



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 07:26 PM UTC

PDB ID : 4UM9 / pdb_00004um9
Title : Crystal structure of alpha V beta 6 with peptide
Authors : Dong, X.; Springer, T.A.
Deposited on : 2014-05-15
Resolution : 2.50 Å(reported)

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

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A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

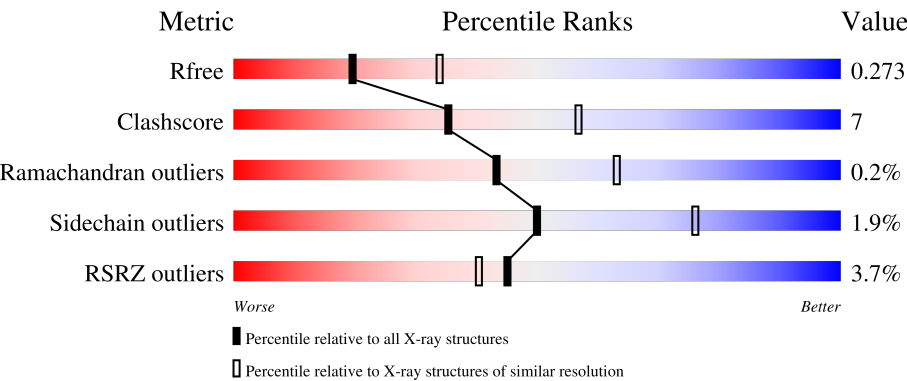
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

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 2.50 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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 $\leq 5\%$. 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	604	2% 79% 19% .
1	C	604	2% 79% 17% ..
2	B	483	5% 72% 19% . 7%
3	D	483	6% 72% 21% . 6%
4	E	13	69% 15% 15%

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Mol	Chain	Length	Quality of chain
4	F	13	 8% 85% 15%
5	G	2	 100%
5	J	2	 100%
5	K	2	 50% 50%
5	L	2	 100%
5	M	2	 100%
5	P	2	 100%
6	H	4	 75% 25%
7	I	6	 33% 67%
7	N	6	 33% 33% 33%
8	O	3	 100%

2 Entry composition

There are 14 unique types of molecules in this entry. The entry contains 17049 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 Integrin alpha-V heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	594	Total	C	N	O	S	0	0	0
			4601	2917	781	882	21			
1	C	589	Total	C	N	O	S	0	0	0
			4566	2897	774	874	21			

There are 20 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	400	CYS	MET	conflict	UNP P06756
A	596	THR	-	expression tag	UNP P06756
A	597	GLY	-	expression tag	UNP P06756
A	598	GLY	-	expression tag	UNP P06756
A	599	LEU	-	expression tag	UNP P06756
A	600	GLU	-	expression tag	UNP P06756
A	601	VAL	-	expression tag	UNP P06756
A	602	LEU	-	expression tag	UNP P06756
A	603	PHE	-	expression tag	UNP P06756
A	604	GLN	-	expression tag	UNP P06756
C	400	CYS	MET	conflict	UNP P06756
C	596	THR	-	expression tag	UNP P06756
C	597	GLY	-	expression tag	UNP P06756
C	598	GLY	-	expression tag	UNP P06756
C	599	LEU	-	expression tag	UNP P06756
C	600	GLU	-	expression tag	UNP P06756
C	601	VAL	-	expression tag	UNP P06756
C	602	LEU	-	expression tag	UNP P06756
C	603	PHE	-	expression tag	UNP P06756
C	604	GLN	-	expression tag	UNP P06756

- Molecule 2 is a protein called Integrin beta-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	448	Total	C	N	O	S	0	0	0
			3433	2155	586	662	30			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	270	CYS	ILE	conflict	UNP P18564
B	452	ASN	HIS	conflict	UNP P18564
B	475	SER	-	expression tag	UNP P18564
B	476	ARG	-	expression tag	UNP P18564
B	477	GLY	-	expression tag	UNP P18564
B	478	LEU	-	expression tag	UNP P18564
B	479	GLN	-	expression tag	UNP P18564
B	480	THR	-	expression tag	UNP P18564
B	481	LEU	-	expression tag	UNP P18564
B	482	PHE	-	expression tag	UNP P18564
B	483	GLN	-	expression tag	UNP P18564

- Molecule 3 is a protein called Integrin beta-6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	455	Total	C	N	O	S	0	0	0
			3485	2185	593	677	30			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
D	29	THR	GLN	conflict	UNP P18564
D	270	CYS	ILE	conflict	UNP P18564
D	452	ASN	HIS	conflict	UNP P18564
D	475	SER	-	expression tag	UNP P18564
D	476	ARG	-	expression tag	UNP P18564
D	477	GLY	-	expression tag	UNP P18564
D	478	LEU	-	expression tag	UNP P18564
D	479	GLN	-	expression tag	UNP P18564
D	480	THR	-	expression tag	UNP P18564
D	481	LEU	-	expression tag	UNP P18564
D	482	PHE	-	expression tag	UNP P18564
D	483	GLN	-	expression tag	UNP P18564

- Molecule 4 is a protein called Latency-associated peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	11	Total	C	N	O	0	0	1
			77	46	19	12			
4	F	13	Total	C	N	O	0	0	2
			89	53	22	14			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	238	ACE	-	expression tag	UNP P10600
E	250	NH2	-	expression tag	UNP P10600
F	238	ACE	-	expression tag	UNP P10600
F	250	NH2	-	expression tag	UNP P10600

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



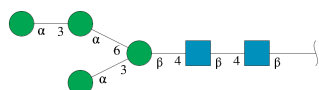
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	J	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	M	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	P	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 6 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.



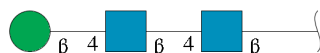
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	H	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-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.



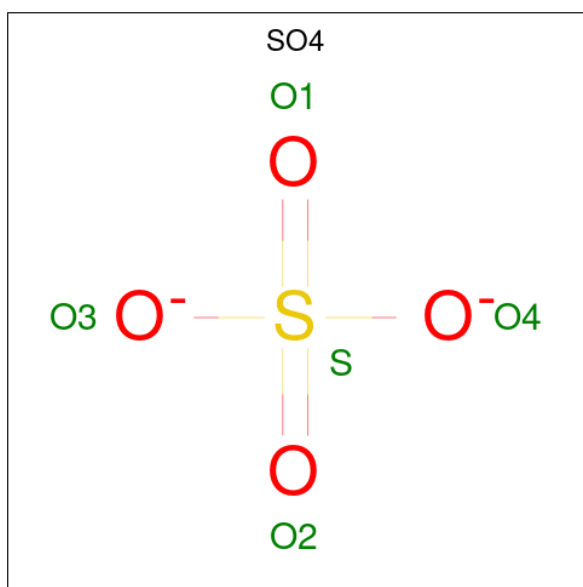
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
7	I	6	Total	C	N	O	0	0	0
			72	40	2	30			
7	N	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 8 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
8	O	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 9 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	O	S	0	0
			5	4	1		
9	A	1	Total	O	S	0	0
			5	4	1		
9	A	1	Total	O	S	0	0
			5	4	1		
9	B	1	Total	O	S	0	0
			5	4	1		
9	C	1	Total	O	S	0	0
			5	4	1		
9	C	1	Total	O	S	0	0
			5	4	1		

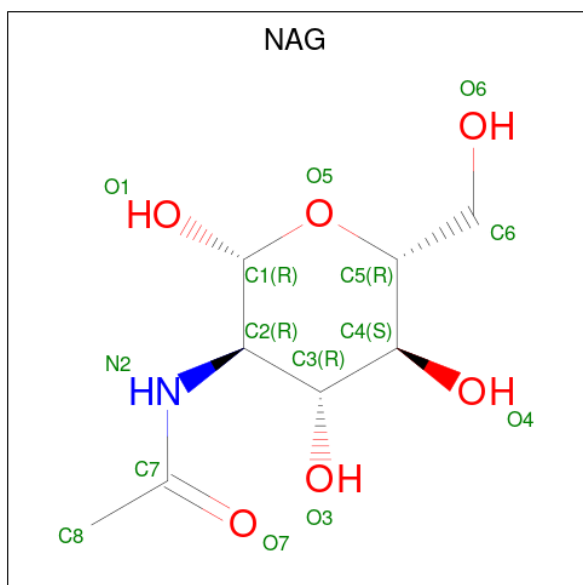
- Molecule 10 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	4	Total	Ca	0	0
			4	4		
10	B	2	Total	Ca	0	0
			2	2		
10	C	4	Total	Ca	0	0
			4	4		
10	D	2	Total	Ca	0	0
			2	2		

- Molecule 11 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

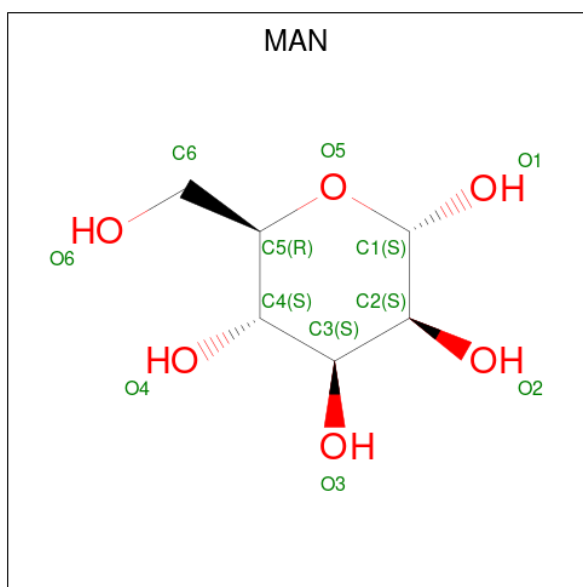
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
11	B	1	Total	Mg	0	0
			1	1		
11	D	1	Total	Mg	0	0
			1	1		

- Molecule 12 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 13 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
13	C	1	Total	C	O	0	0
			11	6	5		

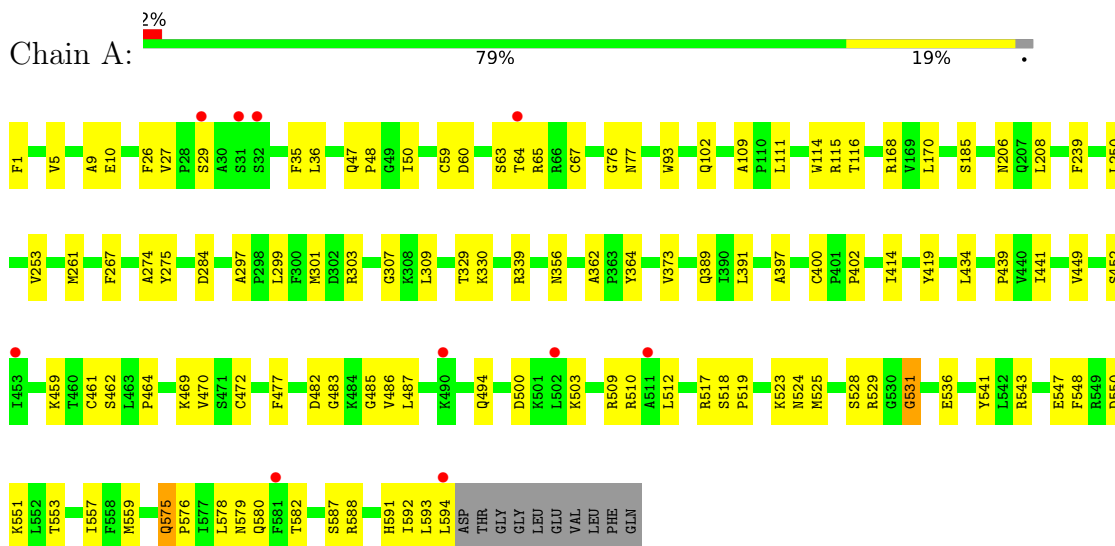
- Molecule 14 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
14	A	85	Total	O	0	0
			85	85		
14	B	65	Total	O	0	0
			65	65		
14	C	52	Total	O	0	0
			52	52		
14	D	25	Total	O	0	0
			25	25		
14	E	3	Total	O	0	0
			3	3		

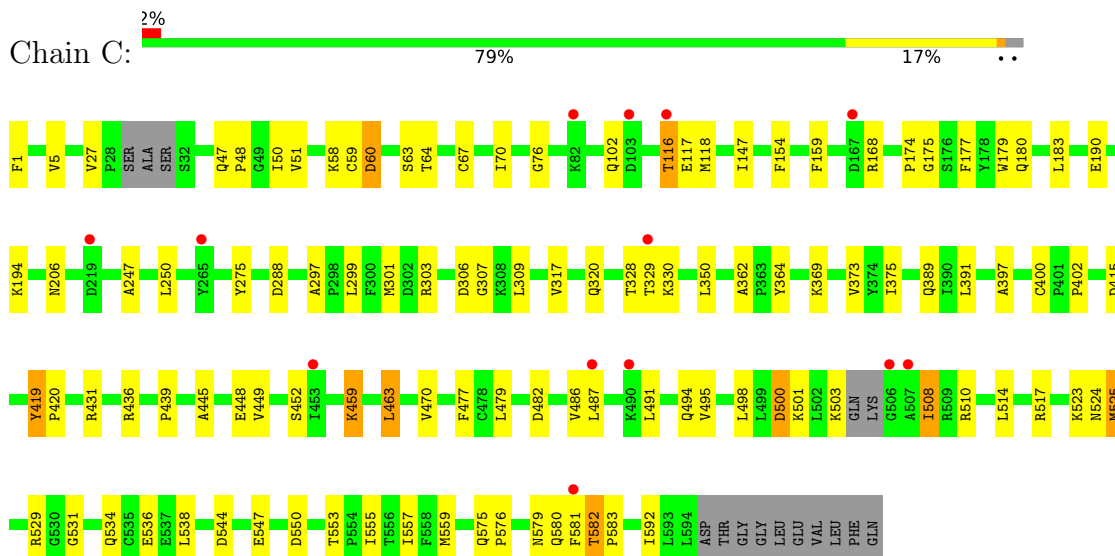
3 Residue-property plots

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: Integrin alpha-V heavy chain

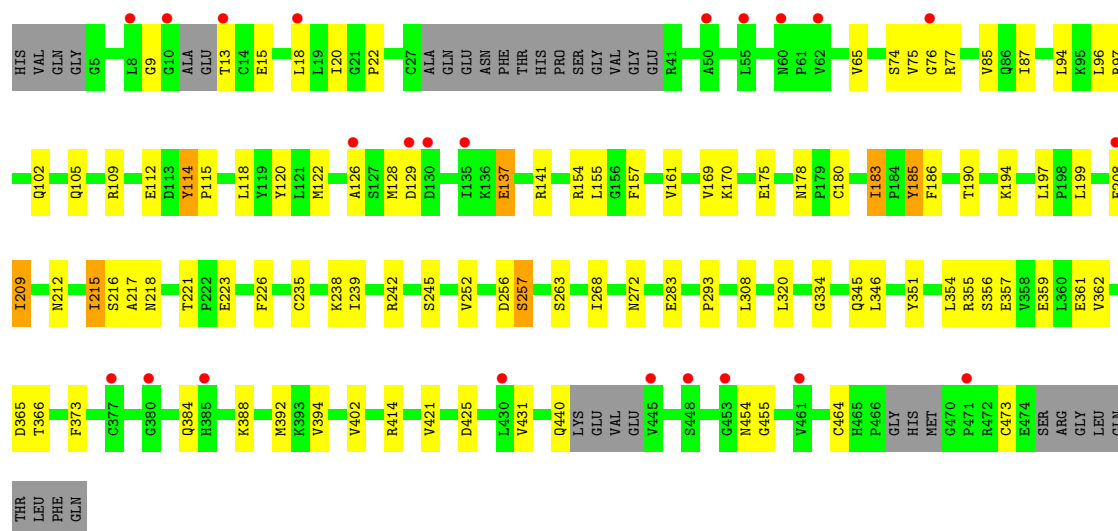


- Molecule 1: Integrin alpha-V heavy chain

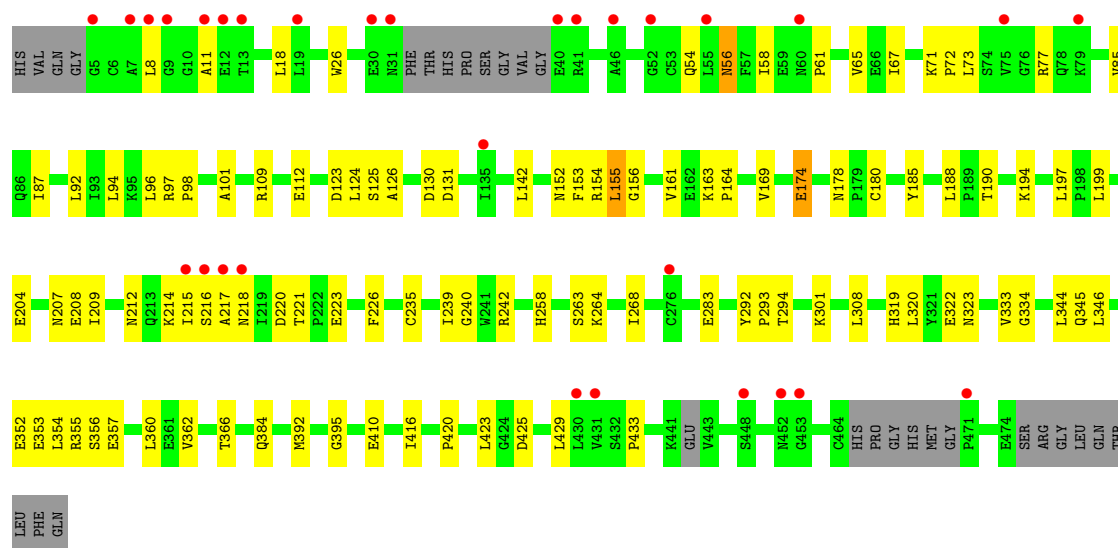
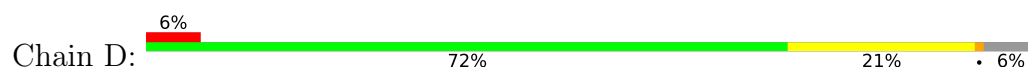


- Molecule 2: Integrin beta-6

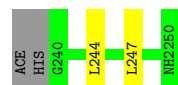




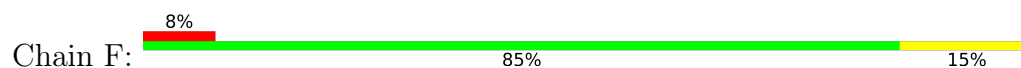
• Molecule 3: Integrin beta-6

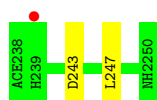


• Molecule 4: Latency-associated peptide



• Molecule 4: Latency-associated peptide





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

Chain G: 100%



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

Chain J: 100%



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

Chain K: 50% 50%



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

Chain L: 100%



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

Chain M: 100%



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

Chain P: 100%



- Molecule 6: 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:  75% 25%

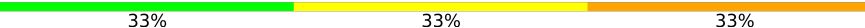


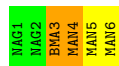
- Molecule 7: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  33% 67%



- Molecule 7: alpha-D-mannopyranose-(1-3)-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 N:  33% 33% 33%



- Molecule 8: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	185.02Å 168.09Å 101.80Å 90.00° 98.95° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 50.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	94.4 (50.00-2.50) 94.4 (50.00-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.35 (at 2.51Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.222 , 0.266 0.227 , 0.273	Depositor DCC
R_{free} test set	1416 reflections (1.42%)	wwPDB-VP
Wilson B-factor (Å ²)	54.5	Xtriage
Anisotropy	0.124	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 76.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	17049	wwPDB-VP
Average B, all atoms (Å ²)	74.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NH2, BMA, MAN, NAG, CA, ACE, MG, SO4

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	0.41	0/4705	1.01	20/6370 (0.3%)
1	C	0.41	0/4668	1.06	27/6318 (0.4%)
2	B	0.40	0/3493	0.93	15/4728 (0.3%)
3	D	0.38	0/3544	0.92	10/4797 (0.2%)
4	E	0.24	0/75	0.67	0/95
4	F	0.25	0/87	0.73	0/112
All	All	0.40	0/16572	0.99	72/22420 (0.3%)

There are no bond length outliers.

All (72) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	582	THR	CA-C-N	12.39	133.21	119.83
1	A	582	THR	C-N-CA	12.39	133.21	119.83
1	C	116	THR	N-CA-C	-11.94	95.13	110.53
1	C	400	CYS	CA-C-N	-9.30	110.80	120.38
1	C	400	CYS	C-N-CA	-9.30	110.80	120.38
1	A	400	CYS	CA-C-N	-8.72	111.39	120.38
1	A	400	CYS	C-N-CA	-8.72	111.39	120.38
1	C	582	THR	CA-C-N	8.63	129.99	119.98
1	C	582	THR	C-N-CA	8.63	129.99	119.98
3	D	197	LEU	CA-C-N	8.24	128.00	119.76
3	D	197	LEU	C-N-CA	8.24	128.00	119.76
2	B	114	TYR	CA-C-N	8.23	128.24	119.76
2	B	114	TYR	C-N-CA	8.23	128.24	119.76
1	A	400	CYS	N-CA-C	-7.84	98.79	110.24
2	B	137	GLU	CB-CA-C	-7.81	98.99	111.18
1	A	531	GLY	N-CA-C	7.43	124.79	112.30
1	C	463	LEU	CA-C-N	7.42	128.59	119.98
1	C	463	LEU	C-N-CA	7.42	128.59	119.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	508	ILE	N-CA-C	7.21	118.41	108.89
3	D	221	THR	CA-C-N	-7.20	112.22	119.92
3	D	221	THR	C-N-CA	-7.20	112.22	119.92
1	C	400	CYS	N-CA-C	-7.03	99.98	110.24
1	A	297	ALA	CA-C-N	6.67	126.36	119.82
1	A	297	ALA	C-N-CA	6.67	126.36	119.82
1	C	117	GLU	N-CA-C	-6.58	104.88	113.17
2	B	215	ILE	N-CA-C	6.44	117.89	109.58
2	B	185	TYR	N-CA-C	-6.34	99.44	109.07
1	C	174	PRO	CB-CA-C	6.34	120.93	111.68
1	C	486	VAL	N-CA-C	6.28	116.87	107.51
2	B	180	CYS	N-CA-C	-6.25	103.06	111.87
2	B	178	ASN	CA-C-N	-6.16	112.14	119.84
2	B	178	ASN	C-N-CA	-6.16	112.14	119.84
2	B	238	LYS	CG-CD-CE	6.15	125.44	111.30
1	C	500	ASP	CA-C-N	6.14	128.51	120.28
1	C	500	ASP	C-N-CA	6.14	128.51	120.28
3	D	112	GLU	CA-C-N	-6.05	114.46	122.99
3	D	112	GLU	C-N-CA	-6.05	114.46	122.99
1	A	486	VAL	N-CA-C	6.02	116.47	107.51
1	A	102	GLN	CB-CA-C	-5.99	109.09	117.23
3	D	180	CYS	N-CA-C	-5.96	103.47	111.87
1	C	362	ALA	CA-C-N	5.95	125.63	119.56
1	C	362	ALA	C-N-CA	5.95	125.63	119.56
1	A	592	ILE	N-CA-C	-5.95	99.85	108.89
1	C	531	GLY	N-CA-C	5.95	122.29	112.30
1	A	575	GLN	CA-C-N	5.86	127.17	119.84
1	A	575	GLN	C-N-CA	5.86	127.17	119.84
1	C	320	GLN	N-CA-C	5.79	117.90	108.63
1	C	592	ILE	N-CA-C	-5.78	100.10	108.89
2	B	221	THR	CA-C-N	-5.76	113.76	119.92
2	B	221	THR	C-N-CA	-5.76	113.76	119.92
2	B	183	ILE	CA-C-N	5.75	140.79	127.00
2	B	183	ILE	C-N-CA	5.75	140.79	127.00
1	A	485	GLY	N-CA-C	5.61	122.46	112.02
1	C	491	LEU	N-CA-C	5.58	117.82	108.73
2	B	183	ILE	C-N-CD	-5.52	108.45	120.60
1	C	495	VAL	N-CA-C	5.38	115.70	108.17
3	D	178	ASN	CA-C-N	-5.33	113.18	119.84
3	D	178	ASN	C-N-CA	-5.33	113.18	119.84
1	C	297	ALA	CA-C-N	5.21	124.92	119.82
1	C	297	ALA	C-N-CA	5.21	124.92	119.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	362	ALA	CA-C-N	5.20	124.86	119.56
1	A	362	ALA	C-N-CA	5.20	124.86	119.56
1	A	500	ASP	CA-C-N	5.16	127.72	120.28
1	A	500	ASP	C-N-CA	5.16	127.72	120.28
1	C	419	TYR	CA-C-N	5.16	124.92	119.76
1	C	419	TYR	C-N-CA	5.16	124.92	119.76
3	D	185	TYR	N-CA-C	-5.15	101.25	109.07
1	A	109	ALA	CA-C-N	5.08	124.87	119.28
1	A	109	ALA	C-N-CA	5.08	124.87	119.28
1	C	102	GLN	CB-CA-C	-5.07	110.33	117.23
1	C	60	ASP	N-CA-C	-5.03	101.50	109.76
2	B	128	MET	N-CA-C	5.00	118.43	112.72

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	4601	0	4452	60	1
1	C	4566	0	4414	58	1
2	B	3433	0	3381	64	0
3	D	3485	0	3431	61	0
4	E	77	0	86	2	0
4	F	89	0	94	2	0
5	G	28	0	25	0	0
5	J	28	0	25	0	0
5	K	28	0	25	0	0
5	L	28	0	25	0	0
5	M	28	0	25	1	0
5	P	28	0	25	1	0
6	H	50	0	43	0	0
7	I	72	0	61	0	0
7	N	72	0	61	1	0
8	O	39	0	34	2	0
9	A	15	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
9	B	5	0	0	0	0
9	C	10	0	0	0	0
10	A	4	0	0	0	0
10	B	2	0	0	0	0
10	C	4	0	0	0	0
10	D	2	0	0	0	0
11	B	1	0	0	0	0
11	D	1	0	0	0	0
12	B	42	0	39	0	0
12	C	28	0	26	0	0
12	D	42	0	39	0	0
13	C	11	0	10	1	0
14	A	85	0	0	0	0
14	B	65	0	0	2	0
14	C	52	0	0	3	0
14	D	25	0	0	4	0
14	E	3	0	0	0	0
All	All	17049	0	16321	245	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 7.

All (245) 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:509:ARG:O	1:A:543:ARG:NH1	2.03	0.91
1:A:494:GLN:NE2	1:A:524:ASN:OD1	2.11	0.84
1:A:10:GLU:H	1:A:65:ARG:HH11	1.35	0.75
3:D:154:ARG:NH2	3:D:240:GLY:O	2.20	0.75
2:B:208:GLU:O	2:B:212:ASN:ND2	2.19	0.74
3:D:152:ASN:N	14:D:2102:HOH:O	2.22	0.72
1:C:494:GLN:NE2	1:C:524:ASN:OD1	2.23	0.72
2:B:170:LYS:HB2	2:B:175:GLU:HG2	1.70	0.71
1:A:487:LEU:HD11	1:A:529:ARG:HD3	1.73	0.70
3:D:71:LYS:HD2	3:D:72:PRO:HD2	1.73	0.69
1:C:397:ALA:HB2	1:C:402:PRO:HD3	1.73	0.69
1:A:10:GLU:H	1:A:65:ARG:NH1	1.90	0.69
1:A:303:ARG:HD2	1:A:307:GLY:O	1.93	0.69
1:C:190:GLU:OE2	1:C:206:ASN:ND2	2.25	0.68
1:C:415:ASP:OD2	14:C:801:HOH:O	2.12	0.67
3:D:420:PRO:HG2	3:D:423:LEU:HB2	1.76	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:161:VAL:HG21	2:B:216:SER:HB2	1.78	0.66
2:B:186:PHE:HB2	3:D:101:ALA:HB2	1.76	0.65
2:B:15:GLU:OE1	2:B:440:GLN:NE2	2.30	0.65
3:D:125:SER:HB2	4:F:243:ASP:OD1	1.97	0.64
1:A:168:ARG:NH2	1:A:206:ASN:O	2.31	0.63
1:C:58:LYS:HB2	1:C:70:ILE:HD11	1.80	0.63
3:D:263:SER:HB2	3:D:268:ILE:HB	1.79	0.63
2:B:351:TYR:OH	2:B:355:ARG:NH1	2.31	0.62
2:B:129:ASP:OD2	2:B:185:TYR:OH	2.11	0.62
1:C:168:ARG:NH2	1:C:206:ASN:O	2.33	0.62
2:B:18:LEU:HD22	2:B:97:ARG:HH22	1.65	0.62
1:C:503:LYS:NZ	1:C:550:ASP:OD2	2.32	0.61
1:A:63:SER:OG	1:A:64:THR:N	2.32	0.61
1:C:303:ARG:NH1	1:C:307:GLY:O	2.34	0.61
3:D:161:VAL:HG21	3:D:216:SER:HB2	1.82	0.60
1:C:59:CYS:HA	1:C:67:CYS:HA	1.83	0.60
1:C:364:TYR:HB3	1:C:369:LYS:HE2	1.83	0.59
1:C:116:THR:HG22	1:C:118:MET:H	1.67	0.59
3:D:194:LYS:HG2	3:D:283:GLU:HG2	1.83	0.59
1:C:1:PHE:CE1	1:C:580:GLN:HG2	2.38	0.58
1:C:1:PHE:HA	1:C:389:GLN:HB2	1.85	0.58
2:B:455:GLY:HA3	2:B:464:CYS:HA	1.86	0.58
1:C:482:ASP:HA	1:C:529:ARG:HG3	1.87	0.57
1:A:441:ILE:HB	1:A:578:LEU:HD23	1.87	0.57
1:A:548:PHE:O	1:A:551:LYS:NZ	2.28	0.56
3:D:226:PHE:HB3	3:D:293:PRO:HG2	1.86	0.56
2:B:65:VAL:HG12	2:B:94:LEU:HD13	1.87	0.56
1:C:436:ARG:HB3	14:C:801:HOH:O	2.05	0.56
1:C:445:ALA:HB1	1:C:559:MET:HE1	1.88	0.56
1:C:47:GLN:HB3	1:C:50:ILE:HD12	1.88	0.56
2:B:235:CYS:HB2	2:B:239:ILE:HD13	1.88	0.55
2:B:190:THR:OG1	2:B:217:ALA:O	2.24	0.55
2:B:256:ASP:OD1	2:B:257:SER:N	2.36	0.55
3:D:156:GLY:HA3	3:D:199:LEU:HD12	1.89	0.55
2:B:226:PHE:HB3	2:B:293:PRO:HG2	1.89	0.54
3:D:319:HIS:O	3:D:323:ASN:ND2	2.41	0.54
2:B:96:LEU:HD21	2:B:102:GLN:HB2	1.90	0.54
1:A:47:GLN:HB3	1:A:50:ILE:HD12	1.90	0.53
3:D:362:VAL:HG12	3:D:366:THR:HG21	1.90	0.53
3:D:235:CYS:HB2	3:D:239:ILE:HD13	1.91	0.53
2:B:154:ARG:HH12	2:B:242:ARG:CZ	2.22	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:359:GLU:HB3	2:B:421:VAL:CG2	2.38	0.53
1:A:208:LEU:HB3	1:A:261:MET:HE2	1.91	0.53
3:D:56:ASN:N	3:D:56:ASN:OD1	2.42	0.53
3:D:190:THR:OG1	3:D:217:ALA:O	2.27	0.52
8:O:1:NAG:H61	8:O:2:NAG:N2	2.25	0.52
3:D:97:ARG:HD2	3:D:98:PRO:HD2	1.92	0.52
2:B:197:LEU:HB2	2:B:209:ILE:HD13	1.92	0.52
3:D:18:LEU:HA	3:D:97:ARG:HH21	1.75	0.52
1:C:63:SER:OG	1:C:64:THR:N	2.41	0.51
1:C:301:MET:HE2	1:C:309:LEU:HB3	1.92	0.51
2:B:263:SER:HB2	2:B:268:ILE:HB	1.91	0.51
2:B:126:ALA:HB1	4:E:247:LEU:HD22	1.92	0.51
1:C:459:LYS:HA	1:C:470:VAL:C	2.36	0.51
1:C:159:PHE:CZ	3:D:264:LYS:HE2	2.46	0.51
1:A:373:VAL:HB	1:A:391:LEU:HB2	1.92	0.51
1:C:498:LEU:HD23	1:C:501:LYS:HD2	1.93	0.51
8:O:2:NAG:H61	8:O:3:BMA:O5	2.11	0.50
2:B:215:ILE:HD12	2:B:215:ILE:H	1.76	0.50
1:C:544:ASP:HB3	1:C:547:GLU:HG3	1.94	0.50
3:D:87:ILE:HD12	3:D:420:PRO:HD3	1.93	0.50
3:D:188:LEU:HD12	3:D:214:LYS:HB2	1.93	0.50
1:A:551:LYS:HB3	1:A:594:LEU:HD11	1.92	0.50
3:D:174:GLU:CD	3:D:174:GLU:H	2.20	0.50
2:B:137:GLU:O	2:B:141:ARG:HG3	2.12	0.49
1:A:1:PHE:HA	1:A:389:GLN:HB2	1.95	0.49
2:B:77:ARG:HB2	2:B:112:GLU:HB2	1.95	0.49
2:B:115:PRO:HB2	2:B:245:SER:CB	2.42	0.49
1:A:414:ILE:HG21	1:A:434:LEU:HD21	1.93	0.49
1:C:50:ILE:HD11	1:C:76:GLY:HA2	1.93	0.49
3:D:142:LEU:HA	3:D:344:LEU:HD11	1.94	0.49
3:D:204:GLU:HA	3:D:207:ASN:HD22	1.77	0.49
3:D:126:ALA:HB1	4:F:247:LEU:HD12	1.94	0.49
3:D:208:GLU:O	3:D:212:ASN:ND2	2.40	0.49
1:C:463:LEU:HD21	1:C:517:ARG:HE	1.76	0.49
3:D:322:GLU:CG	3:D:333:VAL:HG21	2.43	0.49
1:A:523:LYS:HG2	1:A:536:GLU:OE1	2.13	0.49
3:D:353:GLU:O	3:D:357:GLU:HG3	2.13	0.48
1:A:9:ALA:HB3	1:A:434:LEU:HB3	1.95	0.48
2:B:105:GLN:NE2	14:B:606:HOH:O	2.45	0.48
1:A:575:GLN:HA	1:A:576:PRO:HD3	1.73	0.48
1:A:114:TRP:CZ2	1:A:116:THR:HA	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:543:ARG:HD3	1:A:547:GLU:OE1	2.13	0.48
1:C:350:LEU:HA	1:C:420:PRO:HG2	1.96	0.48
1:A:487:LEU:CD1	1:A:529:ARG:HD3	2.43	0.48
3:D:161:VAL:HG21	3:D:216:SER:CB	2.44	0.48
2:B:94:LEU:HG	2:B:96:LEU:HG	1.96	0.48
1:C:579:ASN:OD1	1:C:580:GLN:N	2.47	0.48
3:D:416:ILE:HB	3:D:429:LEU:HB2	1.96	0.48
1:A:529:ARG:C	1:A:531:GLY:H	2.21	0.47
1:A:472:CYS:HA	1:A:541:TYR:HA	1.96	0.47
1:C:431:ARG:NH2	14:C:805:HOH:O	2.41	0.47
1:A:459:LYS:HA	1:A:470:VAL:C	2.40	0.47
3:D:73:LEU:HD23	3:D:85:VAL:HG12	1.95	0.47
1:C:523:LYS:HG2	1:C:536:GLU:OE1	2.14	0.47
1:A:77:ASN:ND2	9:A:703:SO4:O2	2.45	0.47
3:D:322:GLU:HG2	3:D:333:VAL:HG21	1.97	0.47
1:A:397:ALA:HB2	1:A:402:PRO:HD3	1.97	0.47
2:B:9:GLY:HA3	2:B:20:ILE:HD13	1.97	0.47
1:C:154:PHE:O	1:C:175:GLY:HA3	2.15	0.47
1:C:250:LEU:HD12	1:C:250:LEU:HA	1.79	0.47
1:A:10:GLU:N	1:A:65:ARG:HH11	2.06	0.47
2:B:13:THR:HG22	2:B:15:GLU:H	1.79	0.47
2:B:115:PRO:HB2	2:B:245:SER:HB3	1.96	0.47
2:B:362:VAL:HG12	2:B:366:THR:HG21	1.97	0.47
2:B:118:LEU:O	2:B:155:LEU:HA	2.15	0.46
3:D:65:VAL:HG12	3:D:94:LEU:HD13	1.97	0.46
1:A:509:ARG:HG2	1:A:519:PRO:HG3	1.98	0.46
2:B:87:ILE:HG22	2:B:425:ASP:HB3	1.98	0.46
3:D:94:LEU:HG	3:D:96:LEU:HG	1.96	0.46
1:A:36:LEU:HB2	1:A:59:CYS:HB2	1.97	0.46
1:A:115:ARG:O	1:A:116:THR:OG1	2.29	0.46
3:D:308:LEU:HD13	3:D:354:LEU:HD13	1.98	0.46
2:B:357:GLU:OE1	2:B:388:LYS:NZ	2.49	0.46
1:A:329:THR:HG22	1:A:330:LYS:H	1.81	0.46
2:B:154:ARG:HH12	2:B:242:ARG:NH2	2.14	0.46
1:C:329:THR:HG22	1:C:330:LYS:H	1.81	0.46
1:A:26:PHE:HB3	1:A:35:PHE:HB2	1.98	0.46
1:C:116:THR:HG23	1:C:147:ILE:HG21	1.97	0.46
3:D:410:GLU:HB2	3:D:433:PRO:HG2	1.98	0.46
2:B:308:LEU:HD13	2:B:354:LEU:HD13	1.98	0.45
1:A:449:VAL:HG21	1:A:557:ILE:HD13	1.98	0.45
1:A:464:PRO:O	1:A:517:ARG:NH2	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:247:ALA:O	1:C:250:LEU:HB2	2.17	0.45
3:D:264:LYS:HG3	14:D:2101:HOH:O	2.16	0.45
3:D:356:SER:HB2	3:D:392:MET:H	1.79	0.45
3:D:220:ASP:OD2	3:D:258:HIS:NE2	2.43	0.45
2:B:22:PRO:HA	2:B:97:ARG:HH11	1.81	0.45
1:A:482:ASP:HA	1:A:529:ARG:HG3	1.98	0.45
2:B:359:GLU:HG3	2:B:388:LYS:HB2	1.97	0.45
1:C:580:GLN:C	1:C:582:THR:H	2.25	0.45
5:P:1:NAG:H61	5:P:2:NAG:H82	1.99	0.45
2:B:454:ASN:HB3	2:B:473:CYS:SG	2.57	0.45
1:C:498:LEU:HD23	1:C:501:LYS:CD	2.47	0.45
3:D:87:ILE:HG22	3:D:425:ASP:HB3	1.98	0.45
1:C:448:GLU:OE1	13:C:708:MAN:O3	2.24	0.45
1:C:449:VAL:HG21	1:C:557:ILE:HD13	1.99	0.45
2:B:197:LEU:HB2	2:B:209:ILE:CD1	2.47	0.44
1:A:587:SER:O	1:A:588:ARG:HD3	2.17	0.44
3:D:54:GLN:O	3:D:58:ILE:HG13	2.16	0.44
1:C:175:GLY:HA2	1:C:179:TRP:CD1	2.52	0.44
1:A:250:LEU:HD23	1:A:250:LEU:HA	1.85	0.44
1:A:419:TYR:CE1	1:A:439:PRO:HA	2.53	0.44
2:B:120:TYR:CZ	2:B:252:VAL:HG21	2.52	0.44
1:A:111:LEU:HD23	1:A:111:LEU:HA	1.88	0.43
1:A:274:ALA:HA	1:A:299:LEU:HB2	2.00	0.43
1:A:509:ARG:HG2	1:A:519:PRO:CG	2.48	0.43
2:B:414:ARG:HG3	2:B:431:VAL:HB	2.00	0.43
1:C:190:GLU:HG2	1:C:194:LYS:HD3	1.99	0.43
3:D:109:ARG:NH1	3:D:395:GLY:O	2.47	0.43
2:B:154:ARG:HH21	2:B:199:LEU:CD1	2.31	0.43
1:A:579:ASN:OD1	1:A:580:GLN:N	2.51	0.43
2:B:351:TYR:O	2:B:355:ARG:HG2	2.19	0.43
3:D:130:ASP:OD1	3:D:131:ASP:N	2.52	0.43
2:B:359:GLU:HB3	2:B:421:VAL:HG21	1.99	0.43
3:D:11:ALA:HA	3:D:26:TRP:HZ3	1.83	0.43
1:C:47:GLN:HA	1:C:48:PRO:HD3	1.85	0.43
3:D:301:LYS:HD3	3:D:301:LYS:HA	1.84	0.43
7:N:3:BMA:H62	7:N:4:MAN:H2	1.75	0.43
1:A:477:PHE:HB2	1:A:559:MET:SD	2.59	0.43
1:A:339:ARG:HD2	1:A:364:TYR:CE2	2.53	0.43
3:D:123:ASP:OD2	3:D:125:SER:OG	2.37	0.43
1:A:253:VAL:HB	1:A:267:PHE:HB2	2.00	0.42
1:A:284:ASP:O	1:A:356:ASN:HB2	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:74:SER:OG	2:B:85:VAL:N	2.46	0.42
2:B:114:TYR:HA	2:B:115:PRO:HD3	1.75	0.42
1:C:373:VAL:HB	1:C:391:LEU:HB2	2.00	0.42
1:A:170:LEU:HD22	1:A:239:PHE:CD1	2.54	0.42
2:B:161:VAL:HG21	2:B:216:SER:CB	2.47	0.42
2:B:356:SER:O	2:B:392:MET:HG2	2.19	0.42
3:D:352:GLU:HA	3:D:355:ARG:HB2	2.01	0.42
1:A:59:CYS:HA	1:A:67:CYS:HA	2.01	0.42
1:A:483:GLY:HA3	1:A:487:LEU:HD21	1.99	0.42
2:B:183:ILE:HD11	4:E:244:LEU:HD22	2.01	0.42
2:B:362:VAL:HG21	2:B:373:PHE:HZ	1.84	0.42
3:D:18:LEU:HD13	3:D:61:PRO:HG3	2.01	0.42
1:C:487:LEU:HG	1:C:529:ARG:HD3	2.01	0.42
3:D:152:ASN:CB	14:D:2102:HOH:O	2.68	0.42
3:D:163:LYS:HE2	3:D:292:TYR:CE1	2.55	0.42
3:D:360:LEU:HD23	3:D:360:LEU:HA	1.70	0.42
2:B:361:GLU:OE2	2:B:421:VAL:HG22	2.20	0.42
1:C:510:ARG:NH1	1:C:553:THR:O	2.47	0.42
1:A:518:SER:HA	1:A:519:PRO:HD3	1.88	0.42
2:B:122:MET:HE2	2:B:157:PHE:HZ	1.85	0.42
1:C:477:PHE:HB2	1:C:559:MET:SD	2.60	0.42
1:C:500:ASP:OD1	1:C:503:LYS:HG2	2.20	0.42
1:A:301:MET:HE2	1:A:309:LEU:HB3	2.02	0.42
1:A:462:SER:HA	1:A:469:LYS:HG2	2.02	0.42
1:C:575:GLN:HA	1:C:576:PRO:HD3	1.79	0.42
3:D:223:GLU:HB2	14:D:2115:HOH:O	2.20	0.42
2:B:76:GLY:O	2:B:77:ARG:HB3	2.19	0.42
1:C:582:THR:HA	1:C:583:PRO:HD3	1.86	0.42
3:D:67:ILE:HA	3:D:92:LEU:HD23	2.02	0.42
1:A:503:LYS:NZ	1:A:550:ASP:OD2	2.40	0.42
1:C:306:ASP:OD1	1:C:306:ASP:N	2.36	0.42
1:C:177:PHE:O	1:C:180:GLN:HG3	2.19	0.41
1:A:47:GLN:HA	1:A:48:PRO:HD3	1.89	0.41
1:C:580:GLN:O	1:C:581:PHE:HB3	2.20	0.41
3:D:154:ARG:HH12	3:D:242:ARG:CZ	2.33	0.41
1:A:50:ILE:HD11	1:A:76:GLY:HA2	2.02	0.41
2:B:365:ASP:HB3	2:B:414:ARG:HB2	2.01	0.41
1:C:500:ASP:HB2	1:C:555:ILE:HG23	2.01	0.41
1:C:525:MET:HE2	1:C:534:GLN:HB3	2.01	0.41
2:B:373:PHE:CD2	2:B:402:VAL:HG22	2.56	0.41
1:C:317:VAL:O	1:C:328:THR:HA	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:M:1:NAG:H61	5:M:2:NAG:C7	2.49	0.41
1:A:93:TRP:CD1	1:A:111:LEU:HD12	2.56	0.41
1:A:461:CYS:N	1:A:472:CYS:SG	2.92	0.41
1:A:510:ARG:NH1	1:A:553:THR:O	2.48	0.41
2:B:308:LEU:HD21	2:B:421:VAL:HG12	2.02	0.41
2:B:355:ARG:HH21	2:B:394:VAL:HG11	1.86	0.41
1:C:299:LEU:HD22	3:D:294:THR:HG21	2.03	0.41
3:D:345:GLN:OE1	3:D:345:GLN:N	2.39	0.41
2:B:223:GLU:HB2	14:B:628:HOH:O	2.20	0.41
2:B:334:GLY:HA3	2:B:346:LEU:HD22	2.02	0.41
3:D:124:LEU:HD23	3:D:124:LEU:HA	1.94	0.41
3:D:153:PHE:CE2	3:D:155:LEU:HG	2.56	0.41
1:C:288:ASP:OD1	1:C:288:ASP:N	2.54	0.41
2:B:154:ARG:NH2	2:B:242:ARG:HG2	2.36	0.41
3:D:334:GLY:HA3	3:D:346:LEU:HD22	2.03	0.41
1:A:185:SER:HB3	1:A:208:LEU:HB2	2.02	0.41
1:C:419:TYR:CE1	1:C:439:PRO:HA	2.56	0.41
2:B:75:VAL:HB	2:B:109:ARG:NH1	2.36	0.40
2:B:194:LYS:HG2	2:B:283:GLU:HG2	2.02	0.40
1:C:375:ILE:CD1	1:C:391:LEU:HD13	2.51	0.40
2:B:272:ASN:HD22	2:B:293:PRO:HD3	1.87	0.40
3:D:163:LYS:HA	3:D:164:PRO:HD3	1.87	0.40
3:D:126:ALA:HA	3:D:215:ILE:CG2	2.51	0.40
2:B:345:GLN:OE1	2:B:345:GLN:N	2.52	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:452:SER:OG	1:C:452:SER:OG[4_555]	2.19	0.01

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	592/604 (98%)	567 (96%)	24 (4%)	1 (0%)	43	63
1	C	583/604 (96%)	564 (97%)	18 (3%)	1 (0%)	43	63
2	B	438/483 (91%)	418 (95%)	20 (5%)	0	100	100
3	D	447/483 (92%)	424 (95%)	21 (5%)	2 (0%)	30	49
4	E	9/13 (69%)	9 (100%)	0	0	100	100
4	F	11/13 (85%)	11 (100%)	0	0	100	100
All	All	2080/2200 (94%)	1993 (96%)	83 (4%)	4 (0%)	43	63

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	593	LEU
1	C	459	LYS
3	D	8	LEU
3	D	77	ARG

5.3.2 Protein sidechains ⓘ

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	487/495 (98%)	478 (98%)	9 (2%)	51	77
1	C	483/495 (98%)	472 (98%)	11 (2%)	44	72
2	B	393/421 (93%)	387 (98%)	6 (2%)	57	80
3	D	398/421 (94%)	390 (98%)	8 (2%)	48	75
4	E	7/8 (88%)	7 (100%)	0	100	100
4	F	8/8 (100%)	8 (100%)	0	100	100
All	All	1776/1848 (96%)	1742 (98%)	34 (2%)	50	76

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	VAL
1	A	27	VAL
1	A	29	SER
1	A	60	ASP
1	A	275	TYR
1	A	512	LEU
1	A	525	MET
1	A	528	SER
1	A	591	HIS
2	B	169	VAL
2	B	209	ILE
2	B	218	ASN
2	B	257	SER
2	B	320	LEU
2	B	384	GLN
1	C	5	VAL
1	C	27	VAL
1	C	51	VAL
1	C	60	ASP
1	C	183	LEU
1	C	275	TYR
1	C	479	LEU
1	C	508	ILE
1	C	514	LEU
1	C	525	MET
1	C	538	LEU
3	D	56	ASN
3	D	155	LEU
3	D	169	VAL
3	D	174	GLU
3	D	209	ILE
3	D	218	ASN
3	D	320	LEU
3	D	384	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	HIS
1	A	145	GLN
2	B	64	GLN
2	B	78	GLN
2	B	133	ASN

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Mol	Chain	Res	Type
2	B	317	GLN
2	B	436	ASN
2	B	454	ASN
1	C	77	ASN
1	C	91	HIS
1	C	327	GLN
1	C	494	GLN
3	D	110	GLN
3	D	436	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

31 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	G	1	5,1	14,14,15	0.27	0	17,19,21	0.51	0
5	NAG	G	2	5	14,14,15	0.32	0	17,19,21	0.56	0
6	NAG	H	1	6,1	14,14,15	0.20	0	17,19,21	0.41	0
6	NAG	H	2	6	14,14,15	0.19	0	17,19,21	0.45	0
6	BMA	H	3	6	11,11,12	0.71	0	15,15,17	0.85	0
6	MAN	H	4	6	11,11,12	0.67	0	15,15,17	0.97	2 (13%)
7	NAG	I	1	1,7	14,14,15	0.49	0	17,19,21	0.43	0
7	NAG	I	2	7	14,14,15	0.39	0	17,19,21	0.43	0
7	BMA	I	3	7	11,11,12	0.86	1 (9%)	15,15,17	0.74	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
7	MAN	I	4	7	11,11,12	0.77	1 (9%)	15,15,17	0.87	1 (6%)
7	MAN	I	5	7	11,11,12	0.69	0	15,15,17	1.00	2 (13%)
7	MAN	I	6	7	11,11,12	0.72	0	15,15,17	0.91	2 (13%)
5	NAG	J	1	5,1	14,14,15	0.71	0	17,19,21	0.49	0
5	NAG	J	2	5	14,14,15	0.31	0	17,19,21	0.48	0
5	NAG	K	1	5,1	14,14,15	0.57	0	17,19,21	0.67	1 (5%)
5	NAG	K	2	5	14,14,15	0.33	0	17,19,21	0.42	0
5	NAG	L	1	5,1	14,14,15	0.36	0	17,19,21	0.43	0
5	NAG	L	2	5	14,14,15	0.34	0	17,19,21	0.47	0
5	NAG	M	1	5,1	14,14,15	0.20	0	17,19,21	0.51	0
5	NAG	M	2	5	14,14,15	0.23	0	17,19,21	0.50	0
7	NAG	N	1	1,7	14,14,15	0.47	0	17,19,21	0.52	0
7	NAG	N	2	7	14,14,15	0.27	0	17,19,21	0.42	0
7	BMA	N	3	7	11,11,12	0.78	1 (9%)	15,15,17	0.77	1 (6%)
7	MAN	N	4	7	11,11,12	0.63	0	15,15,17	1.01	2 (13%)
7	MAN	N	5	7	11,11,12	0.68	0	15,15,17	0.99	2 (13%)
7	MAN	N	6	7	11,11,12	0.68	0	15,15,17	0.92	2 (13%)
8	NAG	O	1	8,1	14,14,15	0.71	1 (7%)	17,19,21	0.46	0
8	NAG	O	2	8	14,14,15	0.71	1 (7%)	17,19,21	0.98	1 (5%)
8	BMA	O	3	8	11,11,12	0.95	1 (9%)	15,15,17	2.34	4 (26%)
5	NAG	P	1	5,1	14,14,15	0.48	0	17,19,21	0.42	0
5	NAG	P	2	5	14,14,15	0.48	0	17,19,21	0.53	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	G	1	5,1	-	2/6/23/26	0/1/1/1
5	NAG	G	2	5	-	0/6/23/26	0/1/1/1
6	NAG	H	1	6,1	-	0/6/23/26	0/1/1/1
6	NAG	H	2	6	-	4/6/23/26	0/1/1/1
6	BMA	H	3	6	-	1/2/19/22	0/1/1/1
6	MAN	H	4	6	-	0/2/19/22	0/1/1/1
7	NAG	I	1	1,7	-	0/6/23/26	0/1/1/1
7	NAG	I	2	7	-	0/6/23/26	0/1/1/1
7	BMA	I	3	7	-	1/2/19/22	0/1/1/1
7	MAN	I	4	7	-	0/2/19/22	0/1/1/1
7	MAN	I	5	7	-	0/2/19/22	0/1/1/1

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	O	2	NAG	O5-C1	-2.52	1.39	1.43
7	I	3	BMA	O5-C1	-2.41	1.39	1.43
8	O	1	NAG	O5-C1	-2.37	1.39	1.43
8	O	3	BMA	C1-C2	2.32	1.57	1.52
7	I	4	MAN	O5-C1	-2.14	1.40	1.43
7	N	3	BMA	O5-C1	-2.06	1.40	1.43

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	O	3	BMA	C1-C2-C3	-5.35	101.85	109.64
8	O	3	BMA	O3-C3-C2	5.18	120.63	110.05
8	O	3	BMA	C3-C4-C5	-3.15	104.52	110.23
8	O	2	NAG	O4-C4-C3	-2.82	103.72	110.38
7	I	5	MAN	C1-O5-C5	2.66	115.75	112.19
7	N	4	MAN	C1-O5-C5	2.54	115.58	112.19
6	H	4	MAN	C1-O5-C5	2.49	115.52	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	N	5	MAN	C1-O5-C5	2.45	115.47	112.19
7	N	4	MAN	O2-C2-C3	-2.35	105.29	110.15
7	N	6	MAN	C1-O5-C5	2.24	115.19	112.19
5	K	1	NAG	C1-O5-C5	2.18	115.11	112.19
7	I	6	MAN	O2-C2-C3	-2.14	105.71	110.15
7	I	4	MAN	O2-C2-C3	-2.13	105.75	110.15
7	I	6	MAN	C1-O5-C5	2.12	115.03	112.19
7	N	6	MAN	O2-C2-C3	-2.11	105.78	110.15
7	I	5	MAN	O2-C2-C3	-2.11	105.79	110.15
6	H	4	MAN	O2-C2-C3	-2.08	105.85	110.15
7	N	3	BMA	O2-C2-C3	-2.07	105.86	110.15
8	O	3	BMA	O5-C5-C4	-2.06	105.81	110.83
7	N	5	MAN	O2-C2-C3	-2.02	105.96	110.15

There are no chirality outliers.

All (21) torsion outliers are listed below:

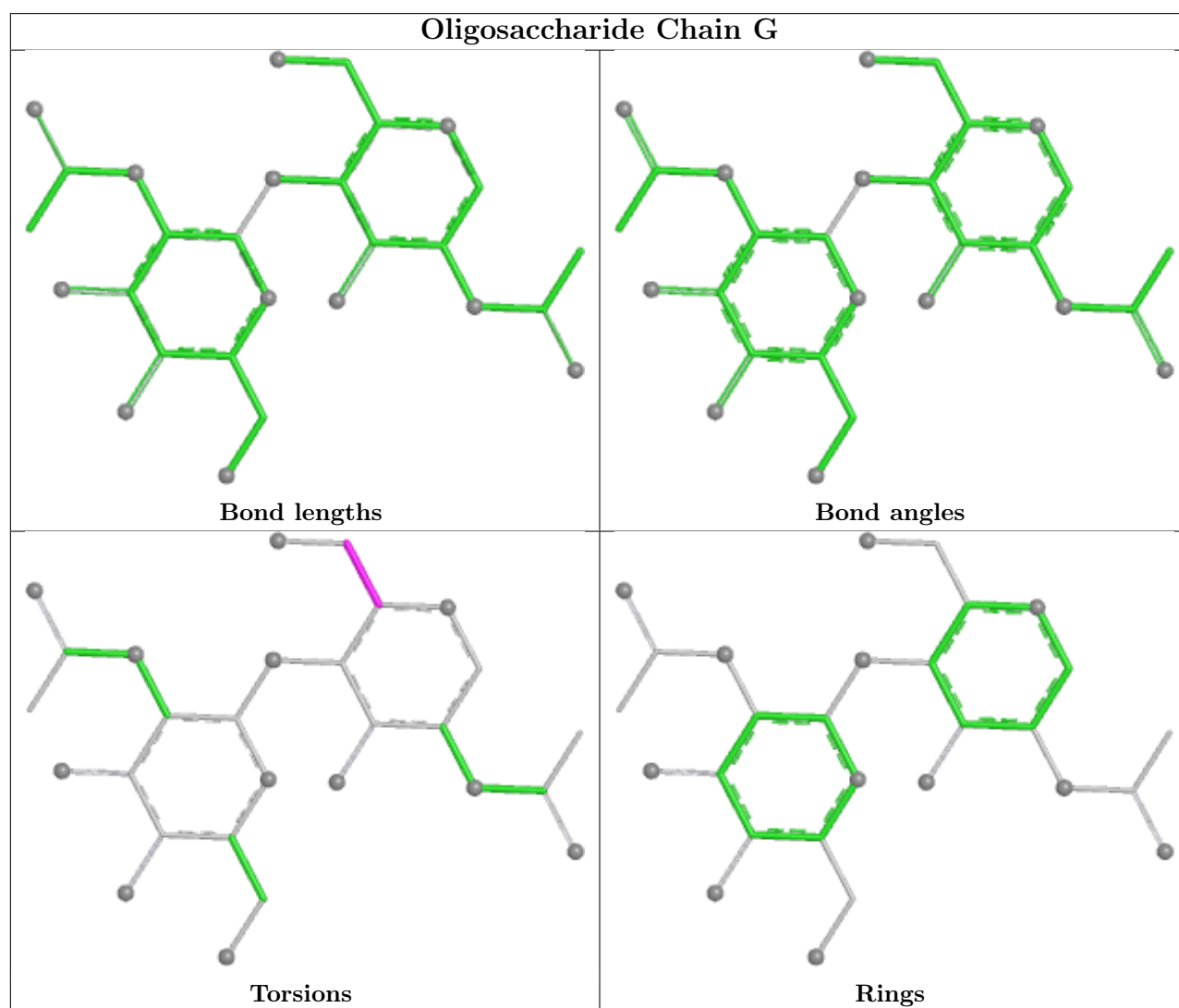
Mol	Chain	Res	Type	Atoms
6	H	2	NAG	O5-C5-C6-O6
5	G	1	NAG	O5-C5-C6-O6
5	K	1	NAG	O5-C5-C6-O6
6	H	2	NAG	C8-C7-N2-C2
6	H	2	NAG	O7-C7-N2-C2
6	H	2	NAG	C4-C5-C6-O6
5	G	1	NAG	C4-C5-C6-O6
5	K	1	NAG	C4-C5-C6-O6
7	N	3	BMA	O5-C5-C6-O6
7	N	5	MAN	O5-C5-C6-O6
5	L	1	NAG	C4-C5-C6-O6
7	I	6	MAN	O5-C5-C6-O6
7	N	6	MAN	O5-C5-C6-O6
7	I	3	BMA	O5-C5-C6-O6
5	P	2	NAG	O5-C5-C6-O6
5	K	2	NAG	O5-C5-C6-O6
5	M	1	NAG	C4-C5-C6-O6
5	L	1	NAG	O5-C5-C6-O6
5	M	1	NAG	O5-C5-C6-O6
5	M	2	NAG	O5-C5-C6-O6
6	H	3	BMA	C4-C5-C6-O6

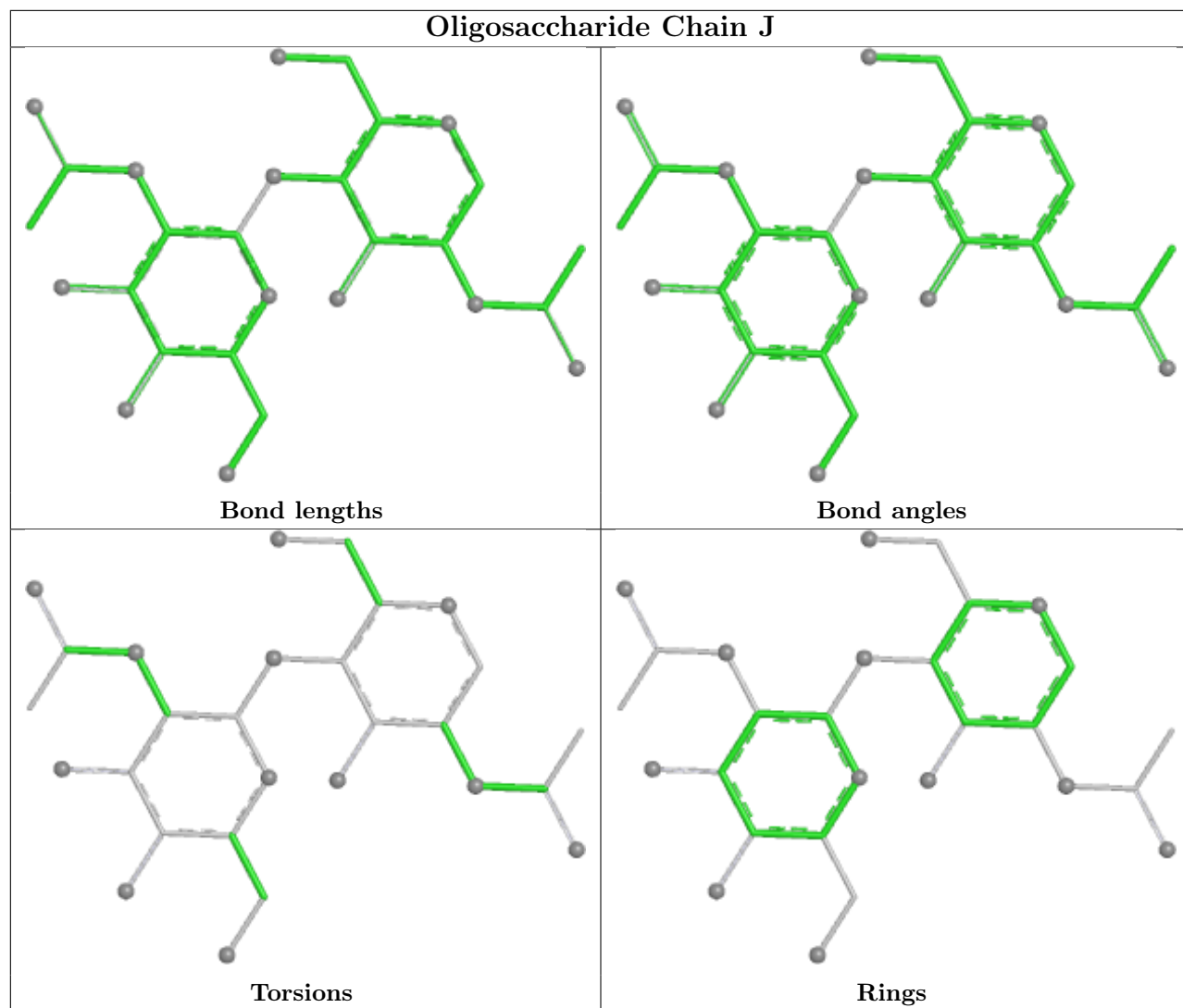
There are no ring outliers.

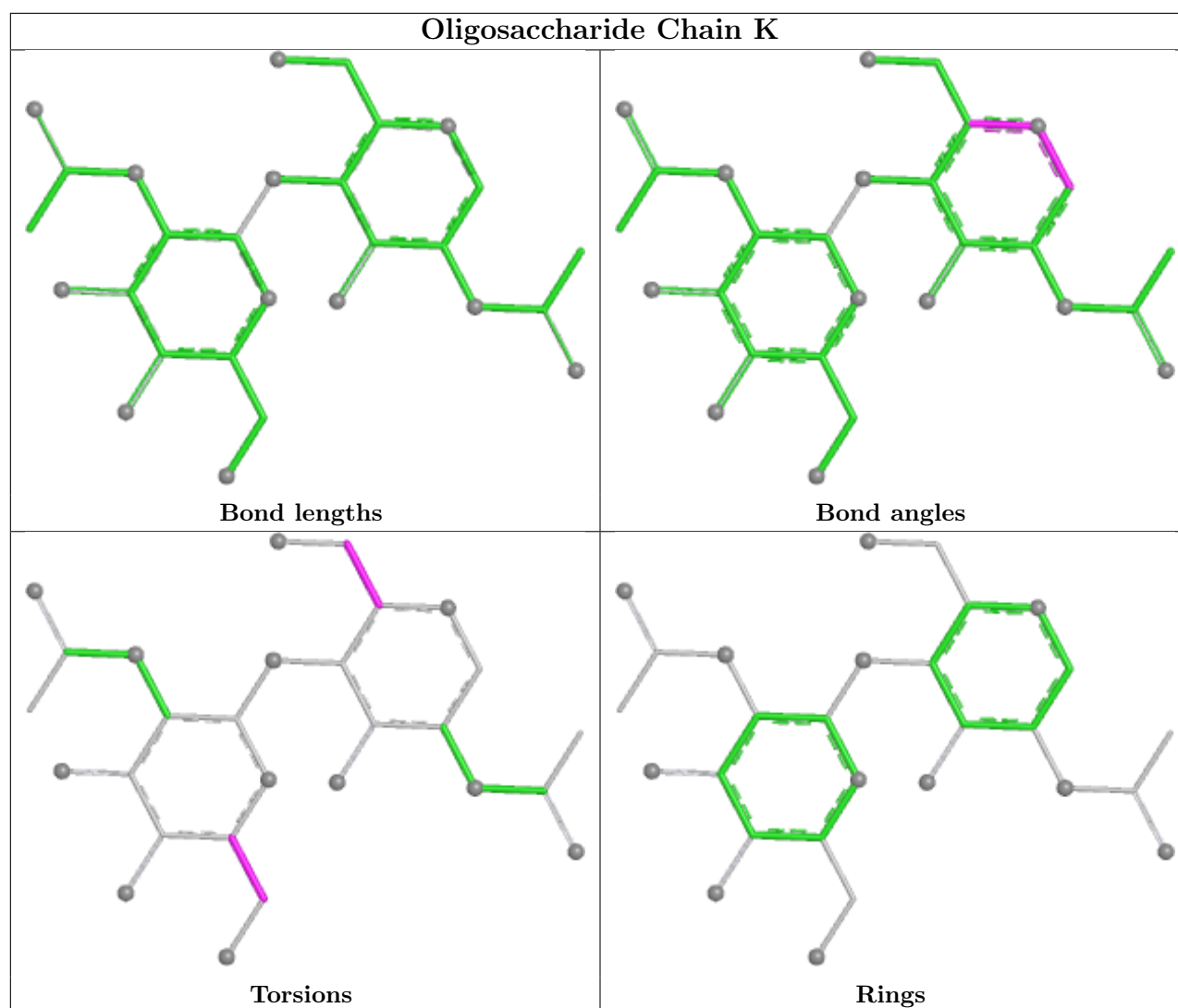
9 monomers are involved in 5 short contacts:

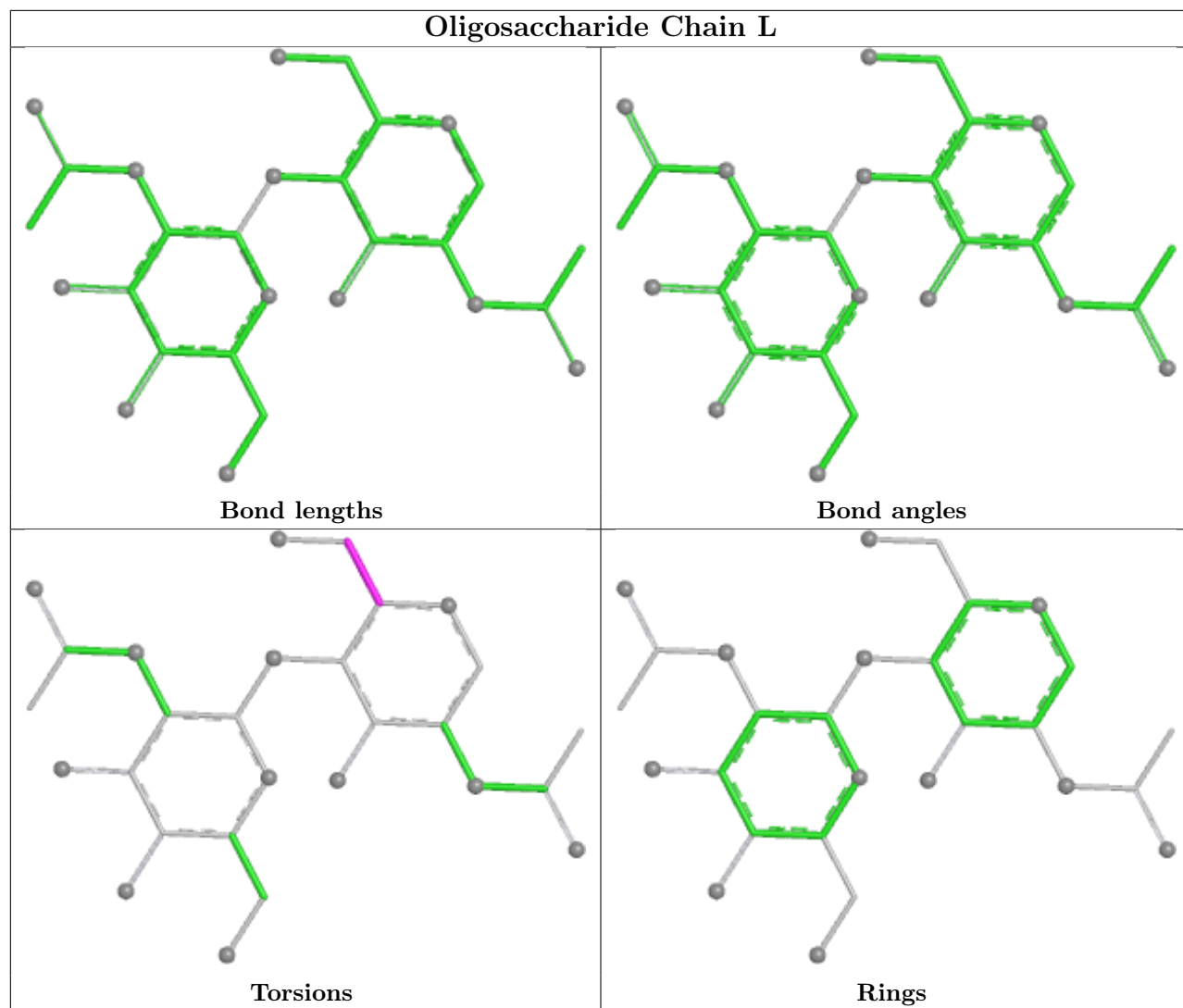
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	P	2	NAG	1	0
8	O	2	NAG	2	0
8	O	3	BMA	1	0
7	N	3	BMA	1	0
8	O	1	NAG	1	0
7	N	4	MAN	1	0
5	M	2	NAG	1	0
5	P	1	NAG	1	0
5	M	1	NAG	1	0

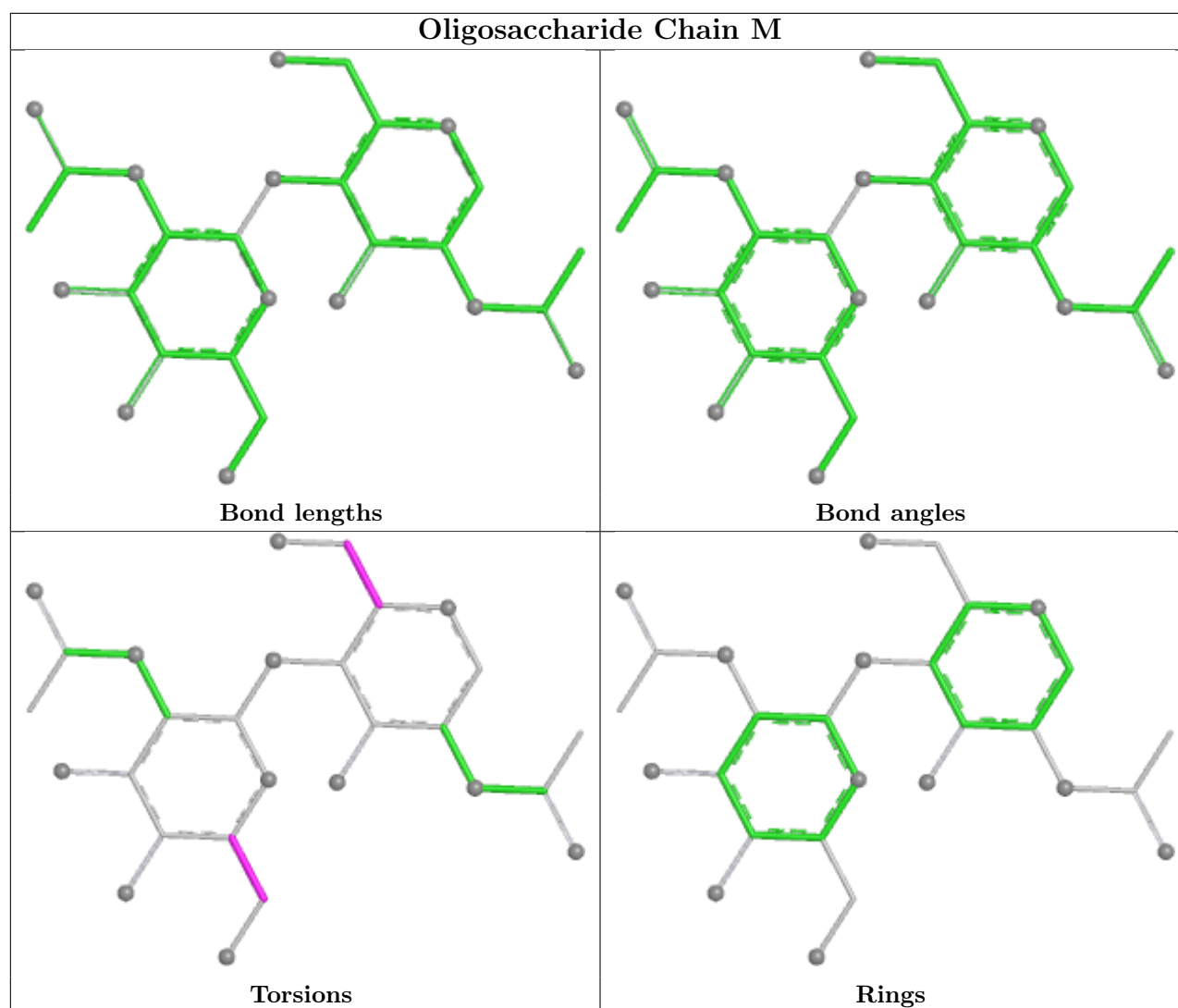
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

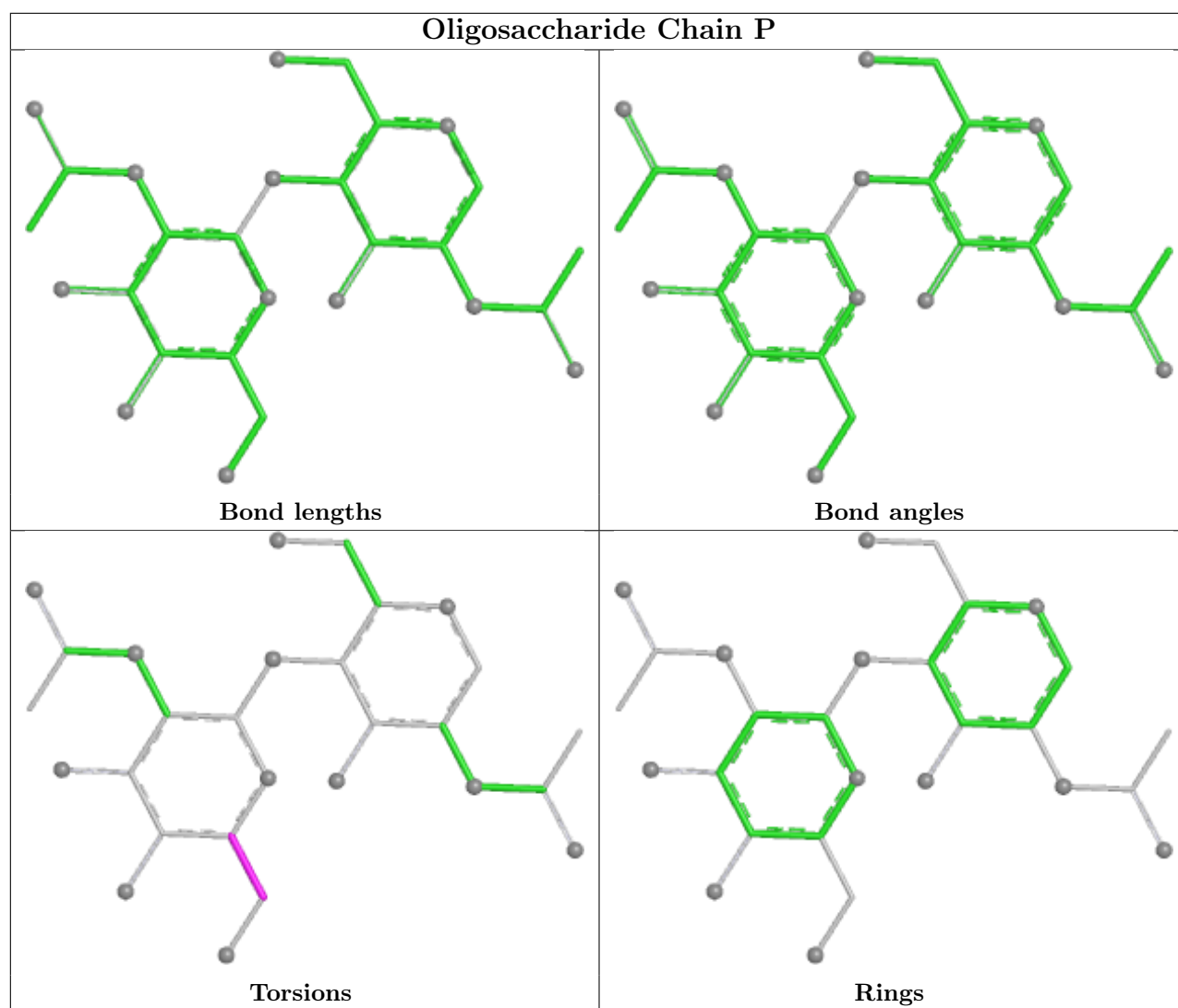




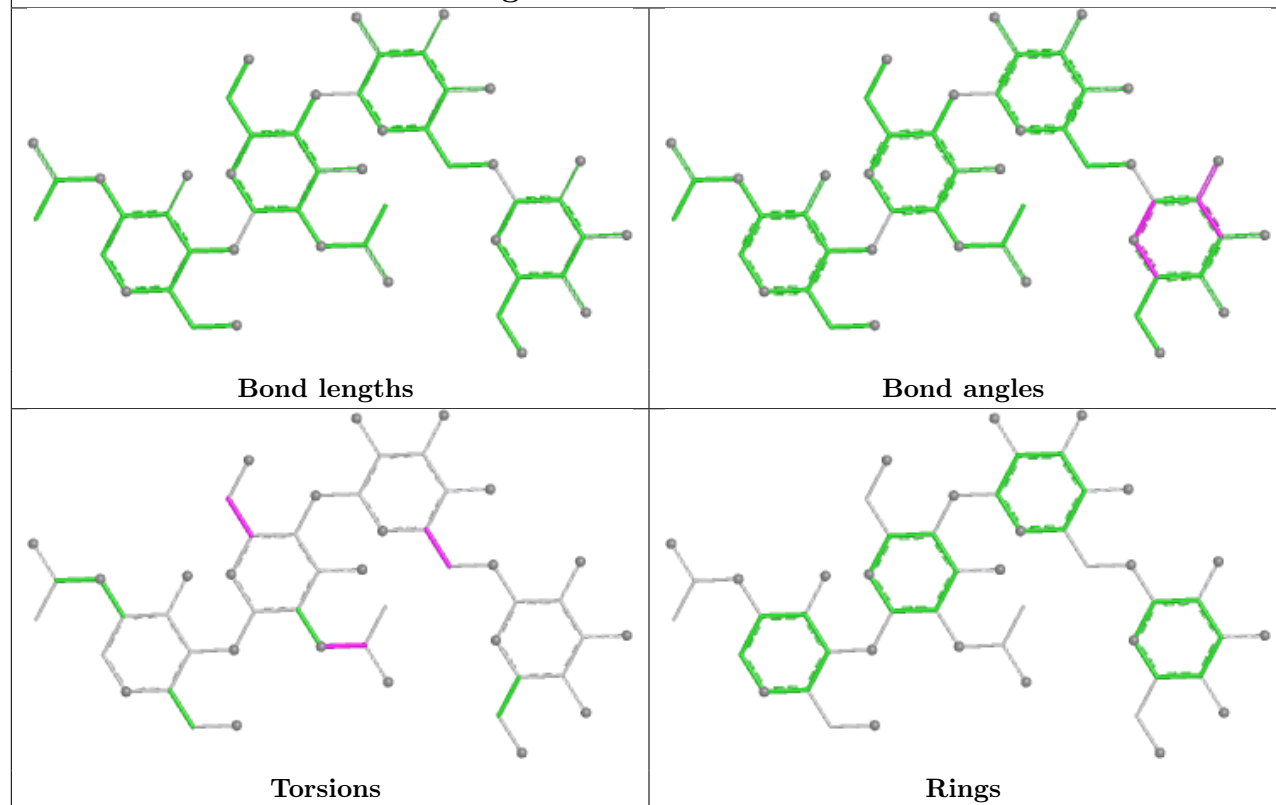




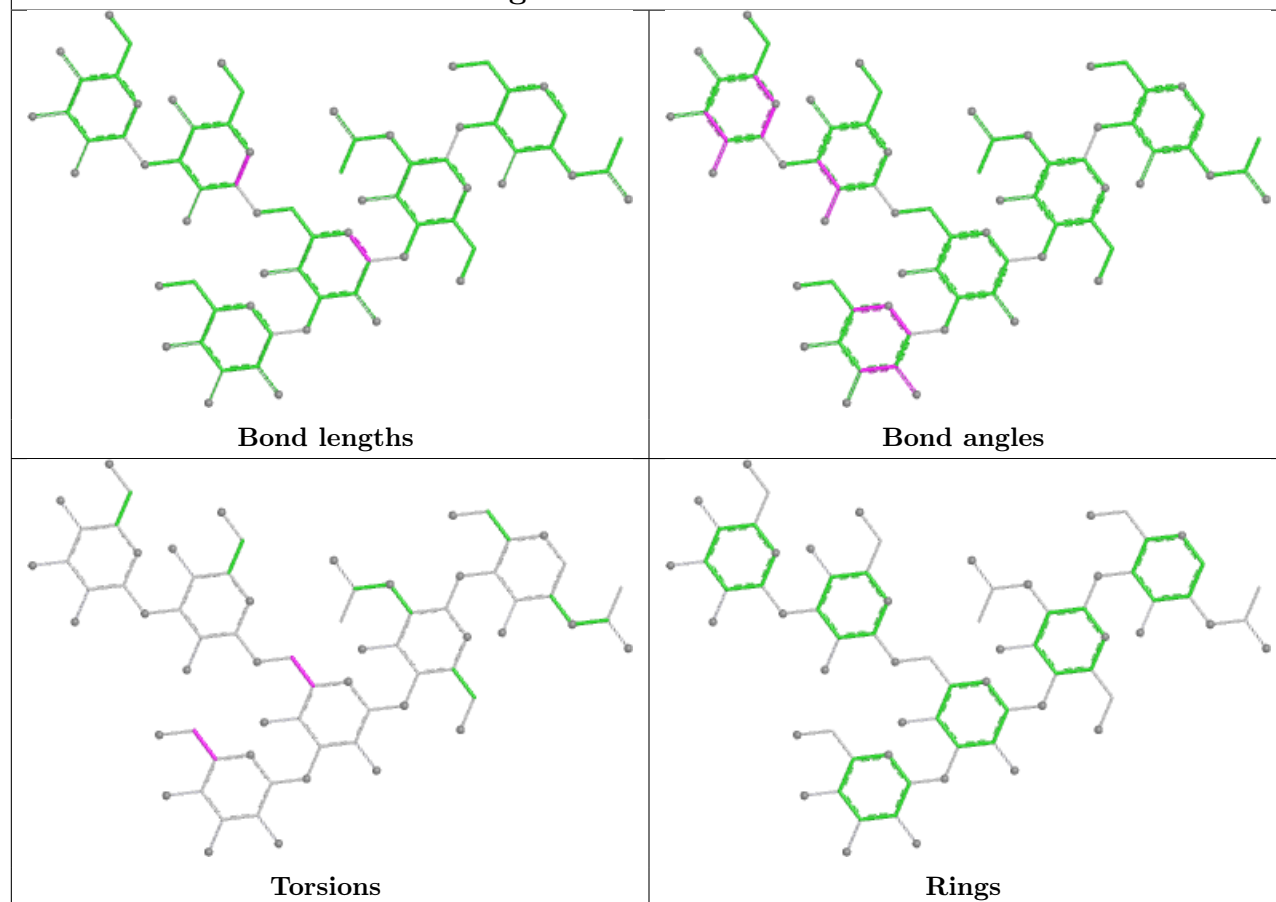




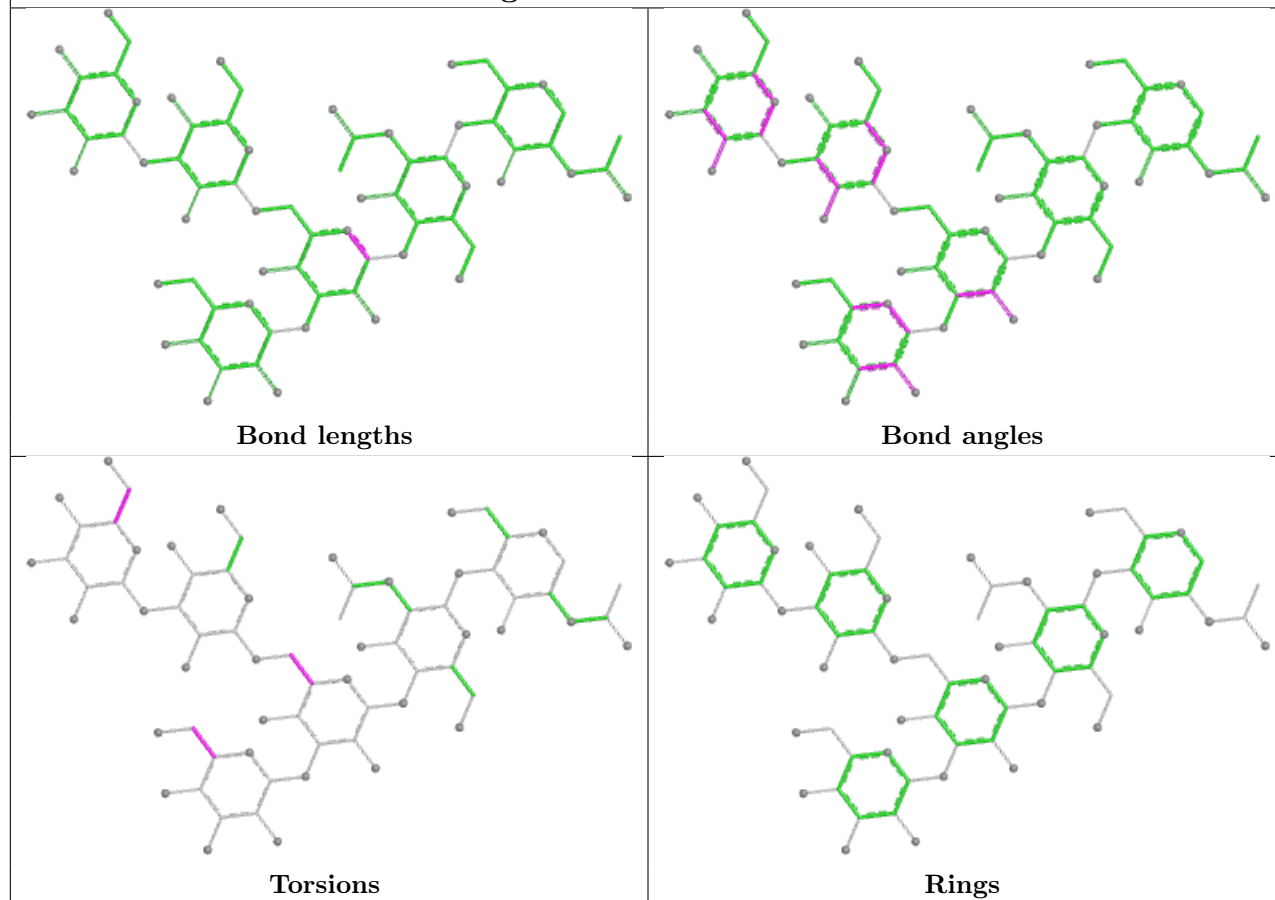
Oligosaccharide Chain H



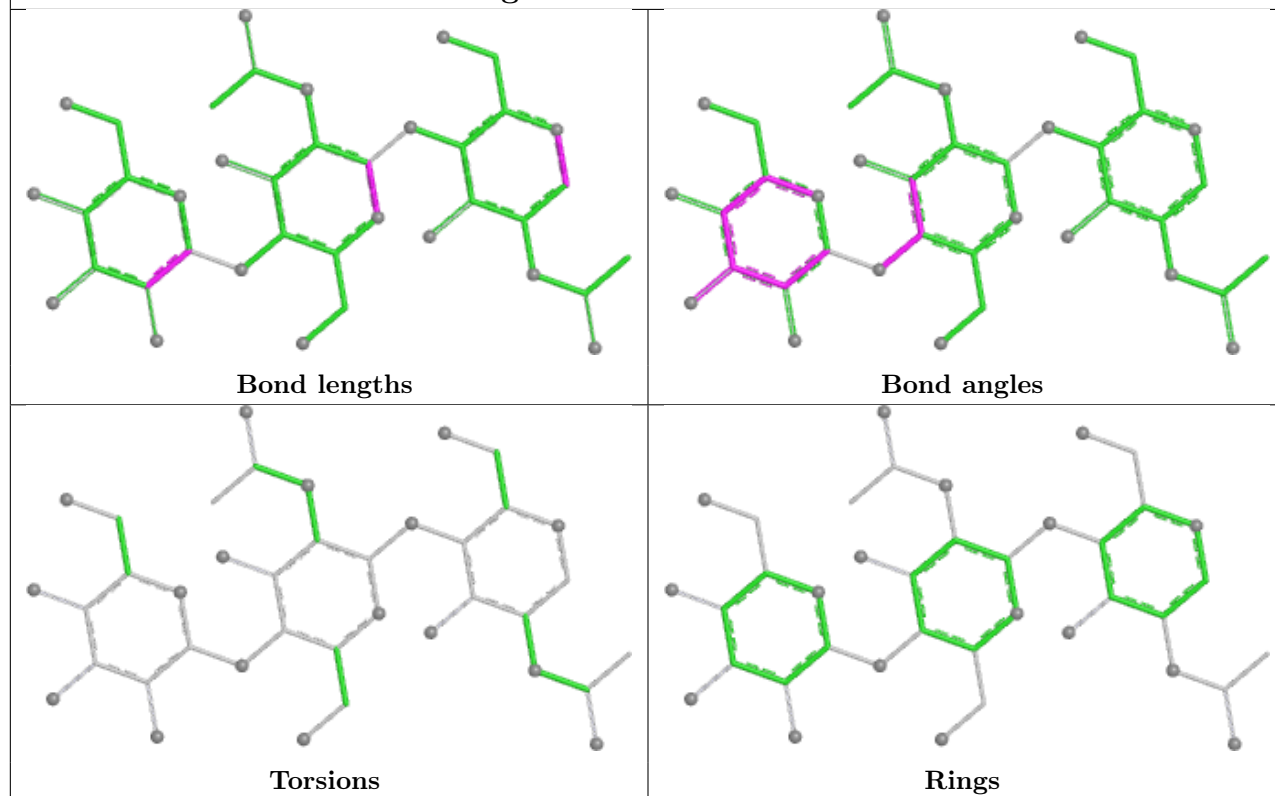
Oligosaccharide Chain I



Oligosaccharide Chain N



Oligosaccharide Chain O



5.6 Ligand geometry

Of 29 ligands modelled in this entry, 14 are monoatomic - leaving 15 for Mogul analysis.

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
12	NAG	C	707	1	14,14,15	0.20	0	17,19,21	0.38	0
9	SO4	A	702	-	4,4,4	0.23	0	6,6,6	0.07	0
9	SO4	A	703	-	4,4,4	0.22	0	6,6,6	0.16	0
12	NAG	C	709	1	14,14,15	0.46	0	17,19,21	0.43	0
9	SO4	A	701	-	4,4,4	0.24	0	6,6,6	0.10	0
9	SO4	B	501	-	4,4,4	0.23	0	6,6,6	0.08	0
9	SO4	C	702	-	4,4,4	0.24	0	6,6,6	0.05	0
9	SO4	C	701	-	4,4,4	0.24	0	6,6,6	0.15	0
13	MAN	C	708	-	11,11,12	1.48	3 (27%)	15,15,17	1.57	3 (20%)
12	NAG	D	2004	3	14,14,15	0.35	0	17,19,21	0.59	0
12	NAG	B	506	2	14,14,15	0.34	0	17,19,21	0.52	0
12	NAG	D	2005	3	14,14,15	0.37	0	17,19,21	0.46	0
12	NAG	B	507	2	14,14,15	0.68	1 (7%)	17,19,21	0.56	0
12	NAG	D	2006	3	14,14,15	0.44	0	17,19,21	0.54	0
12	NAG	B	505	2	14,14,15	0.28	0	17,19,21	0.38	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
12	NAG	C	707	1	-	1/6/23/26	0/1/1/1
12	NAG	C	709	1	-	2/6/23/26	0/1/1/1
13	MAN	C	708	-	-	0/2/19/22	0/1/1/1
12	NAG	D	2004	3	-	0/6/23/26	0/1/1/1
12	NAG	B	506	2	-	1/6/23/26	0/1/1/1
12	NAG	D	2005	3	-	2/6/23/26	0/1/1/1
12	NAG	B	507	2	-	3/6/23/26	0/1/1/1
12	NAG	D	2006	3	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
12	NAG	B	505	2	-	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
13	C	708	MAN	O5-C1	-2.66	1.39	1.43
13	C	708	MAN	C1-C2	-2.61	1.45	1.52
12	B	507	NAG	C1-C2	2.34	1.55	1.52
13	C	708	MAN	O5-C5	2.28	1.47	1.43

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
13	C	708	MAN	C1-C2-C3	-4.00	103.81	109.64
13	C	708	MAN	O2-C2-C3	-2.99	103.95	110.15
13	C	708	MAN	O5-C1-C2	2.07	115.73	110.79

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
12	D	2005	NAG	O5-C5-C6-O6
12	B	507	NAG	O5-C5-C6-O6
12	D	2006	NAG	O5-C5-C6-O6
12	B	507	NAG	C4-C5-C6-O6
12	D	2005	NAG	C4-C5-C6-O6
12	D	2006	NAG	C4-C5-C6-O6
12	C	709	NAG	O5-C5-C6-O6
12	C	707	NAG	O5-C5-C6-O6
12	C	709	NAG	C4-C5-C6-O6
12	B	506	NAG	C1-C2-N2-C7
12	B	507	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	703	SO4	1	0
13	C	708	MAN	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 ⓘ

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	594/604 (98%)	0.14	10 (1%) 69 65	34, 60, 103, 176	0
1	C	589/604 (97%)	0.25	13 (2%) 62 58	41, 71, 110, 204	0
2	B	448/483 (92%)	0.25	23 (5%) 33 29	25, 61, 148, 228	0
3	D	455/483 (94%)	0.45	30 (6%) 24 21	29, 66, 174, 227	0
4	E	10/13 (76%)	0.42	0 100 100	57, 61, 107, 111	0
4	F	11/13 (84%)	0.38	1 (9%) 15 13	53, 67, 103, 157	0
All	All	2107/2200 (95%)	0.26	77 (3%) 45 40	25, 65, 141, 228	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	D	60	ASN	5.2
3	D	8	LEU	4.3
1	A	453	ILE	4.2
1	A	490	LYS	3.9
2	B	126	ALA	3.7
1	C	453	ILE	3.6
3	D	75	VAL	3.5
2	B	471	PRO	3.4
2	B	10	GLY	3.4
2	B	430	LEU	3.3
3	D	46	ALA	3.2
3	D	430	LEU	3.2
1	C	265	TYR	3.1
2	B	385	HIS	3.1
2	B	13	THR	3.1
3	D	11	ALA	3.0
1	C	506	GLY	3.0
3	D	448	SER	3.0
1	A	594	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
3	D	30	GLU	2.9
2	B	55	LEU	2.8
3	D	215	ILE	2.8
2	B	377	CYS	2.8
3	D	55	LEU	2.8
1	C	329	THR	2.7
3	D	452	ASN	2.7
3	D	471	PRO	2.7
3	D	276	CYS	2.6
3	D	7	ALA	2.6
1	A	502	LEU	2.6
3	D	12	GLU	2.6
2	B	380	GLY	2.6
1	A	31	SER	2.6
2	B	135	ILE	2.5
2	B	8	LEU	2.5
2	B	445	VAL	2.5
1	C	507	ALA	2.5
1	C	219	ASP	2.4
3	D	5	GLY	2.4
3	D	135	ILE	2.4
2	B	50	ALA	2.4
3	D	217	ALA	2.4
2	B	448	SER	2.3
2	B	76	GLY	2.3
1	A	511	ALA	2.3
3	D	52	GLY	2.3
1	A	29	SER	2.3
2	B	18	LEU	2.3
1	A	581	PHE	2.3
3	D	79	LYS	2.2
3	D	31	ASN	2.2
1	A	64	THR	2.2
2	B	461	VAL	2.2
1	C	487	LEU	2.2
2	B	62	VAL	2.2
1	C	116	THR	2.2
3	D	453	GLY	2.2
2	B	129	ASP	2.2
3	D	41	ARG	2.2
2	B	453	GLY	2.1
1	C	103	ASP	2.1

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Mol	Chain	Res	Type	RSRZ
3	D	9	GLY	2.1
3	D	19	LEU	2.1
3	D	431	VAL	2.1
3	D	218	ASN	2.1
1	C	490	LYS	2.1
4	F	239	HIS	2.1
1	C	581	PHE	2.1
2	B	60	ASN	2.1
1	A	32	SER	2.1
1	C	82	LYS	2.0
3	D	13	THR	2.0
2	B	130	ASP	2.0
2	B	208	GLU	2.0
3	D	40	GLU	2.0
1	C	167	ASP	2.0
3	D	216	SER	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.

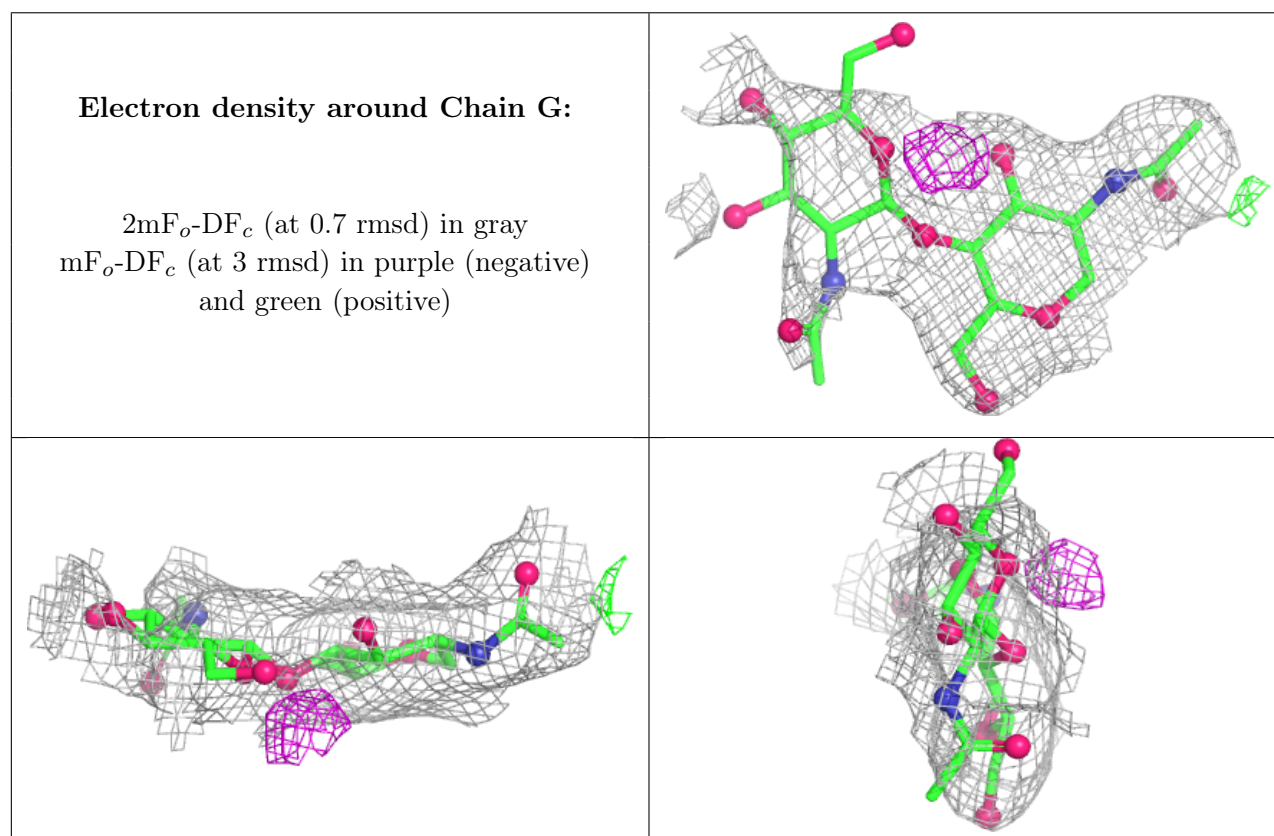
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	L	2	14/15	0.41	0.16	166,175,183,187	0
7	MAN	I	5	11/12	0.41	0.14	153,161,166,166	0
7	MAN	N	5	11/12	0.41	0.18	155,163,173,177	0
7	MAN	N	6	11/12	0.51	0.13	120,148,159,168	0
6	MAN	H	4	11/12	0.54	0.14	129,146,154,158	0
5	NAG	P	2	14/15	0.56	0.15	174,184,193,200	0
6	BMA	H	3	11/12	0.58	0.14	142,152,156,157	0
5	NAG	M	2	14/15	0.60	0.13	76,126,134,141	0
5	NAG	G	2	14/15	0.62	0.16	110,137,151,156	0
5	NAG	K	2	14/15	0.63	0.13	133,157,165,169	0
7	MAN	I	6	11/12	0.63	0.10	100,135,158,161	0
7	MAN	I	4	11/12	0.64	0.12	159,163,172,181	0

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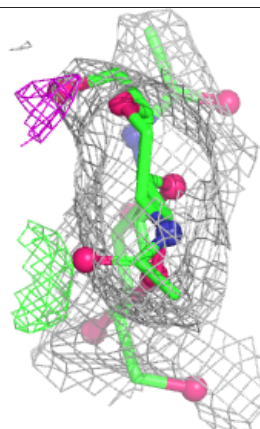
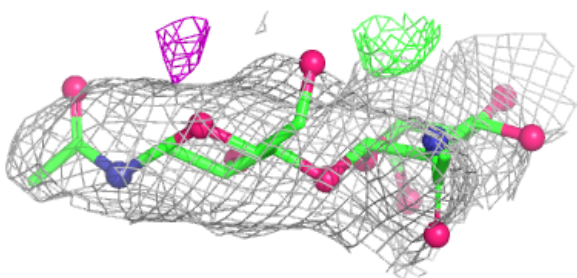
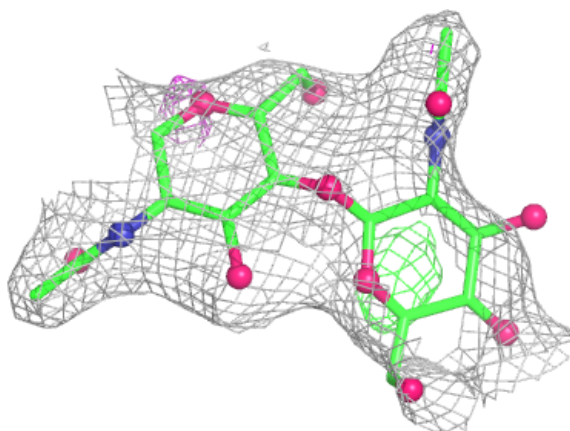
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	NAG	O	2	14/15	0.64	0.15	106,123,142,150	0
8	BMA	O	3	11/12	0.64	0.14	113,119,152,154	0
6	NAG	H	2	14/15	0.68	0.13	76,98,120,144	0
5	NAG	K	1	14/15	0.74	0.14	119,129,160,162	0
7	MAN	N	4	11/12	0.74	0.12	149,154,169,176	0
5	NAG	J	2	14/15	0.76	0.15	94,125,134,134	0
7	BMA	I	3	11/12	0.76	0.10	107,119,144,153	0
7	BMA	N	3	11/12	0.80	0.10	55,96,131,138	0
5	NAG	P	1	14/15	0.80	0.14	91,107,140,155	0
5	NAG	L	1	14/15	0.81	0.11	80,106,126,149	0
5	NAG	M	1	14/15	0.87	0.10	68,90,100,117	0
5	NAG	G	1	14/15	0.91	0.08	45,54,69,72	0
5	NAG	J	1	14/15	0.91	0.10	67,82,97,103	0
8	NAG	O	1	14/15	0.92	0.08	45,78,89,102	0
7	NAG	I	2	14/15	0.93	0.08	47,80,87,98	0
7	NAG	N	2	14/15	0.93	0.07	39,65,89,97	0
7	NAG	I	1	14/15	0.93	0.08	42,49,60,64	0
6	NAG	H	1	14/15	0.95	0.07	46,63,72,79	0
7	NAG	N	1	14/15	0.96	0.08	41,55,65,71	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.



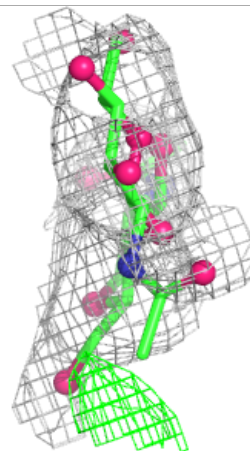
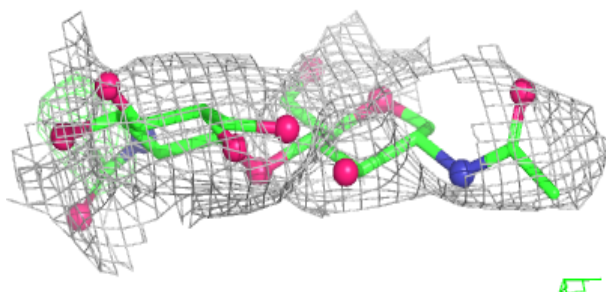
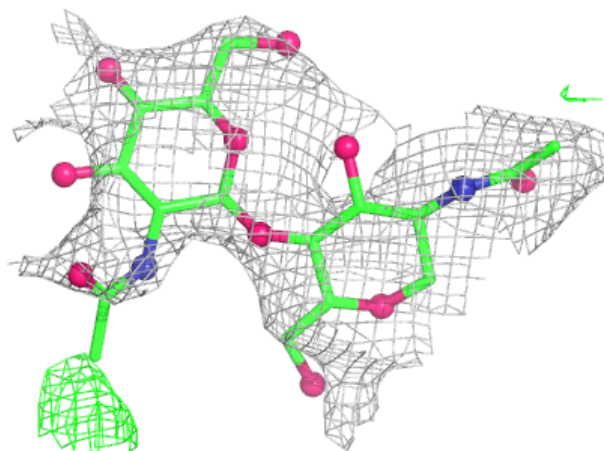
Electron density around Chain J:

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



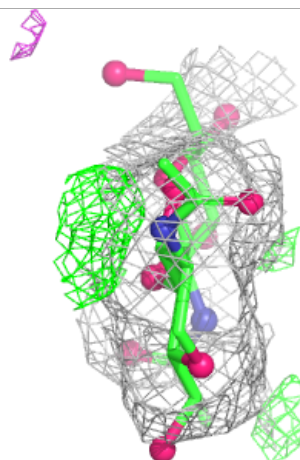
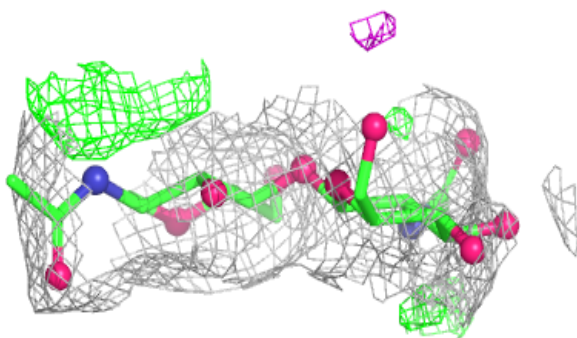
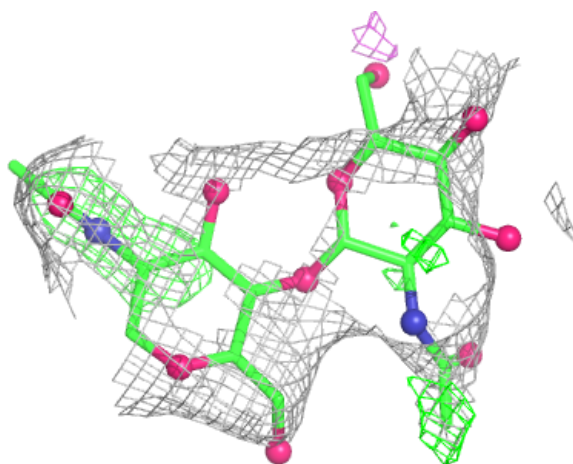
Electron density around Chain K:

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



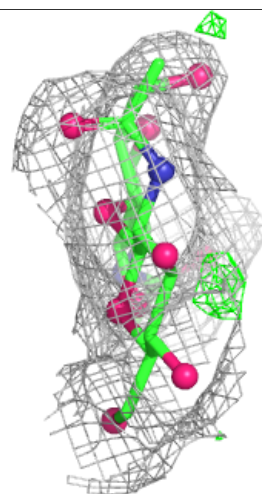
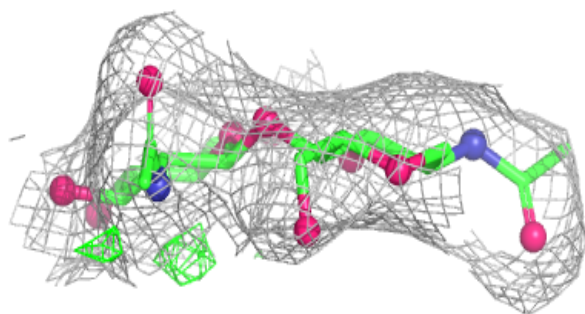
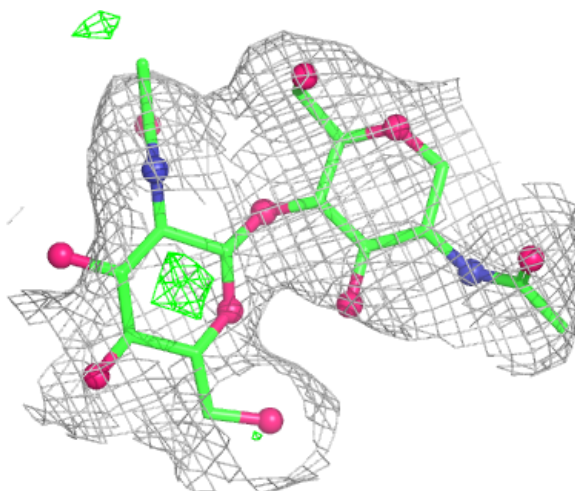
Electron density around Chain L:

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



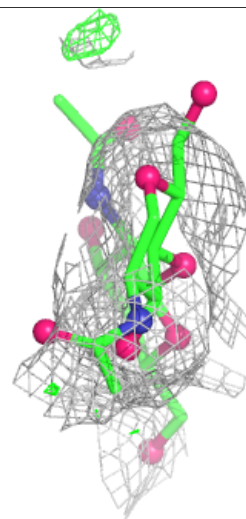
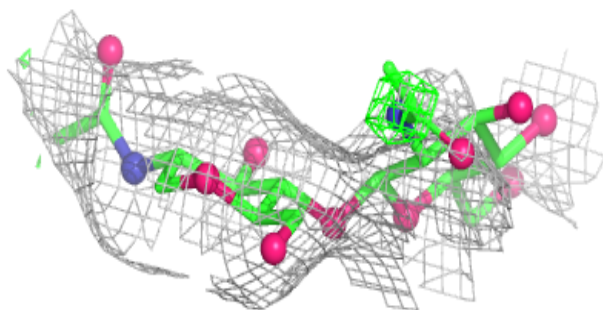
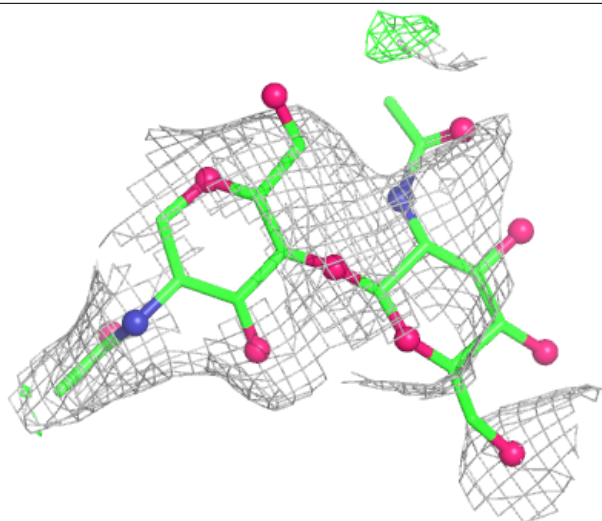
Electron density around Chain M:

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



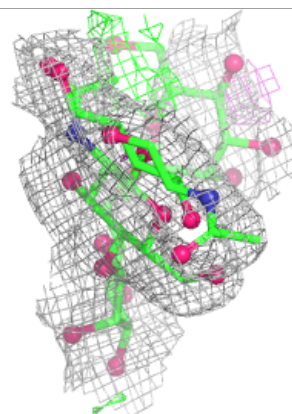
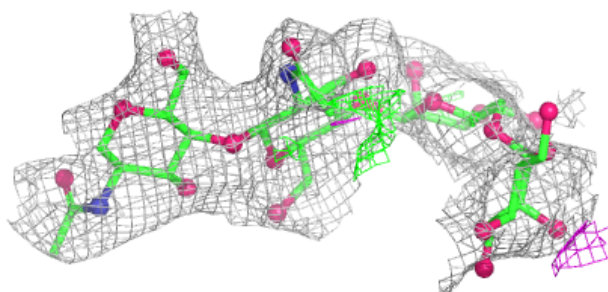
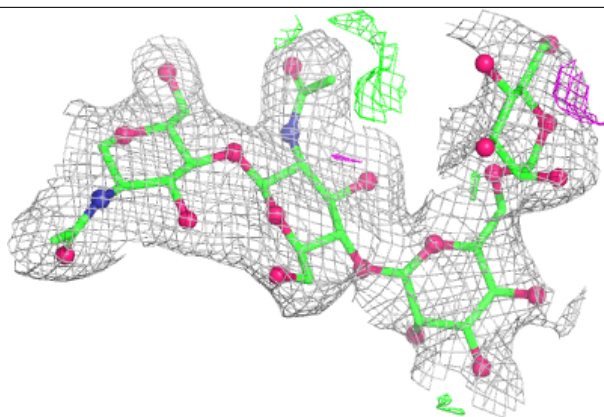
Electron density around Chain P:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

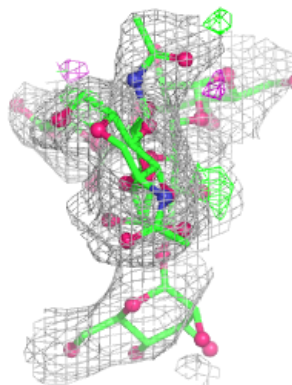
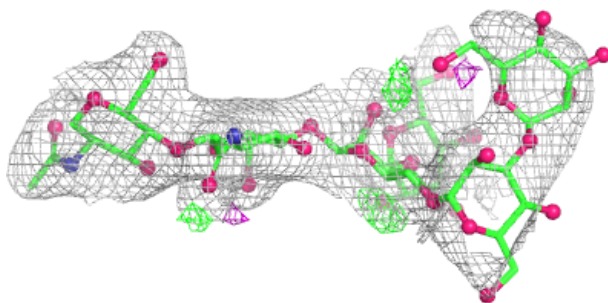
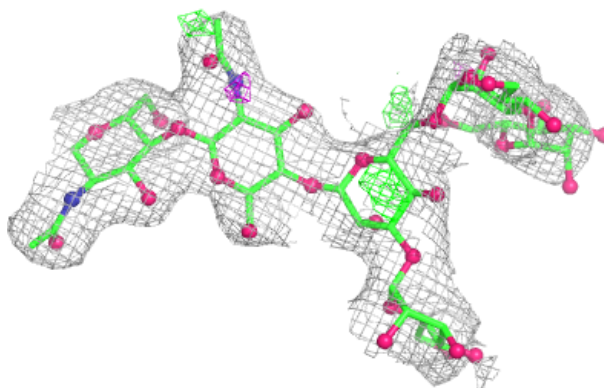


Electron density around Chain H:

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

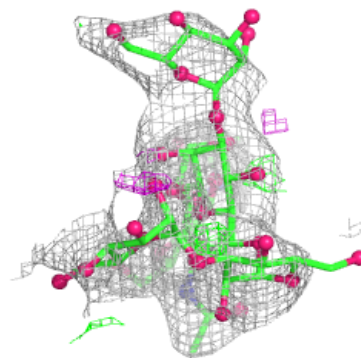
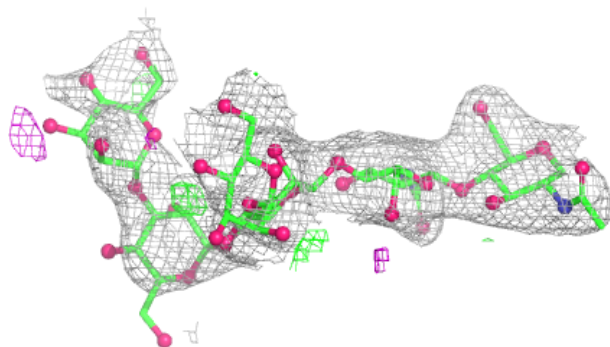
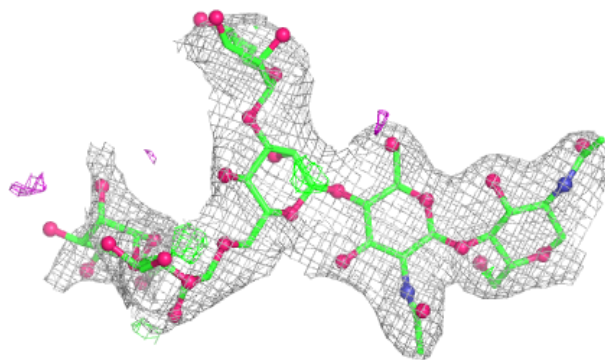
**Electron density around Chain I:**

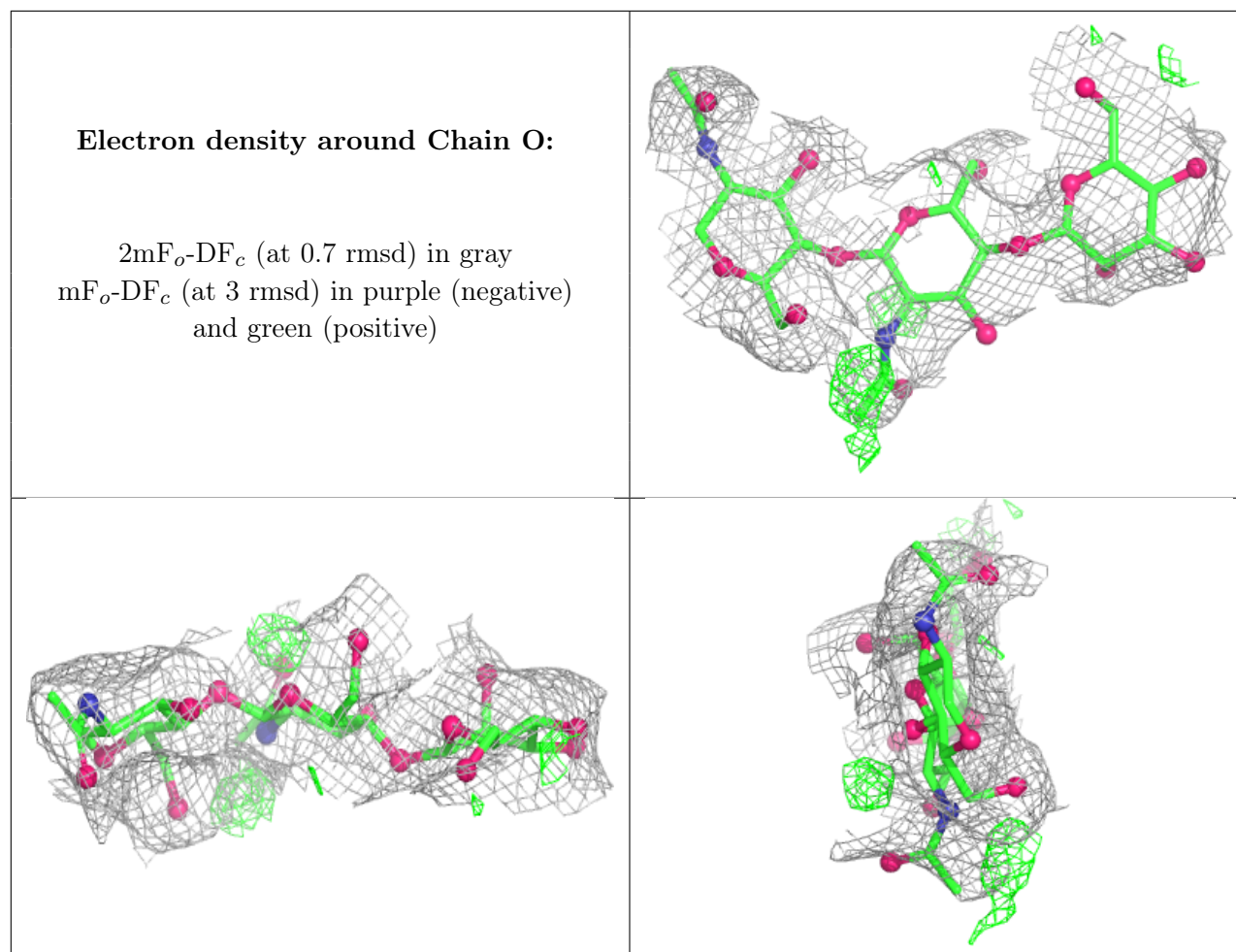
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.4 Ligands [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($\text{\AA}^2$)	Q<0.9
12	NAG	B	507	14/15	0.42	0.17	118,145,153,162	0
12	NAG	D	2004	14/15	0.45	0.16	116,141,156,158	0
12	NAG	C	709	14/15	0.57	0.13	116,136,158,159	0
12	NAG	B	505	14/15	0.60	0.15	109,130,164,168	0
12	NAG	D	2006	14/15	0.60	0.14	100,134,147,151	0
12	NAG	D	2005	14/15	0.63	0.13	101,123,128,129	0
9	SO4	A	702	5/5	0.63	0.22	198,205,206,226	0
9	SO4	B	501	5/5	0.67	0.23	235,237,244,245	0
13	MAN	C	708	11/12	0.67	0.15	98,105,113,115	0
9	SO4	C	702	5/5	0.69	0.11	107,107,109,109	5
12	NAG	B	506	14/15	0.78	0.11	105,111,118,120	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
10	CA	D	2002	1/1	0.84	0.10	140,140,140,140	0
10	CA	B	504	1/1	0.88	0.18	120,120,120,120	0
12	NAG	C	707	14/15	0.90	0.08	55,65,78,85	0
10	CA	C	703	1/1	0.91	0.08	73,73,73,73	0
9	SO4	A	703	5/5	0.92	0.13	82,89,99,101	0
9	SO4	C	701	5/5	0.93	0.07	66,69,86,93	0
9	SO4	A	701	5/5	0.97	0.07	47,47,52,53	5
10	CA	C	705	1/1	0.97	0.04	66,66,66,66	0
10	CA	A	704	1/1	0.98	0.04	67,67,67,67	0
10	CA	B	503	1/1	0.98	0.04	47,47,47,47	0
10	CA	D	2003	1/1	0.98	0.05	49,49,49,49	0
10	CA	C	704	1/1	0.98	0.05	68,68,68,68	0
10	CA	A	705	1/1	0.99	0.03	58,58,58,58	0
11	MG	B	502	1/1	0.99	0.05	61,61,61,61	0
11	MG	D	2001	1/1	0.99	0.04	81,81,81,81	0
10	CA	A	706	1/1	0.99	0.02	53,53,53,53	0
10	CA	C	706	1/1	0.99	0.02	62,62,62,62	0
10	CA	A	707	1/1	0.99	0.04	45,45,45,45	0

6.5 Other polymers

There are no such residues in this entry.