wwPDB X-ray Structure Validation Summary Report (i)

Jun 12, 2024 – 04:35 PM EDT

PDB ID	:	1CVI
Title	:	CRYSTAL STRUCTURE OF HUMAN PROSTATIC ACID PHOSPHATASE
Authors	:	Jakob, C.G.; Lewinski, K.; Kuciel, R.; Ostrowski, W.; Lebioda, L.
Deposited on	:	1999-08-23
Resolution	:	3.20 Å(reported)

This is a wwPDB X-ray Structure 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/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:

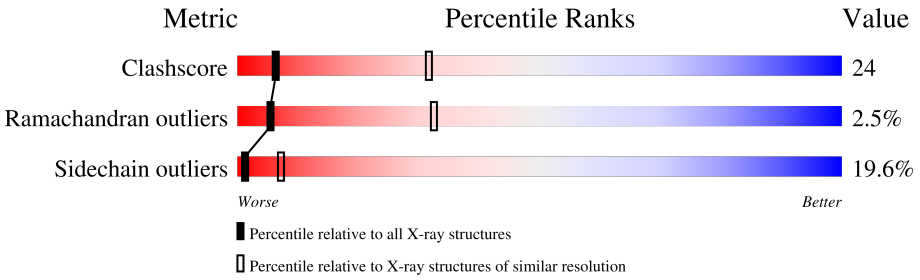
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 3.20 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	342	46% 43% 11%
1	B	342	46% 42% 12%
1	C	342	43% 47% 10%
1	D	342	45% 44% 11%
2	E	8	25% 25% 50%
3	F	4	50% 50%
4	G	3	33% 33% 33%
4	J	3	67% 33%

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Mol	Chain	Length	Quality of chain
4	K	3	33% 67%
4	M	3	33% 67%
5	H	4	25% 75%
5	L	4	25% 75%
6	I	6	33% 50% 17%

2 Entry composition

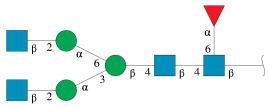
There are 9 unique types of molecules in this entry. The entry contains 12376 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 PROSTATIC ACID PHOSPHATASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	342	Total	C	N	O	S	0	0	0
			2801	1807	461	517	16			
1	B	342	Total	C	N	O	S	0	0	0
			2801	1807	461	517	16			
1	C	342	Total	C	N	O	S	0	0	0
			2801	1807	461	517	16			
1	D	342	Total	C	N	O	S	0	0	0
			2801	1807	461	517	16			

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

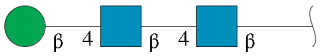


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	8	Total	C	N	O	0	0	0
			99	56	4	39			

- Molecule 3 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	ZeroOcc	AltConf	Trace
3	F	4	Total C N O 50 28 2 20	0	0	0

- Molecule 4 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.

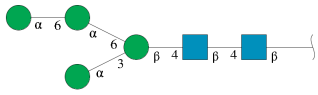


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
4	G	3	Total C N O 39 22 2 15	0	0	0
4	J	3	Total C N O 39 22 2 15	0	0	0
4	K	3	Total C N O 39 22 2 15	0	0	0
4	M	3	Total C N O 39 22 2 15	0	0	0

- Molecule 5 is an oligosaccharide called 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.

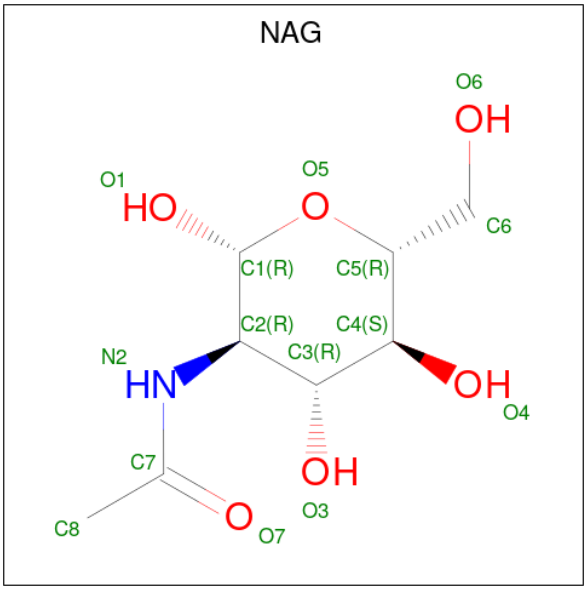
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf	Trace
5	H	4	Total C N O 50 28 2 20	0	0	0
5	L	4	Total C N O 50 28 2 20	0	0	0

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



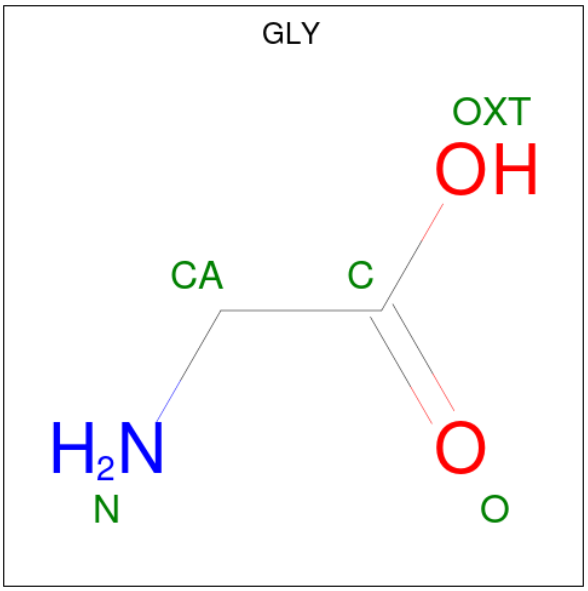
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	I	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	A	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		
7	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 8 is GLYCINE (three-letter code: GLY) (formula: $C_2H_5NO_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	1	Total C N O 5 2 1 2	0	0
8	B	1	Total C N O 5 2 1 2	0	0
8	C	1	Total C N O 5 2 1 2	0	0
8	D	1	Total C N O 5 2 1 2	0	0

- Molecule 9 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
9	A	163	Total O 163 163	0	0
9	B	188	Total O 188 188	0	0
9	C	144	Total O 144 144	0	0
9	D	138	Total O 138 138	0	0

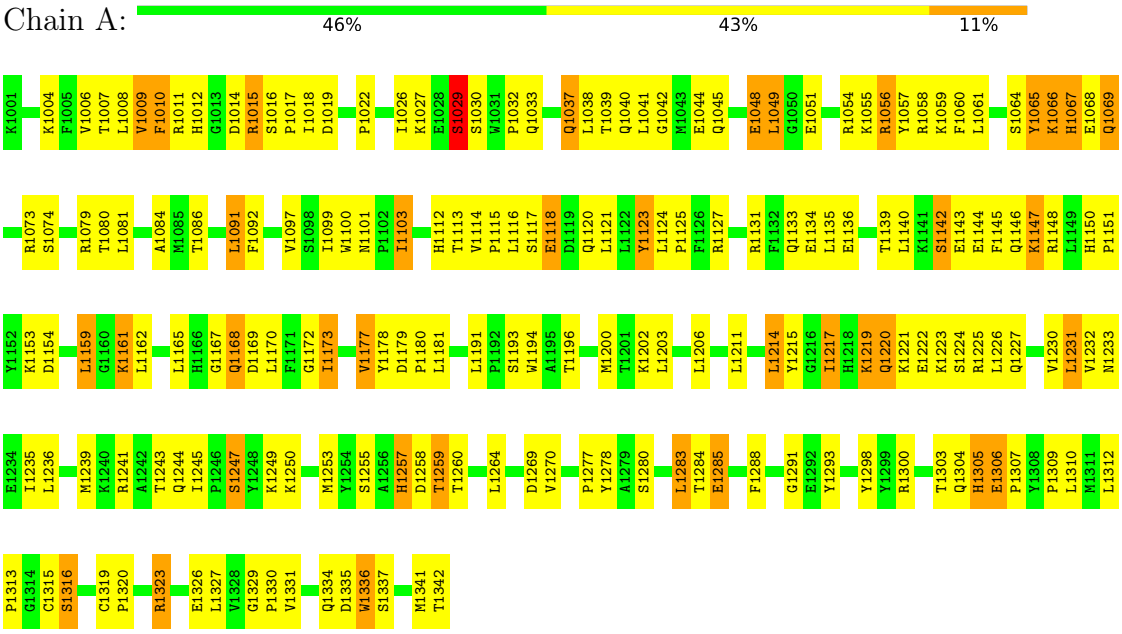
3 Residue-property plots (i)

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. 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. 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.

Note EDS was not executed.

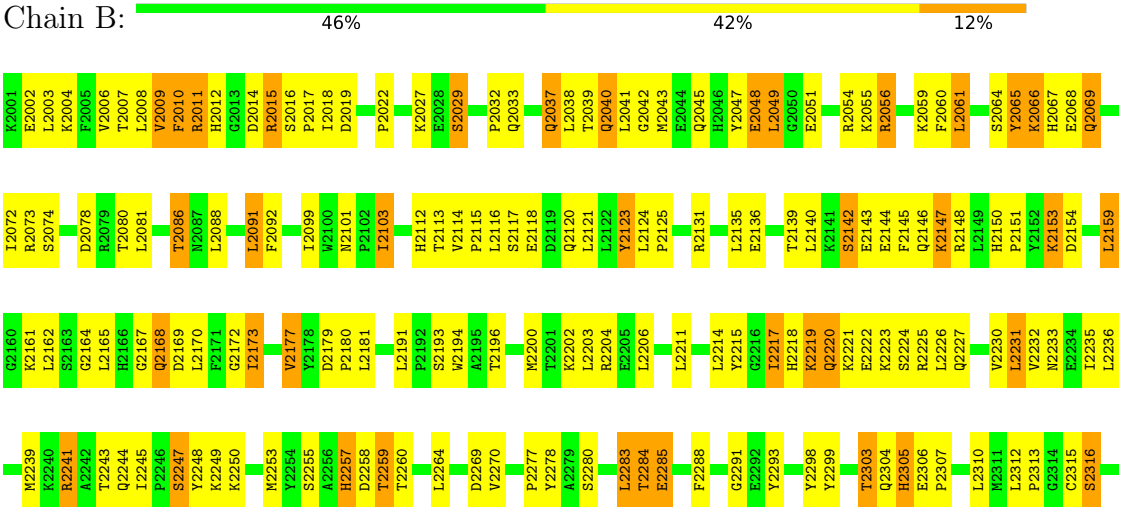
• Molecule 1: PROSTATIC ACID PHOSPHATASE

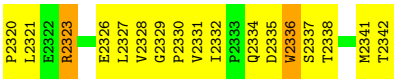
Chain A:



• Molecule 1: PROSTATIC ACID PHOSPHATASE

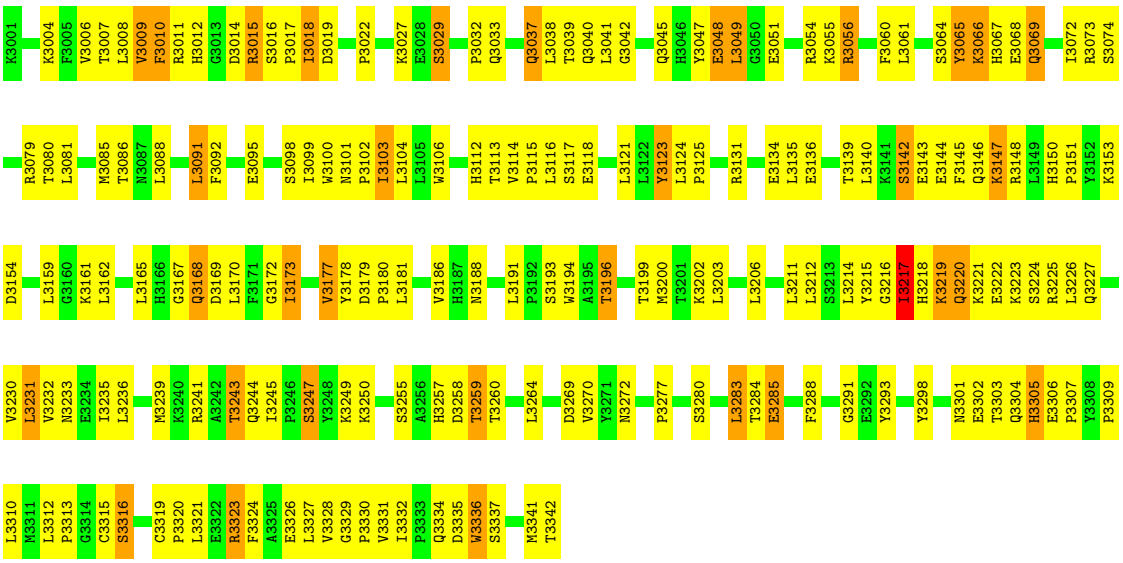
Chain B:





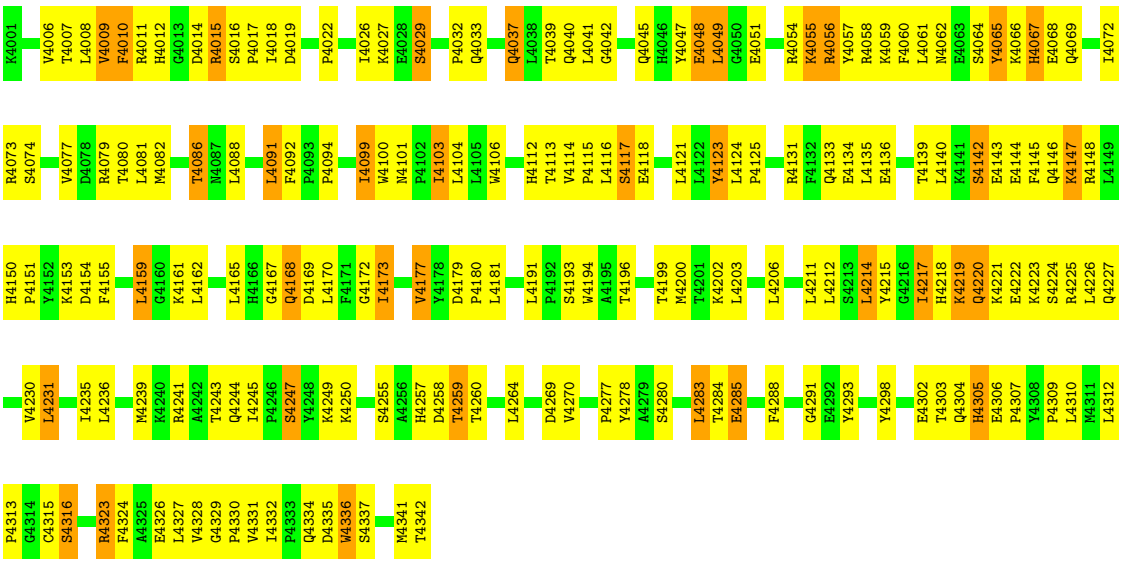
• Molecule 1: PROSTATIC ACID PHOSPHATASE

Chain C:



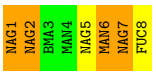
• Molecule 1: PROSTATIC ACID PHOSPHATASE

Chain D:



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

Chain E:

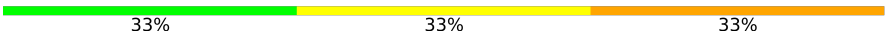


• Molecule 3: alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:

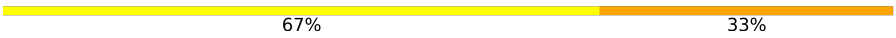
• Molecule 4: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G:



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

Chain J:

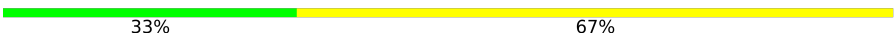


• Molecule 4: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:

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

Chain M:

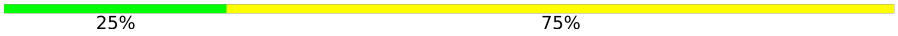


• Molecule 5: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:

• Molecule 5: 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 6: alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)-[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 I:

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	126.33Å 207.96Å 73.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	8.00 – 3.20	Depositor
% Data completeness (in resolution range)	91.7 (8.00-3.20)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.157 , 0.274	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	12376	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| > 5$	RMSZ	$\#\|Z\| > 5$
1	A	0.33	0/2883	0.53	0/3914
1	B	0.32	0/2883	0.53	0/3914
1	C	0.33	0/2883	0.53	0/3914
1	D	0.33	0/2883	0.54	0/3914
All	All	0.33	0/11532	0.53	0/15656

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2801	0	2745	141	0
1	B	2801	0	2745	139	0
1	C	2801	0	2745	154	0
1	D	2801	0	2745	135	0
2	E	99	0	85	9	0
3	F	50	0	43	1	0
4	G	39	0	34	4	0
4	J	39	0	34	3	0
4	K	39	0	34	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	M	39	0	34	1	0
5	H	50	0	43	3	0
5	L	50	0	43	2	0
6	I	72	0	61	2	0
7	A	14	0	13	0	0
7	C	14	0	13	0	0
7	D	14	0	13	0	0
8	A	5	0	2	1	0
8	B	5	0	2	0	0
8	C	5	0	2	1	0
8	D	5	0	2	2	0
9	A	163	0	0	11	0
9	B	188	0	0	12	0
9	C	144	0	0	10	0
9	D	138	0	0	11	0
All	All	12376	0	11438	561	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 24.

The worst 5 of 561 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:C:3159:LEU:HD11	1:C:3203:LEU:HD21	1.39	1.05
1:D:4159:LEU:HD11	1:D:4203:LEU:HD21	1.42	1.02
1:B:2200:MET:HE2	1:B:2200:MET:HA	1.44	0.99
1:A:1159:LEU:HD11	1:A:1203:LEU:HD21	1.44	0.99
1:B:2159:LEU:HD11	1:B:2203:LEU:HD21	1.41	0.98

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	340/342 (99%)	298 (88%)	33 (10%)	9 (3%)	5 31
1	B	340/342 (99%)	298 (88%)	34 (10%)	8 (2%)	6 34
1	C	340/342 (99%)	298 (88%)	34 (10%)	8 (2%)	6 34
1	D	340/342 (99%)	301 (88%)	30 (9%)	9 (3%)	5 31
All	All	1360/1368 (99%)	1195 (88%)	131 (10%)	34 (2%)	5 32

5 of 34 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1065	TYR
1	A	1217	ILE
1	A	1220	GLN
1	B	2065	TYR
1	B	2217	ILE

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	314/314 (100%)	252 (80%)	62 (20%)	1 7
1	B	314/314 (100%)	251 (80%)	63 (20%)	1 6
1	C	314/314 (100%)	255 (81%)	59 (19%)	1 8
1	D	314/314 (100%)	252 (80%)	62 (20%)	1 7
All	All	1256/1256 (100%)	1010 (80%)	246 (20%)	1 7

5 of 246 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	2283	LEU
1	D	4193	SER
1	C	3068	GLU
1	D	4177	VAL
1	D	4259	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
1	D	4037	GLN
1	D	4334	GLN
1	D	4265	GLN
1	B	2265	GLN
1	C	3334	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)

38 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	NAG	E	1	2,1	14,14,15	0.51	0	17,19,21	1.01	1 (5%)
2	NAG	E	2	2	14,14,15	0.61	0	17,19,21	0.80	1 (5%)
2	BMA	E	3	2	11,11,12	0.57	0	15,15,17	0.58	0
2	MAN	E	4	2	11,11,12	0.79	0	15,15,17	0.82	0
2	NAG	E	5	2	14,14,15	0.50	0	17,19,21	0.75	1 (5%)
2	MAN	E	6	2	11,11,12	0.72	0	15,15,17	0.63	1 (6%)
2	NAG	E	7	2	14,14,15	0.47	0	17,19,21	0.76	1 (5%)
2	FUC	E	8	2	10,10,11	0.60	0	14,14,16	0.45	0
3	NAG	F	1	1,3	14,14,15	0.57	0	17,19,21	0.64	0
3	NAG	F	2	3	14,14,15	0.79	0	17,19,21	1.32	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
3	BMA	F	3	3	11,11,12	0.58	0	15,15,17	0.98	1 (6%)
3	MAN	F	4	3	11,11,12	0.73	0	15,15,17	0.59	0
4	NAG	G	1	1,4	14,14,15	0.49	0	17,19,21	0.99	1 (5%)
4	NAG	G	2	4	14,14,15	0.59	0	17,19,21	0.59	0
4	BMA	G	3	4	11,11,12	0.54	0	15,15,17	0.21	0
5	NAG	H	1	1,5	14,14,15	0.64	0	17,19,21	0.62	0
5	NAG	H	2	5	14,14,15	0.64	0	17,19,21	0.77	0
5	BMA	H	3	5	11,11,12	0.66	0	15,15,17	0.34	0
5	MAN	H	4	5	11,11,12	0.66	0	15,15,17	0.65	1 (6%)
6	NAG	I	1	1,6	14,14,15	0.46	0	17,19,21	0.64	0
6	NAG	I	2	6	14,14,15	0.64	0	17,19,21	0.93	2 (11%)
6	BMA	I	3	6	11,11,12	0.73	0	15,15,17	0.44	0
6	MAN	I	4	6	11,11,12	0.74	0	15,15,17	0.77	1 (6%)
6	MAN	I	5	6	11,11,12	0.69	0	15,15,17	0.69	0
6	MAN	I	6	6	11,11,12	0.61	0	15,15,17	0.90	1 (6%)
4	NAG	J	1	1,4	14,14,15	0.65	0	17,19,21	0.82	1 (5%)
4	NAG	J	2	4	14,14,15	0.66	0	17,19,21	0.83	0
4	BMA	J	3	4	11,11,12	0.53	0	15,15,17	0.33	0
4	NAG	K	1	1,4	14,14,15	0.67	0	17,19,21	1.24	3 (17%)
4	NAG	K	2	4	14,14,15	0.89	0	17,19,21	1.01	2 (11%)
4	BMA	K	3	4	11,11,12	0.58	0	15,15,17	0.59	0
5	NAG	L	1	1,5	14,14,15	0.87	0	17,19,21	0.94	0
5	NAG	L	2	5	14,14,15	0.67	0	17,19,21	0.90	0
5	BMA	L	3	5	11,11,12	0.79	0	15,15,17	0.52	0
5	MAN	L	4	5	11,11,12	0.60	0	15,15,17	0.81	1 (6%)
4	NAG	M	1	1,4	14,14,15	0.63	0	17,19,21	0.99	1 (5%)
4	NAG	M	2	4	14,14,15	0.69	0	17,19,21	0.71	0
4	BMA	M	3	4	11,11,12	0.57	0	15,15,17	0.44	0

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

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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	5	2	-	2/6/23/26	0/1/1/1
2	MAN	E	6	2	-	1/2/19/22	0/1/1/1
2	NAG	E	7	2	-	2/6/23/26	0/1/1/1
2	FUC	E	8	2	-	-	0/1/1/1
3	NAG	F	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	BMA	F	3	3	-	2/2/19/22	0/1/1/1
3	MAN	F	4	3	-	0/2/19/22	0/1/1/1
4	NAG	G	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	G	2	4	-	0/6/23/26	0/1/1/1
4	BMA	G	3	4	-	2/2/19/22	0/1/1/1
5	NAG	H	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	H	2	5	-	2/6/23/26	0/1/1/1
5	BMA	H	3	5	-	0/2/19/22	0/1/1/1
5	MAN	H	4	5	-	2/2/19/22	0/1/1/1
6	NAG	I	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	I	2	6	-	0/6/23/26	0/1/1/1
6	BMA	I	3	6	-	2/2/19/22	0/1/1/1
6	MAN	I	4	6	-	2/2/19/22	0/1/1/1
6	MAN	I	5	6	-	0/2/19/22	0/1/1/1
6	MAN	I	6	6	-	2/2/19/22	0/1/1/1
4	NAG	J	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	J	2	4	-	2/6/23/26	0/1/1/1
4	BMA	J	3	4	-	2/2/19/22	0/1/1/1
4	NAG	K	1	1,4	-	2/6/23/26	0/1/1/1
4	NAG	K	2	4	-	3/6/23/26	0/1/1/1
4	BMA	K	3	4	-	0/2/19/22	0/1/1/1
5	NAG	L	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	L	2	5	-	2/6/23/26	0/1/1/1
5	BMA	L	3	5	-	2/2/19/22	0/1/1/1
5	MAN	L	4	5	-	0/2/19/22	0/1/1/1
4	NAG	M	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	M	2	4	-	1/6/23/26	0/1/1/1
4	BMA	M	3	4	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

The worst 5 of 22 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	I	6	MAN	C1-O5-C5	2.98	116.23	112.19
3	F	2	NAG	C2-N2-C7	-2.97	118.68	122.90
4	G	1	NAG	C4-C3-C2	-2.91	106.75	111.02
3	F	2	NAG	C4-C3-C2	-2.80	106.92	111.02
5	L	4	MAN	C1-O5-C5	2.69	115.83	112.19

There are no chirality outliers.

5 of 53 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C3-C2-N2-C7
4	G	3	BMA	O5-C5-C6-O6
2	E	4	MAN	O5-C5-C6-O6
5	L	3	BMA	C4-C5-C6-O6
2	E	5	NAG	O5-C5-C6-O6

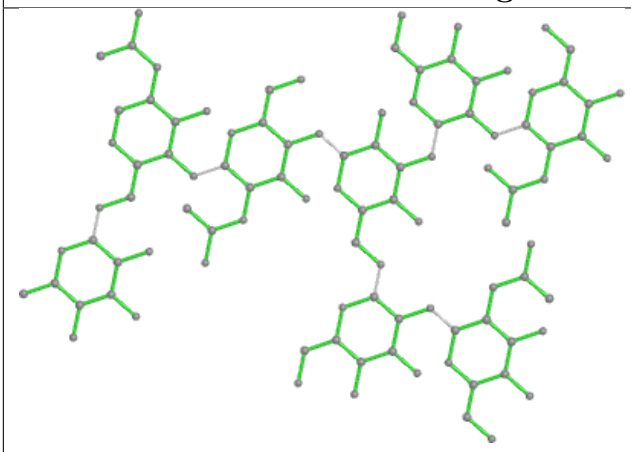
There are no ring outliers.

19 monomers are involved in 25 short contacts:

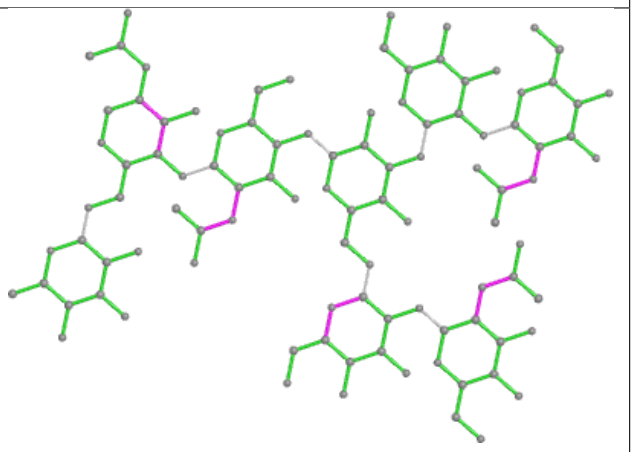
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	3	BMA	1	0
4	G	2	NAG	3	0
2	E	8	FUC	3	0
2	E	2	NAG	4	0
2	E	7	NAG	3	0
5	L	2	NAG	2	0
3	F	3	BMA	1	0
3	F	2	NAG	1	0
4	G	1	NAG	1	0
4	J	2	NAG	1	0
5	H	3	BMA	2	0
6	I	2	NAG	2	0
4	M	2	NAG	1	0
2	E	6	MAN	1	0
6	I	1	NAG	2	0
5	L	1	NAG	2	0
5	H	2	NAG	3	0
4	J	1	NAG	2	0
2	E	1	NAG	4	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.

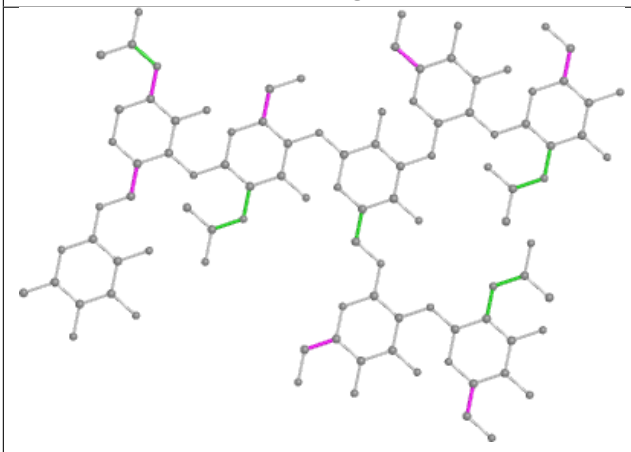
Oligosaccharide Chain E



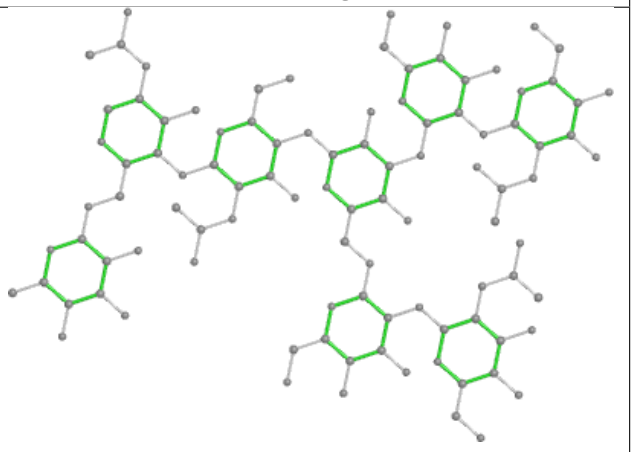
Bond lengths



Bond angles

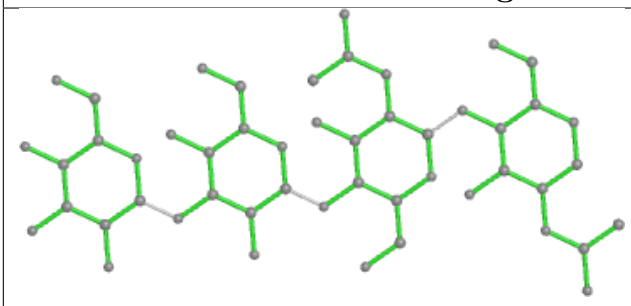


Torsions

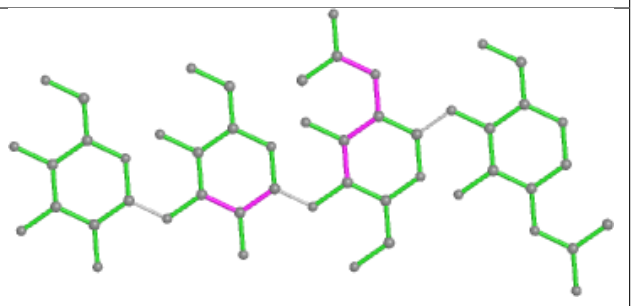


Rings

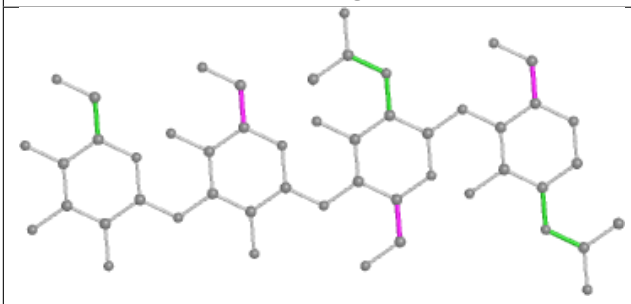
Oligosaccharide Chain F



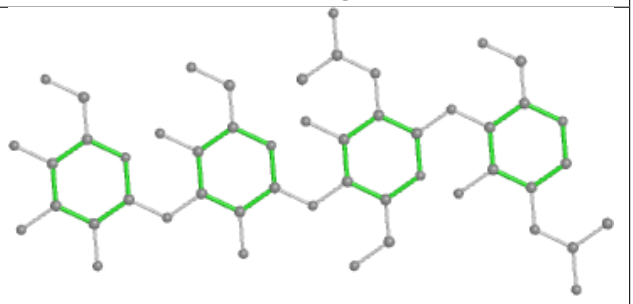
Bond lengths



Bond angles

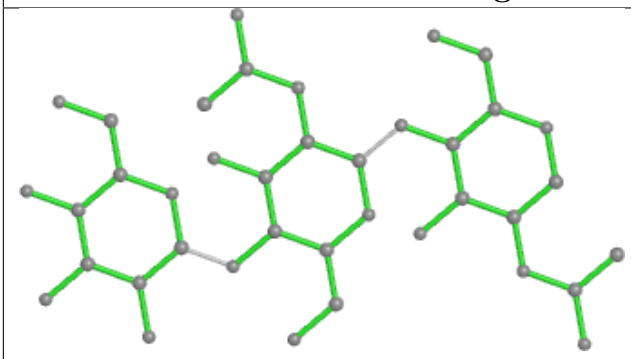


Torsions

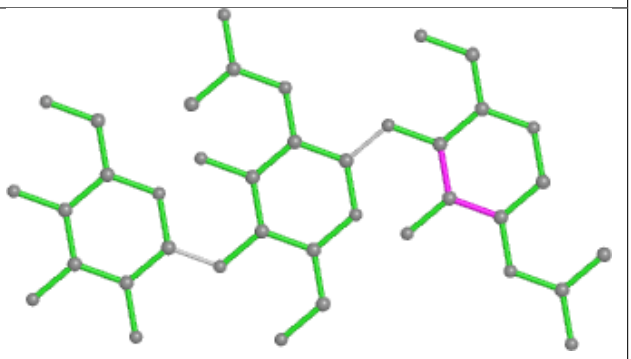


Rings

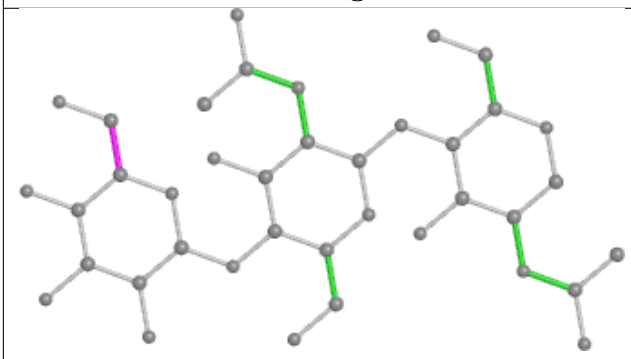
Oligosaccharide Chain G



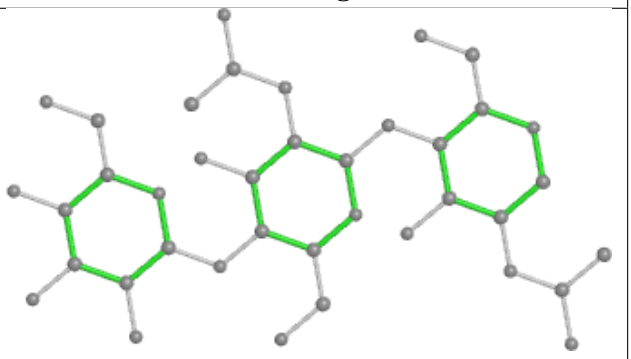
Bond lengths



Bond angles

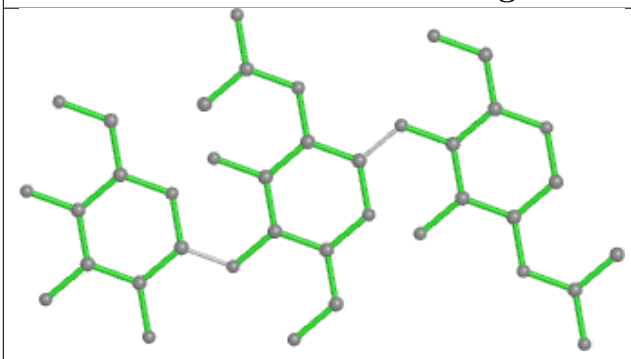


Torsions

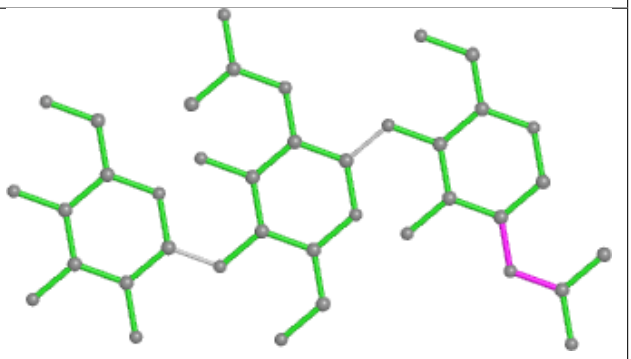


Rings

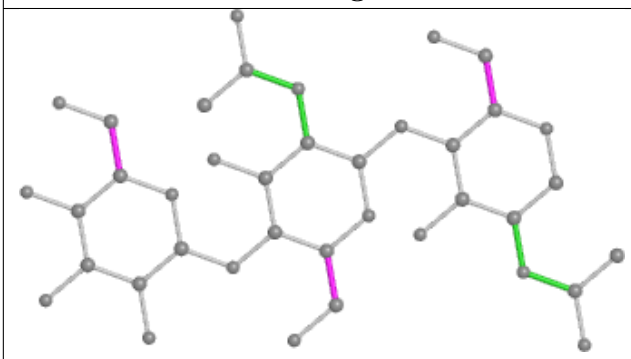
Oligosaccharide Chain J



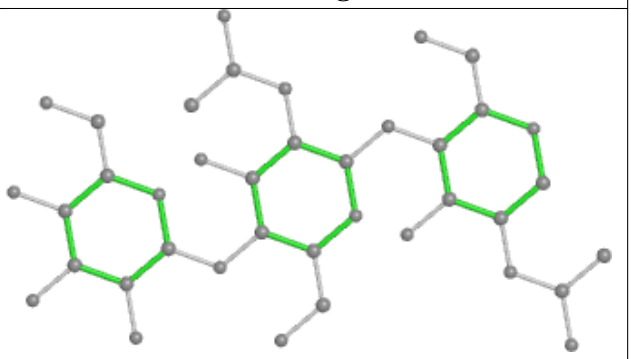
Bond lengths



Bond angles

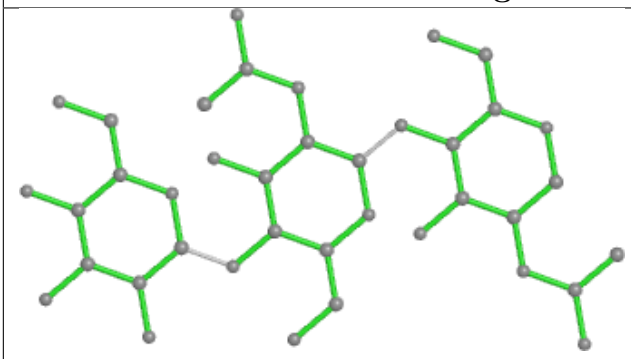


Torsions

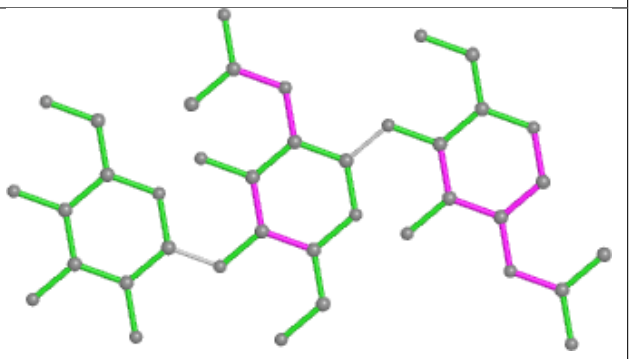


Rings

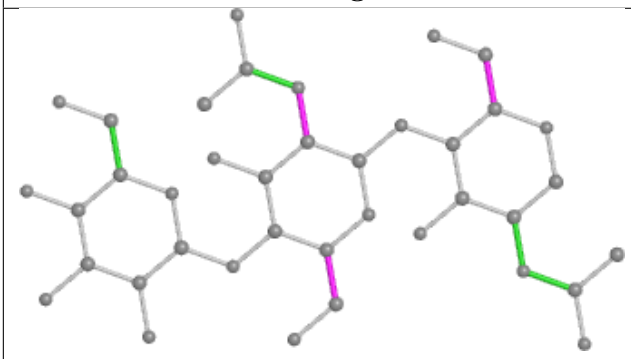
Oligosaccharide Chain K



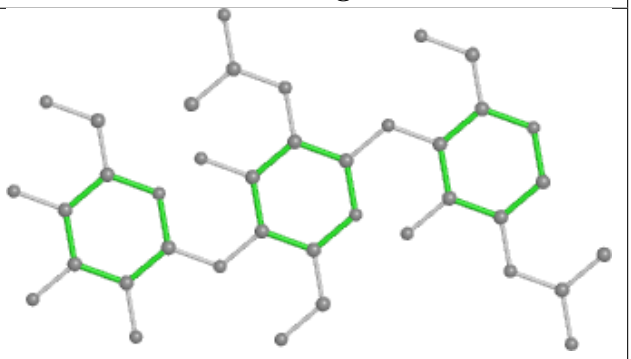
Bond lengths



Bond angles

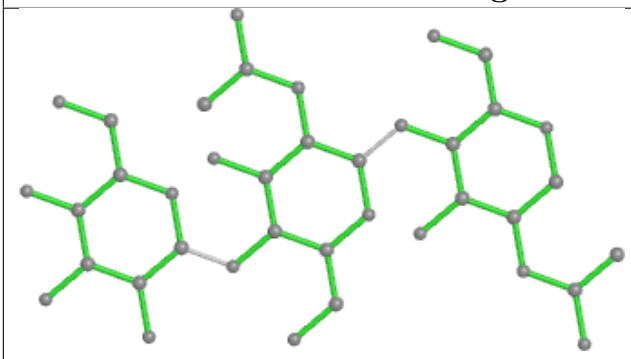


Torsions

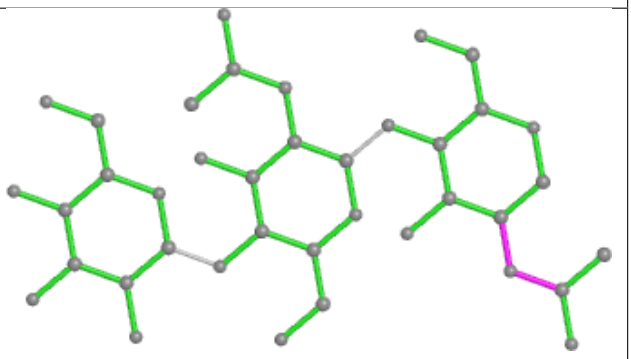


Rings

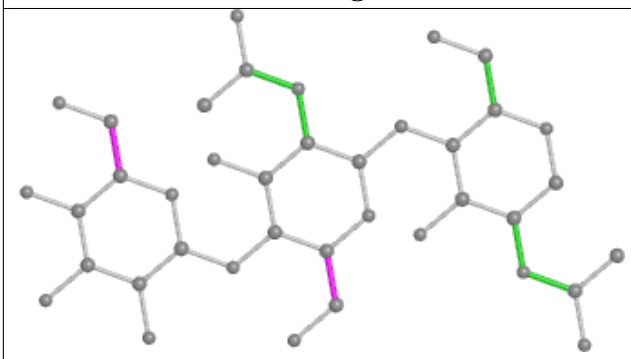
Oligosaccharide Chain M



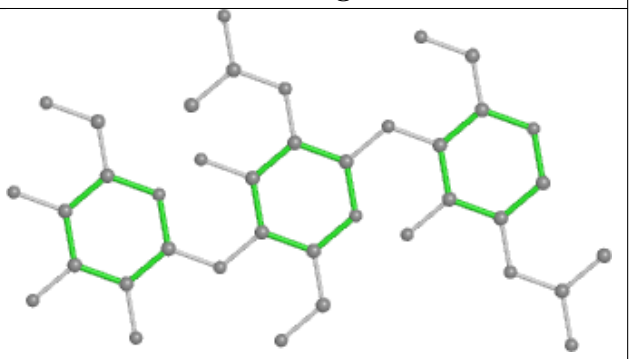
Bond lengths



Bond angles

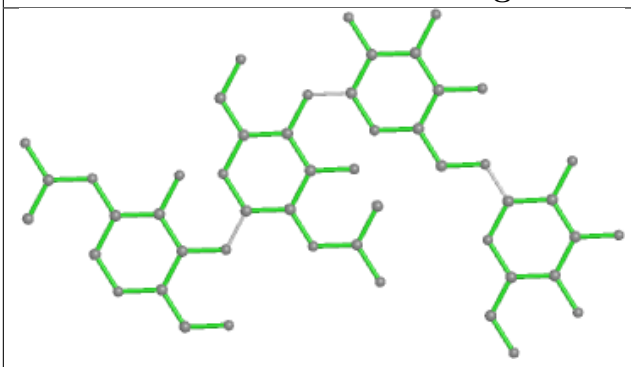


Torsions

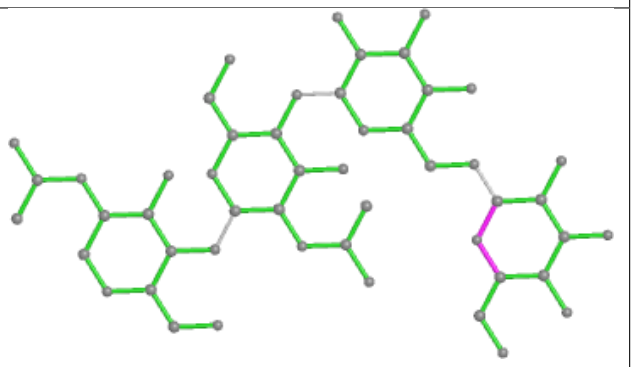


Rings

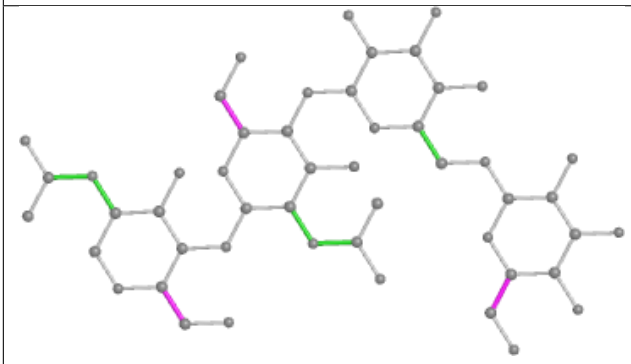
Oligosaccharide Chain H



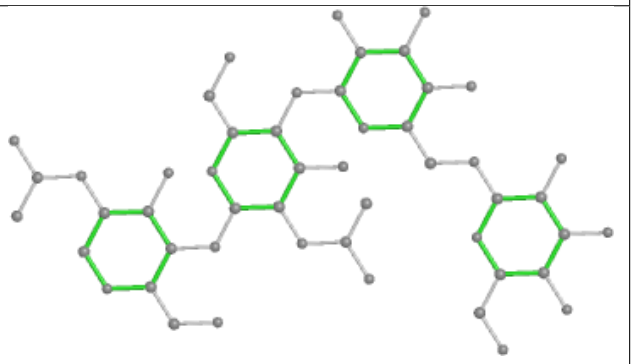
Bond lengths



Bond angles

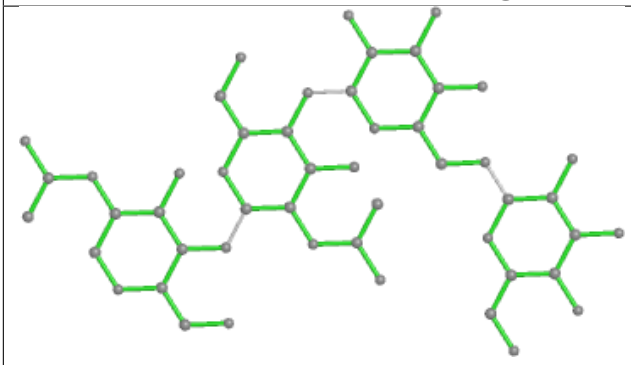


Torsions

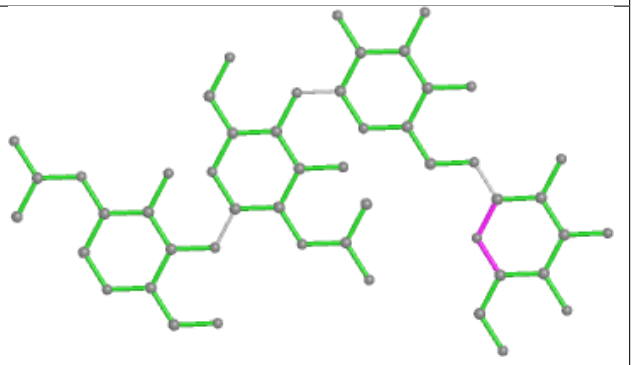


Rings

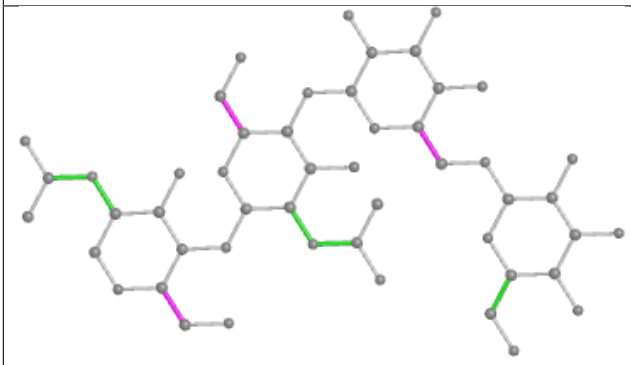
Oligosaccharide Chain L



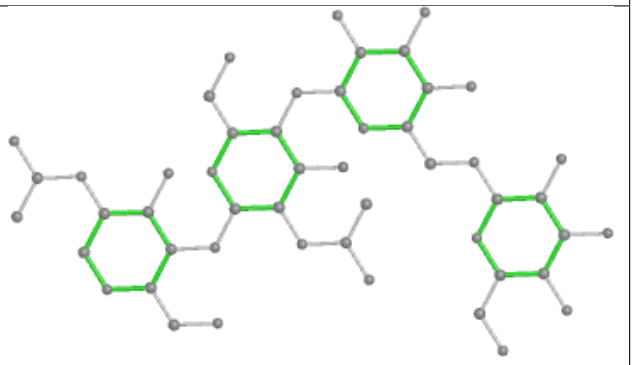
Bond lengths



Bond angles

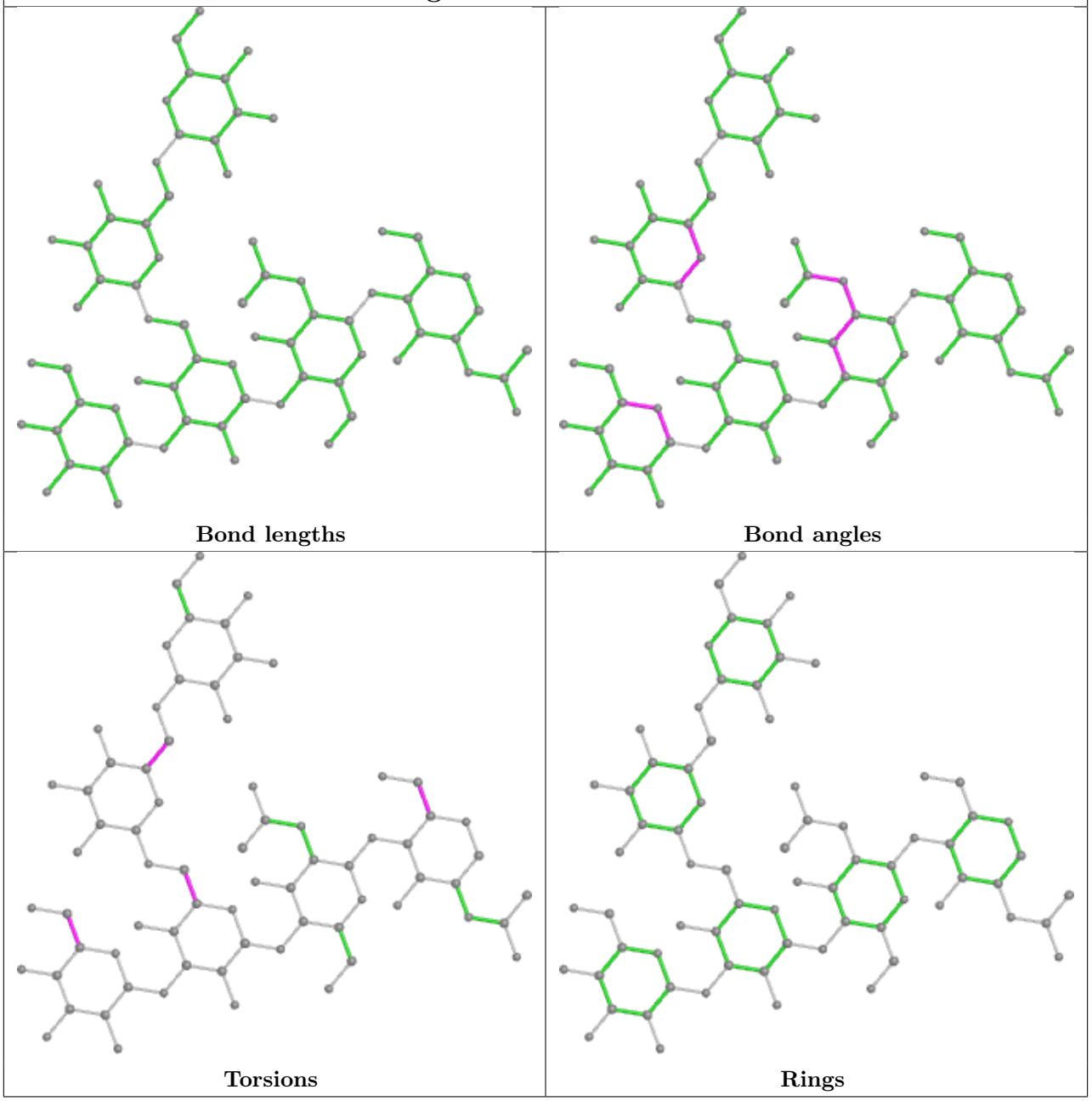


Torsions



Rings

Oligosaccharide Chain I



5.6 Ligand geometry (i)

7 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
8	GLY	D	4350	-	4,4,4	1.07	0	3,4,4	0.83	0
8	GLY	B	2350	-	4,4,4	1.06	0	3,4,4	0.90	0
7	NAG	A	1361	1	14,14,15	0.57	0	17,19,21	0.72	1 (5%)
8	GLY	C	3350	-	4,4,4	1.04	0	3,4,4	0.81	0
7	NAG	C	3361	1	14,14,15	0.58	0	17,19,21	0.80	1 (5%)
8	GLY	A	1350	-	4,4,4	1.04	0	3,4,4	1.00	0
7	NAG	D	4371	1	14,14,15	0.53	0	17,19,21	0.64	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	GLY	D	4350	-	-	1/2/2/2	-
8	GLY	B	2350	-	-	2/2/2/2	-
7	NAG	A	1361	1	-	2/6/23/26	0/1/1/1
8	GLY	C	3350	-	-	2/2/2/2	-
7	NAG	C	3361	1	-	2/6/23/26	0/1/1/1
8	GLY	A	1350	-	-	2/2/2/2	-
7	NAG	D	4371	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
7	C	3361	NAG	C2-N2-C7	-2.41	119.48	122.90
7	A	1361	NAG	C2-N2-C7	-2.16	119.83	122.90

There are no chirality outliers.

5 of 13 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	D	4371	NAG	O5-C5-C6-O6
7	D	4371	NAG	C4-C5-C6-O6
8	A	1350	GLY	OXT-C-CA-N
8	C	3350	GLY	OXT-C-CA-N
7	A	1361	NAG	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	D	4350	GLY	2	0
8	C	3350	GLY	1	0
8	A	1350	GLY	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 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.