Full wwPDB X-ray Structure Validation Report (i)

Sep 28, 2024 – 06:51 AM EDT

PDB ID	:	1WGC
Title	:	2.2 ANGSTROMS RESOLUTION STRUCTURE ANALYSIS OF TWO REFINED N-ACETYLNEURAMINYLLACTOSE-WHEAT GERM AGGLUTININ ISOLECTIN COMPLEXES
Authors	:	Wright, C.S.
Deposited on	:	1990-04-03
Resolution	:	2.20 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:

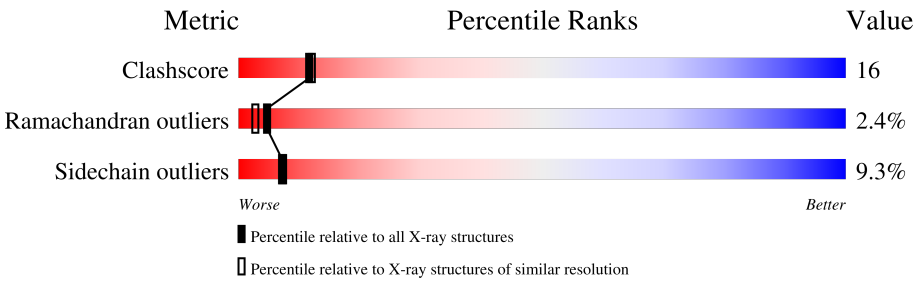
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.20 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	180529	6634 (2.20-2.20)
Ramachandran outliers	177936	6560 (2.20-2.20)
Sidechain outliers	177891	6561 (2.20-2.20)

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	171	60% 32% 7% •
1	B	171	59% 35% 5% •
2	C	3	100%
2	D	3	100%

2 Entry composition (i)

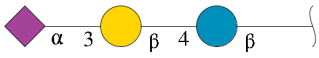
There are 3 unique types of molecules in this entry. The entry contains 2626 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 WHEAT GERM LECTIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	171	1163	678	213	238	34	0	0	0
1	B	171	1163	678	213	238	34	0	0	0

- Molecule 2 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galacto pyranose-(1-4)-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	C	3	43	23	1	19	0	0	0
2	D	3	43	23	1	19	0	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	115	115	115	0	0
3	B	99	99	99	0	0

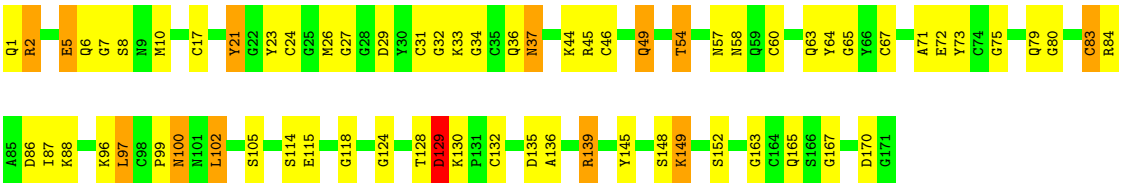
3 Residue-property plots (i)

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.

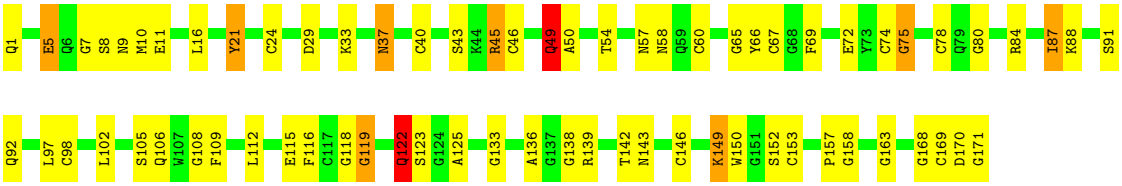
• Molecule 1: WHEAT GERM LECTIN

Chain A: 60% 32% 7%



• Molecule 1: WHEAT GERM LECTIN

Chain B: 59% 35% 5%



• Molecule 2: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain C: 100%

• Molecule 2: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-4)-beta-D-glucopyranose

Chain D: 100%

4 Data and refinement statistics (i)

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

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	51.00Å 73.64Å 91.41Å 90.00° 97.91° 90.00°	Depositor
Resolution (Å)	8.00 – 2.20	Depositor
% Data completeness (in resolution range)	(Not available) (8.00-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	PROLSQ	Depositor
R, R_{free}	0.172 , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2626	wwPDB-VP
Average B, all atoms (Å^2)	16.0	wwPDB-VP

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: SIA, PCA, BGC, GAL

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	1.16	5/1177 (0.4%)	1.87	25/1576 (1.6%)
1	B	1.18	4/1177 (0.3%)	1.88	22/1576 (1.4%)
All	All	1.17	9/2354 (0.4%)	1.87	47/3152 (1.5%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	5	GLU	CD-OE2	7.56	1.33	1.25
1	A	5	GLU	CD-OE2	7.40	1.33	1.25
1	B	11	GLU	CD-OE2	7.38	1.33	1.25
1	A	97	LEU	N-CA	-6.89	1.32	1.46
1	B	72	GLU	CD-OE2	6.58	1.32	1.25
1	A	83	CYS	CB-SG	-6.41	1.71	1.82
1	A	72	GLU	CD-OE2	6.00	1.32	1.25
1	A	24	CYS	CB-SG	-5.90	1.72	1.81
1	B	24	CYS	CB-SG	-5.37	1.73	1.81

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	84	ARG	NE-CZ-NH1	-13.19	113.70	120.30
1	B	84	ARG	NE-CZ-NH2	12.89	126.74	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	2	ARG	NE-CZ-NH1	11.02	125.81	120.30
1	B	139	ARG	CD-NE-CZ	10.55	138.37	123.60
1	B	170	ASP	CB-CG-OD2	-9.02	110.19	118.30
1	A	21	TYR	CB-CG-CD1	8.74	126.25	121.00
1	B	45	ARG	NE-CZ-NH2	-8.60	116.00	120.30
1	A	21	TYR	CB-CG-CD2	-8.41	115.95	121.00
1	A	73	TYR	CB-CG-CD2	-8.31	116.02	121.00
1	B	21	TYR	CB-CG-CD1	8.11	125.86	121.00
1	B	29	ASP	CB-CG-OD2	-7.95	111.14	118.30
1	A	2	ARG	NE-CZ-NH2	-7.92	116.34	120.30
1	A	29	ASP	CB-CG-OD2	-7.84	111.24	118.30
1	B	21	TYR	CB-CG-CD2	-7.81	116.31	121.00
1	A	86	ASP	CB-CG-OD2	-7.72	111.35	118.30
1	B	45	ARG	NE-CZ-NH1	7.68	124.14	120.30
1	A	170	ASP	CB-CG-OD1	7.50	125.05	118.30
1	A	49	GLN	CA-CB-CG	7.43	129.74	113.40
1	B	84	ARG	NH1-CZ-NH2	-7.42	111.23	119.40
1	B	66	TYR	CB-CG-CD2	-7.34	116.59	121.00
1	B	45	ARG	CD-NE-CZ	7.16	133.62	123.60
1	A	170	ASP	CB-CG-OD2	-6.99	112.01	118.30
1	B	84	ARG	CD-NE-CZ	6.65	132.91	123.60
1	A	102	LEU	CB-CG-CD2	6.64	122.30	111.00
1	B	139	ARG	NE-CZ-NH2	-6.58	117.01	120.30
1	A	129	ASP	CB-CG-OD2	-6.50	112.45	118.30
1	A	114	SER	O-C-N	6.43	132.98	122.70
1	A	26	MET	CG-SD-CE	6.42	110.47	100.20
1	B	66	TYR	CB-CG-CD1	6.31	124.79	121.00
1	A	102	LEU	CA-CB-CG	6.13	129.41	115.30
1	A	149	LYS	CA-CB-CG	6.13	126.88	113.40
1	A	5	GLU	OE1-CD-OE2	-6.06	116.03	123.30
1	A	135	ASP	CB-CG-OD2	-6.01	112.89	118.30
1	A	73	TYR	CB-CG-CD1	5.98	124.58	121.00
1	A	129	ASP	CB-CG-OD1	5.81	123.53	118.30
1	A	17	CYS	O-C-N	5.79	131.96	122.70
1	B	122	GLN	CB-CA-C	-5.58	99.23	110.40
1	A	84	ARG	NE-CZ-NH2	5.51	123.05	120.30
1	B	40	CYS	CA-CB-SG	5.46	123.82	114.00
1	B	9	ASN	CA-CB-CG	-5.45	101.42	113.40
1	B	139	ARG	NE-CZ-NH1	5.39	122.99	120.30
1	B	49	GLN	CA-CB-CG	5.29	125.04	113.40
1	B	67	CYS	CA-CB-SG	-5.20	104.64	114.00
1	B	24	CYS	CA-CB-SG	5.18	123.32	114.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	54	THR	N-CA-CB	-5.12	100.56	110.30
1	A	114	SER	N-CA-CB	5.10	118.15	110.50
1	B	49	GLN	CB-CG-CD	5.09	124.83	111.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	122	GLN	Mainchain
1	B	21	TYR	Sidechain

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	1163	0	988	32	0
1	B	1163	0	988	38	0
2	C	43	0	37	0	0
2	D	43	0	37	0	0
3	A	115	0	0	5	0
3	B	99	0	0	4	0
All	All	2626	0	2050	70	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 16.

All (70) 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:A:2:ARG:O	1:A:7:GLY:HA3	1.80	0.81
1:A:37:ASN:HA	1:A:44:LYS:NZ	2.00	0.75
1:B:88:LYS:HG2	1:B:109:PHE:HE1	1.55	0.71
1:B:7:GLY:O	1:B:10:MET:HB2	1.89	0.71
1:A:145:TYR:O	1:A:167:GLY:HA3	1.92	0.69
1:B:60:CYS:HA	1:B:80:GLY:O	1.95	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:97:LEU:HD23	1:B:123:SER:OG	1.95	0.66
1:A:57:ASN:O	1:A:58:ASN:HB2	1.94	0.66
1:A:148:SER:O	1:A:165:GLN:NE2	2.31	0.63
1:A:115:GLU:OE1	1:A:115:GLU:N	2.30	0.63
1:A:99:PRO:O	1:A:100:ASN:HB2	2.00	0.61
1:B:105:SER:O	1:B:122:GLN:NE2	2.33	0.61
1:A:149:LYS:HD3	1:A:163:GLY:HA3	1.82	0.60
1:B:74:CYS:O	1:B:78:CYS:HB2	2.01	0.60
1:A:37:ASN:HA	1:A:44:LYS:HZ2	1.66	0.60
1:B:92:GLN:NE2	1:B:106:GLN:O	2.35	0.60
1:B:102:LEU:CD2	1:B:112:LEU:HG	2.32	0.59
1:A:128:THR:O	1:A:130:LYS:N	2.35	0.59
1:B:92:GLN:OE1	1:B:92:GLN:N	2.35	0.58
1:B:157:PRO:HD2	3:B:326:HOH:O	2.06	0.56
1:B:133:GLY:HA2	3:B:325:HOH:O	2.05	0.56
1:B:98:CYS:HB3	1:B:102:LEU:HB2	1.87	0.56
1:B:88:LYS:HG2	1:B:109:PHE:CE1	2.40	0.55
1:B:136:ALA:C	1:B:138:GLY:H	2.10	0.55
1:A:33:LYS:HD2	1:A:34:GLY:N	2.21	0.55
1:A:37:ASN:HA	1:A:44:LYS:HZ1	1.69	0.54
1:B:142:THR:O	1:B:143:ASN:HB2	2.08	0.52
1:A:63:GLN:HB2	3:A:359:HOH:O	2.08	0.52
1:B:43:SER:HA	3:B:388:HOH:O	2.09	0.51
1:A:6:GLN:NE2	1:A:21:TYR:C	2.64	0.51
1:B:146:CYS:O	1:B:153:CYS:HA	2.11	0.50
1:A:33:LYS:HD2	1:A:34:GLY:H	1.75	0.50
1:B:149:LYS:HG3	1:B:150:TRP:CD1	2.48	0.49
1:B:37:ASN:ND2	3:B:319:HOH:O	2.46	0.49
1:B:169:CYS:O	1:B:171:GLY:OXT	2.30	0.49
1:B:146:CYS:SG	1:B:168:GLY:O	2.71	0.49
1:A:36:GLN:O	3:A:364:HOH:O	2.20	0.49
1:A:97:LEU:HB2	3:A:387:HOH:O	2.13	0.48
1:B:149:LYS:HA	1:B:163:GLY:O	2.14	0.48
1:B:88:LYS:HA	1:B:108:GLY:O	2.14	0.48
1:B:157:PRO:CD	1:B:158:GLY:H	2.27	0.48
1:A:149:LYS:HB3	1:A:163:GLY:HA3	1.96	0.47
1:B:75:GLY:O	1:B:78:CYS:HB2	2.15	0.47
1:B:74:CYS:O	1:B:75:GLY:O	2.32	0.46
1:B:105:SER:HB3	1:B:116:PHE:HA	1.96	0.46
1:A:139:ARG:NH1	3:A:393:HOH:O	2.49	0.45
1:A:71:ALA:O	1:A:75:GLY:HA3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:ASP:OD2	3:A:344:HOH:O	2.21	0.45
1:B:157:PRO:HD2	1:B:158:GLY:H	1.82	0.45
1:A:45:ARG:HH11	1:A:45:ARG:HG3	1.82	0.45
1:A:46:CYS:HB3	1:A:67:CYS:SG	2.57	0.45
1:A:2:ARG:HB3	1:A:6:GLN:O	2.16	0.45
1:A:27:GLY:O	1:A:31:CYS:N	2.37	0.45
1:B:57:ASN:O	1:B:58:ASN:HB2	2.16	0.45
1:A:49:GLN:HG2	1:A:64:TYR:O	2.17	0.44
1:B:87:ILE:O	1:B:109:PHE:HA	2.17	0.44
1:A:2:ARG:NE	1:A:23:TYR:CZ	2.86	0.43
1:B:58:ASN:O	1:B:69:PHE:CZ	2.72	0.43
1:B:58:ASN:O	1:B:69:PHE:HZ	2.01	0.43
1:A:49:GLN:HG2	1:A:65:GLY:HA3	2.00	0.43
1:A:83:CYS:HB2	1:A:87:ILE:HD11	2.01	0.43
1:A:132:CYS:C	1:A:136:ALA:HB3	2.39	0.43
1:B:102:LEU:HD22	1:B:112:LEU:HG	2.00	0.43
1:B:45:ARG:HA	1:B:45:ARG:HD2	1.90	0.42
1:A:102:LEU:O	1:A:124:GLY:HA3	2.20	0.42
1:B:118:GLY:O	1:B:119:GLY:C	2.58	0.41
1:B:49:GLN:HG2	1:B:65:GLY:HA3	2.03	0.41
1:B:102:LEU:HD23	1:B:112:LEU:HG	2.02	0.41
1:B:46:CYS:CA	1:B:50:ALA:HB3	2.52	0.40
1:A:60:CYS:HA	1:A:80:GLY:O	2.21	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	169/171 (99%)	149 (88%)	16 (10%)	4 (2%)	5 3
1	B	169/171 (99%)	145 (86%)	20 (12%)	4 (2%)	5 3

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	338/342 (99%)	294 (87%)	36 (11%)	8 (2%)	5 3

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	129	ASP
1	A	32	GLY
1	B	75	GLY
1	B	125	ALA
1	B	115	GLU
1	A	118	GLY
1	A	105	SER
1	B	119	GLY

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	118/119 (99%)	107 (91%)	11 (9%)	7 7
1	B	118/119 (99%)	107 (91%)	11 (9%)	7 7
All	All	236/238 (99%)	214 (91%)	22 (9%)	7 7

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	GLU
1	A	8	SER
1	A	10	MET
1	A	37	ASN
1	A	54	THR
1	A	79	GLN
1	A	88	LYS
1	A	96	LYS
1	A	100	ASN
1	A	139	ARG

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Mol	Chain	Res	Type
1	A	152	SER
1	B	5	GLU
1	B	8	SER
1	B	16	LEU
1	B	33	LYS
1	B	37	ASN
1	B	49	GLN
1	B	54	THR
1	B	87	ILE
1	B	91	SER
1	B	149	LYS
1	B	152	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	14	ASN
1	A	37	ASN
1	A	100	ASN
1	B	9	ASN
1	B	37	ASN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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	$\#\|Z\| > 2$
1	PCA	A	1	1	7,8,9	0.81	0	9,10,12	0.93	1 (11%)
1	PCA	B	1	1	7,8,9	0.91	1 (14%)	9,10,12	2.92	2 (22%)

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	PCA	A	1	1	-	0/0/11/13	0/1/1/1
1	PCA	B	1	1	-	0/0/11/13	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1	PCA	O-C	2.35	1.28	1.20

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	1	PCA	CB-CA-C	7.22	122.56	112.66
1	B	1	PCA	O-C-CA	-4.36	113.56	124.77
1	A	1	PCA	O-C-CA	-2.18	119.15	124.77

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)

6 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	C	1	2	12,12,12	1.24	2 (16%)	17,17,17	0.86	0
2	GAL	C	2	2	11,11,12	0.94	1 (9%)	15,15,17	1.19	2 (13%)
2	SIA	C	3	2	20,20,21	1.47	5 (25%)	21,28,31	2.42	9 (42%)
2	BGC	D	1	2	12,12,12	1.11	2 (16%)	17,17,17	0.82	0
2	GAL	D	2	2	11,11,12	0.77	1 (9%)	15,15,17	0.92	0
2	SIA	D	3	2	20,20,21	1.32	3 (15%)	21,28,31	2.03	5 (23%)

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	C	1	2	-	1/2/22/22	0/1/1/1
2	GAL	C	2	2	-	2/2/19/22	0/1/1/1
2	SIA	C	3	2	-	3/18/34/38	0/1/1/1
2	BGC	D	1	2	-	2/2/22/22	0/1/1/1
2	GAL	D	2	2	-	0/2/19/22	0/1/1/1
2	SIA	D	3	2	-	2/18/34/38	0/1/1/1

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	3	SIA	C4-C5	-3.15	1.50	1.53
2	C	3	SIA	O1B-C1	-3.12	1.20	1.30
2	C	3	SIA	O1A-C1	3.07	1.31	1.22
2	D	3	SIA	O1B-C1	-2.90	1.21	1.30
2	D	3	SIA	O1A-C1	2.87	1.30	1.22
2	C	1	BGC	O1-C1	2.74	1.48	1.39
2	C	1	BGC	O2-C2	2.49	1.49	1.43
2	D	1	BGC	O2-C2	2.39	1.48	1.43
2	C	3	SIA	C8-C7	-2.25	1.49	1.53
2	C	3	SIA	O6-C6	-2.20	1.40	1.44
2	C	2	GAL	O5-C1	-2.16	1.40	1.43
2	C	3	SIA	C4-C5	-2.12	1.51	1.53
2	D	2	GAL	O5-C1	-2.06	1.40	1.43
2	D	1	BGC	O1-C1	2.05	1.46	1.39

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	D	3	SIA	O1A-C1-C2	-5.13	111.76	122.85
2	D	3	SIA	O1B-C1-C2	4.95	125.58	112.71
2	C	3	SIA	O1B-C1-C2	4.50	124.41	112.71
2	C	3	SIA	O6-C2-C1	4.34	115.91	107.72
2	C	3	SIA	O1A-C1-C2	-4.33	113.50	122.85
2	C	3	SIA	C9-C8-C7	3.95	120.23	112.17
2	C	3	SIA	C6-C5-N5	-3.23	105.76	110.91
2	C	3	SIA	C11-C10-N5	-2.98	111.17	116.12
2	D	3	SIA	C8-C7-C6	-2.75	107.90	113.05
2	C	2	GAL	C1-O5-C5	2.62	115.70	112.19
2	D	3	SIA	O6-C2-C1	2.43	112.30	107.72
2	C	3	SIA	O7-C7-C8	2.42	114.42	108.93
2	C	3	SIA	O10-C10-C11	2.39	126.31	122.05
2	D	3	SIA	C9-C8-C7	2.28	116.81	112.17
2	C	3	SIA	O9-C9-C8	-2.19	106.56	111.16
2	C	2	GAL	C1-C2-C3	-2.05	106.66	109.64

There are no chirality outliers.

All (10) torsion outliers are listed below:

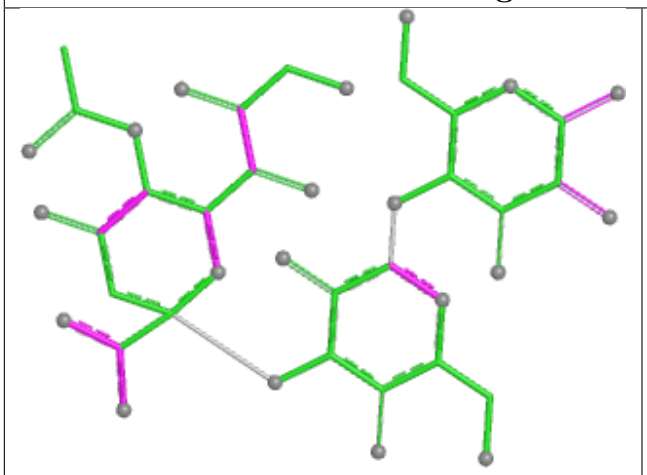
Mol	Chain	Res	Type	Atoms
2	D	3	SIA	O8-C8-C9-O9
2	D	3	SIA	C7-C8-C9-O9
2	C	2	GAL	O5-C5-C6-O6
2	C	2	GAL	C4-C5-C6-O6
2	D	1	BGC	O5-C5-C6-O6
2	D	1	BGC	C4-C5-C6-O6
2	C	3	SIA	C6-C7-C8-O8
2	C	1	BGC	O5-C5-C6-O6
2	C	3	SIA	C5-C6-C7-C8
2	C	3	SIA	C7-C8-C9-O9

There are no ring outliers.

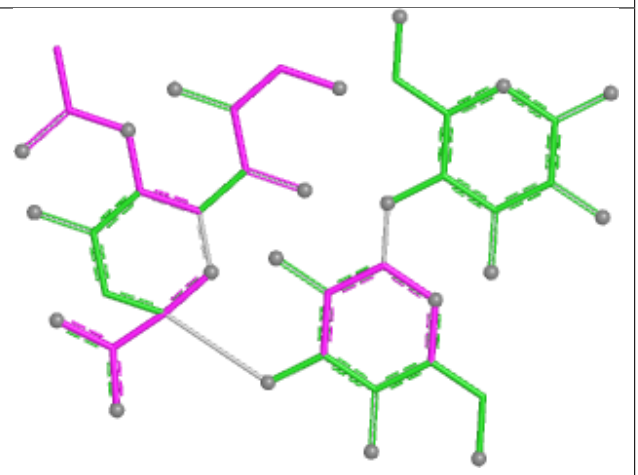
No monomer is involved in short contacts.

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

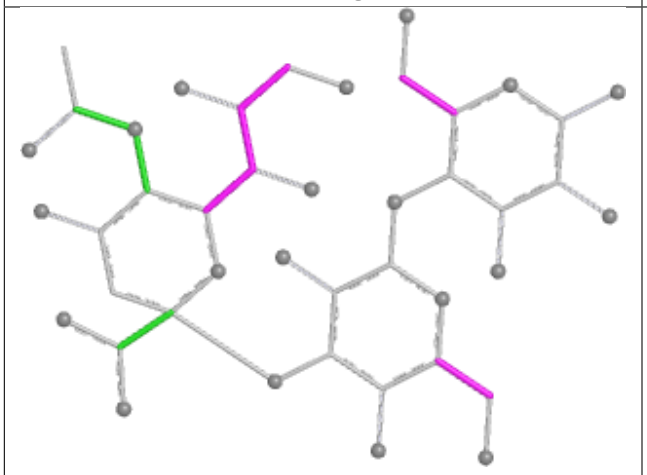
Oligosaccharide Chain C



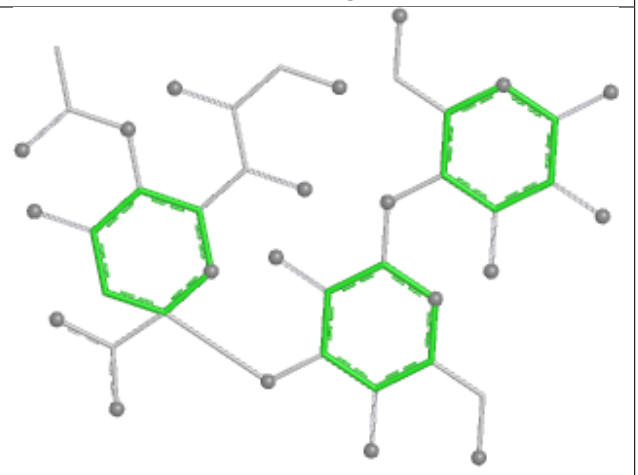
Bond lengths



Bond angles

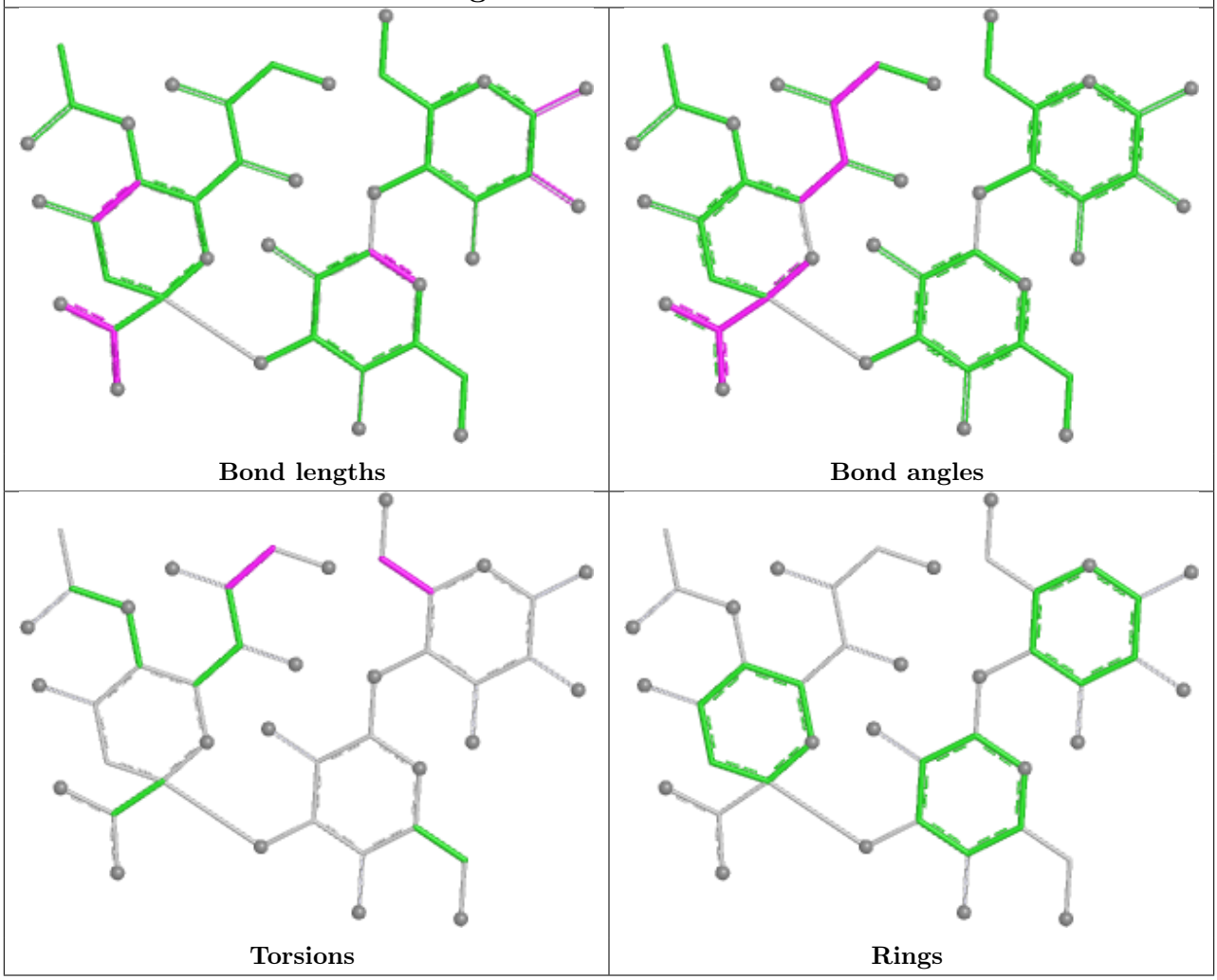


Torsions



Rings

Oligosaccharide Chain D



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)

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.