Full wwPDB EM Validation Report (i)

Nov 6, 2022 – 10:36 AM EST

PDB ID	:	6CRX
EMDB ID	:	EMD-7575
Title	:	SARS Spike Glycoprotein, Stabilized variant, two S1 CTDs in the upwards conformation
Authors	:	Kirchdoerfer, R.N.; Wang, N.; Pallesen, J.; Turner, H.L.; Cottrell, C.A.; McLellan, J.S.; Ward, A.B.
Deposited on	:	2018-03-19
Resolution	:	3.90 Å(reported)
Based on initial models	:	5X4S, 2AJF, 5I08

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 (i) 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 (i)) were used in the production of this report:

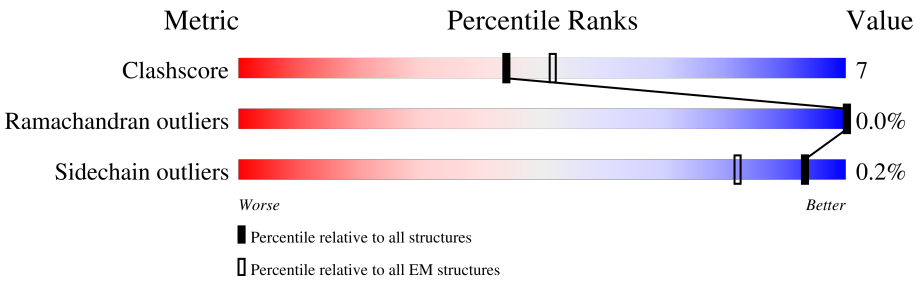
EMDB validation analysis	:	0.0.1.dev43
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.9
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.31.2

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

The reported resolution of this entry is 3.90 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 <=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	1215	22% / 71% / 17% / 12%
1	B	1215	22% / 73% / 15% / 12%
1	C	1215	19% / 70% / 18% / 12%
2	D	2	50% / 100%
2	E	2	100% / 100%
2	H	2	50% / 100%
2	I	2	100% / 100%
2	K	2	100% / 50% / 50%

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Mol	Chain	Length	Quality of chain
2	L	2	100% / 100%
2	M	2	100% / 100%
2	P	2	100%
2	Q	2	100% / 50% / 50%
2	R	2	50% / 50% / 50%
2	S	2	50% / 50% / 50%
2	T	2	50% / 50% / 50%
2	U	2	50% / 50% / 50%
2	V	2	100% / 100%
2	W	2	100% / 100%
2	Y	2	100% / 50% / 50%
2	b	2	100% / 100%
2	c	2	50% / 50% / 50%
2	d	2	50% / 50% / 50%
3	F	3	67% / 67% / 33%
3	J	3	67% / 33% / 67%
3	Z	3	67% / 67% / 33%
4	G	4	50% / 25% / 75%
4	N	4	25% / 25% / 75%
4	O	4	50% / 75% / 25%
4	X	4	75% / 50% / 50%
4	a	4	50% / 25% / 75%

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 26012 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 Spike glycoprotein,Fibritin.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	B	1068	8342	5331	1379	1586	46	0	0
1	C	1069	8346	5333	1380	1587	46	0	0
1	A	1066	8327	5320	1377	1584	46	0	0

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	577	ALA	SER	conflict	UNP P59594
B	968	PRO	LYS	engineered mutation	UNP P59594
B	969	PRO	VAL	engineered mutation	UNP P59594
B	1191	GLY	-	linker	UNP P59594
B	1192	SER	ALA	linker	UNP D9IEJ2
B	1220	GLY	-	expression tag	UNP D9IEJ2
B	1221	ARG	-	expression tag	UNP D9IEJ2
B	1222	SER	-	expression tag	UNP D9IEJ2
B	1223	LEU	-	expression tag	UNP D9IEJ2
B	1224	GLU	-	expression tag	UNP D9IEJ2
B	1225	VAL	-	expression tag	UNP D9IEJ2
B	1226	LEU	-	expression tag	UNP D9IEJ2
B	1227	PHE	-	expression tag	UNP D9IEJ2
B	1228	GLN	-	expression tag	UNP D9IEJ2
C	577	ALA	SER	conflict	UNP P59594
C	968	PRO	LYS	engineered mutation	UNP P59594
C	969	PRO	VAL	engineered mutation	UNP P59594
C	1191	GLY	-	linker	UNP P59594
C	1192	SER	ALA	linker	UNP D9IEJ2
C	1220	GLY	-	expression tag	UNP D9IEJ2
C	1221	ARG	-	expression tag	UNP D9IEJ2
C	1222	SER	-	expression tag	UNP D9IEJ2
C	1223	LEU	-	expression tag	UNP D9IEJ2
C	1224	GLU	-	expression tag	UNP D9IEJ2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	1225	VAL	-	expression tag	UNP D9IEJ2
C	1226	LEU	-	expression tag	UNP D9IEJ2
C	1227	PHE	-	expression tag	UNP D9IEJ2
C	1228	GLN	-	expression tag	UNP D9IEJ2
A	577	ALA	SER	conflict	UNP P59594
A	968	PRO	LYS	engineered mutation	UNP P59594
A	969	PRO	VAL	engineered mutation	UNP P59594
A	1191	GLY	-	linker	UNP P59594
A	1192	SER	ALA	linker	UNP D9IEJ2
A	1220	GLY	-	expression tag	UNP D9IEJ2
A	1221	ARG	-	expression tag	UNP D9IEJ2
A	1222	SER	-	expression tag	UNP D9IEJ2
A	1223	LEU	-	expression tag	UNP D9IEJ2
A	1224	GLU	-	expression tag	UNP D9IEJ2
A	1225	VAL	-	expression tag	UNP D9IEJ2
A	1226	LEU	-	expression tag	UNP D9IEJ2
A	1227	PHE	-	expression tag	UNP D9IEJ2
A	1228	GLN	-	expression tag	UNP D9IEJ2

- Molecule 2 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
			Total	C	N	O		
2	D	2	28	16	2	10	0	0
2	E	2	28	16	2	10	0	0
2	H	2	28	16	2	10	0	0
2	I	2	28	16	2	10	0	0
2	K	2	28	16	2	10	0	0
2	L	2	28	16	2	10	0	0
2	M	2	28	16	2	10	0	0

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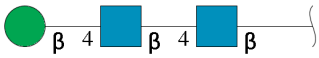
Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
2	P	2	28	16	2	10	0	0
2	Q	2	28	16	2	10	0	0
2	R	2	28	16	2	10	0	0
2	S	2	28	16	2	10	0	0
2	T	2	28	16	2	10	0	0
2	U	2	28	16	2	10	0	0
2	V	2	28	16	2	10	0	0
2	W	2	28	16	2	10	0	0
2	Y	2	28	16	2	10	0	0
2	b	2	28	16	2	10	0	0
2	c	2	28	16	2	10	0	0
2	d	2	28	16	2	10	0	0

- Molecule 3 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.

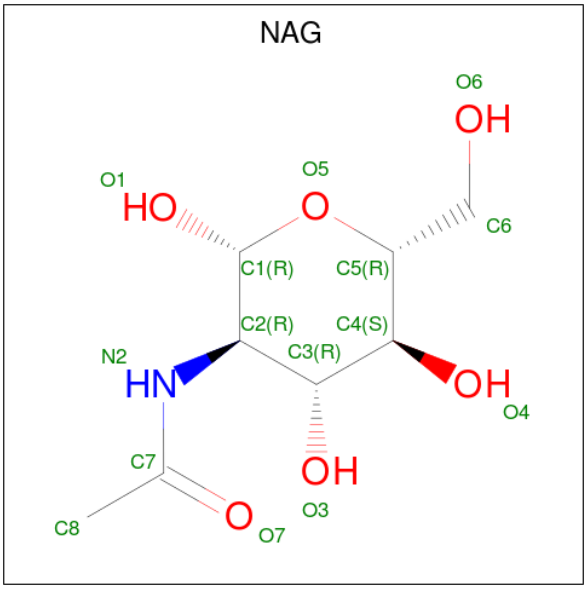


Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
3	F	3	39	22	2	15	0	0
3	J	3	39	22	2	15	0	0
3	Z	3	39	22	2	15	0	0

- Molecule 4 is an oligosaccharide called 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
4	G	4	Total	C	N	O	0	0
			50	28	2	20		
4	N	4	Total	C	N	O	0	0
			50	28	2	20		
4	O	4	Total	C	N	O	0	0
			50	28	2	20		
4	X	4	Total	C	N	O	0	0
			50	28	2	20		
4	a	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				AltConf
5	B	1	Total	C	N	O	0
			56	32	4	20	
5	B	1	Total	C	N	O	0
			56	32	4	20	
5	B	1	Total	C	N	O	0
			56	32	4	20	
5	B	1	Total	C	N	O	0
			56	32	4	20	

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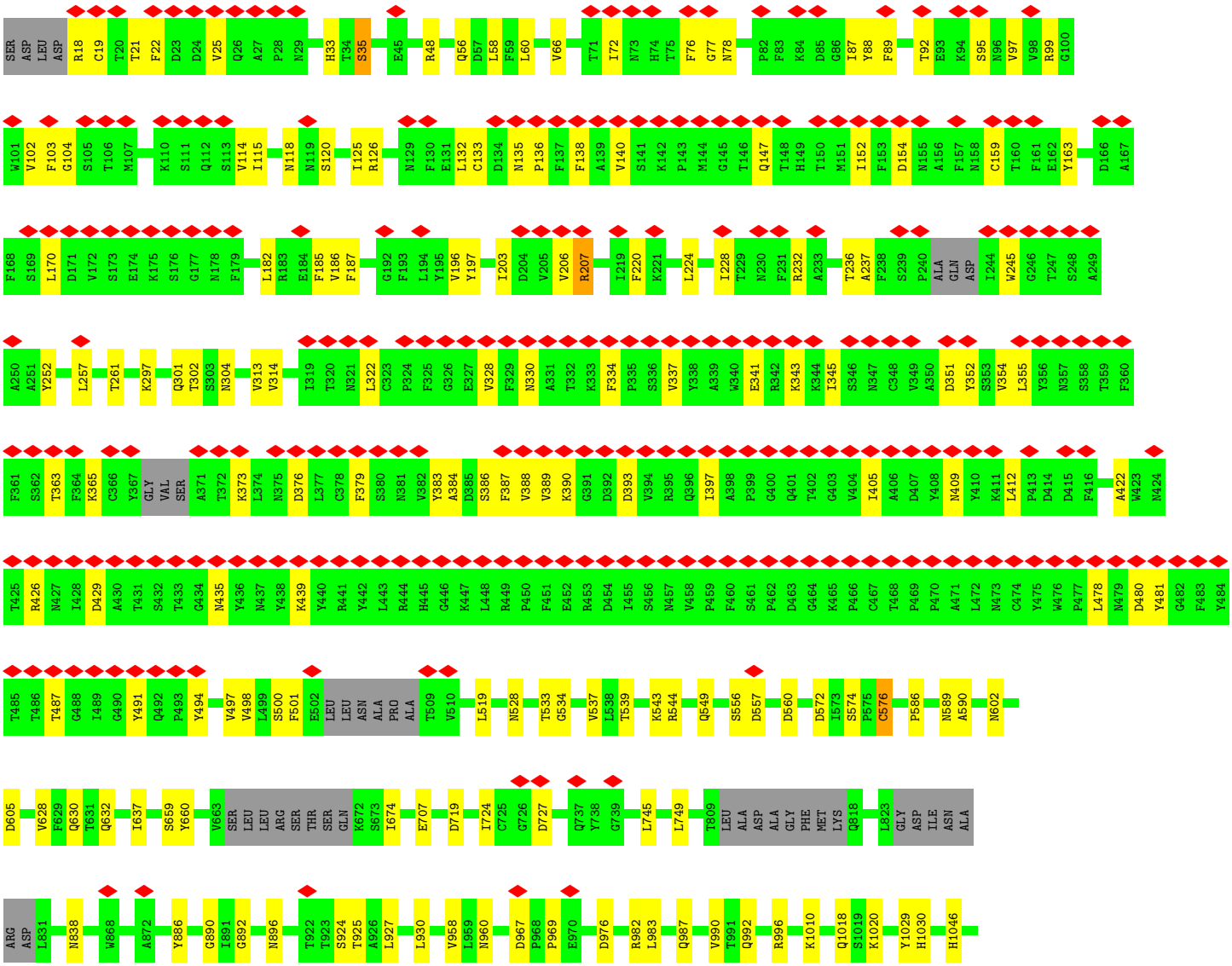
Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
5	C	1	28	16	2	10	0
			Total	C	N	O	
5	C	1	28	16	2	10	0
			Total	C	N	O	
5	A	1	14	8	1	5	0

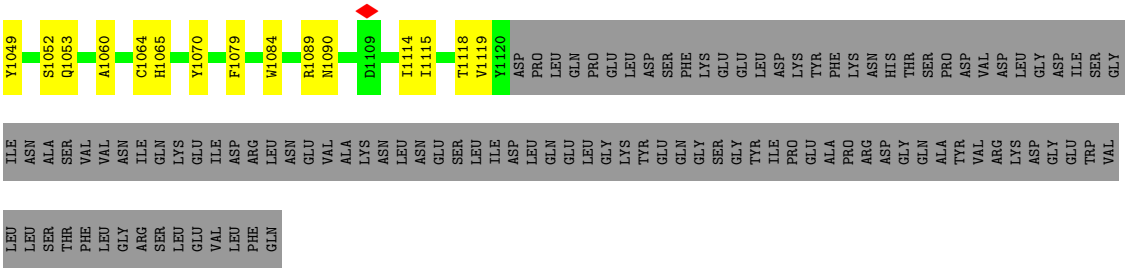
3 Residue-property plots (i)

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: Spike glycoprotein,Fibritin

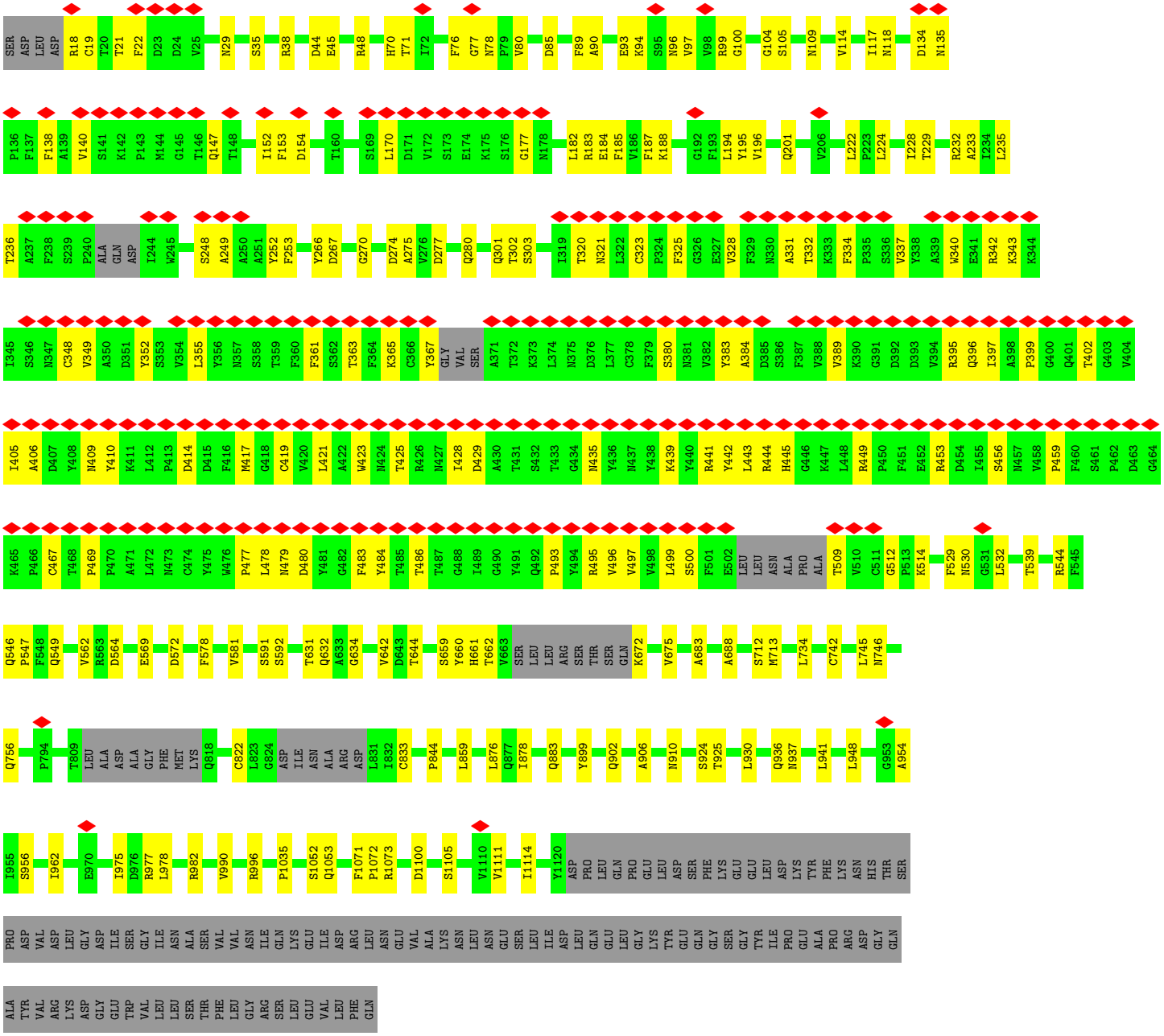
Chain B: 22% 73% 15% 12%





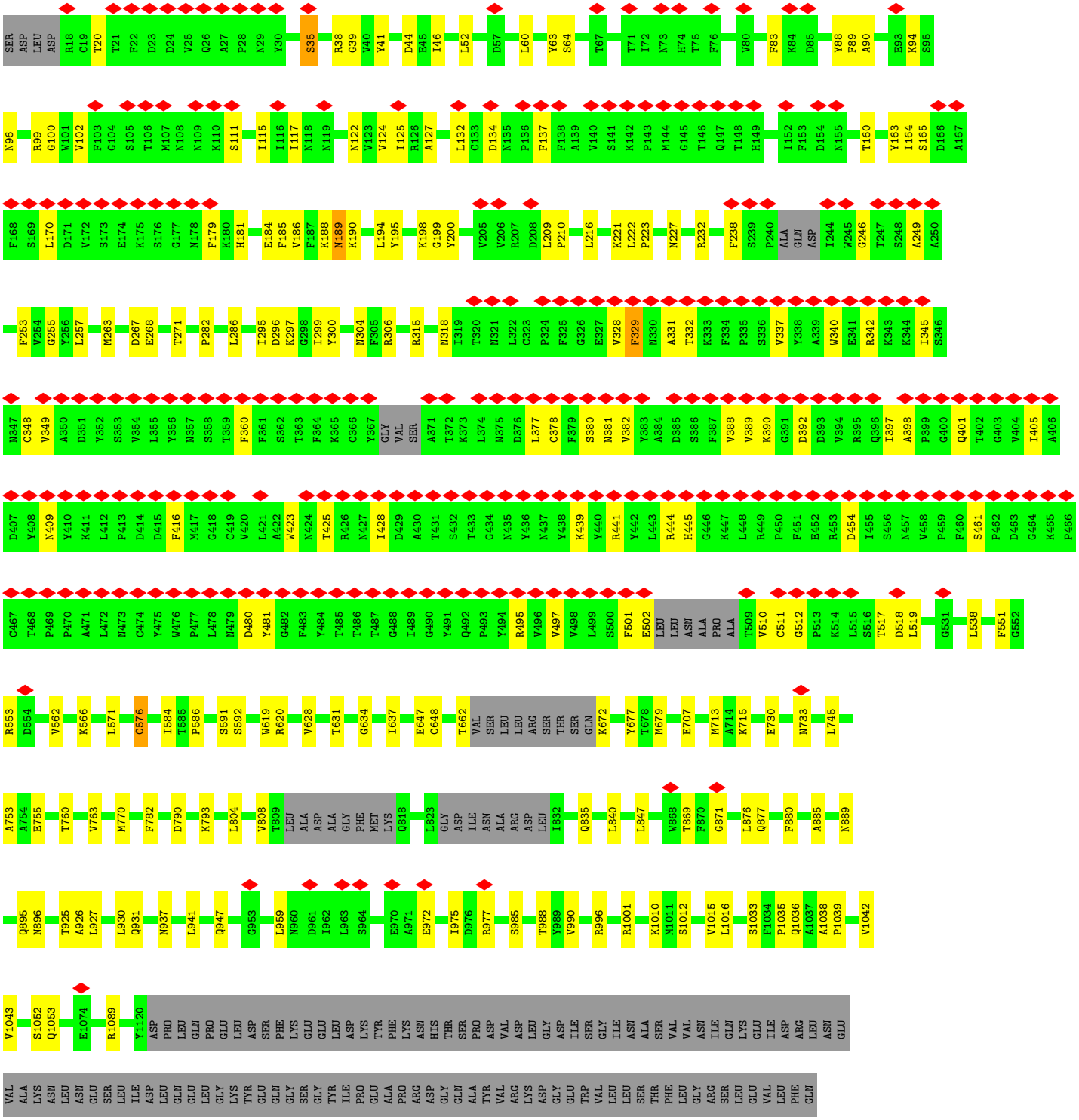
- Molecule 1: Spike glycoprotein,Fibritin

Chain C: 19% 70% 18% 12%

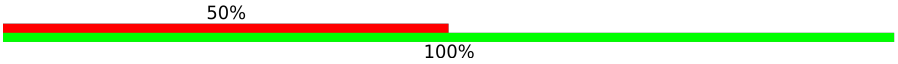


• Molecule 1: Spike glycoprotein,Fibritin

Chain A: 22% 71% 17% 12%



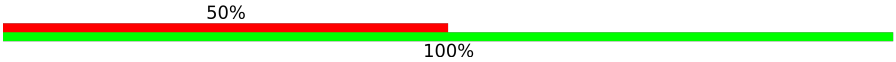
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D: 

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

Chain E:

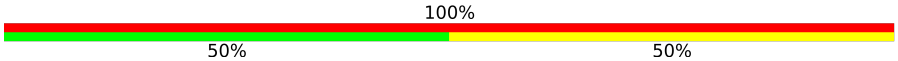
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H: 

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

Chain I:

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

Chain K: 

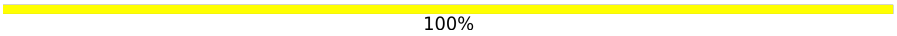
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:

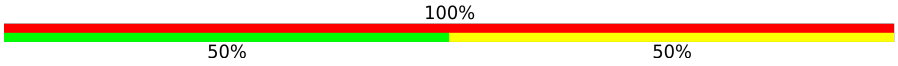
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:

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

Chain P: 

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

Chain Q: 

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

Chain R:

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

Chain S:

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

Chain T:

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

Chain U:

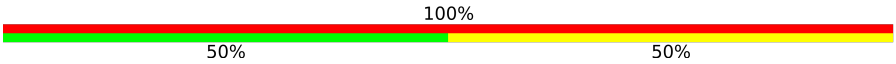
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain V:

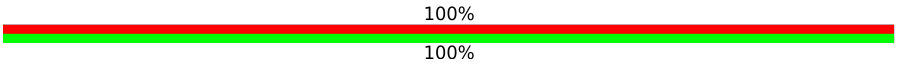
• Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain W:

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

Chain Y: 

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

Chain b: 

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

Chain c:

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

Chain d:

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

Chain F:

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

Chain J:

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

Chain Z:

• Molecule 4: 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

Chain G:

• Molecule 4: 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

Chain N:

• Molecule 4: 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

Chain O:

• Molecule 4: 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

Chain X:

• Molecule 4: 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

Chain a:

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	38603	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 ($e^-/Å^2$)	65	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	29000	Depositor
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.158	Depositor
Minimum map value	-0.075	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.0328	Depositor
Map size (Å)	329.59998, 329.59998, 329.59998	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.03, 1.03, 1.03	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, NAG, BMA

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	0.49	0/8526	0.63	2/11601 (0.0%)
1	B	0.50	0/8541	0.63	1/11622 (0.0%)
1	C	0.50	0/8545	0.63	0/11627
All	All	0.50	0/25612	0.63	3/34850 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	3
1	B	0	4
1	C	0	2
All	All	0	9

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	329	PHE	CB-CG-CD1	6.25	125.17	120.80
1	A	329	PHE	CB-CG-CD2	-5.12	117.21	120.80
1	B	1070	TYR	C-N-CA	5.00	134.20	121.70

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	35	SER	Peptide

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Mol	Chain	Res	Type	Group
1	A	576	CYS	Peptide
1	A	83	PHE	Peptide
1	B	207	ARG	Peptide
1	B	35	SER	Peptide
1	B	576	CYS	Peptide
1	B	72	ILE	Peptide
1	C	35	SER	Peptide
1	C	85	ASP	Peptide

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	8327	0	8068	133	0
1	B	8342	0	8088	113	0
1	C	8346	0	8090	127	0
2	D	28	0	25	0	0
2	E	28	0	25	0	0
2	H	28	0	25	0	0
2	I	28	0	25	0	0
2	K	28	0	25	0	0
2	L	28	0	25	0	0
2	M	28	0	25	0	0
2	P	28	0	25	0	0
2	Q	28	0	25	0	0
2	R	28	0	25	0	0
2	S	28	0	25	0	0
2	T	28	0	25	1	0
2	U	28	0	25	2	0
2	V	28	0	25	0	0
2	W	28	0	25	0	0
2	Y	28	0	25	0	0
2	b	28	0	25	0	0
2	c	28	0	25	0	0
2	d	28	0	25	0	0
3	F	39	0	34	0	0
3	J	39	0	34	0	0
3	Z	39	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	G	50	0	43	0	0
4	N	50	0	43	0	0
4	O	50	0	43	0	0
4	X	50	0	43	0	0
4	a	50	0	43	0	0
5	A	14	0	13	0	0
5	B	56	0	52	0	0
5	C	28	0	26	0	0
All	All	26012	0	25129	354	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.

All (354) 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:B:125:ILE:HB	1:B:163:TYR:HB3	1.70	0.73
1:B:58:LEU:HB3	1:B:257:LEU:HB3	1.72	0.72
1:B:1089:ARG:HH21	1:C:876:LEU:HB2	1.55	0.71
1:A:871:GLY:HA3	1:A:1016:LEU:HB3	1.74	0.68
1:B:334:PHE:HB2	1:B:388:VAL:HG23	1.75	0.67
1:B:586:PRO:HD3	1:B:674:ILE:HD11	1.78	0.65
1:C:29:ASN:HD22	1:C:70:HIS:HB2	1.61	0.65
1:B:1029:TYR:HB2	1:B:1049:TYR:HB3	1.79	0.64
1:A:89:PHE:HB3	1:A:185:PHE:HB2	1.80	0.64
1:B:549:GLN:HA	1:C:45:GLU:HB2	1.79	0.64
1:C:712:SER:OG	1:C:756:GLN:NE2	2.27	0.64
1:C:661:HIS:ND1	1:C:675:VAL:HG21	2.12	0.64
1:C:22:PHE:H	1:C:135:ASN:HB2	1.63	0.63
1:C:182:LEU:HD23	1:C:201:GLN:HE21	1.63	0.62
1:C:96:ASN:O	1:C:99:ARG:NH1	2.32	0.62
1:C:529:PHE:HB2	1:C:532:LEU:HB2	1.81	0.62
1:B:383:TYR:HB2	1:B:500:SER:HB3	1.82	0.61
1:C:688:ALA:O	1:A:877:GLN:NE2	2.33	0.61
1:A:713:MET:SD	1:A:937:ASN:ND2	2.74	0.60
1:C:713:MET:O	1:C:756:GLN:NE2	2.34	0.60
1:B:890:GLY:O	1:B:1020:LYS:NZ	2.34	0.60
1:B:138:PHE:HB2	1:B:236:THR:HA	1.83	0.60
1:C:1105:SER:OG	1:A:896:ASN:ND2	2.34	0.60
1:B:439:LYS:NZ	1:B:480:ASP:OD1	2.32	0.60
1:A:745:LEU:HD22	1:A:990:VAL:HG21	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:348:CYS:HB3	1:A:510:VAL:HG12	1.85	0.59
1:C:459:PRO:HG2	1:C:469:PRO:HG3	1.84	0.59
1:A:390:LYS:HE3	1:A:392:ASP:HB2	1.84	0.59
1:B:21:THR:HB	1:B:135:ASN:HD22	1.68	0.58
1:B:102:VAL:HB	1:B:232:ARG:HB3	1.84	0.58
1:B:1064:CYS:HB2	1:B:1114:ILE:HD12	1.84	0.58
1:C:99:ARG:HE	1:C:170:LEU:HD23	1.69	0.58
1:C:18:ARG:N	1:C:153:PHE:O	2.36	0.58
1:C:342:ARG:NH2	1:A:223:PRO:O	2.37	0.58
1:C:449:ARG:NH2	1:A:189:ASN:O	2.34	0.57
1:C:662:THR:HA	1:C:672:LYS:HG2	1.85	0.57
1:A:405:ILE:HA	1:A:409:ASN:HB2	1.86	0.57
1:B:328:VAL:HG22	1:B:343:LYS:HD3	1.85	0.57
1:C:114:VAL:HG21	1:C:224:LEU:HD13	1.86	0.57
1:B:136:PRO:HB3	1:B:152:ILE:HA	1.87	0.57
1:A:755:GLU:OE2	1:A:1001:ARG:NH1	2.38	0.57
1:B:322:LEU:HD22	1:B:351:ASP:HB3	1.86	0.56
1:C:564:ASP:HB3	1:C:569:GLU:H	1.70	0.56
1:C:591:SER:OG	1:C:592:SER:N	2.37	0.56
1:B:389:VAL:HG11	1:B:397:ILE:HD11	1.87	0.56
1:B:390:LYS:HG3	1:B:481:TYR:HD1	1.68	0.56
1:A:445:HIS:NE2	1:A:461:SER:O	2.33	0.56
1:B:103:PHE:HB2	1:B:114:VAL:HB	1.87	0.56
1:A:194:LEU:HB3	1:A:222:LEU:HB2	1.87	0.56
1:A:342:ARG:NH2	1:A:502:GLU:OE1	2.39	0.56
1:B:89:PHE:HB3	1:B:185:PHE:HB2	1.87	0.55
1:C:936:GLN:O	1:C:996:ARG:NH1	2.39	0.55
1:B:95:SER:O	1:B:99:ARG:NH2	2.39	0.55
1:B:363:THR:HB	1:B:365:LYS:HG3	1.88	0.55
1:B:186:VAL:HB	1:B:197:TYR:HB2	1.89	0.55
1:B:719:ASP:OD1	1:A:304:ASN:ND2	2.40	0.55
1:C:367:TYR:HB2	1:C:417:MET:HB3	1.88	0.55
1:B:1018:GLN:NE2	1:B:1030:HIS:O	2.39	0.55
1:B:126:ARG:NH2	1:B:159:CYS:SG	2.80	0.55
1:A:38:ARG:NH1	1:A:184:GLU:OE2	2.40	0.55
1:B:314:VAL:HG22	1:B:528:ASN:HB3	1.89	0.54
1:B:429:ASP:O	1:B:435:ASN:ND2	2.40	0.54
1:C:277:ASP:HB3	1:C:280:GLN:HB2	1.89	0.54
1:C:1072:PRO:O	1:A:895:GLN:NE2	2.41	0.54
1:C:734:LEU:HD12	1:C:975:ILE:HD11	1.89	0.54
1:C:38:ARG:NH1	1:C:184:GLU:OE2	2.40	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:LYS:NZ	1:A:249:ALA:O	2.40	0.54
1:B:838:ASN:OD1	1:B:838:ASN:N	2.39	0.54
1:C:456:SER:HB2	1:A:111:SER:HB3	1.90	0.54
1:C:1071:PHE:HZ	1:C:1111:VAL:HG21	1.71	0.54
1:C:1073:ARG:NH1	1:C:1100:ASP:O	2.41	0.54
1:A:100:GLY:HA3	1:A:117:ILE:HG22	1.90	0.54
1:B:745:LEU:HD22	1:B:990:VAL:HG21	1.90	0.54
1:C:383:TYR:HB2	1:C:500:SER:HB3	1.88	0.54
1:A:41:TYR:HB3	1:A:216:LEU:HB2	1.89	0.54
1:A:416:PHE:HB3	1:A:501:PHE:HB2	1.88	0.54
1:B:724:ILE:O	1:B:982:ARG:NH1	2.36	0.54
1:A:52:LEU:HD11	1:A:263:MET:HB3	1.89	0.54
1:C:509:THR:OG1	1:C:514:LYS:NZ	2.37	0.54
1:C:421:LEU:HB3	1:C:497:VAL:HB	1.89	0.53
1:B:22:PHE:HE2	1:B:78:ASN:HD21	1.56	0.53
1:C:80:VAL:HG22	1:C:232:ARG:HH11	1.73	0.53
1:C:425:THR:HB	1:C:428:ILE:HB	1.90	0.53
1:C:578:PHE:H	1:A:835:GLN:HE22	1.56	0.53
1:A:707:GLU:OE2	1:A:1010:LYS:NZ	2.42	0.53
1:A:337:VAL:HG13	1:A:340:TRP:HE1	1.73	0.53
1:B:393:ASP:HB3	1:B:405:ILE:HD11	1.90	0.53
1:A:195:TYR:HE1	1:A:221:LYS:HG3	1.74	0.53
1:C:352:TYR:HB3	1:C:355:LEU:HB3	1.90	0.53
1:A:125:ILE:HB	1:A:163:TYR:HB3	1.90	0.53
1:A:972:GLU:HA	1:A:975:ILE:HG22	1.91	0.53
1:B:35:SER:HB2	1:B:60:LEU:HD21	1.91	0.53
1:C:94:LYS:NZ	1:C:177:GLY:O	2.40	0.53
1:B:412:LEU:HD21	1:B:498:VAL:HG11	1.91	0.53
1:C:93:GLU:HG3	1:C:183:ARG:HH21	1.74	0.53
1:A:885:ALA:HB1	1:A:895:GLN:HB2	1.91	0.52
1:C:444:ARG:NH1	1:C:445:HIS:O	2.41	0.52
1:A:331:ALA:O	1:A:495:ARG:NH1	2.41	0.52
1:B:33:HIS:HB2	1:B:66:VAL:HB	1.91	0.52
1:A:35:SER:OG	1:A:64:SER:N	2.39	0.52
1:B:328:VAL:HG21	1:B:345:ILE:HD11	1.91	0.52
1:C:99:ARG:HB2	1:C:138:PHE:HE2	1.74	0.52
1:A:662:THR:HA	1:A:672:LYS:HG2	1.92	0.52
1:C:104:GLY:H	1:C:228:ILE:HG23	1.75	0.52
1:A:39:GLY:HA3	1:A:60:LEU:HB3	1.91	0.52
1:C:444:ARG:HH21	1:A:227:ASN:HD21	1.58	0.52
1:A:591:SER:OG	1:A:592:SER:N	2.43	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:318:ASN:OD1	1:A:566:LYS:NZ	2.43	0.51
1:A:628:VAL:HG22	1:A:637:ILE:HG12	1.91	0.51
1:C:742:CYS:O	1:C:746:ASN:ND2	2.44	0.51
1:C:248:SER:OG	1:C:249:ALA:N	2.41	0.51
1:C:429:ASP:O	1:C:435:ASN:ND2	2.44	0.51
1:A:122:ASN:ND2	1:A:165:SER:O	2.43	0.51
1:A:127:ALA:O	1:A:160:THR:N	2.39	0.51
1:C:549:GLN:NE2	1:A:46:ILE:O	2.44	0.51
1:A:185:PHE:HE1	1:A:198:LYS:HG3	1.76	0.51
1:C:331:ALA:HB3	1:C:334:PHE:HE1	1.76	0.51
1:C:956:SER:H	1:C:962:ILE:HD11	1.76	0.51
1:B:337:VAL:HG22	1:B:387:PHE:HB2	1.93	0.50
1:C:399:PRO:HB2	1:C:414:ASP:HA	1.92	0.50
1:A:44:ASP:OD1	1:A:44:ASP:N	2.44	0.50
1:A:425:THR:HB	1:A:428:ILE:HB	1.93	0.50
1:A:840:LEU:HD21	1:A:941:LEU:HD22	1.93	0.50
1:B:892:GLY:O	1:B:1090:ASN:ND2	2.39	0.50
1:A:268:GLU:HB2	2:U:1:NAG:H82	1.93	0.50
1:B:97:VAL:O	1:B:236:THR:N	2.44	0.50
1:B:104:GLY:H	1:B:228:ILE:HG23	1.77	0.50
1:B:405:ILE:HA	1:B:409:ASN:HB2	1.92	0.50
1:B:537:VAL:HB	1:B:574:SER:HB2	1.93	0.50
1:C:443:LEU:HB3	1:C:477:PRO:HA	1.92	0.50
1:A:297:LYS:HG3	1:A:586:PRO:HA	1.92	0.50
1:C:509:THR:HG21	1:C:530:ASN:HD21	1.75	0.50
1:A:125:ILE:O	1:A:163:TYR:N	2.38	0.50
1:C:419:CYS:H	1:C:499:LEU:HB2	1.77	0.50
1:A:538:LEU:HD22	1:A:571:LEU:HD13	1.94	0.50
1:B:206:VAL:HG23	1:B:207:ARG:HG3	1.93	0.50
1:B:373:LYS:HA	1:B:376:ASP:HB2	1.94	0.50
1:C:337:VAL:HG11	1:C:405:ILE:HD12	1.94	0.50
1:C:859:LEU:HD23	1:C:1035:PRO:HG3	1.94	0.50
1:B:386:SER:HA	1:B:497:VAL:HG22	1.94	0.49
1:C:713:MET:HB2	1:C:937:ASN:HD21	1.77	0.49
1:A:63:TYR:O	1:A:620:ARG:NH2	2.45	0.49
1:A:808:VAL:HA	1:A:931:GLN:HE21	1.77	0.49
1:A:382:VAL:HG22	1:A:501:PHE:HA	1.94	0.49
1:B:707:GLU:OE2	1:B:1010:LYS:NZ	2.35	0.49
1:B:99:ARG:HB3	1:B:118:ASN:HB3	1.94	0.49
1:B:967:ASP:HB3	1:B:969:PRO:HD2	1.93	0.49
1:A:730:GLU:HA	1:A:733:ASN:HD22	1.77	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:102:VAL:HG11	1:B:132:LEU:HD11	1.95	0.49
1:C:948:LEU:O	1:C:982:ARG:NH2	2.46	0.49
1:C:100:GLY:HA3	1:C:117:ILE:HG22	1.96	0.48
1:C:109:ASN:OD1	1:C:109:ASN:N	2.45	0.48
1:C:683:ALA:O	1:A:770:MET:N	2.46	0.48
1:B:77:GLY:O	1:B:252:TYR:OH	2.21	0.48
1:B:543:LYS:NZ	1:B:560:ASP:OD2	2.46	0.48
1:C:409:ASN:OD1	1:C:441:ARG:N	2.45	0.48
1:B:628:VAL:HG22	1:B:637:ILE:HG12	1.95	0.48
1:A:90:ALA:HB3	1:A:253:PHE:HB2	1.96	0.48
1:A:380:SER:OG	1:A:381:ASN:N	2.46	0.48
1:B:992:GLN:HB3	1:B:996:ARG:HH12	1.79	0.48
1:B:56:GLN:HB2	1:B:261:THR:HG22	1.94	0.48
1:A:209:LEU:HD12	1:A:210:PRO:HD2	1.95	0.48
1:B:976:ASP:OD2	1:A:977:ARG:NH2	2.47	0.48
1:C:140:VAL:HG22	1:C:147:GLN:HA	1.94	0.48
1:A:286:LEU:HD21	1:A:300:TYR:HD2	1.79	0.48
1:A:377:LEU:O	1:A:512:GLY:N	2.47	0.48
1:C:78:ASN:ND2	1:C:233:ALA:O	2.46	0.47
1:C:445:HIS:CE1	1:C:467:CYS:H	2.32	0.47
1:C:899:TYR:HA	1:C:902:GLN:HE21	1.79	0.47
1:C:105:SER:HA	1:C:229:THR:HG22	1.95	0.47
1:A:190:LYS:HD3	1:A:195:TYR:HE2	1.80	0.47
1:A:378:CYS:HA	1:A:511:CYS:HA	1.96	0.47
1:A:425:THR:HG21	1:A:495:ARG:HD2	1.95	0.47
1:A:647:GLU:O	1:A:677:TYR:OH	2.30	0.47
1:C:152:ILE:HB	1:C:153:PHE:HD1	1.79	0.47
1:C:19:CYS:N	1:C:154:ASP:OD1	2.47	0.47
1:A:360:PHE:HB3	1:A:423:TRP:HB2	1.95	0.47
1:A:102:VAL:HG22	1:A:115:ILE:HG12	1.96	0.47
1:B:341:GLU:HB3	1:B:386:SER:HB3	1.96	0.47
1:B:958:VAL:HG12	1:B:960:ASN:H	1.80	0.47
1:C:97:VAL:HA	1:C:236:THR:HB	1.96	0.47
1:C:188:LYS:HB2	1:C:195:TYR:HB2	1.96	0.47
1:C:194:LEU:HB3	1:C:222:LEU:HB2	1.97	0.47
1:C:320:THR:OG1	1:C:321:ASN:N	2.48	0.47
1:C:363:THR:HB	1:C:365:LYS:HG3	1.96	0.47
1:C:395:ARG:HH12	1:C:396:GLN:HE21	1.62	0.47
1:B:140:VAL:HG22	1:B:147:GLN:HA	1.97	0.47
1:C:659:SER:OG	1:C:660:TYR:N	2.48	0.47
1:C:878:ILE:HG23	1:C:883:GLN:HE21	1.80	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:ARG:HH21	1:A:553:ARG:HD2	1.79	0.46
1:B:927:LEU:HD23	1:B:930:LEU:HD12	1.97	0.46
1:B:182:LEU:HB2	1:B:203:ILE:HD13	1.97	0.46
1:B:196:VAL:HB	1:B:220:PHE:HB2	1.97	0.46
1:B:544:ARG:NH1	2:T:1:NAG:O5	2.48	0.46
1:B:727:ASP:OD1	1:A:306:ARG:NH1	2.49	0.46
1:A:88:TYR:N	1:A:255:GLY:O	2.47	0.46
1:B:18:ARG:N	1:B:154:ASP:OD1	2.48	0.46
1:B:707:GLU:OE1	1:B:1046:HIS:NE2	2.46	0.46
1:B:120:SER:HA	1:B:170:LEU:HD23	1.97	0.46
1:A:124:VAL:HG22	1:A:164:ILE:HG12	1.96	0.46
1:A:181:HIS:CE1	1:A:200:TYR:HH	2.33	0.46
1:A:295:ILE:HG23	1:A:300:TYR:HE2	1.81	0.46
1:B:328:VAL:HG11	1:B:384:ALA:HB1	1.98	0.46
1:B:602:ASN:OD1	1:B:630:GLN:NE2	2.48	0.46
1:A:188:LYS:HB2	1:A:195:TYR:HB2	1.98	0.46
1:A:390:LYS:HB2	1:A:481:TYR:HD1	1.81	0.46
1:B:992:GLN:O	1:B:996:ARG:NH1	2.49	0.46
1:C:76:PHE:CZ	1:C:78:ASN:HB3	2.50	0.46
1:C:484:TYR:HB3	1:C:486:THR:H	1.80	0.45
1:B:313:VAL:HG11	1:B:519:LEU:HD12	1.99	0.45
1:A:1038:ALA:HB2	1:A:1043:VAL:HG23	1.97	0.45
1:B:1060:ALA:HB2	1:B:1115:ILE:HD12	1.97	0.45
1:A:238:PHE:N	1:A:246:GLY:O	2.49	0.45
1:A:389:VAL:HG11	1:A:397:ILE:HD11	1.98	0.45
1:C:361:PHE:CD1	1:C:423:TRP:HB3	2.52	0.45
1:C:439:LYS:HB3	1:C:478:LEU:HB3	1.99	0.45
1:B:115:ILE:HD11	1:B:132:LEU:HD21	1.99	0.45
1:B:487:THR:HG21	1:B:491:TYR:HB2	1.99	0.45
1:B:605:ASP:N	1:B:605:ASP:OD1	2.48	0.45
1:A:1012:SER:HA	1:A:1016:LEU:HD12	1.98	0.45
1:B:533:THR:OG1	1:B:534:GLY:N	2.50	0.45
1:C:89:PHE:HB3	1:C:185:PHE:HB2	1.98	0.45
1:C:302:THR:OG1	1:C:303:SER:N	2.50	0.45
1:C:405:ILE:HD13	1:C:409:ASN:HD22	1.81	0.45
1:B:659:SER:OG	1:B:660:TYR:N	2.50	0.45
1:C:546:GLN:HA	1:C:547:PRO:HD3	1.85	0.45
1:C:924:SER:OG	1:C:925:THR:N	2.49	0.45
1:A:441:ARG:NH2	1:A:454:ASP:O	2.39	0.45
1:A:869:THR:HG21	1:A:876:LEU:HD12	1.99	0.45
1:A:551:PHE:HB3	1:A:562:VAL:HG23	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:VAL:HA	1:B:76:PHE:HD2	1.82	0.44
1:B:337:VAL:HG21	1:B:389:VAL:HG12	1.99	0.44
1:B:422:ALA:HB1	1:B:494:TYR:HD2	1.82	0.44
1:C:325:PHE:HB3	1:C:328:VAL:HB	1.99	0.44
1:C:906:ALA:O	1:C:910:ASN:ND2	2.50	0.44
1:A:715:LYS:HD2	1:A:753:ALA:HB1	2.00	0.44
1:A:1015:VAL:HG11	1:A:1035:PRO:HD3	1.99	0.44
1:B:87:ILE:HD11	1:B:187:PHE:HD2	1.83	0.44
1:B:886:TYR:HB2	1:A:1089:ARG:NH1	2.32	0.44
1:A:315:ARG:HH21	1:A:566:LYS:HB3	1.81	0.44
1:A:332:THR:HA	1:A:495:ARG:HH12	1.83	0.44
1:C:77:GLY:O	1:C:252:TYR:OH	2.24	0.44
1:C:301:GLN:NE2	1:C:581:VAL:O	2.51	0.44
1:A:102:VAL:HG11	1:A:132:LEU:HD11	2.00	0.44
1:A:348:CYS:SG	1:A:349:VAL:N	2.90	0.44
1:A:631:THR:HG23	1:A:634:GLY:H	1.82	0.44
1:A:648:CYS:HB2	1:A:679:MET:HG3	2.00	0.44
1:B:330:ASN:HB2	1:B:354:VAL:HG21	2.00	0.44
1:C:954:ALA:HB2	1:C:977:ARG:HD3	1.99	0.44
1:A:388:VAL:HG22	1:A:495:ARG:HG2	2.00	0.44
1:C:402:THR:HA	1:C:406:ALA:HB3	1.99	0.44
1:A:328:VAL:HG21	1:A:345:ILE:HD11	1.98	0.44
1:B:114:VAL:HG21	1:B:224:LEU:HD13	2.00	0.43
1:C:99:ARG:HB3	1:C:118:ASN:HB3	2.00	0.43
1:A:267:ASP:OD1	1:A:271:THR:N	2.34	0.43
1:B:439:LYS:HB3	1:B:478:LEU:HB3	1.98	0.43
1:B:632:GLN:HB3	1:C:844:PRO:HG3	1.99	0.43
1:A:808:VAL:HG21	1:A:1039:PRO:HG3	1.99	0.43
1:C:745:LEU:HD22	1:C:990:VAL:HG21	2.01	0.43
1:C:439:LYS:NZ	1:C:480:ASP:OD1	2.46	0.43
1:A:96:ASN:O	1:A:99:ARG:NE	2.52	0.43
1:C:187:PHE:HE1	1:C:196:VAL:HG13	1.82	0.43
1:C:332:THR:O	1:C:495:ARG:NH1	2.49	0.43
1:C:631:THR:OG1	1:C:632:GLN:N	2.51	0.43
1:C:337:VAL:HG13	1:C:340:TRP:HE1	1.82	0.43
1:B:78:ASN:O	1:B:232:ARG:NE	2.51	0.43
1:B:301:GLN:NE2	1:B:302:THR:O	2.52	0.43
1:B:337:VAL:HG12	1:B:409:ASN:HB3	1.99	0.43
1:B:1065:HIS:ND1	1:B:1118:THR:HA	2.33	0.43
1:C:529:PHE:CE2	1:C:562:VAL:HG21	2.53	0.43
1:C:1052:SER:OG	1:C:1053:GLN:N	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:20:THR:HB	1:A:137:PHE:HE1	1.83	0.43
1:A:299:ILE:HG12	1:A:584:ILE:HG12	2.01	0.43
1:B:1079:PHE:HB2	1:B:1084:TRP:NE1	2.34	0.43
1:C:642:VAL:HG12	1:C:644:THR:HG22	2.00	0.43
1:B:1079:PHE:HB2	1:B:1084:TRP:CD1	2.54	0.42
1:C:343:LYS:O	1:C:384:ALA:N	2.45	0.42
1:A:398:ALA:HB3	1:A:401:GLN:HG3	2.01	0.42
1:C:274:ASP:OD1	1:C:275:ALA:N	2.52	0.42
1:C:539:THR:O	1:C:572:ASP:N	2.48	0.42
1:A:1036:GLN:HB2	1:A:1043:VAL:HB	2.01	0.42
1:B:924:SER:OG	1:B:925:THR:N	2.52	0.42
1:B:99:ARG:HE	1:B:170:LEU:HD21	1.85	0.42
1:C:21:THR:HG21	1:C:134:ASP:HB2	2.02	0.42
1:C:267:ASP:OD1	1:C:270:GLY:N	2.53	0.42
1:A:337:VAL:HG11	1:A:405:ILE:HG12	2.00	0.42
1:C:76:PHE:HE1	1:C:235:LEU:HD12	1.84	0.42
1:C:483:PHE:HB3	1:C:493:PRO:HD3	2.02	0.42
1:C:544:ARG:NH1	2:U:1:NAG:H5	2.35	0.42
1:A:186:VAL:HG13	1:A:257:LEU:HD11	2.00	0.42
1:A:184:GLU:O	1:A:199:GLY:N	2.50	0.42
1:A:329:PHE:HA	1:A:497:VAL:HG21	2.01	0.42
1:A:439:LYS:NZ	1:A:480:ASP:OD1	2.48	0.42
1:B:237:ALA:HB1	1:B:245:TRP:HB3	2.01	0.42
1:C:442:TYR:CD1	1:C:479:ASN:HB2	2.54	0.42
1:C:978:LEU:HD23	1:C:978:LEU:HA	1.89	0.42
1:A:296:ASP:OD1	1:A:297:LYS:N	2.53	0.42
1:A:790:ASP:OD2	1:A:793:LYS:N	2.53	0.42
1:B:983:LEU:HD23	1:B:983:LEU:HA	1.90	0.42
1:C:1111:VAL:HB	1:C:1114:ILE:HD11	2.02	0.42
1:A:517:THR:OG1	1:A:518:ASP:N	2.49	0.42
1:B:352:TYR:HB3	1:B:355:LEU:HG	2.01	0.41
1:B:1052:SER:OG	1:B:1053:GLN:N	2.53	0.41
1:C:410:TYR:HE1	1:C:453:ARG:HD2	1.85	0.41
1:A:760:THR:HG22	1:A:847:LEU:HD23	2.03	0.41
1:B:589:ASN:OD1	1:B:590:ALA:N	2.54	0.41
1:C:44:ASP:N	1:C:44:ASP:OD1	2.52	0.41
1:C:512:GLY:O	1:C:514:LYS:N	2.52	0.41
1:A:194:LEU:HD12	1:A:194:LEU:HA	1.91	0.41
1:A:763:VAL:HG11	1:A:1042:VAL:HG21	2.02	0.41
1:A:782:PHE:CD2	1:A:880:PHE:HE2	2.38	0.41
1:A:889:ASN:HD21	1:A:895:GLN:HB3	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:LEU:HB2	1:B:88:TYR:CZ	2.54	0.41
1:B:556:SER:OG	1:B:557:ASP:N	2.53	0.41
1:C:930:LEU:HD23	1:C:930:LEU:HA	1.91	0.41
1:B:390:LYS:HG3	1:B:481:TYR:CD1	2.53	0.41
1:B:539:THR:HG23	1:B:572:ASP:HB2	2.02	0.41
1:C:822:CYS:HA	1:C:833:CYS:HA	2.03	0.41
1:A:1010:LYS:HE3	1:A:1010:LYS:HB2	1.92	0.41
1:B:19:CYS:SG	1:B:133:CYS:N	2.93	0.41
1:C:941:LEU:HD23	1:C:941:LEU:HA	1.93	0.41
1:A:94:LYS:HB3	1:A:179:PHE:HA	2.01	0.41
1:C:71:THR:HG22	1:C:76:PHE:HA	2.03	0.41
1:C:397:ILE:HG21	1:C:496:VAL:HG11	2.02	0.41
1:A:134:ASP:HA	1:A:232:ARG:HH12	1.86	0.41
1:A:441:ARG:HD3	1:A:444:ARG:HB2	2.02	0.41
1:A:947:GLN:OE1	1:A:985:SER:OG	2.23	0.41
1:B:1065:HIS:CD2	1:B:1119:VAL:H	2.39	0.41
1:C:90:ALA:HB3	1:C:253:PHE:HB2	2.03	0.41
1:A:35:SER:HB2	1:A:60:LEU:HD21	2.02	0.41
1:A:89:PHE:N	1:A:185:PHE:O	2.49	0.41
1:B:92:THR:HG22	1:B:182:LEU:HD13	2.03	0.41
1:C:323:CYS:N	1:C:348:CYS:SG	2.94	0.41
1:A:804:LEU:HA	1:A:804:LEU:HD23	1.89	0.41
1:A:804:LEU:HD13	1:A:1043:VAL:HG21	2.02	0.41
1:A:927:LEU:HD22	1:A:930:LEU:HD12	2.03	0.41
1:A:937:ASN:OD1	1:A:996:ARG:NE	2.53	0.41
1:B:304:ASN:OD1	1:B:304:ASN:N	2.53	0.41
1:A:282:PRO:HG3	1:A:619:TRP:CE2	2.56	0.41
1:A:519:LEU:HD22	1:A:571:LEU:HD11	2.03	0.41
1:B:379:PHE:CD2	1:B:501:PHE:HA	2.56	0.40
1:C:48:ARG:HB2	1:C:266:TYR:CD2	2.56	0.40
1:C:631:THR:HG23	1:C:634:GLY:H	1.85	0.40
1:A:390:LYS:HB2	1:A:481:TYR:CD1	2.56	0.40
1:C:29:ASN:ND2	1:C:70:HIS:HB2	2.32	0.40
1:C:320:THR:HG23	1:C:349:VAL:HG23	2.02	0.40
1:B:297:LYS:HG3	1:B:586:PRO:HA	2.03	0.40
1:B:987:GLN:NE2	1:A:988:THR:OG1	2.54	0.40
1:A:1015:VAL:HG12	1:A:1033:SER:HB3	2.02	0.40
1:B:749:LEU:HD23	1:B:749:LEU:HA	1.89	0.40
1:A:99:ARG:HH12	1:A:170:LEU:HD23	1.86	0.40
1:A:959:LEU:HA	1:A:959:LEU:HD23	1.88	0.40
1:A:925:THR:OG1	1:A:926:ALA:N	2.54	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1052:SER:OG	1:A:1053:GLN:N	2.54	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	1052/1215 (87%)	981 (93%)	71 (7%)	0	100	100
1	B	1054/1215 (87%)	981 (93%)	73 (7%)	0	100	100
1	C	1055/1215 (87%)	983 (93%)	71 (7%)	1 (0%)	51	84
All	All	3161/3645 (87%)	2945 (93%)	215 (7%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	380	SER

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	925/1053 (88%)	923 (100%)	2 (0%)	93	96
1	B	927/1053 (88%)	924 (100%)	3 (0%)	92	95

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	927/1053 (88%)	926 (100%)	1 (0%)	93	97
All	All	2779/3159 (88%)	2773 (100%)	6 (0%)	93	96

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	426	ARG
1	B	576	CYS
1	B	896	ASN
1	C	389	VAL
1	A	189	ASN
1	A	576	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	29	ASN
1	B	550	GLN
1	B	599	GLN
1	B	733	ASN
1	B	987	GLN
1	B	993	GLN
1	B	1088	GLN
1	C	29	ASN
1	C	158	ASN
1	C	201	GLN
1	C	396	GLN
1	C	445	HIS
1	C	530	ASN
1	C	746	ASN
1	C	902	GLN
1	C	937	ASN
1	C	984	GLN
1	A	189	ASN
1	A	227	ASN
1	A	733	ASN
1	A	835	GLN
1	A	877	GLN
1	A	896	ASN
1	A	907	ASN
1	A	931	GLN

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Mol	Chain	Res	Type
1	A	1030	HIS

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)

67 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	$\#\|Z\| > 2$
2	NAG	D	1	1,2	14,14,15	0.25	0	17,19,21	0.48	0
2	NAG	D	2	2	14,14,15	0.28	0	17,19,21	0.50	0
2	NAG	E	1	1,2	14,14,15	0.48	0	17,19,21	0.61	0
2	NAG	E	2	2	14,14,15	0.35	0	17,19,21	0.37	0
3	NAG	F	1	1,3	14,14,15	0.25	0	17,19,21	0.62	0
3	NAG	F	2	3	14,14,15	0.29	0	17,19,21	0.63	0
3	BMA	F	3	3	11,11,12	0.75	0	15,15,17	0.98	1 (6%)
4	NAG	G	1	1,4	14,14,15	1.36	1 (7%)	17,19,21	1.01	1 (5%)
4	NAG	G	2	4	14,14,15	0.23	0	17,19,21	0.57	0
4	BMA	G	3	4	11,11,12	1.09	1 (9%)	15,15,17	1.56	4 (26%)
4	MAN	G	4	4	11,11,12	0.72	0	15,15,17	1.38	2 (13%)
2	NAG	H	1	1,2	14,14,15	0.48	0	17,19,21	0.53	0
2	NAG	H	2	2	14,14,15	0.33	0	17,19,21	0.46	0
2	NAG	I	1	1,2	14,14,15	0.66	0	17,19,21	0.64	0
2	NAG	I	2	2	14,14,15	0.21	0	17,19,21	0.61	0
3	NAG	J	1	1,3	14,14,15	0.57	0	17,19,21	0.79	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	NAG	J	2	3	14,14,15	0.23	0	17,19,21	0.46	0
3	BMA	J	3	3	11,11,12	0.75	0	15,15,17	0.94	1 (6%)
2	NAG	K	1	1,2	14,14,15	0.49	0	17,19,21	0.57	0
2	NAG	K	2	2	14,14,15	0.60	0	17,19,21	0.72	1 (5%)
2	NAG	L	1	1,2	14,14,15	0.35	0	17,19,21	0.63	0
2	NAG	L	2	2	14,14,15	0.40	0	17,19,21	0.56	0
2	NAG	M	1	1,2	14,14,15	0.27	0	17,19,21	0.44	0
2	NAG	M	2	2	14,14,15	0.43	0	17,19,21	0.45	0
4	NAG	N	1	1,4	14,14,15	0.65	1 (7%)	17,19,21	0.50	0
4	NAG	N	2	4	14,14,15	0.38	0	17,19,21	0.58	0
4	BMA	N	3	4	11,11,12	1.01	1 (9%)	15,15,17	0.93	0
4	MAN	N	4	4	11,11,12	0.91	1 (9%)	15,15,17	1.12	2 (13%)
4	NAG	O	1	1,4	14,14,15	0.50	0	17,19,21	0.51	0
4	NAG	O	2	4	14,14,15	0.29	0	17,19,21	0.67	0
4	BMA	O	3	4	11,11,12	0.71	0	15,15,17	0.85	0
4	MAN	O	4	4	11,11,12	0.86	0	15,15,17	0.97	1 (6%)
2	NAG	P	1	1,2	14,14,15	0.33	0	17,19,21	0.94	1 (5%)
2	NAG	P	2	2	14,14,15	0.44	0	17,19,21	0.69	1 (5%)
2	NAG	Q	1	1,2	14,14,15	0.25	0	17,19,21	0.55	0
2	NAG	Q	2	2	14,14,15	0.35	0	17,19,21	0.74	1 (5%)
2	NAG	R	1	1,2	14,14,15	0.64	1 (7%)	17,19,21	0.57	0
2	NAG	R	2	2	14,14,15	0.45	0	17,19,21	0.52	0
2	NAG	S	1	1,2	14,14,15	0.57	0	17,19,21	0.69	0
2	NAG	S	2	2	14,14,15	0.21	0	17,19,21	0.88	1 (5%)
2	NAG	T	1	1,2	14,14,15	0.44	0	17,19,21	0.63	0
2	NAG	T	2	2	14,14,15	0.36	0	17,19,21	0.46	0
2	NAG	U	1	1,2	14,14,15	0.24	0	17,19,21	0.57	0
2	NAG	U	2	2	14,14,15	0.22	0	17,19,21	0.56	0
2	NAG	V	1	1,2	14,14,15	0.22	0	17,19,21	0.44	0
2	NAG	V	2	2	14,14,15	0.28	0	17,19,21	0.48	0
2	NAG	W	1	1,2	14,14,15	0.29	0	17,19,21	0.71	0
2	NAG	W	2	2	14,14,15	0.36	0	17,19,21	0.41	0
4	NAG	X	1	1,4	14,14,15	0.48	0	17,19,21	0.44	0
4	NAG	X	2	4	14,14,15	0.34	0	17,19,21	0.64	1 (5%)
4	BMA	X	3	4	11,11,12	0.96	0	15,15,17	0.93	0
4	MAN	X	4	4	11,11,12	1.47	3 (27%)	15,15,17	1.95	4 (26%)
2	NAG	Y	1	1,2	14,14,15	0.68	1 (7%)	17,19,21	0.81	0
2	NAG	Y	2	2	14,14,15	0.38	0	17,19,21	0.44	0
3	NAG	Z	1	1,3	14,14,15	0.34	0	17,19,21	0.56	0
3	NAG	Z	2	3	14,14,15	0.33	0	17,19,21	0.65	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	BMA	Z	3	3	11,11,12	0.95	1 (9%)	15,15,17	1.98	4 (26%)
4	NAG	a	1	1,4	14,14,15	0.90	1 (7%)	17,19,21	0.66	0
4	NAG	a	2	4	14,14,15	0.21	0	17,19,21	0.58	0
4	BMA	a	3	4	11,11,12	0.94	1 (9%)	15,15,17	2.04	4 (26%)
4	MAN	a	4	4	11,11,12	0.88	0	15,15,17	1.27	2 (13%)
2	NAG	b	1	1,2	14,14,15	0.54	0	17,19,21	0.66	0
2	NAG	b	2	2	14,14,15	0.35	0	17,19,21	0.50	0
2	NAG	c	1	1,2	14,14,15	0.26	0	17,19,21	0.40	0
2	NAG	c	2	2	14,14,15	0.21	0	17,19,21	0.63	1 (5%)
2	NAG	d	1	1,2	14,14,15	0.92	1 (7%)	17,19,21	0.59	0
2	NAG	d	2	2	14,14,15	0.25	0	17,19,21	0.42	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
2	NAG	D	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	D	2	2	-	0/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	2/2/19/22	0/1/1/1
4	NAG	G	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
4	BMA	G	3	4	-	1/2/19/22	0/1/1/1
4	MAN	G	4	4	-	2/2/19/22	0/1/1/1
2	NAG	H	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
3	NAG	J	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	J	2	3	-	0/6/23/26	0/1/1/1
3	BMA	J	3	3	-	2/2/19/22	0/1/1/1
2	NAG	K	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	K	2	2	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	L	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	L	2	2	-	2/6/23/26	0/1/1/1
2	NAG	M	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	M	2	2	-	1/6/23/26	0/1/1/1
4	NAG	N	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	N	2	4	-	2/6/23/26	0/1/1/1
4	BMA	N	3	4	-	1/2/19/22	0/1/1/1
4	MAN	N	4	4	-	0/2/19/22	0/1/1/1
4	NAG	O	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	O	2	4	-	2/6/23/26	0/1/1/1
4	BMA	O	3	4	-	0/2/19/22	0/1/1/1
4	MAN	O	4	4	-	0/2/19/22	0/1/1/1
2	NAG	P	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	P	2	2	-	0/6/23/26	0/1/1/1
2	NAG	Q	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	Q	2	2	-	2/6/23/26	0/1/1/1
2	NAG	R	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	R	2	2	-	0/6/23/26	0/1/1/1
2	NAG	S	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	S	2	2	-	2/6/23/26	0/1/1/1
2	NAG	T	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	T	2	2	-	2/6/23/26	0/1/1/1
2	NAG	U	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	U	2	2	-	2/6/23/26	0/1/1/1
2	NAG	V	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	V	2	2	-	2/6/23/26	0/1/1/1
2	NAG	W	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	W	2	2	-	2/6/23/26	0/1/1/1
4	NAG	X	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	X	2	4	-	0/6/23/26	0/1/1/1
4	BMA	X	3	4	-	2/2/19/22	0/1/1/1
4	MAN	X	4	4	-	2/2/19/22	0/1/1/1
2	NAG	Y	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	Y	2	2	-	2/6/23/26	0/1/1/1
3	NAG	Z	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	Z	2	3	-	2/6/23/26	0/1/1/1
3	BMA	Z	3	3	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	a	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	a	2	4	-	0/6/23/26	0/1/1/1
4	BMA	a	3	4	-	1/2/19/22	0/1/1/1
4	MAN	a	4	4	-	2/2/19/22	0/1/1/1
2	NAG	b	1	1,2	-	1/6/23/26	0/1/1/1
2	NAG	b	2	2	-	2/6/23/26	0/1/1/1
2	NAG	c	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	c	2	2	-	2/6/23/26	0/1/1/1
2	NAG	d	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	d	2	2	-	2/6/23/26	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	1	NAG	O5-C1	-4.96	1.35	1.43
4	X	4	MAN	C1-C2	3.30	1.59	1.52
2	d	1	NAG	O5-C1	-3.29	1.38	1.43
4	a	1	NAG	O5-C1	-3.17	1.38	1.43
4	G	3	BMA	C2-C3	2.65	1.56	1.52
4	N	3	BMA	O5-C5	2.52	1.48	1.43
3	Z	3	BMA	C1-C2	2.51	1.57	1.52
4	X	4	MAN	O5-C1	2.33	1.47	1.43
2	R	1	NAG	O5-C1	-2.30	1.40	1.43
2	Y	1	NAG	O5-C1	-2.29	1.40	1.43
4	N	1	NAG	O5-C1	-2.26	1.40	1.43
4	X	4	MAN	O5-C5	2.20	1.47	1.43
4	N	4	MAN	C1-C2	2.01	1.56	1.52
4	a	3	BMA	C2-C3	2.00	1.55	1.52

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	X	4	MAN	C1-O5-C5	5.86	120.14	112.19
4	a	3	BMA	C1-C2-C3	4.64	115.37	109.67
3	Z	3	BMA	C1-O5-C5	4.32	118.05	112.19
3	Z	3	BMA	C1-C2-C3	4.13	114.75	109.67
4	G	4	MAN	C1-O5-C5	4.05	117.69	112.19
4	a	4	MAN	C1-O5-C5	3.75	117.28	112.19
4	a	3	BMA	O5-C1-C2	3.59	116.31	110.77
4	G	3	BMA	C1-C2-C3	3.50	113.97	109.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	a	3	BMA	C1-O5-C5	3.49	116.93	112.19
2	S	2	NAG	C1-O5-C5	3.31	116.68	112.19
2	P	1	NAG	C1-O5-C5	3.12	116.41	112.19
3	Z	3	BMA	O5-C1-C2	3.03	115.45	110.77
4	G	3	BMA	C2-C3-C4	2.81	115.77	110.89
4	N	4	MAN	C1-O5-C5	2.76	115.93	112.19
3	J	1	NAG	C1-O5-C5	2.71	115.86	112.19
2	Q	2	NAG	C1-O5-C5	2.66	115.80	112.19
4	X	4	MAN	C1-C2-C3	2.64	112.91	109.67
4	G	1	NAG	O4-C4-C3	-2.53	104.50	110.35
4	G	4	MAN	O2-C2-C3	-2.49	105.16	110.14
4	N	4	MAN	O2-C2-C3	-2.48	105.17	110.14
3	J	3	BMA	C1-O5-C5	2.39	115.44	112.19
3	F	3	BMA	C1-O5-C5	2.39	115.43	112.19
2	K	2	NAG	C1-O5-C5	2.33	115.34	112.19
4	X	4	MAN	O5-C1-C2	2.23	114.22	110.77
4	X	2	NAG	C1-O5-C5	2.19	115.16	112.19
2	c	2	NAG	C1-O5-C5	2.18	115.15	112.19
4	O	4	MAN	O2-C2-C3	-2.17	105.78	110.14
4	G	3	BMA	O5-C1-C2	2.17	114.12	110.77
4	G	3	BMA	O5-C5-C4	-2.14	105.62	110.83
4	X	4	MAN	O2-C2-C3	-2.05	106.02	110.14
4	a	3	BMA	C2-C3-C4	2.05	114.44	110.89
2	P	2	NAG	C1-O5-C5	2.04	114.96	112.19
4	a	4	MAN	O2-C2-C3	-2.03	106.07	110.14
3	Z	3	BMA	O2-C2-C3	-2.02	106.09	110.14

There are no chirality outliers.

All (94) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	Q	2	NAG	O5-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
3	J	1	NAG	O5-C5-C6-O6
4	N	1	NAG	O5-C5-C6-O6
2	S	1	NAG	O5-C5-C6-O6
2	V	1	NAG	O5-C5-C6-O6
3	Z	2	NAG	O5-C5-C6-O6
2	D	1	NAG	O5-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	U	2	NAG	O5-C5-C6-O6
2	c	2	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
3	F	3	BMA	O5-C5-C6-O6
4	X	4	MAN	O5-C5-C6-O6
4	a	4	MAN	O5-C5-C6-O6
4	a	1	NAG	C4-C5-C6-O6
2	b	2	NAG	O5-C5-C6-O6
2	d	1	NAG	O5-C5-C6-O6
2	d	2	NAG	O5-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
2	U	2	NAG	C4-C5-C6-O6
2	V	1	NAG	C4-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
4	G	4	MAN	O5-C5-C6-O6
4	N	2	NAG	O5-C5-C6-O6
4	a	1	NAG	O5-C5-C6-O6
2	S	1	NAG	C4-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
3	F	3	BMA	C4-C5-C6-O6
3	J	1	NAG	C4-C5-C6-O6
3	Z	2	NAG	C4-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
2	W	2	NAG	O5-C5-C6-O6
2	Y	2	NAG	O5-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
2	Q	2	NAG	C4-C5-C6-O6
2	Y	1	NAG	C4-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6
4	N	1	NAG	C4-C5-C6-O6
4	X	3	BMA	C4-C5-C6-O6
2	Q	1	NAG	O5-C5-C6-O6
2	c	1	NAG	O5-C5-C6-O6
2	b	2	NAG	C4-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
4	O	1	NAG	O5-C5-C6-O6
2	D	1	NAG	C4-C5-C6-O6
4	N	2	NAG	C4-C5-C6-O6
2	d	1	NAG	C4-C5-C6-O6
4	a	4	MAN	C4-C5-C6-O6
2	Y	2	NAG	C4-C5-C6-O6
4	O	1	NAG	C4-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	K	2	NAG	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
2	c	1	NAG	C4-C5-C6-O6
4	X	4	MAN	C4-C5-C6-O6
2	c	2	NAG	C4-C5-C6-O6
2	W	1	NAG	C4-C5-C6-O6
2	Y	1	NAG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
2	Q	1	NAG	C4-C5-C6-O6
2	W	2	NAG	C4-C5-C6-O6
2	d	2	NAG	C4-C5-C6-O6
2	T	2	NAG	O5-C5-C6-O6
2	V	2	NAG	O5-C5-C6-O6
2	S	2	NAG	O5-C5-C6-O6
3	J	3	BMA	O5-C5-C6-O6
2	S	2	NAG	C4-C5-C6-O6
2	T	2	NAG	C4-C5-C6-O6
2	L	1	NAG	C4-C5-C6-O6
3	J	3	BMA	C4-C5-C6-O6
4	X	3	BMA	O5-C5-C6-O6
2	L	1	NAG	O5-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
4	a	3	BMA	O5-C5-C6-O6
4	G	4	MAN	C4-C5-C6-O6
4	G	3	BMA	O5-C5-C6-O6
2	W	1	NAG	O5-C5-C6-O6
4	N	3	BMA	O5-C5-C6-O6
2	V	2	NAG	C4-C5-C6-O6
2	b	1	NAG	O5-C5-C6-O6
4	O	2	NAG	C4-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	M	1	NAG	O5-C5-C6-O6
2	P	1	NAG	O5-C5-C6-O6
2	K	1	NAG	C4-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	L	2	NAG	C4-C5-C6-O6
4	O	2	NAG	O5-C5-C6-O6
2	E	2	NAG	O5-C5-C6-O6
2	L	2	NAG	O5-C5-C6-O6
2	T	1	NAG	C4-C5-C6-O6
2	K	1	NAG	O5-C5-C6-O6
2	M	2	NAG	C4-C5-C6-O6
2	R	1	NAG	C4-C5-C6-O6

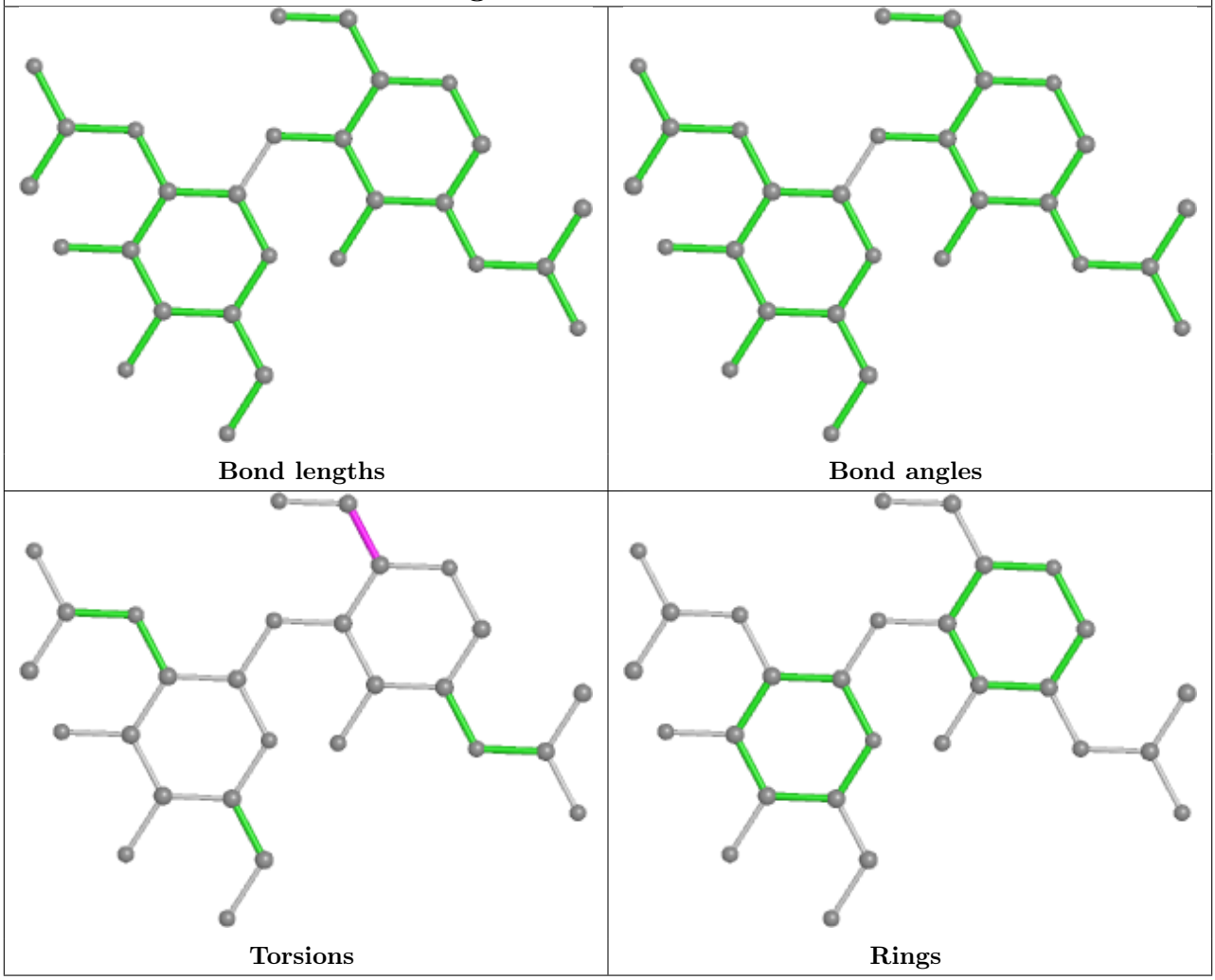
There are no ring outliers.

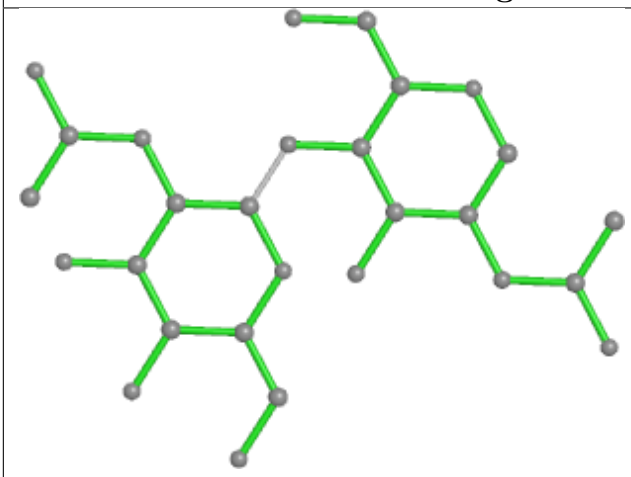
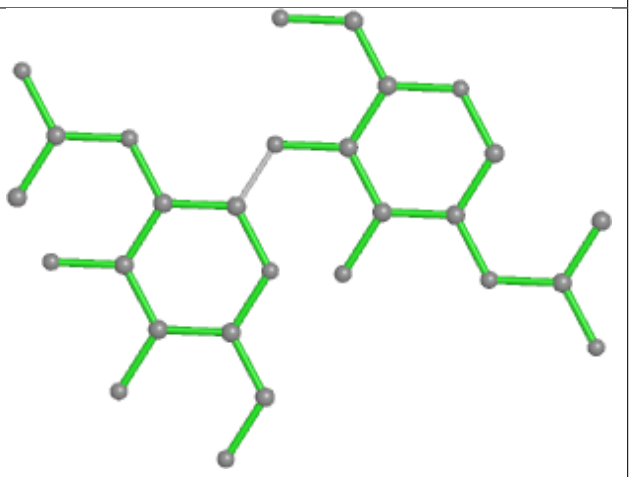
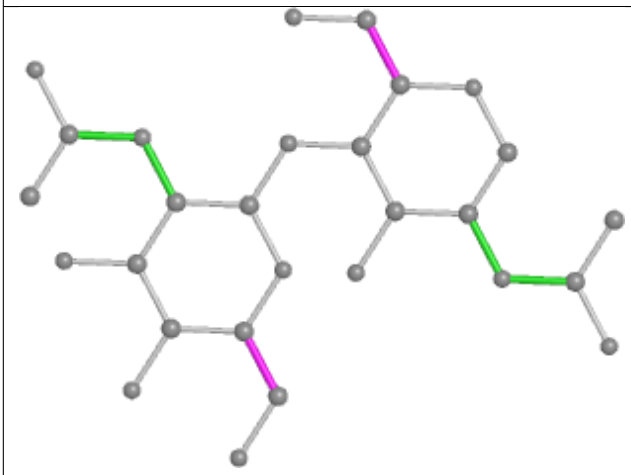
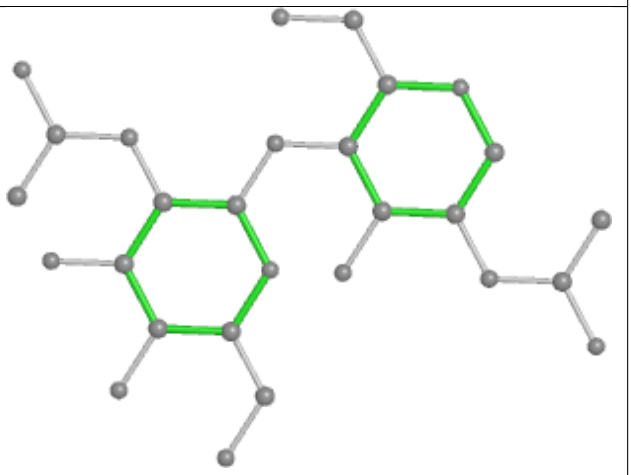
2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	T	1	NAG	1	0
2	U	1	NAG	2	0

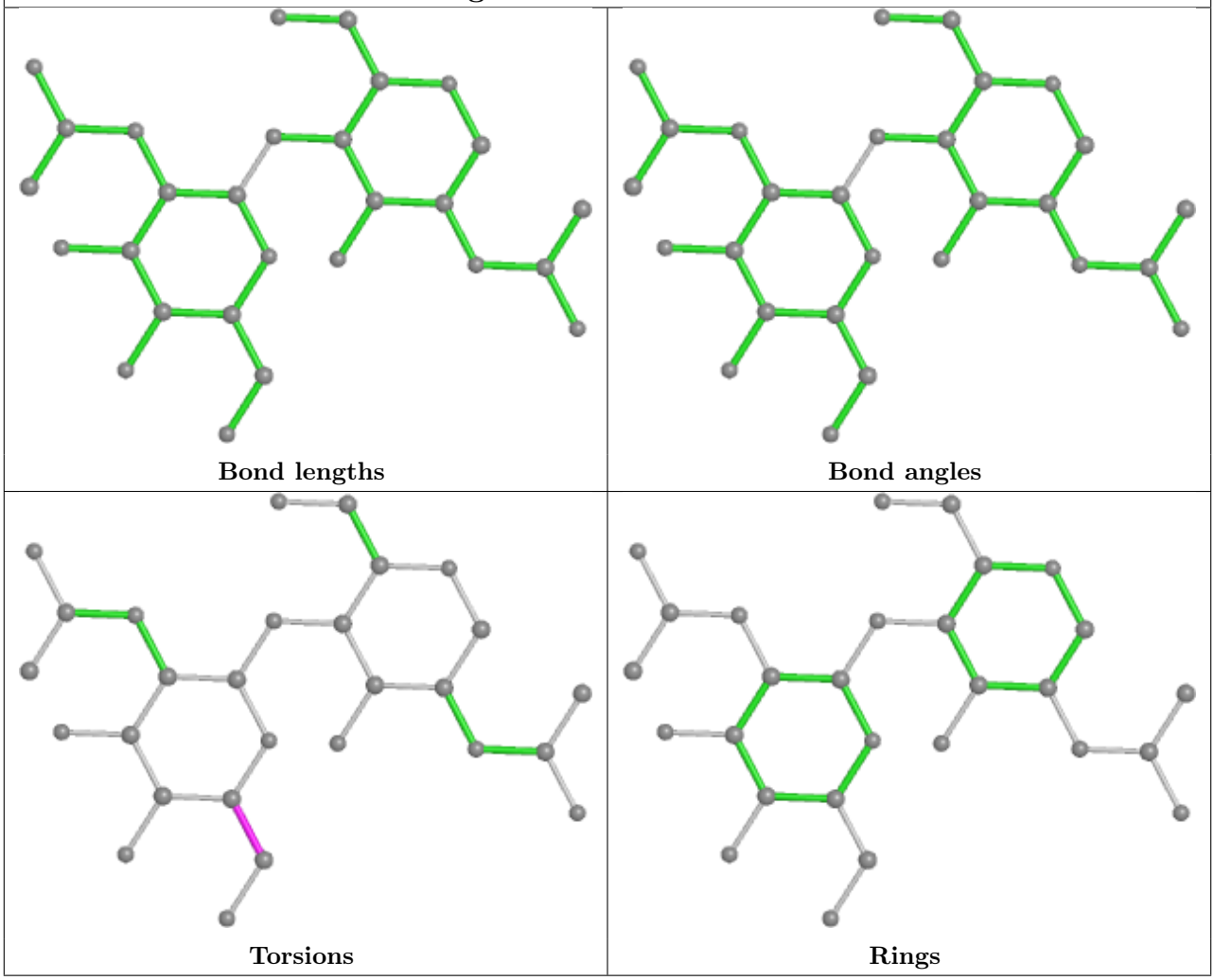
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

Oligosaccharide Chain D

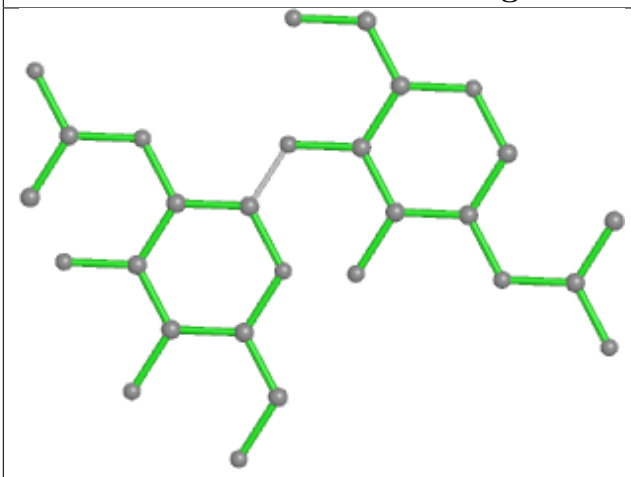


Oligosaccharide Chain E	
 Bond lengths	 Bond angles
 Torsions	 Rings

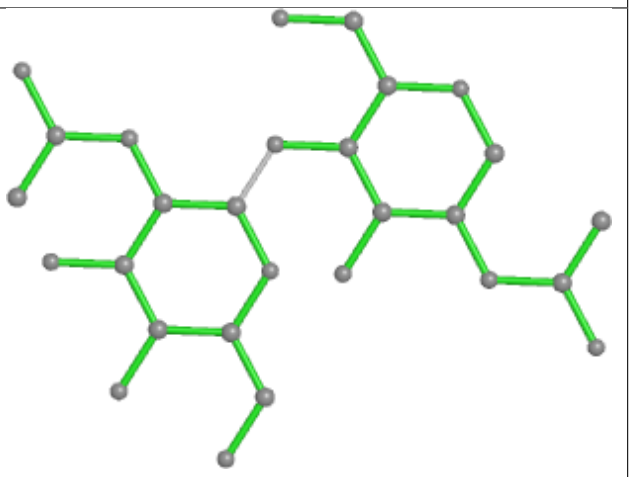
Oligosaccharide Chain H



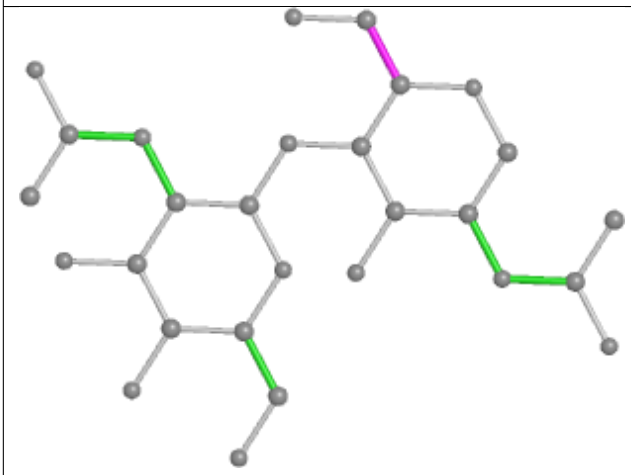
Oligosaccharide Chain I



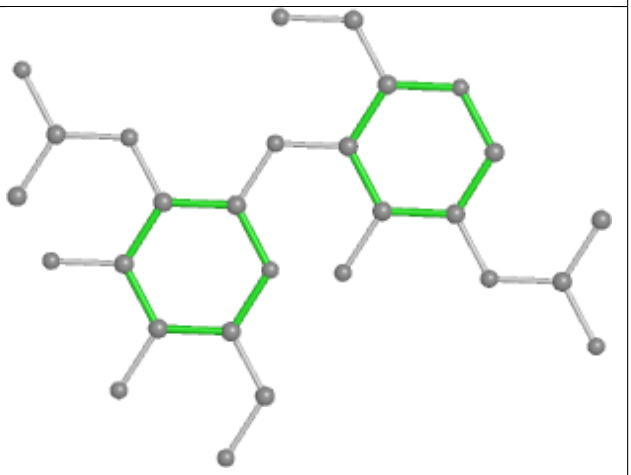
Bond lengths



Bond angles

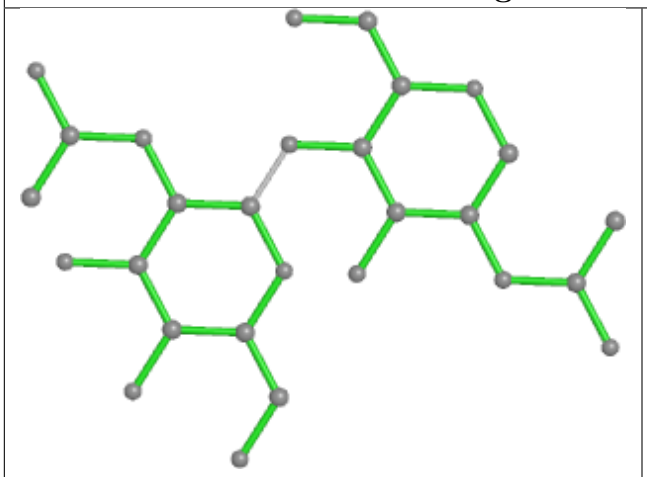


Torsions

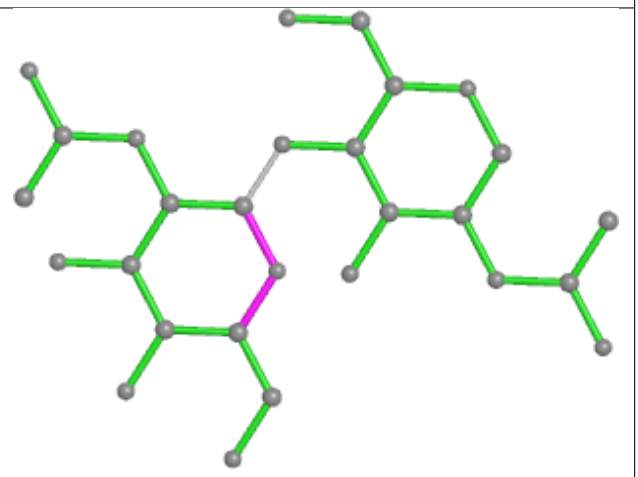


Rings

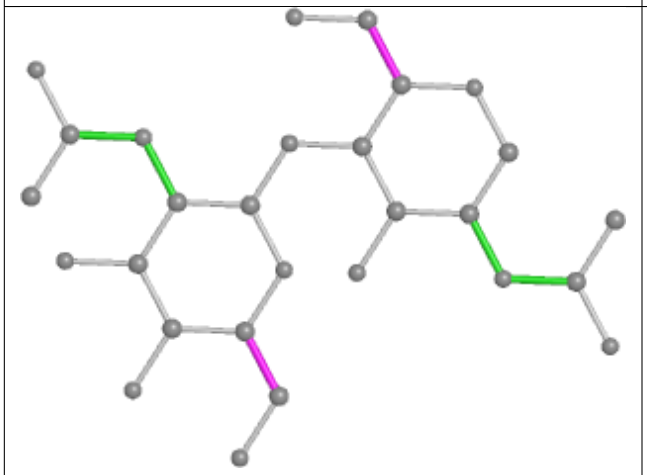
Oligosaccharide Chain K



Bond lengths



Bond angles



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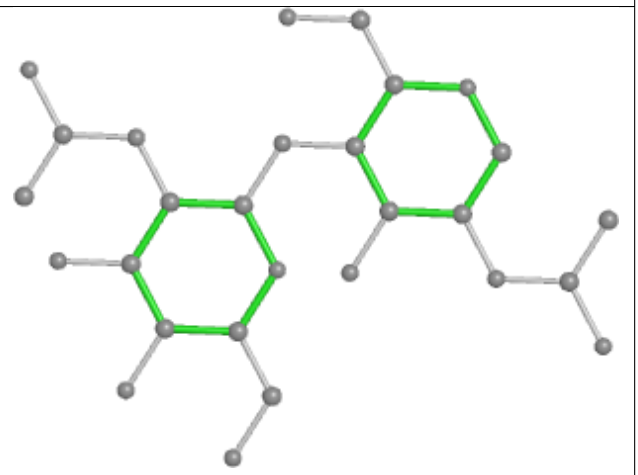
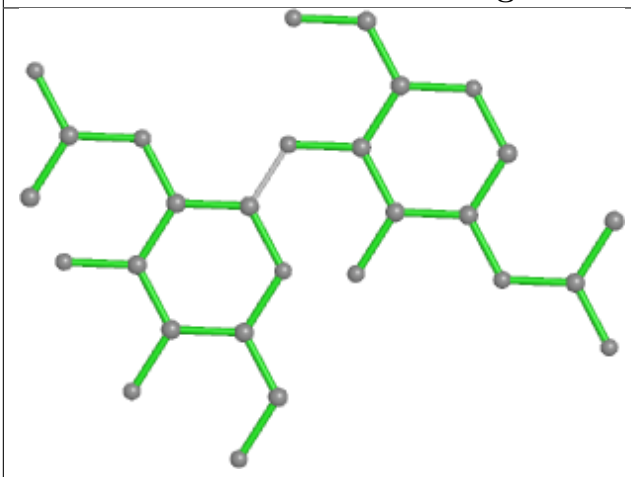
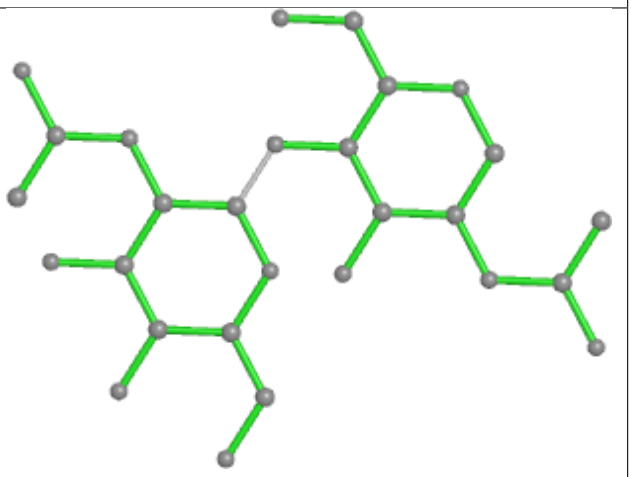
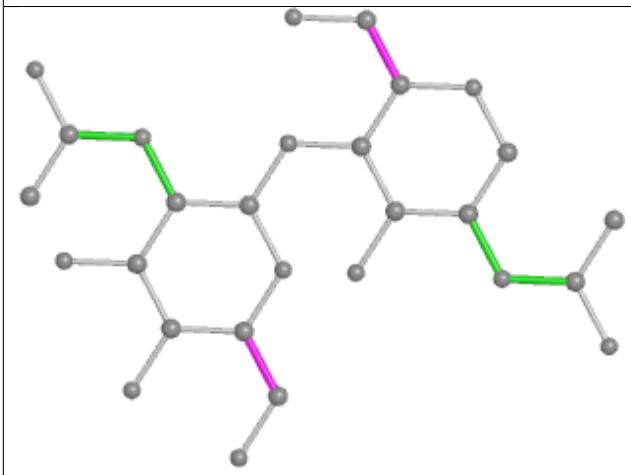
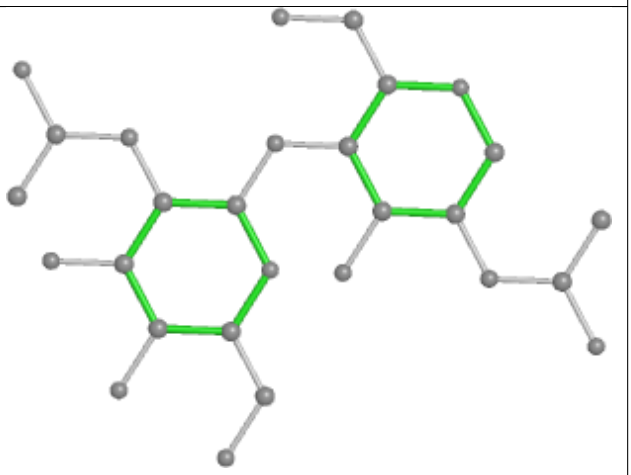


Rings

Oligosaccharide Chain L	
 Bond lengths	 Bond angles
 Torsions	 Rings

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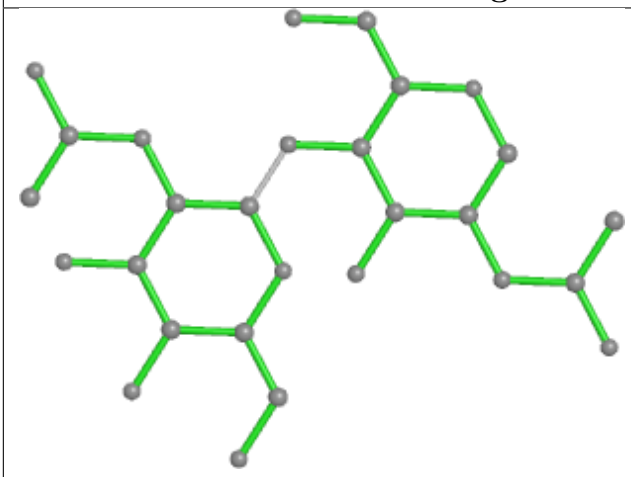
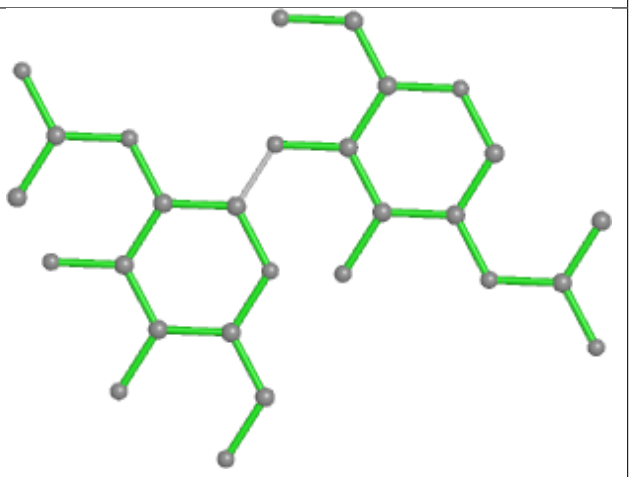
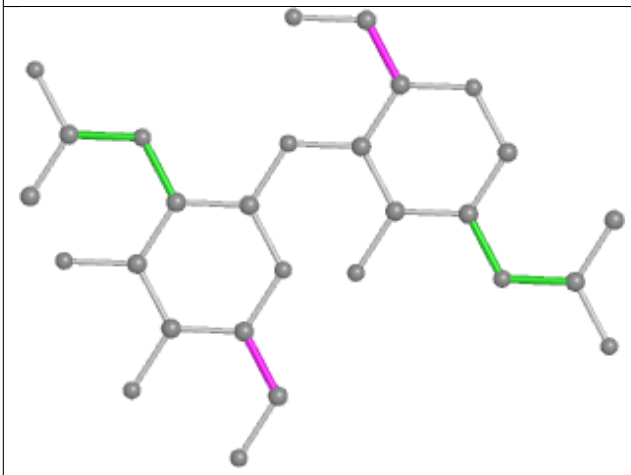
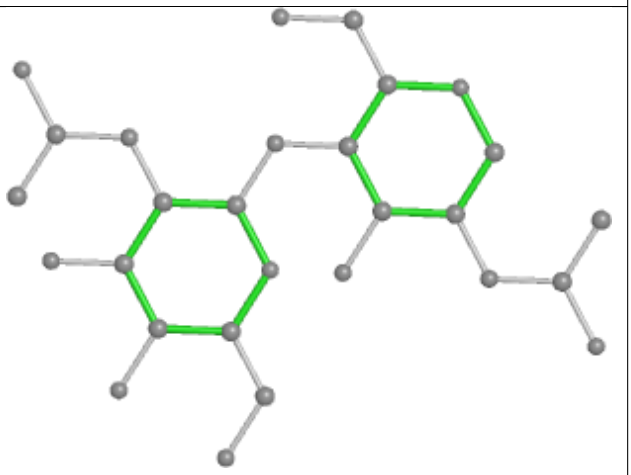
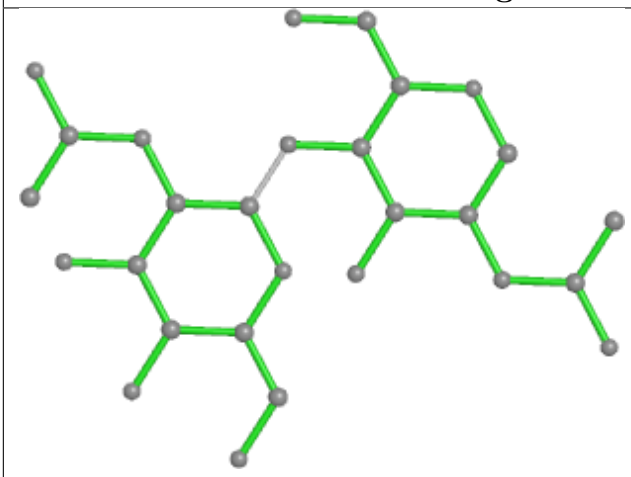
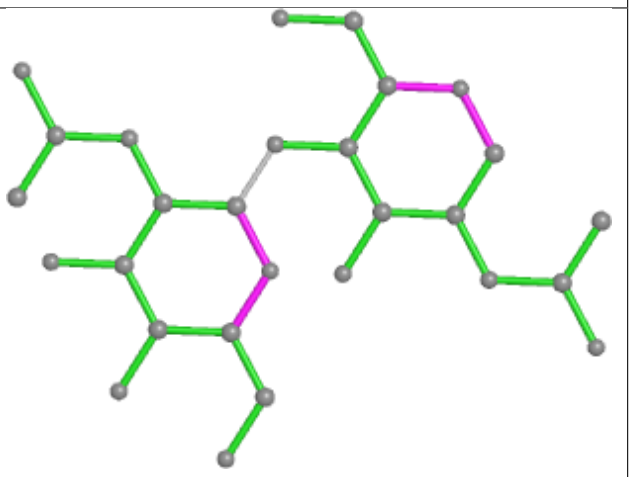
 Bond lengths	 Bond angles
 Torsions	 Rings

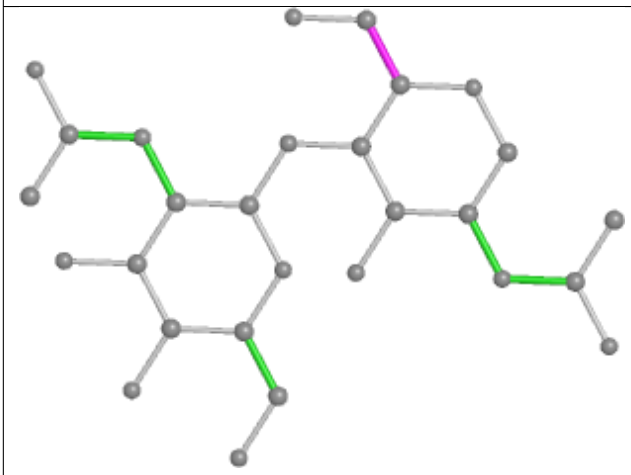
Oligosaccharide Chain P



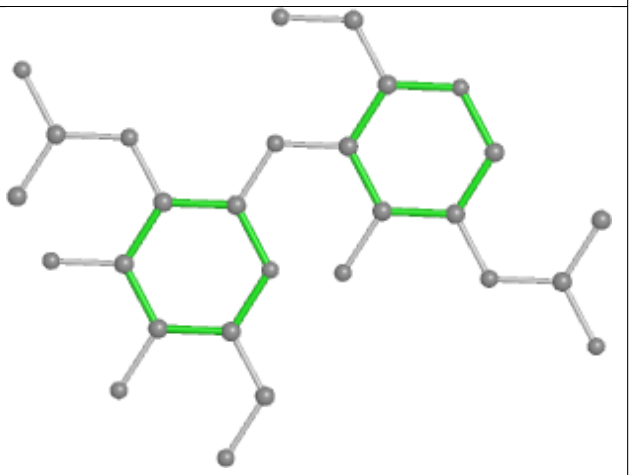
Bond lengths



Bond angles

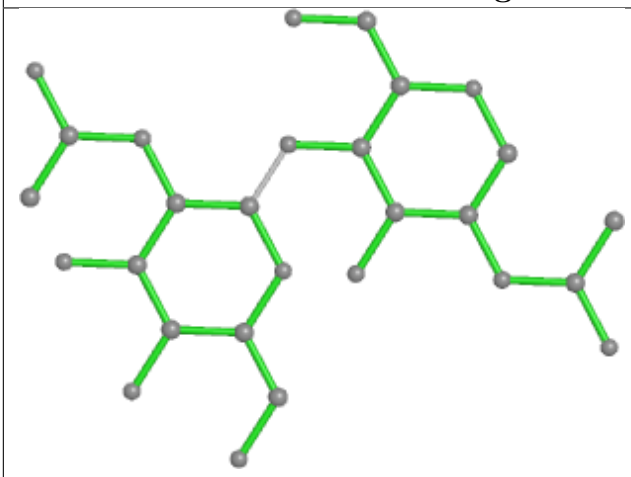


Torsions



Rings

Oligosaccharide Chain Q



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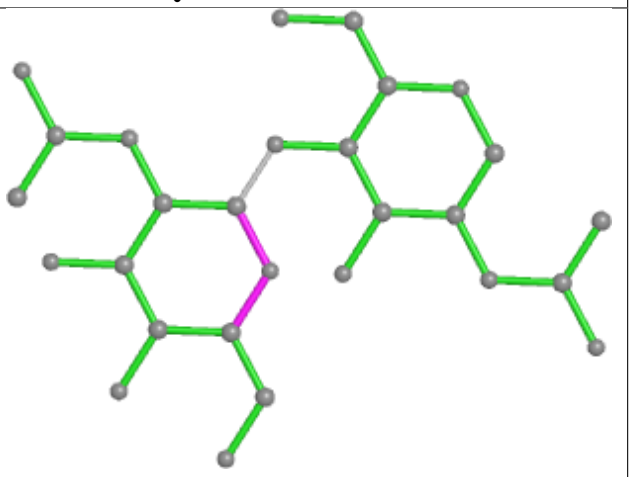

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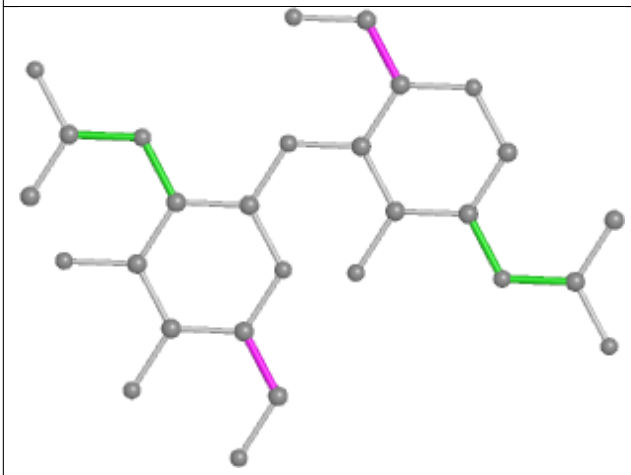

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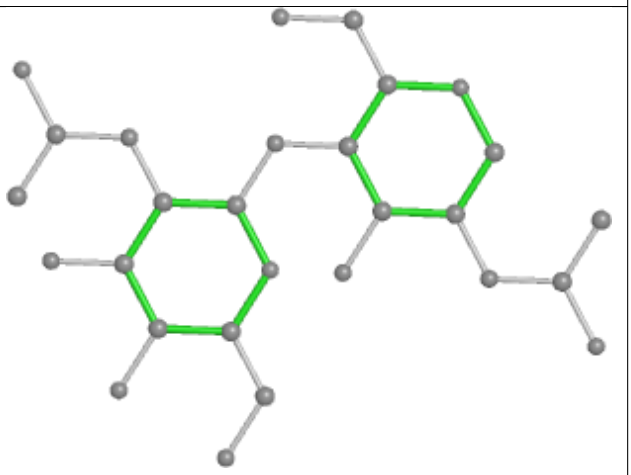
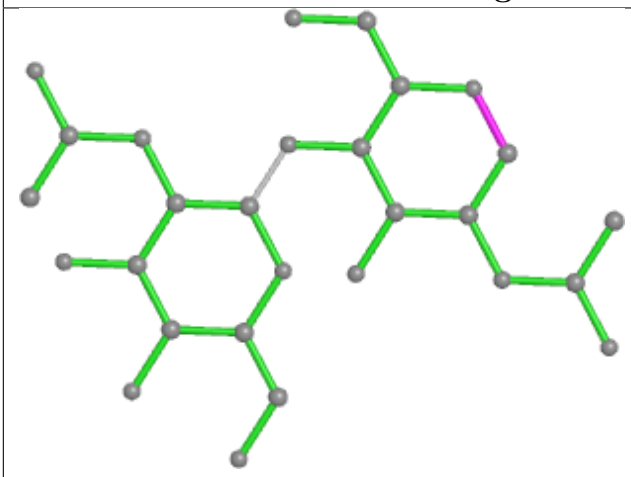
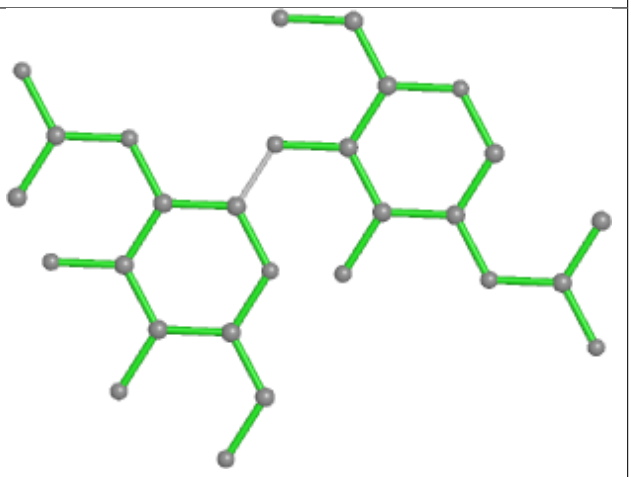
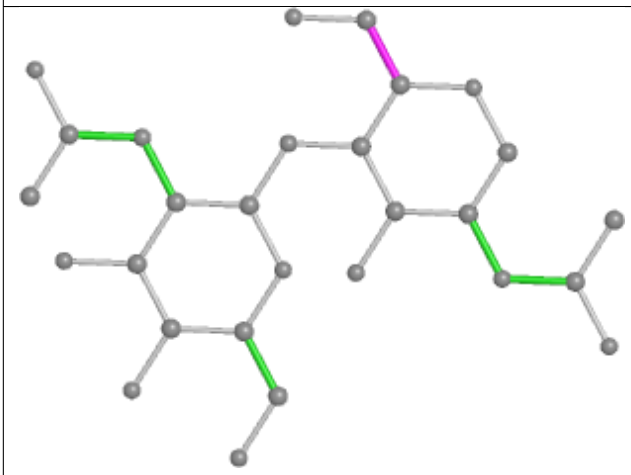
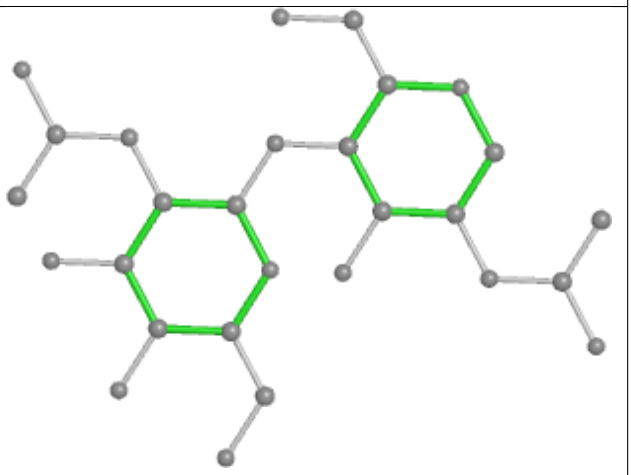
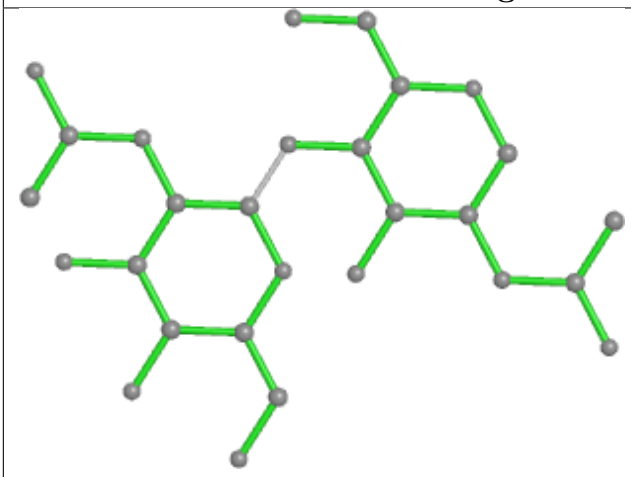


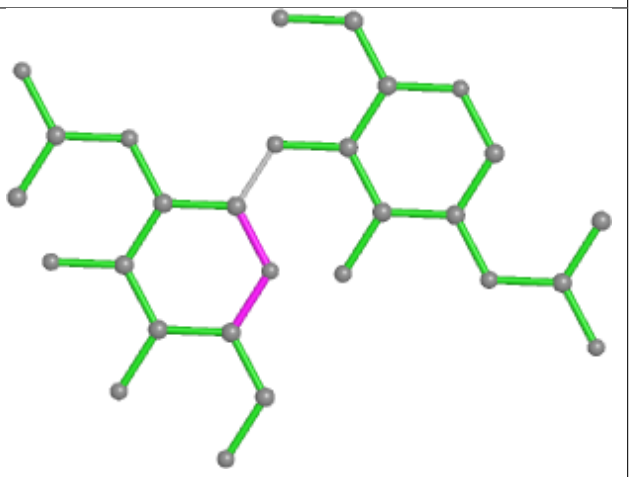
Rings

Oligosaccharide Chain R	
 Bond lengths	 Bond angles
 Torsions	 Rings

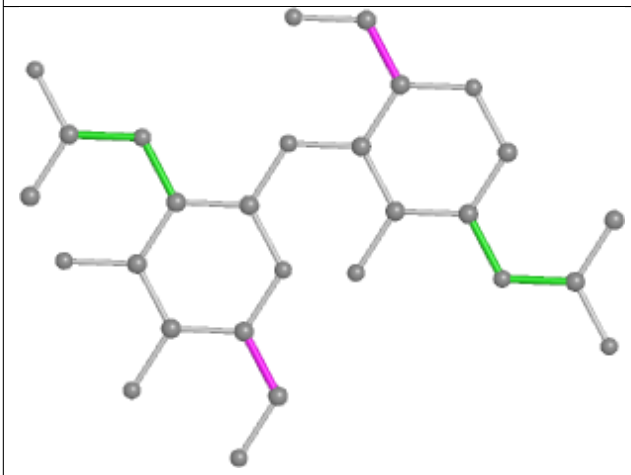
Oligosaccharide Chain S



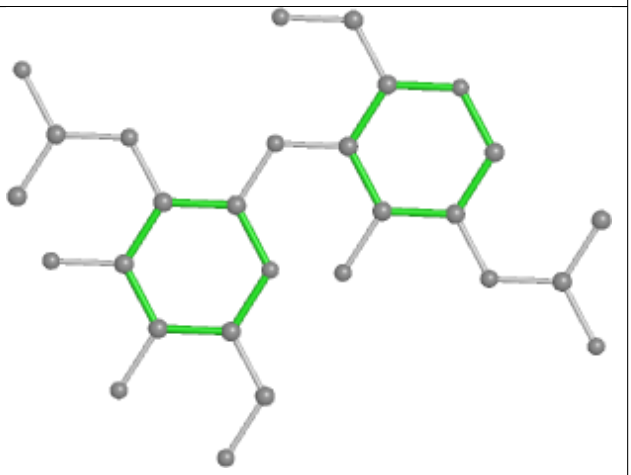
Bond lengths



Bond angles

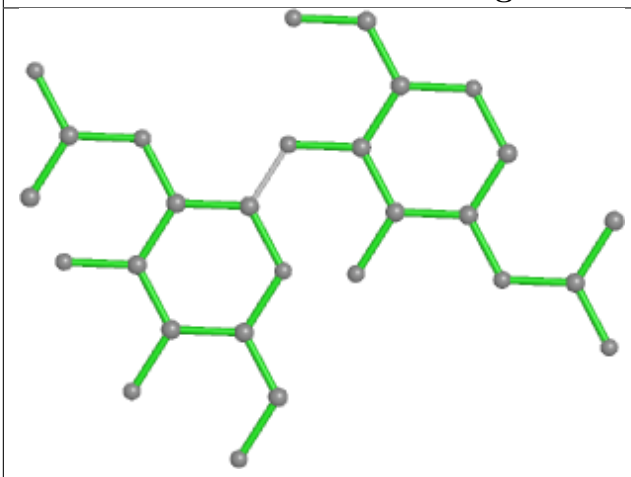


Torsions

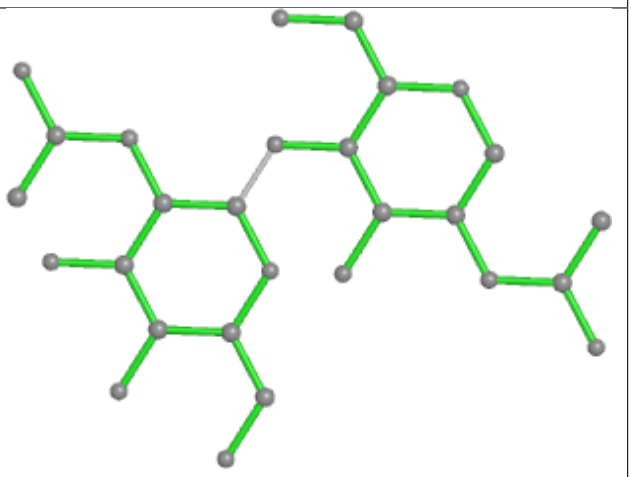


Rings

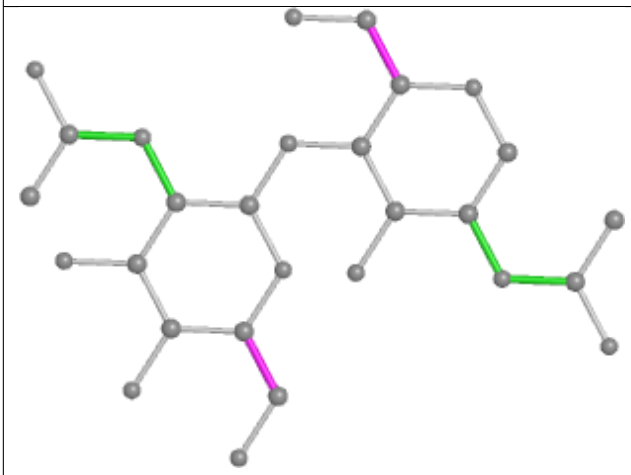
Oligosaccharide Chain T



Bond lengths



Bond angles



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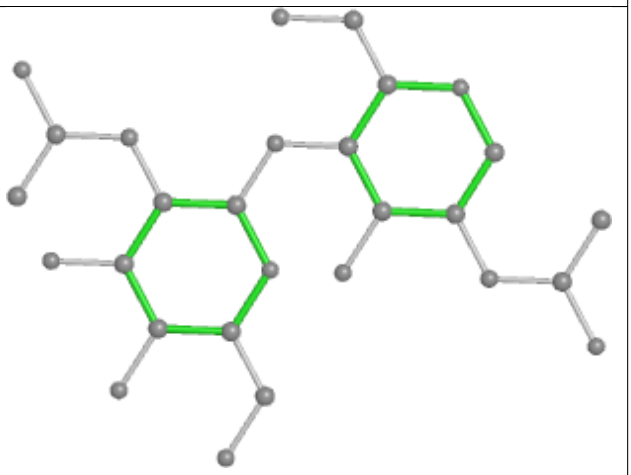
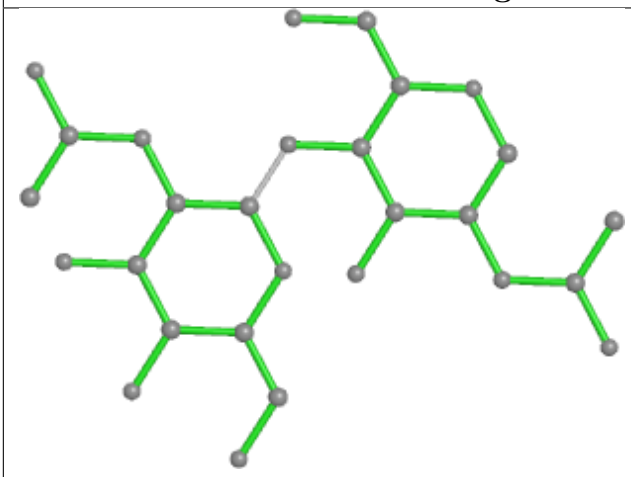
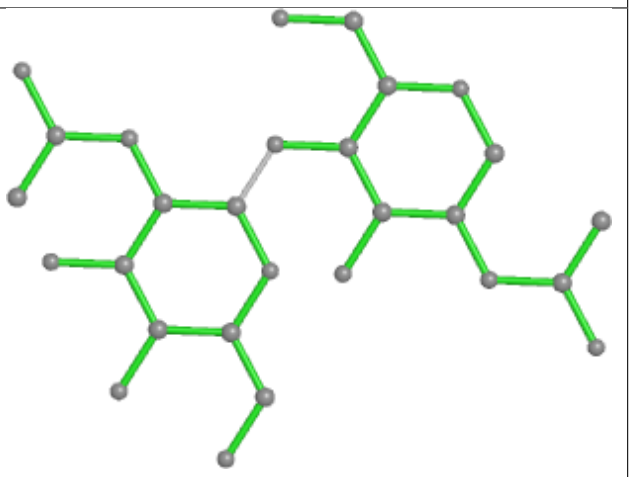
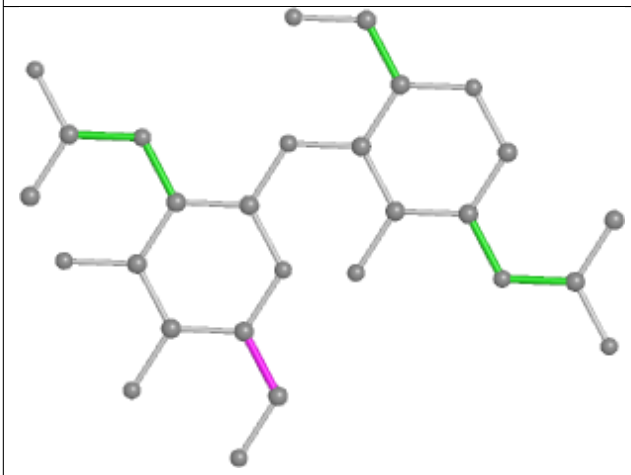
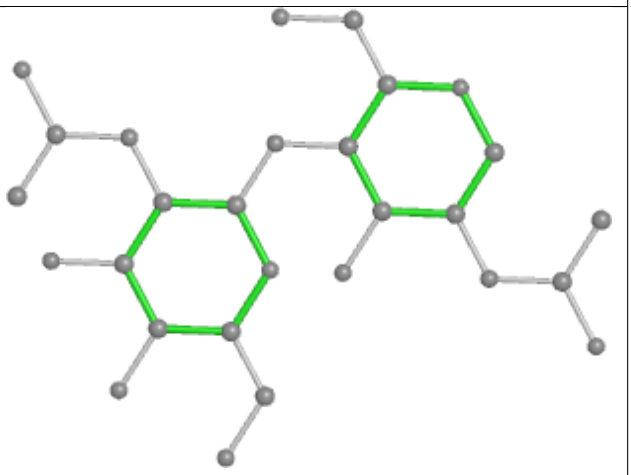
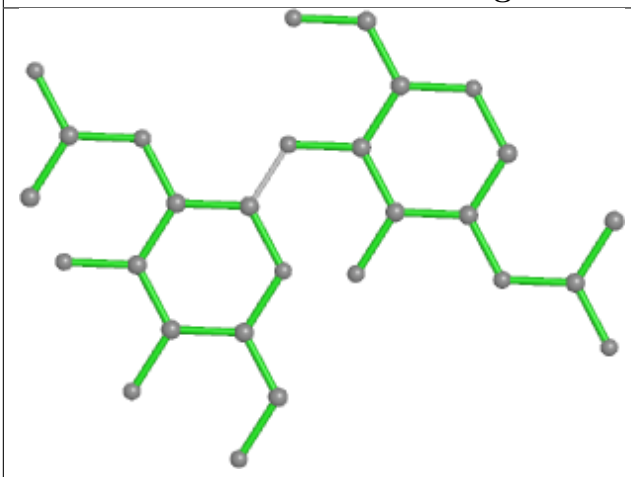


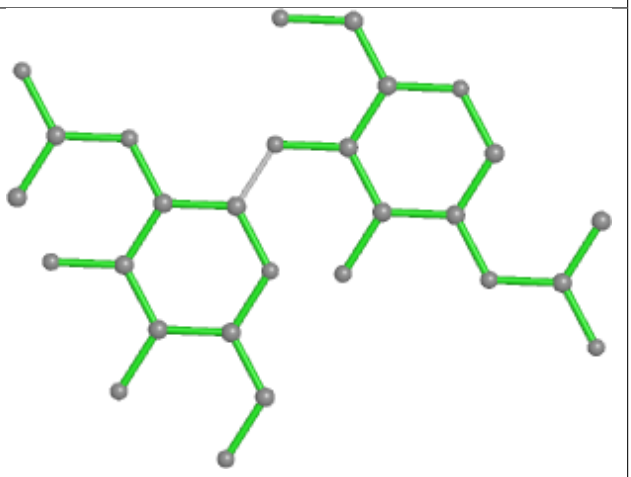
Rings

Oligosaccharide Chain U	
 Bond lengths	 Bond angles
 Torsions	 Rings

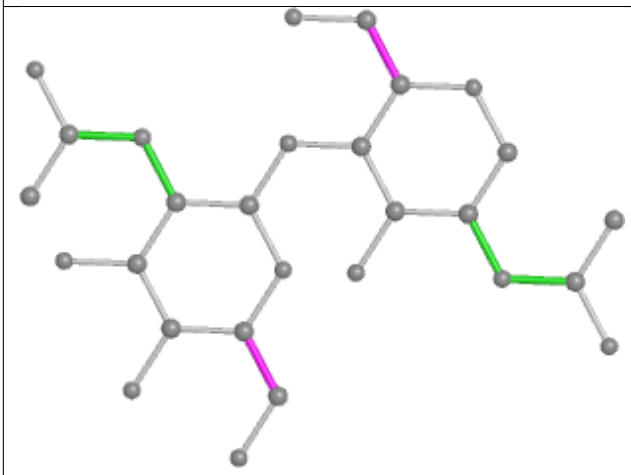
Oligosaccharide Chain V



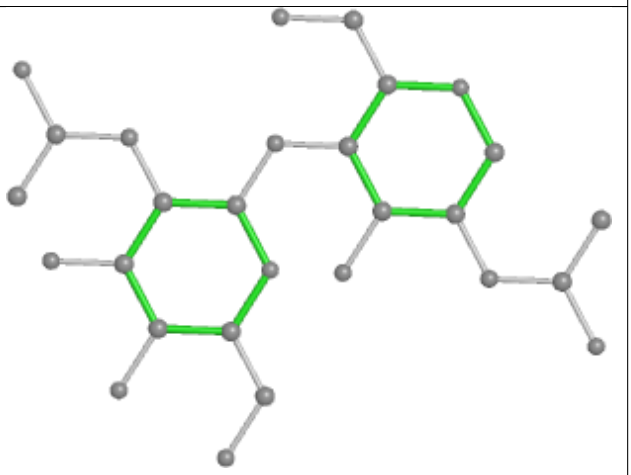
Bond lengths



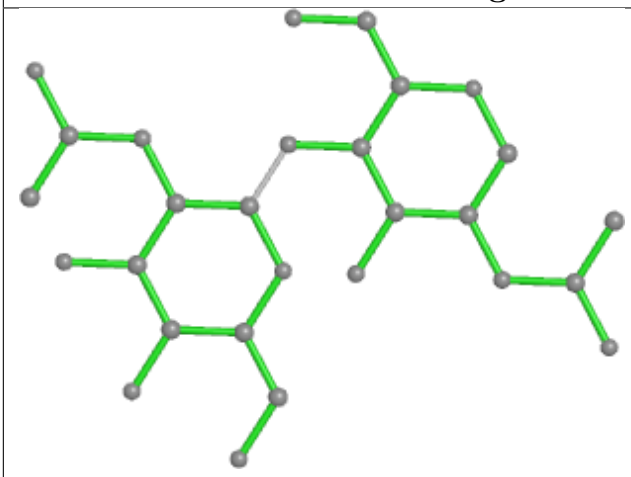
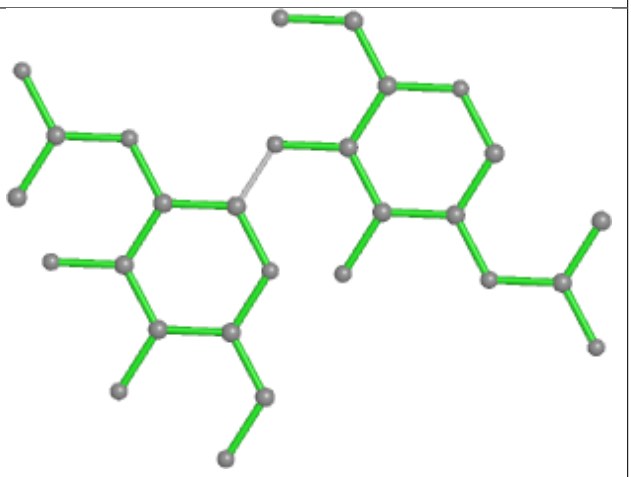
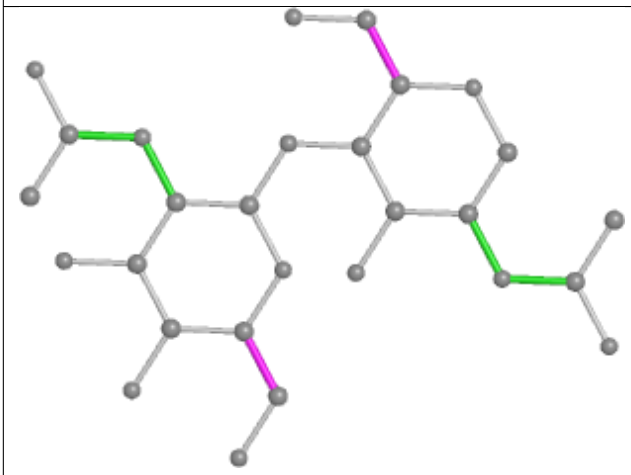
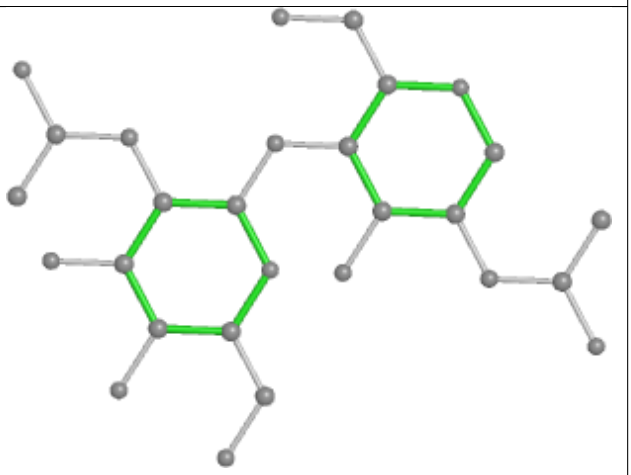
Bond angles

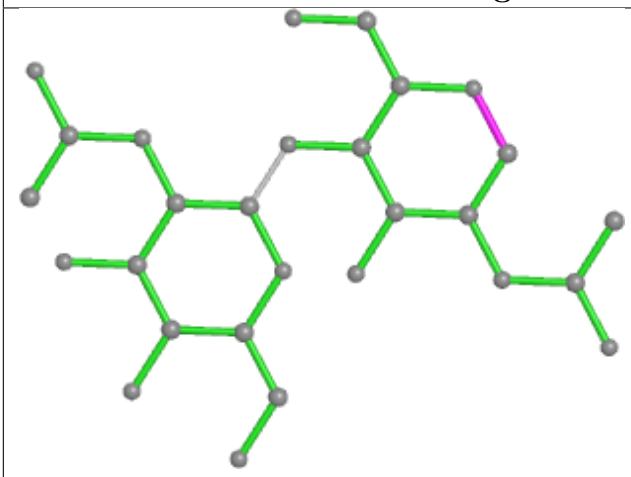
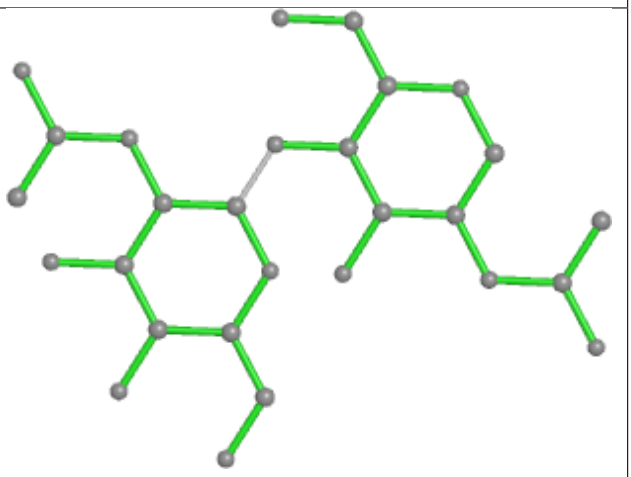
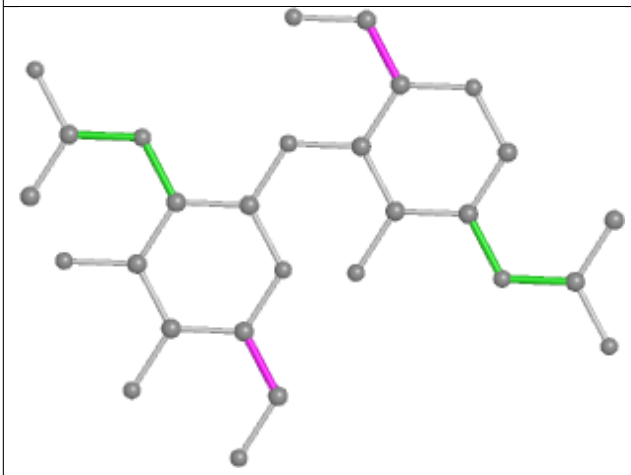
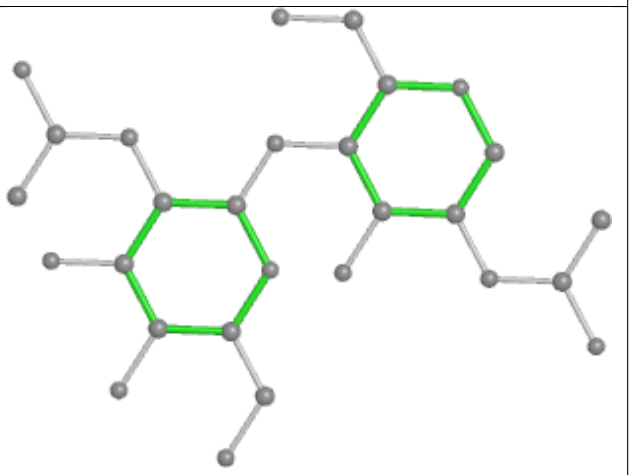


Torsions

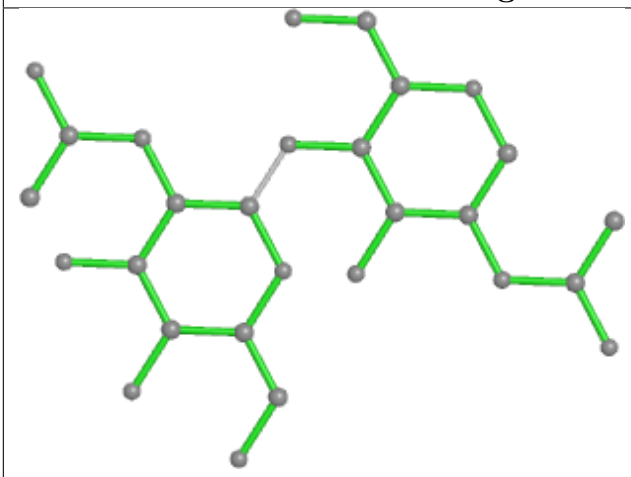


Rings

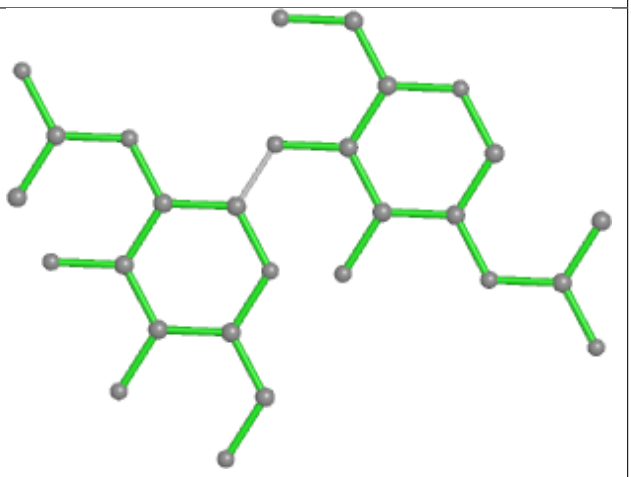
Oligosaccharide Chain W	
 Bond lengths	 Bond angles
 Torsions	 Rings

Oligosaccharide Chain Y	
 Bond lengths	 Bond angles
 Torsions	 Rings

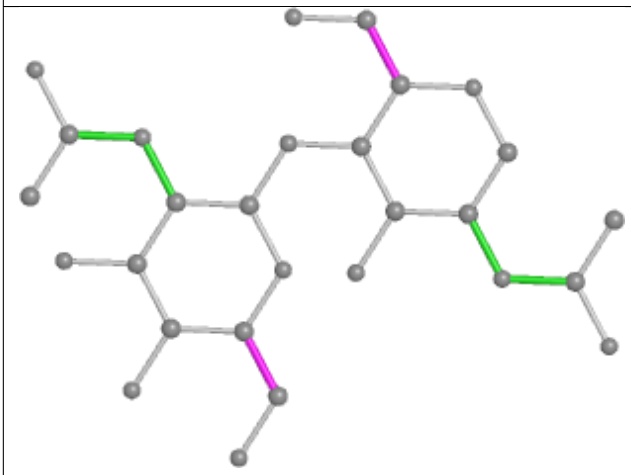
Oligosaccharide Chain b



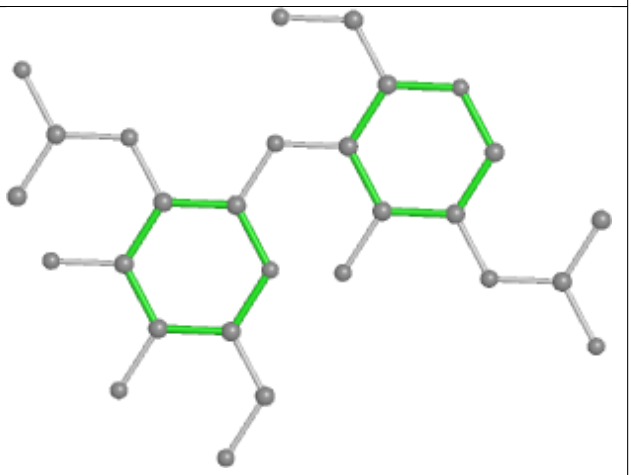
Bond lengths



Bond angles

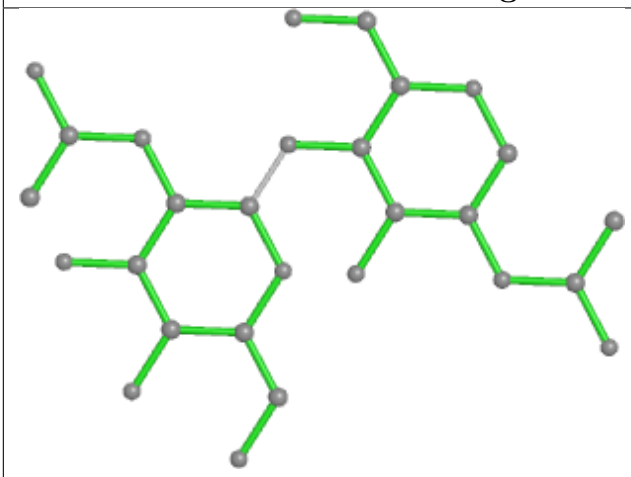


Torsions

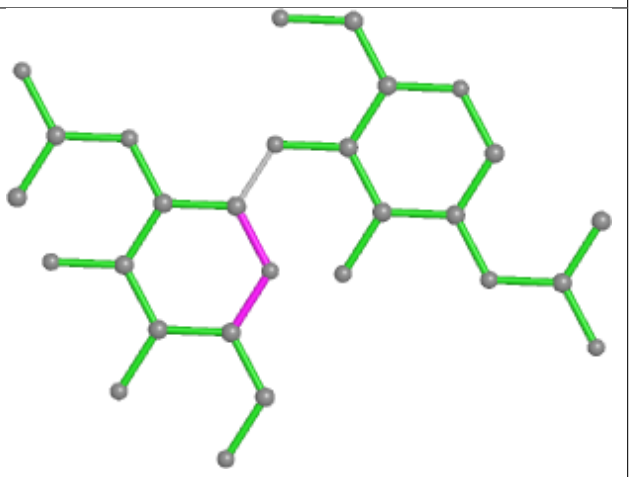


Rings

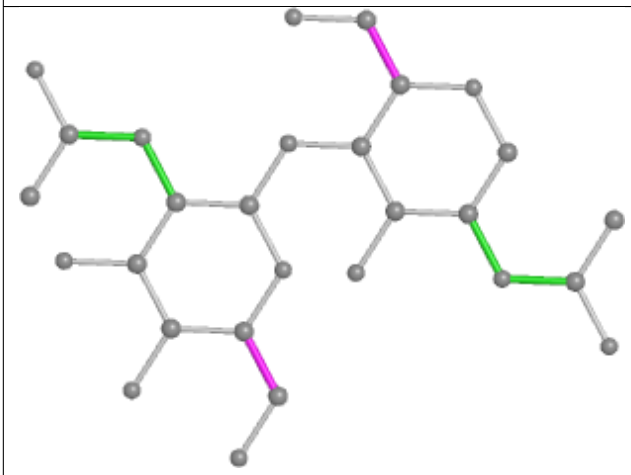
Oligosaccharide Chain c



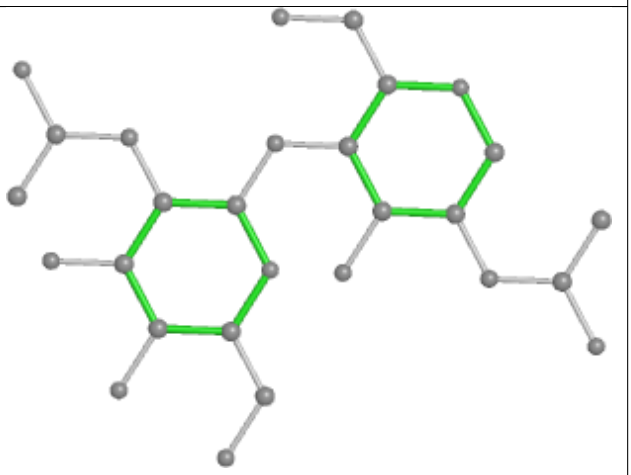
Bond lengths



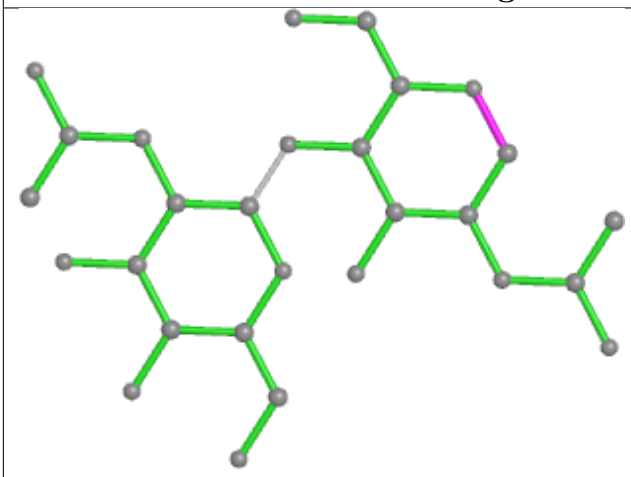
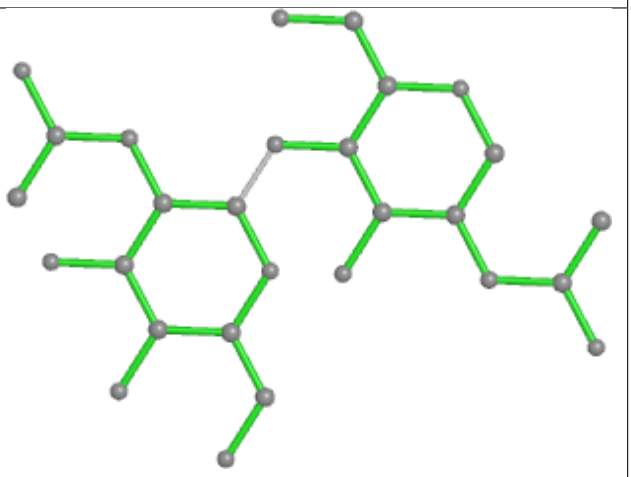
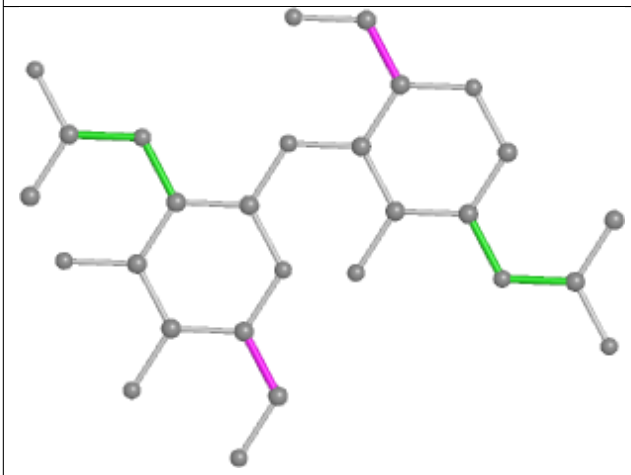
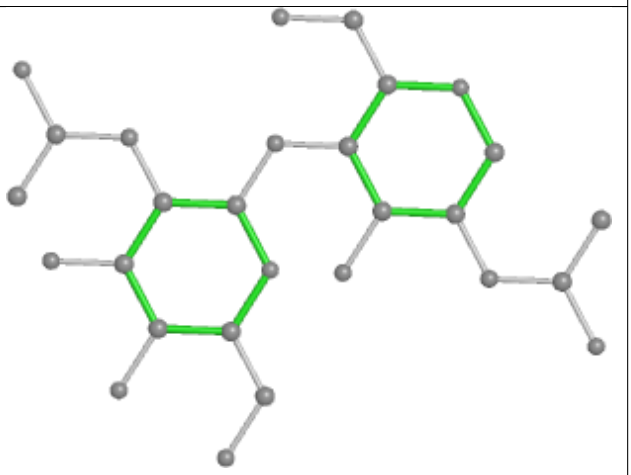
Bond angles



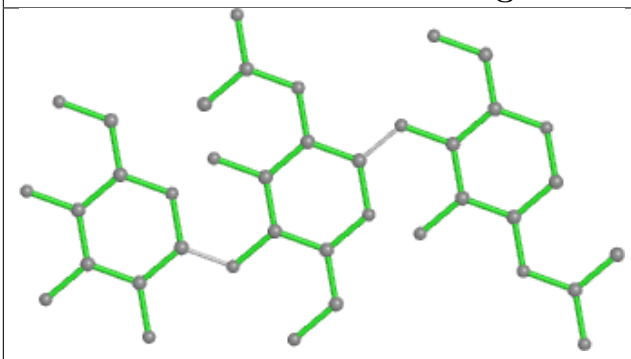
Torsions



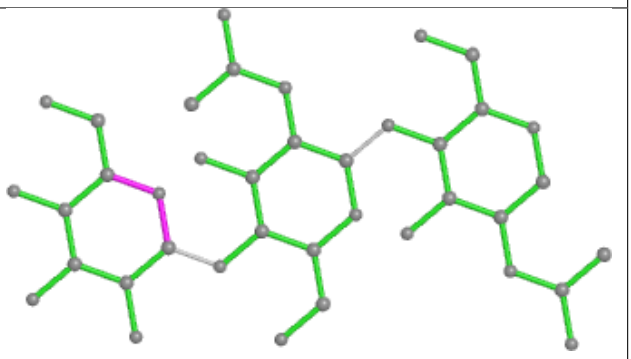
Rings

Oligosaccharide Chain d	
 Bond lengths	 Bond angles
 Torsions	 Rings

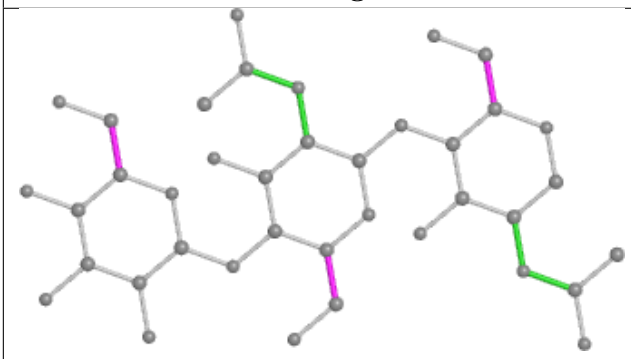
Oligosaccharide Chain F



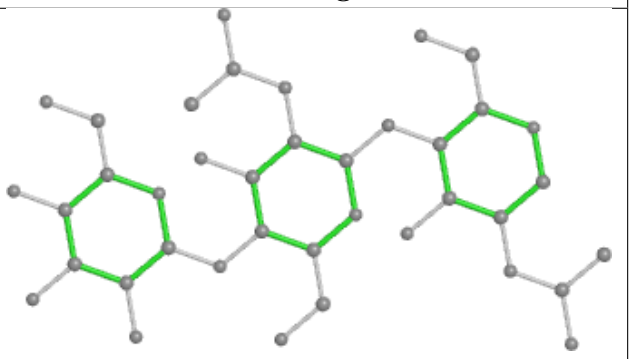
Bond lengths



Bond angles

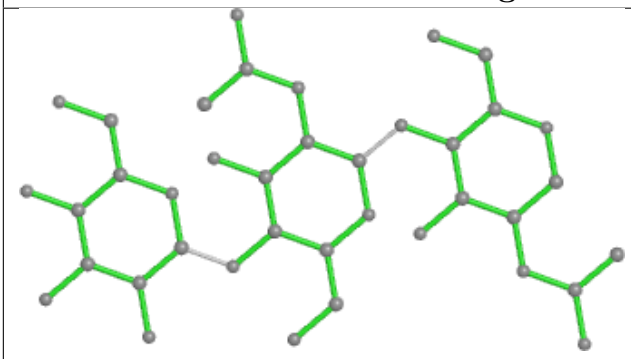


Torsions

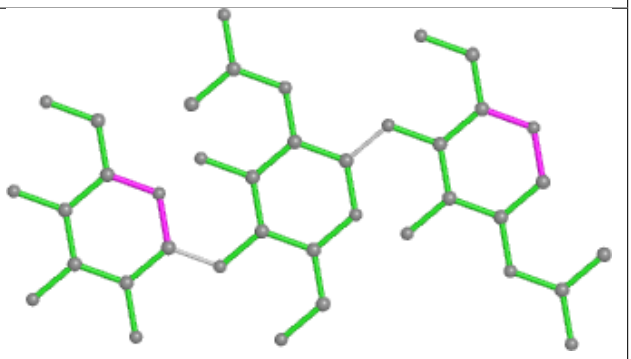


Rings

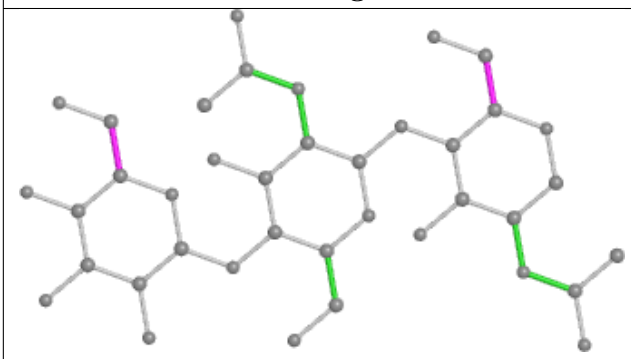
Oligosaccharide Chain J



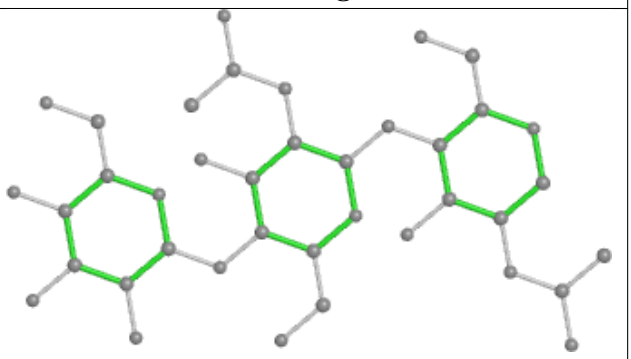
Bond lengths



Bond angles

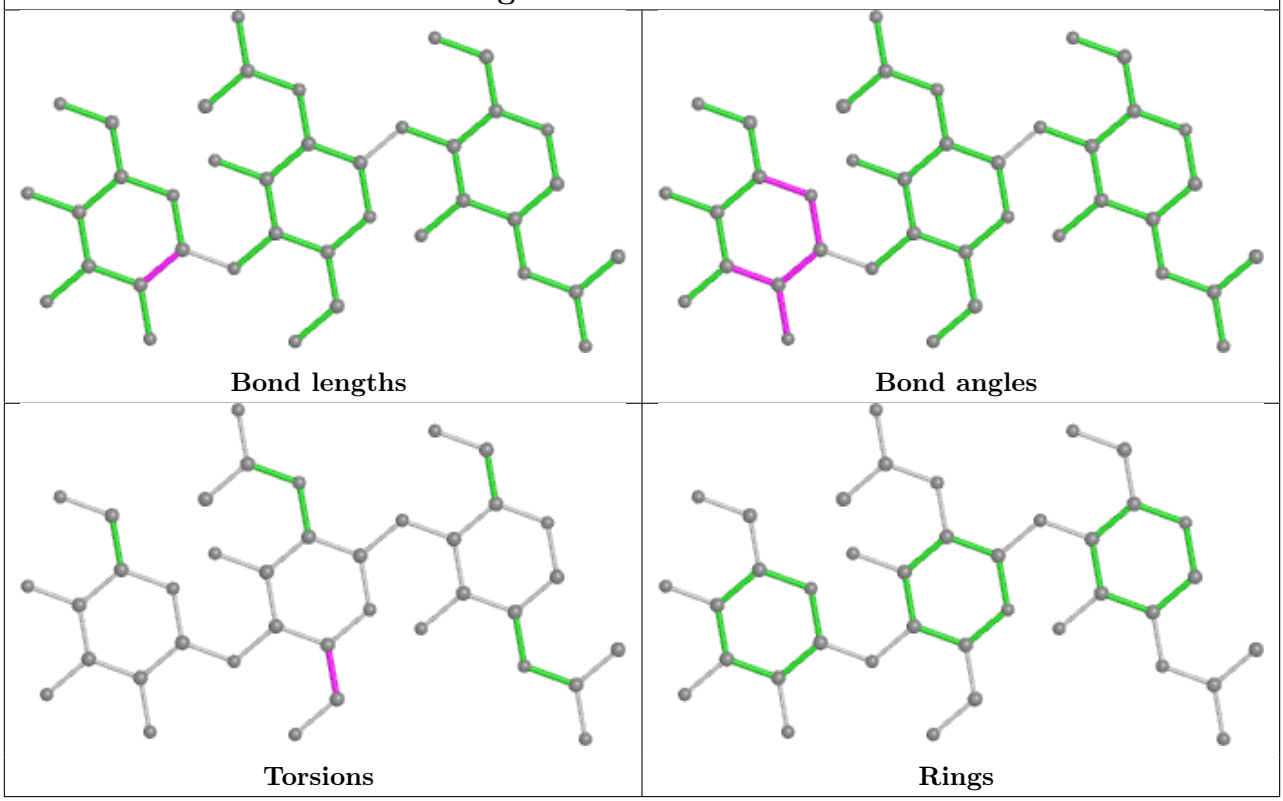


Torsions

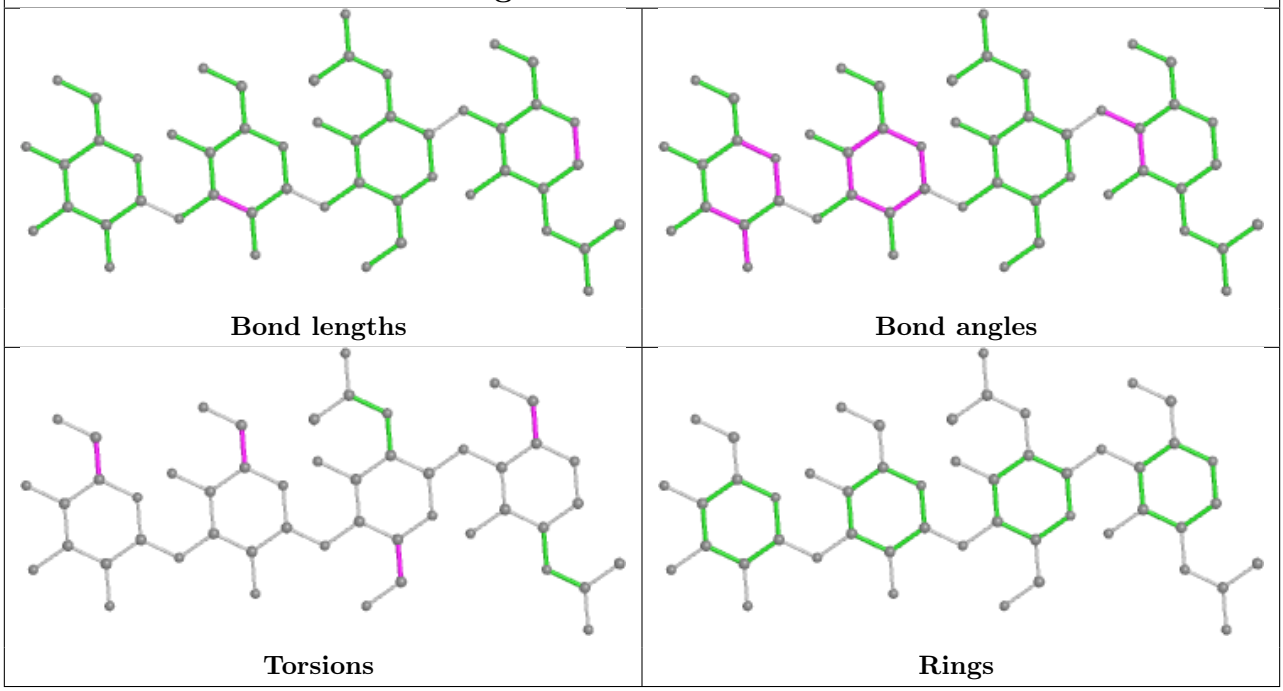


Rings

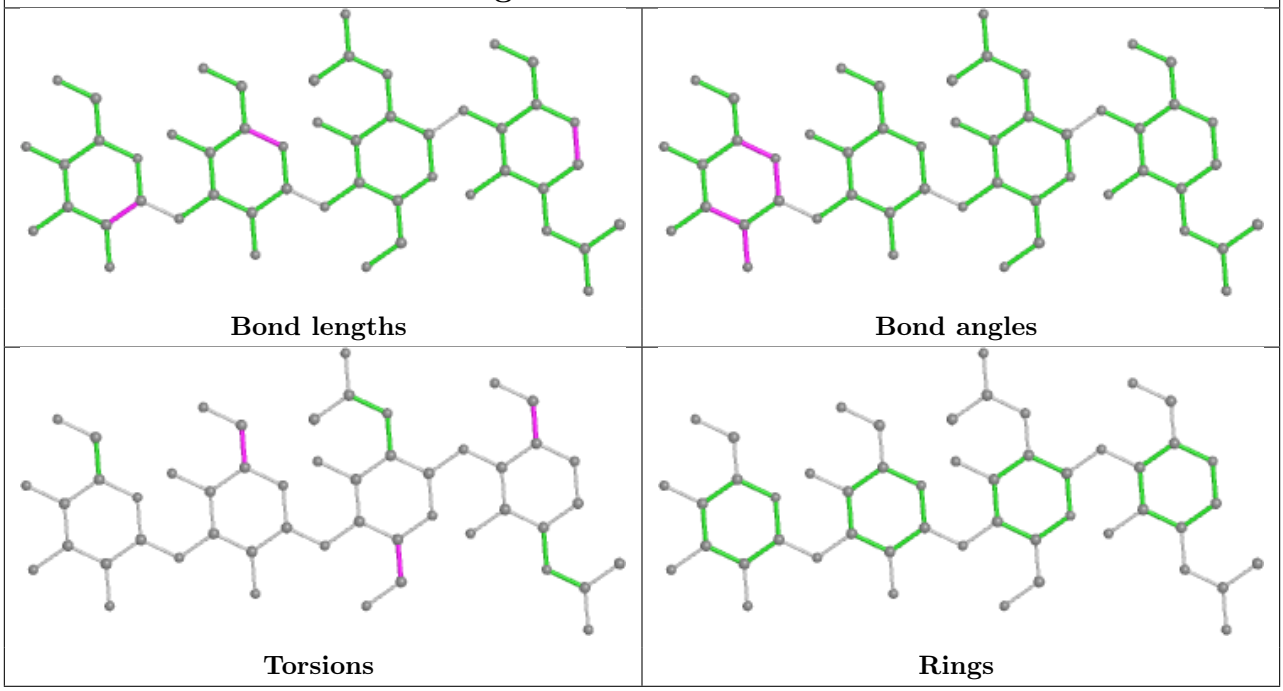
Oligosaccharide Chain Z



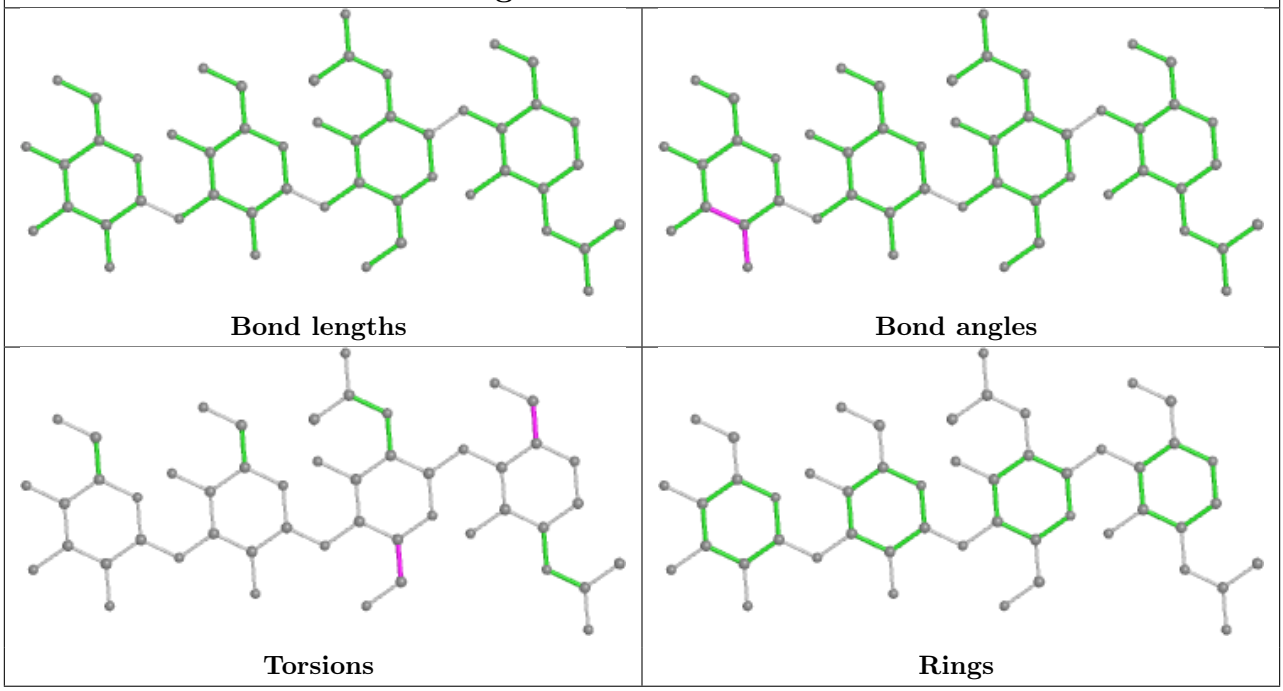
Oligosaccharide Chain G



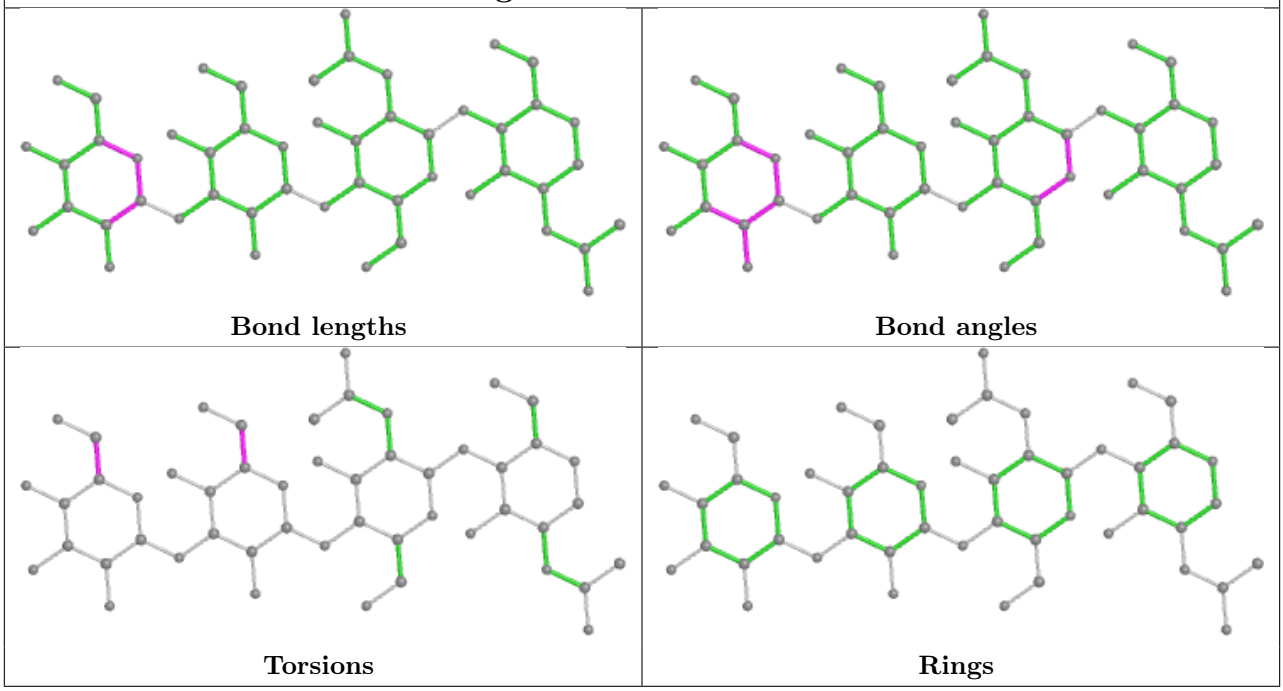
Oligosaccharide Chain N



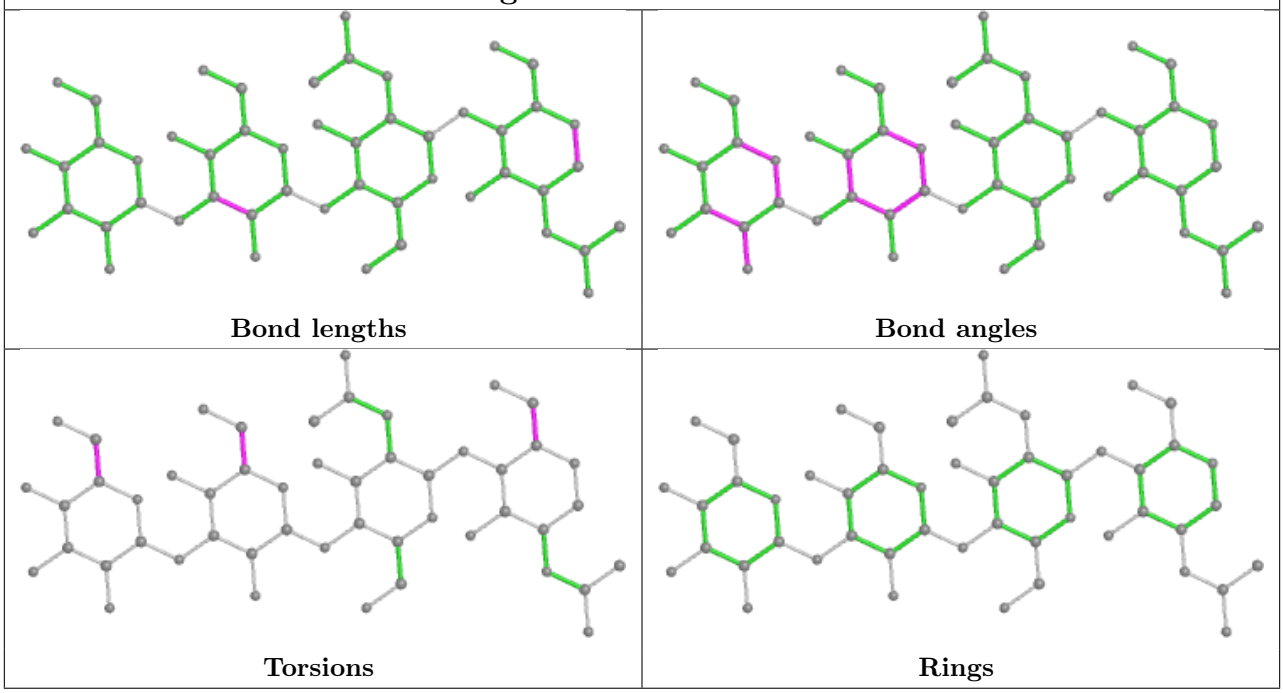
Oligosaccharide Chain O



Oligosaccharide Chain X



Oligosaccharide Chain a



5.6 Ligand geometry (i)

7 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	$\#\|Z\| > 2$
5	NAG	C	1307	1	14,14,15	0.45	0	17,19,21	0.48	0
5	NAG	A	1301	1	14,14,15	0.27	0	17,19,21	0.47	0
5	NAG	B	1308	1	14,14,15	0.47	0	17,19,21	0.49	0
5	NAG	C	1312	1	14,14,15	0.32	0	17,19,21	0.46	0
5	NAG	B	1309	1	14,14,15	0.24	0	17,19,21	0.48	0
5	NAG	B	1317	1	14,14,15	0.36	0	17,19,21	0.44	0
5	NAG	B	1314	1	14,14,15	0.20	0	17,19,21	0.56	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	C	1307	1	-	0/6/23/26	0/1/1/1
5	NAG	A	1301	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1308	1	-	2/6/23/26	0/1/1/1
5	NAG	C	1312	1	-	2/6/23/26	0/1/1/1
5	NAG	B	1309	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1317	1	-	0/6/23/26	0/1/1/1
5	NAG	B	1314	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	1301	NAG	O5-C5-C6-O6
5	C	1312	NAG	O5-C5-C6-O6
5	A	1301	NAG	C4-C5-C6-O6
5	C	1312	NAG	C4-C5-C6-O6
5	B	1308	NAG	O5-C5-C6-O6
5	B	1308	NAG	C4-C5-C6-O6

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.

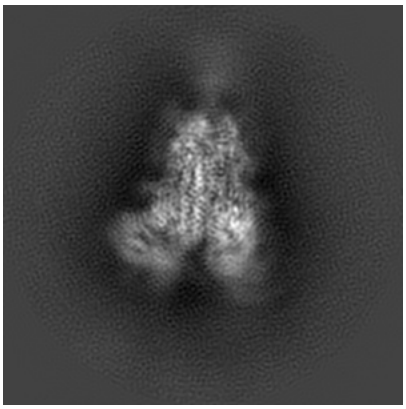
6 Map visualisation (i)

This section contains visualisations of the EMDB entry EMD-7575. These allow visual inspection of the internal detail of the map and identification of artifacts.

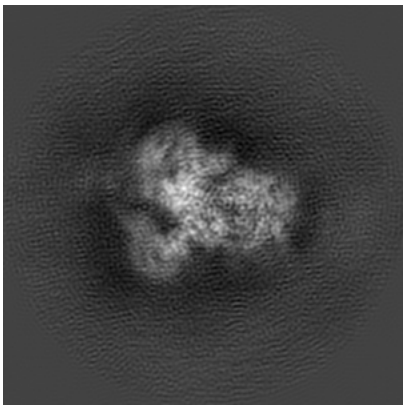
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections (i)

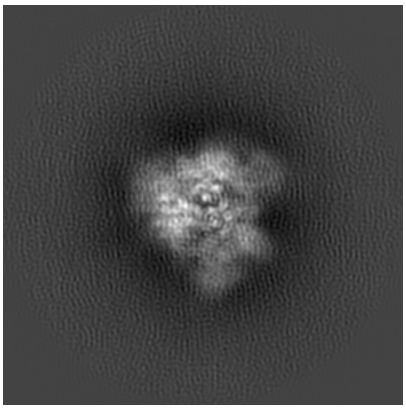
6.1.1 Primary map



X

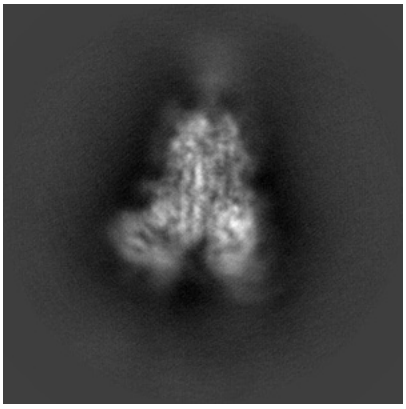


Y

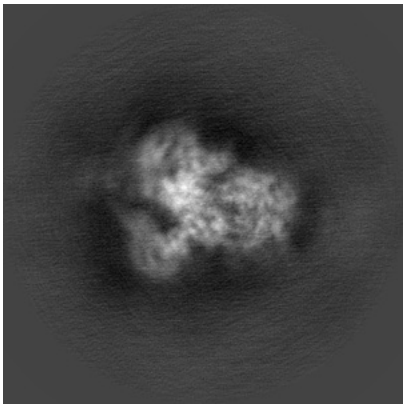


Z

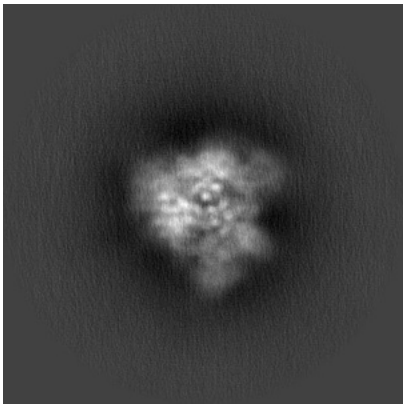
6.1.2 Raw map



X



Y

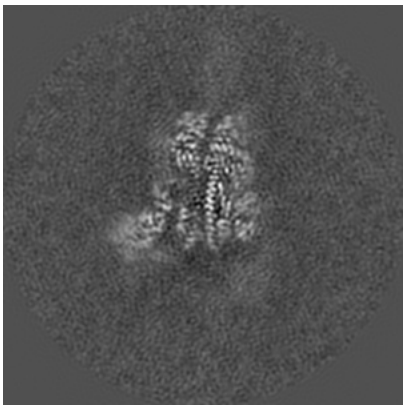


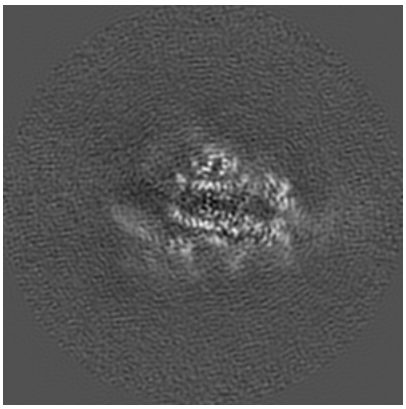
Z

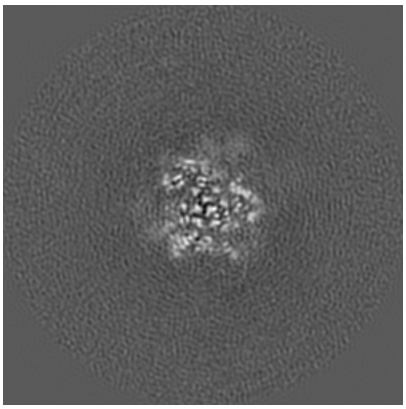
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

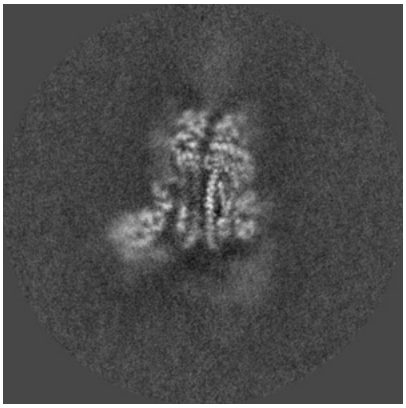
6.2.1 Primary map

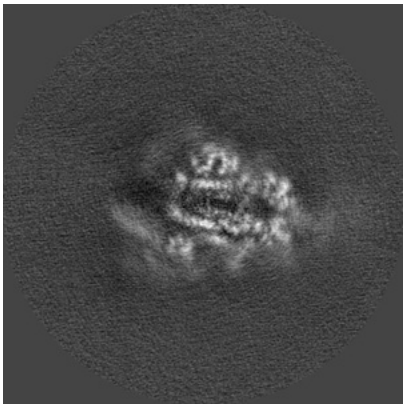

X Index: 160

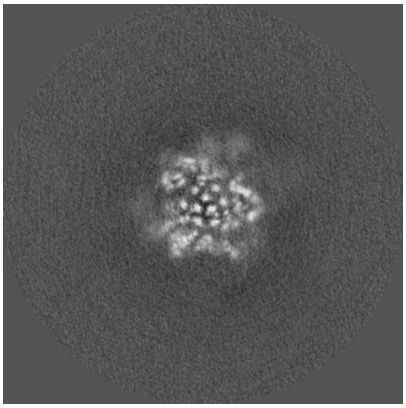

Y Index: 160


Z Index: 160

6.2.2 Raw map


X Index: 160

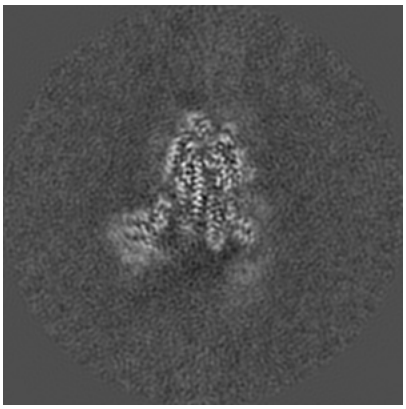

Y Index: 160

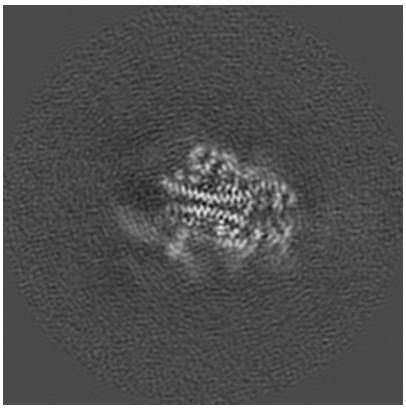

Z Index: 160

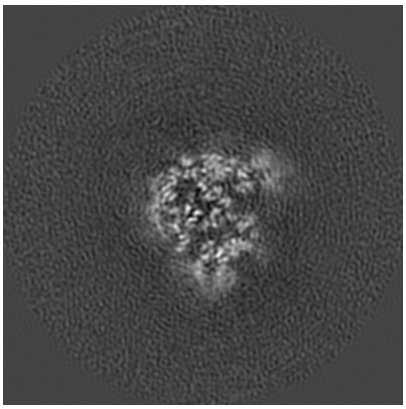
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

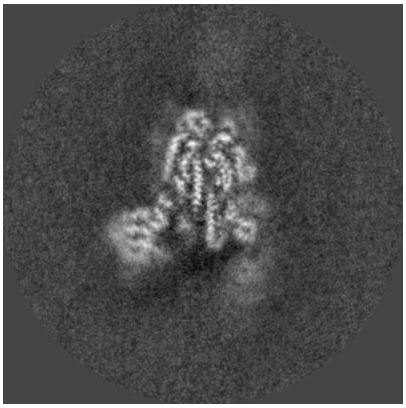
6.3.1 Primary map

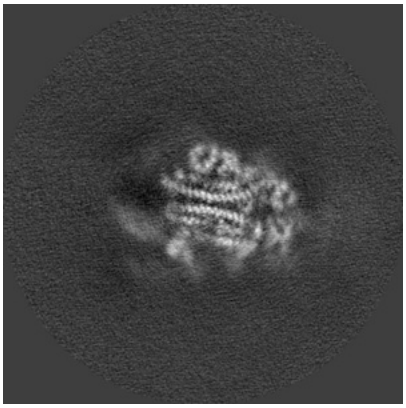

X Index: 165

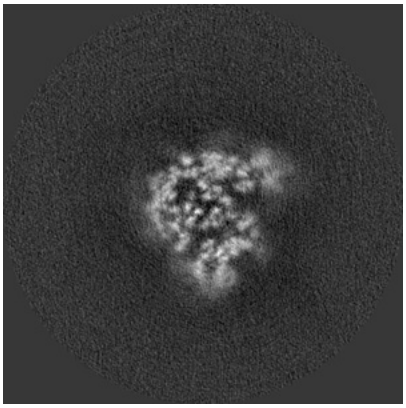

Y Index: 155


Z Index: 145

6.3.2 Raw map


X Index: 164


Y Index: 156


Z Index: 145

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 0.0328. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.4.2 Raw map

X Y Z

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

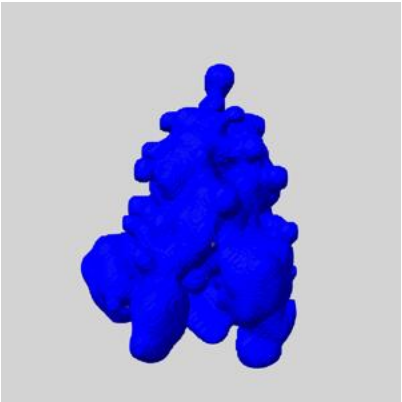
6.5 Mask visualisation (i)

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

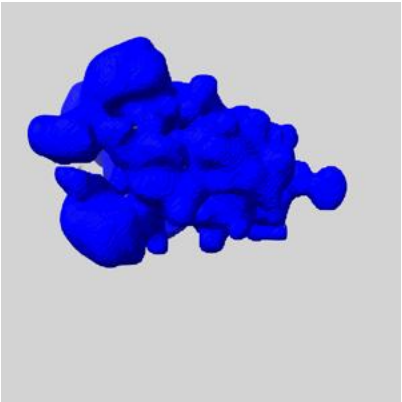
A mask typically either:

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

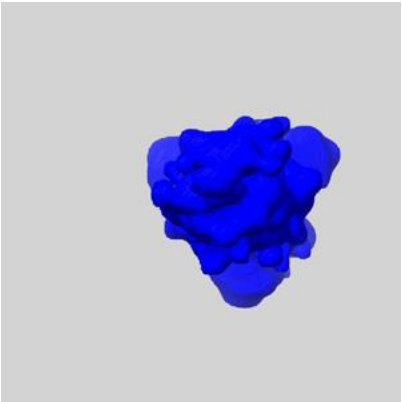
6.5.1 emd_7575_msk_1.map (i)



X



Y

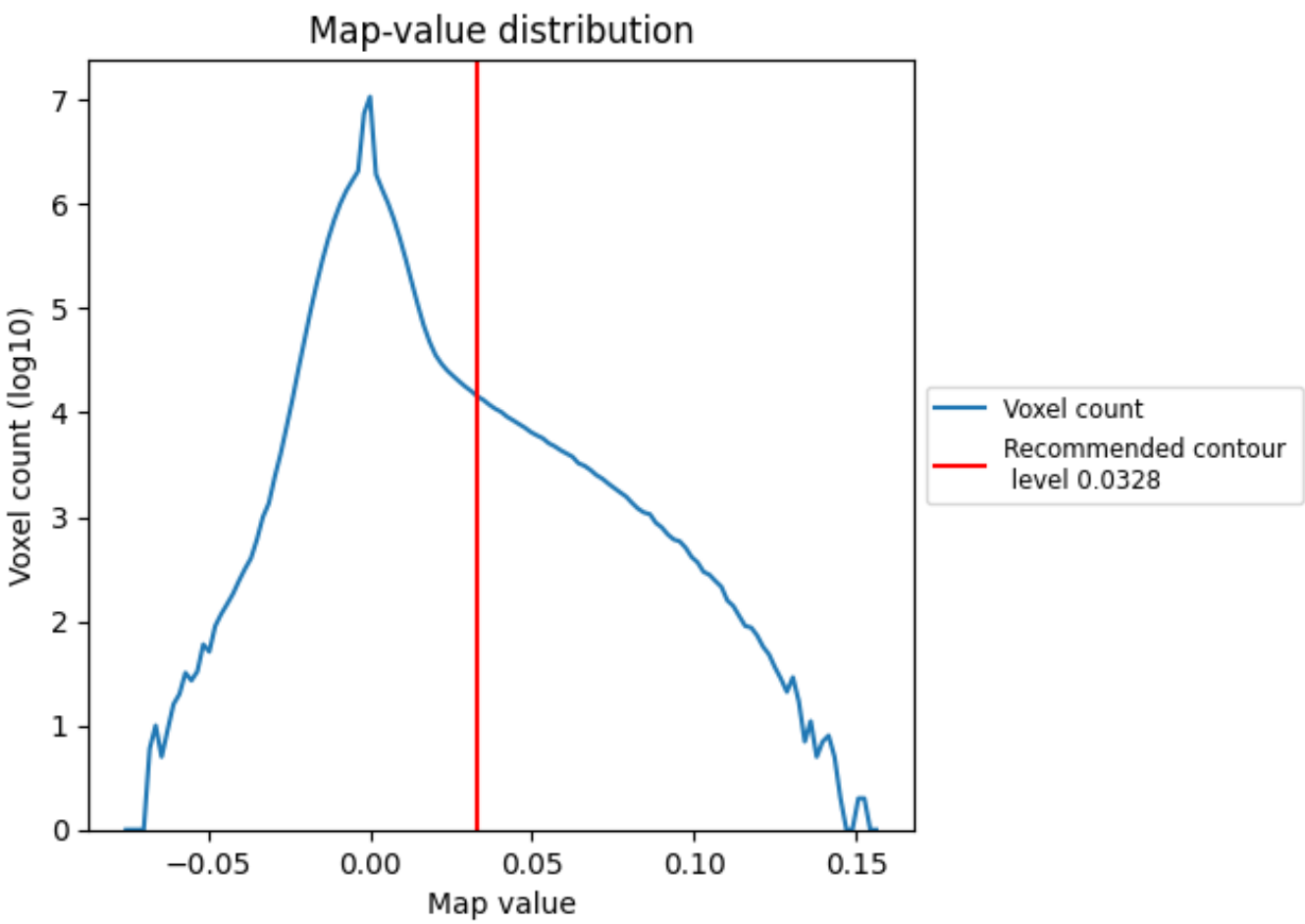


Z

7 Map analysis (i)

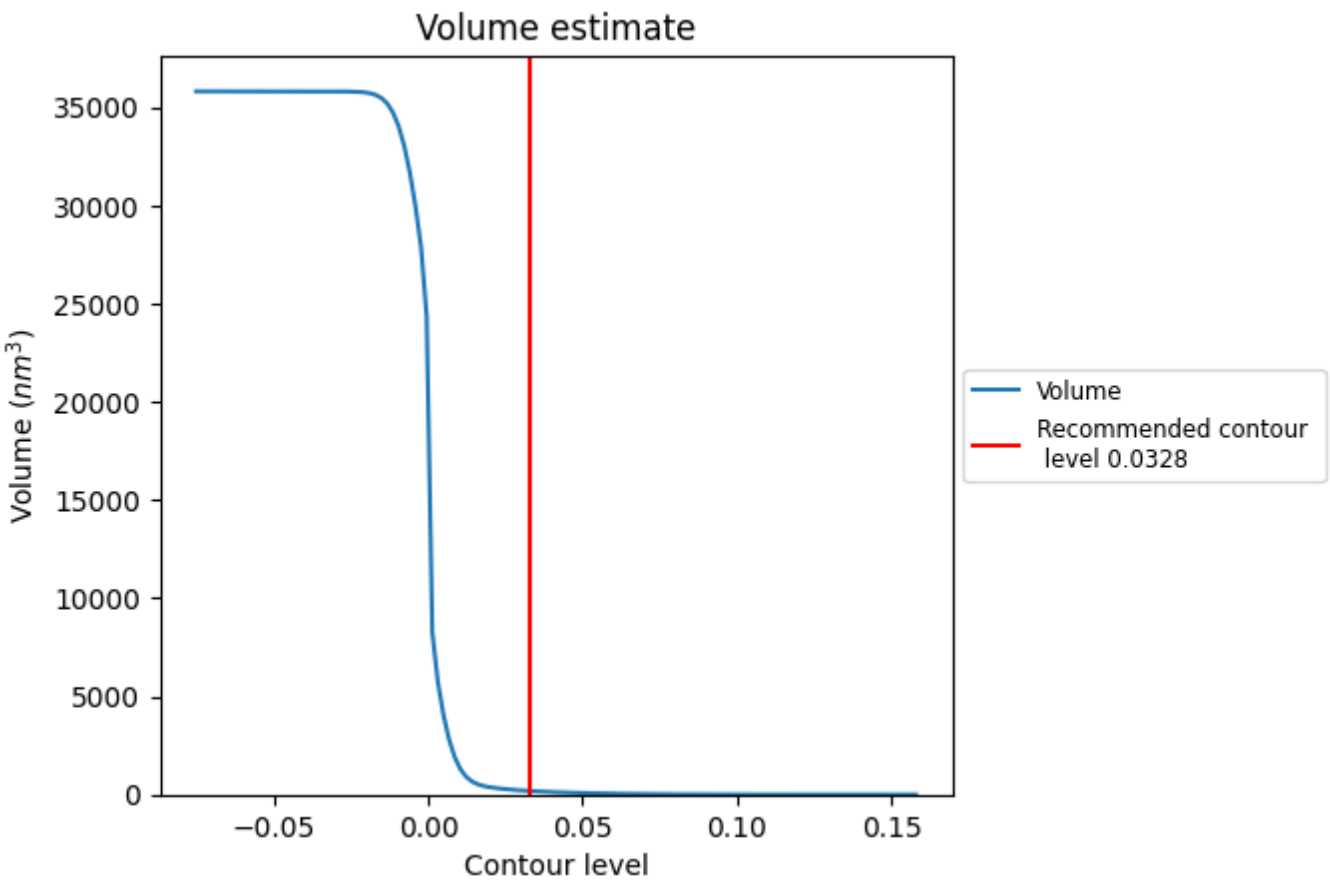
This section contains the results of statistical analysis of the map.

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.

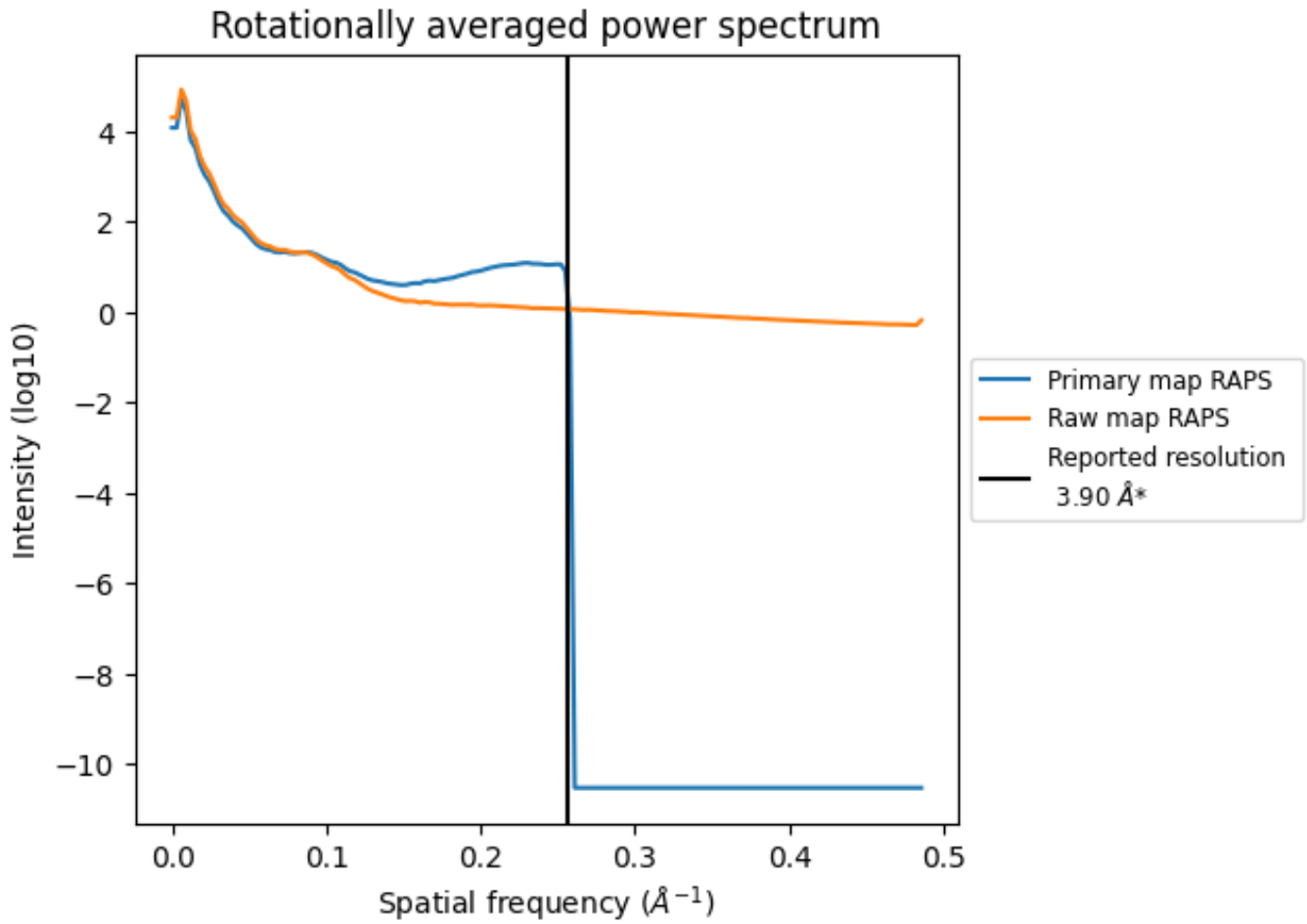
7.2 Volume estimate (i)



The volume at the recommended contour level is 185 nm^3; this corresponds to an approximate mass of 167 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 (i)

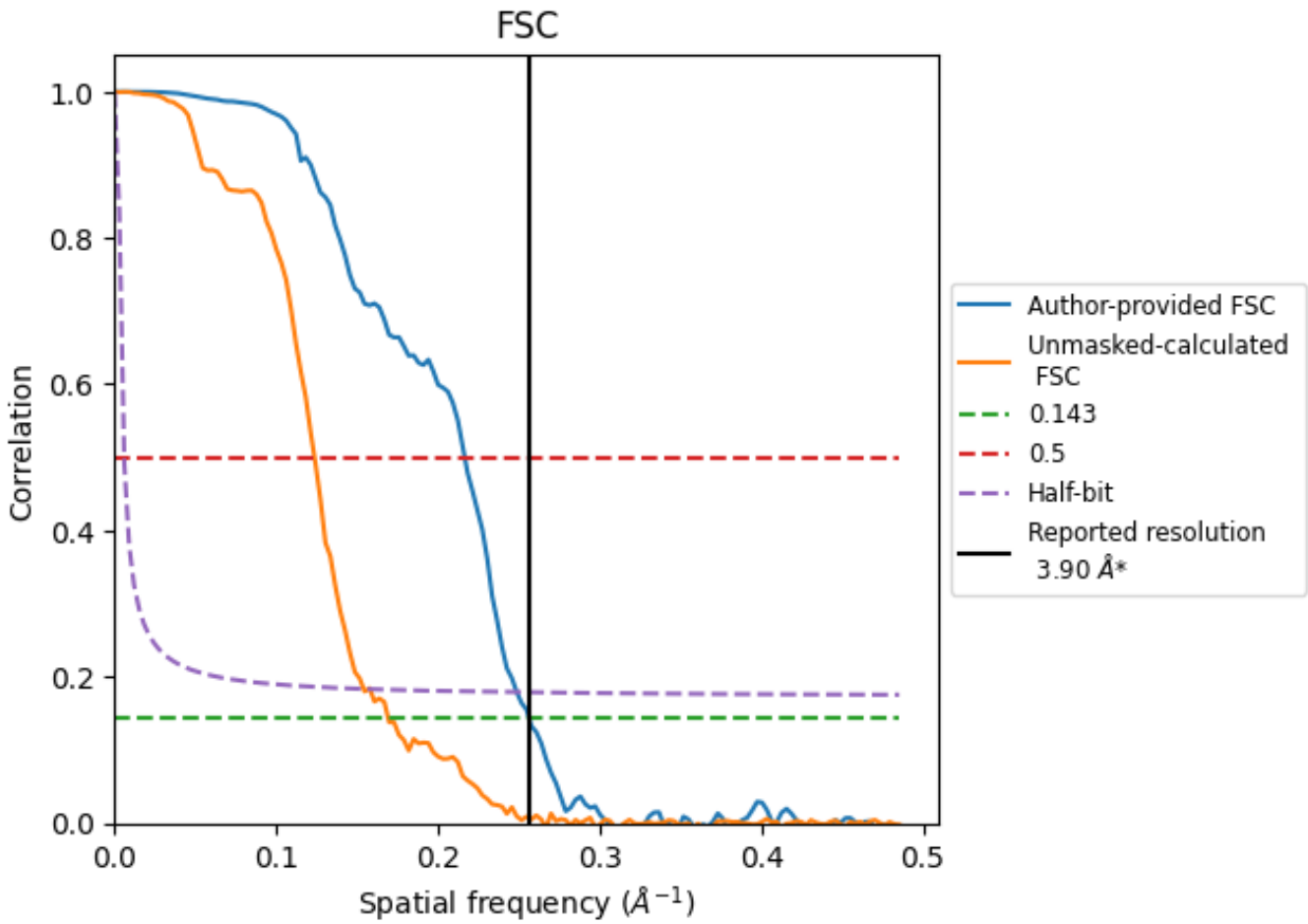


*Reported resolution corresponds to spatial frequency of 0.256 Å^{-1}

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 0.256 Å^{-1}

8.2 Resolution estimates (i)

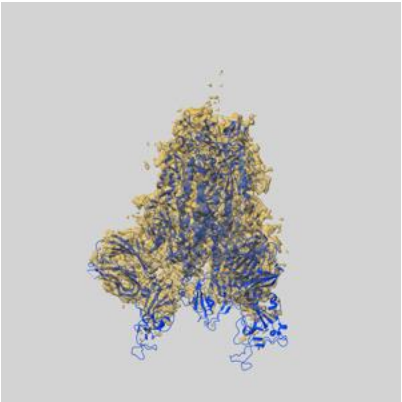
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.90	-	-
Author-provided FSC curve	3.90	4.62	4.02
Unmasked-calculated*	5.91	8.08	6.49

*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 5.91 differs from the reported value 3.9 by more than 10 %

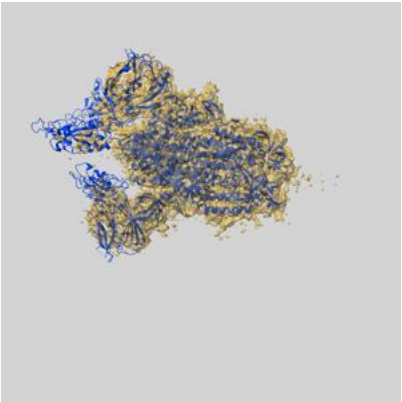
9 Map-model fit (i)

This section contains information regarding the fit between EMDB map EMD-7575 and PDB model 6CRX. Per-residue inclusion information can be found in section 3 on page 9.

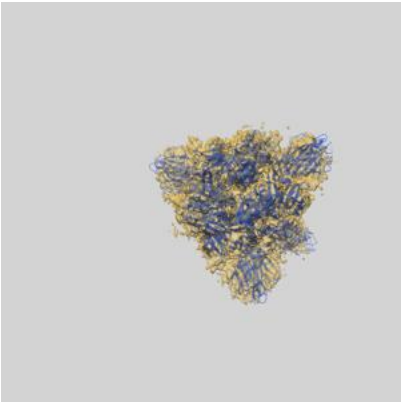
9.1 Map-model overlay (i)



X



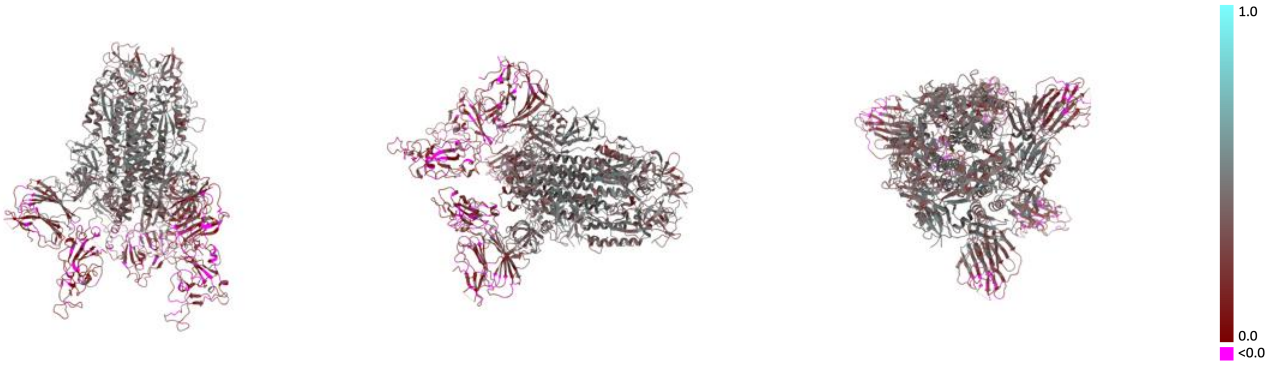
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.0328 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.

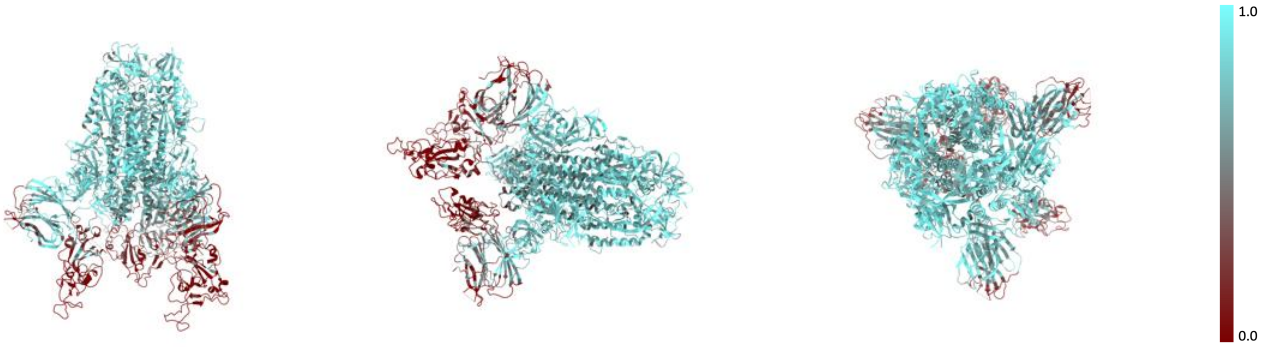
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

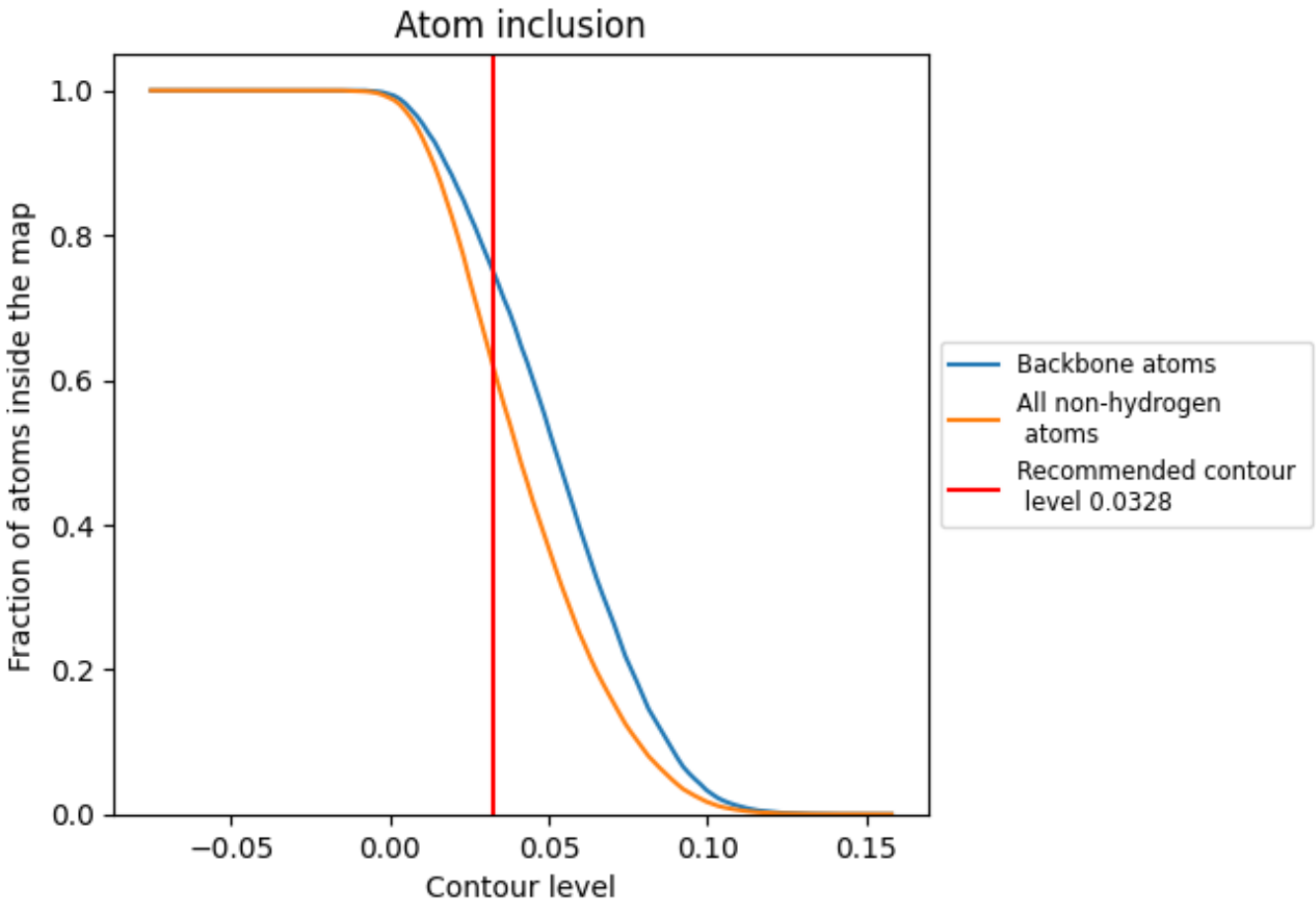
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.0328).

9.4 Atom inclusion (i)



At the recommended contour level, 75% of all backbone atoms, 61% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (0.0328) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6142	0.3240
A	0.6047	0.3200
B	0.6137	0.3230
C	0.6565	0.3300
D	0.4286	0.3220
E	0.1071	0.3380
F	0.2564	0.2650
G	0.4400	0.3120
H	0.5000	0.4390
I	0.2143	0.3140
J	0.3590	0.3190
K	0.2143	0.2510
L	0.0357	0.1380
M	0.1071	0.1480
N	0.4800	0.3230
O	0.2800	0.3070
P	0.6071	0.3660
Q	0.1071	0.2180
R	0.3571	0.1670
S	0.4286	0.4250
T	0.3571	0.3390
U	0.5000	0.3790
V	0.2500	0.3010
W	0.0000	0.3340
X	0.3400	0.2900
Y	0.0714	0.3160
Z	0.4359	0.4180
a	0.4200	0.3160
b	0.1786	0.3370
c	0.5000	0.3500
d	0.2500	0.2870