wwPDB EM Validation Summary Report

Aug 22, 2026 – 01:13 PM EDT

PDB ID	:	9ZRX / pdb_00009zrx
EMDB ID	:	EMD-74654
Title	:	Cryo-EM structure of SHIV-elicited CE79-1571 Fab in complex with HIV Env trimer Q23-SCT27
Authors	:	Roark, R.S.; Shapiro, L.; Kwong, P.D.
Deposited on	:	2025-12-21
Resolution	:	3.30 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:

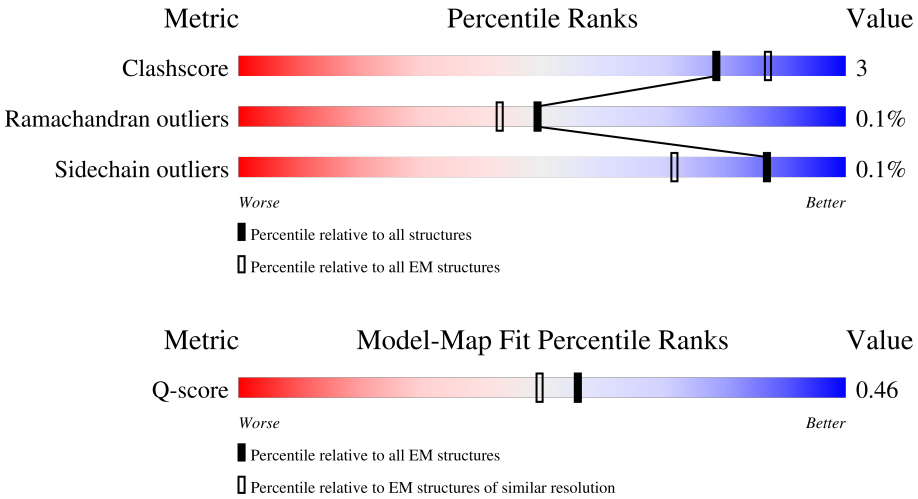
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	2022.3.0, CSD as543be (2022)
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

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 3.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	15087 (2.80 - 3.80)

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	248	13% 46% 9% 46%
2	B	221	13% 45% 6% 48%
3	b	503	82% 7% 12%
3	d	503	86% 12%

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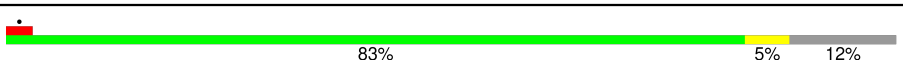
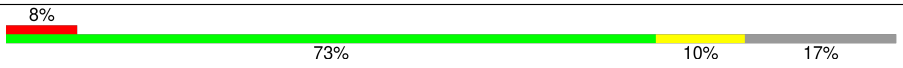
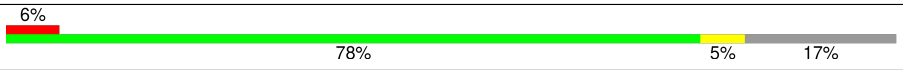
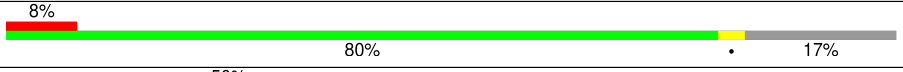
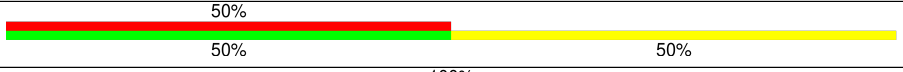
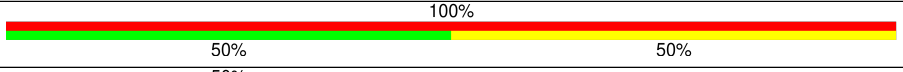
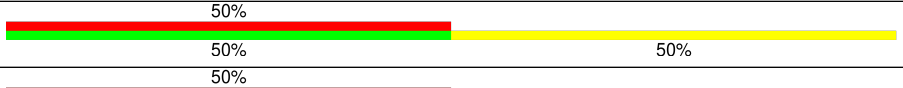
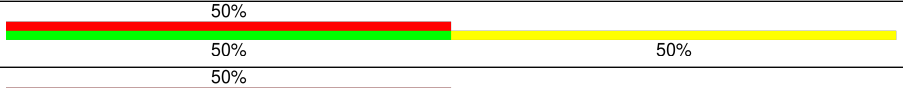
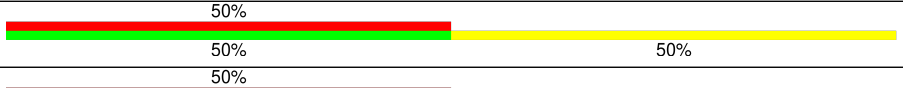
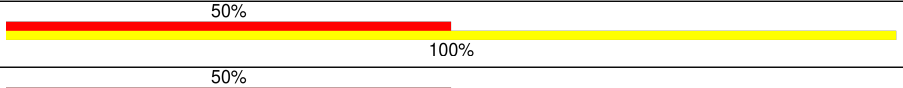
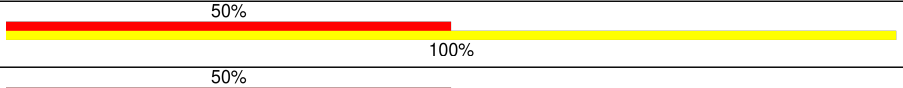
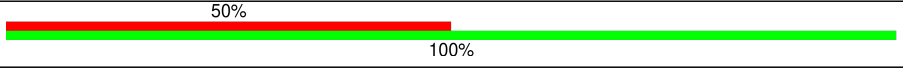
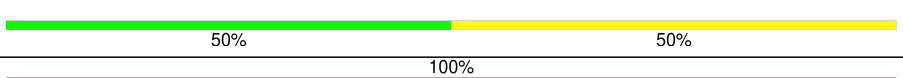
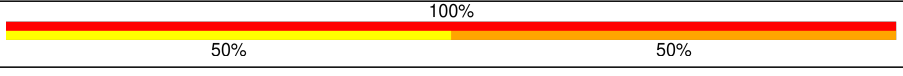
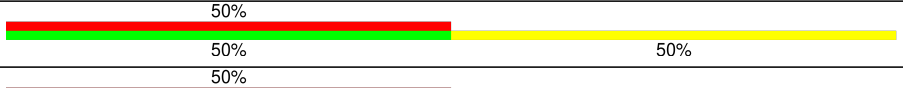
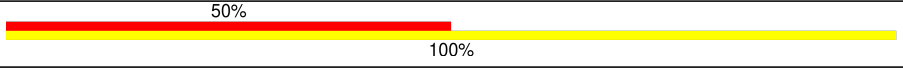
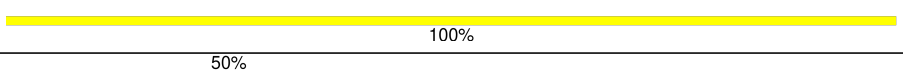
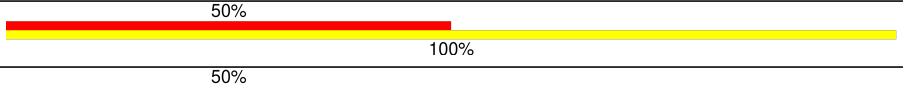
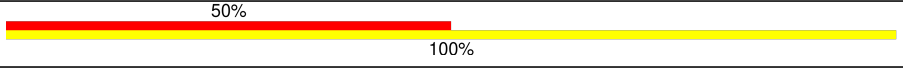
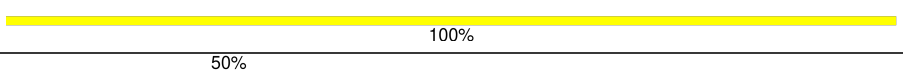
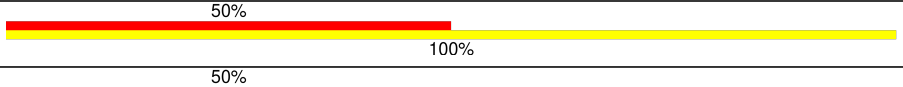
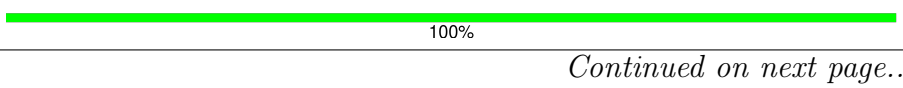
Mol	Chain	Length	Quality of chain
3	f	503	
4	c	153	
4	e	153	
4	g	153	
5	C	2	
5	F	2	
5	G	2	
5	H	2	
5	I	2	
5	K	2	
5	L	2	
5	M	2	
5	N	2	
5	O	2	
5	R	2	
5	T	2	
5	U	2	
5	V	2	
5	W	2	
5	X	2	
5	Y	2	
5	Z	2	
5	a	2	
5	h	2	
5	j	2	

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Mol	Chain	Length	Quality of chain
5	m	2	50% 100%
5	o	2	50% 50%
5	p	2	50% 50%
5	q	2	100%
5	r	2	50% 50%
5	s	2	50% 100%
5	t	2	50% 100%
5	u	2	50% 50% 50%
6	D	3	33% 67% 33%
6	J	3	67% 100%
6	P	3	33% 33% 67%
6	k	3	33% 67% 33%
7	E	5	40% 100%
7	Q	5	40% 20% 80%
7	l	5	40% 100%
7	n	5	20% 20% 60% 20%
8	S	4	75% 25%
9	i	3	100%

2 Entry composition (i)

There are 10 unique types of molecules in this entry. The entry contains 17291 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 CE79-1571 heavy chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	135	Total	C	N	O	S	0	0
			1036	646	171	216	3		

- Molecule 2 is a protein called CE79-1571 light chain.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	114	Total	C	N	O	S	0	0
			873	547	157	166	3		

- Molecule 3 is a protein called Envelope gp120.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	b	445	Total	C	N	O	S	0	0
			3525	2218	621	657	29		
3	d	445	Total	C	N	O	S	0	0
			3525	2218	621	657	29		
3	f	445	Total	C	N	O	S	0	0
			3525	2218	621	657	29		

There are 135 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
b	-4	MET	-	initiating methionine	UNP O55774
b	-3	ASP	-	expression tag	UNP O55774
b	-2	ALA	-	expression tag	UNP O55774
b	-1	MET	-	expression tag	UNP O55774
b	0	LYS	-	expression tag	UNP O55774
b	1	ARG	-	expression tag	UNP O55774
b	2	GLY	-	expression tag	UNP O55774
b	3	LEU	-	expression tag	UNP O55774
b	4	CYS	-	expression tag	UNP O55774
b	5	CYS	-	expression tag	UNP O55774
b	6	VAL	-	expression tag	UNP O55774

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Chain	Residue	Modelled	Actual	Comment	Reference
b	7	LEU	-	expression tag	UNP O55774
b	8	LEU	-	expression tag	UNP O55774
b	9	LEU	-	expression tag	UNP O55774
b	10	CYS	-	expression tag	UNP O55774
b	11	GLY	-	expression tag	UNP O55774
b	12	ALA	-	expression tag	UNP O55774
b	13	VAL	-	expression tag	UNP O55774
b	14	PHE	-	expression tag	UNP O55774
b	15	VAL	-	expression tag	UNP O55774
b	16	SER	-	expression tag	UNP O55774
b	17	PRO	-	expression tag	UNP O55774
b	18	SER	-	expression tag	UNP O55774
b	19	GLN	-	expression tag	UNP O55774
b	20	GLU	-	expression tag	UNP O55774
b	21	ILE	-	expression tag	UNP O55774
b	22	HIS	-	expression tag	UNP O55774
b	23	ALA	-	expression tag	UNP O55774
b	24	ARG	-	expression tag	UNP O55774
b	25	PHE	-	expression tag	UNP O55774
b	26	ARG	-	expression tag	UNP O55774
b	27	ARG	-	expression tag	UNP O55774
b	28	GLY	-	expression tag	UNP O55774
b	29	ALA	-	expression tag	UNP O55774
b	30	ARG	-	expression tag	UNP O55774
b	31	ALA	VAL	conflict	UNP O55774
b	49	GLU	ASP	conflict	UNP O55774
b	153	GLU	GLY	conflict	UNP O55774
b	201	CYS	ILE	conflict	UNP O55774
b	219	ALA	THR	conflict	UNP O55774
b	302	TYR	ASN	conflict	UNP O55774
b	320	MET	THR	conflict	UNP O55774
b	334	SER	THR	conflict	UNP O55774
b	433	CYS	ALA	conflict	UNP O55774
b	501	CYS	ALA	conflict	UNP O55774
d	-4	MET	-	initiating methionine	UNP O55774
d	-3	ASP	-	expression tag	UNP O55774
d	-2	ALA	-	expression tag	UNP O55774
d	-1	MET	-	expression tag	UNP O55774
d	0	LYS	-	expression tag	UNP O55774
d	1	ARG	-	expression tag	UNP O55774
d	2	GLY	-	expression tag	UNP O55774
d	3	LEU	-	expression tag	UNP O55774

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Chain	Residue	Modelled	Actual	Comment	Reference
d	4	CYS	-	expression tag	UNP O55774
d	5	CYS	-	expression tag	UNP O55774
d	6	VAL	-	expression tag	UNP O55774
d	7	LEU	-	expression tag	UNP O55774
d	8	LEU	-	expression tag	UNP O55774
d	9	LEU	-	expression tag	UNP O55774
d	10	CYS	-	expression tag	UNP O55774
d	11	GLY	-	expression tag	UNP O55774
d	12	ALA	-	expression tag	UNP O55774
d	13	VAL	-	expression tag	UNP O55774
d	14	PHE	-	expression tag	UNP O55774
d	15	VAL	-	expression tag	UNP O55774
d	16	SER	-	expression tag	UNP O55774
d	17	PRO	-	expression tag	UNP O55774
d	18	SER	-	expression tag	UNP O55774
d	19	GLN	-	expression tag	UNP O55774
d	20	GLU	-	expression tag	UNP O55774
d	21	ILE	-	expression tag	UNP O55774
d	22	HIS	-	expression tag	UNP O55774
d	23	ALA	-	expression tag	UNP O55774
d	24	ARG	-	expression tag	UNP O55774
d	25	PHE	-	expression tag	UNP O55774
d	26	ARG	-	expression tag	UNP O55774
d	27	ARG	-	expression tag	UNP O55774
d	28	GLY	-	expression tag	UNP O55774
d	29	ALA	-	expression tag	UNP O55774
d	30	ARG	-	expression tag	UNP O55774
d	31	ALA	VAL	conflict	UNP O55774
d	49	GLU	ASP	conflict	UNP O55774
d	153	GLU	GLY	conflict	UNP O55774
d	201	CYS	ILE	conflict	UNP O55774
d	219	ALA	THR	conflict	UNP O55774
d	302	TYR	ASN	conflict	UNP O55774
d	320	MET	THR	conflict	UNP O55774
d	334	SER	THR	conflict	UNP O55774
d	433	CYS	ALA	conflict	UNP O55774
d	501	CYS	ALA	conflict	UNP O55774
f	-4	MET	-	initiating methionine	UNP O55774
f	-3	ASP	-	expression tag	UNP O55774
f	-2	ALA	-	expression tag	UNP O55774
f	-1	MET	-	expression tag	UNP O55774
f	0	LYS	-	expression tag	UNP O55774

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Chain	Residue	Modelled	Actual	Comment	Reference
f	1	ARG	-	expression tag	UNP O55774
f	2	GLY	-	expression tag	UNP O55774
f	3	LEU	-	expression tag	UNP O55774
f	4	CYS	-	expression tag	UNP O55774
f	5	CYS	-	expression tag	UNP O55774
f	6	VAL	-	expression tag	UNP O55774
f	7	LEU	-	expression tag	UNP O55774
f	8	LEU	-	expression tag	UNP O55774
f	9	LEU	-	expression tag	UNP O55774
f	10	CYS	-	expression tag	UNP O55774
f	11	GLY	-	expression tag	UNP O55774
f	12	ALA	-	expression tag	UNP O55774
f	13	VAL	-	expression tag	UNP O55774
f	14	PHE	-	expression tag	UNP O55774
f	15	VAL	-	expression tag	UNP O55774
f	16	SER	-	expression tag	UNP O55774
f	17	PRO	-	expression tag	UNP O55774
f	18	SER	-	expression tag	UNP O55774
f	19	GLN	-	expression tag	UNP O55774
f	20	GLU	-	expression tag	UNP O55774
f	21	ILE	-	expression tag	UNP O55774
f	22	HIS	-	expression tag	UNP O55774
f	23	ALA	-	expression tag	UNP O55774
f	24	ARG	-	expression tag	UNP O55774
f	25	PHE	-	expression tag	UNP O55774
f	26	ARG	-	expression tag	UNP O55774
f	27	ARG	-	expression tag	UNP O55774
f	28	GLY	-	expression tag	UNP O55774
f	29	ALA	-	expression tag	UNP O55774
f	30	ARG	-	expression tag	UNP O55774
f	31	ALA	VAL	conflict	UNP O55774
f	49	GLU	ASP	conflict	UNP O55774
f	153	GLU	GLY	conflict	UNP O55774
f	201	CYS	ILE	conflict	UNP O55774
f	219	ALA	THR	conflict	UNP O55774
f	302	TYR	ASN	conflict	UNP O55774
f	320	MET	THR	conflict	UNP O55774
f	334	SER	THR	conflict	UNP O55774
f	433	CYS	ALA	conflict	UNP O55774
f	501	CYS	ALA	conflict	UNP O55774

- Molecule 4 is a protein called Envelope gp41 ectodomain.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	c	127	Total	C	N	O	S	0	0
			1026	650	184	187	5		
4	e	127	Total	C	N	O	S	0	0
			1026	650	184	187	5		
4	g	127	Total	C	N	O	S	0	0
			1026	650	184	187	5		

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
c	519	ARG	PHE	conflict	UNP O55774
c	520	ARG	LEU	conflict	UNP O55774
c	533	ALA	THR	conflict	UNP O55774
c	551	PRO	GLN	conflict	UNP O55774
c	559	PRO	ILE	conflict	UNP O55774
c	569	GLY	THR	conflict	UNP O55774
c	605	CYS	THR	conflict	UNP O55774
c	636	GLY	ASN	conflict	UNP O55774
c	662	ALA	GLU	conflict	UNP O55774
e	519	ARG	PHE	conflict	UNP O55774
e	520	ARG	LEU	conflict	UNP O55774
e	533	ALA	THR	conflict	UNP O55774
e	551	PRO	GLN	conflict	UNP O55774
e	559	PRO	ILE	conflict	UNP O55774
e	569	GLY	THR	conflict	UNP O55774
e	605	CYS	THR	conflict	UNP O55774
e	636	GLY	ASN	conflict	UNP O55774
e	662	ALA	GLU	conflict	UNP O55774
g	519	ARG	PHE	conflict	UNP O55774
g	520	ARG	LEU	conflict	UNP O55774
g	533	ALA	THR	conflict	UNP O55774
g	551	PRO	GLN	conflict	UNP O55774
g	559	PRO	ILE	conflict	UNP O55774
g	569	GLY	THR	conflict	UNP O55774
g	605	CYS	THR	conflict	UNP O55774
g	636	GLY	ASN	conflict	UNP O55774
g	662	ALA	GLU	conflict	UNP O55774

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



Mol	Chain	Residues	Atoms				AltConf	Trace
5	C	2	Total	C	N	O	0	0
			28	16	2	10		
5	F	2	Total	C	N	O	0	0
			28	16	2	10		
5	G	2	Total	C	N	O	0	0
			28	16	2	10		
5	H	2	Total	C	N	O	0	0
			28	16	2	10		
5	I	2	Total	C	N	O	0	0
			28	16	2	10		
5	K	2	Total	C	N	O	0	0
			28	16	2	10		
5	L	2	Total	C	N	O	0	0
			28	16	2	10		
5	M	2	Total	C	N	O	0	0
			28	16	2	10		
5	N	2	Total	C	N	O	0	0
			28	16	2	10		
5	O	2	Total	C	N	O	0	0
			28	16	2	10		
5	R	2	Total	C	N	O	0	0
			28	16	2	10		
5	T	2	Total	C	N	O	0	0
			28	16	2	10		
5	U	2	Total	C	N	O	0	0
			28	16	2	10		
5	V	2	Total	C	N	O	0	0
			28	16	2	10		
5	W	2	Total	C	N	O	0	0
			28	16	2	10		
5	X	2	Total	C	N	O	0	0
			28	16	2	10		
5	Y	2	Total	C	N	O	0	0
			28	16	2	10		
5	Z	2	Total	C	N	O	0	0
			28	16	2	10		
5	a	2	Total	C	N	O	0	0
			28	16	2	10		

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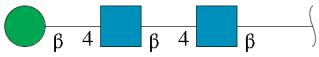
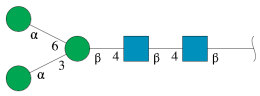
Mol	Chain	Residues	Atoms				AltConf	Trace
5	h	2	Total	C	N	O	0	0
			28	16	2	10		
5	j	2	Total	C	N	O	0	0
			28	16	2	10		
5	m	2	Total	C	N	O	0	0
			28	16	2	10		
5	o	2	Total	C	N	O	0	0
			28	16	2	10		
5	p	2	Total	C	N	O	0	0
			28	16	2	10		
5	q	2	Total	C	N	O	0	0
			28	16	2	10		
5	r	2	Total	C	N	O	0	0
			28	16	2	10		
5	s	2	Total	C	N	O	0	0
			28	16	2	10		
5	t	2	Total	C	N	O	0	0
			28	16	2	10		
5	u	2	Total	C	N	O	0	0
			28	16	2	10		

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



Mol	Chain	Residues	Atoms				AltConf	Trace
6	D	3	Total	C	N	O	0	0
			39	22	2	15		
6	J	3	Total	C	N	O	0	0
			39	22	2	15		
6	P	3	Total	C	N	O	0	0
			39	22	2	15		
6	k	3	Total	C	N	O	0	0
			39	22	2	15		

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



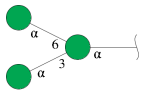
Mol	Chain	Residues	Atoms				AltConf	Trace
7	E	5	Total	C	N	O	0	0
			61	34	2	25		
7	Q	5	Total	C	N	O	0	0
			61	34	2	25		
7	l	5	Total	C	N	O	0	0
			61	34	2	25		
7	n	5	Total	C	N	O	0	0
			61	34	2	25		

- Molecule 8 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.



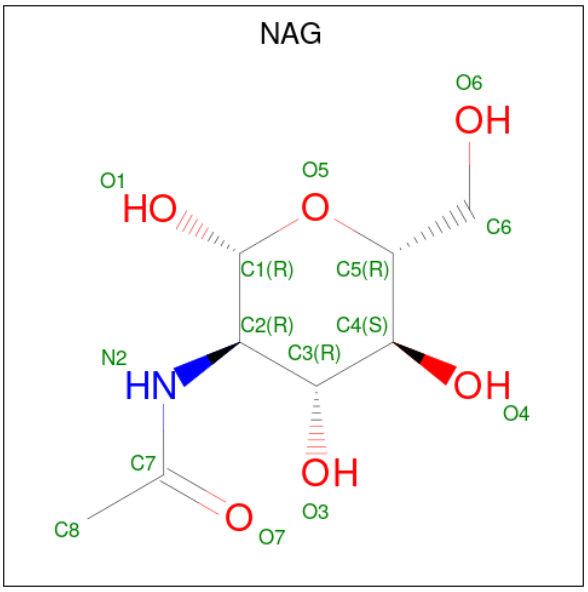
Mol	Chain	Residues	Atoms				AltConf	Trace
8	S	4	Total	C	N	O	0	0
			50	28	2	20		

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose.



Mol	Chain	Residues	Atoms			AltConf	Trace
9	i	3	Total	C	O	0	0
			33	18	15		

- Molecule 10 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				AltConf
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	b	1	Total	C	N	O	0
			14	8	1	5	
10	c	1	Total	C	N	O	0
			14	8	1	5	
10	c	1	Total	C	N	O	0
			14	8	1	5	
10	c	1	Total	C	N	O	0
			14	8	1	5	
10	d	1	Total	C	N	O	0
			14	8	1	5	
10	d	1	Total	C	N	O	0
			14	8	1	5	
10	d	1	Total	C	N	O	0
			14	8	1	5	

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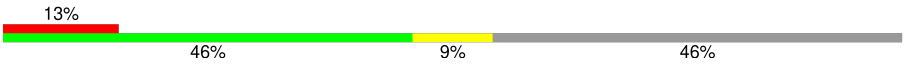
Mol	Chain	Residues	Atoms				AltConf
10	d	1	Total	C	N	O	0
			14	8	1	5	
10	d	1	Total	C	N	O	0
			14	8	1	5	
10	d	1	Total	C	N	O	0
			14	8	1	5	
10	e	1	Total	C	N	O	0
			14	8	1	5	
10	e	1	Total	C	N	O	0
			14	8	1	5	
10	e	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	f	1	Total	C	N	O	0
			14	8	1	5	
10	g	1	Total	C	N	O	0
			14	8	1	5	
10	g	1	Total	C	N	O	0
			14	8	1	5	
10	g	1	Total	C	N	O	0
			14	8	1	5	

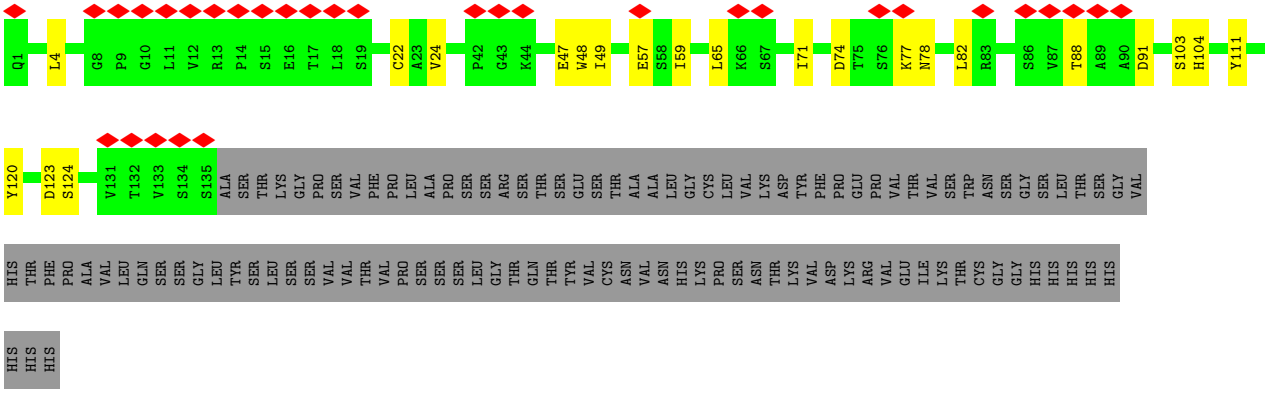
3 Residue-property plots (i)

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: CE79-1571 heavy chain

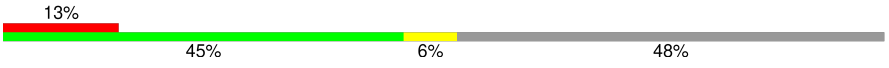
Chain A:

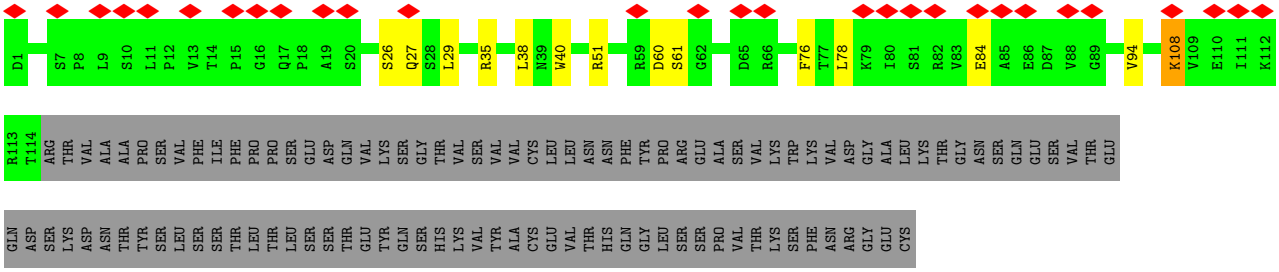




• Molecule 2: CE79-1571 light chain

Chain B:

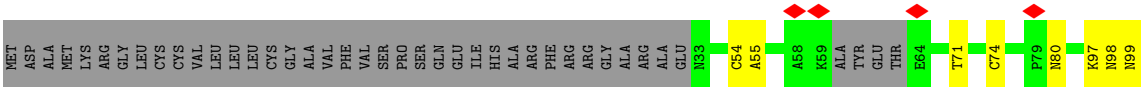


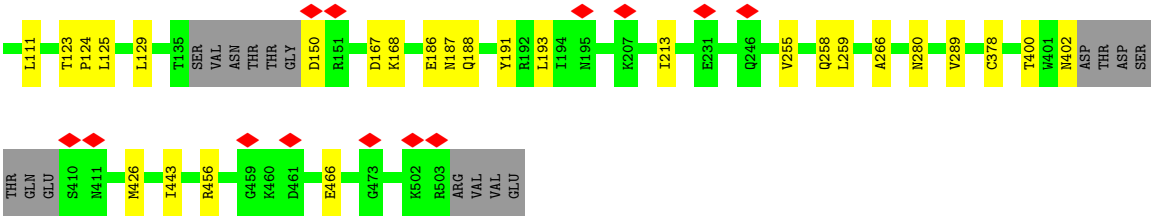


• Molecule 3: Envelope gp120

Chain b:



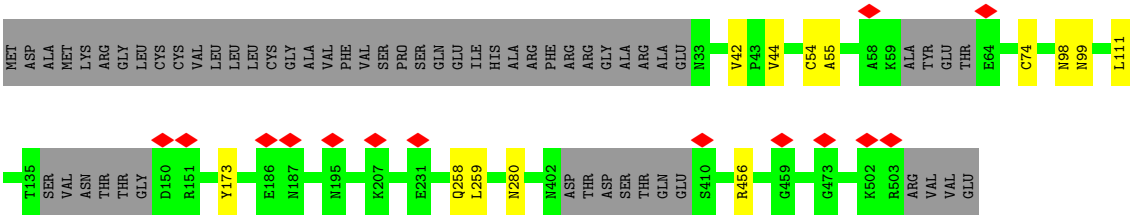




• Molecule 3: Envelope gp120

Chain d:

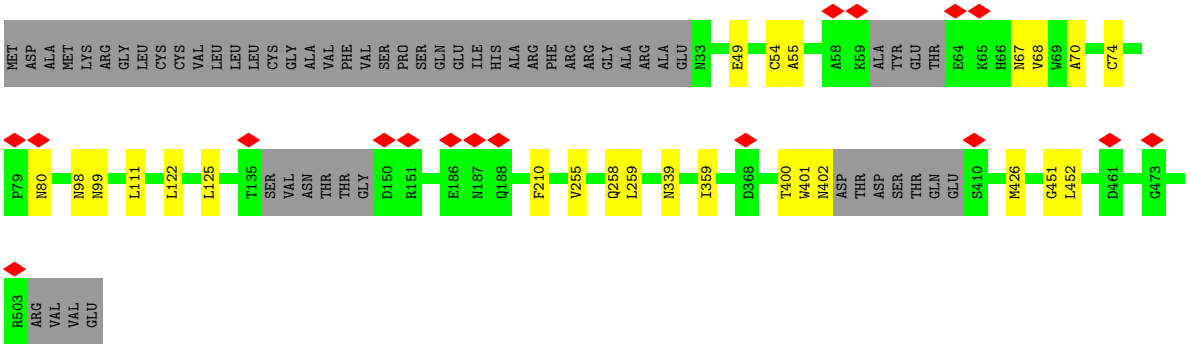




• Molecule 3: Envelope gp120

Chain f:

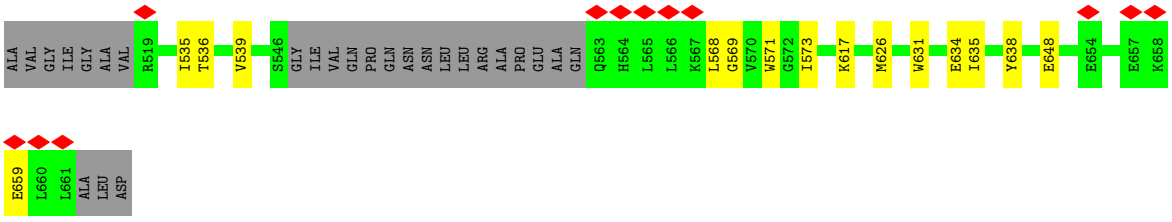




• Molecule 4: Envelope gp41 ectodomain

Chain c:

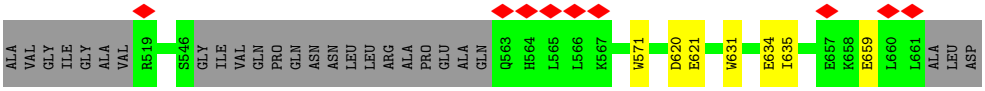




• Molecule 4: Envelope gp41 ectodomain

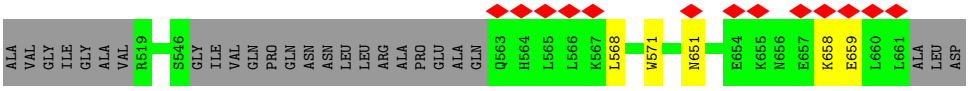
Chain e:





• Molecule 4: Envelope gp41 ectodomain





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





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





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





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





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





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

Chain K:





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

Chain L:





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

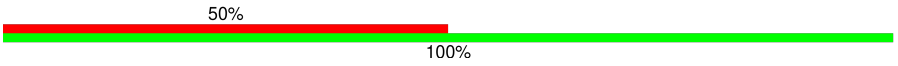
Chain M:





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

Chain N:





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

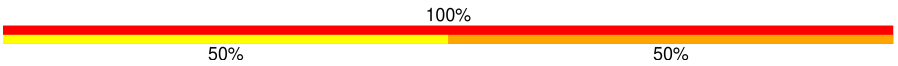
Chain O:





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

Chain R:





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

Chain T:





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

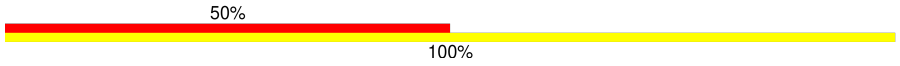
Chain U:





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

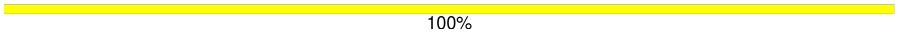
Chain V:





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

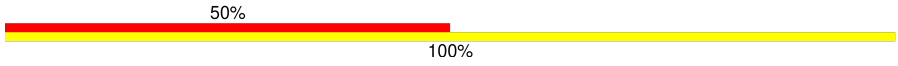
Chain W:





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

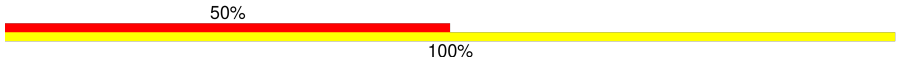
Chain X:





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

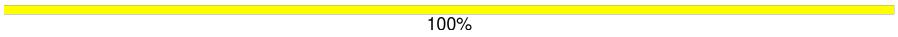
Chain Y:





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

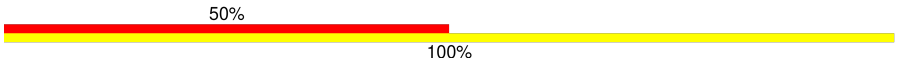
Chain Z:





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

Chain a:





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

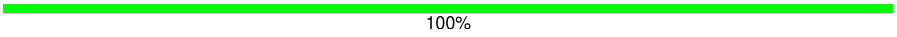
Chain h:





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

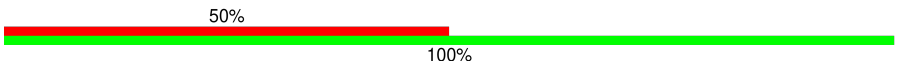
Chain j:





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

Chain m:





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

Chain o:





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

Chain p:





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

Chain q:





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

Chain r:





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

Chain s:





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

Chain t:





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

Chain u:





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

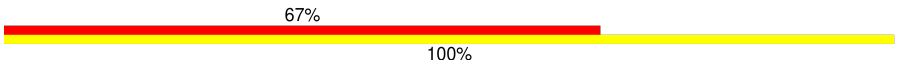
Chain D:





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

Chain J:





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

Chain P:





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

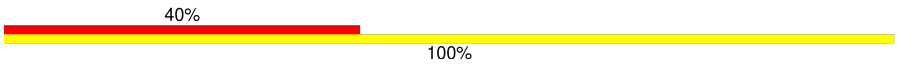
Chain k:





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

Chain E:



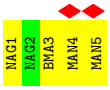


• Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyra

nose

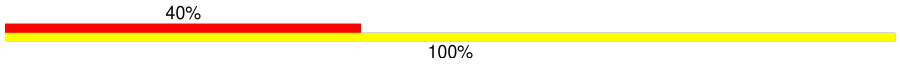
Chain Q:





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

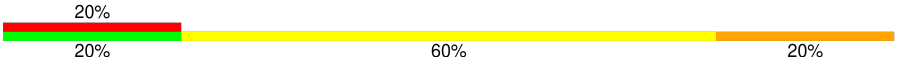
Chain l:

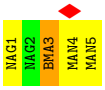




• Molecule 7: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain n:





• Molecule 8: 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

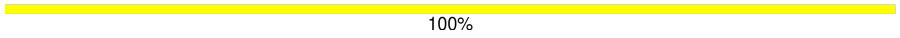
Chain S:





• Molecule 9: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose

Chain i:





4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, POINT, POINT, POINT	Depositor
Number of particles used	204457, 204457, 204457, 204457	Depositor
Resolution determination method	FSC 0.143 CUT-OFF, FSC 0.143 CUT-OFF, FSC 0.143 CUT-OFF, FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION, PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	58	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.412	Depositor
Minimum map value	-0.208	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.011	Depositor
Recommended contour level	0.09	Depositor
Map size (Å)	318.72, 318.72, 318.72	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83, 0.83, 0.83	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, MAN, NAG, TYS

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	0.21	0/1043	0.54	0/1422
2	B	0.19	0/895	0.55	1/1217 (0.1%)
3	b	0.18	0/3601	0.51	0/4888
3	d	0.17	0/3601	0.49	0/4888
3	f	0.18	0/3601	0.52	0/4888
4	c	0.21	0/1044	0.54	0/1410
4	e	0.22	0/1044	0.55	0/1410
4	g	0.19	0/1044	0.50	0/1410
All	All	0.19	0/15873	0.52	1/21533 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	108	LYS	CB-CG-CD	5.33	123.55	111.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1036	0	972	16	0
2	B	873	0	863	11	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	b	3525	0	3445	24	0
3	d	3525	0	3446	8	0
3	f	3525	0	3445	16	0
4	c	1026	0	1025	13	0
4	e	1026	0	1025	5	0
4	g	1026	0	1025	4	0
5	C	28	0	25	0	0
5	F	28	0	25	0	0
5	G	28	0	25	0	0
5	H	28	0	25	0	0
5	I	28	0	25	0	0
5	K	28	0	25	0	0
5	L	28	0	25	0	0
5	M	28	0	25	0	0
5	N	28	0	25	0	0
5	O	28	0	25	0	0
5	R	28	0	25	1	0
5	T	28	0	25	0	0
5	U	28	0	25	0	0
5	V	28	0	25	0	0
5	W	28	0	25	0	0
5	X	28	0	25	0	0
5	Y	28	0	25	0	0
5	Z	28	0	25	0	0
5	a	28	0	25	0	0
5	h	28	0	25	0	0
5	j	28	0	25	0	0
5	m	28	0	25	0	0
5	o	28	0	25	0	0
5	p	28	0	25	0	0
5	q	28	0	25	1	0
5	r	28	0	25	0	0
5	s	28	0	25	0	0
5	t	28	0	25	0	0
5	u	28	0	25	1	0
6	D	39	0	34	0	0
6	J	39	0	34	0	0
6	P	39	0	34	0	0
6	k	39	0	34	0	0
7	E	61	0	52	0	0
7	Q	61	0	52	0	0
7	l	61	0	52	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	n	61	0	52	1	0
8	S	50	0	43	3	0
9	i	33	0	28	0	0
10	b	112	0	104	0	0
10	c	42	0	39	1	0
10	d	84	0	78	0	0
10	e	42	0	39	1	0
10	f	112	0	104	1	0
10	g	42	0	39	0	0
All	All	17291	0	16789	95	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

The worst 5 of 95 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:B:26:SER:OG	2:B:27:GLN:OE1	1.94	0.86
8:S:3:BMA:HO6	8:S:3:BMA:HO2	1.41	0.67
4:e:620:ASP:OD2	4:e:621:GLU:N	2.28	0.66
1:A:4:LEU:HD12	1:A:22:CYS:SG	2.36	0.66
3:d:54:CYS:SG	3:d:55:ALA:N	2.68	0.65

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	132/248 (53%)	126 (96%)	6 (4%)	0	100	100
2	B	112/221 (51%)	110 (98%)	2 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
3	b	437/503 (87%)	412 (94%)	24 (6%)	1 (0%)	43 71
3	d	437/503 (87%)	416 (95%)	21 (5%)	0	100 100
3	f	437/503 (87%)	416 (95%)	21 (5%)	0	100 100
4	c	123/153 (80%)	118 (96%)	5 (4%)	0	100 100
4	e	123/153 (80%)	119 (97%)	4 (3%)	0	100 100
4	g	123/153 (80%)	119 (97%)	4 (3%)	0	100 100
All	All	1924/2437 (79%)	1836 (95%)	87 (4%)	1 (0%)	49 75

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	b	187	ASN

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	113/213 (53%)	113 (100%)	0	100 100
2	B	99/196 (50%)	99 (100%)	0	100 100
3	b	398/446 (89%)	398 (100%)	0	100 100
3	d	398/446 (89%)	398 (100%)	0	100 100
3	f	398/446 (89%)	397 (100%)	1 (0%)	86 86
4	c	110/128 (86%)	109 (99%)	1 (1%)	70 78
4	e	110/128 (86%)	110 (100%)	0	100 100
4	g	110/128 (86%)	110 (100%)	0	100 100
All	All	1736/2131 (82%)	1734 (100%)	2 (0%)	87 90

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
4	c	638	TYR

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Mol	Chain	Res	Type
3	f	339	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 8 such sidechains are listed below:

Mol	Chain	Res	Type
4	g	656	ASN
4	g	652	GLN
4	e	607	ASN
3	d	425	ASN
3	f	85	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)

1 non-standard protein/DNA/RNA residue is 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	$\#\|Z\| > 2$
1	TYS	A	111	1	15,16,17	1.59	2 (13%)	15,22,24	0.87	1 (6%)

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
1	TYS	A	111	1	-	0/10/11/13	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	111	TYS	OH-S	5.05	1.68	1.58
1	A	111	TYS	OH-CZ	-3.10	1.37	1.42

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	111	TYS	O3-S-O2	2.01	115.57	108.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates (i)

97 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	NAG	C	1	3,5	14,14,15	0.70	0	17,19,21	0.92	1 (5%)
5	NAG	C	2	5	14,14,15	0.68	0	17,19,21	0.84	0
6	NAG	D	1	3,6	14,14,15	0.70	0	17,19,21	0.89	0
6	NAG	D	2	6	14,14,15	0.69	0	17,19,21	0.88	0
6	BMA	D	3	6	11,11,12	0.85	0	15,15,17	2.76	6 (40%)
7	NAG	E	1	3,7	14,14,15	0.73	0	17,19,21	1.12	1 (5%)
7	NAG	E	2	7	14,14,15	0.71	0	17,19,21	0.96	1 (5%)
7	BMA	E	3	7	11,11,12	0.82	0	15,15,17	2.49	6 (40%)
7	MAN	E	4	7	11,11,12	0.64	0	15,15,17	1.27	1 (6%)
7	MAN	E	5	7	11,11,12	0.73	0	15,15,17	0.98	1 (6%)
5	NAG	F	1	3,5	14,14,15	0.73	0	17,19,21	1.01	1 (5%)
5	NAG	F	2	5	14,14,15	0.68	0	17,19,21	0.89	0
5	NAG	G	1	3,5	14,14,15	0.71	0	17,19,21	0.88	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	NAG	G	2	5	14,14,15	0.72	0	17,19,21	0.96	0
5	NAG	H	1	3,5	14,14,15	0.71	0	17,19,21	1.08	1 (5%)
5	NAG	H	2	5	14,14,15	0.70	0	17,19,21	0.97	0
5	NAG	I	1	3,5	14,14,15	0.70	0	17,19,21	0.96	1 (5%)
5	NAG	I	2	5	14,14,15	0.68	0	17,19,21	0.87	0
6	NAG	J	1	3,6	14,14,15	0.82	0	17,19,21	1.09	1 (5%)
6	NAG	J	2	6	14,14,15	0.79	0	17,19,21	1.33	1 (5%)
6	BMA	J	3	6	11,11,12	0.87	0	15,15,17	2.91	6 (40%)
5	NAG	K	1	3,5	14,14,15	0.73	0	17,19,21	1.10	2 (11%)
5	NAG	K	2	5	14,14,15	0.76	0	17,19,21	0.91	0
5	NAG	L	1	3,5	14,14,15	0.67	0	17,19,21	0.91	1 (5%)
5	NAG	L	2	5	14,14,15	0.73	0	17,19,21	1.20	1 (5%)
5	NAG	M	1	3,5	14,14,15	0.78	0	17,19,21	1.17	1 (5%)
5	NAG	M	2	5	14,14,15	0.76	0	17,19,21	1.38	1 (5%)
5	NAG	N	1	3,5	14,14,15	0.65	0	17,19,21	0.85	0
5	NAG	N	2	5	14,14,15	0.71	0	17,19,21	0.98	0
5	NAG	O	1	3,5	14,14,15	0.69	0	17,19,21	0.91	2 (11%)
5	NAG	O	2	5	14,14,15	0.67	0	17,19,21	0.87	0
6	NAG	P	1	3,6	14,14,15	0.71	0	17,19,21	0.86	1 (5%)
6	NAG	P	2	6	14,14,15	0.70	0	17,19,21	0.90	0
6	BMA	P	3	6	11,11,12	0.85	0	15,15,17	2.73	6 (40%)
7	NAG	Q	1	3,7	14,14,15	0.74	0	17,19,21	1.17	1 (5%)
7	NAG	Q	2	7	14,14,15	0.70	0	17,19,21	0.91	0
7	BMA	Q	3	7	11,11,12	0.81	0	15,15,17	2.28	5 (33%)
7	MAN	Q	4	7	11,11,12	0.63	0	15,15,17	1.32	1 (6%)
7	MAN	Q	5	7	11,11,12	0.67	0	15,15,17	1.19	1 (6%)
5	NAG	R	1	3,5	14,14,15	0.74	0	17,19,21	0.95	1 (5%)
5	NAG	R	2	5	14,14,15	0.70	0	17,19,21	0.93	1 (5%)
8	NAG	S	1	8,3	14,14,15	0.76	0	17,19,21	1.10	1 (5%)
8	NAG	S	2	8	14,14,15	0.71	0	17,19,21	0.79	0
8	BMA	S	3	8	11,11,12	0.71	0	15,15,17	2.59	4 (26%)
8	MAN	S	4	8	11,11,12	0.67	0	15,15,17	1.17	1 (6%)
5	NAG	T	1	3,5	14,14,15	0.72	0	17,19,21	1.13	1 (5%)
5	NAG	T	2	5	14,14,15	0.70	0	17,19,21	0.94	0
5	NAG	U	1	3,5	14,14,15	0.70	0	17,19,21	0.98	1 (5%)
5	NAG	U	2	5	14,14,15	0.68	0	17,19,21	0.89	0
5	NAG	V	1	3,5	14,14,15	0.80	0	17,19,21	1.03	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	NAG	V	2	5	14,14,15	0.79	0	17,19,21	1.40	1 (5%)
5	NAG	W	1	3,5	14,14,15	0.69	0	17,19,21	1.22	2 (11%)
5	NAG	W	2	5	14,14,15	0.77	0	17,19,21	0.93	1 (5%)
5	NAG	X	1	3,5	14,14,15	0.70	0	17,19,21	0.94	1 (5%)
5	NAG	X	2	5	14,14,15	0.72	0	17,19,21	1.29	1 (5%)
5	NAG	Y	1	3,5	14,14,15	0.74	0	17,19,21	1.67	4 (23%)
5	NAG	Y	2	5	14,14,15	0.67	0	17,19,21	0.92	1 (5%)
5	NAG	Z	1	3,5	14,14,15	0.71	0	17,19,21	1.03	1 (5%)
5	NAG	Z	2	5	14,14,15	0.69	0	17,19,21	0.93	1 (5%)
5	NAG	a	1	3,5	14,14,15	0.79	0	17,19,21	1.27	1 (5%)
5	NAG	a	2	5	14,14,15	0.74	0	17,19,21	1.16	1 (5%)
5	NAG	h	1	3,5	14,14,15	0.67	0	17,19,21	0.86	1 (5%)
5	NAG	h	2	5	14,14,15	0.69	0	17,19,21	0.98	0
9	MAN	i	1	9	11,11,12	0.85	1 (9%)	15,15,17	1.09	1 (6%)
9	MAN	i	2	9	11,11,12	0.65	0	15,15,17	1.31	1 (6%)
9	MAN	i	3	9	11,11,12	0.76	0	15,15,17	0.90	1 (6%)
5	NAG	j	1	3,5	14,14,15	0.69	0	17,19,21	0.85	0
5	NAG	j	2	5	14,14,15	0.68	0	17,19,21	0.89	0
6	NAG	k	1	3,6	14,14,15	0.70	0	17,19,21	0.91	0
6	NAG	k	2	6	14,14,15	0.70	0	17,19,21	0.92	0
6	BMA	k	3	6	11,11,12	0.85	0	15,15,17	2.79	6 (40%)
7	NAG	l	1	3,7	14,14,15	0.76	0	17,19,21	1.10	1 (5%)
7	NAG	l	2	7	14,14,15	0.70	0	17,19,21	1.44	2 (11%)
7	BMA	l	3	7	11,11,12	0.87	0	15,15,17	2.28	4 (26%)
7	MAN	l	4	7	11,11,12	0.68	0	15,15,17	1.23	1 (6%)
7	MAN	l	5	7	11,11,12	0.72	0	15,15,17	1.00	1 (6%)
5	NAG	m	1	3,5	14,14,15	0.71	0	17,19,21	0.84	0
5	NAG	m	2	5	14,14,15	0.69	0	17,19,21	0.86	0
7	NAG	n	1	3,7	14,14,15	0.72	0	17,19,21	1.07	2 (11%)
7	NAG	n	2	7	14,14,15	0.68	0	17,19,21	1.01	0
7	BMA	n	3	7	11,11,12	0.88	0	15,15,17	2.40	5 (33%)
7	MAN	n	4	7	11,11,12	0.72	0	15,15,17	1.14	1 (6%)
7	MAN	n	5	7	11,11,12	0.66	0	15,15,17	1.36	1 (6%)
5	NAG	o	1	3,5	14,14,15	0.70	0	17,19,21	1.10	1 (5%)
5	NAG	o	2	5	14,14,15	0.69	0	17,19,21	1.02	0
5	NAG	p	1	3,5	14,14,15	0.72	0	17,19,21	0.92	1 (5%)
5	NAG	p	2	5	14,14,15	0.69	0	17,19,21	0.83	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	NAG	q	1	3,5	14,14,15	0.65	0	17,19,21	1.32	3 (17%)
5	NAG	q	2	5	14,14,15	0.80	1 (7%)	17,19,21	1.06	1 (5%)
5	NAG	r	1	3,5	14,14,15	0.69	0	17,19,21	1.21	2 (11%)
5	NAG	r	2	5	14,14,15	0.76	0	17,19,21	0.88	0
5	NAG	s	1	3,5	14,14,15	0.69	0	17,19,21	0.88	1 (5%)
5	NAG	s	2	5	14,14,15	0.67	0	17,19,21	1.03	1 (5%)
5	NAG	t	1	3,5	14,14,15	0.77	0	17,19,21	1.13	1 (5%)
5	NAG	t	2	5	14,14,15	0.77	0	17,19,21	1.18	1 (5%)
5	NAG	u	1	3,5	14,14,15	0.73	0	17,19,21	0.92	1 (5%)
5	NAG	u	2	5	14,14,15	0.70	0	17,19,21	0.92	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	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	C	2	5	-	0/6/23/26	0/1/1/1
6	NAG	D	1	3,6	-	0/6/23/26	0/1/1/1
6	NAG	D	2	6	-	0/6/23/26	0/1/1/1
6	BMA	D	3	6	-	0/2/19/22	0/1/1/1
7	NAG	E	1	3,7	-	0/6/23/26	0/1/1/1
7	NAG	E	2	7	-	0/6/23/26	0/1/1/1
7	BMA	E	3	7	-	0/2/19/22	0/1/1/1
7	MAN	E	4	7	-	0/2/19/22	0/1/1/1
7	MAN	E	5	7	-	0/2/19/22	0/1/1/1
5	NAG	F	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	F	2	5	-	0/6/23/26	0/1/1/1
5	NAG	G	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	G	2	5	-	2/6/23/26	0/1/1/1
5	NAG	H	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	H	2	5	-	3/6/23/26	0/1/1/1
5	NAG	I	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	1/6/23/26	0/1/1/1
6	NAG	J	1	3,6	-	1/6/23/26	0/1/1/1
6	NAG	J	2	6	-	0/6/23/26	0/1/1/1
6	BMA	J	3	6	-	1/2/19/22	0/1/1/1
5	NAG	K	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	K	2	5	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	L	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	L	2	5	-	3/6/23/26	0/1/1/1
5	NAG	M	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	M	2	5	-	1/6/23/26	0/1/1/1
5	NAG	N	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	N	2	5	-	3/6/23/26	0/1/1/1
5	NAG	O	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	O	2	5	-	0/6/23/26	0/1/1/1
6	NAG	P	1	3,6	-	0/6/23/26	0/1/1/1
6	NAG	P	2	6	-	0/6/23/26	0/1/1/1
6	BMA	P	3	6	-	0/2/19/22	0/1/1/1
7	NAG	Q	1	3,7	-	0/6/23/26	0/1/1/1
7	NAG	Q	2	7	-	0/6/23/26	0/1/1/1
7	BMA	Q	3	7	-	1/2/19/22	0/1/1/1
7	MAN	Q	4	7	-	0/2/19/22	0/1/1/1
7	MAN	Q	5	7	-	0/2/19/22	0/1/1/1
5	NAG	R	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	R	2	5	-	0/6/23/26	0/1/1/1
8	NAG	S	1	8,3	-	0/6/23/26	0/1/1/1
8	NAG	S	2	8	-	0/6/23/26	0/1/1/1
8	BMA	S	3	8	-	0/2/19/22	0/1/1/1
8	MAN	S	4	8	-	1/2/19/22	0/1/1/1
5	NAG	T	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	T	2	5	-	3/6/23/26	0/1/1/1
5	NAG	U	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	U	2	5	-	1/6/23/26	0/1/1/1
5	NAG	V	1	3,5	-	1/6/23/26	0/1/1/1
5	NAG	V	2	5	-	0/6/23/26	0/1/1/1
5	NAG	W	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	W	2	5	-	0/6/23/26	0/1/1/1
5	NAG	X	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	X	2	5	-	3/6/23/26	0/1/1/1
5	NAG	Y	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	Y	2	5	-	0/6/23/26	0/1/1/1
5	NAG	Z	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	Z	2	5	-	1/6/23/26	0/1/1/1
5	NAG	a	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	a	2	5	-	1/6/23/26	0/1/1/1
5	NAG	h	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	h	2	5	-	3/6/23/26	0/1/1/1
9	MAN	i	1	9	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	MAN	i	2	9	-	1/2/19/22	0/1/1/1
9	MAN	i	3	9	-	1/2/19/22	0/1/1/1
5	NAG	j	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	j	2	5	-	0/6/23/26	0/1/1/1
6	NAG	k	1	3,6	-	0/6/23/26	0/1/1/1
6	NAG	k	2	6	-	2/6/23/26	0/1/1/1
6	BMA	k	3	6	-	0/2/19/22	0/1/1/1
7	NAG	l	1	3,7	-	0/6/23/26	0/1/1/1
7	NAG	l	2	7	-	0/6/23/26	0/1/1/1
7	BMA	l	3	7	-	0/2/19/22	0/1/1/1
7	MAN	l	4	7	-	0/2/19/22	0/1/1/1
7	MAN	l	5	7	-	0/2/19/22	0/1/1/1
5	NAG	m	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	m	2	5	-	0/6/23/26	0/1/1/1
7	NAG	n	1	3,7	-	0/6/23/26	0/1/1/1
7	NAG	n	2	7	-	2/6/23/26	0/1/1/1
7	BMA	n	3	7	-	0/2/19/22	0/1/1/1
7	MAN	n	4	7	-	0/2/19/22	0/1/1/1
7	MAN	n	5	7	-	1/2/19/22	0/1/1/1
5	NAG	o	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	o	2	5	-	3/6/23/26	0/1/1/1
5	NAG	p	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	p	2	5	-	1/6/23/26	0/1/1/1
5	NAG	q	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	q	2	5	-	0/6/23/26	0/1/1/1
5	NAG	r	1	3,5	-	2/6/23/26	0/1/1/1
5	NAG	r	2	5	-	0/6/23/26	0/1/1/1
5	NAG	s	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	s	2	5	-	3/6/23/26	0/1/1/1
5	NAG	t	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	t	2	5	-	1/6/23/26	0/1/1/1
5	NAG	u	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	u	2	5	-	3/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	i	1	MAN	O5-C1	-2.30	1.39	1.43
5	q	2	NAG	C1-C2	2.08	1.55	1.52

The worst 5 of 119 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
8	S	3	BMA	C1-O5-C5	8.79	123.96	112.19
6	J	3	BMA	C1-O5-C5	8.51	123.59	112.19
6	k	3	BMA	C1-O5-C5	7.88	122.75	112.19
6	D	3	BMA	C1-O5-C5	7.72	122.53	112.19
6	P	3	BMA	C1-O5-C5	7.61	122.39	112.19

There are no chirality outliers.

5 of 63 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	r	1	NAG	C1-C2-N2-C7
5	q	1	NAG	O5-C5-C6-O6
6	J	3	BMA	O5-C5-C6-O6
5	I	2	NAG	O5-C5-C6-O6
5	X	2	NAG	O5-C5-C6-O6

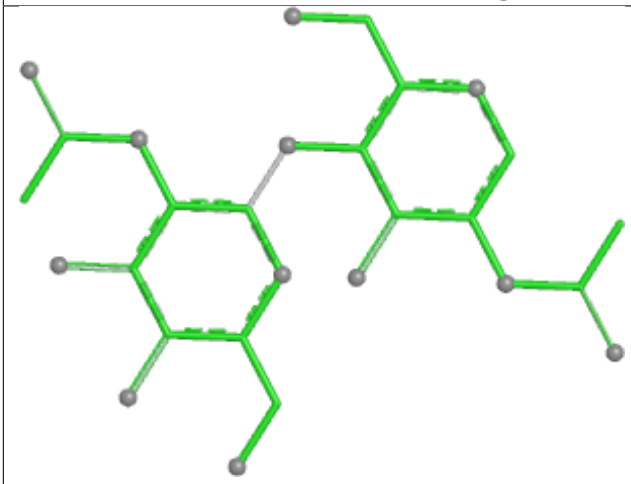
There are no ring outliers.

7 monomers are involved in 7 short contacts:

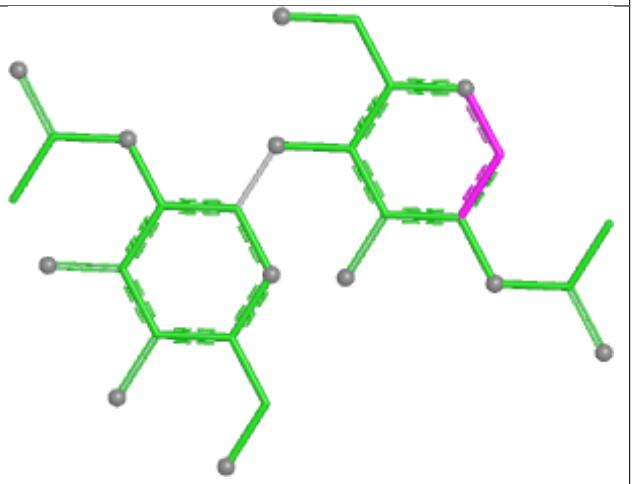
Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	S	3	BMA	2	0
5	R	2	NAG	1	0
5	u	1	NAG	1	0
5	q	2	NAG	1	0
7	n	3	BMA	1	0
8	S	2	NAG	1	0
5	q	1	NAG	1	0

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

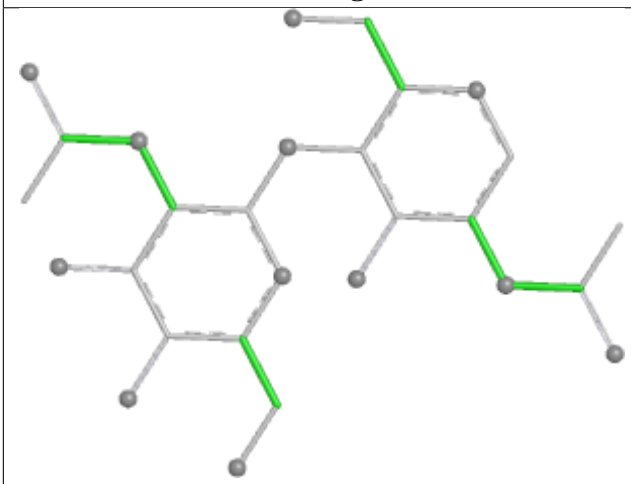
Oligosaccharide Chain C



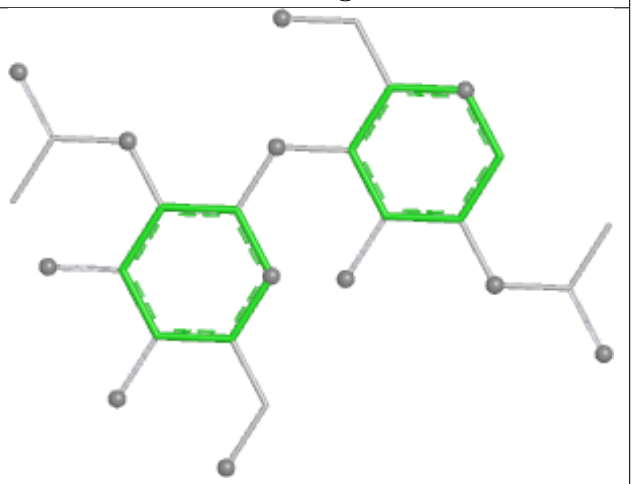
Bond lengths



Bond angles

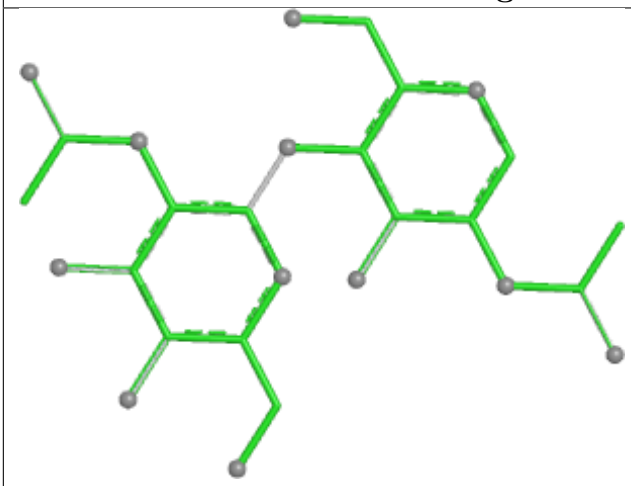


Torsions

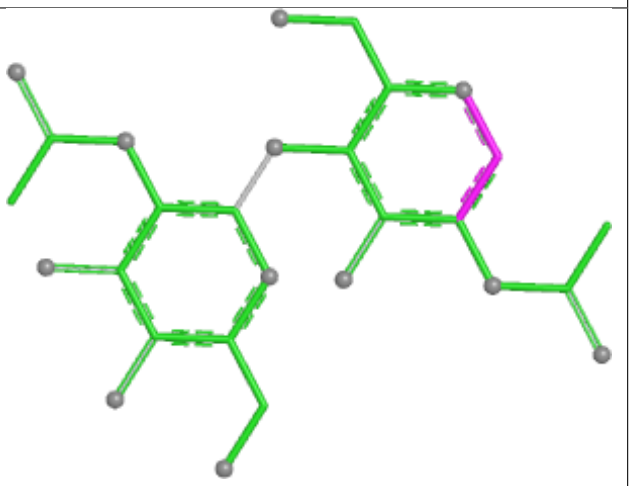


Rings

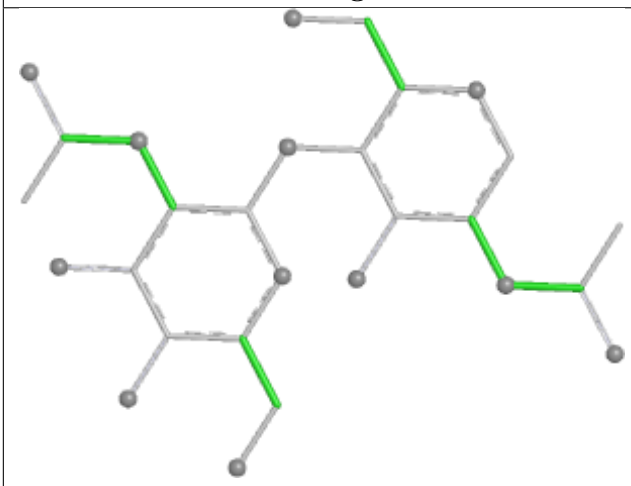
Oligosaccharide Chain F



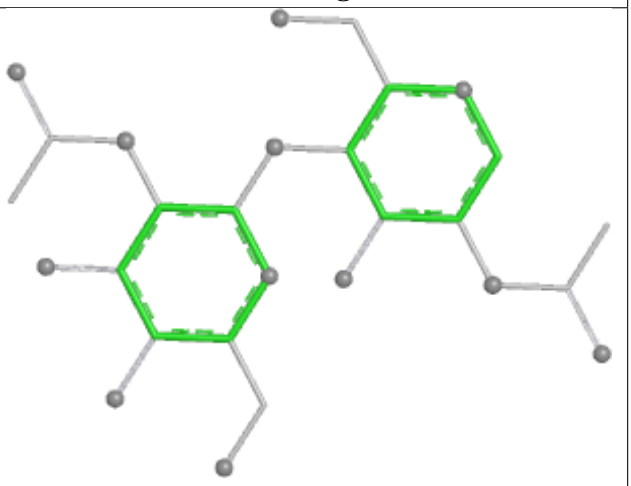
Bond lengths



Bond angles

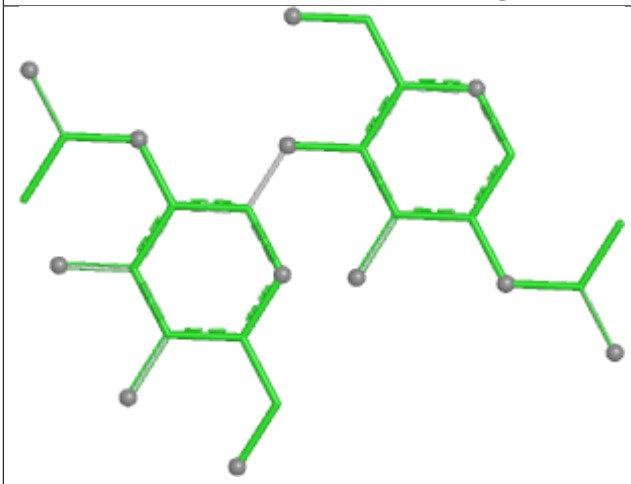


Torsions

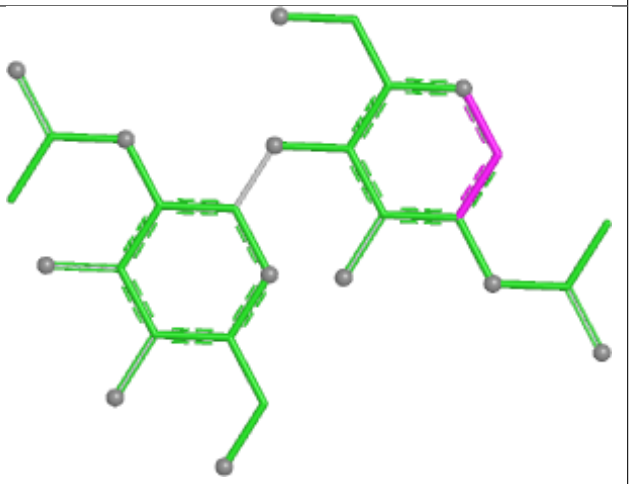


Rings

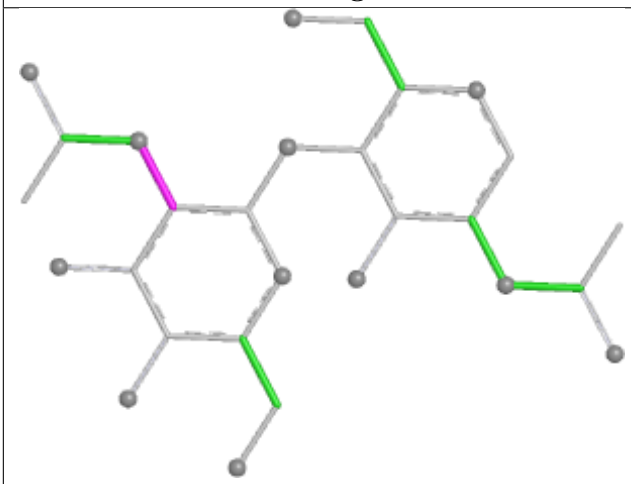
Oligosaccharide Chain G



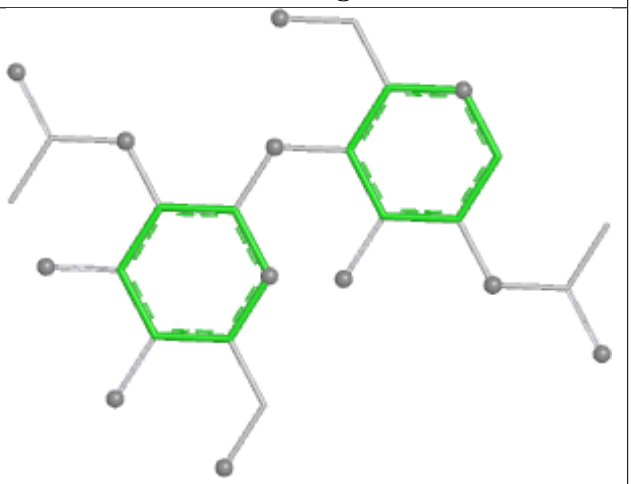
Bond lengths



Bond angles

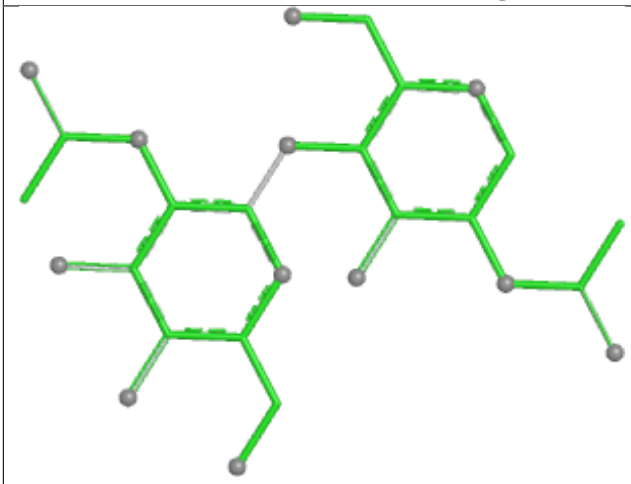


Torsions

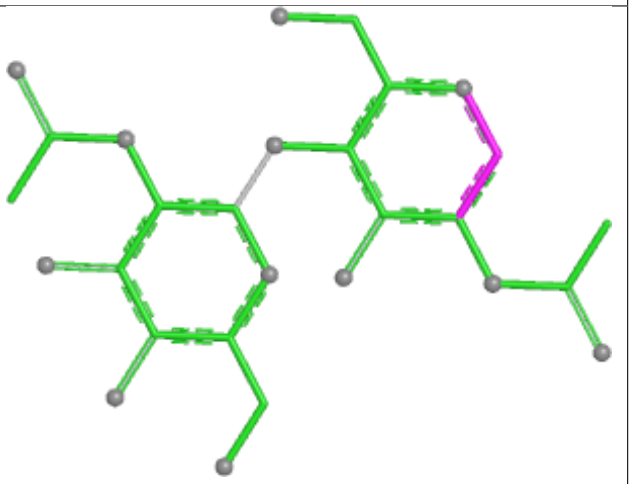


Rings

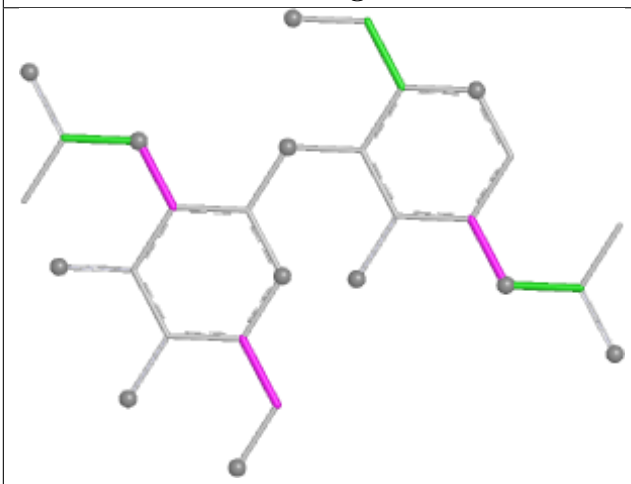
Oligosaccharide Chain H



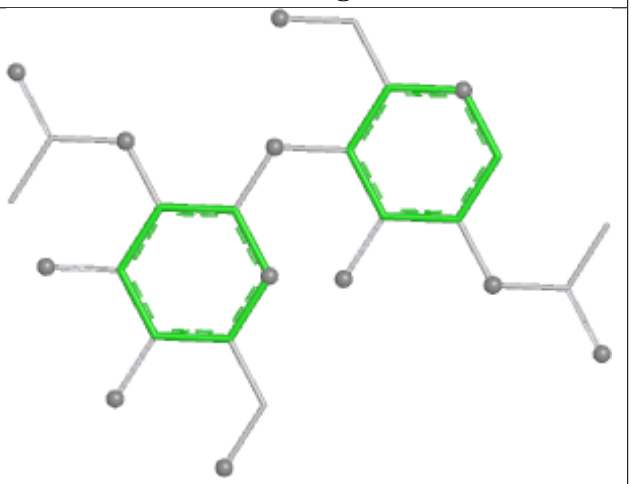
Bond lengths



Bond angles

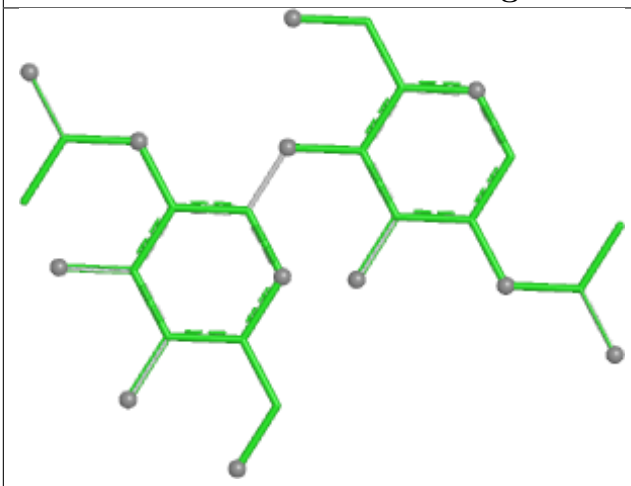


Torsions

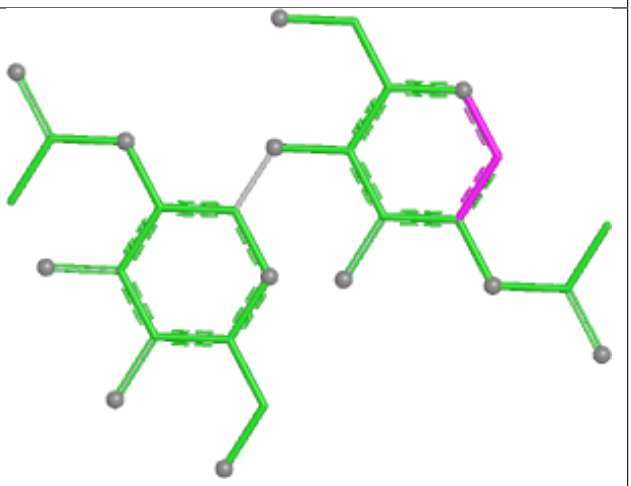


Rings

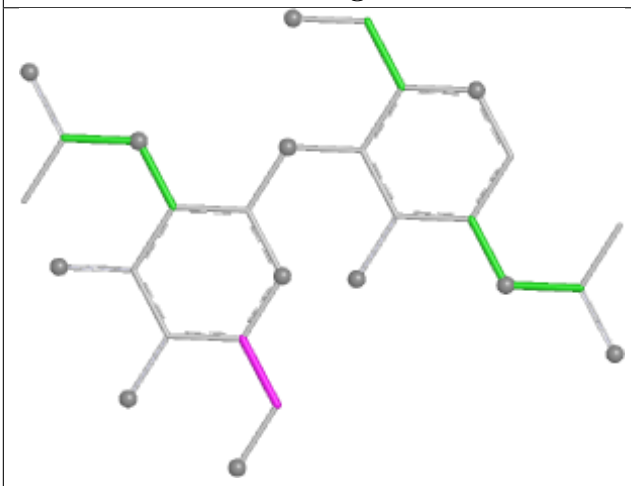
Oligosaccharide Chain I



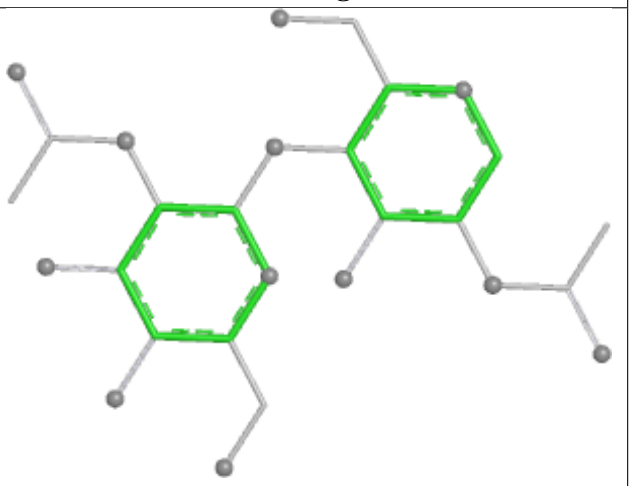
Bond lengths



Bond angles

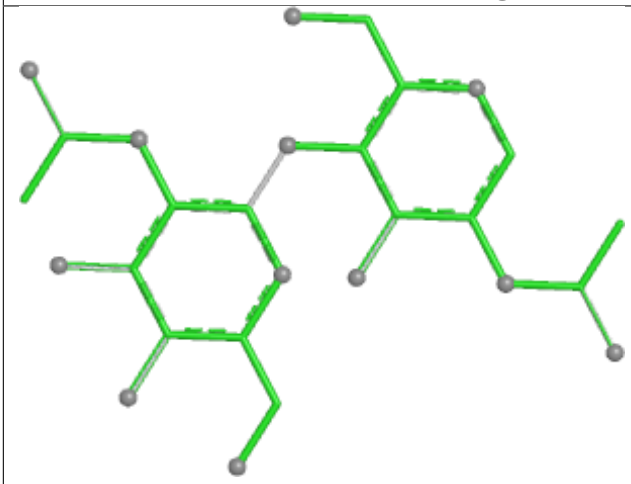


Torsions

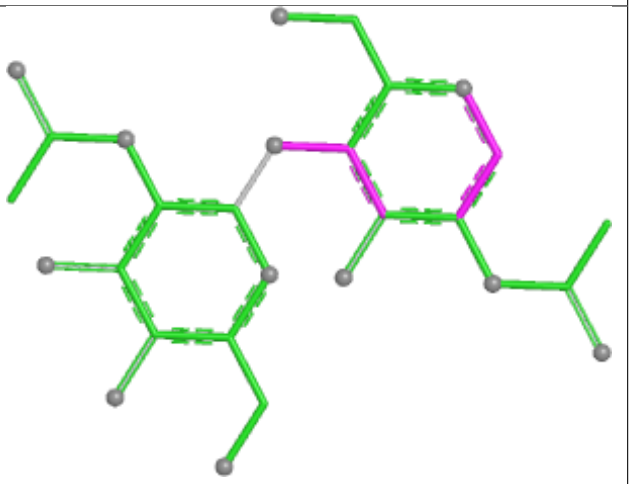


Rings

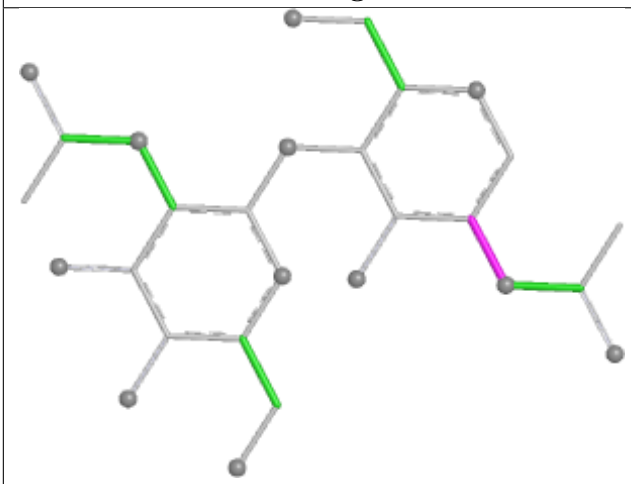
Oligosaccharide Chain K



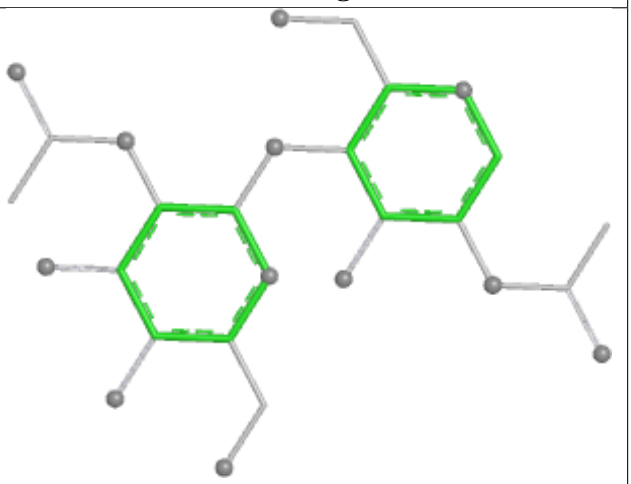
Bond lengths



Bond angles

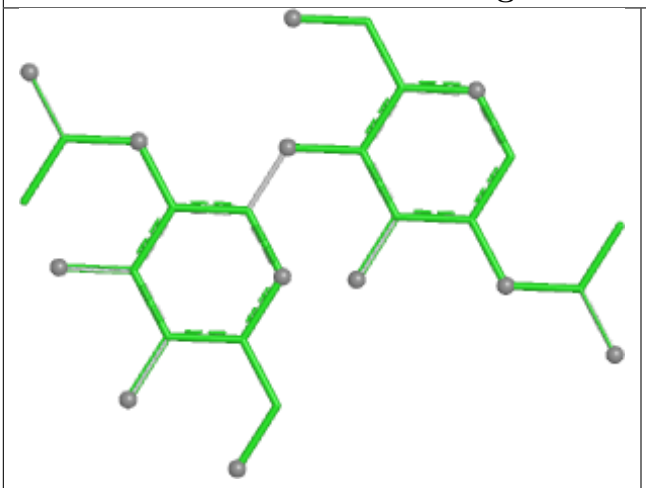


Torsions

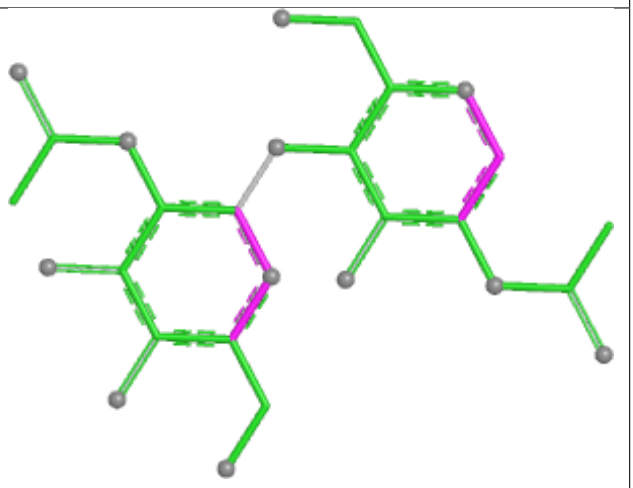


Rings

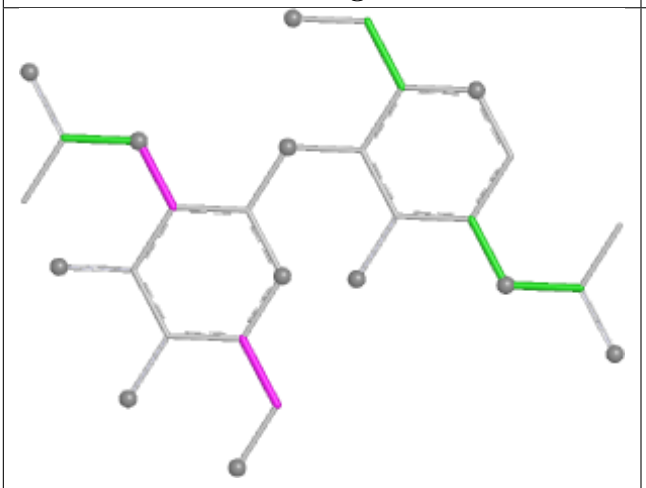
Oligosaccharide Chain L



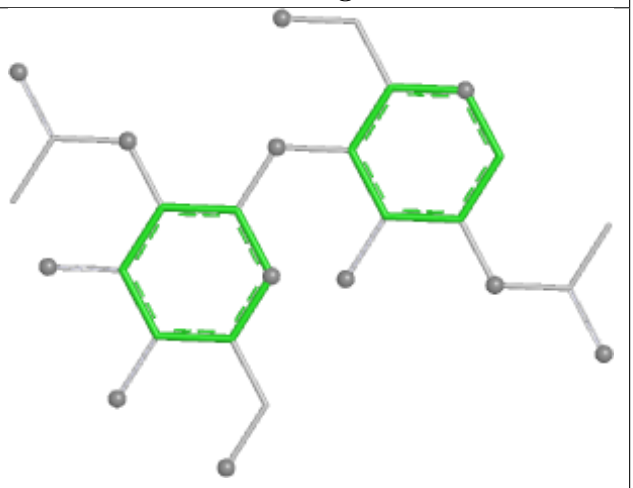
Bond lengths



Bond angles

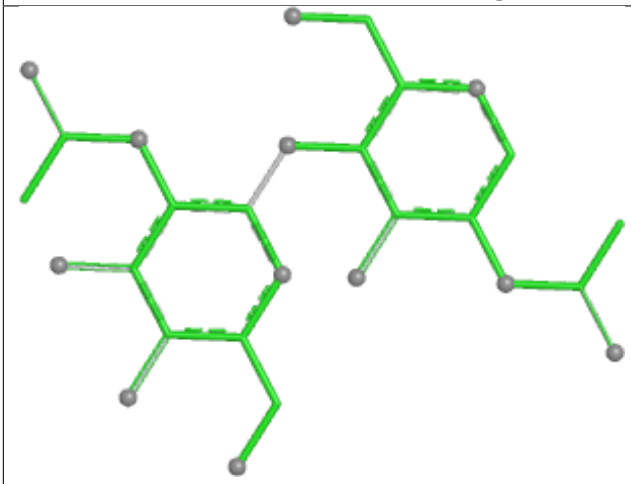


Torsions



Rings

Oligosaccharide Chain M



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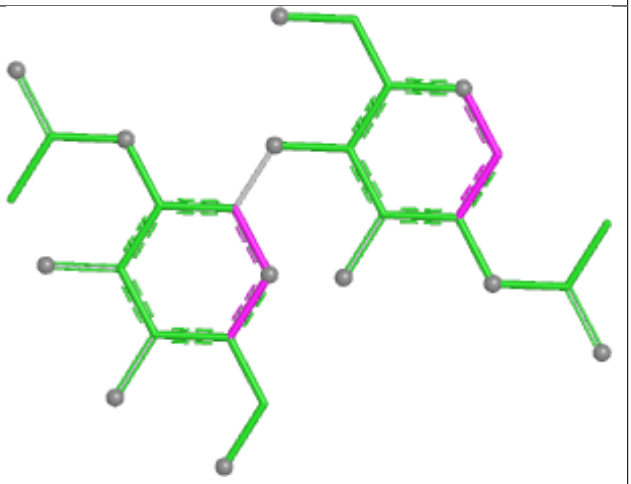

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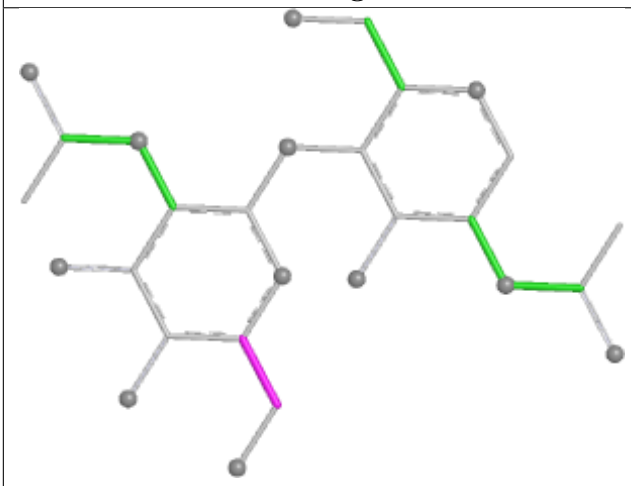

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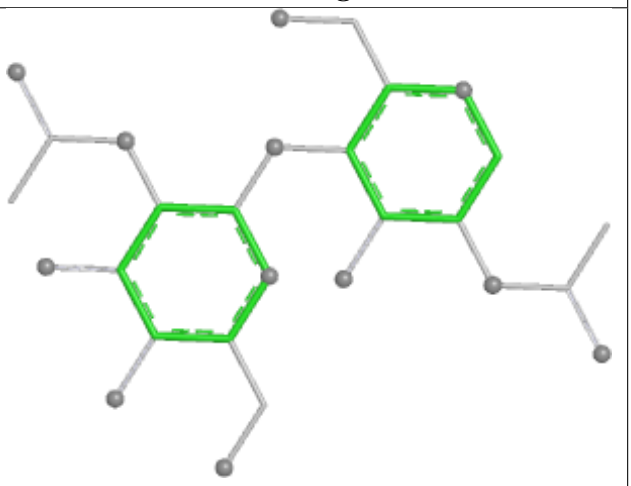
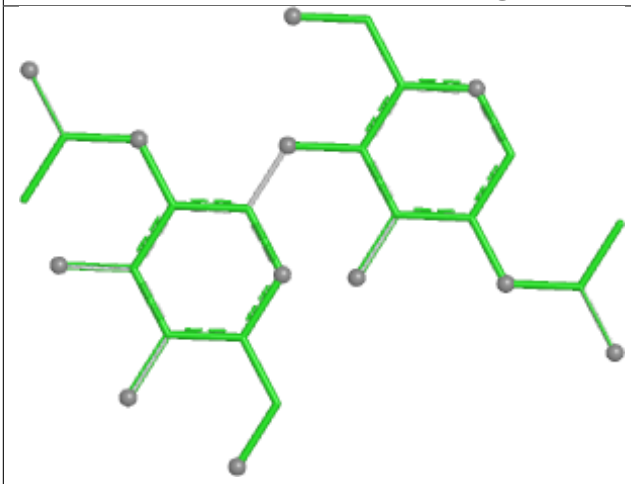


Rings

Oligosaccharide Chain N



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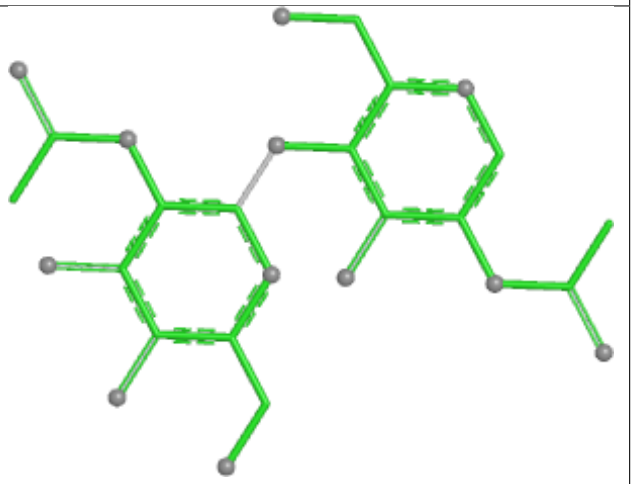

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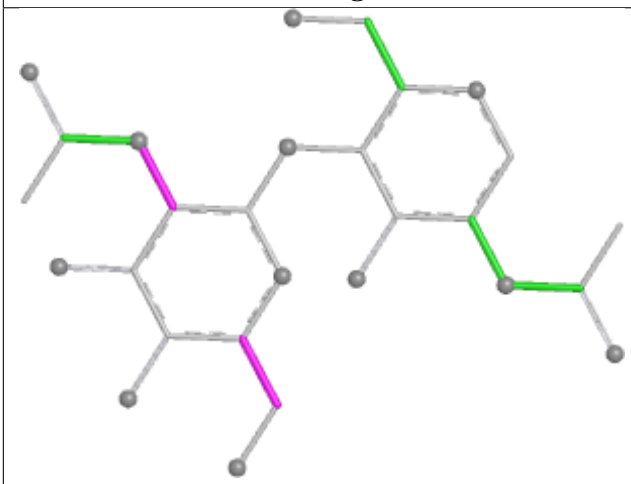

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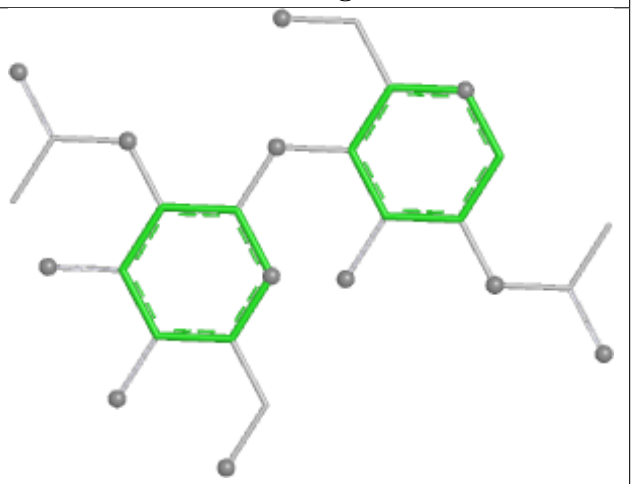
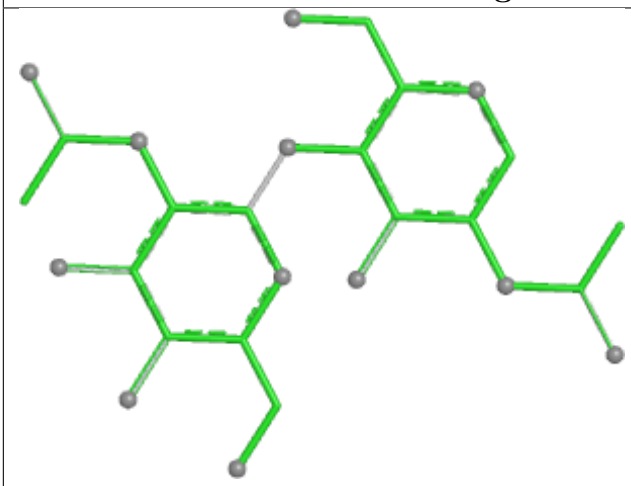


Rings

Oligosaccharide Chain O



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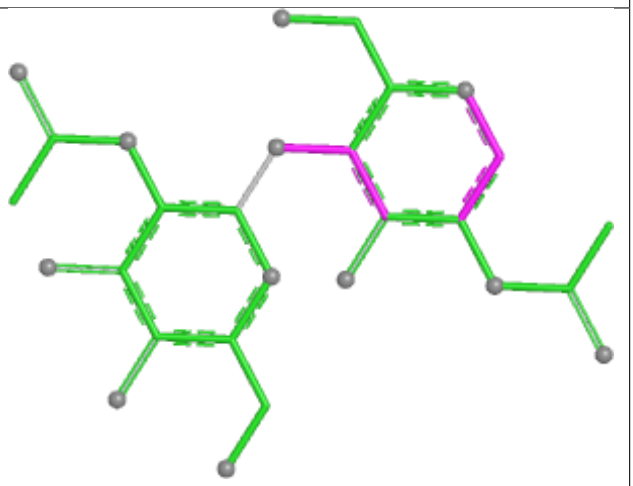

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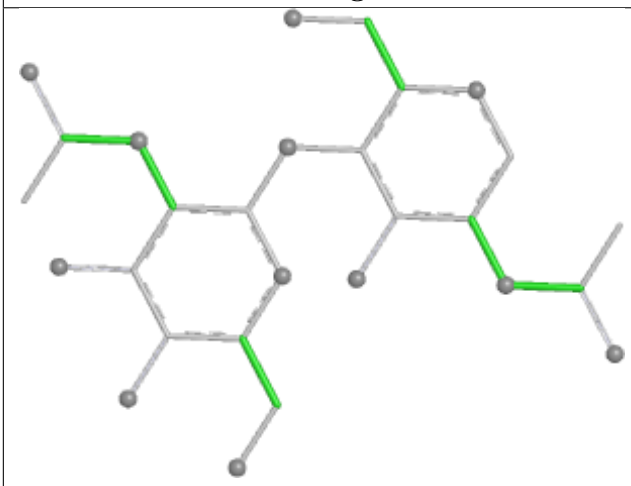

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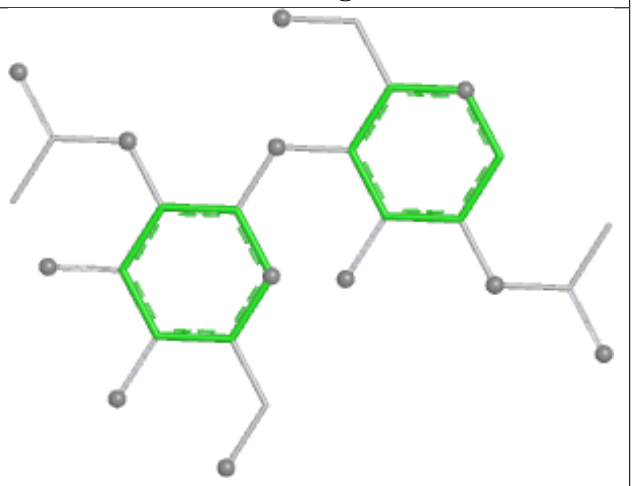
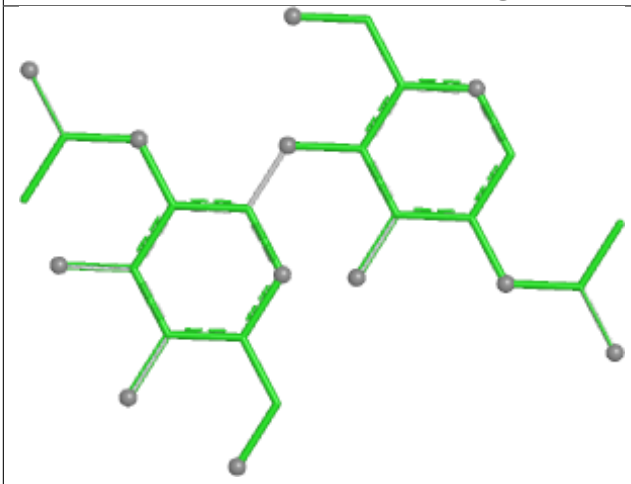


Rings

Oligosaccharide Chain R



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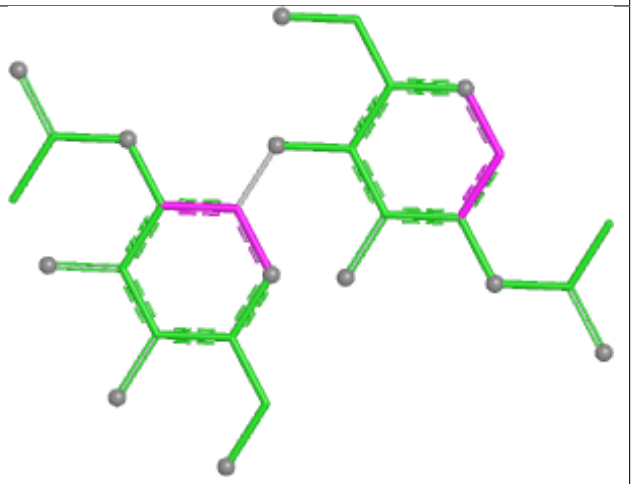
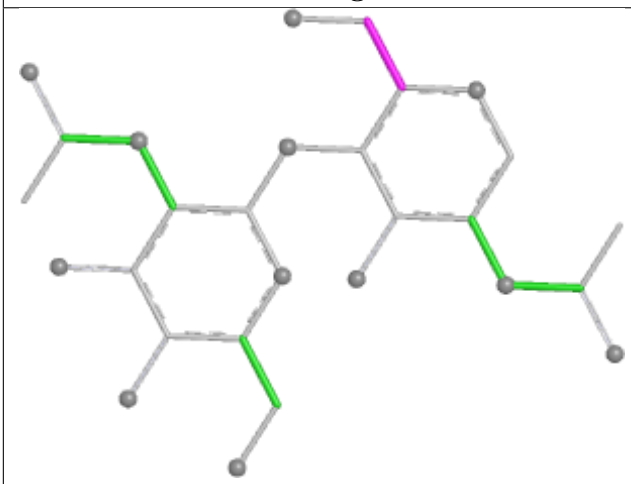


Bond angles



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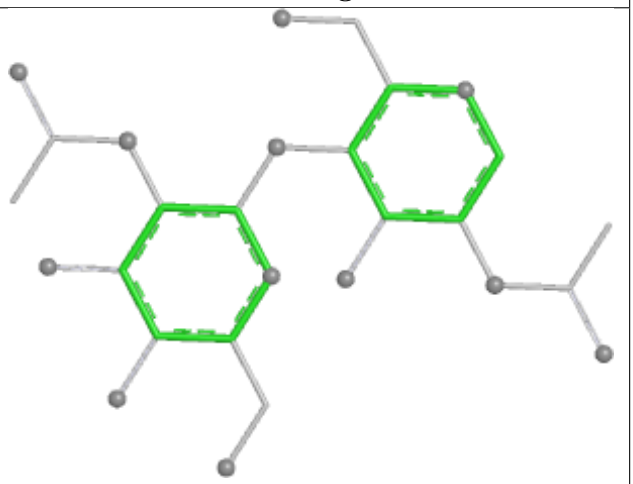
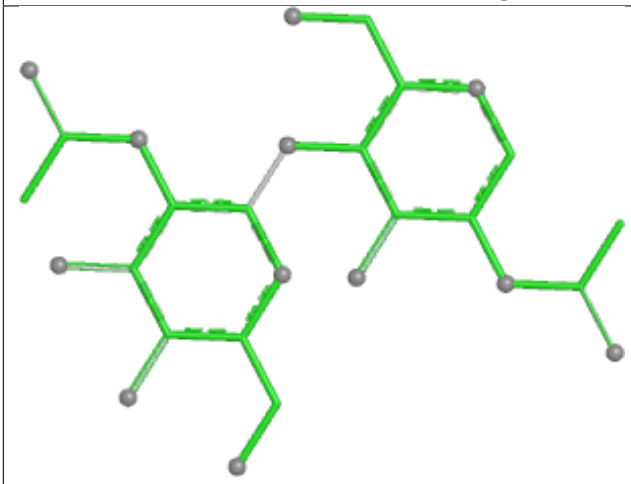


Rings

Oligosaccharide Chain T



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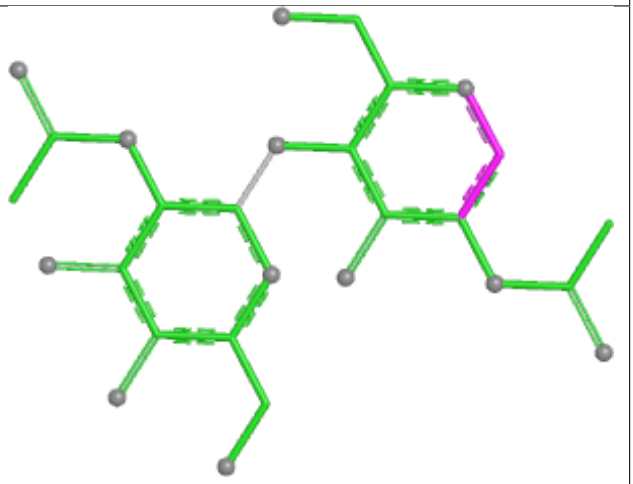

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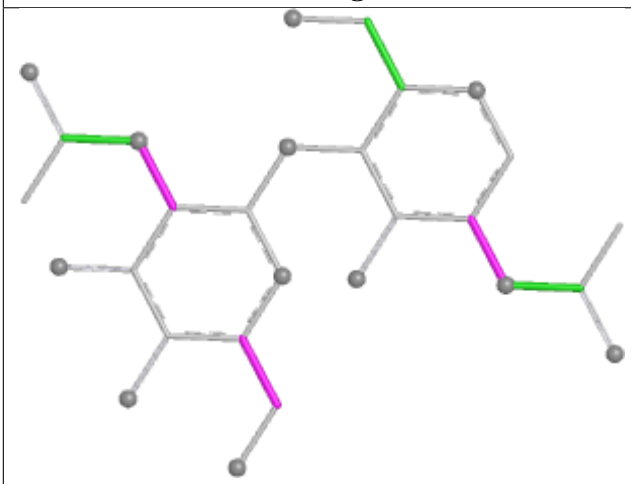

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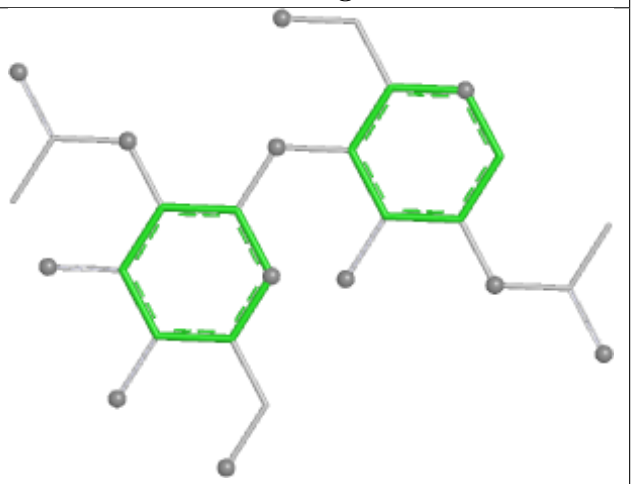
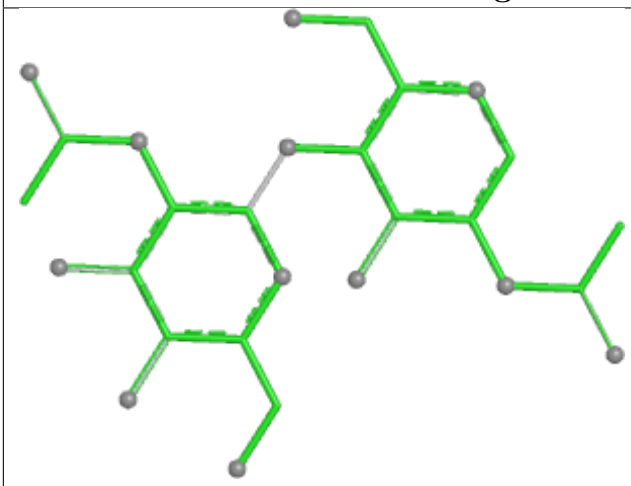


Rings

Oligosaccharide Chain U



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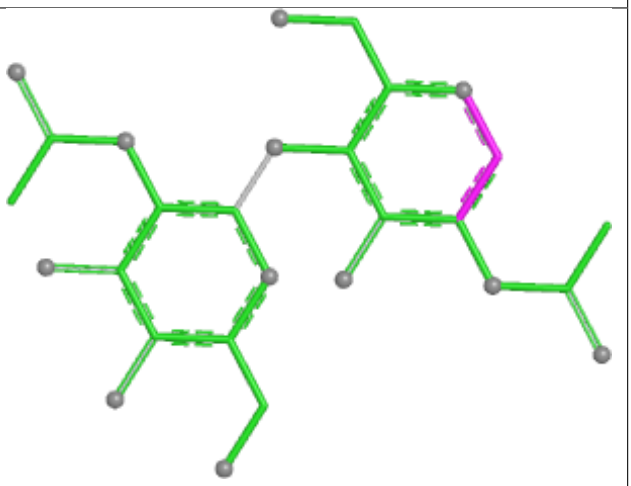

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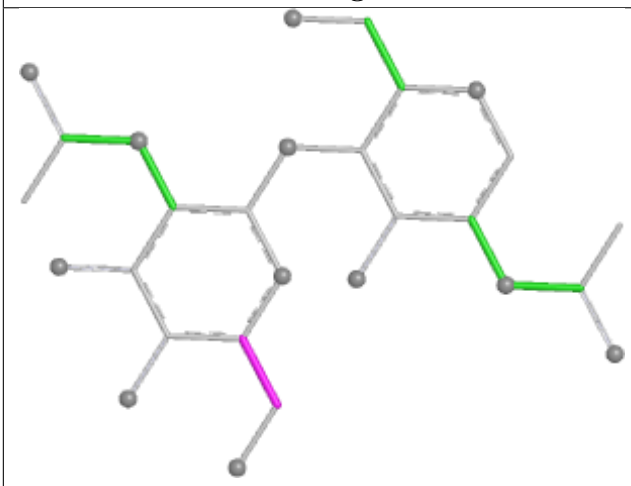

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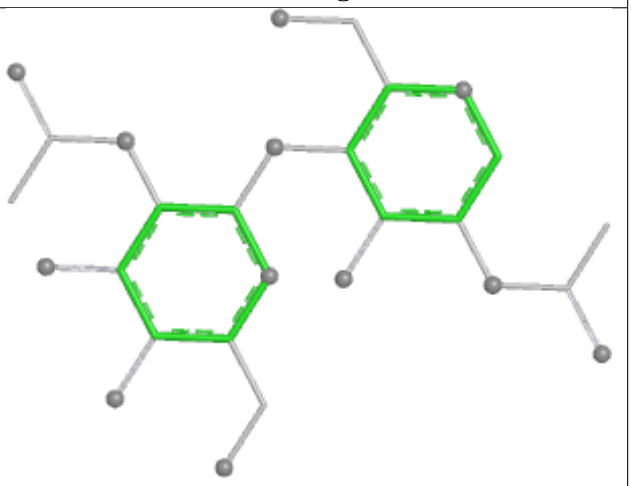
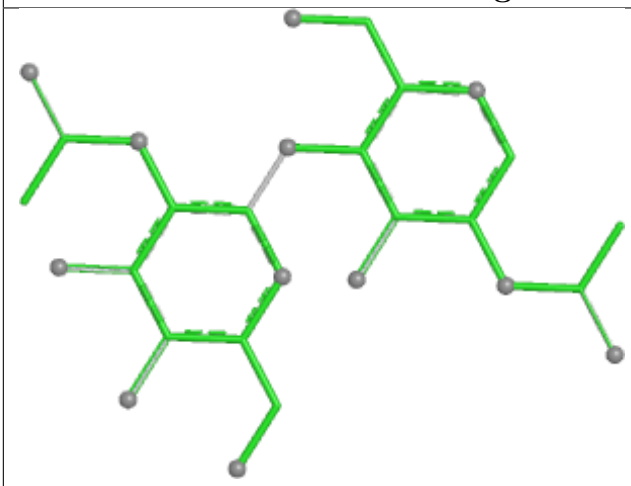


Rings

Oligosaccharide Chain V



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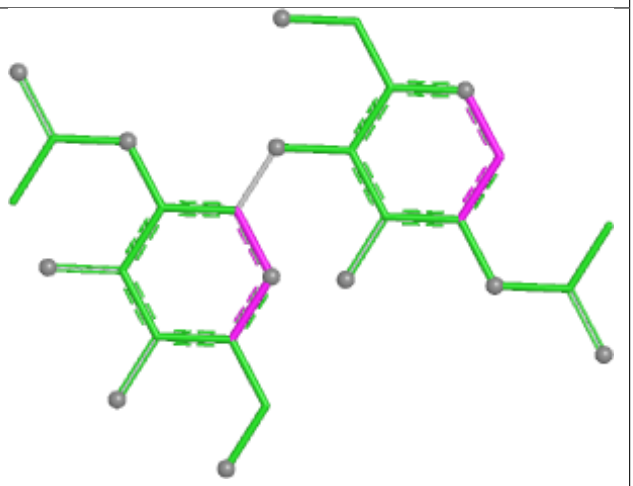

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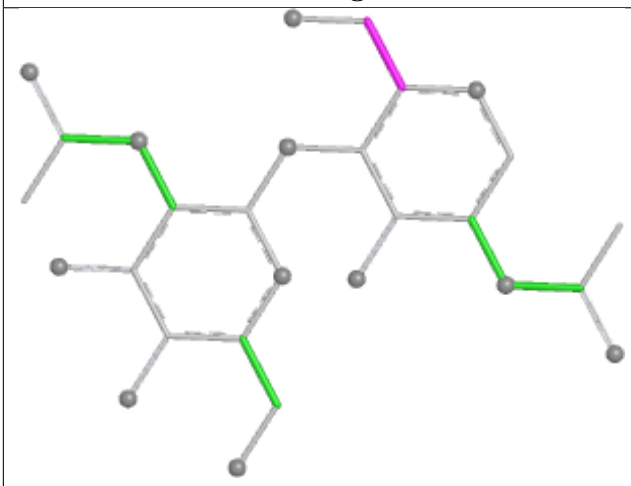

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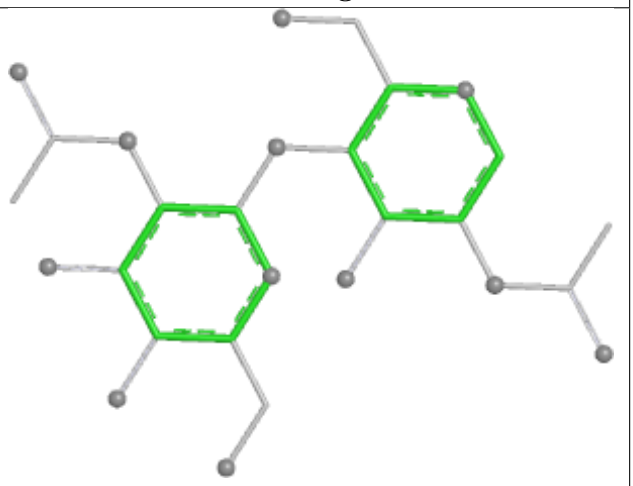
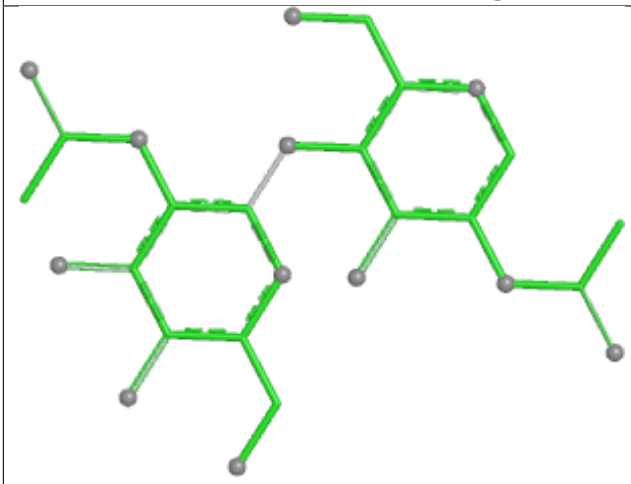


Rings

Oligosaccharide Chain W



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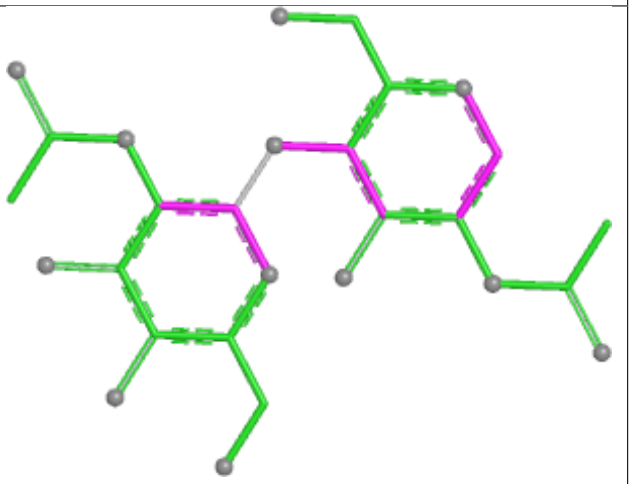

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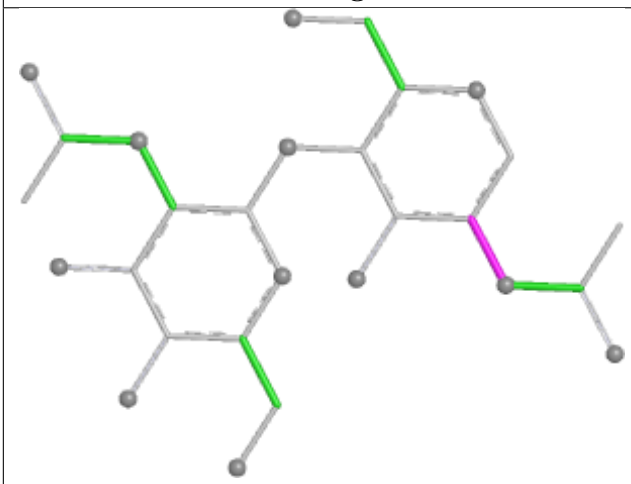

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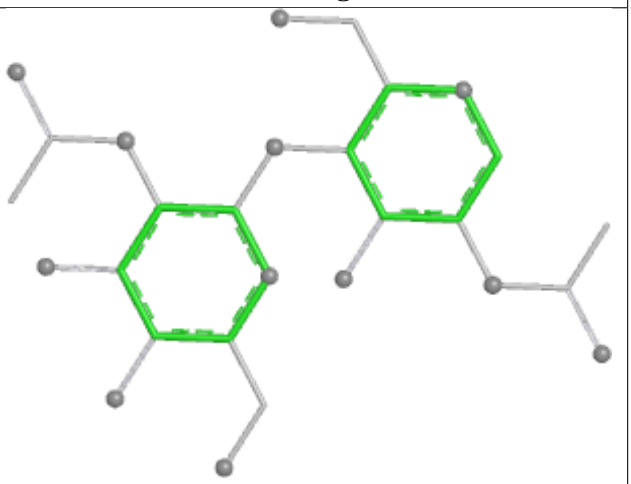
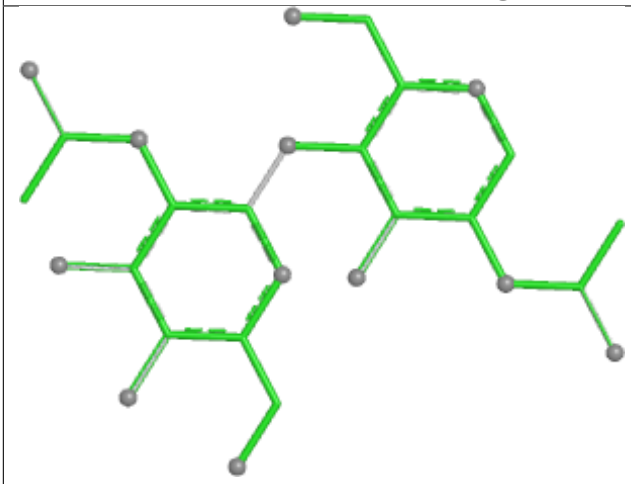


Rings

Oligosaccharide Chain X



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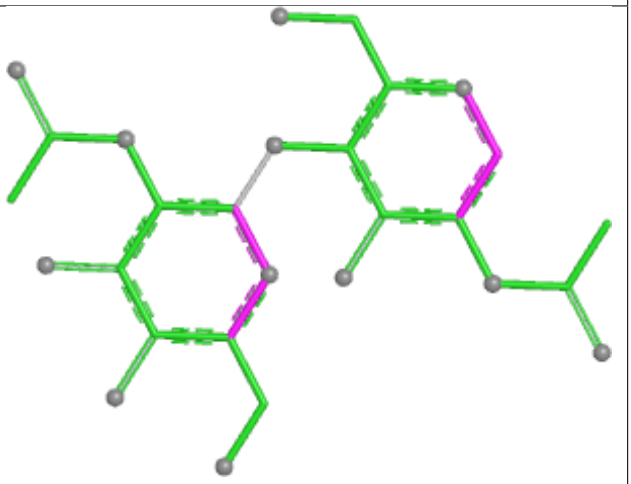

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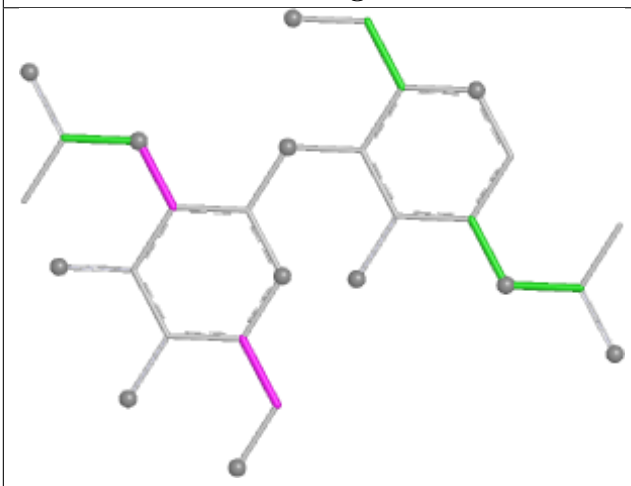

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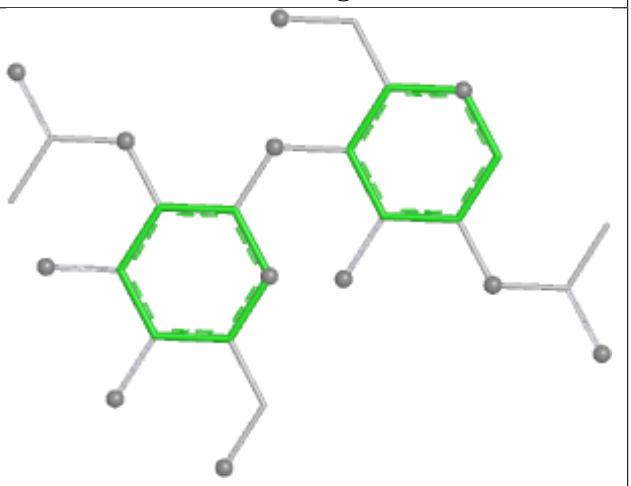
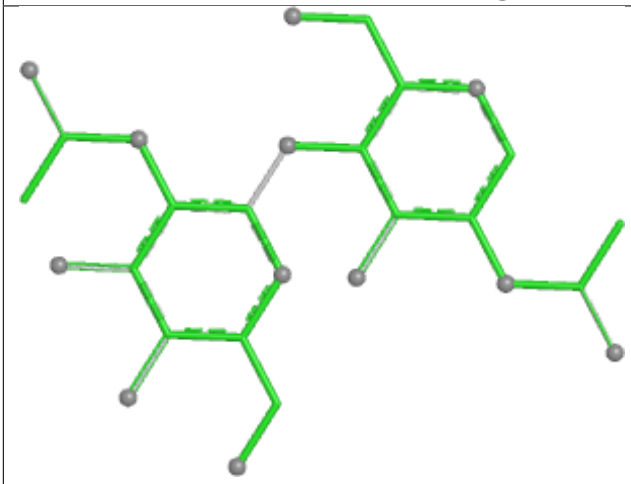


Rings

Oligosaccharide Chain Y



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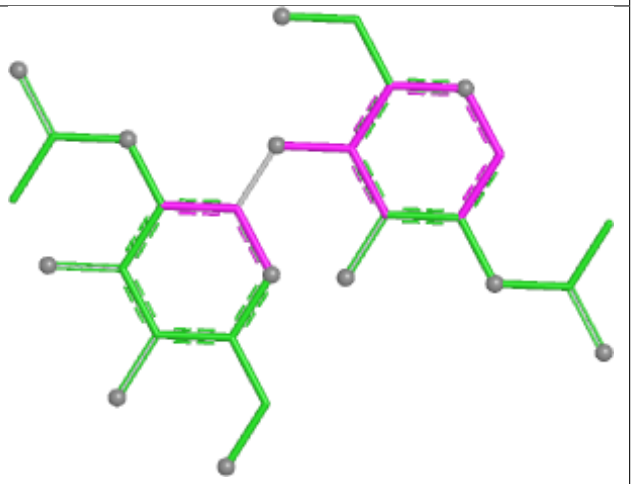

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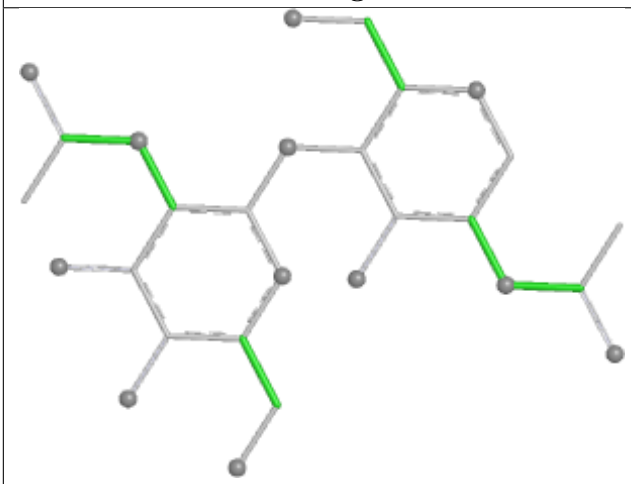

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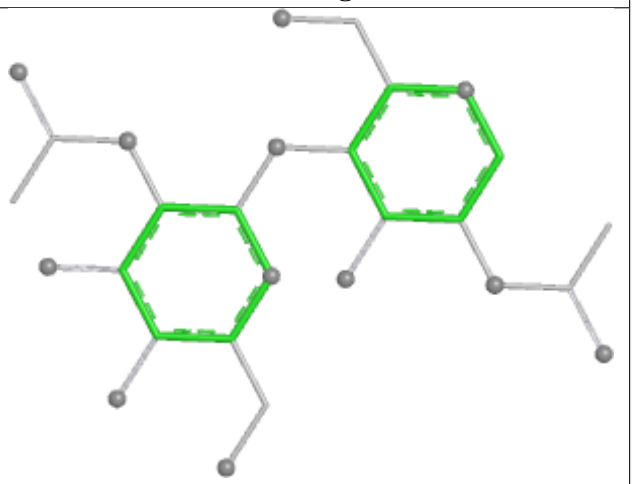
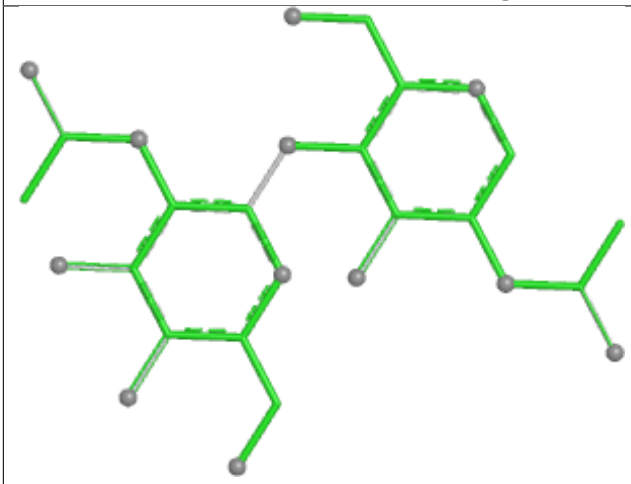


Rings

Oligosaccharide Chain Z



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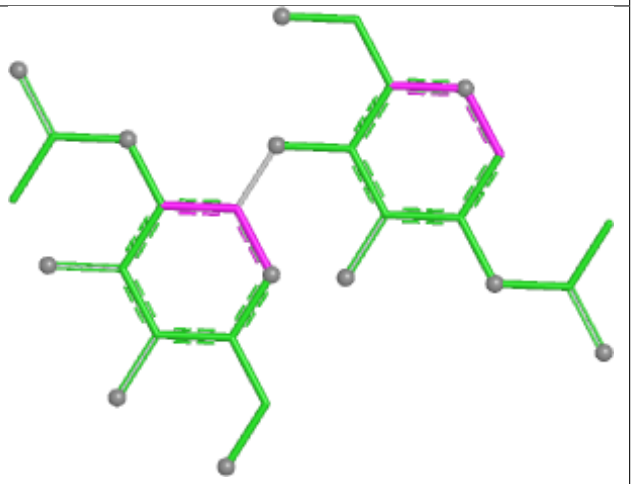
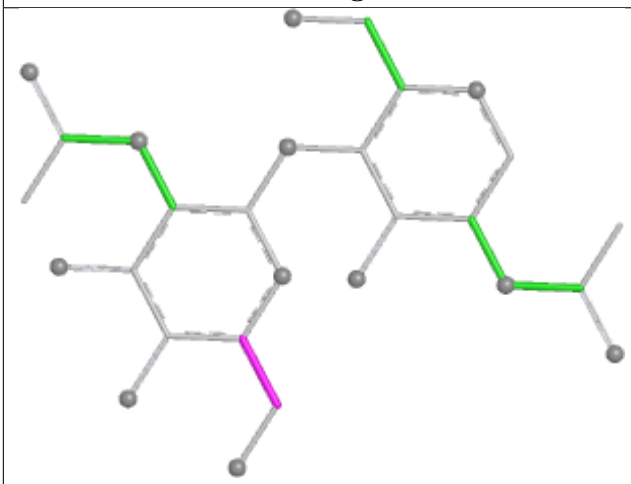


Bond angles



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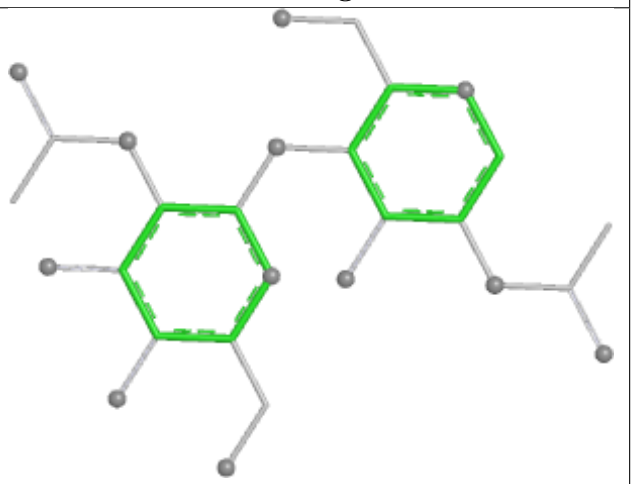
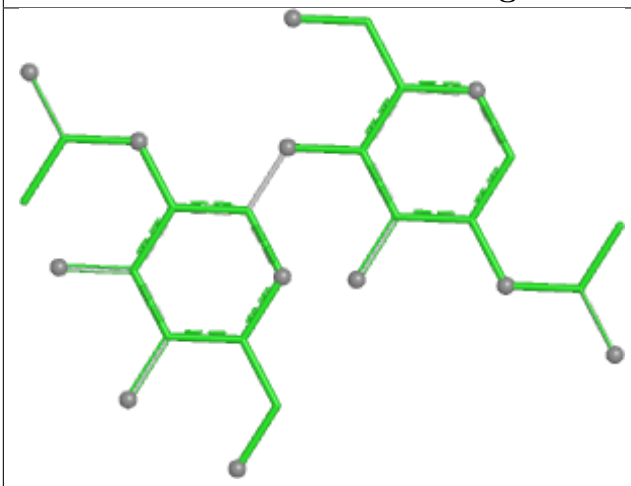


Rings

Oligosaccharide Chain a



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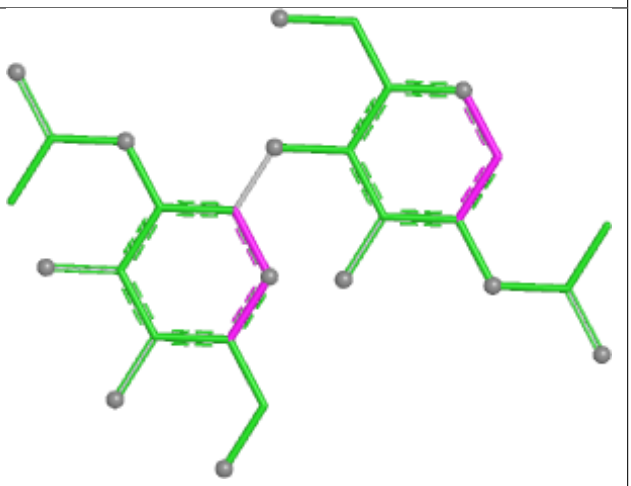
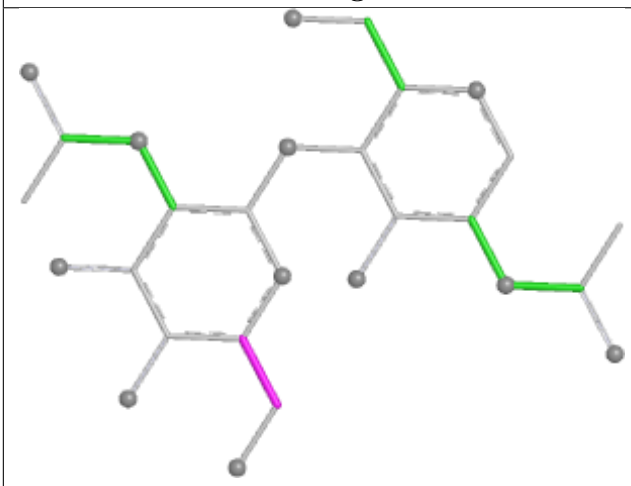


Bond angles



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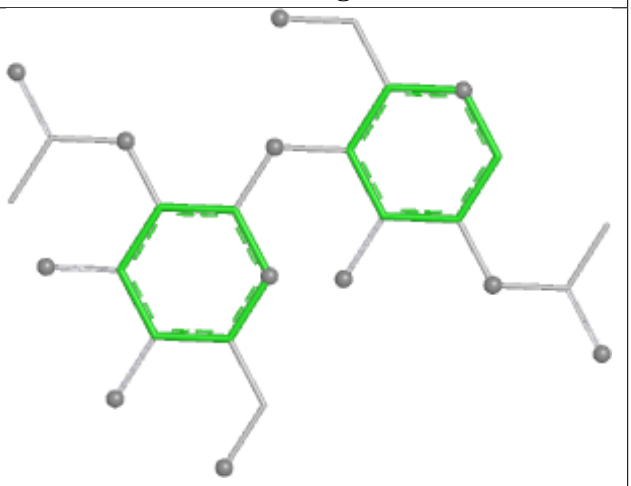
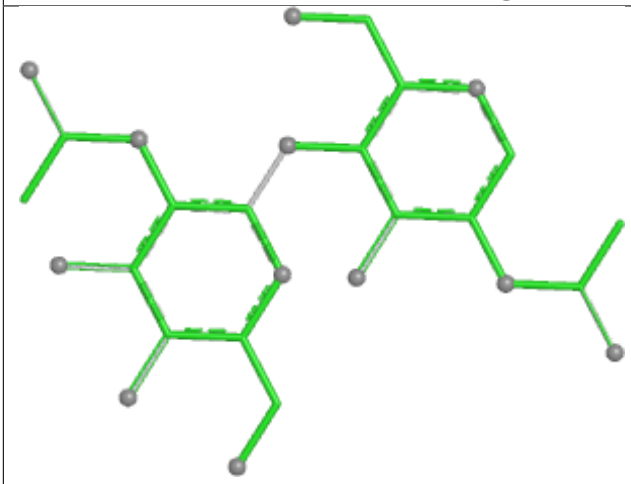
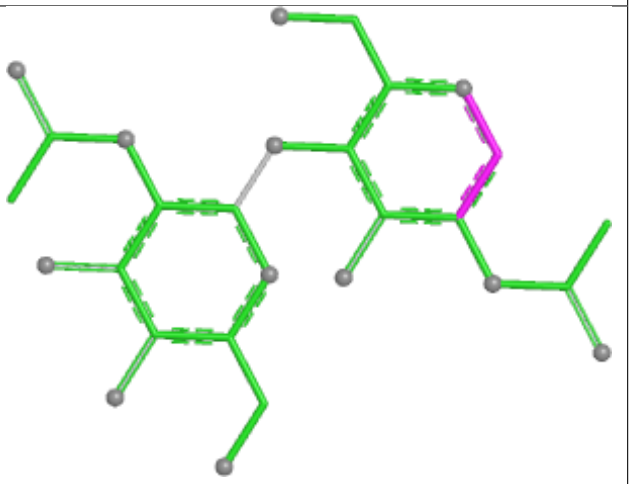


Rings

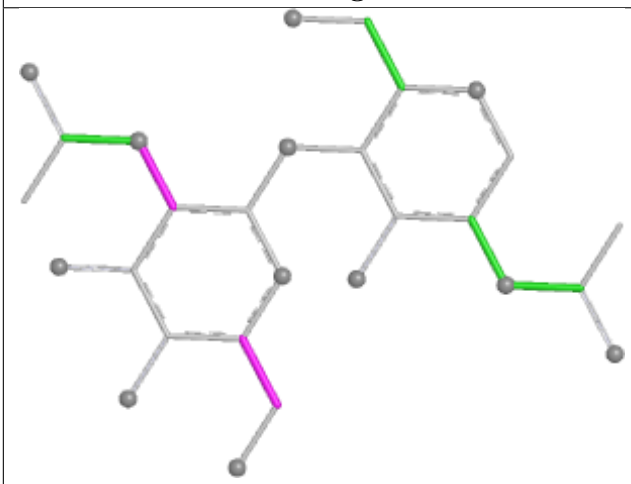
Oligosaccharide Chain h



Bond lengths



Bond angles



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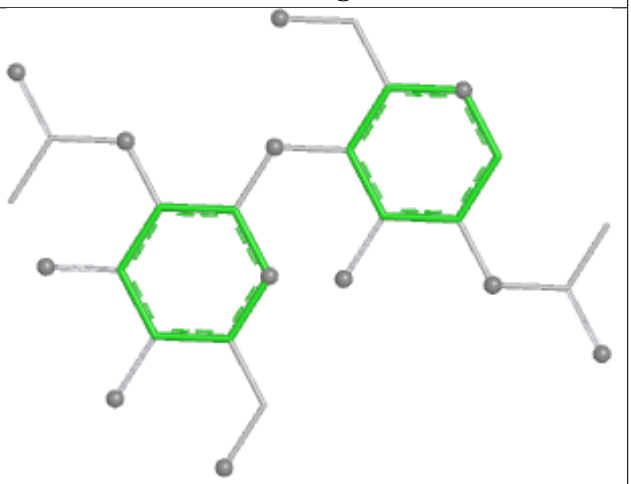
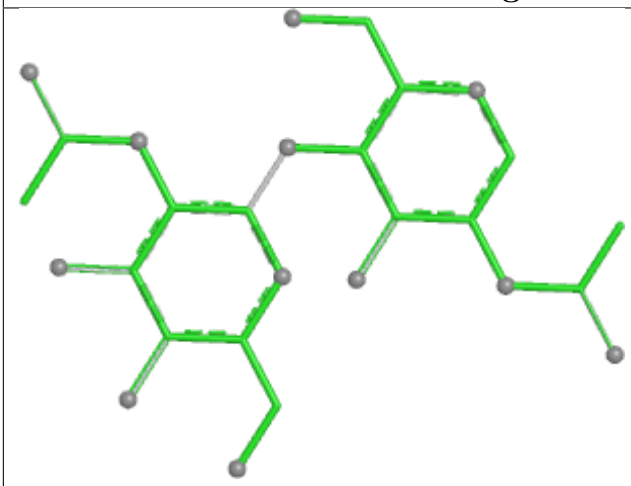
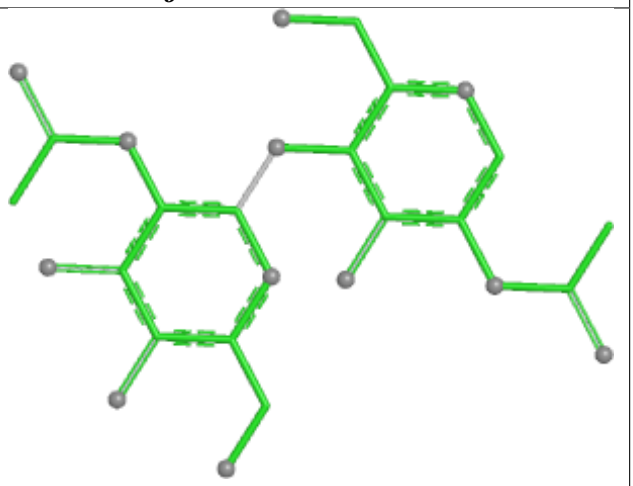


Rings

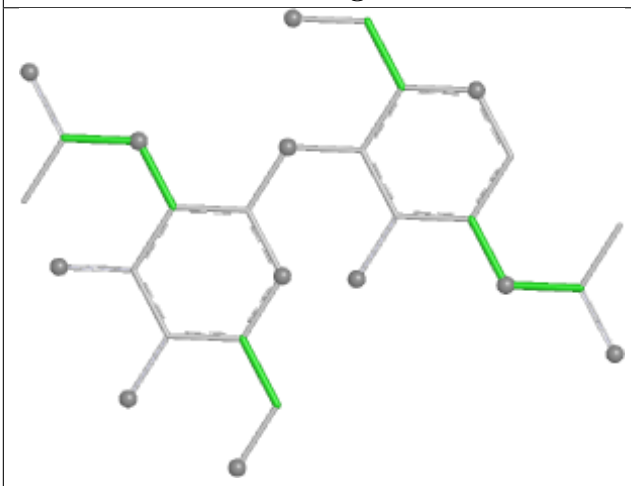
Oligosaccharide Chain j



Bond lengths



Bond angles



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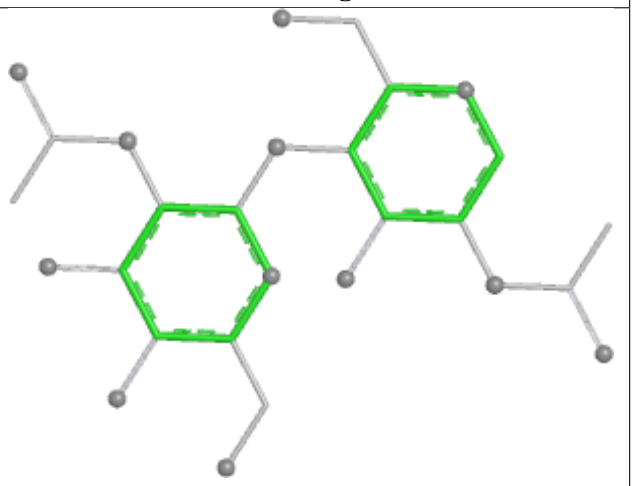
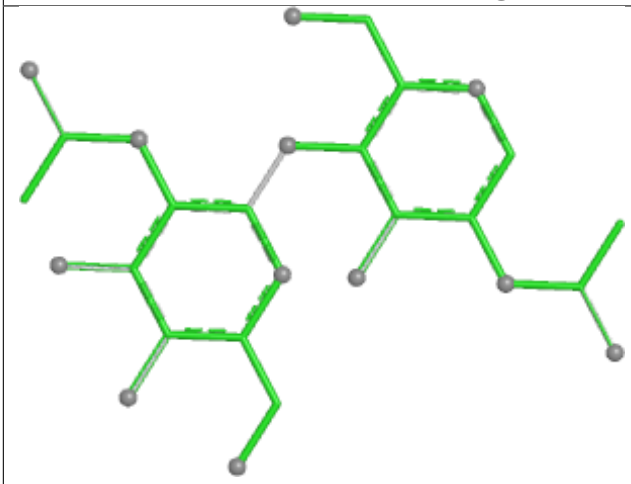
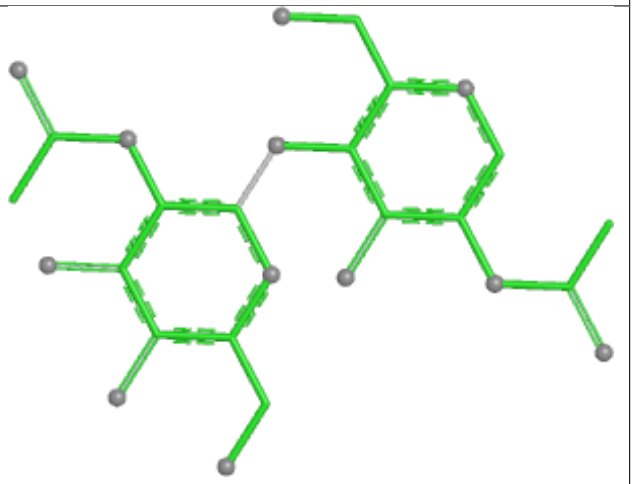


Rings

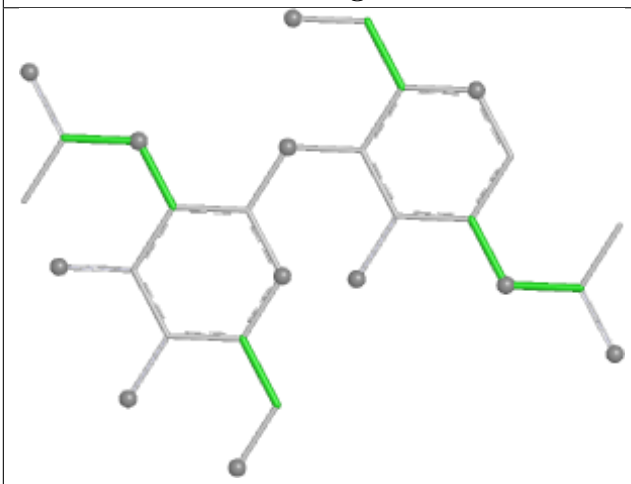
Oligosaccharide Chain m



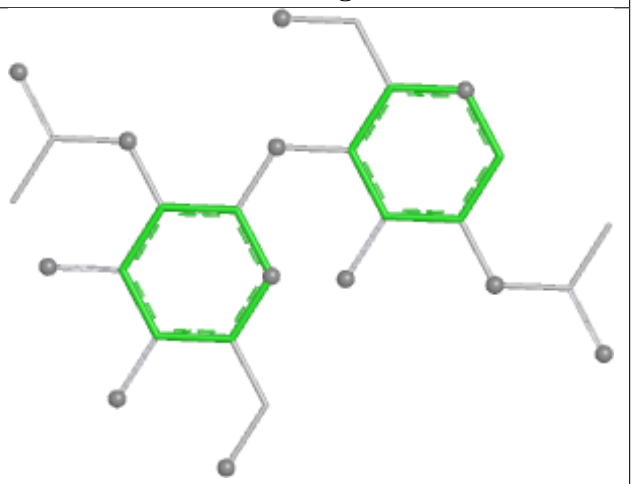
Bond lengths



Bond angles

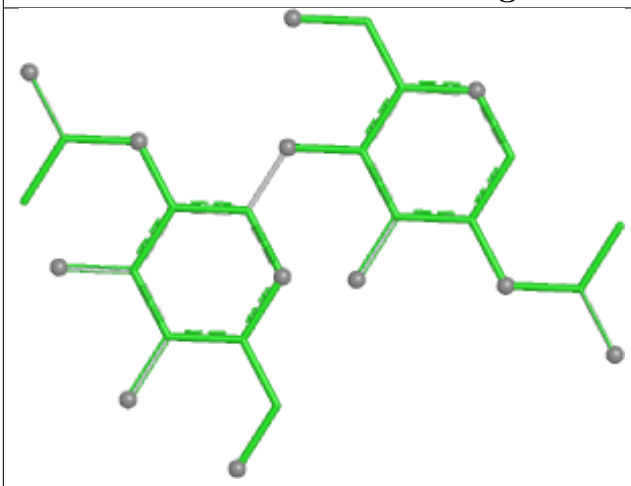


Torsions



Rings

Oligosaccharide Chain o



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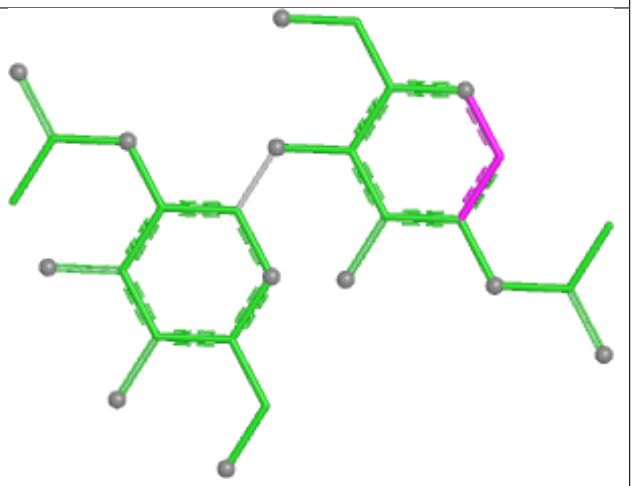

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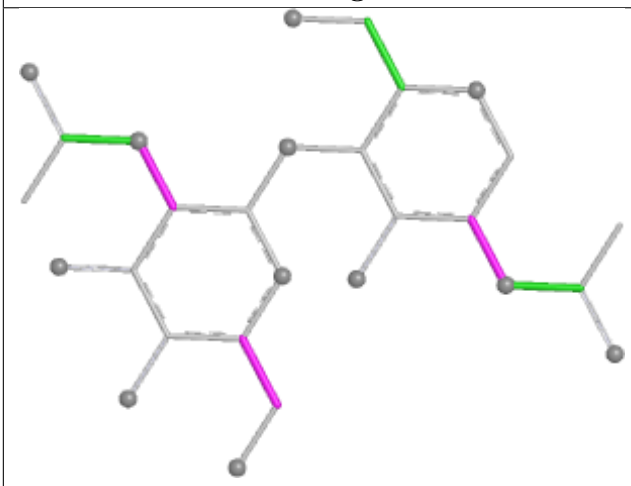

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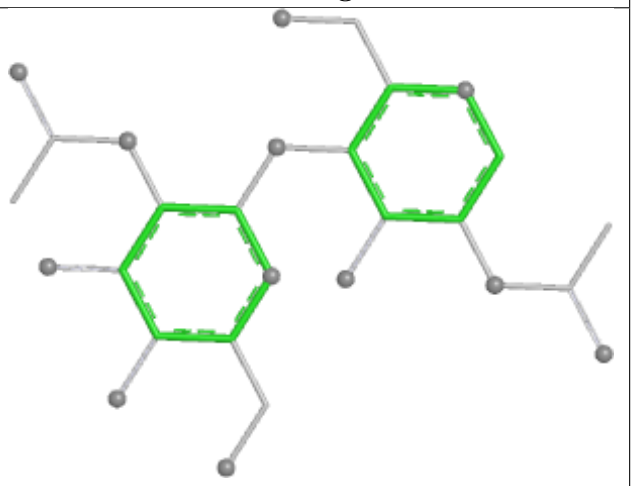
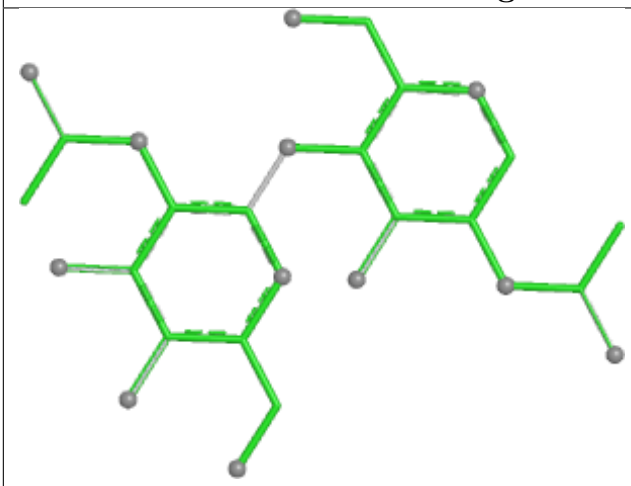


Rings

Oligosaccharide Chain p



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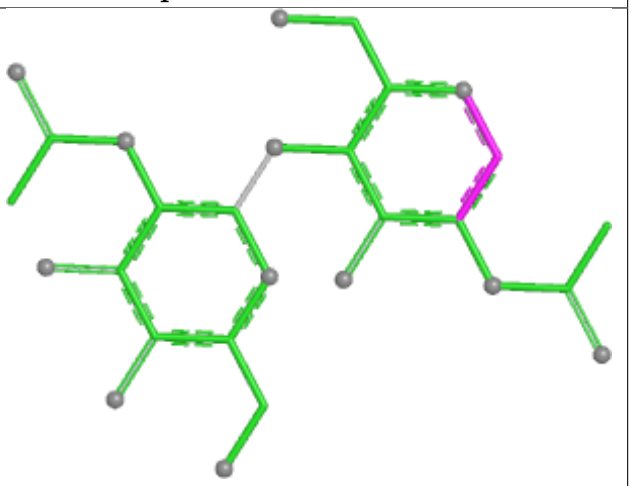

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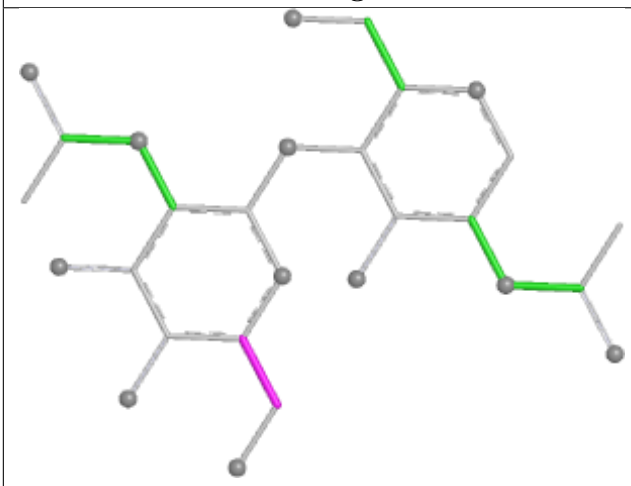

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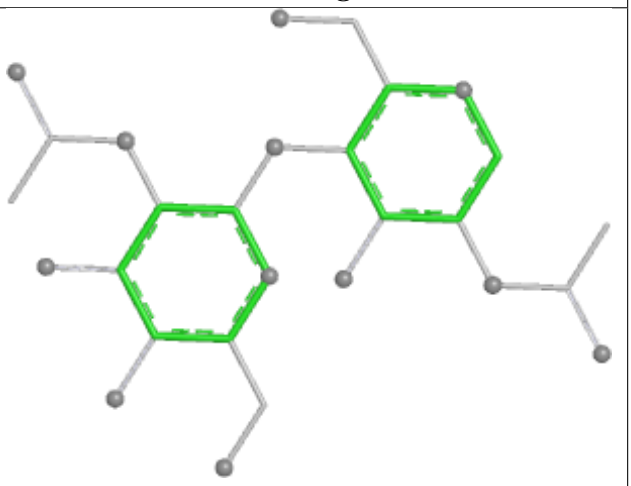
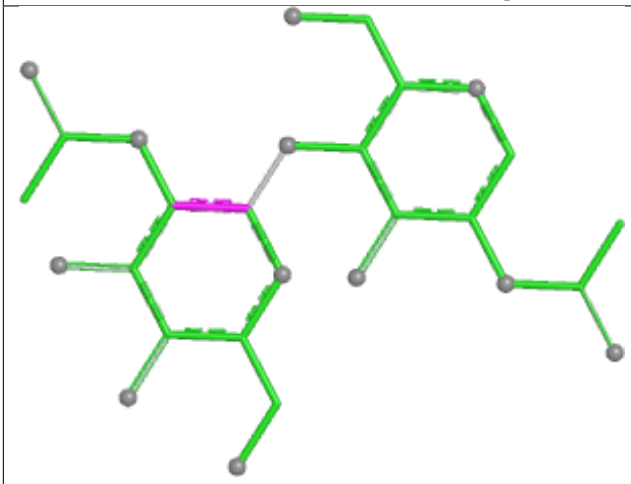


Rings

Oligosaccharide Chain q



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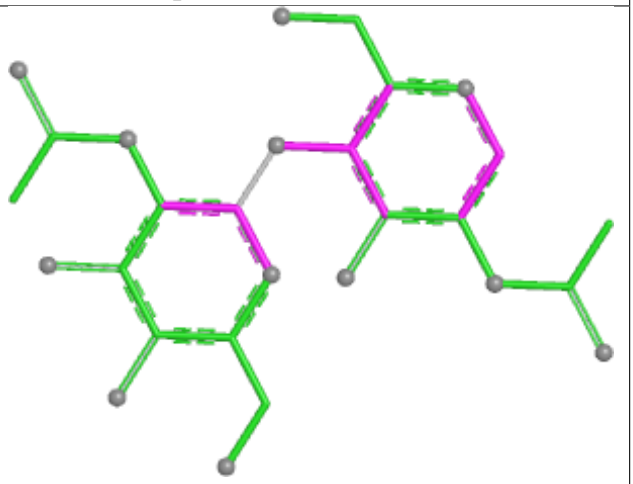

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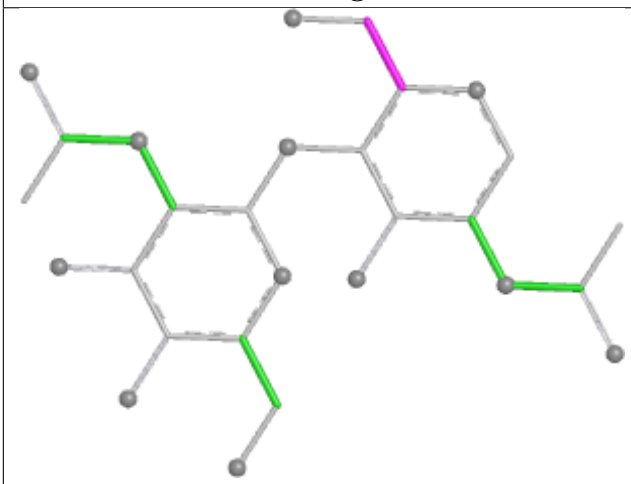

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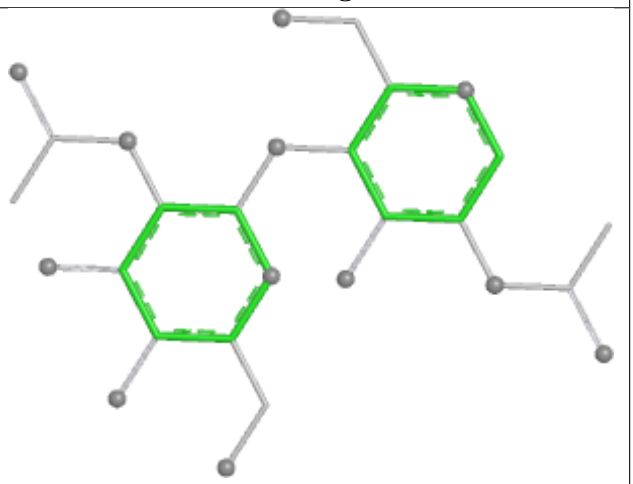
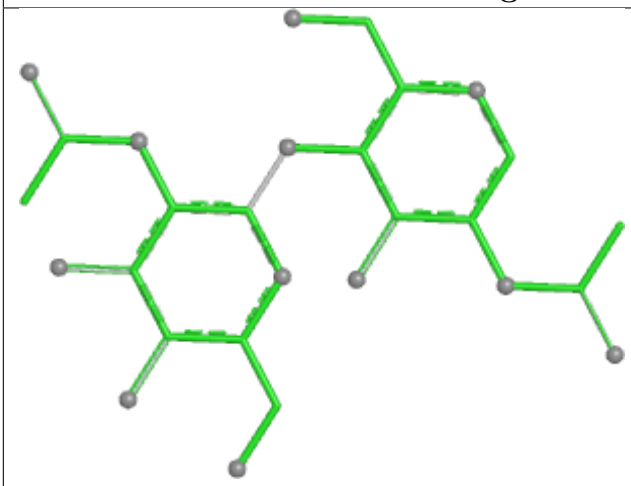


Rings

Oligosaccharide Chain r



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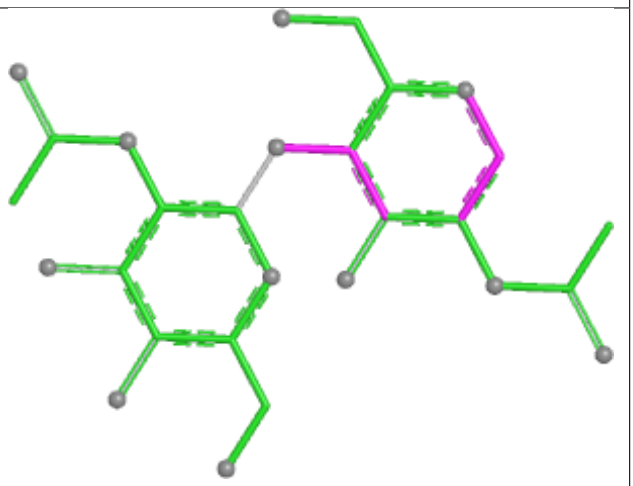

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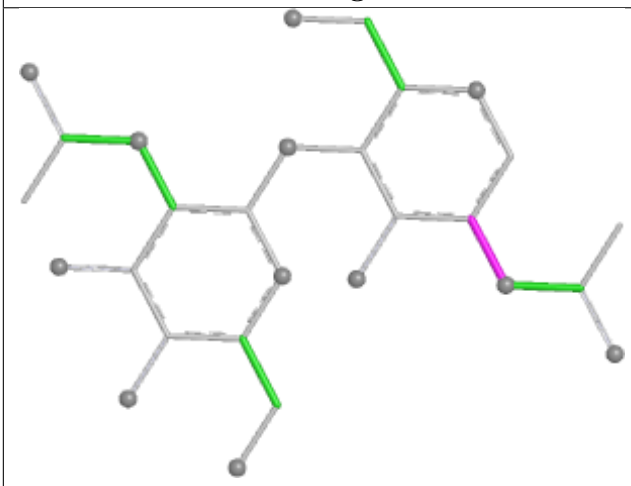

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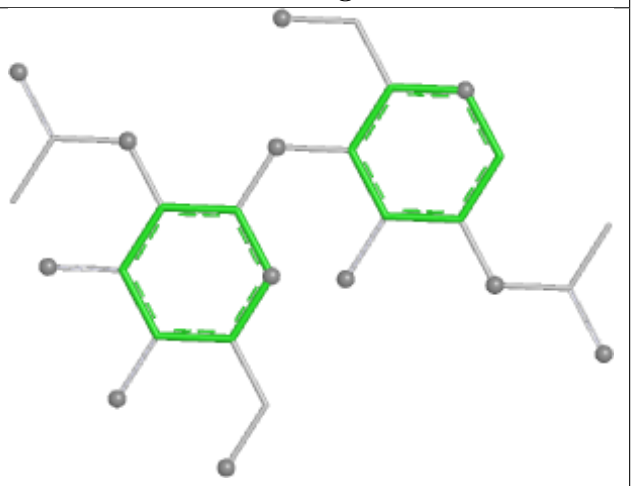
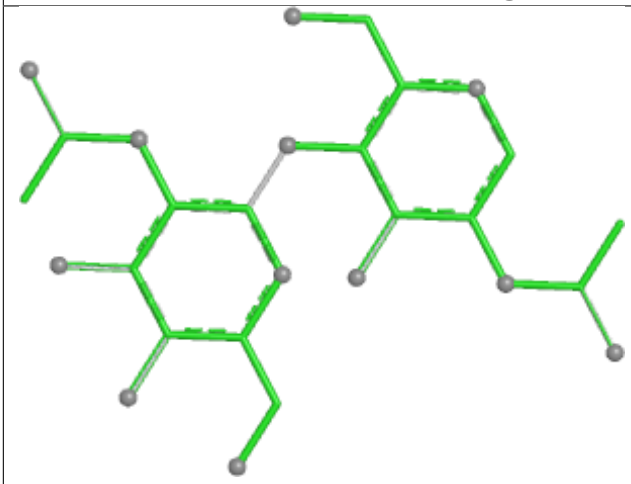


Rings

Oligosaccharide Chain s



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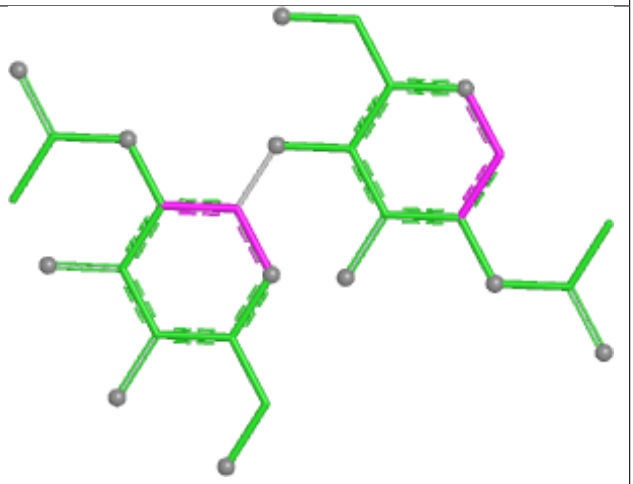

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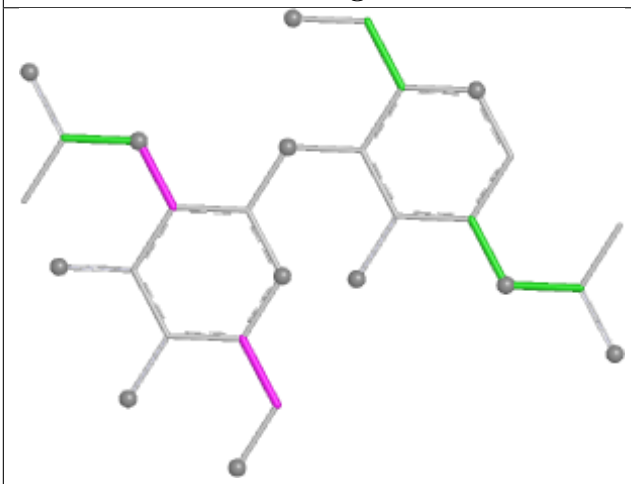

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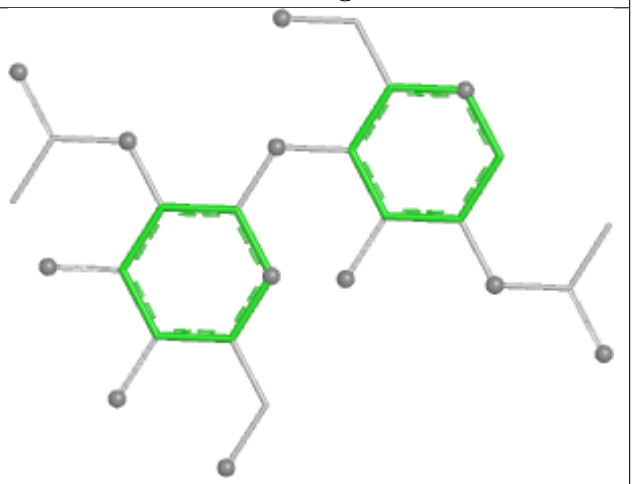
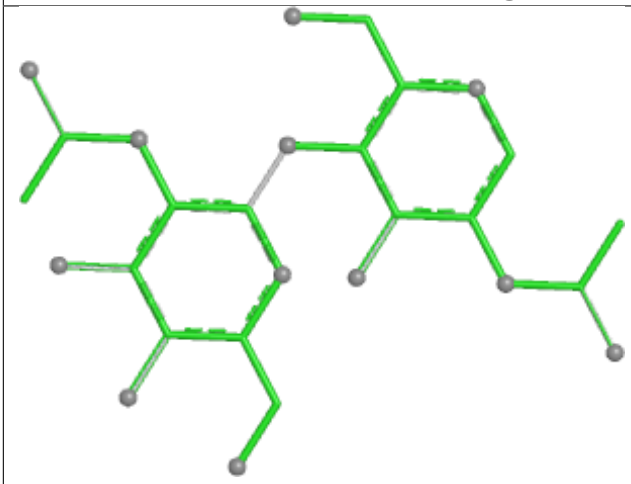


Rings

Oligosaccharide Chain t



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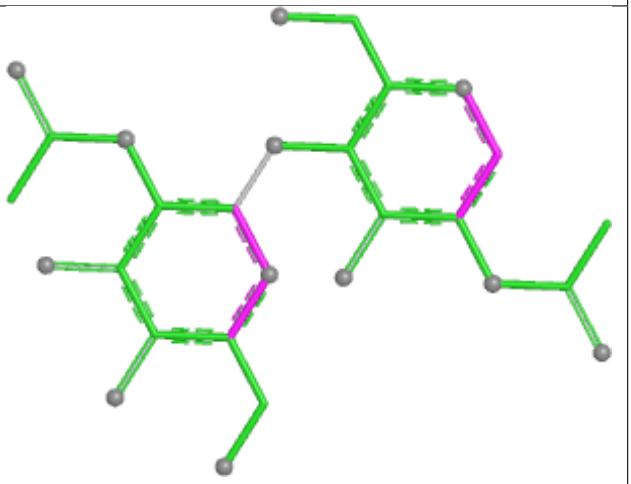

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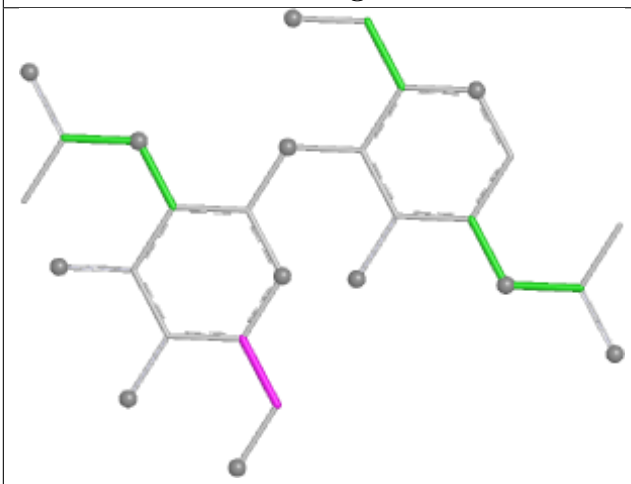

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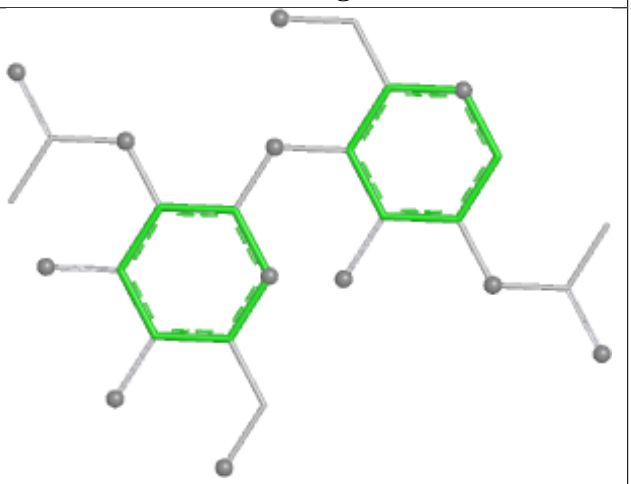
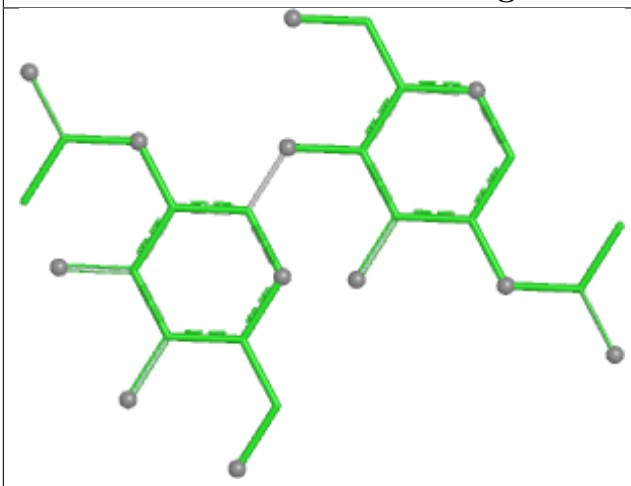


Rings

Oligosaccharide Chain u



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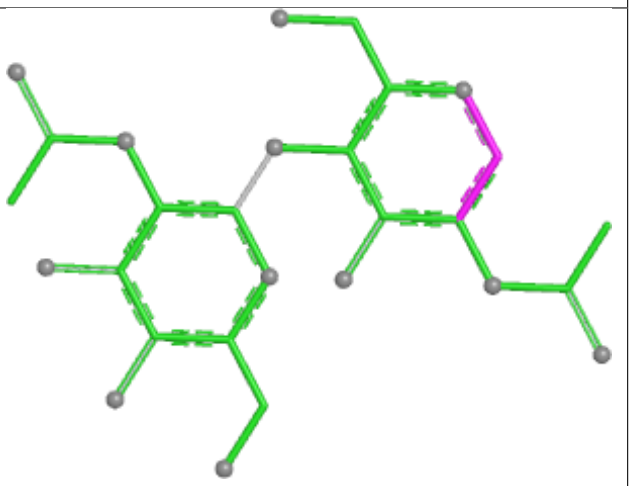

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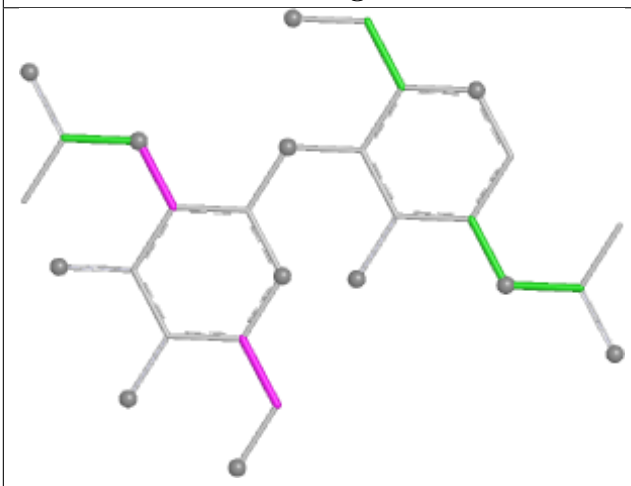

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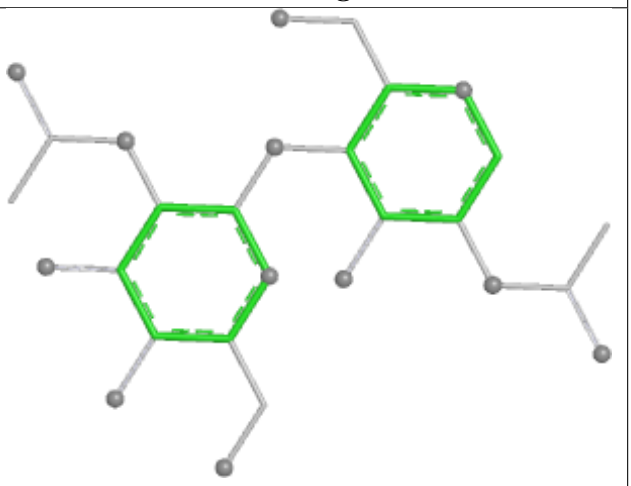
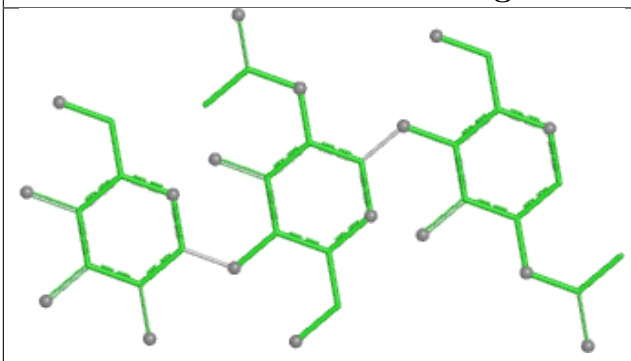


Rings

Oligosaccharide Chain D



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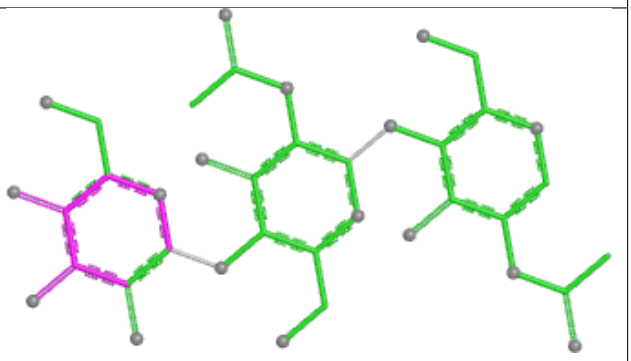

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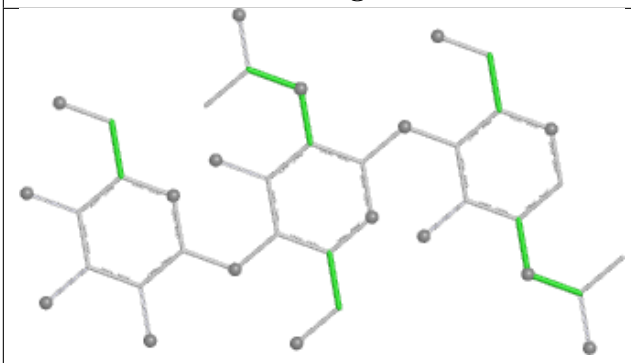

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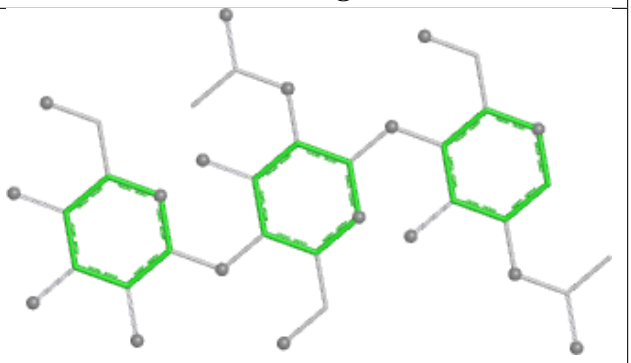
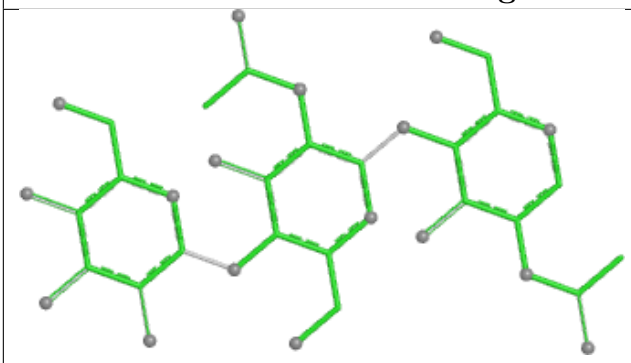


Rings

Oligosaccharide Chain J



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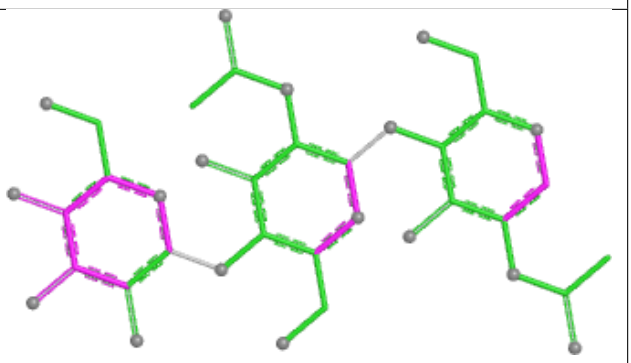

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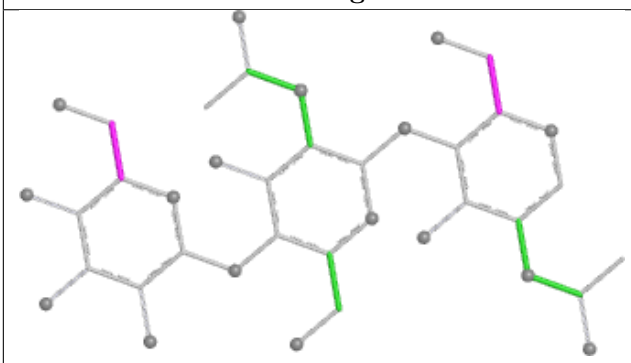

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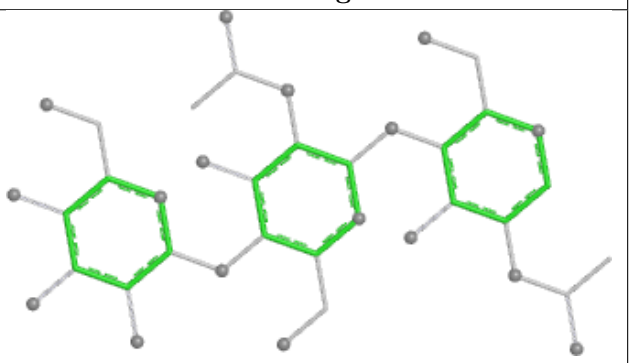
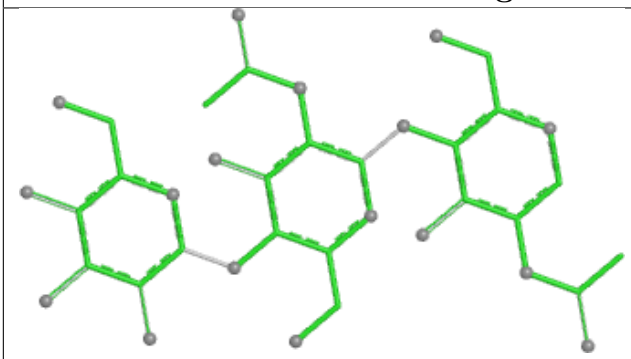


Rings

Oligosaccharide Chain P



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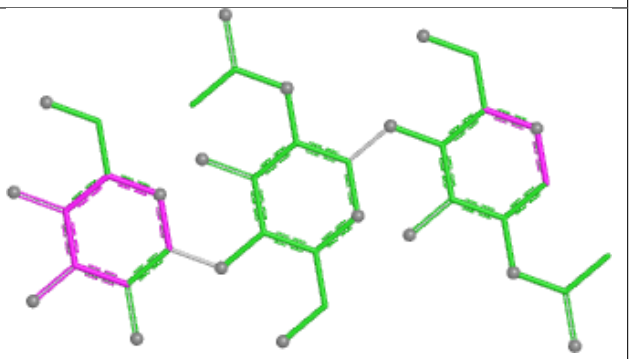

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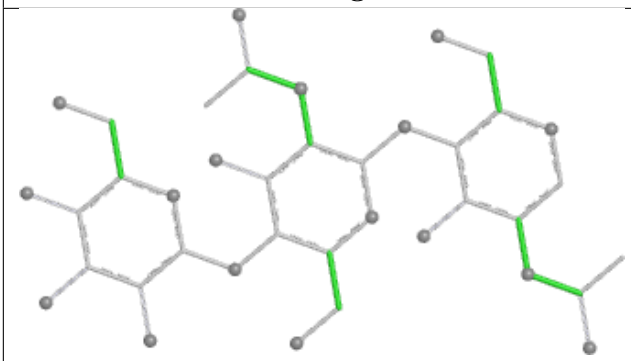

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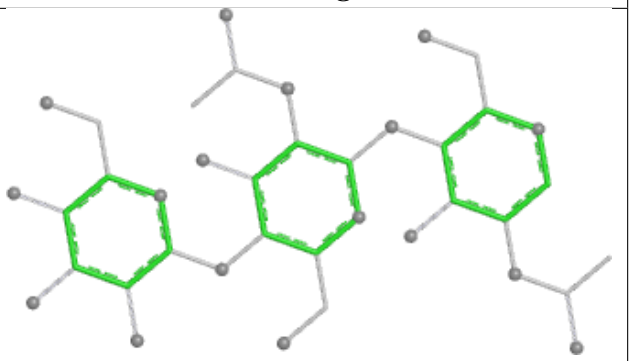
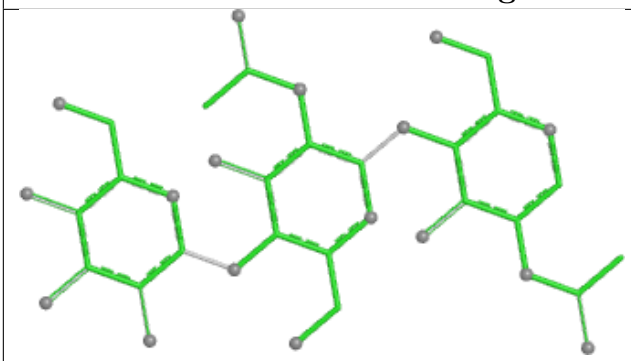


Rings

Oligosaccharide Chain k



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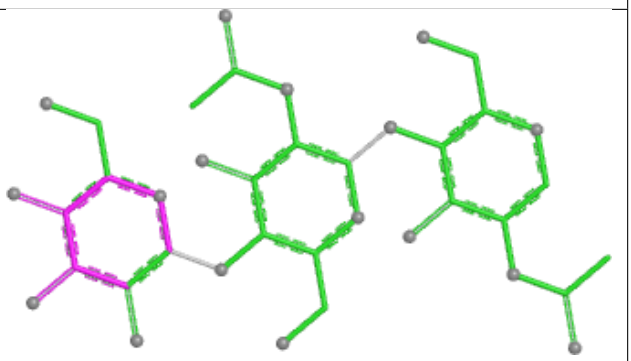
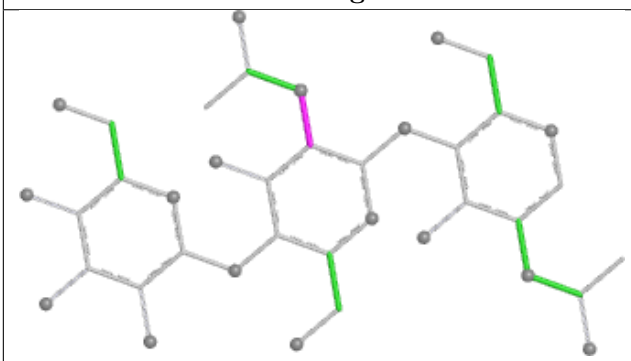


Bond angles



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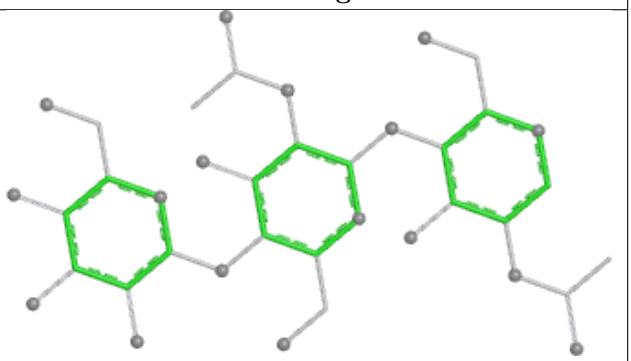
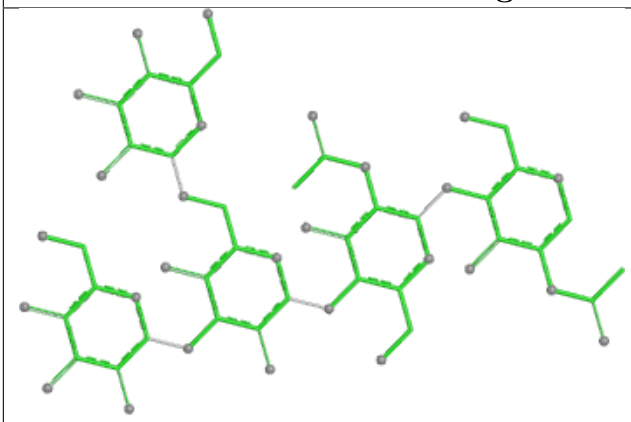


Rings

Oligosaccharide Chain E



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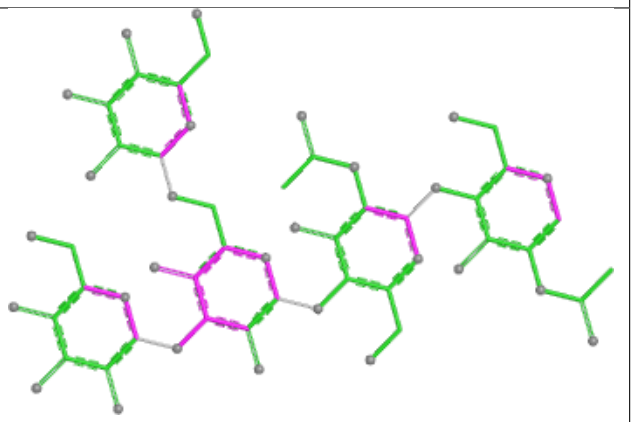
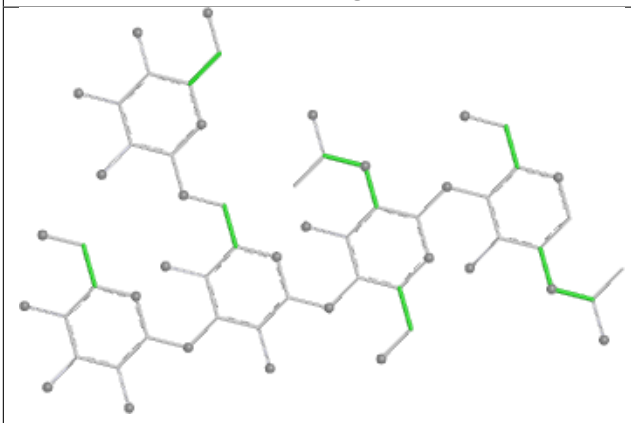


Bond angles



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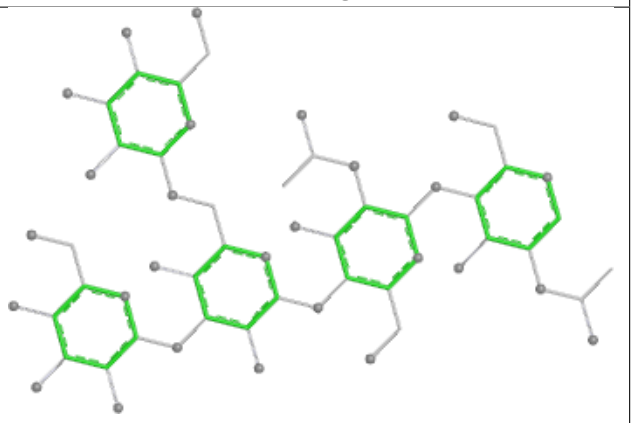
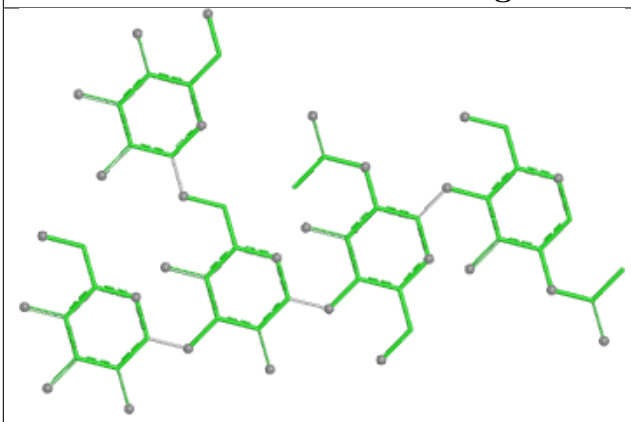


Rings

Oligosaccharide Chain Q



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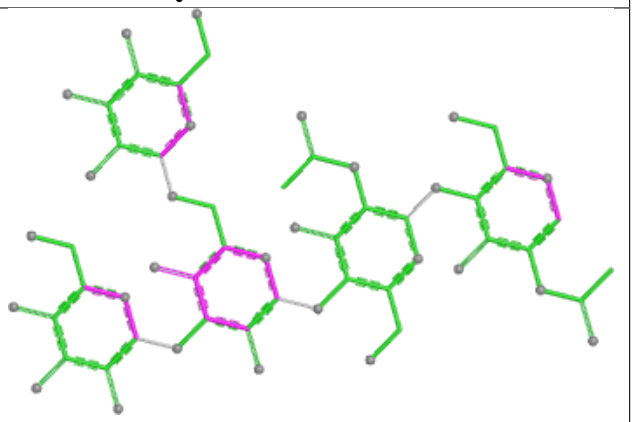
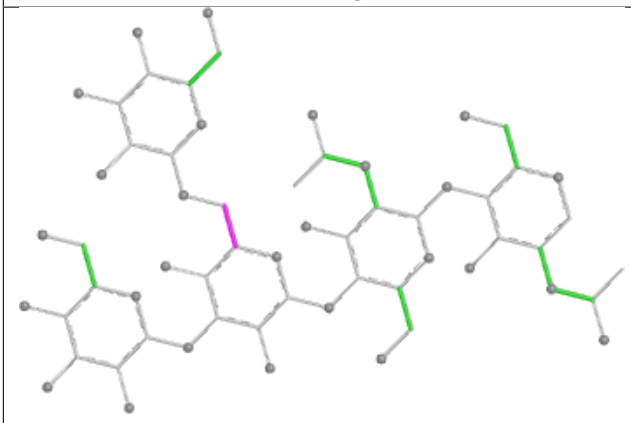


Bond angles



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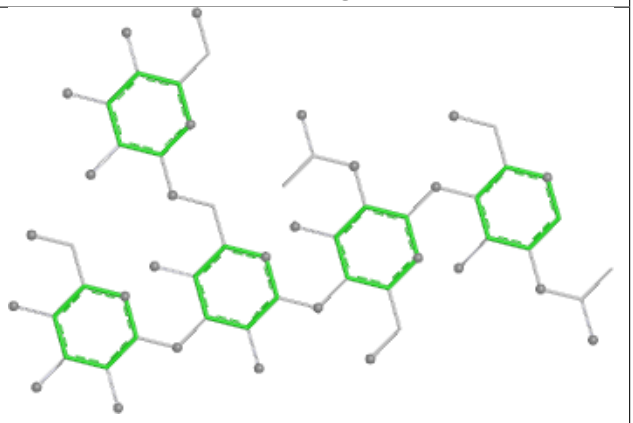
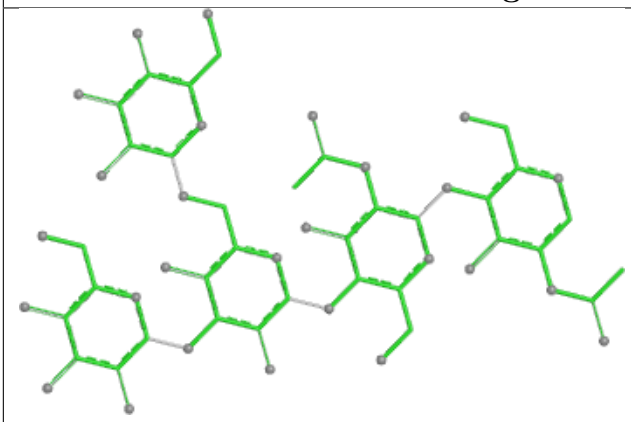


Rings

Oligosaccharide Chain l



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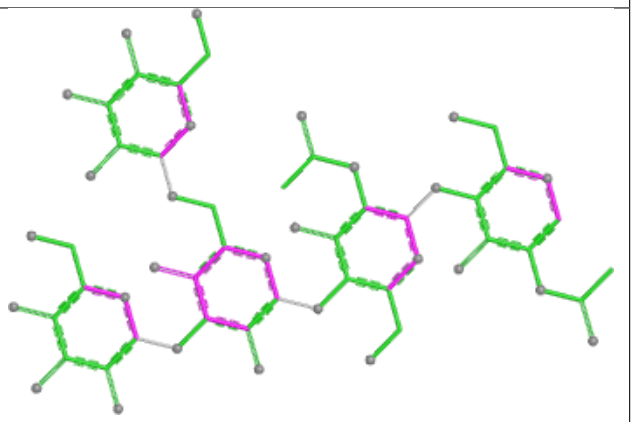
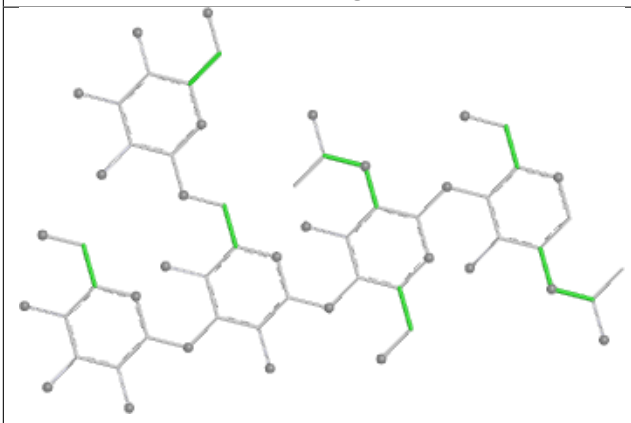
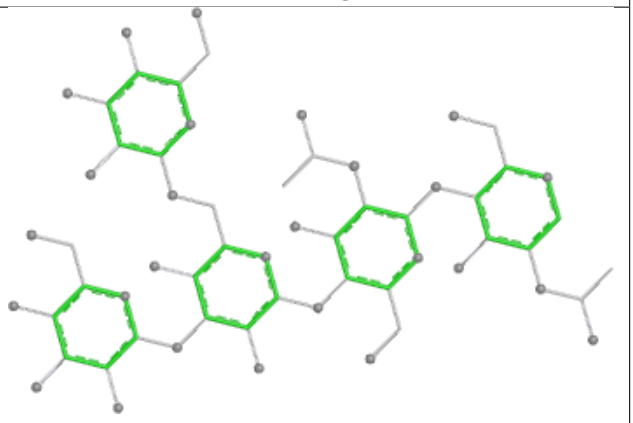


Bond angles

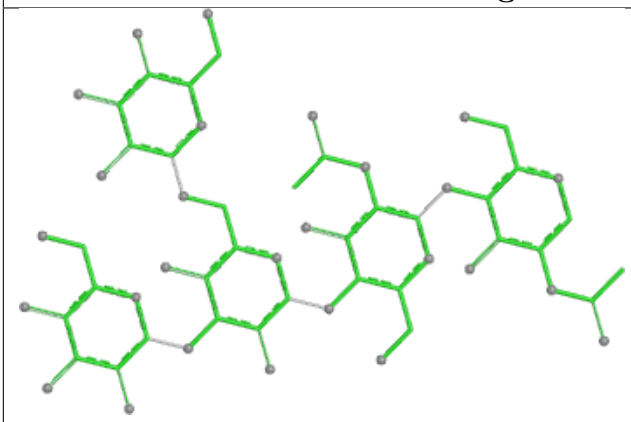


Torsions

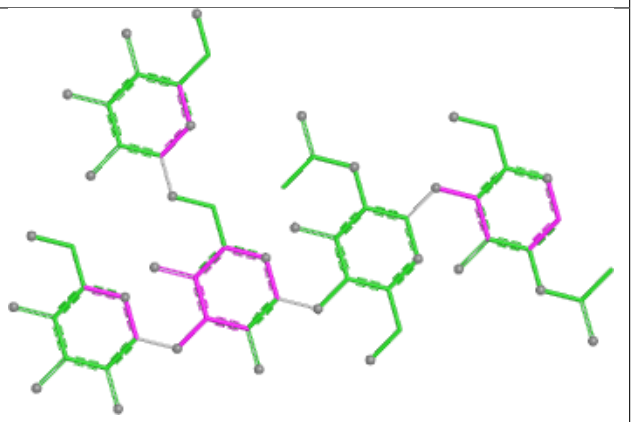


Rings

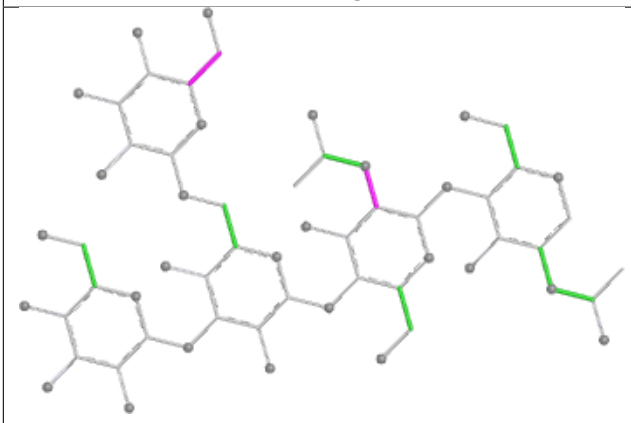
Oligosaccharide Chain n



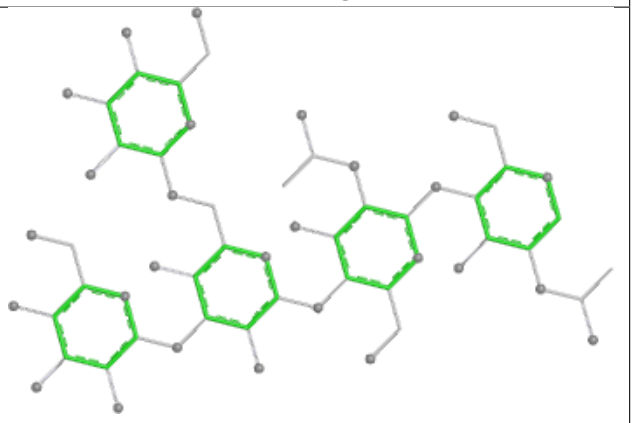
Bond lengths



Bond angles

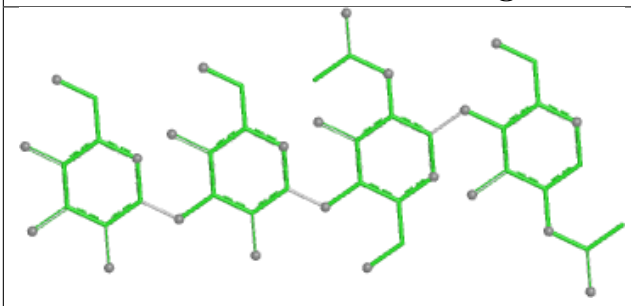


Torsions

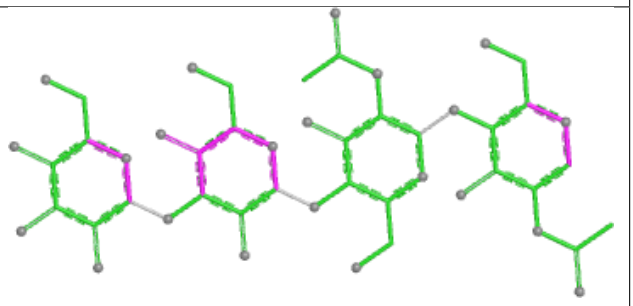


Rings

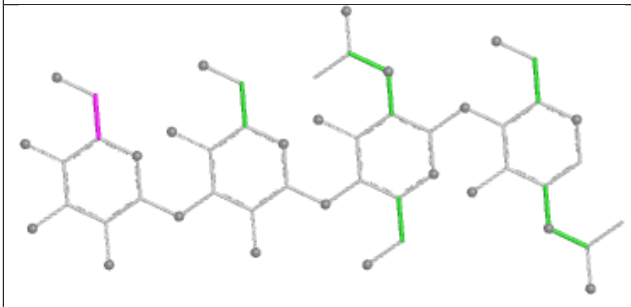
Oligosaccharide Chain S



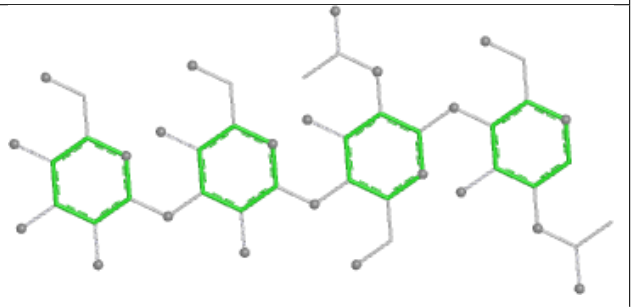
Bond lengths



Bond angles

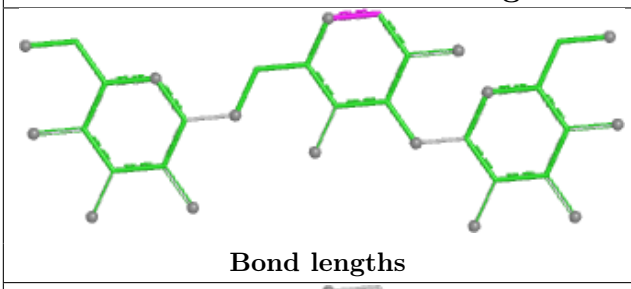


Torsions

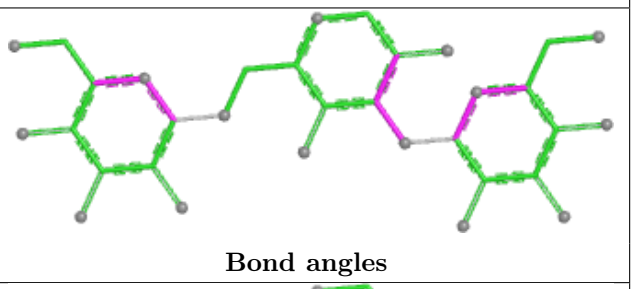


Rings

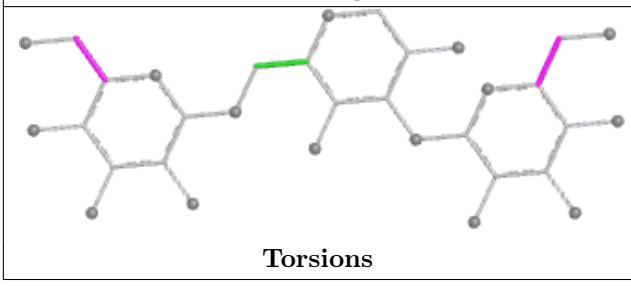
Oligosaccharide Chain i



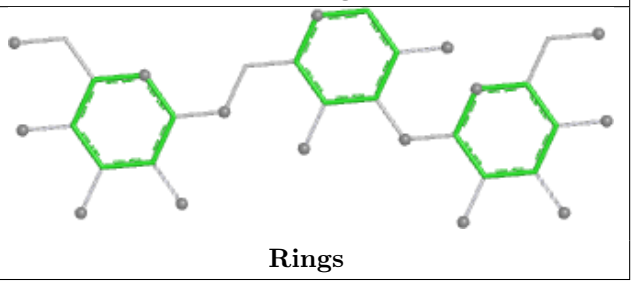
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

31 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	$\#\|Z\| > 2$
10	NAG	c	701	4	14,14,15	0.69	0	17,19,21	0.90	0
10	NAG	b	601	3	14,14,15	0.67	0	17,19,21	0.92	0
10	NAG	b	606	3	14,14,15	0.71	0	17,19,21	0.91	0
10	NAG	c	702	4	14,14,15	0.70	0	17,19,21	0.86	0
10	NAG	d	606	-	14,14,15	0.71	0	17,19,21	0.78	0
10	NAG	f	603	3	14,14,15	0.74	0	17,19,21	1.00	0
10	NAG	f	605	3	14,14,15	0.73	0	17,19,21	0.79	0
10	NAG	d	603	3	14,14,15	0.66	0	17,19,21	0.84	0
10	NAG	f	608	3	14,14,15	0.69	0	17,19,21	1.00	1 (5%)
10	NAG	b	604	3	14,14,15	0.72	0	17,19,21	1.04	1 (5%)
10	NAG	e	701	4	14,14,15	0.67	0	17,19,21	0.88	0
10	NAG	d	601	3	14,14,15	0.64	0	17,19,21	0.89	0
10	NAG	f	602	3	14,14,15	0.65	0	17,19,21	0.88	0
10	NAG	f	601	3	14,14,15	0.68	0	17,19,21	0.88	0
10	NAG	f	606	-	14,14,15	0.73	0	17,19,21	0.79	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
10	NAG	d	602	3	14,14,15	0.69	0	17,19,21	0.82	0
10	NAG	b	605	3	14,14,15	0.72	0	17,19,21	0.88	0
10	NAG	c	703	4	14,14,15	0.69	0	17,19,21	0.79	0
10	NAG	e	702	4	14,14,15	0.72	0	17,19,21	0.82	0
10	NAG	f	604	3	14,14,15	0.67	0	17,19,21	0.82	0
10	NAG	e	703	4	14,14,15	0.70	0	17,19,21	0.85	0
10	NAG	f	607	3	14,14,15	0.72	0	17,19,21	1.09	1 (5%)
10	NAG	g	701	4	14,14,15	0.67	0	17,19,21	0.88	0
10	NAG	b	608	-	14,14,15	0.70	0	17,19,21	0.95	1 (5%)
10	NAG	d	604	3	14,14,15	0.72	0	17,19,21	0.86	0
10	NAG	g	702	4	14,14,15	0.72	0	17,19,21	0.83	0
10	NAG	g	703	4	14,14,15	0.71	0	17,19,21	0.84	0
10	NAG	b	603	3	14,14,15	0.77	0	17,19,21	1.76	2 (11%)
10	NAG	b	607	3	14,14,15	0.73	0	17,19,21	0.91	0
10	NAG	d	605	-	14,14,15	0.75	0	17,19,21	0.72	0
10	NAG	b	602	3	14,14,15	0.69	0	17,19,21	0.84	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
10	NAG	c	701	4	-	0/6/23/26	0/1/1/1
10	NAG	b	601	3	-	0/6/23/26	0/1/1/1
10	NAG	b	606	3	-	0/6/23/26	0/1/1/1
10	NAG	c	702	4	-	0/6/23/26	0/1/1/1
10	NAG	d	606	-	-	0/6/23/26	0/1/1/1
10	NAG	f	603	3	-	2/6/23/26	0/1/1/1
10	NAG	f	605	3	-	0/6/23/26	0/1/1/1
10	NAG	d	603	3	-	0/6/23/26	0/1/1/1
10	NAG	f	608	3	-	0/6/23/26	0/1/1/1
10	NAG	b	604	3	-	0/6/23/26	0/1/1/1
10	NAG	e	701	4	-	0/6/23/26	0/1/1/1
10	NAG	d	601	3	-	0/6/23/26	0/1/1/1
10	NAG	f	602	3	-	0/6/23/26	0/1/1/1
10	NAG	f	601	3	-	0/6/23/26	0/1/1/1
10	NAG	f	606	-	-	0/6/23/26	0/1/1/1
10	NAG	d	602	3	-	0/6/23/26	0/1/1/1
10	NAG	b	605	3	-	0/6/23/26	0/1/1/1
10	NAG	c	703	4	-	0/6/23/26	0/1/1/1
10	NAG	e	702	4	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
10	NAG	f	604	3	-	0/6/23/26	0/1/1/1
10	NAG	e	703	4	-	0/6/23/26	0/1/1/1
10	NAG	f	607	3	-	0/6/23/26	0/1/1/1
10	NAG	g	701	4	-	0/6/23/26	0/1/1/1
10	NAG	b	608	-	-	2/6/23/26	0/1/1/1
10	NAG	d	604	3	-	0/6/23/26	0/1/1/1
10	NAG	g	702	4	-	0/6/23/26	0/1/1/1
10	NAG	g	703	4	-	0/6/23/26	0/1/1/1
10	NAG	b	603	3	-	2/6/23/26	0/1/1/1
10	NAG	b	607	3	-	1/6/23/26	0/1/1/1
10	NAG	d	605	-	-	0/6/23/26	0/1/1/1
10	NAG	b	602	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
10	b	603	NAG	C1-O5-C5	5.61	119.71	112.19
10	f	607	NAG	C1-O5-C5	2.97	116.16	112.19
10	b	604	NAG	C1-O5-C5	2.74	115.85	112.19
10	b	603	NAG	C1-C2-N2	2.46	114.31	110.43
10	b	608	NAG	C2-N2-C7	2.16	125.80	122.90

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	b	607	NAG	O5-C5-C6-O6
10	b	603	NAG	C1-C2-N2-C7
10	b	608	NAG	C1-C2-N2-C7
10	f	603	NAG	C1-C2-N2-C7
10	b	603	NAG	C3-C2-N2-C7

There are no ring outliers.

3 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
10	c	702	NAG	1	0
10	f	603	NAG	1	0
10	e	702	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.

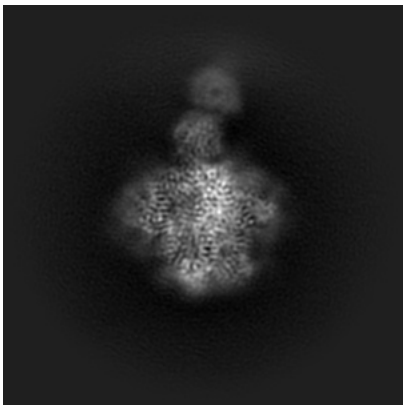
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-74654. These allow visual inspection of the internal detail of the map and identification of artifacts.

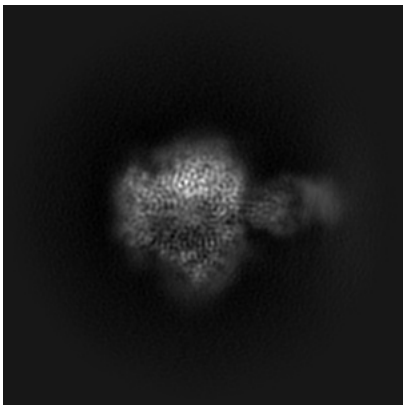
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

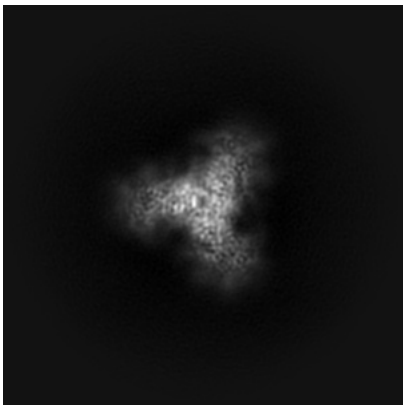
6.1.1 Primary map



X

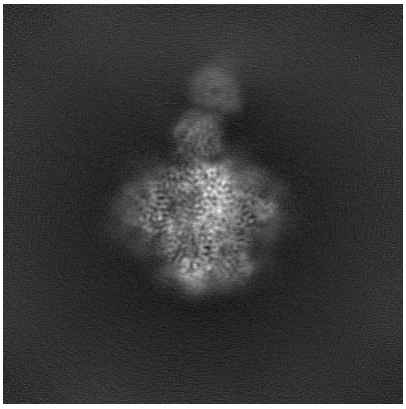


Y

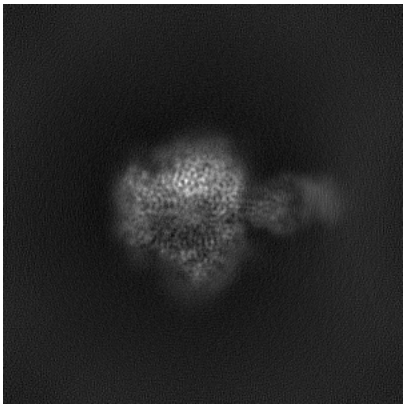


Z

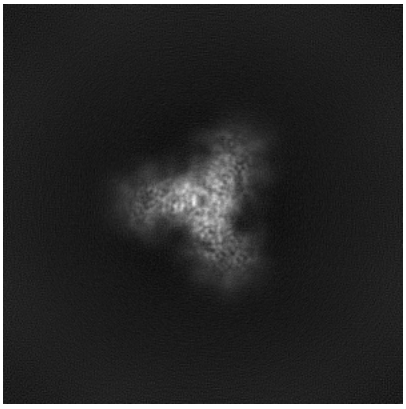
6.1.2 Raw map



X



Y

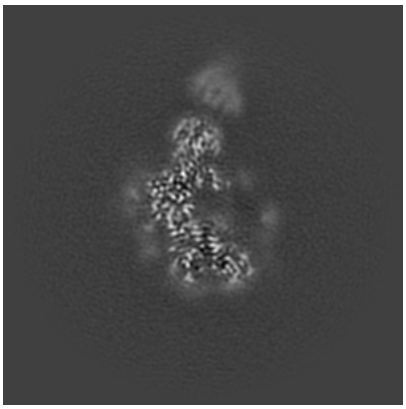


Z

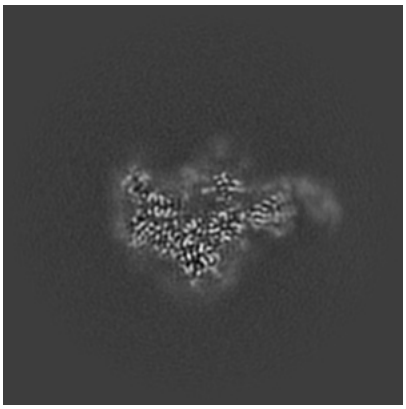
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

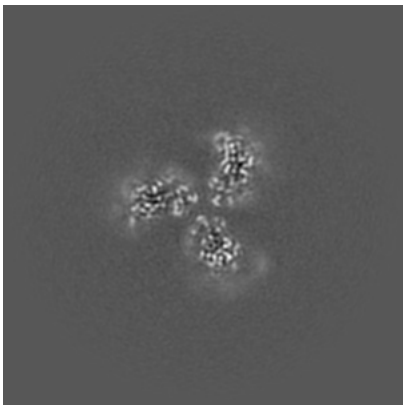
6.2.1 Primary map



X Index: 192

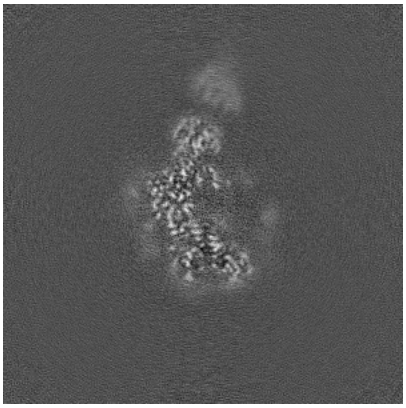


Y Index: 192

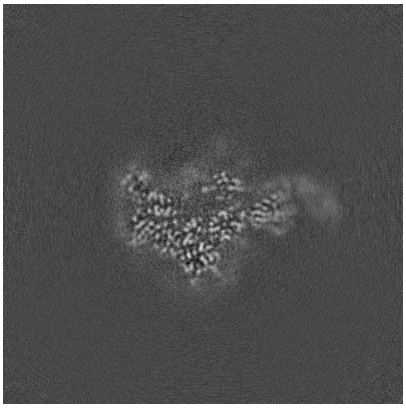


Z Index: 192

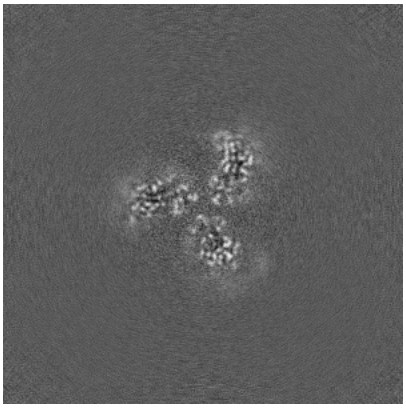
6.2.2 Raw map



X Index: 192



Y Index: 192

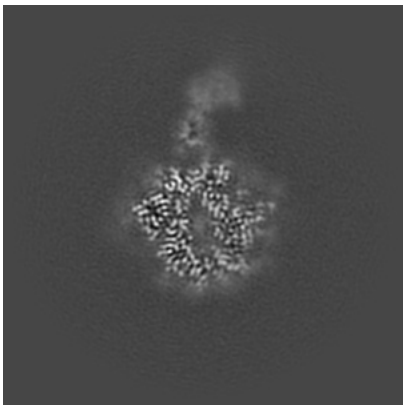


Z Index: 192

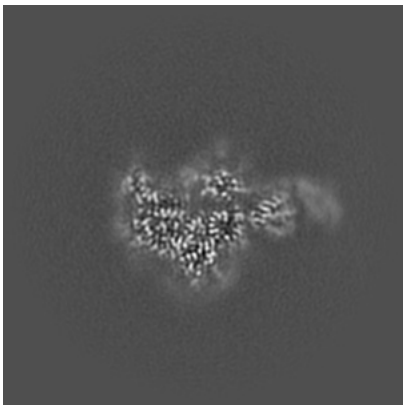
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

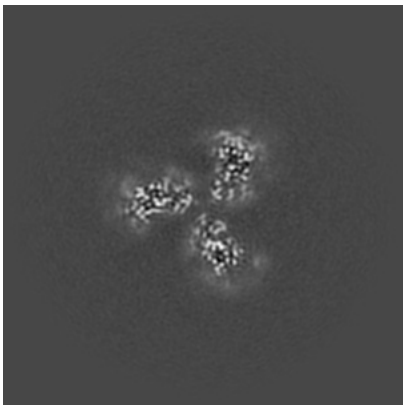
6.3.1 Primary map



X Index: 206

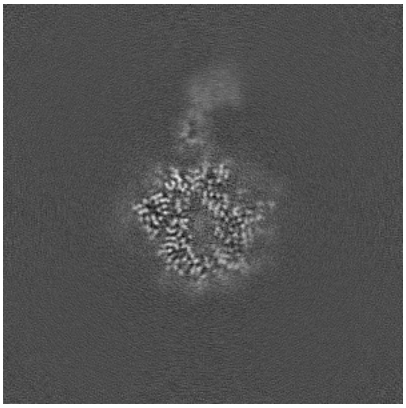


Y Index: 194

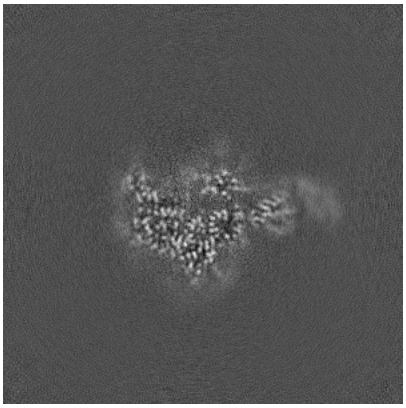


Z Index: 189

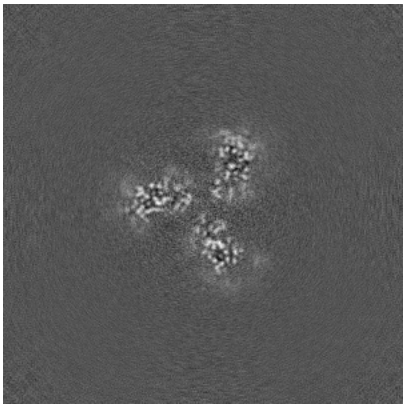
6.3.2 Raw map



X Index: 206



Y Index: 194



Z Index: 189

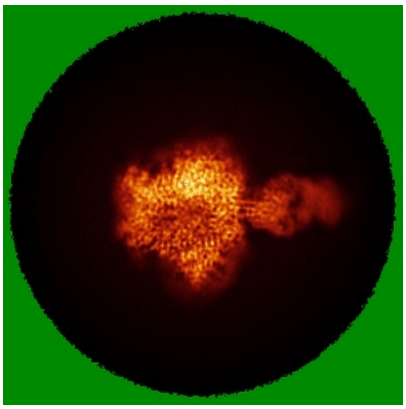
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

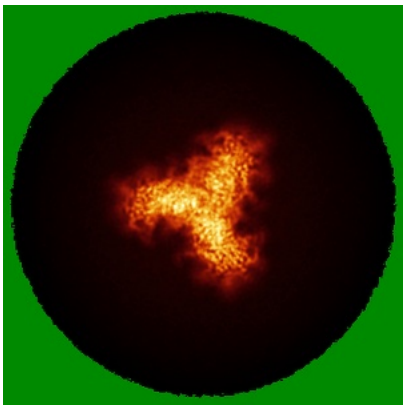
6.4.1 Primary map



X

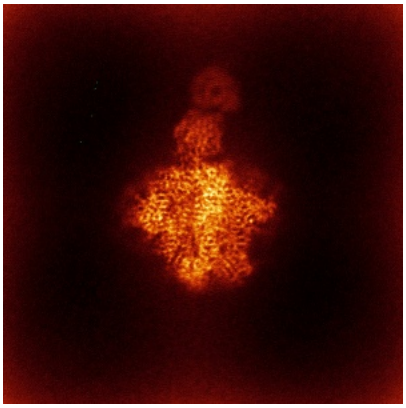


Y

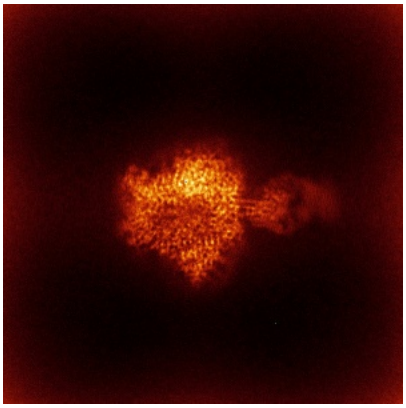


Z

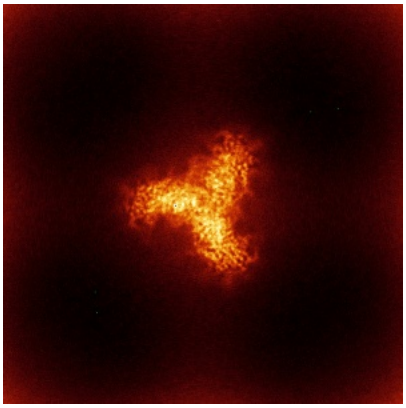
6.4.2 Raw map



X



Y



Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views (i)

6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.09. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X Y Z

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

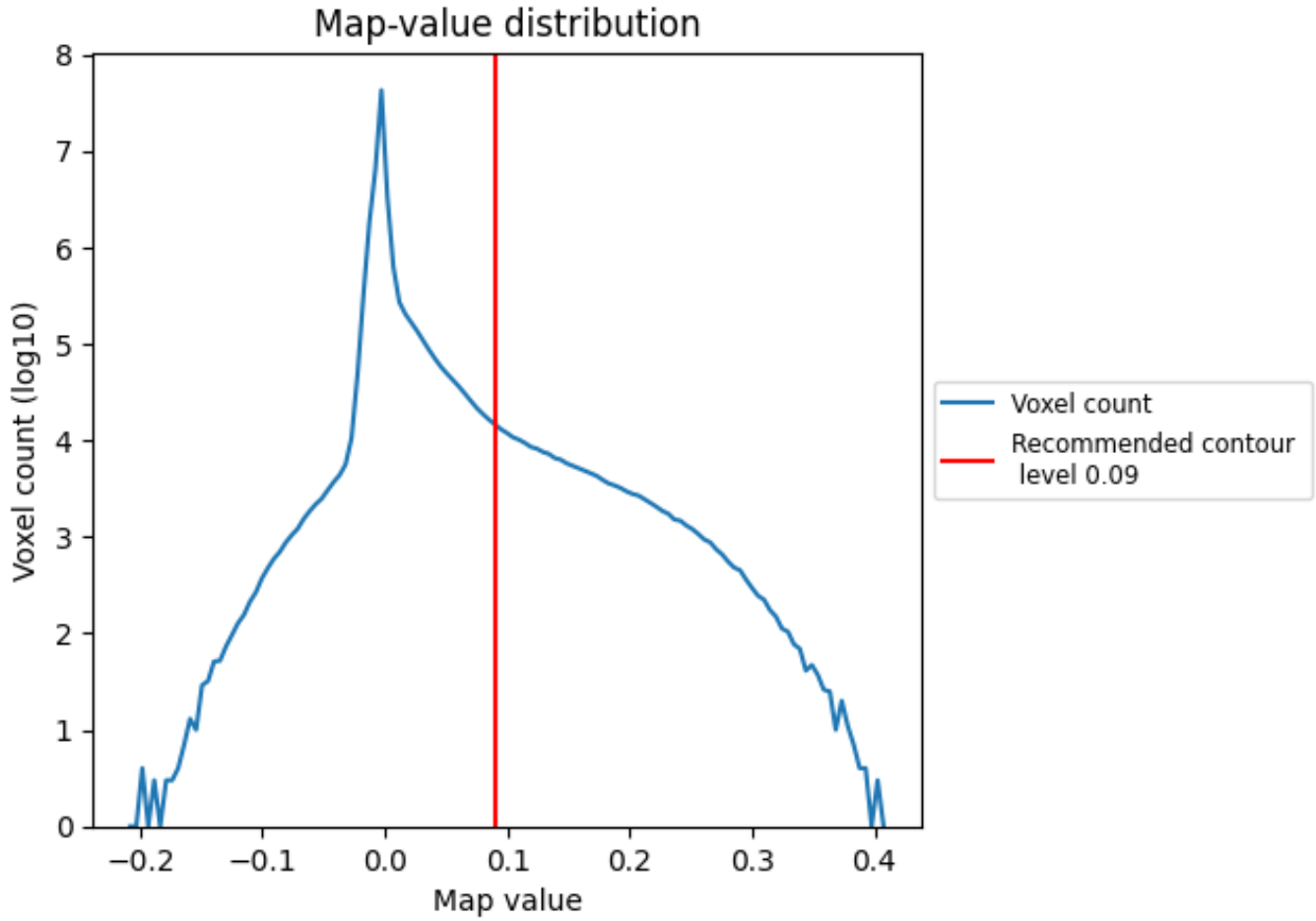
6.6 Mask visualisation (i)

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

7 Map analysis

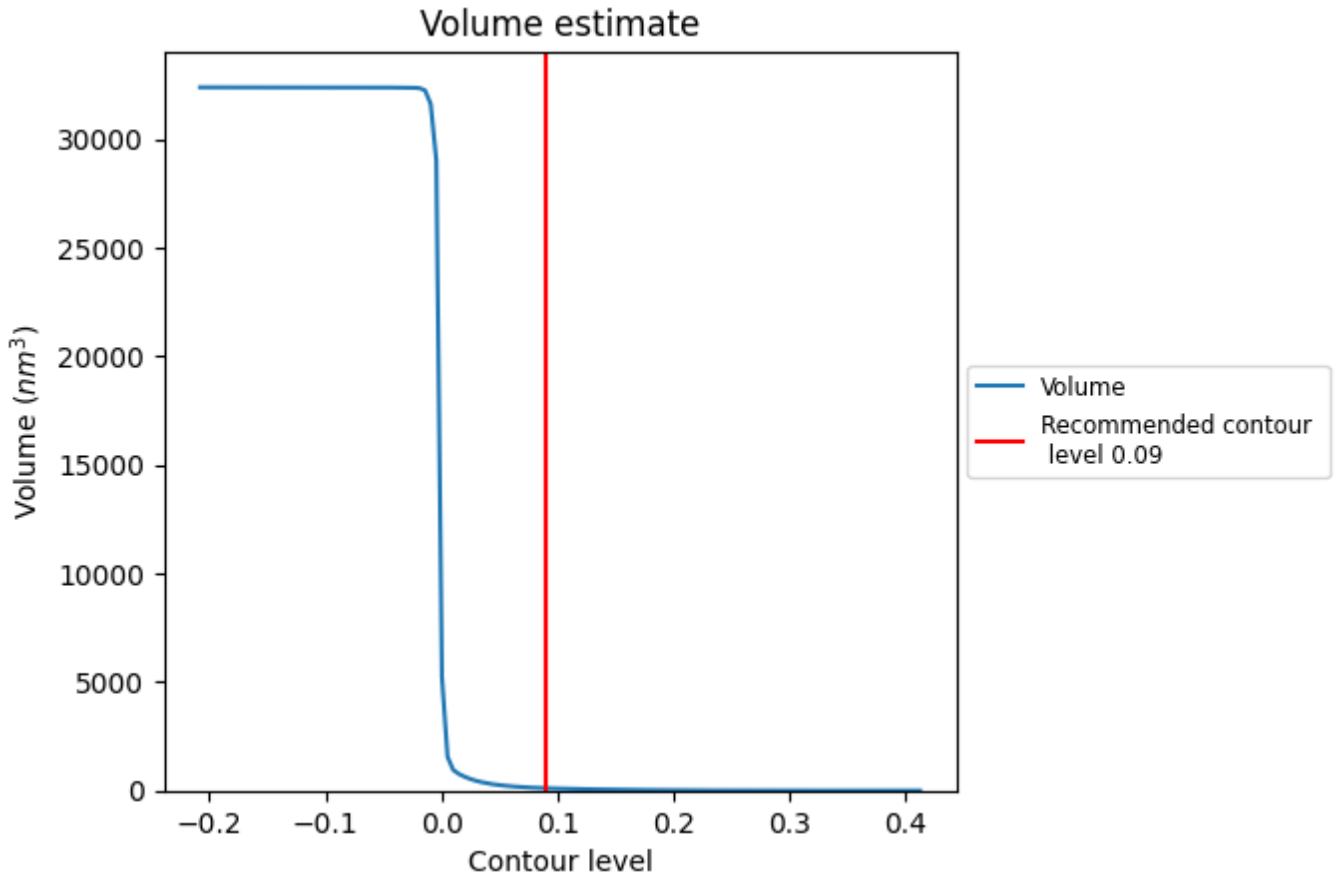
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

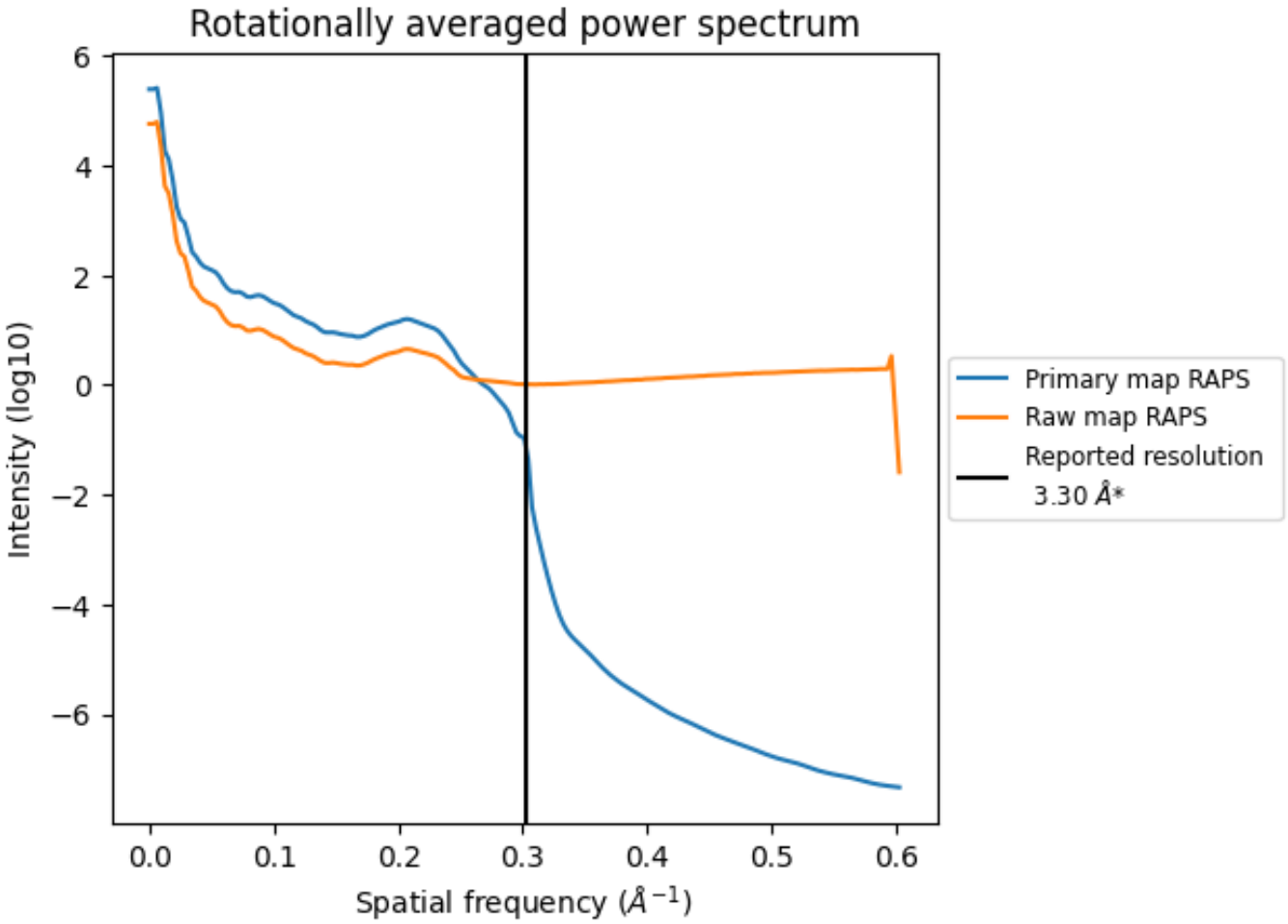
7.2 Volume estimate (i)



The volume at the recommended contour level is 109 nm^3; this corresponds to an approximate mass of 98 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)

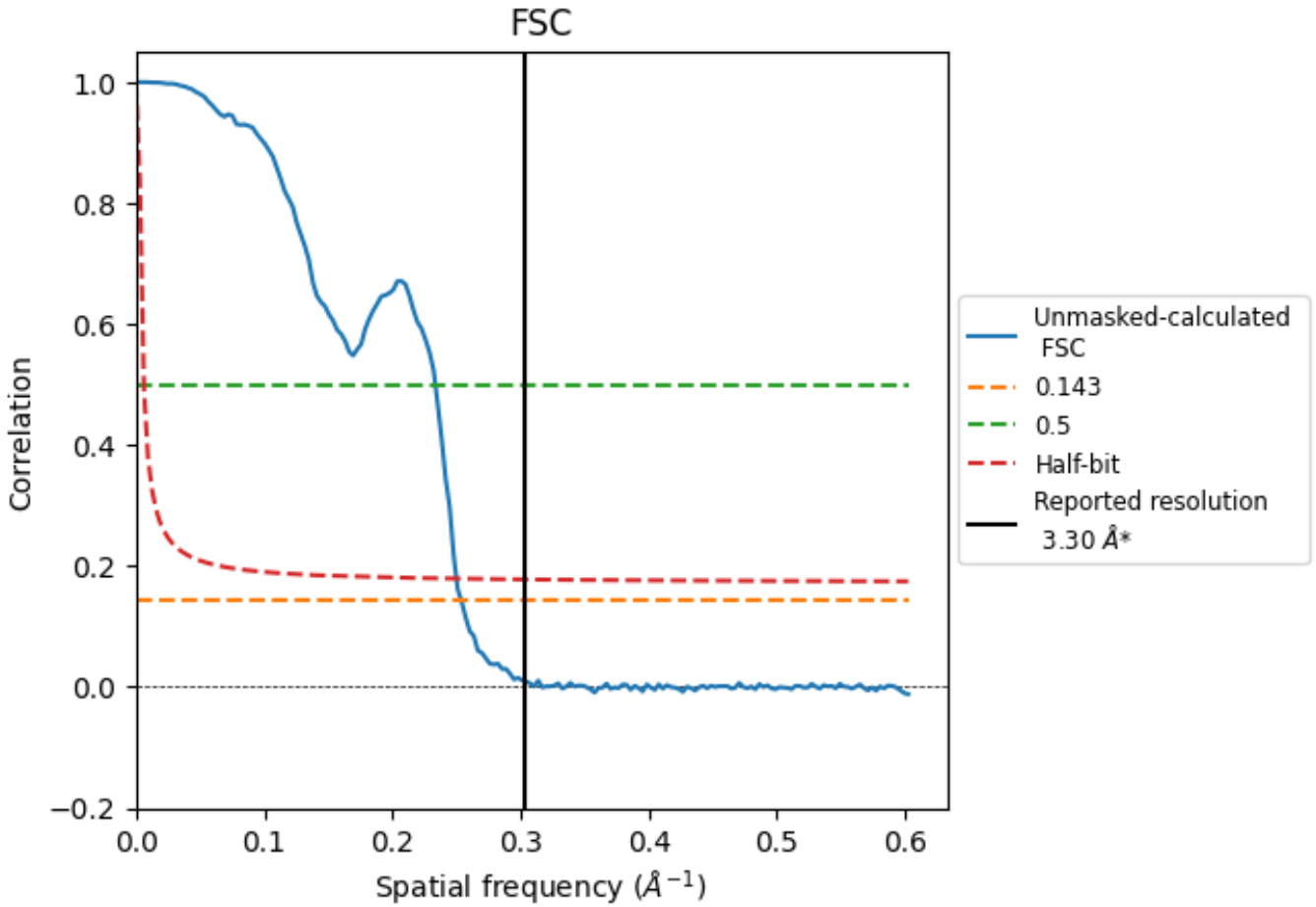


*Reported resolution corresponds to spatial frequency of 0.303 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.303 Å^{-1}

8.2 Resolution estimates (i)

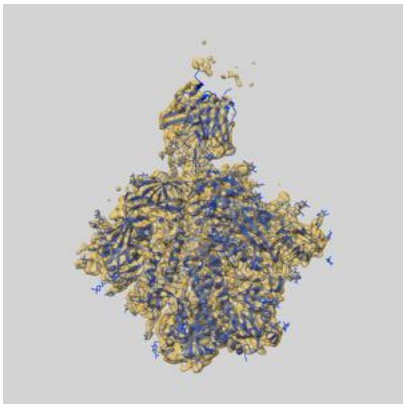
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	-	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.94	4.28	4.00

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

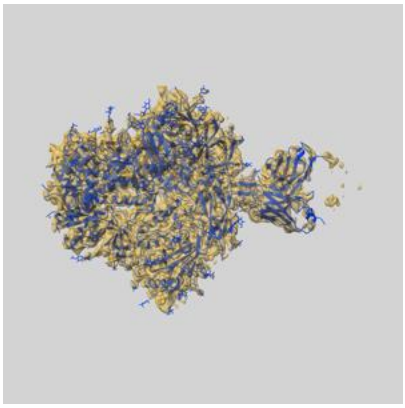
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-74654 and PDB model 9ZRX. Per-residue inclusion information can be found in section 3 on page 15.

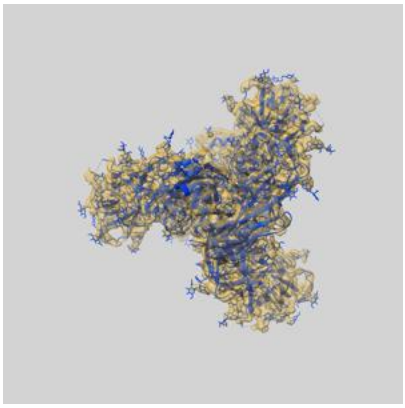
9.1 Map-model overlay



X



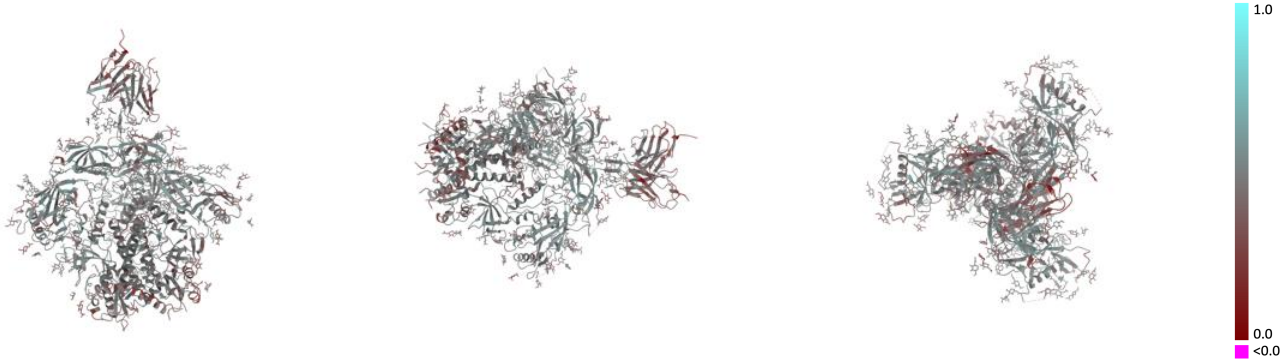
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.09 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.

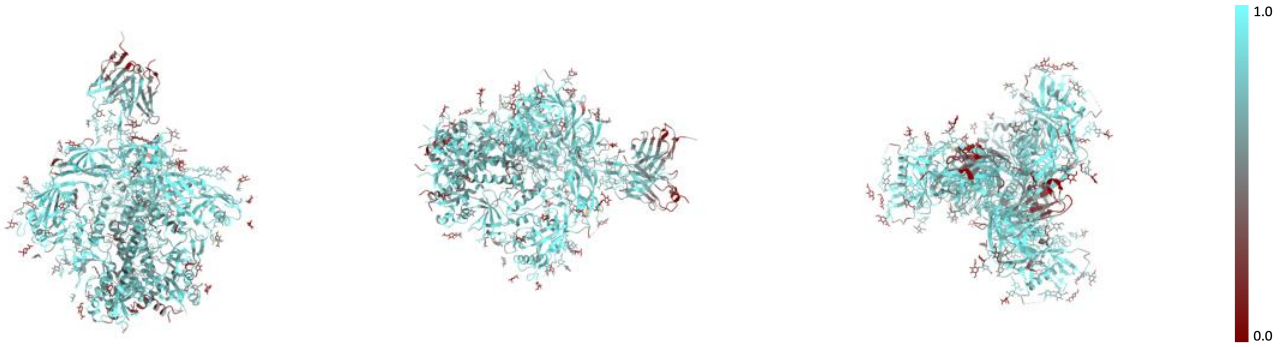
9.2 Q-score mapped to coordinate model (i)



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.

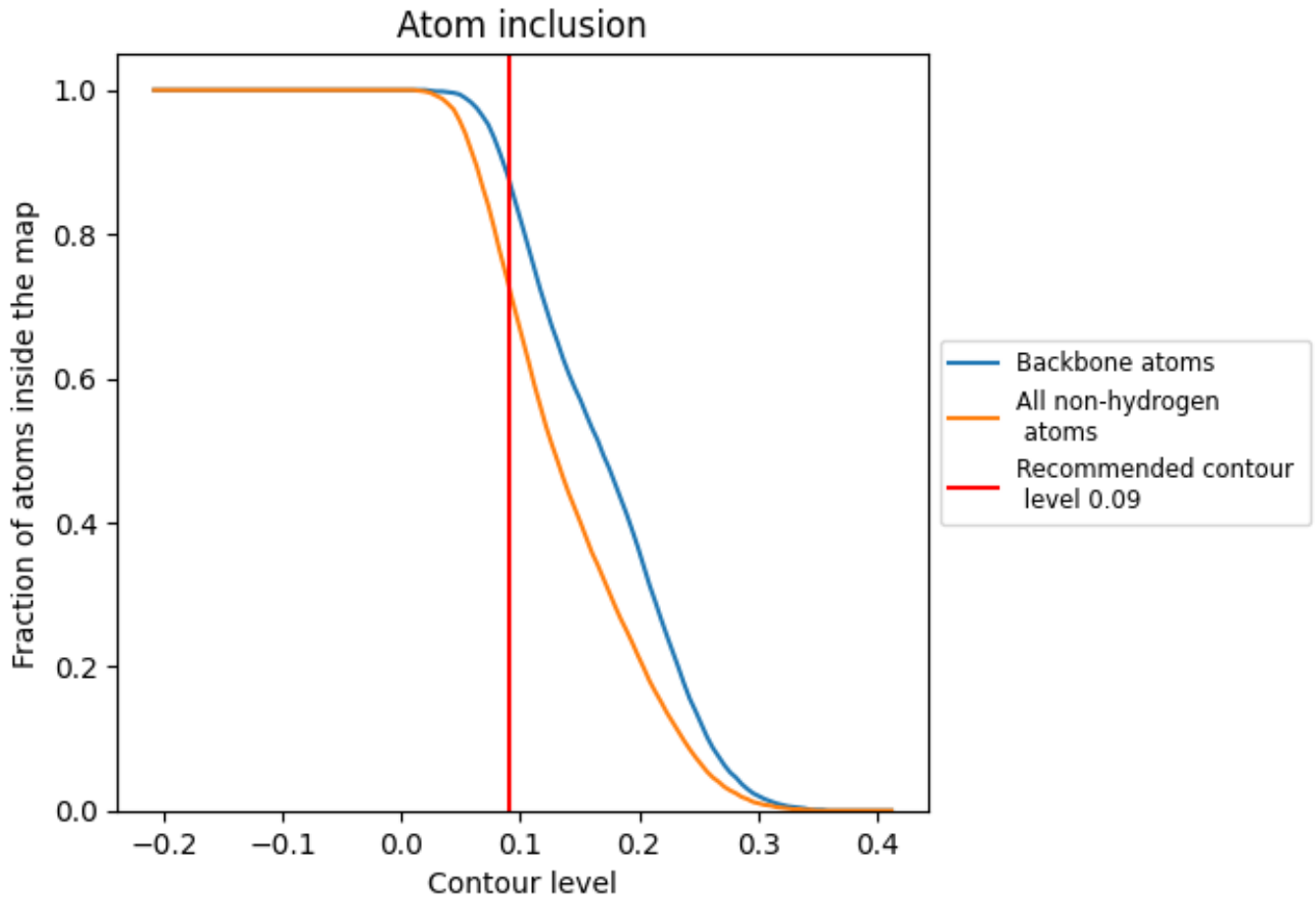
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.09).

9.4 Atom inclusion (i)



At the recommended contour level, 88% of all backbone atoms, 73% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary

The table lists the average atom inclusion at the recommended contour level (0.09) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7310	0.4600
A	0.5910	0.4120
B	0.5760	0.3830
C	0.5360	0.4610
D	0.5900	0.4560
E	0.5740	0.4680
F	0.1430	0.3180
G	0.2860	0.4100
H	0.4640	0.4470
I	0.4290	0.3870
J	0.4100	0.4270
K	0.5360	0.3780
L	0.2860	0.3260
M	0.3570	0.4220
N	0.3570	0.3870
O	0.6070	0.4710
P	0.5900	0.4750
Q	0.5570	0.4630
R	0.0360	0.3240
S	0.7800	0.4760
T	0.6070	0.4930
U	0.5360	0.4410
V	0.5710	0.4700
W	0.5710	0.3370
X	0.3210	0.3870
Y	0.5000	0.3530
Z	0.7140	0.4390
a	0.3570	0.3860
b	0.7910	0.4760
c	0.6690	0.4120
d	0.7980	0.4870
e	0.6990	0.4460
f	0.8050	0.4900
g	0.6640	0.4240
h	0.5360	0.4160



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Chain	Atom inclusion	Q-score
i	0.8180	0.4880
j	0.6070	0.4730
k	0.5640	0.4840
l	0.5410	0.4020
m	0.2140	0.3490
n	0.6560	0.4560
o	0.6790	0.4690
p	0.5710	0.4660
q	0.6430	0.4560
r	0.5710	0.3690
s	0.3930	0.4330
t	0.5000	0.4130
u	0.5000	0.3890