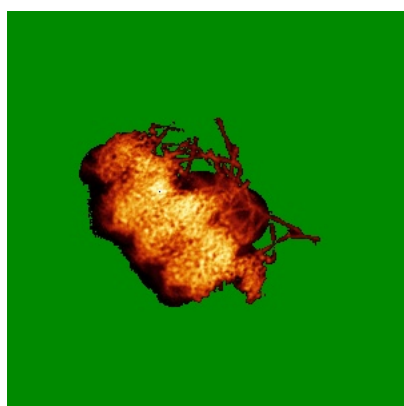


Z Index: 150

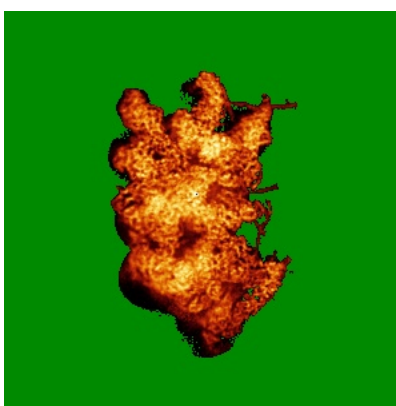
The images above show the largest variance slices of the map in three orthogonal directions.

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

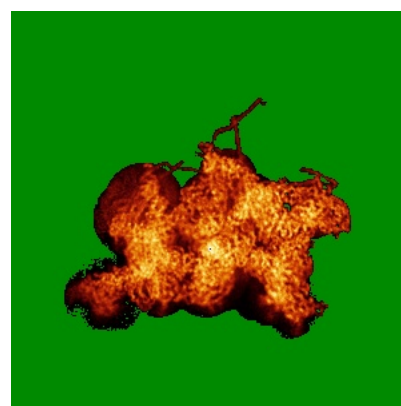
6.4.1 Primary map



X



Y



Z

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 1.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

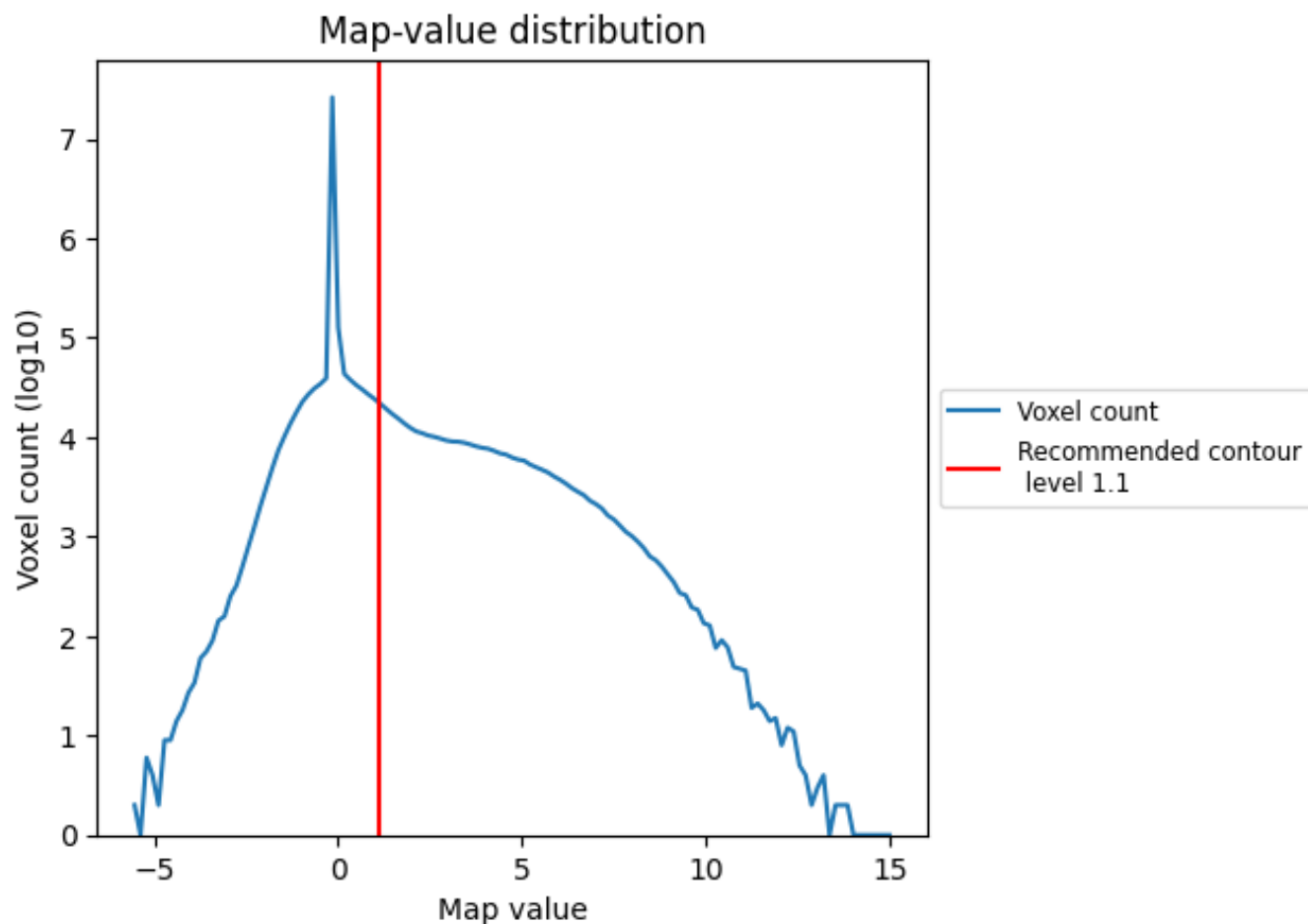
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

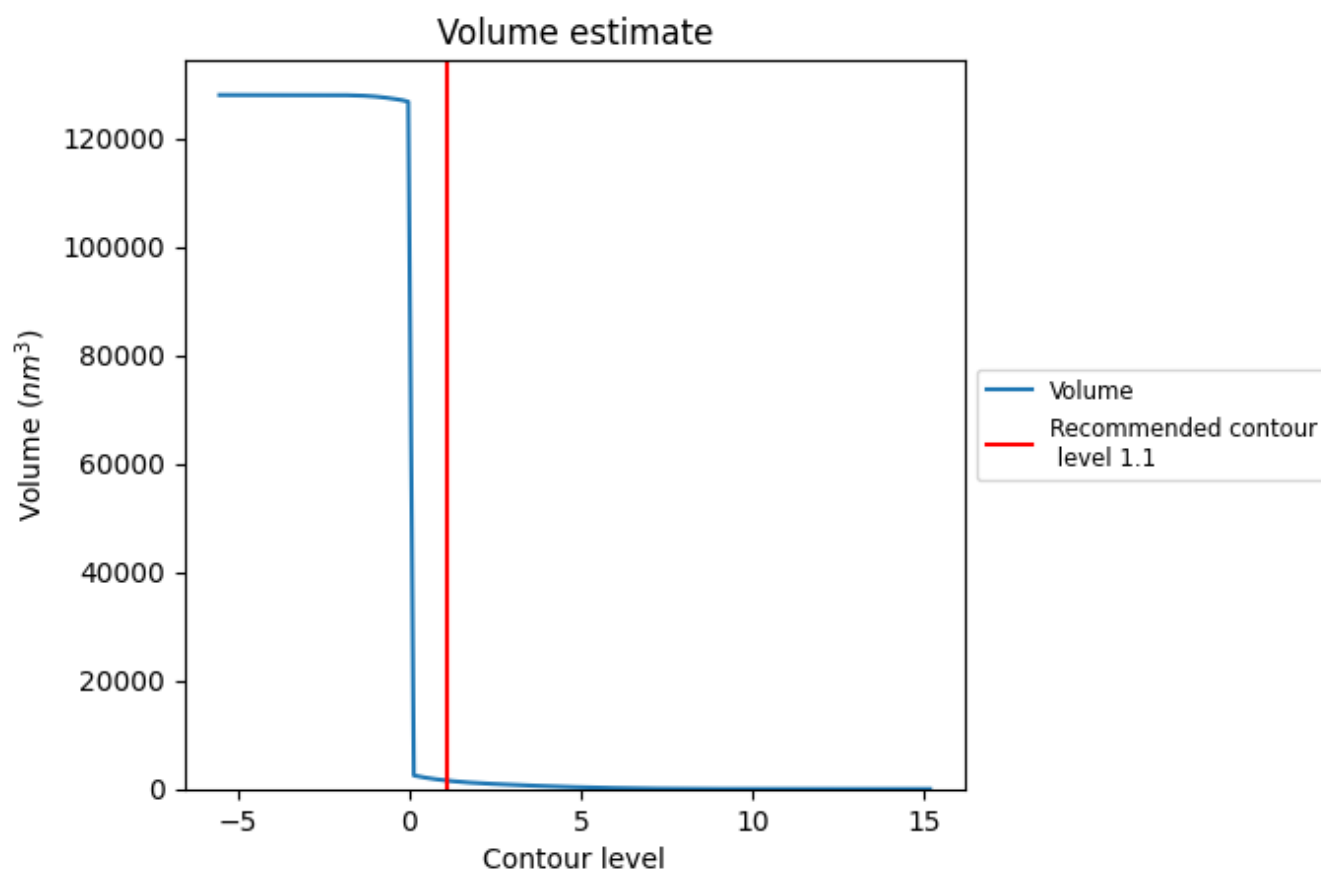
This section contains the results of statistical analysis of the map.

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.

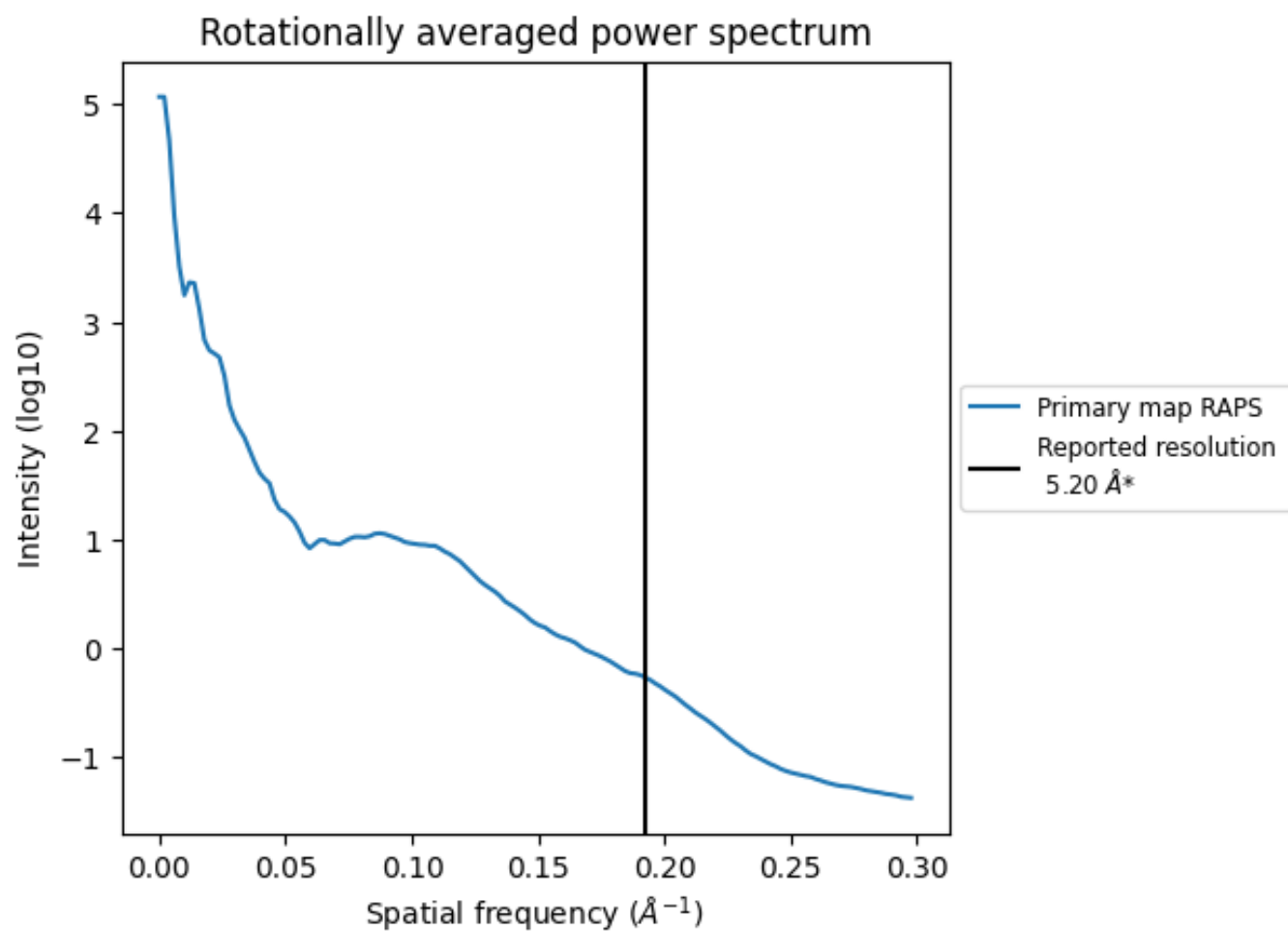
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1538 nm^3 ; this corresponds to an approximate mass of 1390 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 0.192 \AA^{-1}

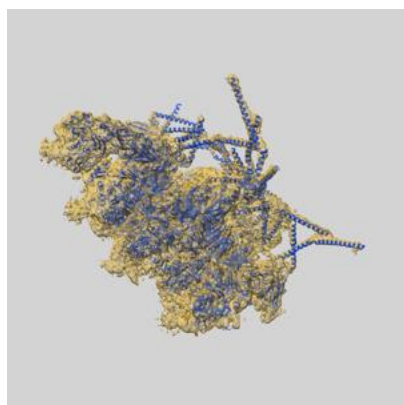
8 Fourier-Shell correlation

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

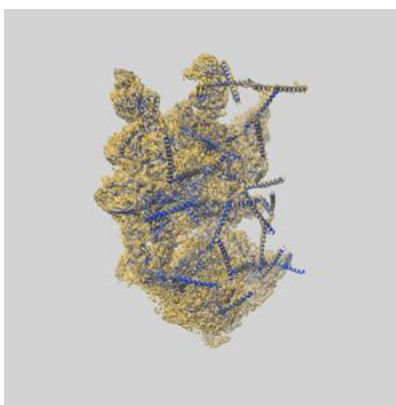
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-35183 and PDB model 8I4T. Per-residue inclusion information can be found in section [3](#) on page [9](#).

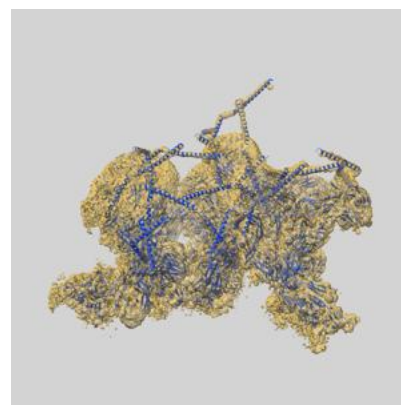
9.1 Map-model overlay [i](#)



X



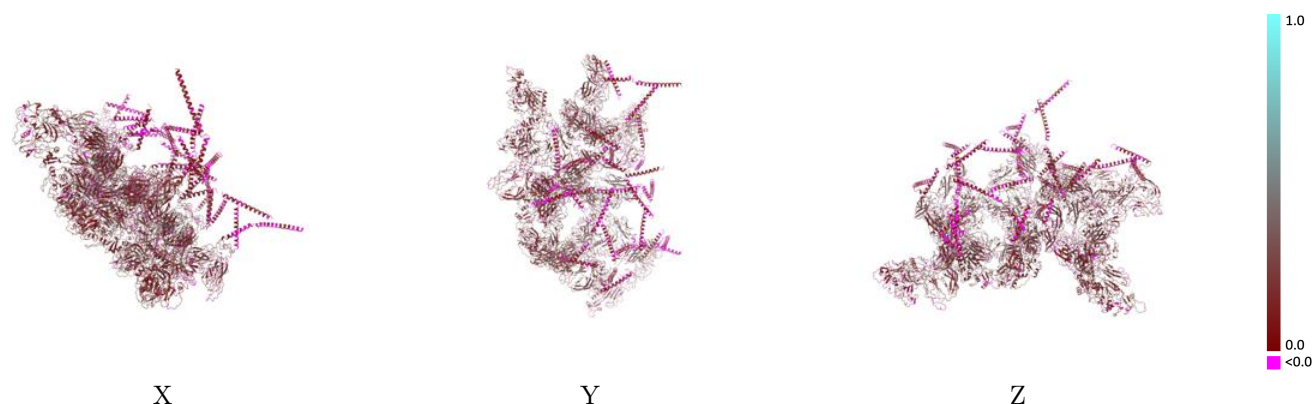
Y



Z

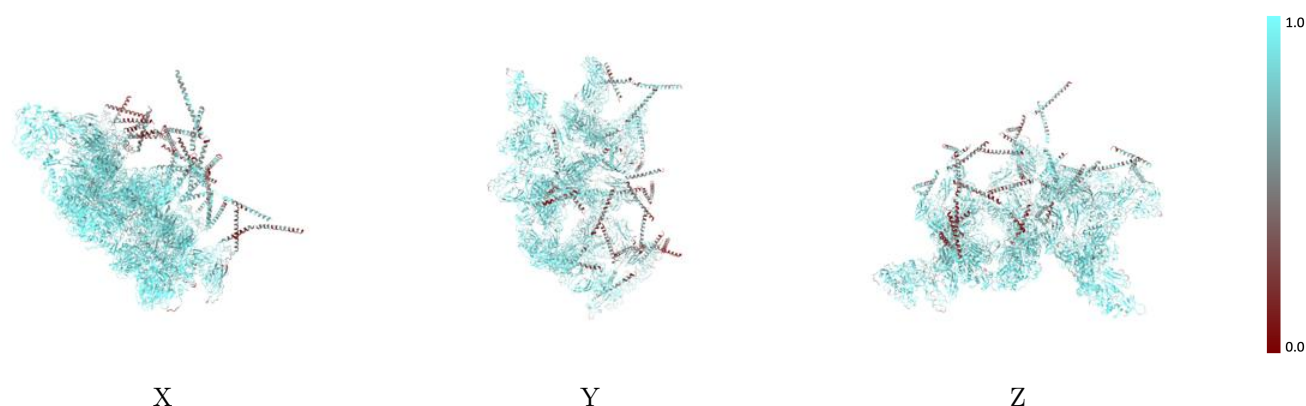
The images above show the 3D surface view of the map at the recommended contour level 1.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



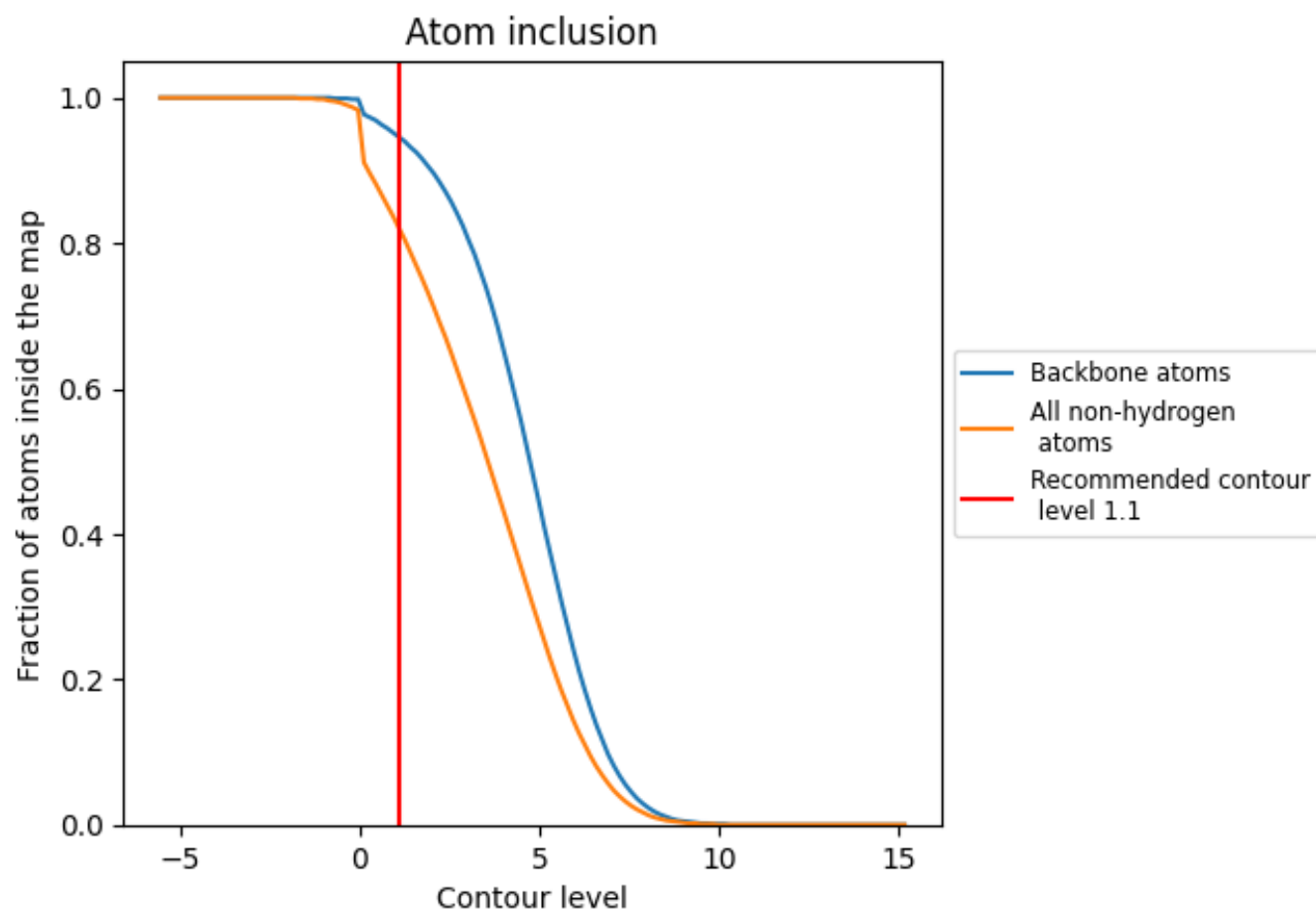
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (1.1).































































9.4 Atom inclusion [i](#)



At the recommended contour level, 95% 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 (1.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8210	 0.2060
A	 0.8070	 0.2060
B	 0.7900	 0.2180
C	 0.8250	 0.2090
D	 0.8560	 0.2180
E	 0.8630	 0.1960
F	 0.7850	 0.1700
G	 0.7310	 0.2060
H	 0.7990	 0.2130
I	 0.8620	 0.2310
J	 0.8360	 0.2110
K	 0.8380	 0.2370
L	 0.8230	 0.2210
M	 0.8650	 0.2190
N	 0.8170	 0.1990
O	 0.8090	 0.2070
P	 0.7740	 0.1800
Q	 0.8230	 0.1850
R	 0.8370	 0.2020
S	 0.8450	 0.2130
T	 0.7780	 0.1760
U	 0.8470	 0.2120
V	 0.7820	 0.1940
W	 0.8380	 0.2150
X	 0.8570	 0.2000
Y	 0.9290	 0.2180
Z	 1.0000	 0.3170
a	 1.0000	 0.2520
b	 1.0000	 0.3330
c	 1.0000	 0.2950
d	 1.0000	 0.2450

