Full wwPDB X-ray Structure Validation Report (i)

Mar 3, 2024 – 11:01 PM EST

PDB ID	:	6UBB
Title	:	Crystal structure of a GH128 (subgroup VI) exo-beta-1,3-glucanase from Aureobasidium namibiae (AnGH128_VI) with laminaribiose at the surface-binding site
Authors	:	Santos, C.R.; Vieira, P.S.; Domingues, M.N.; Cordeiro, R.L.; Tomazini, A.; Murakami, M.T.
Deposited on	:	2019-09-11
Resolution	:	2.35 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

The following versions of software and data (see references (i)) were used in the production of this report:

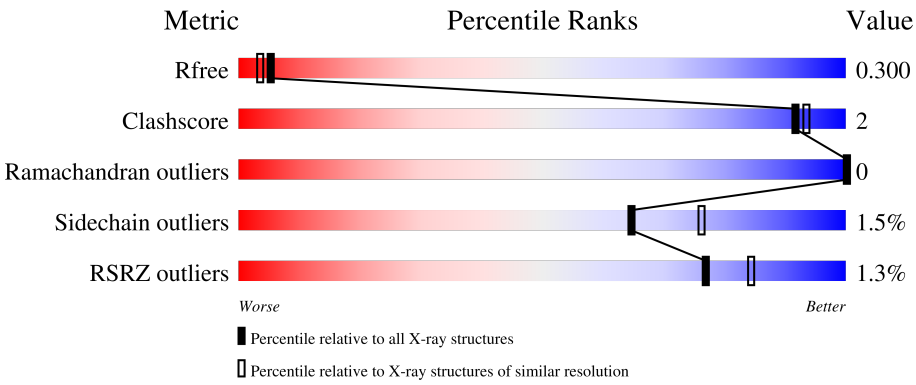
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	1164 (2.36-2.36)
Clashscore	141614	1232 (2.36-2.36)
Ramachandran outliers	138981	1211 (2.36-2.36)
Sidechain outliers	138945	1212 (2.36-2.36)
RSRZ outliers	127900	1150 (2.36-2.36)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	266	88% 5% 7%
1	B	266	88% 5% 7%
1	C	266	.% 91% • 6%
1	D	266	2% 88% 5% 7%

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Mol	Chain	Length	Quality of chain
1	E	266	89% • 7%
1	F	266	.% 85% 8% • 7%
1	G	266	88% 5% 7%
1	H	266	3% 89% • 7%
1	I	266	3% 87% 6% 7%
1	J	266	2% 89% • 7%
2	K	2	100%
2	L	2	50% 50%
2	M	2	100%
2	N	2	100%
2	O	2	100%
2	P	2	100%
2	Q	2	100%
2	R	2	100%
2	S	2	50% 50%
2	T	2	100%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 21017 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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 Glyco_hydro_cc domain-containing protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	247	2030	1300	352	362	16	0	0	0
1	B	247	2030	1300	352	362	16	0	0	0
1	C	249	2040	1305	354	365	16	0	0	0
1	D	247	2030	1300	352	362	16	0	0	0
1	E	247	2030	1300	352	362	16	0	0	0
1	F	247	2030	1300	352	362	16	0	0	0
1	G	247	2030	1300	352	362	16	0	0	0
1	H	247	2030	1300	352	362	16	0	0	0
1	I	247	2030	1300	352	362	16	0	0	0
1	J	247	2030	1300	352	362	16	0	0	0

There are 200 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A074W9U7
A	-18	GLY	-	expression tag	UNP A0A074W9U7
A	-17	SER	-	expression tag	UNP A0A074W9U7
A	-16	SER	-	expression tag	UNP A0A074W9U7
A	-15	HIS	-	expression tag	UNP A0A074W9U7
A	-14	HIS	-	expression tag	UNP A0A074W9U7
A	-13	HIS	-	expression tag	UNP A0A074W9U7
A	-12	HIS	-	expression tag	UNP A0A074W9U7
A	-11	HIS	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-10	HIS	-	expression tag	UNP A0A074W9U7
A	-9	SER	-	expression tag	UNP A0A074W9U7
A	-8	SER	-	expression tag	UNP A0A074W9U7
A	-7	GLY	-	expression tag	UNP A0A074W9U7
A	-6	LEU	-	expression tag	UNP A0A074W9U7
A	-5	VAL	-	expression tag	UNP A0A074W9U7
A	-4	PRO	-	expression tag	UNP A0A074W9U7
A	-3	ALA	-	expression tag	UNP A0A074W9U7
A	-2	GLY	-	expression tag	UNP A0A074W9U7
A	-1	SER	-	expression tag	UNP A0A074W9U7
A	0	HIS	-	expression tag	UNP A0A074W9U7
B	-19	MET	-	initiating methionine	UNP A0A074W9U7
B	-18	GLY	-	expression tag	UNP A0A074W9U7
B	-17	SER	-	expression tag	UNP A0A074W9U7
B	-16	SER	-	expression tag	UNP A0A074W9U7
B	-15	HIS	-	expression tag	UNP A0A074W9U7
B	-14	HIS	-	expression tag	UNP A0A074W9U7
B	-13	HIS	-	expression tag	UNP A0A074W9U7
B	-12	HIS	-	expression tag	UNP A0A074W9U7
B	-11	HIS	-	expression tag	UNP A0A074W9U7
B	-10	HIS	-	expression tag	UNP A0A074W9U7
B	-9	SER	-	expression tag	UNP A0A074W9U7
B	-8	SER	-	expression tag	UNP A0A074W9U7
B	-7	GLY	-	expression tag	UNP A0A074W9U7
B	-6	LEU	-	expression tag	UNP A0A074W9U7
B	-5	VAL	-	expression tag	UNP A0A074W9U7
B	-4	PRO	-	expression tag	UNP A0A074W9U7
B	-3	ALA	-	expression tag	UNP A0A074W9U7
B	-2	GLY	-	expression tag	UNP A0A074W9U7
B	-1	SER	-	expression tag	UNP A0A074W9U7
B	0	HIS	-	expression tag	UNP A0A074W9U7
C	-19	MET	-	initiating methionine	UNP A0A074W9U7
C	-18	GLY	-	expression tag	UNP A0A074W9U7
C	-17	SER	-	expression tag	UNP A0A074W9U7
C	-16	SER	-	expression tag	UNP A0A074W9U7
C	-15	HIS	-	expression tag	UNP A0A074W9U7
C	-14	HIS	-	expression tag	UNP A0A074W9U7
C	-13	HIS	-	expression tag	UNP A0A074W9U7
C	-12	HIS	-	expression tag	UNP A0A074W9U7
C	-11	HIS	-	expression tag	UNP A0A074W9U7
C	-10	HIS	-	expression tag	UNP A0A074W9U7
C	-9	SER	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-8	SER	-	expression tag	UNP A0A074W9U7
C	-7	GLY	-	expression tag	UNP A0A074W9U7
C	-6	LEU	-	expression tag	UNP A0A074W9U7
C	-5	VAL	-	expression tag	UNP A0A074W9U7
C	-4	PRO	-	expression tag	UNP A0A074W9U7
C	-3	ALA	-	expression tag	UNP A0A074W9U7
C	-2	GLY	-	expression tag	UNP A0A074W9U7
C	-1	SER	-	expression tag	UNP A0A074W9U7
C	0	HIS	-	expression tag	UNP A0A074W9U7
D	-19	MET	-	initiating methionine	UNP A0A074W9U7
D	-18	GLY	-	expression tag	UNP A0A074W9U7
D	-17	SER	-	expression tag	UNP A0A074W9U7
D	-16	SER	-	expression tag	UNP A0A074W9U7
D	-15	HIS	-	expression tag	UNP A0A074W9U7
D	-14	HIS	-	expression tag	UNP A0A074W9U7
D	-13	HIS	-	expression tag	UNP A0A074W9U7
D	-12	HIS	-	expression tag	UNP A0A074W9U7
D	-11	HIS	-	expression tag	UNP A0A074W9U7
D	-10	HIS	-	expression tag	UNP A0A074W9U7
D	-9	SER	-	expression tag	UNP A0A074W9U7
D	-8	SER	-	expression tag	UNP A0A074W9U7
D	-7	GLY	-	expression tag	UNP A0A074W9U7
D	-6	LEU	-	expression tag	UNP A0A074W9U7
D	-5	VAL	-	expression tag	UNP A0A074W9U7
D	-4	PRO	-	expression tag	UNP A0A074W9U7
D	-3	ALA	-	expression tag	UNP A0A074W9U7
D	-2	GLY	-	expression tag	UNP A0A074W9U7
D	-1	SER	-	expression tag	UNP A0A074W9U7
D	0	HIS	-	expression tag	UNP A0A074W9U7
E	-19	MET	-	initiating methionine	UNP A0A074W9U7
E	-18	GLY	-	expression tag	UNP A0A074W9U7
E	-17	SER	-	expression tag	UNP A0A074W9U7
E	-16	SER	-	expression tag	UNP A0A074W9U7
E	-15	HIS	-	expression tag	UNP A0A074W9U7
E	-14	HIS	-	expression tag	UNP A0A074W9U7
E	-13	HIS	-	expression tag	UNP A0A074W9U7
E	-12	HIS	-	expression tag	UNP A0A074W9U7
E	-11	HIS	-	expression tag	UNP A0A074W9U7
E	-10	HIS	-	expression tag	UNP A0A074W9U7
E	-9	SER	-	expression tag	UNP A0A074W9U7
E	-8	SER	-	expression tag	UNP A0A074W9U7
E	-7	GLY	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	-6	LEU	-	expression tag	UNP A0A074W9U7
E	-5	VAL	-	expression tag	UNP A0A074W9U7
E	-4	PRO	-	expression tag	UNP A0A074W9U7
E	-3	ALA	-	expression tag	UNP A0A074W9U7
E	-2	GLY	-	expression tag	UNP A0A074W9U7
E	-1	SER	-	expression tag	UNP A0A074W9U7
E	0	HIS	-	expression tag	UNP A0A074W9U7
F	-19	MET	-	initiating methionine	UNP A0A074W9U7
F	-18	GLY	-	expression tag	UNP A0A074W9U7
F	-17	SER	-	expression tag	UNP A0A074W9U7
F	-16	SER	-	expression tag	UNP A0A074W9U7
F	-15	HIS	-	expression tag	UNP A0A074W9U7
F	-14	HIS	-	expression tag	UNP A0A074W9U7
F	-13	HIS	-	expression tag	UNP A0A074W9U7
F	-12	HIS	-	expression tag	UNP A0A074W9U7
F	-11	HIS	-	expression tag	UNP A0A074W9U7
F	-10	HIS	-	expression tag	UNP A0A074W9U7
F	-9	SER	-	expression tag	UNP A0A074W9U7
F	-8	SER	-	expression tag	UNP A0A074W9U7
F	-7	GLY	-	expression tag	UNP A0A074W9U7
F	-6	LEU	-	expression tag	UNP A0A074W9U7
F	-5	VAL	-	expression tag	UNP A0A074W9U7
F	-4	PRO	-	expression tag	UNP A0A074W9U7
F	-3	ALA	-	expression tag	UNP A0A074W9U7
F	-2	GLY	-	expression tag	UNP A0A074W9U7
F	-1	SER	-	expression tag	UNP A0A074W9U7
F	0	HIS	-	expression tag	UNP A0A074W9U7
G	-19	MET	-	initiating methionine	UNP A0A074W9U7
G	-18	GLY	-	expression tag	UNP A0A074W9U7
G	-17	SER	-	expression tag	UNP A0A074W9U7
G	-16	SER	-	expression tag	UNP A0A074W9U7
G	-15	HIS	-	expression tag	UNP A0A074W9U7
G	-14	HIS	-	expression tag	UNP A0A074W9U7
G	-13	HIS	-	expression tag	UNP A0A074W9U7
G	-12	HIS	-	expression tag	UNP A0A074W9U7
G	-11	HIS	-	expression tag	UNP A0A074W9U7
G	-10	HIS	-	expression tag	UNP A0A074W9U7
G	-9	SER	-	expression tag	UNP A0A074W9U7
G	-8	SER	-	expression tag	UNP A0A074W9U7
G	-7	GLY	-	expression tag	UNP A0A074W9U7
G	-6	LEU	-	expression tag	UNP A0A074W9U7
G	-5	VAL	-	expression tag	UNP A0A074W9U7

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-4	PRO	-	expression tag	UNP A0A074W9U7
G	-3	ALA	-	expression tag	UNP A0A074W9U7
G	-2	GLY	-	expression tag	UNP A0A074W9U7
G	-1	SER	-	expression tag	UNP A0A074W9U7
G	0	HIS	-	expression tag	UNP A0A074W9U7
H	-19	MET	-	initiating methionine	UNP A0A074W9U7
H	-18	GLY	-	expression tag	UNP A0A074W9U7
H	-17	SER	-	expression tag	UNP A0A074W9U7
H	-16	SER	-	expression tag	UNP A0A074W9U7
H	-15	HIS	-	expression tag	UNP A0A074W9U7
H	-14	HIS	-	expression tag	UNP A0A074W9U7
H	-13	HIS	-	expression tag	UNP A0A074W9U7
H	-12	HIS	-	expression tag	UNP A0A074W9U7
H	-11	HIS	-	expression tag	UNP A0A074W9U7
H	-10	HIS	-	expression tag	UNP A0A074W9U7
H	-9	SER	-	expression tag	UNP A0A074W9U7
H	-8	SER	-	expression tag	UNP A0A074W9U7
H	-7	GLY	-	expression tag	UNP A0A074W9U7
H	-6	LEU	-	expression tag	UNP A0A074W9U7
H	-5	VAL	-	expression tag	UNP A0A074W9U7
H	-4	PRO	-	expression tag	UNP A0A074W9U7
H	-3	ALA	-	expression tag	UNP A0A074W9U7
H	-2	GLY	-	expression tag	UNP A0A074W9U7
H	-1	SER	-	expression tag	UNP A0A074W9U7
H	0	HIS	-	expression tag	UNP A0A074W9U7
I	-19	MET	-	initiating methionine	UNP A0A074W9U7
I	-18	GLY	-	expression tag	UNP A0A074W9U7
I	-17	SER	-	expression tag	UNP A0A074W9U7
I	-16	SER	-	expression tag	UNP A0A074W9U7
I	-15	HIS	-	expression tag	UNP A0A074W9U7
I	-14	HIS	-	expression tag	UNP A0A074W9U7
I	-13	HIS	-	expression tag	UNP A0A074W9U7
I	-12	HIS	-	expression tag	UNP A0A074W9U7
I	-11	HIS	-	expression tag	UNP A0A074W9U7
I	-10	HIS	-	expression tag	UNP A0A074W9U7
I	-9	SER	-	expression tag	UNP A0A074W9U7
I	-8	SER	-	expression tag	UNP A0A074W9U7
I	-7	GLY	-	expression tag	UNP A0A074W9U7
I	-6	LEU	-	expression tag	UNP A0A074W9U7
I	-5	VAL	-	expression tag	UNP A0A074W9U7
I	-4	PRO	-	expression tag	UNP A0A074W9U7
I	-3	ALA	-	expression tag	UNP A0A074W9U7

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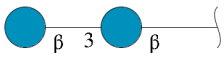
Chain	Residue	Modelled	Actual	Comment	Reference
I	-2	GLY	-	expression tag	UNP A0A074W9U7
I	-1	SER	-	expression tag	UNP A0A074W9U7
I	0	HIS	-	expression tag	UNP A0A074W9U7
J	-19	MET	-	initiating methionine	UNP A0A074W9U7
J	-18	GLY	-	expression tag	UNP A0A074W9U7
J	-17	SER	-	expression tag	UNP A0A074W9U7
J	-16	SER	-	expression tag	UNP A0A074W9U7
J	-15	HIS	-	expression tag	UNP A0A074W9U7
J	-14	HIS	-	expression tag	UNP A0A074W9U7
J	-13	HIS	-	expression tag	UNP A0A074W9U7
J	-12	HIS	-	expression tag	UNP A0A074W9U7
J	-11	HIS	-	expression tag	UNP A0A074W9U7
J	-10	HIS	-	expression tag	UNP A0A074W9U7
J	-9	SER	-	expression tag	UNP A0A074W9U7
J	-8	SER	-	expression tag	UNP A0A074W9U7
J	-7	GLY	-	expression tag	UNP A0A074W9U7
J	-6	LEU	-	expression tag	UNP A0A074W9U7
J	-5	VAL	-	expression tag	UNP A0A074W9U7
J	-4	PRO	-	expression tag	UNP A0A074W9U7
J	-3	ALA	-	expression tag	UNP A0A074W9U7
J	-2	GLY	-	expression tag	UNP A0A074W9U7
J	-1	SER	-	expression tag	UNP A0A074W9U7
J	0	HIS	-	expression tag	UNP A0A074W9U7

- Molecule 2 is an oligosaccharide called beta-D-glucopyranose-(1-3)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	K	2	23	12	11	0	0	0
2	L	2	23	12	11	0	0	0
2	M	2	23	12	11	0	0	0
2	N	2	23	12	11	0	0	0
2	O	2	23	12	11	0	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
			Total	C	O			
2	P	2	23	12	11	0	0	0
2	Q	2	23	12	11	0	0	0
2	R	2	23	12	11	0	0	0
2	S	2	23	12	11	0	0	0
2	T	2	23	12	11	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	63	63	63	0	0
3	B	44	44	44	0	0
3	C	53	53	53	0	0
3	D	40	40	40	0	0
3	E	50	50	50	0	0
3	F	55	55	55	0	0
3	G	61	61	61	0	0
3	H	43	43	43	0	0
3	I	35	35	35	0	0
3	J	33	33	33	0	0

3 Residue-property plots (i)

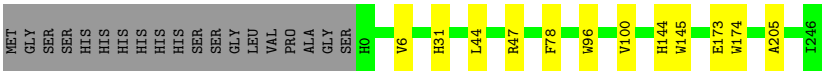
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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 dot above a residue indicates a poor fit to the electron density (RSRZ > 2). 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: Glyco_hydro_cc domain-containing protein

Chain A:

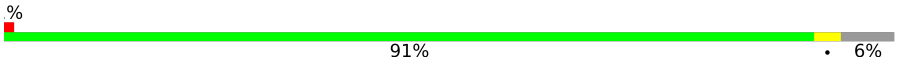
• Molecule 1: Glyco_hydro_cc domain-containing protein

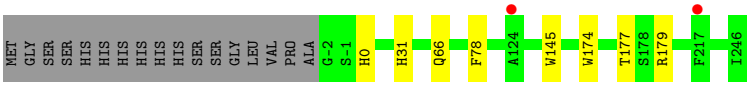
Chain B:



• Molecule 1: Glyco_hydro_cc domain-containing protein

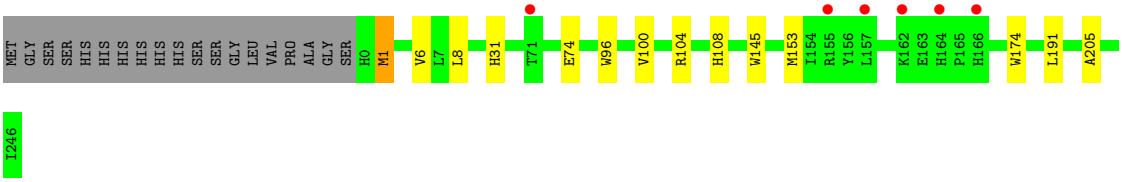
Chain C:





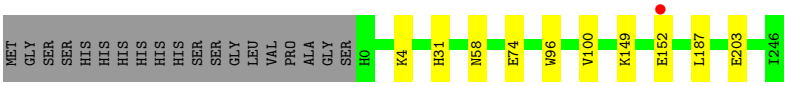
• Molecule 1: Glyco_hydro_cc domain-containing protein

Chain D:



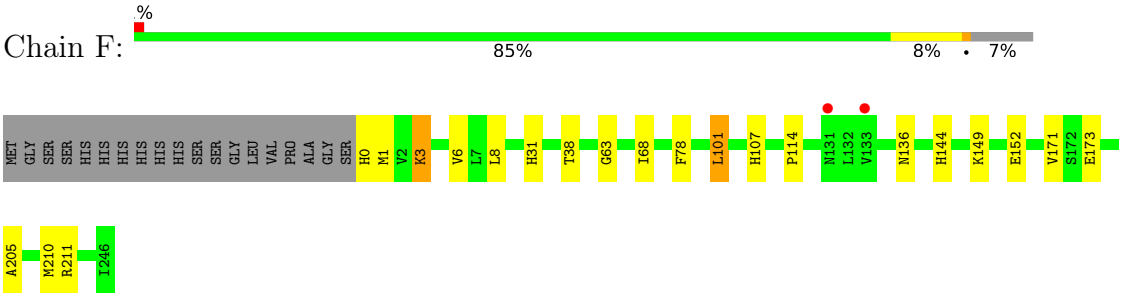
• Molecule 1: Glyco_hydro_cc domain-containing protein

Chain E:



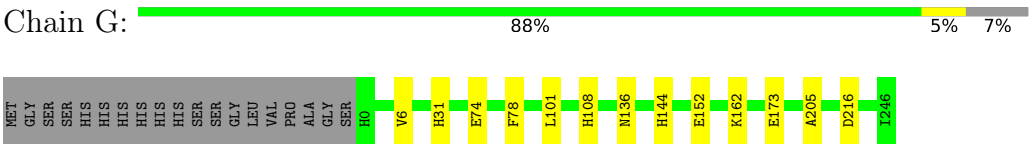
• Molecule 1: Glyco_hydro_cc domain-containing protein

Chain F:



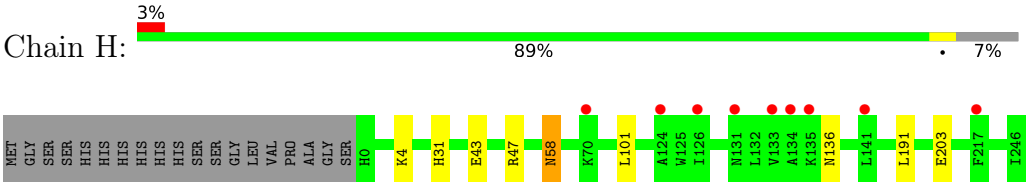
• Molecule 1: Glyco_hydro_cc domain-containing protein

Chain G:



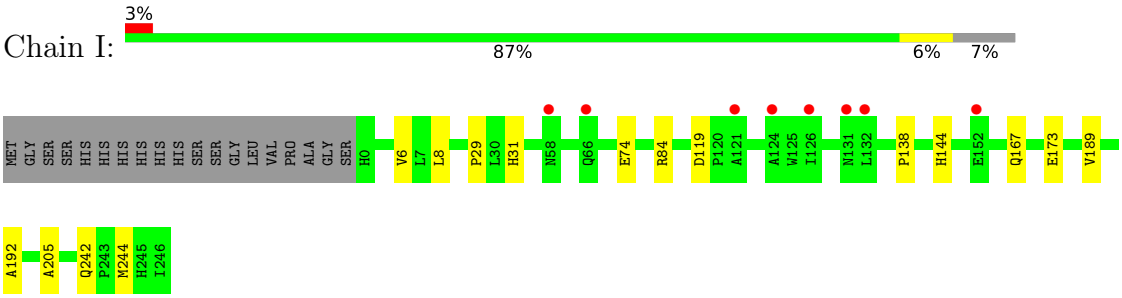
• Molecule 1: Glyco_hydro_cc domain-containing protein

Chain H:



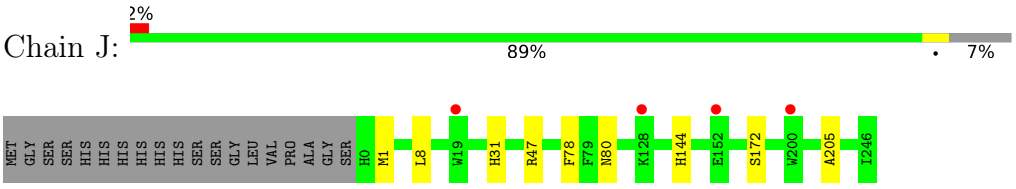
• Molecule 1: Glyco_hydro_cc domain-containing protein

Chain I:



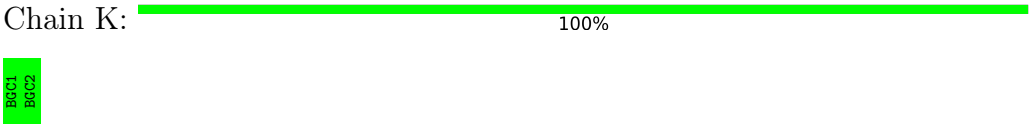
• Molecule 1: Glyco_hydro_cc domain-containing protein

Chain J:



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

Chain K:



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

Chain L: 50% 50%

BGC1 BGC2

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

Chain M: 100%

BGC1 BGC2

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

Chain N: 100%

BGC1 BGC2

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

Chain O: 100%

BGC1 BGC2

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

Chain P: 100%

BGC1 BGC2

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

Chain Q: 100%

BGC1 BGC2

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

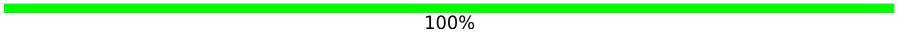
Chain R: 100%

BGC1 BGC2

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

Chain S: 50% 50%

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

Chain T: 

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	121.35Å 152.73Å 169.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.30 – 2.35 49.30 – 2.35	Depositor EDS
% Data completeness (in resolution range)	99.8 (49.30-2.35) 99.9 (49.30-2.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.35 (at 2.34Å)	Xtriage
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.265 , 0.301 0.266 , 0.300	Depositor DCC
R_{free} test set	6541 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	28.5	Xtriage
Anisotropy	0.684	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 33.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	21017	wwPDB-VP
Average B, all atoms (Å^2)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.98% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: BGC

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	0.64	0/2098	0.70	0/2851
1	B	0.64	0/2098	0.69	0/2851
1	C	0.64	0/2108	0.69	0/2864
1	D	0.64	0/2098	0.70	0/2851
1	E	0.64	0/2098	0.70	0/2851
1	F	0.64	0/2098	0.71	0/2851
1	G	0.64	0/2098	0.70	0/2851
1	H	0.64	0/2098	0.69	0/2851
1	I	0.64	0/2098	0.69	0/2851
1	J	0.64	0/2098	0.70	0/2851
All	All	0.64	0/20990	0.70	0/28523

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2030	0	1958	7	0
1	B	2030	0	1958	5	0
1	C	2040	0	1966	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	2030	0	1958	12	0
1	E	2030	0	1958	5	0
1	F	2030	0	1958	19	0
1	G	2030	0	1958	5	0
1	H	2030	0	1958	4	0
1	I	2030	0	1958	9	0
1	J	2030	0	1958	2	0
2	K	23	0	21	0	0
2	L	23	0	21	0	0
2	M	23	0	21	0	0
2	N	23	0	21	0	0
2	O	23	0	21	0	0
2	P	23	0	21	0	0
2	Q	23	0	21	0	0
2	R	23	0	21	0	0
2	S	23	0	21	0	0
2	T	23	0	21	0	0
3	A	63	0	0	0	0
3	B	44	0	0	0	0
3	C	53	0	0	0	0
3	D	40	0	0	0	0
3	E	50	0	0	0	0
3	F	55	0	0	0	0
3	G	61	0	0	0	0
3	H	43	0	0	0	0
3	I	35	0	0	0	0
3	J	33	0	0	0	0
All	All	21017	0	19798	64	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 2.

All (64) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1:MET:HB2	1:F:63:GLY:HA3	1.67	0.74
1:C:66:GLN:HE22	1:F:0:HIS:HE1	1.39	0.69
1:D:1:MET:HB2	1:F:63:GLY:CA	2.27	0.64
1:B:44:LEU:O	1:B:47:ARG:HD3	1.98	0.63
1:I:189:VAL:HG22	1:I:244:MET:HG2	1.79	0.63
1:G:144:HIS:CD2	1:G:173:GLU:HB2	2.39	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:MET:HB3	1:D:191:LEU:HD21	1.89	0.55
1:I:6:VAL:HB	1:I:205:ALA:HB2	1.89	0.54
1:I:84:ARG:NH2	1:I:119:ASP:OD2	2.40	0.54
1:D:1:MET:CE	1:F:63:GLY:HA3	2.39	0.53
1:F:8:LEU:HG	1:F:205:ALA:HB1	1.93	0.51
1:I:138:PRO:HG2	1:I:167:GLN:HE22	1.76	0.50
1:H:43:GLU:O	1:H:47:ARG:NH1	2.43	0.50
1:F:149:LYS:HB2	1:F:152:GLU:HG3	1.94	0.50
1:G:101:LEU:HD11	1:G:136:ASN:CB	2.42	0.50
1:J:80:ASN:ND2	1:J:144:HIS:HE1	2.10	0.49
1:C:177:THR:O	1:C:179:ARG:NH2	2.43	0.48
1:A:8:LEU:HG	1:A:205:ALA:HB1	1.95	0.48
1:C:0:HIS:CE1	1:F:107:HIS:HE2	2.31	0.48
1:I:192:ALA:CB	1:I:244:MET:CE	2.91	0.48
1:J:8:LEU:HG	1:J:205:ALA:HB1	1.95	0.48
1:I:8:LEU:HG	1:I:205:ALA:HB1	1.96	0.47
1:D:1:MET:HE3	1:F:63:GLY:HA3	1.97	0.47
1:F:144:HIS:CD2	1:F:173:GLU:HB2	2.51	0.46
1:F:171:VAL:CG2	1:F:171:VAL:O	2.62	0.46
1:F:1:MET:HG3	1:F:3:LYS:HG2	1.97	0.46
1:F:78:PHE:CD2	1:F:114:PRO:HG3	2.51	0.46
1:C:66:GLN:HE22	1:F:0:HIS:CE1	2.26	0.46
1:A:67:ASN:HA	1:A:70:LYS:HE3	1.98	0.46
1:A:6:VAL:HB	1:A:205:ALA:HB2	1.98	0.45
1:E:149:LYS:HB2	1:E:152:GLU:HG3	1.99	0.45
1:E:187:LEU:C	1:E:187:LEU:HD23	2.36	0.45
1:D:145:TRP:O	1:D:174:TRP:HA	2.17	0.45
1:A:4:LYS:HB2	1:A:203:GLU:HA	1.99	0.44
1:A:145:TRP:O	1:A:174:TRP:HA	2.17	0.44
1:H:58:ASN:H	1:H:58:ASN:HD22	1.65	0.44
1:E:96:TRP:HA	1:E:100:VAL:HB	2.00	0.44
1:H:101:LEU:HD11	1:H:136:ASN:CB	2.47	0.44
1:D:6:VAL:HB	1:D:205:ALA:HB2	2.00	0.44
1:F:6:VAL:HB	1:F:205:ALA:HB2	1.99	0.43
1:D:96:TRP:HA	1:D:100:VAL:HB	2.01	0.43
1:D:1:MET:HE2	1:F:63:GLY:HA3	2.01	0.43
1:F:101:LEU:HD11	1:F:136:ASN:CB	2.48	0.43
1:B:6:VAL:HB	1:B:205:ALA:HB2	2.01	0.42
1:B:144:HIS:CD2	1:B:173:GLU:HB2	2.55	0.42
1:F:171:VAL:O	1:F:171:VAL:HG22	2.19	0.42
1:B:145:TRP:O	1:B:174:TRP:HA	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:8:LEU:HG	1:D:205:ALA:HB1	2.01	0.42
1:H:4:LYS:HB3	1:H:203:GLU:HA	2.01	0.42
1:C:145:TRP:O	1:C:174:TRP:HA	2.19	0.42
1:B:96:TRP:HA	1:B:100:VAL:HB	2.01	0.42
1:I:29:PRO:HG3	1:I:242:GLN:HB2	2.00	0.42
1:D:153:MET:SD	1:D:191:LEU:CD2	3.08	0.42
1:G:101:LEU:HD11	1:G:136:ASN:HB3	2.02	0.42
1:I:192:ALA:CB	1:I:244:MET:HE3	2.50	0.42
1:C:0:HIS:CE1	1:F:107:HIS:NE2	2.89	0.41
1:A:199:PRO:O	1:G:108:HIS:NE2	2.52	0.41
1:F:38:THR:HG22	1:F:68:ILE:HD11	2.02	0.41
1:G:6:VAL:HB	1:G:205:ALA:HB2	2.02	0.41
1:A:170:ILE:HG23	1:A:203:GLU:HG3	2.03	0.41
1:D:104:ARG:O	1:D:108:HIS:HA	2.22	0.40
1:E:4:LYS:HB2	1:E:203:GLU:HA	2.02	0.40
1:I:144:HIS:CD2	1:I:173:GLU:HB2	2.56	0.40
1:E:58:ASN:HD22	1:E:58:ASN:H	1.69	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	245/266 (92%)	232 (95%)	13 (5%)	0	100	100
1	B	245/266 (92%)	232 (95%)	13 (5%)	0	100	100
1	C	247/266 (93%)	234 (95%)	13 (5%)	0	100	100
1	D	245/266 (92%)	234 (96%)	11 (4%)	0	100	100
1	E	245/266 (92%)	234 (96%)	11 (4%)	0	100	100
1	F	245/266 (92%)	233 (95%)	12 (5%)	0	100	100
1	G	245/266 (92%)	229 (94%)	16 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	H	245/266 (92%)	233 (95%)	12 (5%)	0	100	100
1	I	245/266 (92%)	234 (96%)	11 (4%)	0	100	100
1	J	245/266 (92%)	235 (96%)	10 (4%)	0	100	100
All	All	2452/2660 (92%)	2330 (95%)	122 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	220/235 (94%)	217 (99%)	3 (1%)	67	78
1	B	220/235 (94%)	218 (99%)	2 (1%)	78	87
1	C	221/235 (94%)	219 (99%)	2 (1%)	78	87
1	D	220/235 (94%)	217 (99%)	3 (1%)	67	78
1	E	220/235 (94%)	218 (99%)	2 (1%)	78	87
1	F	220/235 (94%)	215 (98%)	5 (2%)	50	61
1	G	220/235 (94%)	214 (97%)	6 (3%)	44	55
1	H	220/235 (94%)	217 (99%)	3 (1%)	67	78
1	I	220/235 (94%)	218 (99%)	2 (1%)	78	87
1	J	220/235 (94%)	215 (98%)	5 (2%)	50	61
All	All	2201/2350 (94%)	2168 (98%)	33 (2%)	65	76

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	74	GLU
1	A	99	GLN
1	A	167	GLN
1	B	31	HIS
1	B	78	PHE

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Mol	Chain	Res	Type
1	C	31	HIS
1	C	78	PHE
1	D	1	MET
1	D	31	HIS
1	D	74	GLU
1	E	31	HIS
1	E	74	GLU
1	F	3	LYS
1	F	31	HIS
1	F	101	LEU
1	F	210	MET
1	F	211	ARG
1	G	31	HIS
1	G	74	GLU
1	G	78	PHE
1	G	152	GLU
1	G	162	LYS
1	G	216	ASP
1	H	31	HIS
1	H	58	ASN
1	H	191	LEU
1	I	31	HIS
1	I	74	GLU
1	J	1	MET
1	J	31	HIS
1	J	47	ARG
1	J	78	PHE
1	J	172	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (23) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	66	GLN
1	A	167	GLN
1	A	212	GLN
1	B	58	ASN
1	C	58	ASN
1	C	66	GLN
1	C	161	HIS
1	D	58	ASN
1	D	66	GLN
1	D	212	GLN

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Mol	Chain	Res	Type
1	D	245	HIS
1	E	58	ASN
1	E	107	HIS
1	E	131	ASN
1	E	212	GLN
1	H	0	HIS
1	H	58	ASN
1	H	107	HIS
1	J	55	HIS
1	J	66	GLN
1	J	80	ASN
1	J	144	HIS
1	J	167	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

20 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	BGC	K	1	2	12,12,12	0.50	0	17,17,17	0.58	0
2	BGC	K	2	2	11,11,12	0.29	0	15,15,17	0.77	0
2	BGC	L	1	2	12,12,12	0.47	0	17,17,17	0.66	0
2	BGC	L	2	2	11,11,12	0.26	0	15,15,17	0.79	1 (6%)
2	BGC	M	1	2	12,12,12	0.51	0	17,17,17	0.69	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	BGC	M	2	2	11,11,12	0.31	0	15,15,17	0.74	0
2	BGC	N	1	2	12,12,12	0.49	0	17,17,17	0.56	0
2	BGC	N	2	2	11,11,12	0.26	0	15,15,17	0.71	0
2	BGC	O	1	2	12,12,12	0.53	0	17,17,17	0.47	0
2	BGC	O	2	2	11,11,12	0.28	0	15,15,17	0.75	0
2	BGC	P	1	2	12,12,12	0.49	0	17,17,17	0.50	0
2	BGC	P	2	2	11,11,12	0.29	0	15,15,17	0.77	0
2	BGC	Q	1	2	12,12,12	0.52	0	17,17,17	0.83	0
2	BGC	Q	2	2	11,11,12	0.25	0	15,15,17	0.83	0
2	BGC	R	1	2	12,12,12	0.49	0	17,17,17	0.58	0
2	BGC	R	2	2	11,11,12	0.27	0	15,15,17	0.80	0
2	BGC	S	1	2	12,12,12	0.55	0	17,17,17	0.82	0
2	BGC	S	2	2	11,11,12	0.26	0	15,15,17	0.84	1 (6%)
2	BGC	T	1	2	12,12,12	0.52	0	17,17,17	0.38	0
2	BGC	T	2	2	11,11,12	0.30	0	15,15,17	0.55	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	BGC	K	1	2	-	0/2/22/22	0/1/1/1
2	BGC	K	2	2	-	0/2/19/22	0/1/1/1
2	BGC	L	1	2	-	0/2/22/22	0/1/1/1
2	BGC	L	2	2	-	0/2/19/22	0/1/1/1
2	BGC	M	1	2	-	0/2/22/22	0/1/1/1
2	BGC	M	2	2	-	0/2/19/22	0/1/1/1
2	BGC	N	1	2	-	0/2/22/22	0/1/1/1
2	BGC	N	2	2	-	0/2/19/22	0/1/1/1
2	BGC	O	1	2	-	2/2/22/22	0/1/1/1
2	BGC	O	2	2	-	2/2/19/22	0/1/1/1
2	BGC	P	1	2	-	0/2/22/22	0/1/1/1
2	BGC	P	2	2	-	0/2/19/22	0/1/1/1
2	BGC	Q	1	2	-	0/2/22/22	0/1/1/1
2	BGC	Q	2	2	-	0/2/19/22	0/1/1/1
2	BGC	R	1	2	-	2/2/22/22	0/1/1/1
2	BGC	R	2	2	-	1/2/19/22	0/1/1/1
2	BGC	S	1	2	-	0/2/22/22	0/1/1/1
2	BGC	S	2	2	-	2/2/19/22	0/1/1/1
2	BGC	T	1	2	-	0/2/22/22	0/1/1/1
2	BGC	T	2	2	-	1/2/19/22	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	S	2	BGC	O5-C5-C6	2.32	110.83	107.20
2	L	2	BGC	C1-C2-C3	2.01	112.13	109.67

There are no chirality outliers.

All (10) torsion outliers are listed below:

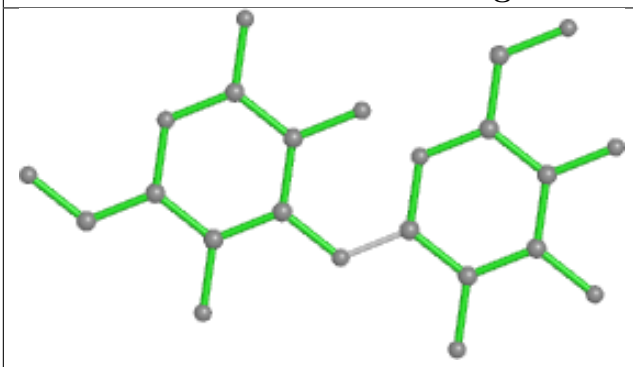
Mol	Chain	Res	Type	Atoms
2	O	1	BGC	O5-C5-C6-O6
2	O	1	BGC	C4-C5-C6-O6
2	O	2	BGC	O5-C5-C6-O6
2	S	2	BGC	O5-C5-C6-O6
2	R	1	BGC	O5-C5-C6-O6
2	O	2	BGC	C4-C5-C6-O6
2	T	2	BGC	O5-C5-C6-O6
2	R	2	BGC	O5-C5-C6-O6
2	S	2	BGC	C4-C5-C6-O6
2	R	1	BGC	C4-C5-C6-O6

There are no ring outliers.

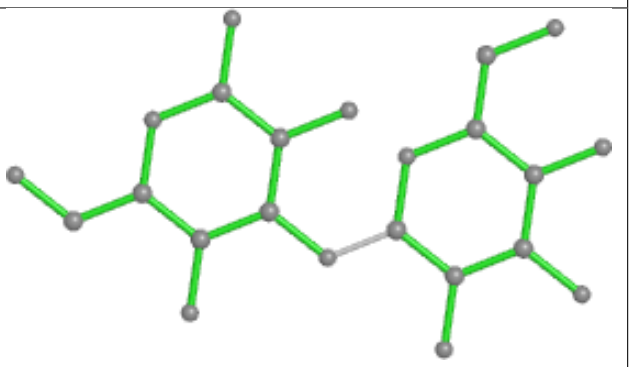
No monomer is involved in short contacts.

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

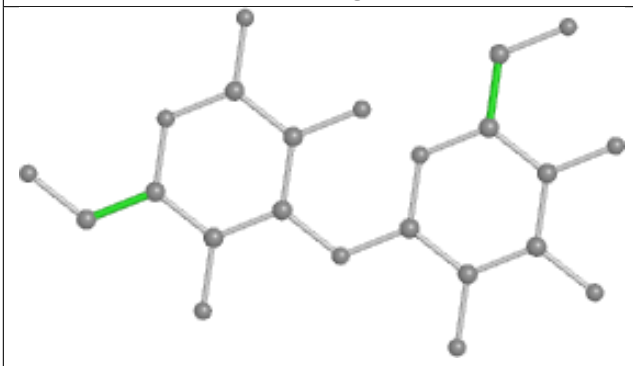
Oligosaccharide Chain K



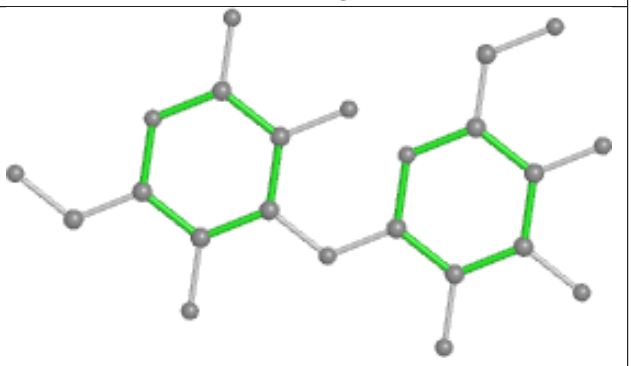
Bond lengths



Bond angles

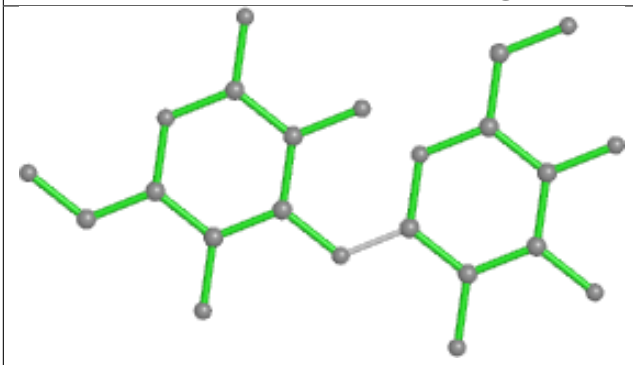


Torsions

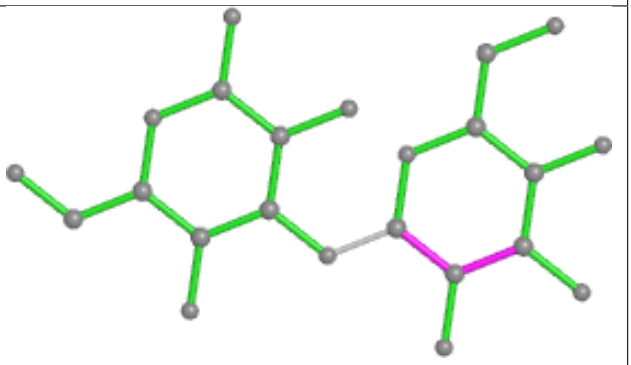


Rings

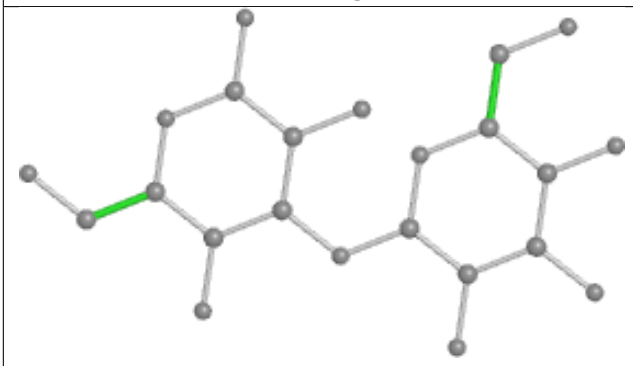
Oligosaccharide Chain L



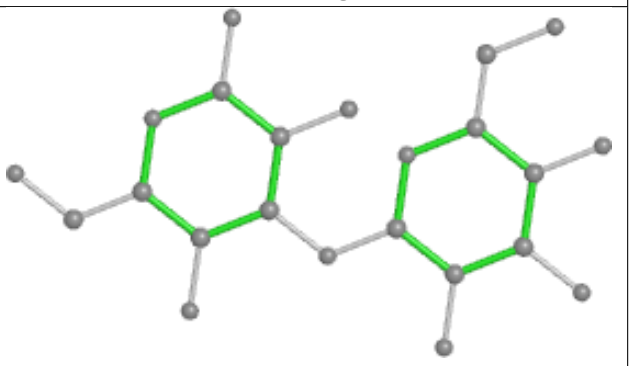
Bond lengths



Bond angles

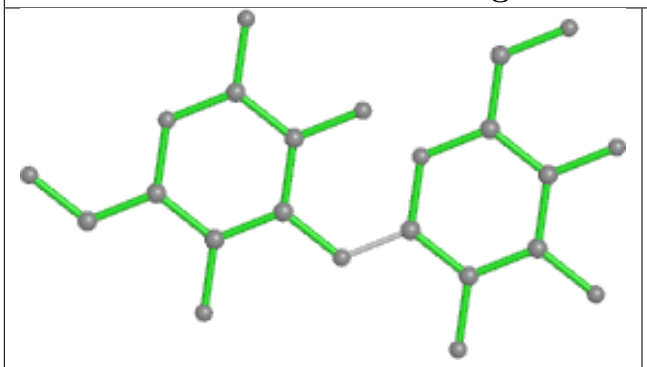


Torsions

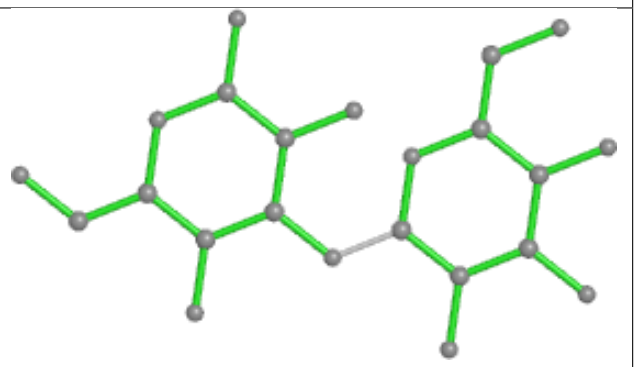


Rings

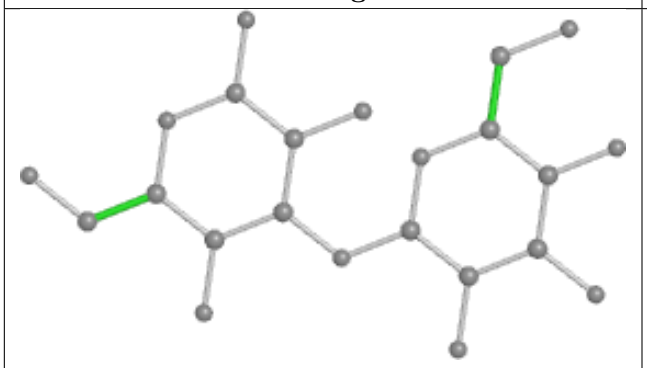
Oligosaccharide Chain M



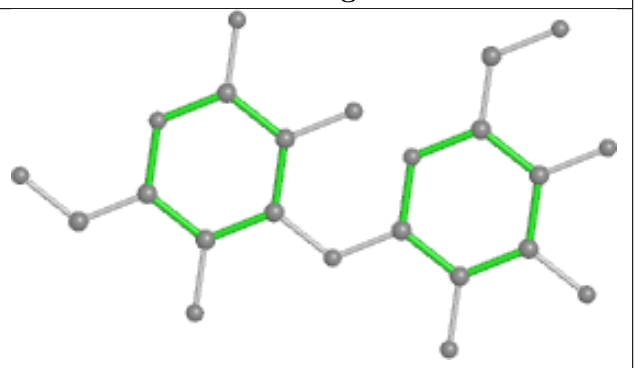
Bond lengths



Bond angles

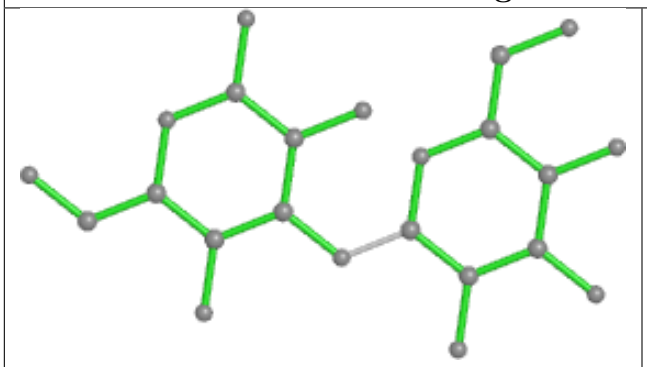


Torsions

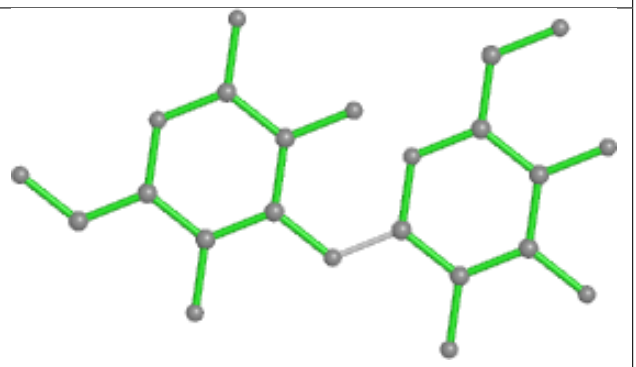


Rings

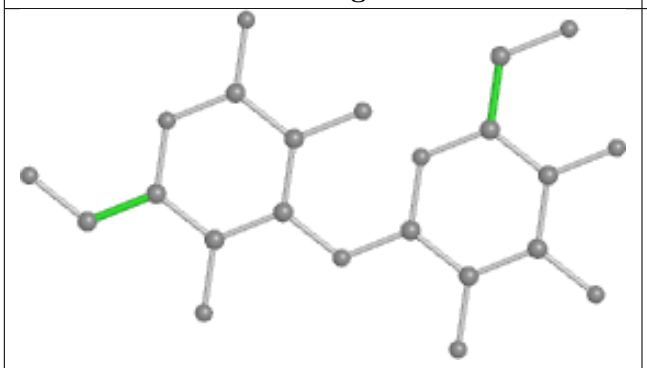
Oligosaccharide Chain N



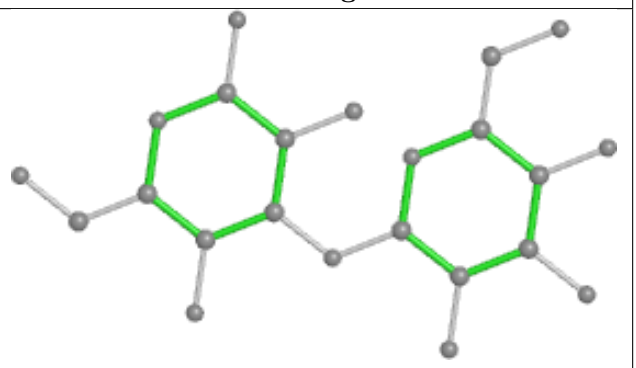
Bond lengths



Bond angles

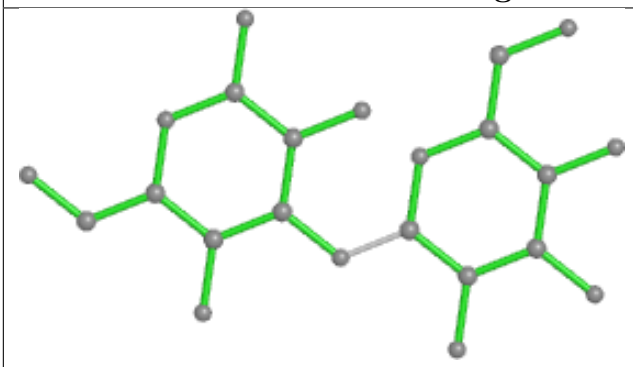


Torsions

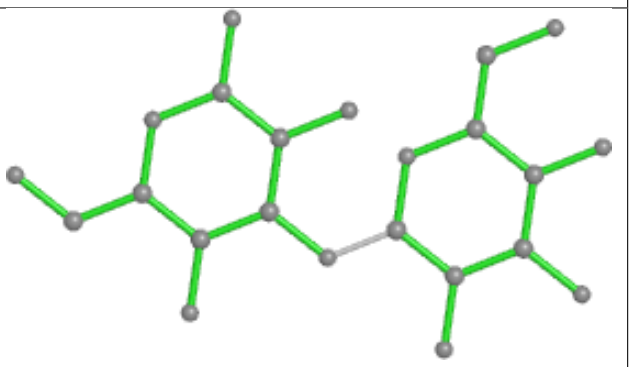


Rings

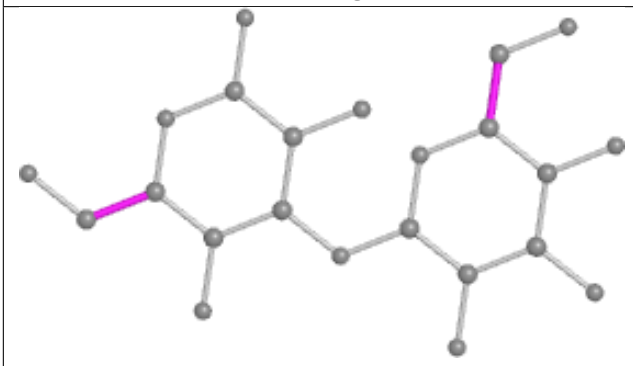
Oligosaccharide Chain O



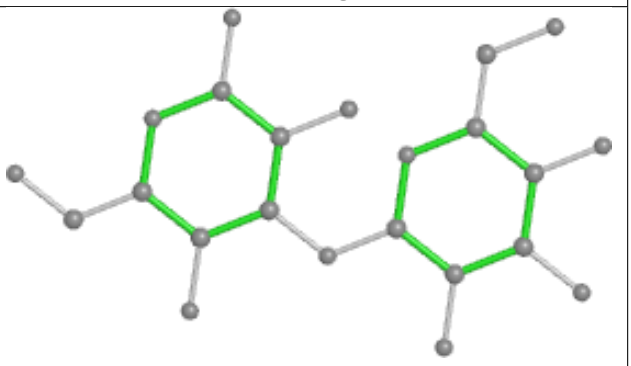
Bond lengths



Bond angles

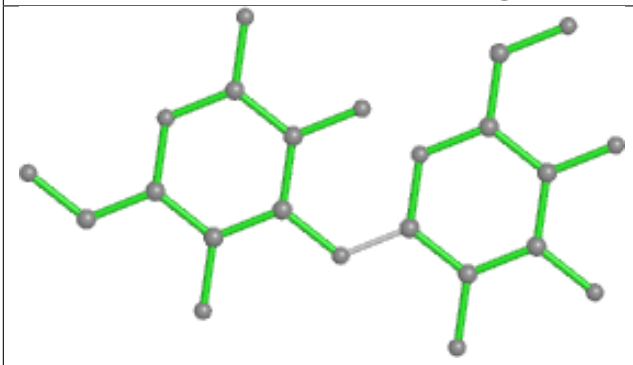


Torsions

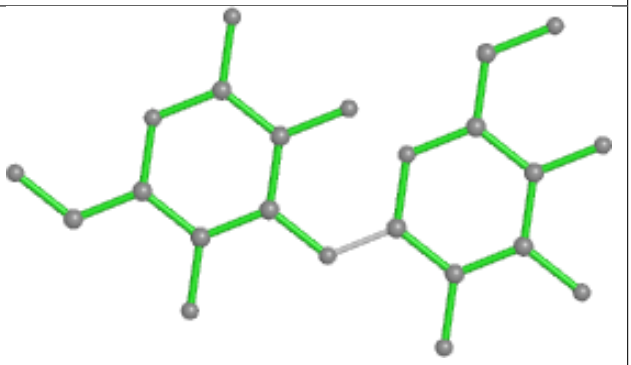


Rings

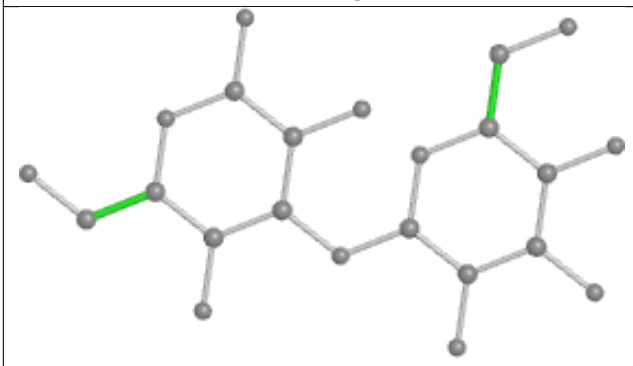
Oligosaccharide Chain P



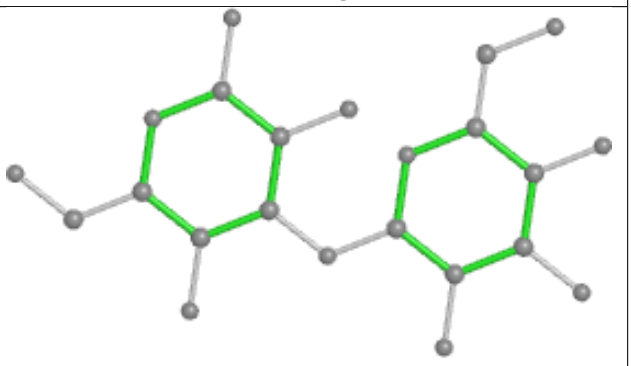
Bond lengths



Bond angles

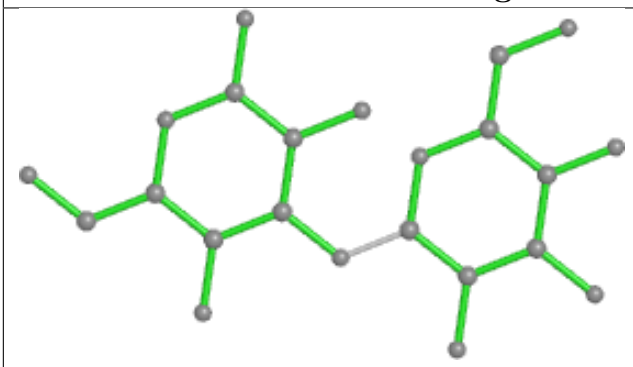


Torsions

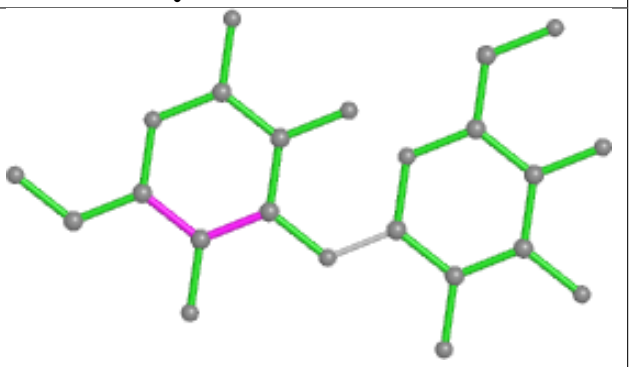


Rings

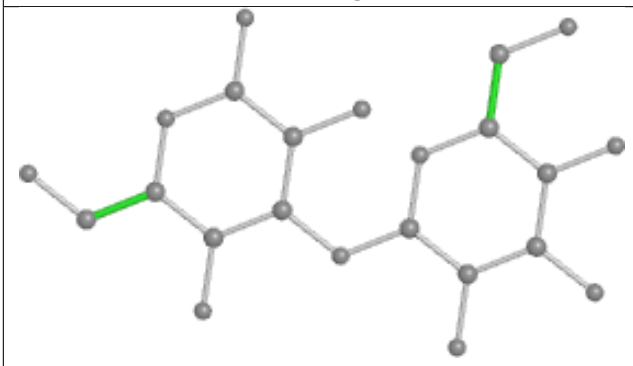
Oligosaccharide Chain Q



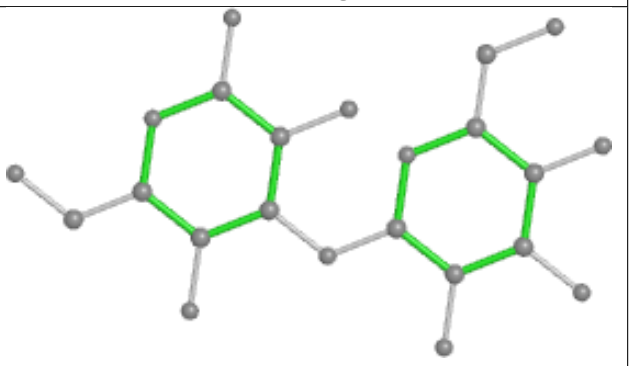
Bond lengths



Bond angles

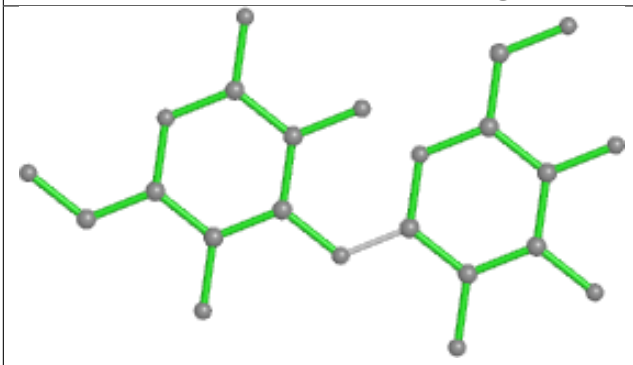


Torsions

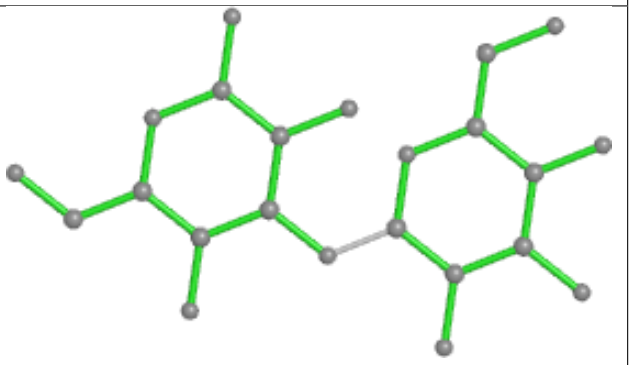


Rings

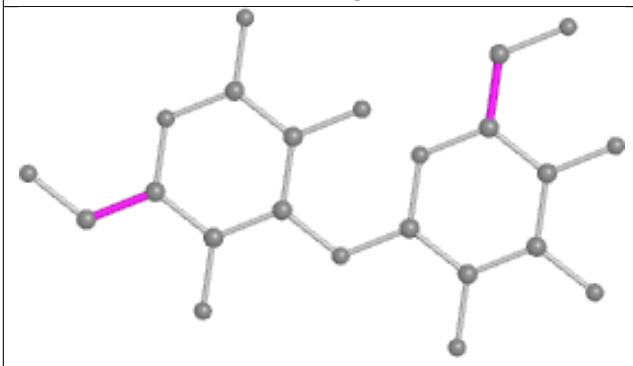
Oligosaccharide Chain R



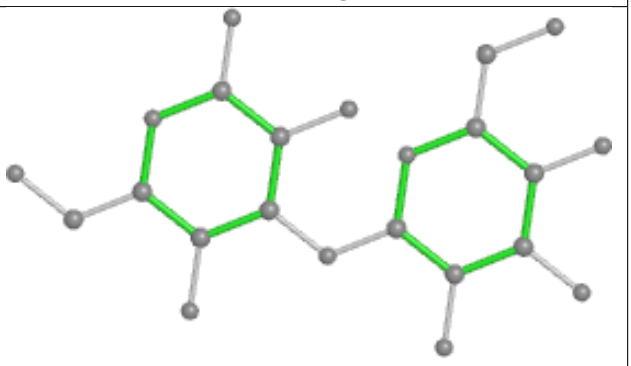
Bond lengths



Bond angles

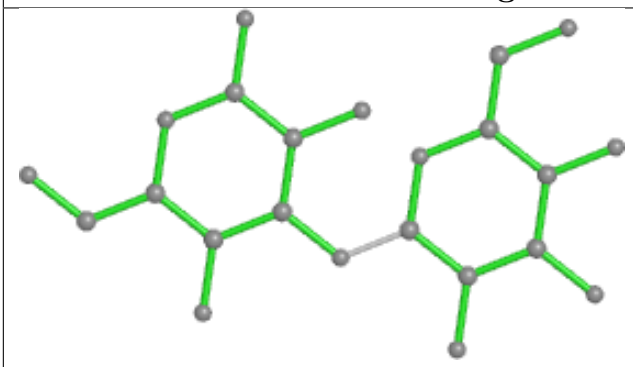


Torsions

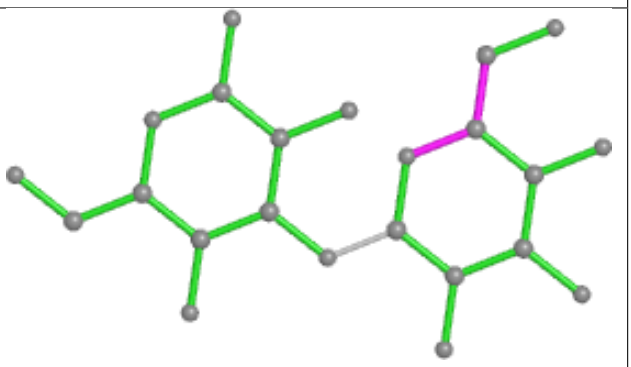


Rings

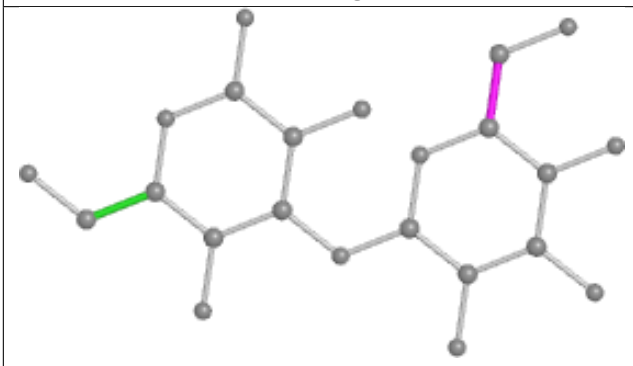
Oligosaccharide Chain S



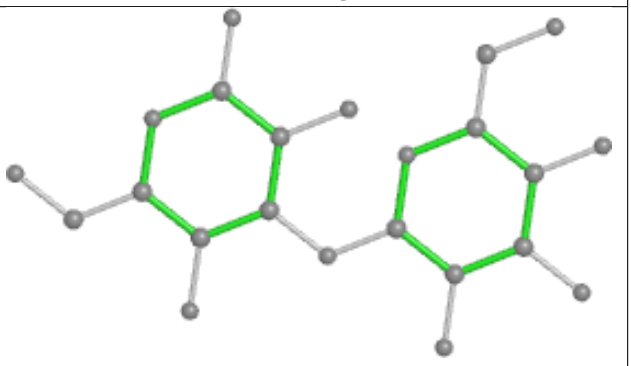
Bond lengths



Bond angles

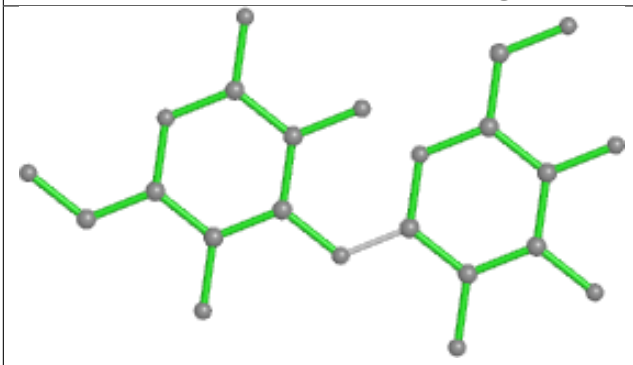


Torsions

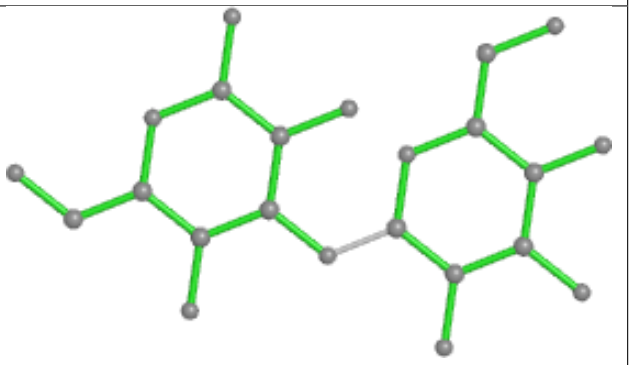


Rings

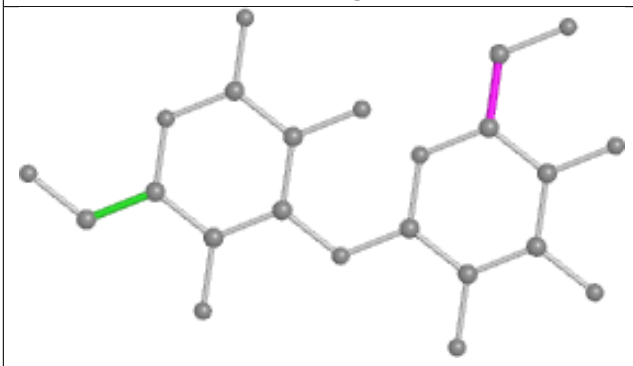
Oligosaccharide Chain T



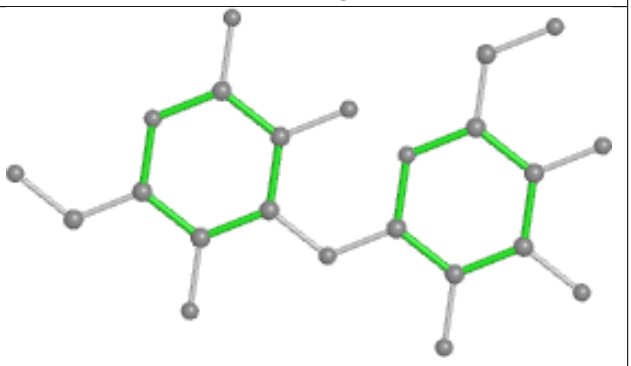
Bond lengths



Bond angles



Torsions



Rings

5.6 Ligand geometry (i)

There are no ligands in this entry.

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 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	247/266 (92%)	0.11	0	100	100	22, 32, 40, 48	0
1	B	247/266 (92%)	0.17	0	100	100	22, 33, 41, 47	0
1	C	249/266 (93%)	0.20	2 (0%)	86	91	22, 33, 44, 49	0
1	D	247/266 (92%)	0.43	6 (2%)	59	68	28, 41, 58, 60	0
1	E	247/266 (92%)	0.17	1 (0%)	92	96	24, 35, 47, 52	0
1	F	247/266 (92%)	0.22	2 (0%)	86	91	22, 34, 44, 47	0
1	G	247/266 (92%)	0.15	0	100	100	22, 33, 42, 46	0
1	H	247/266 (92%)	0.35	9 (3%)	42	55	22, 36, 47, 50	0
1	I	247/266 (92%)	0.43	8 (3%)	47	59	29, 42, 62, 66	0
1	J	247/266 (92%)	0.45	4 (1%)	72	80	30, 44, 68, 77	0
All	All	2472/2660 (92%)	0.27	32 (1%)	77	84	22, 36, 55, 77	0

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	I	58	ASN	3.4
1	H	134	ALA	2.9
1	H	217	PHE	2.9
1	H	131	ASN	2.9
1	I	66	GLN	2.9
1	H	70	LYS	2.7
1	I	131	ASN	2.7
1	C	124	ALA	2.7
1	H	135	LYS	2.7
1	H	124	ALA	2.6
1	D	155	ARG	2.6
1	H	126	ILE	2.6
1	I	126	ILE	2.6

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Mol	Chain	Res	Type	RSRZ
1	I	152	GLU	2.5
1	H	141	LEU	2.5
1	E	152	GLU	2.5
1	J	128	LYS	2.4
1	H	133	VAL	2.4
1	F	133	VAL	2.4
1	D	164	HIS	2.3
1	D	157	LEU	2.3
1	C	217	PHE	2.2
1	J	200	TRP	2.2
1	D	71	THR	2.2
1	I	132	LEU	2.2
1	D	166	HIS	2.2
1	I	121	ALA	2.2
1	F	131	ASN	2.1
1	D	162	LYS	2.1
1	I	124	ALA	2.1
1	J	19	TRP	2.1
1	J	152	GLU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains (i)

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

6.3 Carbohydrates (i)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	BGC	Q	1	12/12	0.81	0.22	26,27,28,28	0
2	BGC	L	2	11/12	0.84	0.15	25,26,26,26	0
2	BGC	R	1	12/12	0.84	0.16	26,28,28,29	0
2	BGC	M	1	12/12	0.86	0.14	25,25,26,26	0
2	BGC	P	1	12/12	0.86	0.20	26,27,27,28	0
2	BGC	T	2	11/12	0.86	0.14	25,26,26,26	0
2	BGC	O	1	12/12	0.87	0.16	25,26,26,27	0
2	BGC	L	1	12/12	0.88	0.22	26,27,27,28	0
2	BGC	Q	2	11/12	0.88	0.17	25,26,26,26	0

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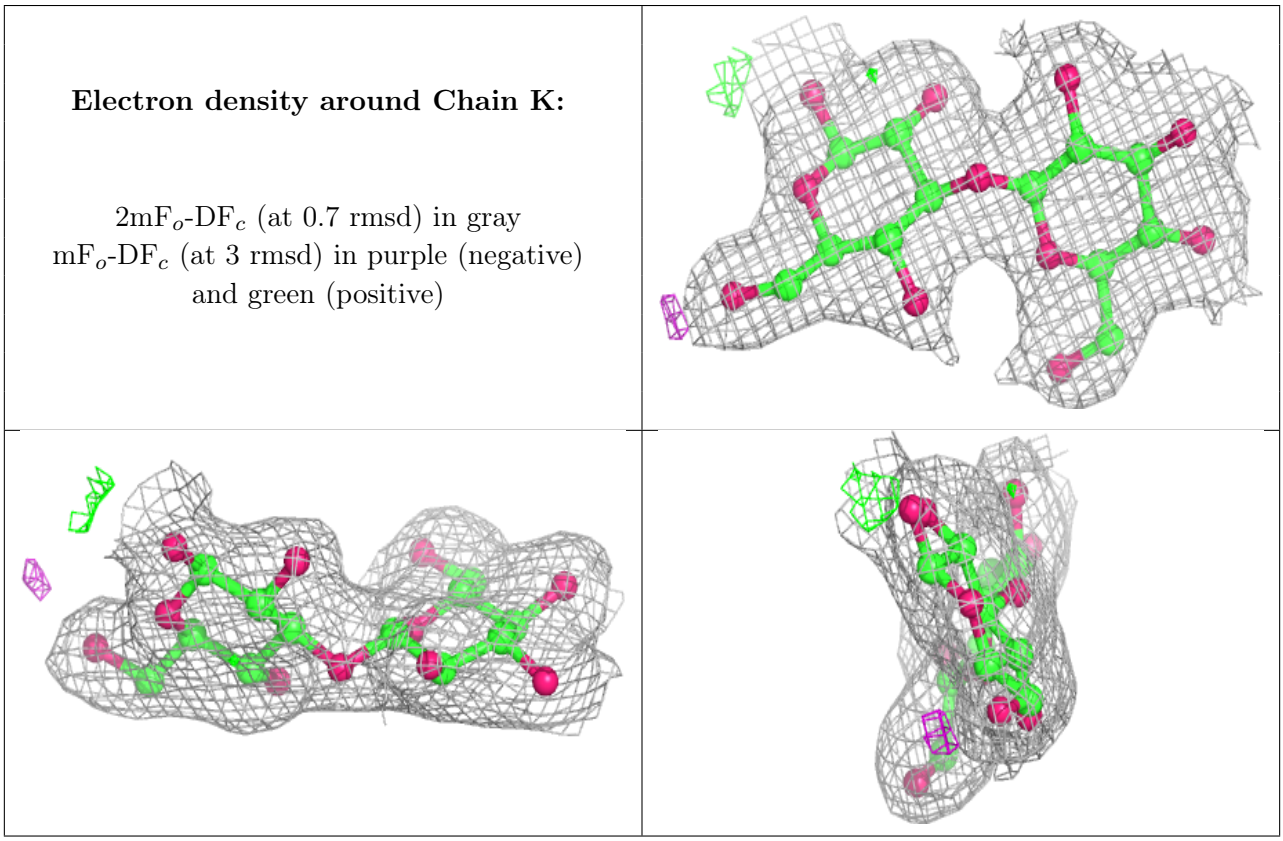
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	BGC	S	1	12/12	0.89	0.15	25,26,26,27	0
2	BGC	T	1	12/12	0.89	0.17	26,27,27,28	0
2	BGC	R	2	11/12	0.89	0.14	26,26,26,27	0
2	BGC	K	2	11/12	0.90	0.12	24,25,25,25	0
2	BGC	M	2	11/12	0.90	0.15	24,24,25,25	0
2	BGC	N	1	12/12	0.90	0.19	27,28,28,29	0
2	BGC	K	1	12/12	0.91	0.15	25,26,26,27	0
2	BGC	N	2	11/12	0.92	0.13	26,26,27,27	0
2	BGC	S	2	11/12	0.93	0.10	24,25,25,25	0
2	BGC	O	2	11/12	0.94	0.09	25,25,25,25	0
2	BGC	P	2	11/12	0.94	0.10	25,25,26,26	0

The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

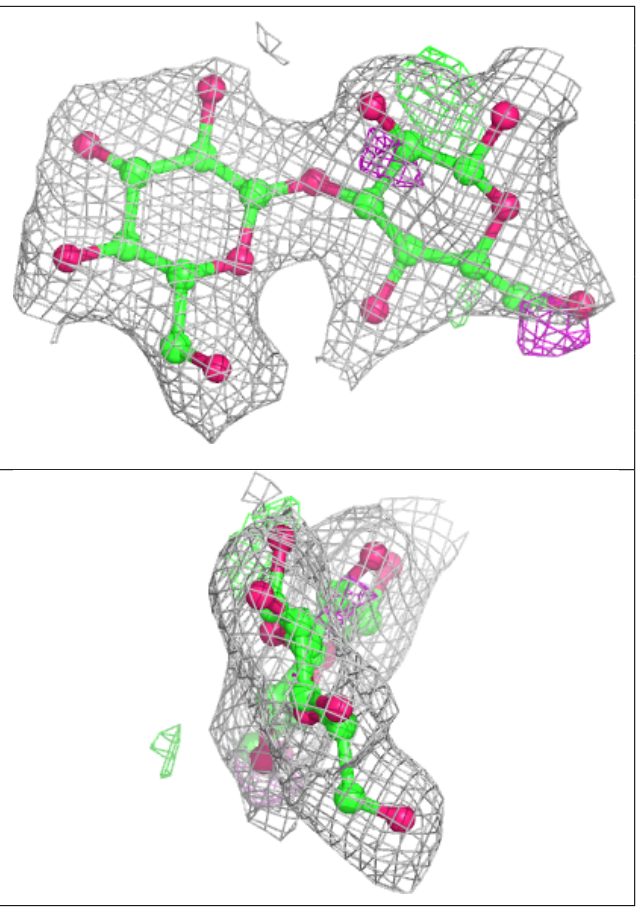
Electron density around Chain K:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)



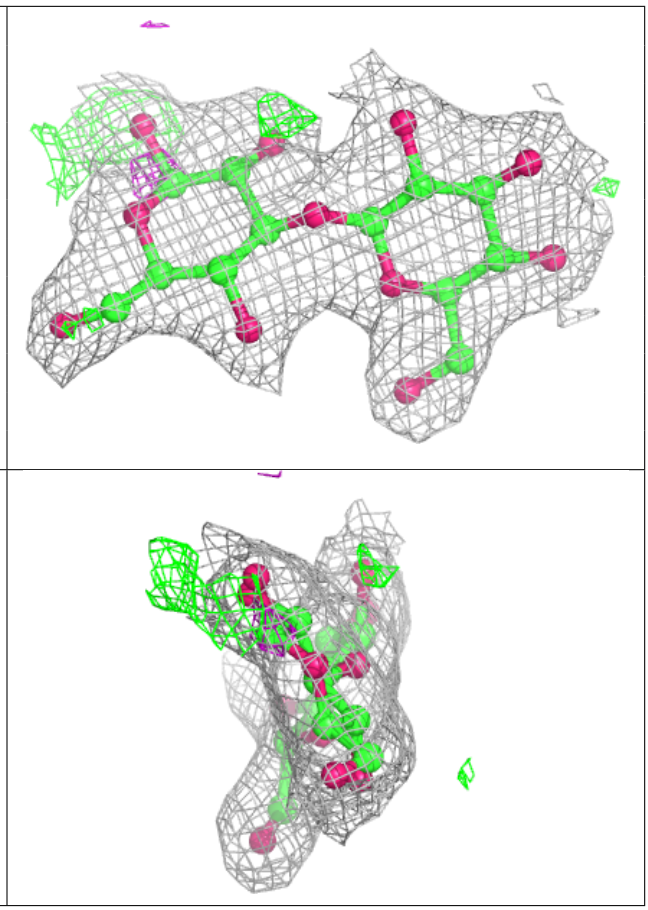
Electron density around Chain L:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



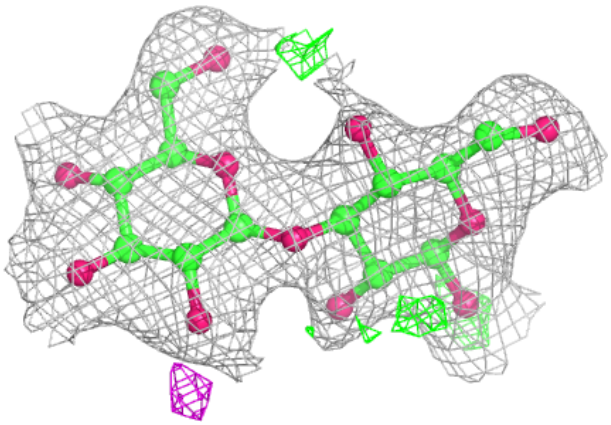
Electron density around Chain M:

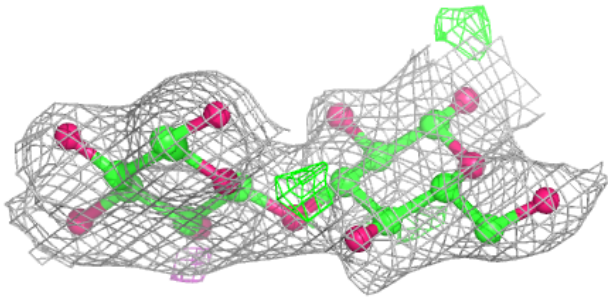
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

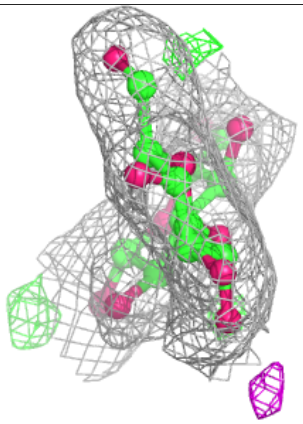


Electron density around Chain N:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

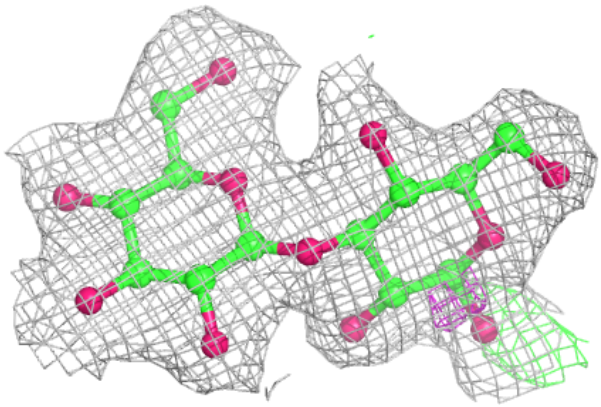


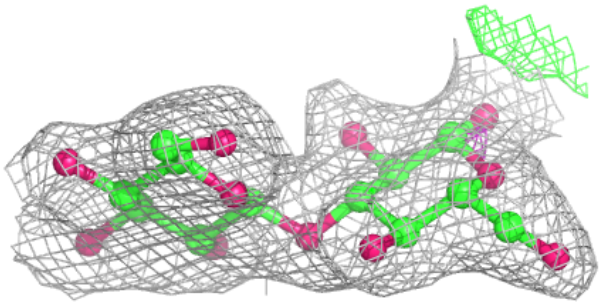


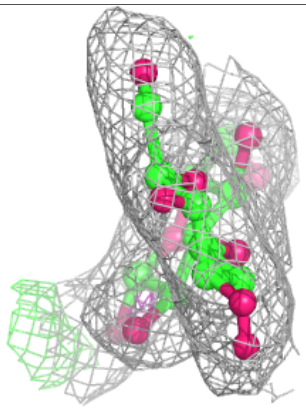


Electron density around Chain O:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

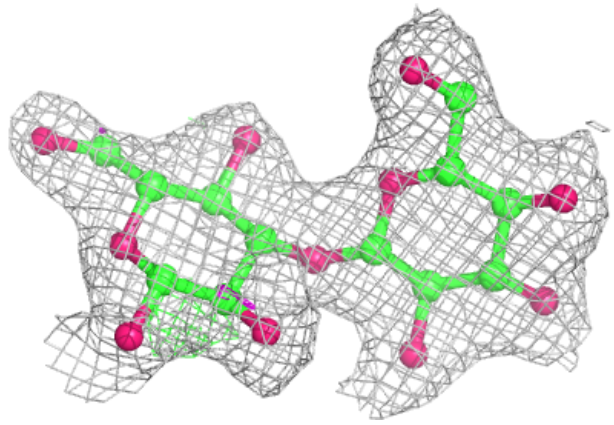


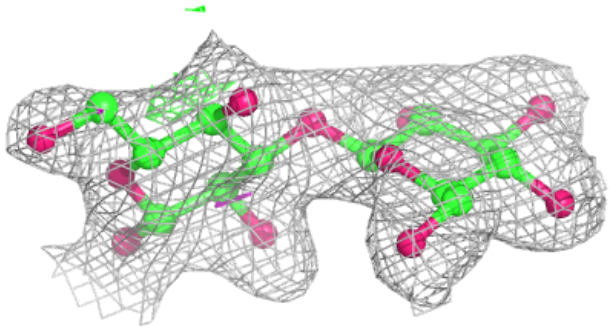


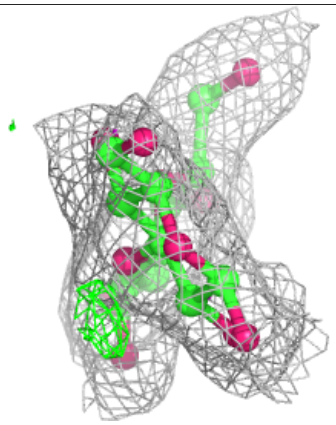


Electron density around Chain P:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

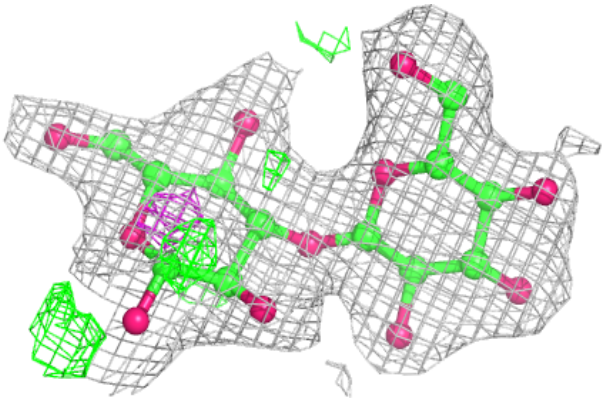


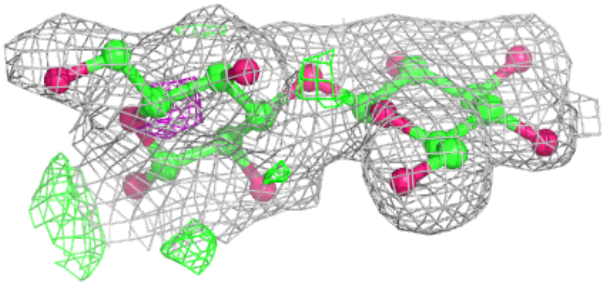


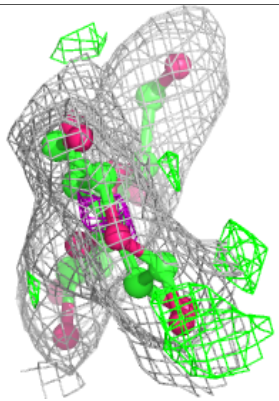


Electron density around Chain Q:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

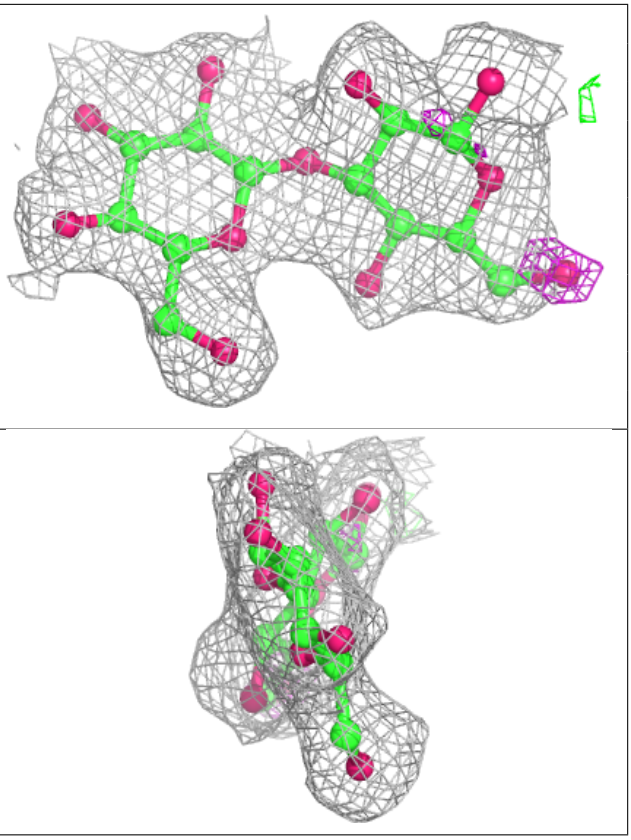


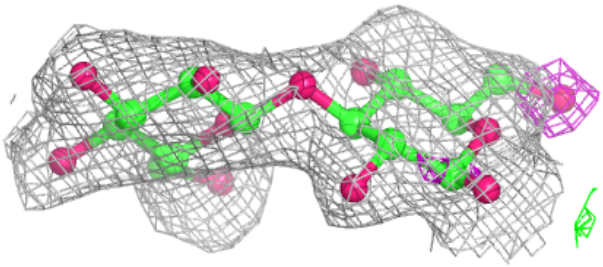




Electron density around Chain R:

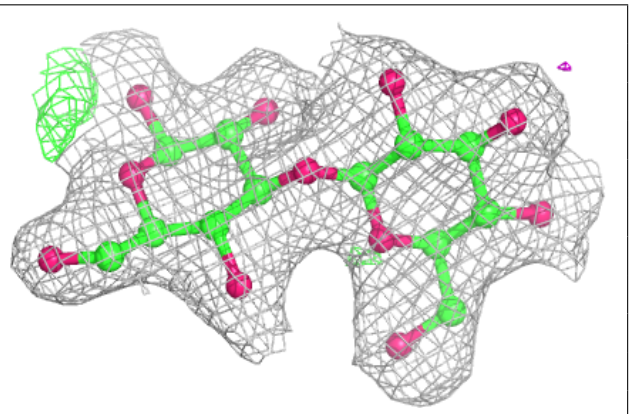
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

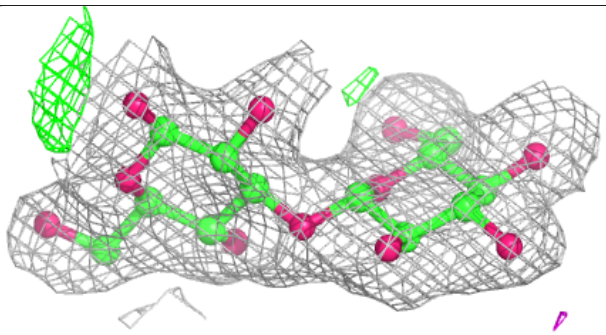


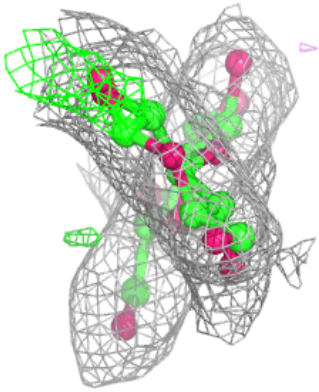


Electron density around Chain S:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

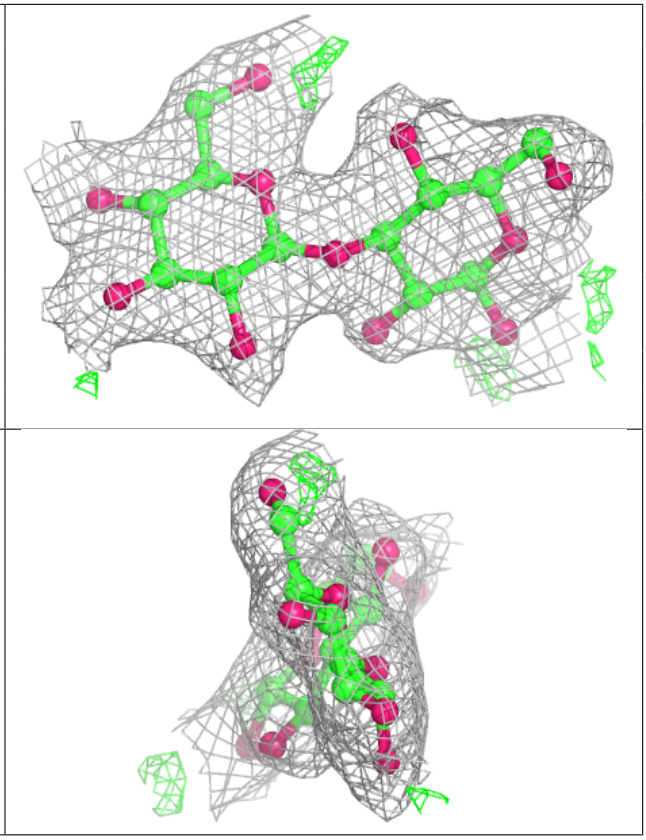






Electron density around Chain T:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.