



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 06:28 PM UTC

PDB ID : 3PPS / pdb_00003pps
Title : Crystal structure of an ascomycete fungal laccase from Thielavia arenaria
Authors : Kallio, J.P.; Rouvinen, J.; Hakulinen, N.
Deposited on : 2010-11-25
Resolution : 2.50 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 ⓘ 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)
Xtrriage (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

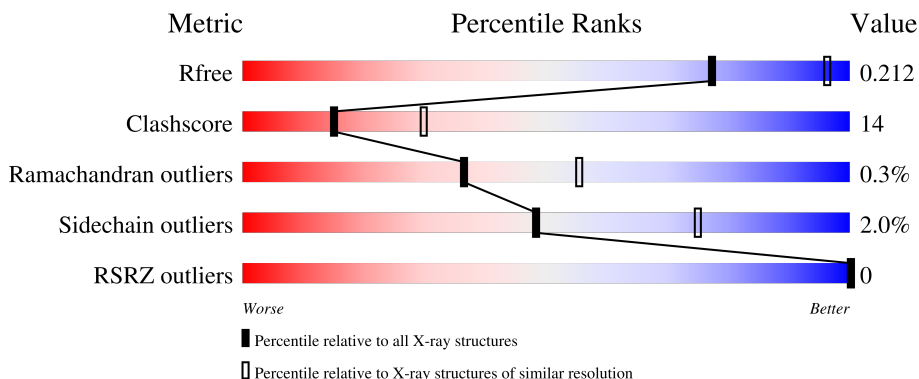
1 Overall quality at a glance i

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.



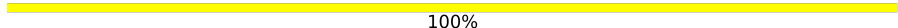
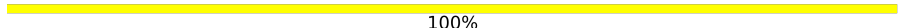
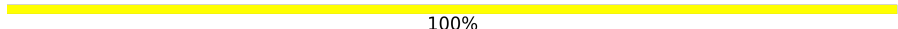
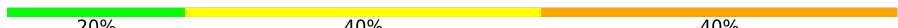
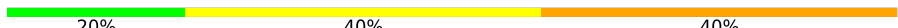

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	
1	B	604	
1	C	604	
1	D	604	
2	E	2	

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Mol	Chain	Length	Quality of chain
2	G	2	 100%
2	H	2	 100%
2	J	2	 50% 50%
2	K	2	 100%
2	M	2	 50% 50%
2	N	2	 100%
2	Q	2	 100%
3	F	6	 33% 50% 17%
4	I	5	 20% 40% 40%
4	P	5	 20% 40% 40%
5	L	3	 33% 67%
5	O	3	 33% 33% 33%

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 18812 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 Laccase.

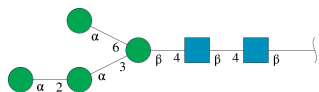
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	564	4406	2793	776	821	16	0	0	0
1	B	564	4406	2793	776	821	16	0	0	0
1	C	564	4406	2793	776	821	16	0	0	0
1	D	564	4406	2793	776	821	16	0	0	0

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



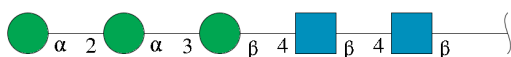
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	E	2	28	16	2	10	0	0	0
2	G	2	28	16	2	10	0	0	0
2	H	2	28	16	2	10	0	0	0
2	J	2	28	16	2	10	0	0	0
2	K	2	28	16	2	10	0	0	0
2	M	2	28	16	2	10	0	0	0
2	N	2	28	16	2	10	0	0	0
2	Q	2	28	16	2	10	0	0	0

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



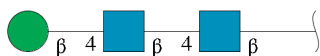
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
3	F	6	72	40	2	30	0	0	0

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-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
			Total	C	N	O			
4	I	5	61	34	2	25	0	0	0
4	P	5	61	34	2	25	0	0	0

- Molecule 5 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.

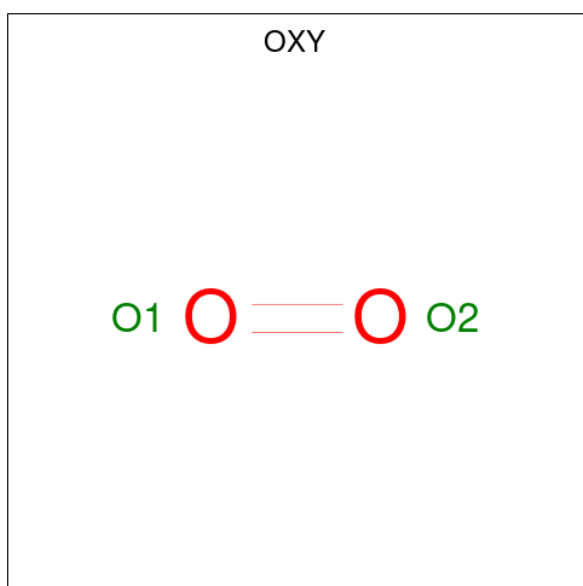


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	L	3	39	22	2	15	0	0	0
5	O	3	39	22	2	15	0	0	0

- Molecule 6 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	4	Total Cu 4 4	0	0
6	B	4	Total Cu 4 4	0	0
6	C	4	Total Cu 4 4	0	0
6	D	4	Total Cu 4 4	0	0

- Molecule 7 is OXYGEN MOLECULE (CCD ID: OXY) (formula: O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total O 2 2	0	0

- Molecule 8 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 9 is water.

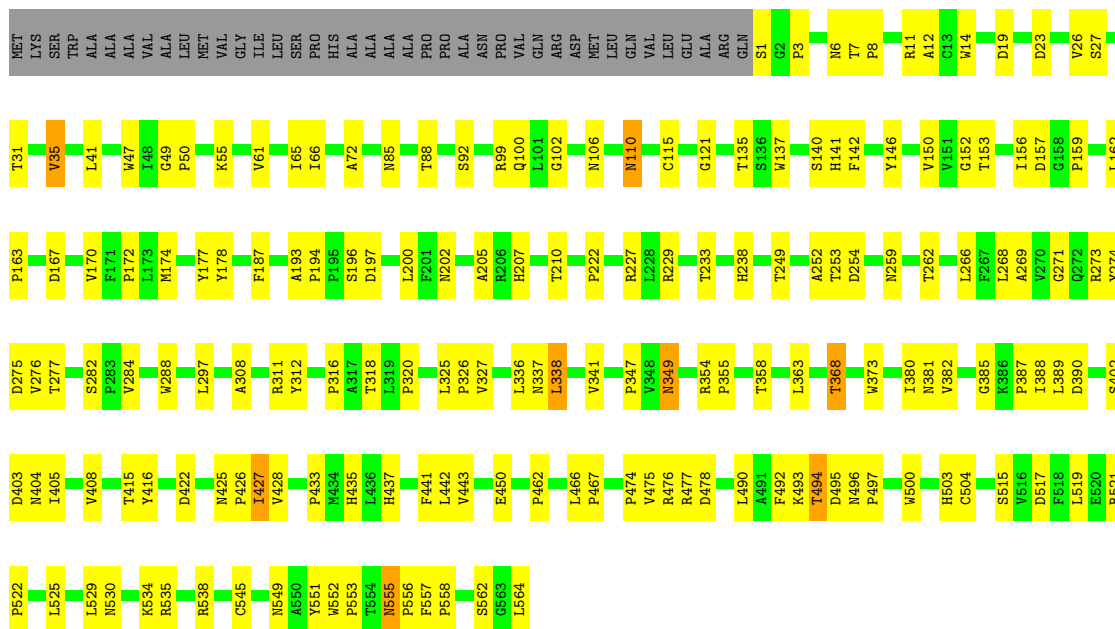
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	159	Total	O	0	0
			159	159		
9	B	138	Total	O	0	0
			138	138		
9	C	124	Total	O	0	0
			124	124		
9	D	141	Total	O	0	0
			141	141		

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.

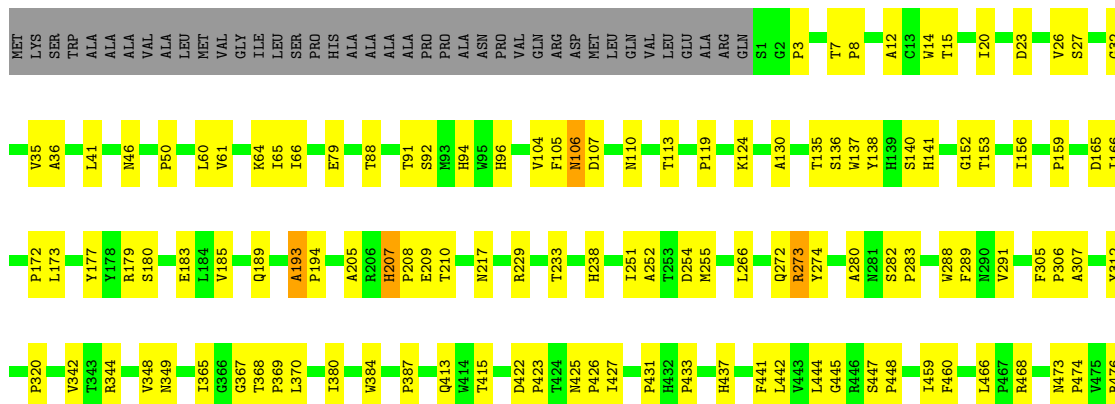
- Molecule 1: Laccase

Chain A: 



- Molecule 1: Laccase

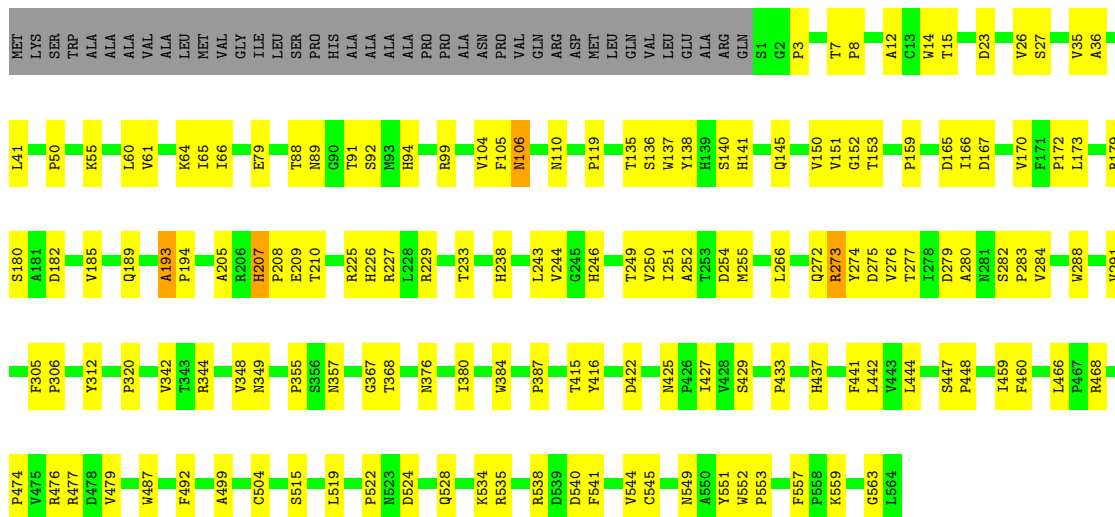
Chain B: 





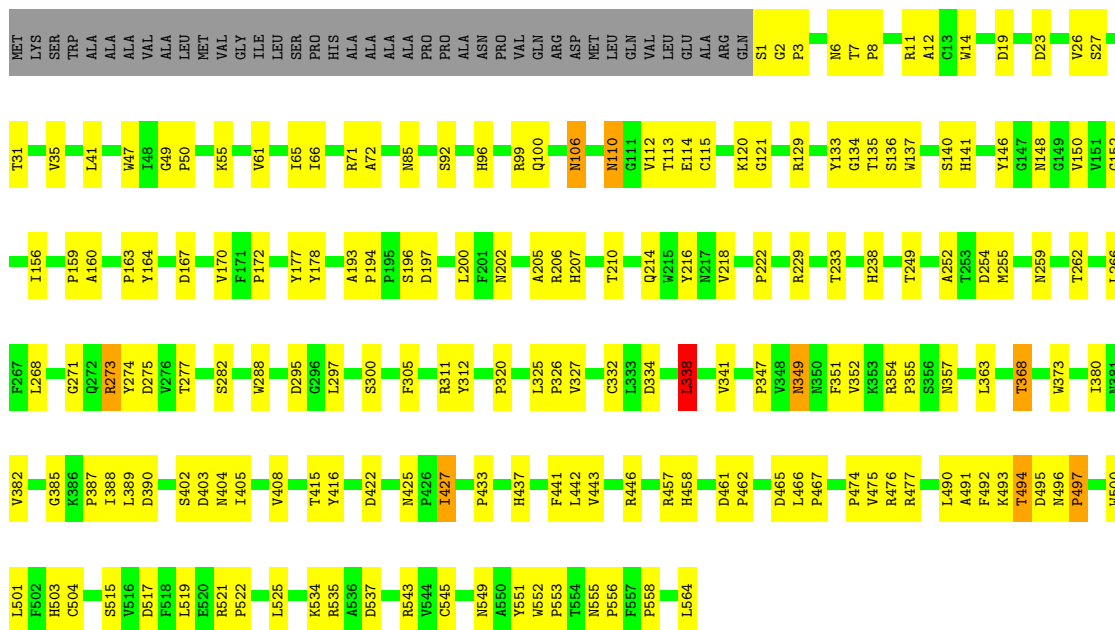
- Molecule 1: Laccase

Chain C: 69% 24% 7%



- Molecule 1: Laccase

Chain D: 63% 28% 7%




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

Chain E: 100%

MAG1
MAG2

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

Chain G:  100%MAG1
MAG2

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

Chain H:  100%MAG1
MAG2


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

Chain J:  50%MAG1
MAG2

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

Chain K:  100%MAG1
MAG2

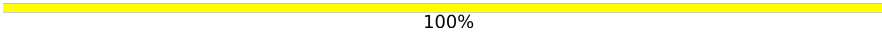
- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  50%MAG1
MAG2

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

Chain N:  100%MAG1
MAG2

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

Chain Q:  100%

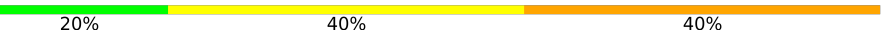
MAG1
MAG2

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

Chain F:  33% 50% 17%

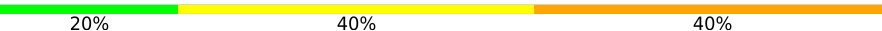
MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

- Molecule 4: alpha-D-mannopyranose-(1-2)-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:  20% 40% 40%

MAG1
MAG2
BMA3
MAN4
MAN5

- Molecule 4: alpha-D-mannopyranose-(1-2)-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 P:  20% 40% 40%

MAG1
MAG2
BMA3
MAN4
MAN5

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

Chain L:  33% 67%

MAG1
MAG2
BMA3

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

Chain O:  33% 33% 33%

MAG1
MAG2
BMA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	61.36Å 178.95Å 118.13Å 90.00° 90.26° 90.00°	Depositor
Resolution (Å)	42.65 – 2.50 42.65 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (42.65-2.50) 98.6 (42.65-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.63 (at 2.51Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.182 , 0.222 0.182 , 0.212	Depositor DCC
R_{free} test set	4402 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	12.6	Xtrriage
Anisotropy	1.837	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 0.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.21$	Xtrriage
Estimated twinning fraction	0.319 for h,-k,-l	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	18812	wwPDB-VP
Average B, all atoms (Å ²)	20.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.22 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.4366e-04.*

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 [i](#)

5.1 Standard geometry [i](#)

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

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	0.46	0/4539	0.82	5/6219 (0.1%)
1	B	0.47	0/4539	0.81	4/6219 (0.1%)
1	C	0.46	0/4539	0.81	5/6219 (0.1%)
1	D	0.44	0/4539	0.82	3/6219 (0.0%)
All	All	0.46	0/18156	0.82	17/24876 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	193	ALA	CA-C-N	6.96	127.55	120.38
1	B	193	ALA	C-N-CA	6.96	127.55	120.38
1	A	338	LEU	CD1-CG-CD2	-6.28	96.99	110.80
1	C	193	ALA	CA-C-N	6.16	126.72	120.38
1	C	193	ALA	C-N-CA	6.16	126.72	120.38
1	A	327	VAL	CA-C-N	6.01	125.97	119.78
1	A	327	VAL	C-N-CA	6.01	125.97	119.78
1	D	338	LEU	CD1-CG-CD2	-5.93	97.75	110.80
1	D	327	VAL	CA-C-N	5.75	125.70	119.78
1	D	327	VAL	C-N-CA	5.75	125.70	119.78
1	C	207	HIS	CA-C-N	5.51	125.86	119.47
1	C	207	HIS	C-N-CA	5.51	125.86	119.47
1	C	150	VAL	N-CA-C	5.35	111.94	106.21
1	B	207	HIS	CA-C-N	5.25	125.57	119.47
1	B	207	HIS	C-N-CA	5.25	125.57	119.47
1	A	555	ASN	CA-C-N	5.02	124.71	119.64
1	A	555	ASN	C-N-CA	5.02	124.71	119.64

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	4406	0	4223	133	0
1	B	4406	0	4223	109	0
1	C	4406	0	4224	107	0
1	D	4406	0	4225	141	0
2	E	28	0	25	4	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
2	J	28	0	25	1	0
2	K	28	0	25	4	0
2	M	28	0	25	1	0
2	N	28	0	25	4	0
2	Q	28	0	25	0	0
3	F	72	0	61	1	0
4	I	61	0	52	2	0
4	P	61	0	52	4	0
5	L	39	0	34	3	0
5	O	39	0	34	1	0
6	A	4	0	0	0	0
6	B	4	0	0	0	0
6	C	4	0	0	0	0
6	D	4	0	0	0	0
7	A	2	0	0	0	0
8	A	42	0	39	1	0
8	B	42	0	39	0	0
8	C	14	0	13	0	0
8	D	14	0	13	1	0
9	A	159	0	0	13	0
9	B	138	0	0	15	0
9	C	124	0	0	14	0
9	D	141	0	0	28	0
All	All	18812	0	17432	489	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 14.

All (489) 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:376:ASN:HD22	2:N:1:NAG:H83	1.35	0.92
1:A:253:THR:HG21	1:A:338:LEU:HD12	1.59	0.83
1:D:207:HIS:HD2	1:D:210:THR:H	1.27	0.82
1:A:207:HIS:HD2	1:A:210:THR:H	1.28	0.81
1:C:104:VAL:HG11	9:C:694:HOH:O	1.79	0.81
1:D:442:LEU:HB2	9:D:663:HOH:O	1.80	0.79
1:D:552:TRP:HB3	1:D:553:PRO:HD3	1.65	0.79
1:A:14:TRP:HB2	1:A:159:PRO:HG3	1.65	0.78
1:D:252:ALA:HB3	1:D:275:ASP:HB2	1.66	0.77
1:D:14:TRP:HB2	1:D:159:PRO:HG3	1.66	0.77
1:B:422:ASP:OD1	1:B:425:ASN:HB2	1.86	0.76
1:C:422:ASP:OD1	1:C:425:ASN:HB2	1.86	0.75
1:A:552:TRP:HB3	1:A:553:PRO:HD3	1.67	0.75
1:D:114:GLU:HB3	9:D:582:HOH:O	1.86	0.74
1:C:167:ASP:OD1	9:C:654:HOH:O	2.05	0.73
1:C:429:SER:HB2	9:C:663:HOH:O	1.89	0.73
1:D:352:VAL:HB	9:D:665:HOH:O	1.87	0.72
1:A:50:PRO:HB2	1:A:205:ALA:HB3	1.72	0.71
1:A:349:ASN:ND2	9:A:635:HOH:O	2.22	0.71
1:C:14:TRP:HB2	1:C:159:PRO:HG3	1.72	0.70
1:D:50:PRO:HB2	1:D:205:ALA:HB3	1.73	0.70
1:D:238:HIS:NE2	1:D:433:PRO:HG3	2.07	0.70
1:C:244:VAL:O	9:C:651:HOH:O	2.08	0.70
1:A:503:HIS:HB3	1:A:515:SER:OG	1.91	0.69
1:B:14:TRP:HB2	1:B:159:PRO:HG3	1.72	0.69
1:A:102:GLY:HA2	9:A:705:HOH:O	1.91	0.69
1:B:487:TRP:O	9:B:631:HOH:O	2.09	0.69
1:A:238:HIS:NE2	1:A:433:PRO:HG3	2.08	0.69
1:C:540:ASP:O	1:C:544:VAL:HG23	1.93	0.68
1:D:462:PRO:HA	1:D:466:LEU:HD13	1.74	0.68
1:B:540:ASP:O	1:B:544:VAL:HG23	1.94	0.67
1:A:252:ALA:HB3	1:A:275:ASP:HB2	1.75	0.67
1:A:426:PRO:HG2	1:B:370:LEU:HD21	1.77	0.67
1:B:14:TRP:CD1	1:B:15:THR:HG23	2.30	0.67
1:C:433:PRO:HD2	1:C:504:CYS:SG	2.35	0.66
1:D:338:LEU:HB2	1:D:475:VAL:HG22	1.78	0.66
1:D:503:HIS:HB3	1:D:515:SER:OG	1.95	0.66
1:C:244:VAL:C	9:C:651:HOH:O	2.37	0.66
1:A:462:PRO:HA	1:A:466:LEU:HD13	1.76	0.66
1:C:376:ASN:ND2	2:N:1:NAG:H83	2.09	0.66
1:B:3:PRO:HG3	1:B:14:TRP:CH2	2.32	0.65
1:B:92:SER:O	1:B:140:SER:HA	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:85:ASN:OD1	1:D:121:GLY:HA2	1.96	0.65
1:A:254:ASP:OD2	1:A:477:ARG:HB2	1.97	0.65
1:C:448:PRO:HD3	1:C:468:ARG:NH2	2.12	0.65
1:B:32:GLY:N	9:B:642:HOH:O	2.27	0.65
1:D:254:ASP:OD2	1:D:477:ARG:HB2	1.96	0.65
1:A:557:PHE:HB2	2:E:1:NAG:O7	1.96	0.64
1:C:3:PRO:HG3	1:C:14:TRP:CH2	2.33	0.64
1:B:106:ASN:N	1:B:106:ASN:HD22	1.95	0.64
1:B:94:HIS:CE1	1:B:437:HIS:CE1	2.86	0.64
1:C:106:ASN:N	1:C:106:ASN:HD22	1.95	0.64
1:C:14:TRP:CD1	1:C:15:THR:HG23	2.33	0.64
1:C:182:ASP:HB2	2:K:2:NAG:H82	1.78	0.64
1:D:446:ARG:NH2	9:D:656:HOH:O	2.31	0.63
1:B:433:PRO:HD2	1:B:504:CYS:SG	2.38	0.63
1:C:442:LEU:HD23	1:C:476:ARG:HB3	1.81	0.63
1:C:477:ARG:NH1	9:C:579:HOH:O	2.32	0.62
1:D:135:THR:HG21	1:D:229:ARG:HB3	1.80	0.62
1:D:214:GLN:CD	8:D:720:NAG:H83	2.23	0.62
1:A:100:GLN:OE1	1:A:100:GLN:HA	1.98	0.62
1:A:408:VAL:HG11	1:A:494:THR:HG21	1.81	0.62
1:A:380:ILE:HG21	1:A:515:SER:HB3	1.81	0.62
1:C:92:SER:O	1:C:140:SER:HA	1.99	0.62
1:D:100:GLN:HA	1:D:100:GLN:OE1	2.00	0.61
1:B:138:TYR:CZ	1:B:152:GLY:HA3	2.34	0.61
1:D:368:THR:HG22	1:D:368:THR:O	2.01	0.61
1:D:519:LEU:HG	1:D:522:PRO:HG3	1.83	0.61
1:D:380:ILE:HG21	1:D:515:SER:HB3	1.83	0.60
1:A:92:SER:OG	1:A:110:ASN:HB3	2.01	0.60
1:C:138:TYR:CZ	1:C:152:GLY:HA3	2.36	0.60
1:A:85:ASN:OD1	1:A:121:GLY:HA2	2.02	0.60
1:D:457:ARG:C	9:D:671:HOH:O	2.43	0.60
1:D:408:VAL:HG11	1:D:494:THR:HG21	1.82	0.60
1:A:519:LEU:HG	1:A:522:PRO:HG3	1.82	0.60
1:B:185:VAL:O	1:B:189:GLN:HG3	2.02	0.60
1:D:92:SER:OG	1:D:110:ASN:HB3	2.02	0.59
1:A:135:THR:HG21	1:A:229:ARG:HB3	1.84	0.59
1:C:185:VAL:O	1:C:189:GLN:HG3	2.02	0.59
1:A:12:ALA:O	1:A:19:ASP:HB2	2.03	0.59
1:D:249:THR:HB	1:D:277:THR:OG1	2.03	0.59
1:D:491:ALA:HB3	9:D:663:HOH:O	2.02	0.59
1:B:425:ASN:OD1	1:B:427:ILE:HG22	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:94:HIS:CE1	1:C:437:HIS:CE1	2.90	0.59
1:D:2:GLY:N	9:D:615:HOH:O	2.35	0.59
1:C:254:ASP:OD2	1:C:477:ARG:HB2	2.02	0.58
1:B:442:LEU:HD23	1:B:476:ARG:HB3	1.86	0.58
1:C:238:HIS:NE2	1:C:433:PRO:HG3	2.18	0.58
1:A:88:THR:HG22	9:A:573:HOH:O	2.04	0.58
1:C:425:ASN:OD1	1:C:427:ILE:HG22	2.04	0.58
1:A:92:SER:O	1:A:140:SER:HA	2.04	0.57
1:B:448:PRO:HD3	1:B:468:ARG:NH2	2.19	0.57
1:B:282:SER:HB3	1:B:283:PRO:HD2	1.86	0.57
1:D:357:ASN:ND2	9:D:699:HOH:O	2.32	0.57
1:B:266:LEU:HD21	1:B:274:TYR:CG	2.39	0.57
1:C:266:LEU:HD21	1:C:274:TYR:CG	2.40	0.57
1:D:415:THR:HG22	1:D:492:PHE:O	2.05	0.57
1:C:415:THR:HG22	1:C:492:PHE:O	2.04	0.57
1:A:249:THR:HB	1:A:277:THR:OG1	2.04	0.57
1:C:448:PRO:HD3	1:C:468:ARG:HH21	1.70	0.57
1:A:253:THR:CG2	1:A:338:LEU:HD12	2.32	0.56
1:A:368:THR:CG2	1:A:368:THR:O	2.52	0.56
1:B:238:HIS:NE2	1:B:433:PRO:HG3	2.19	0.56
1:D:92:SER:O	1:D:140:SER:HA	2.06	0.56
1:D:352:VAL:N	9:D:670:HOH:O	2.24	0.56
1:D:368:THR:O	1:D:368:THR:CG2	2.53	0.56
1:A:170:VAL:HA	1:A:229:ARG:HB2	1.88	0.56
1:C:170:VAL:O	5:L:1:NAG:N2	2.37	0.56
1:D:170:VAL:HA	1:D:229:ARG:HB2	1.88	0.55
1:A:65:ILE:O	1:A:66:ILE:HD13	2.07	0.55
1:A:415:THR:HG22	1:A:492:PHE:O	2.05	0.55
1:A:405:ILE:HD13	1:A:517:ASP:HB3	1.89	0.55
1:B:254:ASP:OD2	1:B:477:ARG:HB2	2.06	0.55
1:A:308:ALA:HA	9:A:617:HOH:O	2.05	0.55
1:B:483:PRO:HG2	9:B:631:HOH:O	2.07	0.55
1:B:415:THR:HG22	1:B:492:PHE:O	2.06	0.55
1:C:7:THR:HB	1:C:8:PRO:HD2	1.88	0.55
1:B:368:THR:HG21	9:B:622:HOH:O	2.06	0.55
1:C:180:SER:OG	2:K:2:NAG:H81	2.07	0.55
1:C:243:LEU:HG	9:C:651:HOH:O	2.07	0.55
1:D:65:ILE:O	1:D:66:ILE:HD13	2.06	0.55
1:A:23:ASP:HB3	1:A:26:VAL:HG22	1.89	0.54
1:C:282:SER:HB3	1:C:283:PRO:HD2	1.87	0.54
1:A:338:LEU:HB2	1:A:475:VAL:HG22	1.87	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:71:ARG:HH12	4:P:2:NAG:H4	1.71	0.54
1:C:65:ILE:O	1:C:66:ILE:HD13	2.08	0.54
1:A:337:ASN:O	1:A:338:LEU:HD23	2.08	0.54
1:D:23:ASP:HB3	1:D:26:VAL:HG22	1.90	0.54
1:D:363:LEU:HB2	1:D:373:TRP:CZ3	2.43	0.54
1:B:23:ASP:HB3	1:B:26:VAL:HG22	1.90	0.54
1:D:200:LEU:HD23	1:D:205:ALA:HB2	1.90	0.54
1:C:524:ASP:O	1:C:528:GLN:HG3	2.08	0.54
1:D:405:ILE:HD13	1:D:517:ASP:HB3	1.90	0.54
1:D:491:ALA:N	9:D:663:HOH:O	2.40	0.54
1:A:1:SER:HB2	1:A:31:THR:HG22	1.90	0.53
1:C:105:PHE:C	1:C:106:ASN:HD22	2.15	0.53
1:A:282:SER:HB2	1:A:312:TYR:OH	2.08	0.53
1:B:105:PHE:C	1:B:106:ASN:HD22	2.14	0.53
1:B:207:HIS:CD2	1:B:209:GLU:H	2.25	0.53
1:C:291:VAL:HB	1:C:306:PRO:HB2	1.91	0.53
1:A:450:GLU:HB2	9:A:670:HOH:O	2.06	0.53
1:D:461:ASP:C	9:D:650:HOH:O	2.51	0.53
1:A:200:LEU:HD23	1:A:205:ALA:HB2	1.91	0.53
1:A:368:THR:O	1:A:368:THR:HG23	2.09	0.53
1:C:23:ASP:HB3	1:C:26:VAL:CG2	2.39	0.53
1:C:246:HIS:N	9:C:651:HOH:O	2.41	0.53
1:D:288:TRP:CE2	1:D:320:PRO:HB2	2.44	0.53
1:C:23:ASP:HB3	1:C:26:VAL:HG22	1.91	0.53
1:D:12:ALA:O	1:D:19:ASP:HB2	2.08	0.53
1:D:282:SER:HB2	1:D:312:TYR:OH	2.08	0.53
1:A:427:ILE:HD13	1:B:365:ILE:HD11	1.91	0.53
1:B:7:THR:HB	1:B:8:PRO:HD2	1.90	0.53
1:B:485:GLY:O	9:B:661:HOH:O	2.19	0.52
1:D:461:ASP:N	9:D:650:HOH:O	2.42	0.52
1:A:194:PRO:HD2	9:A:708:HOH:O	2.09	0.52
1:B:535:ARG:NH2	9:B:569:HOH:O	2.41	0.52
1:C:135:THR:HG21	1:C:229:ARG:HB3	1.92	0.52
1:D:129:ARG:HD3	9:D:661:HOH:O	2.07	0.52
1:B:291:VAL:HB	1:B:306:PRO:HB2	1.92	0.52
1:B:64:LYS:HB2	9:B:688:HOH:O	2.09	0.52
1:C:207:HIS:CD2	1:C:209:GLU:H	2.27	0.52
1:A:363:LEU:HB2	1:A:373:TRP:CZ3	2.44	0.52
5:L:2:NAG:H83	5:L:2:NAG:O3	2.09	0.52
1:D:216:TYR:HE2	4:P:1:NAG:H82	1.74	0.52
1:B:448:PRO:HD3	1:B:468:ARG:HH21	1.75	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:305:PHE:HA	9:D:703:HOH:O	2.10	0.52
1:A:385:GLY:O	1:A:387:PRO:HD3	2.11	0.51
1:D:146:TYR:HB2	1:D:150:VAL:O	2.09	0.51
1:B:519:LEU:HG	1:B:522:PRO:HG3	1.92	0.51
1:D:120:LYS:HA	9:D:595:HOH:O	2.09	0.51
1:A:557:PHE:CD2	2:E:2:NAG:H62	2.45	0.51
1:A:61:VAL:N	9:A:577:HOH:O	2.19	0.51
1:D:1:SER:HB2	1:D:31:THR:HG22	1.92	0.51
1:D:385:GLY:O	1:D:387:PRO:HD3	2.11	0.51
1:A:3:PRO:HG3	1:A:14:TRP:CH2	2.46	0.51
1:B:23:ASP:HB3	1:B:26:VAL:CG2	2.41	0.51
1:B:524:ASP:O	1:B:528:GLN:HG3	2.10	0.51
1:C:119:PRO:HG2	1:C:551:TYR:CZ	2.46	0.51
1:D:47:TRP:CZ3	1:D:49:GLY:HA2	2.46	0.51
1:C:384:TRP:CZ3	1:C:559:LYS:HD2	2.45	0.51
1:D:6:ASN:HD22	1:D:14:TRP:HE3	1.59	0.51
1:D:26:VAL:HG23	1:D:27:SER:N	2.26	0.51
1:D:325:LEU:HB3	1:D:326:PRO:CD	2.41	0.51
1:C:41:LEU:HA	1:C:60:LEU:O	2.11	0.51
1:D:354:ARG:NH1	9:D:665:HOH:O	2.43	0.50
1:A:47:TRP:CZ3	1:A:49:GLY:HA2	2.46	0.50
1:B:153:THR:HG21	1:B:172:PRO:HG3	1.92	0.50
1:D:218:VAL:HG22	4:P:1:NAG:H81	1.93	0.50
1:D:433:PRO:HD2	1:D:504:CYS:SG	2.52	0.50
1:C:380:ILE:HG21	1:C:515:SER:HB3	1.94	0.50
1:B:65:ILE:O	1:B:66:ILE:HD13	2.12	0.50
1:C:172:PRO:O	1:C:173:LEU:HD23	2.11	0.50
1:D:3:PRO:HG3	1:D:14:TRP:CH2	2.47	0.50
1:A:41:LEU:CD2	1:A:61:VAL:HG22	2.42	0.50
1:D:437:HIS:O	1:D:500:TRP:HB3	2.11	0.50
1:B:384:TRP:CZ2	1:B:563:GLY:HA2	2.46	0.50
1:A:207:HIS:CD2	1:A:210:THR:H	2.19	0.50
1:B:384:TRP:CZ3	1:B:559:LYS:HD2	2.46	0.50
1:C:104:VAL:CG1	9:C:694:HOH:O	2.50	0.50
1:B:305:PHE:CZ	2:J:2:NAG:H61	2.47	0.50
1:C:519:LEU:HG	1:C:522:PRO:HG3	1.93	0.50
1:A:288:TRP:CE2	1:A:320:PRO:HB2	2.47	0.49
1:D:496:ASN:O	1:D:497:PRO:C	2.54	0.49
1:B:272:GLN:HG3	1:B:479:VAL:CG1	2.43	0.49
1:C:384:TRP:CE3	1:C:559:LYS:HD2	2.47	0.49
1:B:41:LEU:HD23	1:B:61:VAL:HG22	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:LEU:HD23	1:C:61:VAL:HG22	1.94	0.49
1:A:422:ASP:OD1	1:A:425:ASN:HB2	2.13	0.49
1:B:172:PRO:O	1:B:173:LEU:HD23	2.12	0.49
1:B:384:TRP:CE3	1:B:559:LYS:HD2	2.47	0.49
1:D:177:TYR:CE2	1:D:196:SER:HA	2.48	0.49
1:B:413:GLN:HB2	9:B:580:HOH:O	2.12	0.49
1:A:354:ARG:HB2	1:A:355:PRO:HD2	1.95	0.49
1:B:280:ALA:HB1	1:B:312:TYR:CE2	2.48	0.49
1:C:384:TRP:CZ2	1:C:563:GLY:HA2	2.48	0.49
1:D:347:PRO:HD2	1:D:416:TYR:OH	2.13	0.49
1:C:50:PRO:HB2	1:C:205:ALA:HB3	1.94	0.48
1:A:26:VAL:HG23	1:A:27:SER:N	2.27	0.48
1:B:96:HIS:O	1:B:136:SER:OG	2.26	0.48
1:B:380:ILE:HG21	1:B:515:SER:HB3	1.96	0.48
1:D:216:TYR:CE2	4:P:1:NAG:H82	2.48	0.48
1:C:12:ALA:O	1:C:159:PRO:HB3	2.12	0.48
1:B:46:ASN:ND2	9:B:685:HOH:O	2.36	0.48
1:D:338:LEU:HB2	1:D:475:VAL:CG2	2.43	0.48
1:D:493:LYS:O	1:D:495:ASP:N	2.44	0.48
1:B:119:PRO:HG2	1:B:551:TYR:CZ	2.49	0.48
1:B:135:THR:HG21	1:B:229:ARG:HB3	1.96	0.48
1:C:280:ALA:HB1	1:C:312:TYR:CE2	2.49	0.48
1:D:354:ARG:HB2	1:D:355:PRO:HD2	1.96	0.48
1:D:382:VAL:CG1	1:D:564:LEU:HD11	2.43	0.48
1:A:6:ASN:HD22	1:A:14:TRP:HE3	1.62	0.48
1:A:146:TYR:HB2	1:A:150:VAL:O	2.14	0.48
1:B:124:LYS:NZ	9:B:626:HOH:O	2.46	0.48
1:C:557:PHE:CD2	2:K:2:NAG:H62	2.49	0.48
1:D:172:PRO:HD2	1:D:202:ASN:HA	1.95	0.48
1:C:26:VAL:HG23	1:C:27:SER:N	2.30	0.47
1:D:255:MET:HB2	9:D:566:HOH:O	2.13	0.47
1:A:496:ASN:O	1:A:497:PRO:C	2.57	0.47
1:A:433:PRO:HD2	1:A:504:CYS:SG	2.54	0.47
1:C:368:THR:HG23	1:C:368:THR:O	2.14	0.47
1:D:382:VAL:HG13	1:D:564:LEU:HD11	1.96	0.47
1:D:442:LEU:HD23	1:D:476:ARG:HB3	1.95	0.47
1:A:466:LEU:N	1:A:467:PRO:CD	2.77	0.47
1:C:422:ASP:CG	1:C:425:ASN:HB2	2.40	0.47
1:D:266:LEU:HD21	1:D:274:TYR:CG	2.50	0.47
1:A:55:LYS:HE3	1:A:55:LYS:HB3	1.60	0.47
1:A:325:LEU:HB3	1:A:326:PRO:CD	2.44	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:530:ASN:HA	9:A:653:HOH:O	2.13	0.47
1:B:538:ARG:O	1:B:541:PHE:HB3	2.15	0.47
1:D:334:ASP:HB2	9:D:664:HOH:O	2.15	0.47
1:A:227:ARG:NH2	9:A:678:HOH:O	2.46	0.47
1:A:408:VAL:CG1	1:A:494:THR:HG21	2.44	0.47
1:B:217:ASN:HB2	9:B:612:HOH:O	2.14	0.47
1:C:251:ILE:O	1:C:252:ALA:HB2	2.15	0.47
1:D:249:THR:HA	1:D:262:THR:HA	1.97	0.47
1:D:55:LYS:HB3	1:D:55:LYS:HE3	1.66	0.47
1:D:252:ALA:HB1	9:D:596:HOH:O	2.15	0.47
1:B:534:LYS:O	1:B:535:ARG:C	2.58	0.47
1:C:266:LEU:HD21	1:C:274:TYR:CD1	2.50	0.47
1:C:441:PHE:O	1:C:476:ARG:HB2	2.15	0.47
1:D:26:VAL:HG23	1:D:27:SER:H	1.79	0.47
1:D:26:VAL:HA	9:D:625:HOH:O	2.15	0.47
1:D:425:ASN:HA	9:D:589:HOH:O	2.13	0.47
1:C:88:THR:HG21	1:C:557:PHE:HE1	1.80	0.47
1:A:177:TYR:CE2	1:A:196:SER:HA	2.50	0.46
1:C:99:ARG:HG3	9:C:676:HOH:O	2.14	0.46
1:C:153:THR:HG21	1:C:172:PRO:HG3	1.96	0.46
1:C:545:CYS:O	1:C:549:ASN:ND2	2.44	0.46
1:A:557:PHE:HA	1:A:558:PRO:HD3	1.80	0.46
1:C:165:ASP:HB2	1:C:166:ILE:HD12	1.97	0.46
1:C:474:PRO:HA	9:C:598:HOH:O	2.15	0.46
1:B:445:GLY:HA2	9:B:695:HOH:O	2.14	0.46
1:A:249:THR:HA	1:A:262:THR:HA	1.96	0.46
1:B:88:THR:HG21	1:B:557:PHE:HE1	1.80	0.46
1:D:458:HIS:N	9:D:671:HOH:O	2.48	0.46
1:D:207:HIS:CD2	1:D:210:THR:H	2.18	0.46
1:D:388:ILE:HD12	1:D:405:ILE:HD11	1.98	0.46
1:A:380:ILE:O	1:A:562:SER:HB2	2.16	0.46
1:B:431:PRO:HB3	9:B:629:HOH:O	2.16	0.46
1:D:534:LYS:O	1:D:535:ARG:C	2.59	0.46
1:A:26:VAL:HG23	1:A:27:SER:H	1.80	0.46
1:A:442:LEU:HD23	1:A:476:ARG:HB3	1.97	0.46
1:B:41:LEU:HA	1:B:60:LEU:O	2.15	0.46
1:B:130:ALA:HB1	1:B:156:ILE:HD13	1.98	0.46
1:C:272:GLN:HG3	1:C:479:VAL:CG1	2.46	0.46
2:N:1:NAG:H82	2:N:1:NAG:H2	1.74	0.46
1:D:99:ARG:O	1:D:100:GLN:C	2.59	0.46
1:D:193:ALA:HA	1:D:194:PRO:HD3	1.80	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:PHE:HZ	2:E:1:NAG:H83	1.81	0.46
1:A:428:VAL:CG1	1:B:427:ILE:HD11	2.46	0.46
1:A:437:HIS:O	1:A:500:TRP:HB3	2.16	0.46
1:D:295:ASP:HB3	1:D:332:CYS:SG	2.56	0.46
1:A:493:LYS:O	1:A:495:ASP:N	2.48	0.45
1:C:207:HIS:HD2	1:C:210:THR:H	1.64	0.45
1:A:521:ARG:N	1:A:522:PRO:HD3	2.31	0.45
1:B:36:ALA:HA	1:B:79:GLU:O	2.15	0.45
1:C:255:MET:HG2	1:C:273:ARG:HH21	1.81	0.45
1:D:497:PRO:O	1:D:525:LEU:HD13	2.17	0.45
1:A:349:ASN:OD1	1:A:349:ASN:N	2.48	0.45
1:A:388:ILE:HD12	1:A:405:ILE:HD11	1.98	0.45
1:B:441:PHE:O	1:B:476:ARG:HB2	2.16	0.45
1:C:551:TYR:O	1:C:552:TRP:C	2.59	0.45
1:A:266:LEU:HD21	1:A:274:TYR:CG	2.51	0.45
1:A:358:THR:O	8:A:760:NAG:H82	2.16	0.45
1:A:425:ASN:OD1	1:A:427:ILE:HG23	2.17	0.45
1:A:534:LYS:O	1:A:535:ARG:C	2.59	0.45
1:A:497:PRO:O	1:A:525:LEU:HD13	2.17	0.45
1:B:207:HIS:HD2	1:B:210:THR:H	1.65	0.45
1:C:538:ARG:O	1:C:541:PHE:HB3	2.17	0.45
1:A:388:ILE:CD1	1:A:405:ILE:HD11	2.47	0.45
1:B:255:MET:HG2	1:B:273:ARG:HH21	1.82	0.45
1:B:367:GLY:O	1:B:368:THR:C	2.59	0.45
1:A:7:THR:HB	1:A:8:PRO:HD2	1.99	0.45
1:C:36:ALA:HA	1:C:79:GLU:O	2.16	0.45
1:D:137:TRP:HB2	1:D:152:GLY:O	2.16	0.45
1:D:466:LEU:N	1:D:467:PRO:CD	2.80	0.45
1:A:137:TRP:CZ3	1:A:233:THR:HB	2.52	0.45
1:B:12:ALA:O	1:B:159:PRO:HB3	2.17	0.45
5:L:2:NAG:O3	5:L:3:BMA:O5	2.30	0.45
1:C:534:LYS:O	1:C:535:ARG:C	2.60	0.44
1:D:349:ASN:N	1:D:349:ASN:OD1	2.50	0.44
1:D:408:VAL:CG1	1:D:494:THR:HG21	2.46	0.44
1:A:349:ASN:CG	9:A:682:HOH:O	2.61	0.44
1:B:266:LEU:HD21	1:B:274:TYR:CD1	2.52	0.44
1:D:41:LEU:CD2	1:D:61:VAL:HG22	2.48	0.44
1:D:300:SER:HB2	9:D:668:HOH:O	2.17	0.44
1:A:99:ARG:O	1:A:100:GLN:C	2.60	0.44
1:A:207:HIS:HD2	1:A:210:THR:N	2.07	0.44
1:D:135:THR:O	1:D:273:ARG:HG3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:387:PRO:O	1:D:390:ASP:HB2	2.18	0.44
1:D:23:ASP:HB3	1:D:26:VAL:CG2	2.48	0.44
1:D:137:TRP:CZ3	1:D:233:THR:HB	2.53	0.44
1:D:558:PRO:O	5:O:1:NAG:H83	2.17	0.44
1:A:380:ILE:HA	1:A:402:SER:O	2.18	0.44
1:B:166:ILE:HG23	4:I:4:MAN:H61	2.00	0.44
1:B:422:ASP:CG	1:B:425:ASN:HB2	2.43	0.44
1:A:555:ASN:HA	1:A:556:PRO:HD3	1.75	0.44
1:D:380:ILE:HA	1:D:402:SER:O	2.17	0.44
1:A:172:PRO:HD2	1:A:202:ASN:HA	2.00	0.43
1:C:499:ALA:HB3	9:C:694:HOH:O	2.18	0.43
1:D:422:ASP:OD1	1:D:425:ASN:HB2	2.18	0.43
1:D:442:LEU:HB3	1:D:474:PRO:HG2	1.99	0.43
1:D:545:CYS:O	1:D:549:ASN:ND2	2.51	0.43
1:A:443:VAL:HA	1:A:490:LEU:HD23	2.00	0.43
1:B:65:ILE:HG23	1:B:66:ILE:HG12	2.00	0.43
1:B:459:ILE:O	1:B:460:PHE:C	2.62	0.43
1:A:178:TYR:CE1	1:A:197:ASP:HB3	2.53	0.43
1:A:538:ARG:NE	9:A:653:HOH:O	2.31	0.43
1:C:284:VAL:HG13	1:C:312:TYR:HB3	2.00	0.43
1:A:382:VAL:CG1	1:A:564:LEU:HD11	2.49	0.43
1:B:368:THR:HG23	1:B:368:THR:O	2.18	0.43
1:C:227:ARG:HB3	9:C:654:HOH:O	2.18	0.43
1:C:459:ILE:O	1:C:460:PHE:C	2.62	0.43
1:D:178:TYR:CE1	1:D:197:ASP:HB3	2.54	0.43
1:D:297:LEU:HD23	1:D:297:LEU:HA	1.82	0.43
1:D:388:ILE:CD1	1:D:405:ILE:HD11	2.48	0.43
1:D:465:ASP:HB2	9:D:650:HOH:O	2.18	0.43
1:A:441:PHE:O	1:A:476:ARG:HA	2.18	0.43
1:B:26:VAL:HG23	1:B:27:SER:N	2.34	0.43
1:C:387:PRO:HB2	1:C:541:PHE:HZ	1.83	0.43
1:D:351:PHE:HA	9:D:670:HOH:O	2.18	0.43
1:D:425:ASN:OD1	1:D:427:ILE:HG23	2.19	0.43
1:D:521:ARG:N	1:D:522:PRO:HD3	2.34	0.43
1:B:50:PRO:HB2	1:B:205:ALA:HB3	1.99	0.43
1:B:348:VAL:CG1	1:B:444:LEU:HD22	2.49	0.43
1:D:403:ASP:O	1:D:404:ASN:C	2.61	0.43
1:C:225:ARG:HB2	1:C:279:ASP:OD2	2.19	0.43
1:A:35:VAL:HG21	1:A:72:ALA:HB2	2.00	0.42
1:A:115:CYS:CB	1:A:545:CYS:HA	2.49	0.42
1:A:193:ALA:HA	1:A:194:PRO:HD3	1.81	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:GLN:HB3	1:B:369:PRO:HB3	2.00	0.42
1:B:551:TYR:O	1:B:552:TRP:C	2.62	0.42
1:C:55:LYS:HE3	1:C:55:LYS:HB3	1.75	0.42
1:D:113:THR:HA	1:D:389:LEU:HG	2.01	0.42
1:B:207:HIS:HA	1:B:208:PRO:HD3	1.91	0.42
1:B:423:PRO:HA	9:B:661:HOH:O	2.19	0.42
1:C:207:HIS:HA	1:C:208:PRO:HD3	1.90	0.42
1:C:355:PRO:HB2	2:N:2:NAG:O6	2.20	0.42
1:A:23:ASP:HB3	1:A:26:VAL:CG2	2.49	0.42
1:A:137:TRP:HB2	1:A:152:GLY:O	2.20	0.42
1:C:226:HIS:O	1:C:277:THR:HA	2.19	0.42
1:D:35:VAL:HG21	1:D:72:ALA:HB2	2.01	0.42
1:C:60:LEU:HD22	1:C:64:LYS:O	2.20	0.42
1:C:552:TRP:HB3	1:C:553:PRO:HD3	2.02	0.42
1:A:545:CYS:O	1:A:549:ASN:ND2	2.53	0.42
1:B:41:LEU:CD2	1:B:61:VAL:HG22	2.49	0.42
1:B:370:LEU:N	9:B:624:HOH:O	2.52	0.42
1:D:7:THR:HB	1:D:8:PRO:HD2	2.02	0.42
1:D:96:HIS:O	1:D:136:SER:OG	2.30	0.42
1:D:382:VAL:HG11	1:D:564:LEU:HD21	2.02	0.42
1:A:142:PHE:CZ	2:E:1:NAG:H83	2.54	0.42
1:A:427:ILE:HD11	1:B:427:ILE:HG13	2.00	0.42
1:B:137:TRP:CE3	1:B:233:THR:HB	2.55	0.42
1:D:11:ARG:HG2	1:D:163:PRO:HA	2.01	0.42
1:D:207:HIS:HD2	1:D:210:THR:N	2.06	0.42
1:B:165:ASP:HB2	1:B:166:ILE:HD12	2.00	0.42
1:B:193:ALA:HA	1:B:194:PRO:HD3	1.66	0.42
1:B:288:TRP:CE2	1:B:320:PRO:HB2	2.55	0.42
1:D:115:CYS:CB	1:D:545:CYS:HA	2.49	0.42
1:A:7:THR:O	1:A:11:ARG:HD3	2.20	0.42
1:A:137:TRP:CZ2	1:A:271:GLY:HA2	2.54	0.42
1:A:382:VAL:HG13	1:A:564:LEU:HD11	2.01	0.42
1:A:403:ASP:O	1:A:404:ASN:C	2.63	0.42
1:D:441:PHE:O	1:D:476:ARG:HA	2.20	0.42
1:D:555:ASN:HA	1:D:556:PRO:HD3	1.71	0.42
1:A:387:PRO:O	1:A:390:ASP:HB2	2.20	0.42
1:C:180:SER:CB	2:K:2:NAG:H81	2.50	0.42
1:C:288:TRP:CE2	1:C:320:PRO:HB2	2.55	0.42
1:B:60:LEU:HD22	1:B:64:LYS:O	2.20	0.41
1:B:251:ILE:O	1:B:252:ALA:HB2	2.20	0.41
1:C:136:SER:HA	1:C:273:ARG:HD2	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:499:ALA:N	9:C:694:HOH:O	2.28	0.41
1:B:425:ASN:HA	1:B:426:PRO:HD3	1.88	0.41
1:C:447:SER:HB3	1:C:487:TRP:CD1	2.55	0.41
1:D:543:ARG:NH2	9:D:681:HOH:O	2.37	0.41
1:B:289:PHE:O	1:B:307:ALA:HA	2.20	0.41
1:C:305:PHE:CE1	2:M:2:NAG:H5	2.55	0.41
1:D:146:TYR:C	1:D:148:ASN:H	2.27	0.41
1:D:551:TYR:O	1:D:552:TRP:C	2.63	0.41
1:A:11:ARG:HG2	1:A:163:PRO:HA	2.02	0.41
1:A:551:TYR:O	1:A:552:TRP:C	2.63	0.41
1:C:104:VAL:C	1:C:106:ASN:N	2.79	0.41
1:C:137:TRP:CE3	1:C:233:THR:HB	2.55	0.41
1:C:357:ASN:HB2	1:C:416:TYR:O	2.21	0.41
1:A:174:MET:HB2	1:A:233:THR:OG1	2.21	0.41
1:B:94:HIS:CE1	1:B:437:HIS:ND1	2.88	0.41
1:C:193:ALA:HA	1:C:194:PRO:HD3	1.69	0.41
1:A:153:THR:HG21	1:A:172:PRO:HG3	2.02	0.41
1:A:238:HIS:CD2	1:A:269:ALA:HA	2.56	0.41
1:A:382:VAL:HG11	1:A:564:LEU:HD21	2.02	0.41
1:A:425:ASN:HA	1:A:426:PRO:HD3	1.91	0.41
1:B:254:ASP:HA	1:B:273:ARG:HG2	2.02	0.41
1:B:342:VAL:HG12	1:B:344:ARG:HH12	1.85	0.41
1:C:249:THR:O	1:C:276:VAL:HA	2.21	0.41
1:C:367:GLY:O	1:C:368:THR:C	2.63	0.41
1:D:72:ALA:O	1:D:156:ILE:HA	2.21	0.41
1:D:167:ASP:OD1	1:D:167:ASP:C	2.63	0.41
1:A:297:LEU:HD23	1:A:297:LEU:HA	1.87	0.41
1:D:7:THR:O	1:D:11:ARG:HD3	2.20	0.41
1:D:534:LYS:O	1:D:537:ASP:N	2.54	0.41
1:A:167:ASP:C	1:A:167:ASP:OD1	2.64	0.41
1:A:347:PRO:HD2	1:A:416:TYR:OH	2.20	0.41
1:A:381:ASN:HB2	9:A:736:HOH:O	2.21	0.41
1:A:435:HIS:CE1	1:A:478:ASP:HB2	2.56	0.41
1:B:136:SER:HA	1:B:273:ARG:HD2	2.02	0.41
1:B:447:SER:HB3	1:B:487:TRP:CD1	2.56	0.41
1:C:342:VAL:HG12	1:C:344:ARG:HH12	1.85	0.41
1:D:8:PRO:HA	1:D:164:TYR:O	2.21	0.41
1:D:92:SER:OG	1:D:110:ASN:CB	2.68	0.41
1:A:72:ALA:O	1:A:156:ILE:HA	2.21	0.41
1:A:427:ILE:HG12	1:B:427:ILE:HG12	2.03	0.41
1:B:387:PRO:HB2	1:B:541:PHE:HZ	1.85	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:11:ARG:HB3	1:D:160:ALA:HB3	2.02	0.41
1:D:106:ASN:N	1:D:106:ASN:HD22	2.19	0.41
1:A:259:ASN:OD1	1:A:341:VAL:HG22	2.21	0.40
1:B:20:ILE:O	1:B:20:ILE:HG13	2.22	0.40
1:B:107:ASP:HB3	1:B:113:THR:HG21	2.03	0.40
1:B:180:SER:OG	1:B:183:GLU:HG3	2.21	0.40
1:B:473:ASN:N	1:B:474:PRO:HD3	2.36	0.40
1:B:545:CYS:O	1:B:549:ASN:ND2	2.50	0.40
1:D:443:VAL:HA	1:D:490:LEU:HD23	2.01	0.40
1:A:157:ASP:HA	9:A:652:HOH:O	2.21	0.40
1:A:187:PHE:CE2	1:A:194:PRO:HG3	2.56	0.40
1:B:365:ILE:O	1:B:365:ILE:HG22	2.20	0.40
1:C:106:ASN:N	1:C:106:ASN:ND2	2.66	0.40
1:D:137:TRP:CZ2	1:D:271:GLY:HA2	2.57	0.40
1:D:259:ASN:OD1	1:D:341:VAL:HG22	2.21	0.40
1:A:336:LEU:O	1:A:475:VAL:HG23	2.21	0.40
1:A:389:LEU:HD22	1:A:529:LEU:HD11	2.03	0.40
1:B:104:VAL:C	1:B:106:ASN:N	2.79	0.40
1:C:89:ASN:HB3	1:C:145:GLN:NE2	2.36	0.40
1:C:252:ALA:HB3	1:C:275:ASP:HB2	2.03	0.40
1:C:348:VAL:CG1	1:C:444:LEU:HD22	2.51	0.40
1:D:133:TYR:CE1	1:D:159:PRO:HD2	2.57	0.40
1:D:134:GLY:HA2	9:D:653:HOH:O	2.20	0.40
1:D:325:LEU:HB3	1:D:326:PRO:HD2	2.02	0.40
3:F:2:NAG:H82	3:F:2:NAG:C1	2.52	0.40
4:I:1:NAG:H62	4:I:2:NAG:C7	2.51	0.40
1:A:316:PRO:HB2	1:A:318:THR:HG23	2.04	0.40
1:A:428:VAL:HG11	1:B:427:ILE:HD11	2.02	0.40
1:A:442:LEU:HB3	1:A:474:PRO:HG2	2.04	0.40
1:D:112:VAL:HG21	1:D:501:LEU:HD21	2.03	0.40
1:A:162:LEU:HD12	1:A:162:LEU:HA	1.98	0.40
1:A:249:THR:O	1:A:276:VAL:HA	2.21	0.40
1:B:177:TYR:CD1	1:B:177:TYR:N	2.90	0.40
1:C:41:LEU:CD2	1:C:61:VAL:HG22	2.51	0.40
1:C:151:VAL:HG22	1:C:152:GLY:N	2.36	0.40
1:C:282:SER:HB3	1:C:283:PRO:CD	2.52	0.40
1:D:206:ARG:NE	9:D:660:HOH:O	2.54	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	562/604 (93%)	517 (92%)	42 (8%)	3 (0%)	24	43
1	B	562/604 (93%)	502 (89%)	60 (11%)	0	100	100
1	C	562/604 (93%)	505 (90%)	57 (10%)	0	100	100
1	D	562/604 (93%)	515 (92%)	44 (8%)	3 (0%)	24	43
All	All	2248/2416 (93%)	2039 (91%)	203 (9%)	6 (0%)	36	55

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	494	THR
1	D	494	THR
1	A	284	VAL
1	D	222	PRO
1	D	497	PRO
1	A	222	PRO

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	484/513 (94%)	474 (98%)	10 (2%)	47	74
1	B	484/513 (94%)	475 (98%)	9 (2%)	50	76
1	C	484/513 (94%)	474 (98%)	10 (2%)	47	74
1	D	484/513 (94%)	474 (98%)	10 (2%)	47	74

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1936/2052 (94%)	1897 (98%)	39 (2%)	48 75

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	35	VAL
1	A	106	ASN
1	A	110	ASN
1	A	141	HIS
1	A	268	LEU
1	A	273	ARG
1	A	311	ARG
1	A	349	ASN
1	A	368	THR
1	A	427	ILE
1	B	35	VAL
1	B	91	THR
1	B	106	ASN
1	B	110	ASN
1	B	141	HIS
1	B	179	ARG
1	B	273	ARG
1	B	349	ASN
1	B	466	LEU
1	C	35	VAL
1	C	91	THR
1	C	106	ASN
1	C	110	ASN
1	C	141	HIS
1	C	179	ARG
1	C	250	VAL
1	C	273	ARG
1	C	349	ASN
1	C	466	LEU
1	D	106	ASN
1	D	110	ASN
1	D	141	HIS
1	D	268	LEU
1	D	273	ARG
1	D	311	ARG
1	D	338	LEU
1	D	349	ASN

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Mol	Chain	Res	Type
1	D	368	THR
1	D	427	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	103	ASN
1	A	207	HIS
1	A	237	ASN
1	A	528	GLN
1	B	10	ASN
1	B	38	GLN
1	B	103	ASN
1	B	207	HIS
1	B	240	GLN
1	B	247	ASN
1	B	286	ASN
1	C	10	ASN
1	C	38	GLN
1	C	103	ASN
1	C	207	HIS
1	C	240	GLN
1	C	286	ASN
1	D	38	GLN
1	D	103	ASN
1	D	207	HIS
1	D	237	ASN
1	D	528	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	0.50	0	17,19,21	1.64	4 (23%)
2	NAG	E	2	2	14,14,15	0.42	0	17,19,21	1.61	2 (11%)
3	NAG	F	1	3,1	14,14,15	0.55	0	17,19,21	1.01	0
3	NAG	F	2	3	14,14,15	0.61	0	17,19,21	1.70	4 (23%)
3	BMA	F	3	3	11,11,12	0.51	0	15,15,17	1.01	0
3	MAN	F	4	3	11,11,12	0.83	1 (9%)	15,15,17	1.18	3 (20%)
3	MAN	F	5	3	11,11,12	0.88	0	15,15,17	1.72	4 (26%)
3	MAN	F	6	3	11,11,12	0.46	0	15,15,17	0.81	1 (6%)
2	NAG	G	1	2,1	14,14,15	0.67	0	17,19,21	1.29	2 (11%)
2	NAG	G	2	2	14,14,15	0.54	0	17,19,21	1.02	1 (5%)
2	NAG	H	1	2,1	14,14,15	0.52	0	17,19,21	1.54	3 (17%)
2	NAG	H	2	2	14,14,15	0.51	0	17,19,21	1.29	2 (11%)
4	NAG	I	1	1,4	14,14,15	0.75	0	17,19,21	1.15	2 (11%)
4	NAG	I	2	4	14,14,15	0.64	0	17,19,21	0.93	0
4	BMA	I	3	4	11,11,12	0.67	0	15,15,17	0.90	0
4	MAN	I	4	4	11,11,12	0.65	0	15,15,17	1.33	2 (13%)
4	MAN	I	5	4	11,11,12	0.48	0	15,15,17	1.62	5 (33%)
2	NAG	J	1	2,1	14,14,15	0.45	0	17,19,21	0.85	0
2	NAG	J	2	2	14,14,15	0.56	0	17,19,21	0.94	0
2	NAG	K	1	2,1	14,14,15	0.43	0	17,19,21	1.10	1 (5%)
2	NAG	K	2	2	14,14,15	0.52	0	17,19,21	0.86	0
5	NAG	L	1	5,1	14,14,15	0.55	0	17,19,21	0.83	0
5	NAG	L	2	5	14,14,15	0.43	0	17,19,21	1.23	1 (5%)
5	BMA	L	3	5	11,11,12	0.77	0	15,15,17	1.18	2 (13%)
2	NAG	M	1	2,1	14,14,15	0.67	0	17,19,21	1.26	1 (5%)
2	NAG	M	2	2	14,14,15	0.42	0	17,19,21	1.93	5 (29%)
2	NAG	N	1	2,1	14,14,15	0.53	0	17,19,21	1.19	1 (5%)
2	NAG	N	2	2	14,14,15	0.59	0	17,19,21	1.10	1 (5%)
5	NAG	O	1	5,1	14,14,15	0.44	0	17,19,21	1.45	1 (5%)
5	NAG	O	2	5	14,14,15	0.51	0	17,19,21	0.98	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	BMA	O	3	5	11,11,12	0.60	0	15,15,17	0.60	0
4	NAG	P	1	1,4	14,14,15	0.49	0	17,19,21	1.64	3 (17%)
4	NAG	P	2	4	14,14,15	0.72	0	17,19,21	1.67	3 (17%)
4	BMA	P	3	4	11,11,12	0.66	0	15,15,17	0.81	0
4	MAN	P	4	4	11,11,12	0.74	0	15,15,17	1.50	3 (20%)
4	MAN	P	5	4	11,11,12	0.64	0	15,15,17	1.16	2 (13%)
2	NAG	Q	1	2,1	14,14,15	0.54	0	17,19,21	1.20	1 (5%)
2	NAG	Q	2	2	14,14,15	0.40	0	17,19,21	1.35	2 (11%)

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	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
3	NAG	F	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	F	2	3	-	5/6/23/26	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
3	MAN	F	5	3	-	2/2/19/22	0/1/1/1
3	MAN	F	6	3	-	2/2/19/22	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	1/6/23/26	0/1/1/1
2	NAG	H	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	4/6/23/26	0/1/1/1
4	NAG	I	1	1,4	-	0/6/23/26	0/1/1/1
4	NAG	I	2	4	-	2/6/23/26	0/1/1/1
4	BMA	I	3	4	-	0/2/19/22	0/1/1/1
4	MAN	I	4	4	-	0/2/19/22	0/1/1/1
4	MAN	I	5	4	-	2/2/19/22	0/1/1/1
2	NAG	J	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	J	2	2	-	2/6/23/26	0/1/1/1
2	NAG	K	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	K	2	2	-	2/6/23/26	0/1/1/1
5	NAG	L	1	5,1	-	1/6/23/26	0/1/1/1
5	NAG	L	2	5	-	3/6/23/26	0/1/1/1

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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	4	MAN	O5-C1	-2.05	1.40	1.43

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	2	NAG	C1-O5-C5	4.96	118.83	112.19
5	O	1	NAG	C1-O5-C5	4.62	118.38	112.19
4	P	2	NAG	C4-C3-C2	4.60	117.76	111.02
2	M	2	NAG	C1-O5-C5	4.30	117.95	112.19
5	L	2	NAG	C2-N2-C7	-4.20	117.28	122.90
2	H	1	NAG	C1-O5-C5	4.01	117.56	112.19
2	E	1	NAG	C1-O5-C5	3.99	117.53	112.19
3	F	2	NAG	O5-C1-C2	-3.94	105.19	111.29
3	F	5	MAN	C2-C3-C4	-3.91	103.99	110.86
2	M	1	NAG	C1-O5-C5	3.90	117.41	112.19
2	Q	2	NAG	C1-O5-C5	3.86	117.36	112.19
2	M	2	NAG	C4-C3-C2	-3.82	105.42	111.02
2	N	1	NAG	C2-N2-C7	-3.71	117.92	122.90
3	F	2	NAG	C1-C2-N2	3.52	115.98	110.43
2	Q	1	NAG	C1-O5-C5	3.50	116.88	112.19
4	P	1	NAG	C2-N2-C7	-3.45	118.28	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	2	NAG	C2-N2-C7	-3.39	118.36	122.90
3	F	2	NAG	C2-N2-C7	3.23	127.22	122.90
2	M	2	NAG	C2-N2-C7	-3.22	118.58	122.90
2	H	2	NAG	C1-O5-C5	3.18	116.44	112.19
4	P	4	MAN	C3-C4-C5	3.15	115.95	110.23
4	P	1	NAG	C3-C4-C5	2.96	115.59	110.23
3	F	5	MAN	O2-C2-C1	-2.91	102.56	109.22
5	O	2	NAG	C1-O5-C5	2.88	116.05	112.19
4	P	4	MAN	C2-C3-C4	2.88	115.92	110.86
2	M	2	NAG	O5-C1-C2	2.80	115.62	111.29
2	E	1	NAG	C2-N2-C7	-2.79	119.17	122.90
4	I	4	MAN	C1-O5-C5	2.74	115.86	112.19
4	I	5	MAN	O5-C5-C6	2.67	112.86	107.66
4	P	1	NAG	O5-C1-C2	-2.66	107.17	111.29
4	I	5	MAN	C2-C3-C4	-2.62	106.25	110.86
2	K	1	NAG	C2-N2-C7	-2.60	119.42	122.90
2	Q	2	NAG	C2-N2-C7	-2.53	119.50	122.90
2	E	1	NAG	C6-C5-C4	-2.51	106.86	113.02
3	F	5	MAN	C1-C2-C3	2.48	113.26	109.64
2	G	1	NAG	C1-O5-C5	2.48	115.50	112.19
4	P	5	MAN	C2-C3-C4	2.46	115.19	110.86
2	N	2	NAG	C1-C2-N2	-2.45	106.58	110.43
4	I	5	MAN	O3-C3-C4	2.41	116.06	110.38
4	I	4	MAN	C3-C4-C5	-2.40	105.88	110.23
4	I	5	MAN	C3-C4-C5	-2.39	105.89	110.23
2	H	1	NAG	C6-C5-C4	-2.39	107.16	113.02
3	F	4	MAN	C1-C2-C3	-2.38	106.18	109.64
2	E	2	NAG	O5-C1-C2	2.36	114.94	111.29
4	I	1	NAG	O5-C1-C2	-2.33	107.68	111.29
4	I	1	NAG	C4-C3-C2	2.32	114.42	111.02
2	G	1	NAG	O5-C5-C6	2.32	112.18	107.66
4	P	4	MAN	O2-C2-C1	2.23	114.33	109.22
3	F	6	MAN	C1-O5-C5	2.22	115.16	112.19
4	P	2	NAG	C1-O5-C5	-2.20	109.24	112.19
2	G	2	NAG	C2-N2-C7	-2.20	119.95	122.90
2	E	1	NAG	O6-C6-C5	-2.16	103.98	111.33
2	H	1	NAG	O4-C4-C5	-2.15	104.03	109.32
5	L	3	BMA	C2-C3-C4	2.14	114.63	110.86
3	F	2	NAG	C4-C3-C2	-2.12	107.91	111.02
4	P	5	MAN	C1-C2-C3	2.12	112.73	109.64
5	L	3	BMA	C1-C2-C3	2.12	112.72	109.64
4	P	2	NAG	O7-C7-C8	-2.10	118.31	122.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	I	5	MAN	O4-C4-C3	2.08	115.29	110.38
3	F	5	MAN	C6-C5-C4	-2.07	107.95	113.02
2	M	2	NAG	O5-C5-C6	2.04	111.63	107.66
3	F	4	MAN	O5-C1-C2	-2.03	105.94	110.79
3	F	4	MAN	O3-C3-C2	-2.00	105.97	110.05

There are no chirality outliers.

All (64) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	K	2	NAG	C8-C7-N2-C2
2	K	2	NAG	O7-C7-N2-C2
3	F	2	NAG	C1-C2-N2-C7
3	F	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O7-C7-N2-C2
4	P	1	NAG	O7-C7-N2-C2
5	L	2	NAG	C8-C7-N2-C2
5	L	2	NAG	O7-C7-N2-C2
5	O	1	NAG	O7-C7-N2-C2
2	J	2	NAG	C4-C5-C6-O6
2	Q	1	NAG	C8-C7-N2-C2
2	Q	1	NAG	O7-C7-N2-C2
4	P	1	NAG	C8-C7-N2-C2
5	O	1	NAG	C8-C7-N2-C2
2	H	2	NAG	O5-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	Q	2	NAG	C4-C5-C6-O6
3	F	5	MAN	O5-C5-C6-O6
3	F	5	MAN	C4-C5-C6-O6
3	F	6	MAN	O5-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
4	I	5	MAN	C4-C5-C6-O6
4	P	1	NAG	C4-C5-C6-O6
2	M	1	NAG	C8-C7-N2-C2
2	Q	2	NAG	O5-C5-C6-O6
5	O	1	NAG	O5-C5-C6-O6
2	M	1	NAG	O7-C7-N2-C2
2	N	1	NAG	C8-C7-N2-C2
3	F	1	NAG	C4-C5-C6-O6
4	P	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	I	5	MAN	O5-C5-C6-O6
5	O	1	NAG	C4-C5-C6-O6
2	M	2	NAG	C8-C7-N2-C2
2	N	1	NAG	O7-C7-N2-C2
5	O	2	NAG	C8-C7-N2-C2
3	F	3	BMA	C4-C5-C6-O6
3	F	3	BMA	O5-C5-C6-O6
4	P	4	MAN	C4-C5-C6-O6
2	E	2	NAG	C8-C7-N2-C2
2	H	2	NAG	C8-C7-N2-C2
2	M	2	NAG	O7-C7-N2-C2
5	O	2	NAG	O7-C7-N2-C2
3	F	1	NAG	O5-C5-C6-O6
4	P	4	MAN	O5-C5-C6-O6
2	N	1	NAG	O5-C5-C6-O6
3	F	6	MAN	C4-C5-C6-O6
4	P	5	MAN	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
4	P	3	BMA	O5-C5-C6-O6
2	H	2	NAG	O7-C7-N2-C2
5	O	3	BMA	C4-C5-C6-O6
2	E	2	NAG	O7-C7-N2-C2
5	O	2	NAG	O5-C5-C6-O6
5	L	1	NAG	C4-C5-C6-O6
5	O	3	BMA	O5-C5-C6-O6
2	J	1	NAG	C8-C7-N2-C2
3	F	2	NAG	O5-C5-C6-O6
2	J	1	NAG	O7-C7-N2-C2
5	L	2	NAG	C4-C5-C6-O6
4	I	2	NAG	O5-C5-C6-O6
4	I	2	NAG	C4-C5-C6-O6

There are no ring outliers.

17 monomers are involved in 25 short contacts:

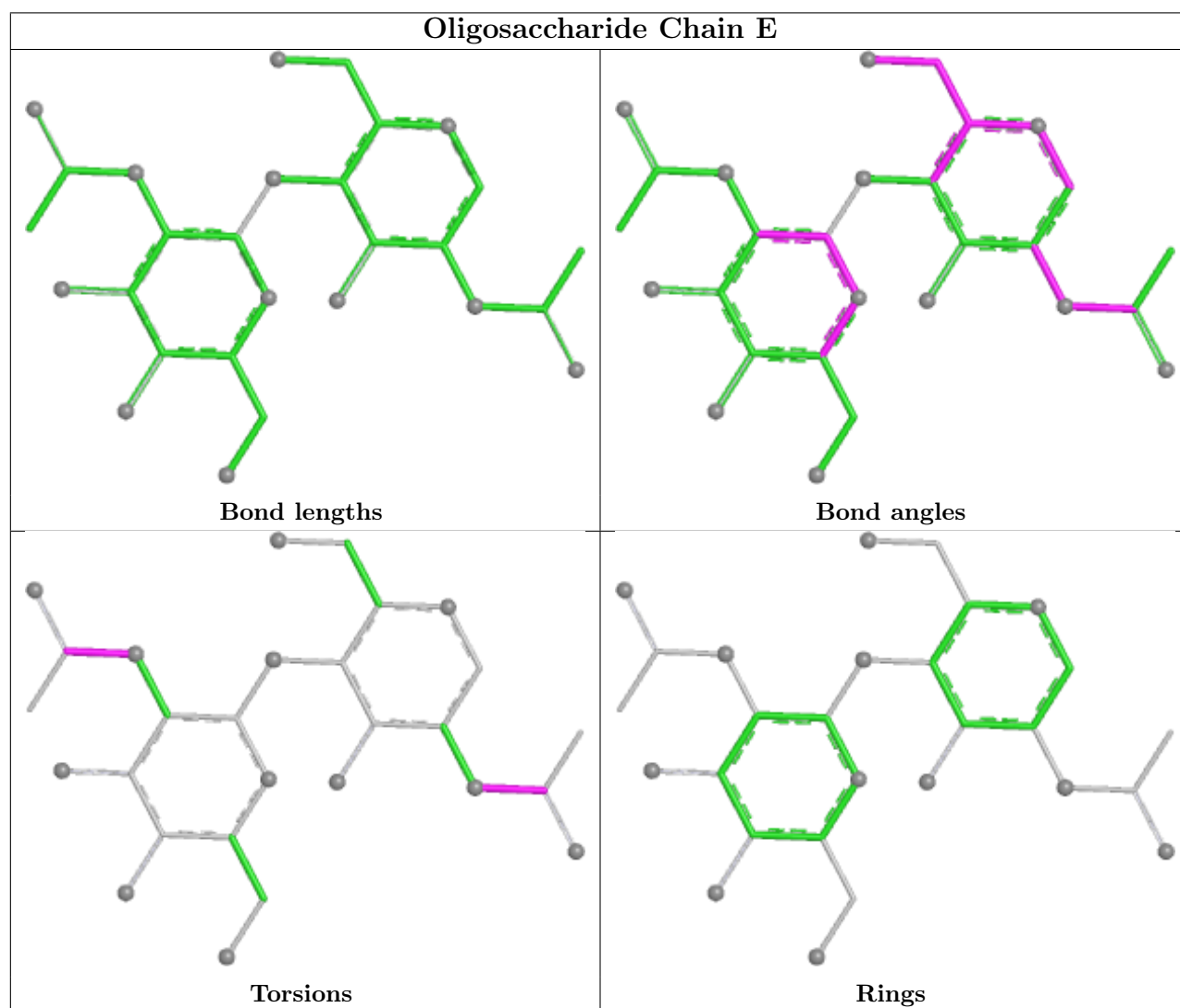
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	1	NAG	1	0
4	I	2	NAG	1	0
5	L	1	NAG	1	0
2	N	1	NAG	3	0
5	L	3	BMA	1	0

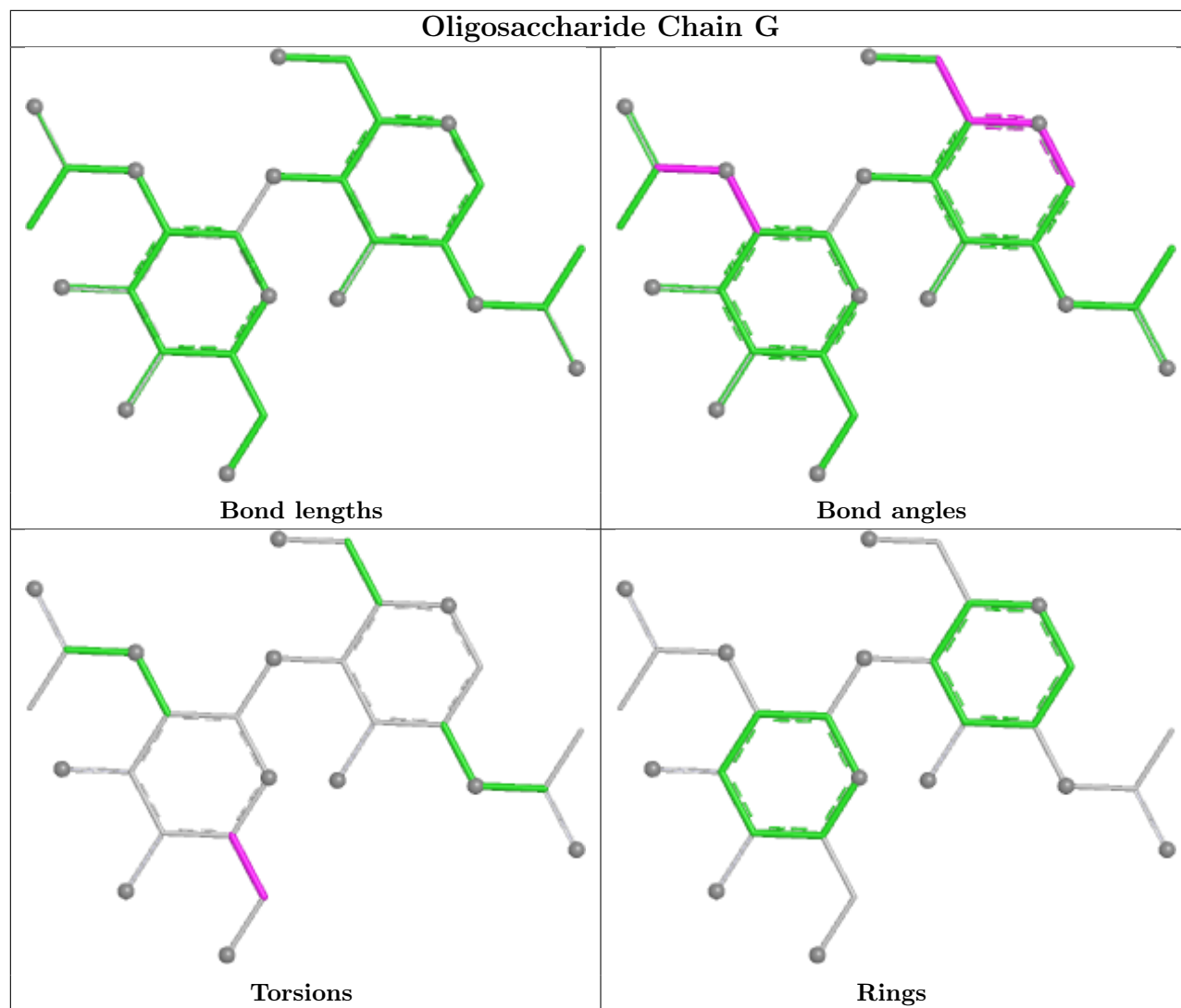
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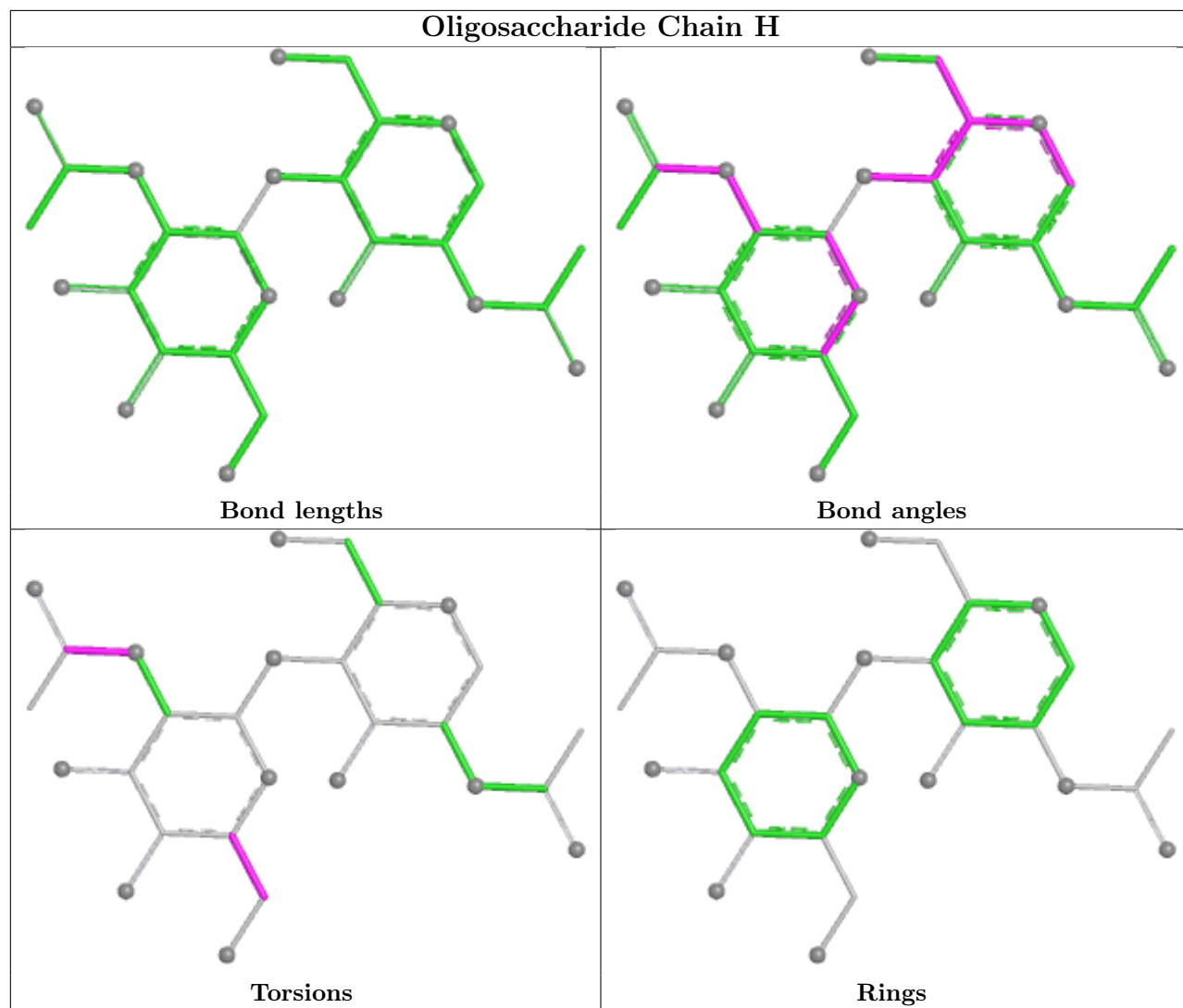
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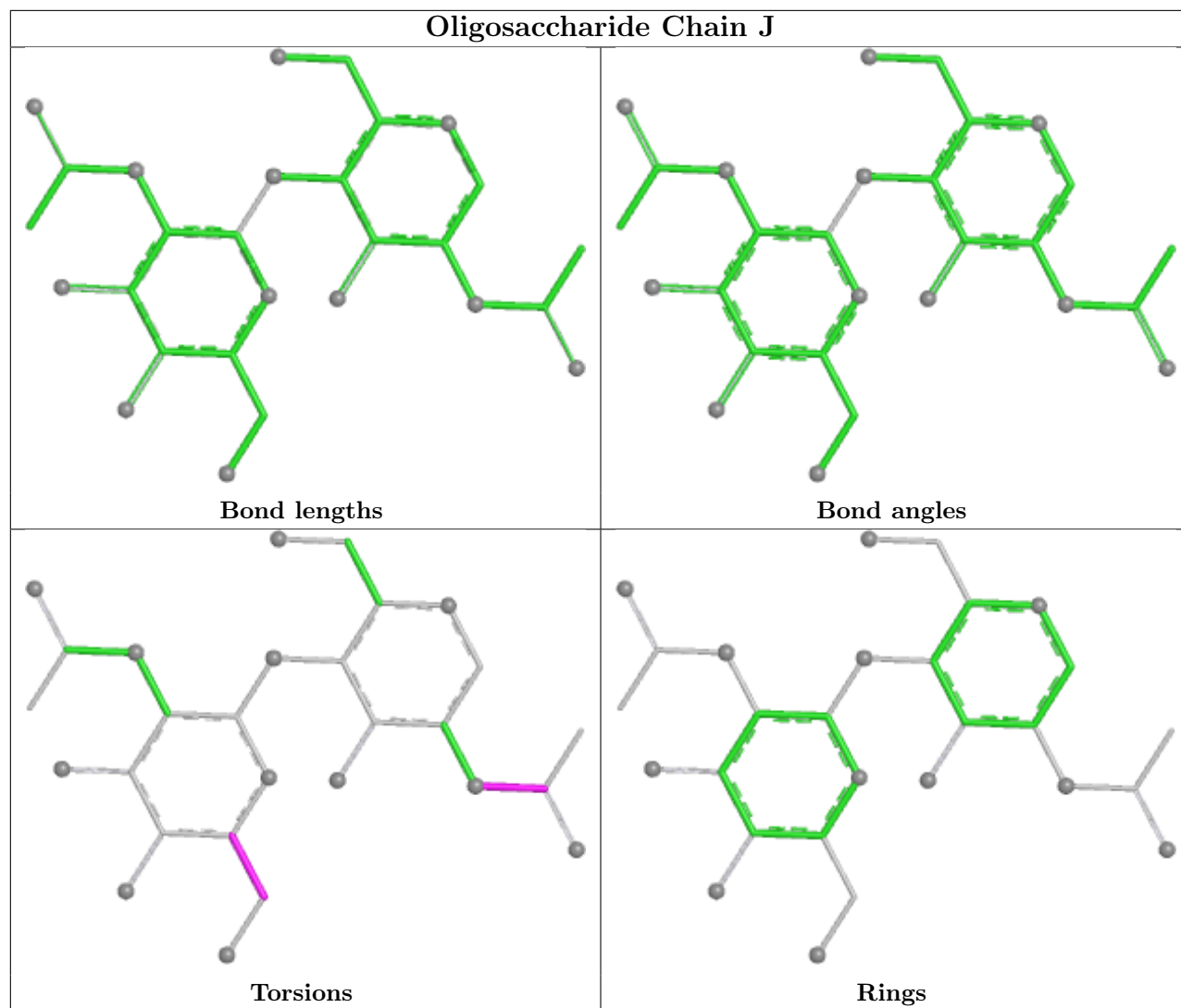
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	4	MAN	1	0
3	F	2	NAG	1	0
2	J	2	NAG	1	0
5	L	2	NAG	2	0
5	O	1	NAG	1	0
4	P	2	NAG	1	0
2	E	1	NAG	3	0
2	E	2	NAG	1	0
2	K	2	NAG	4	0
4	P	1	NAG	3	0
2	M	2	NAG	1	0
2	N	2	NAG	1	0

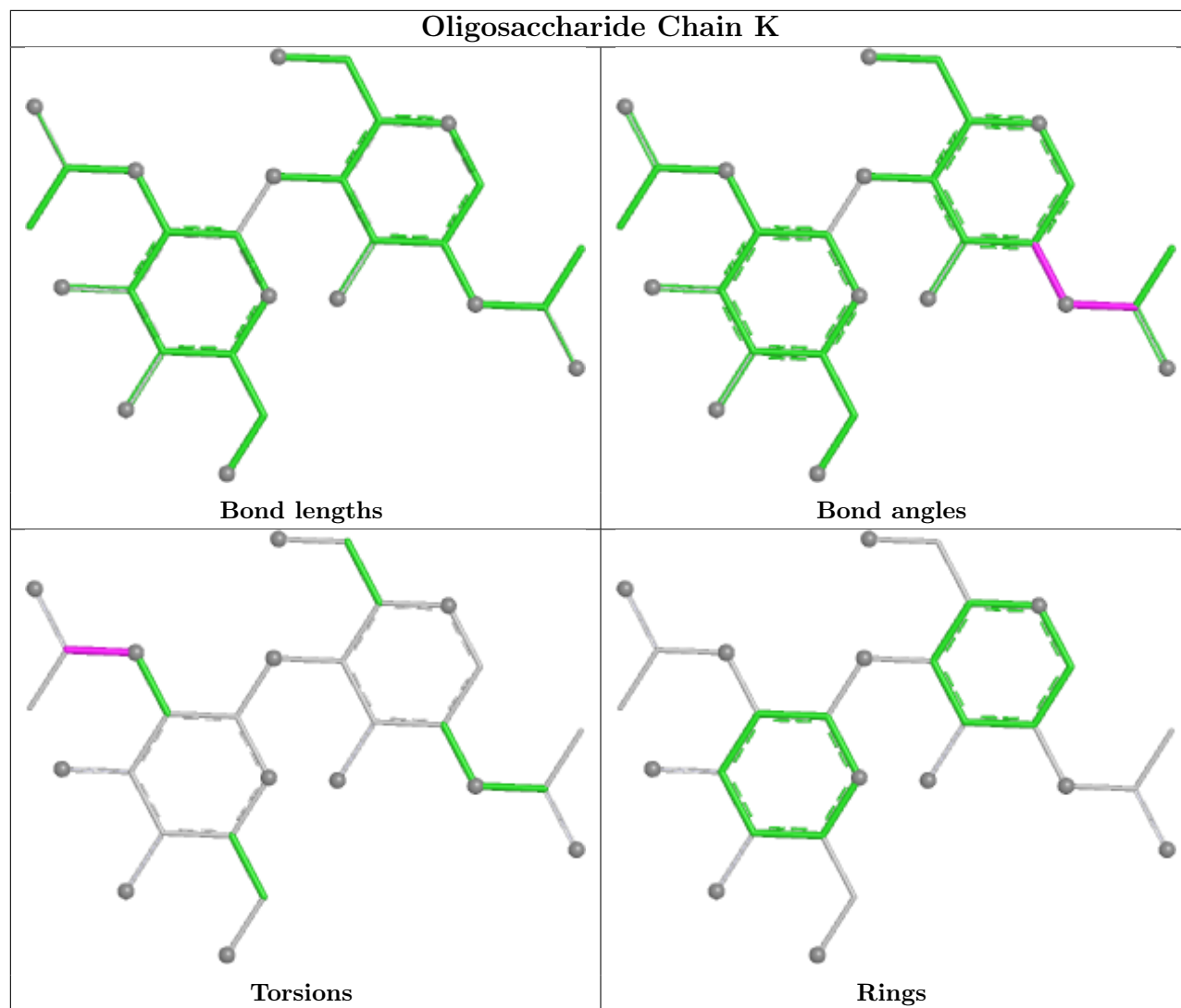
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

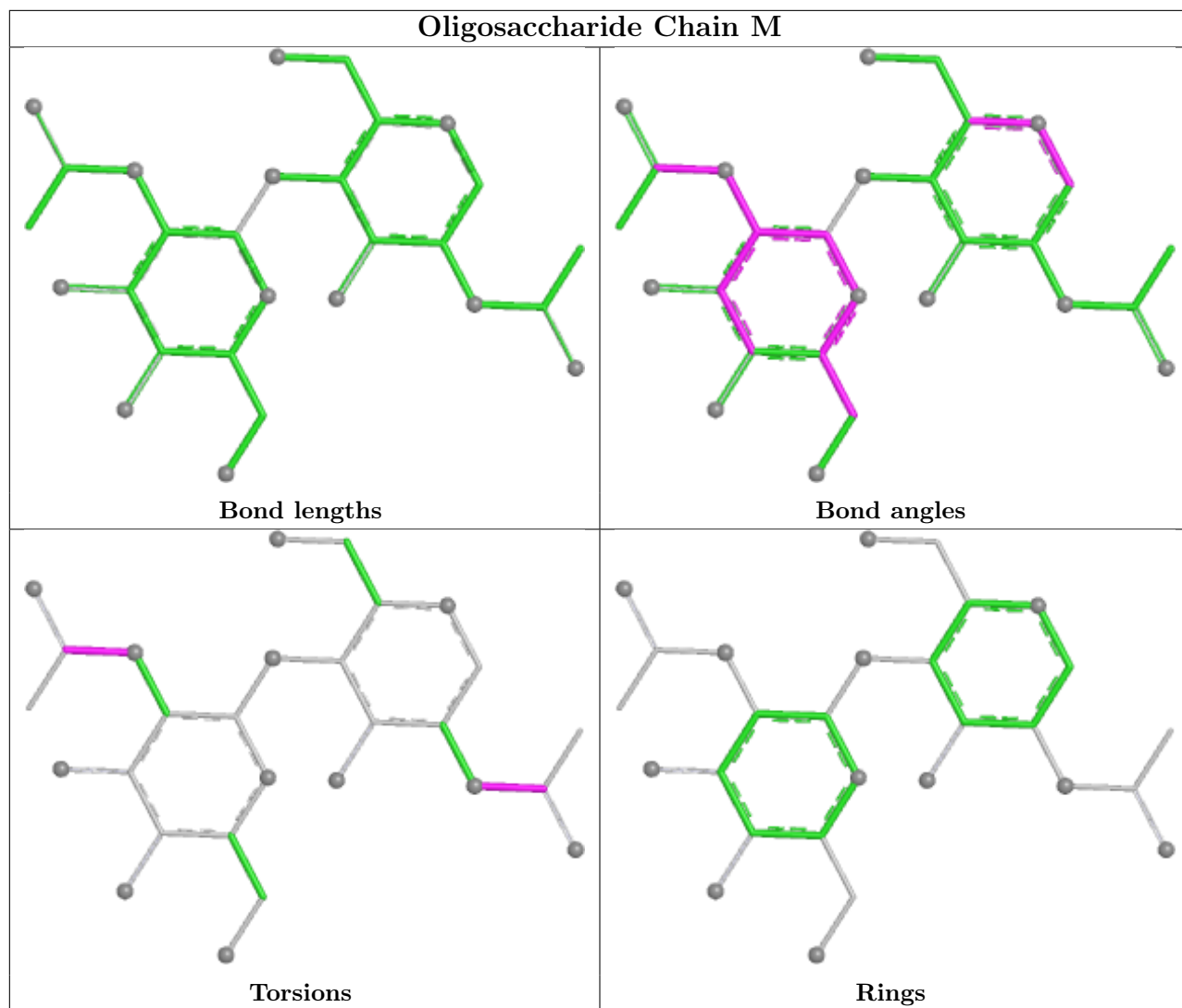


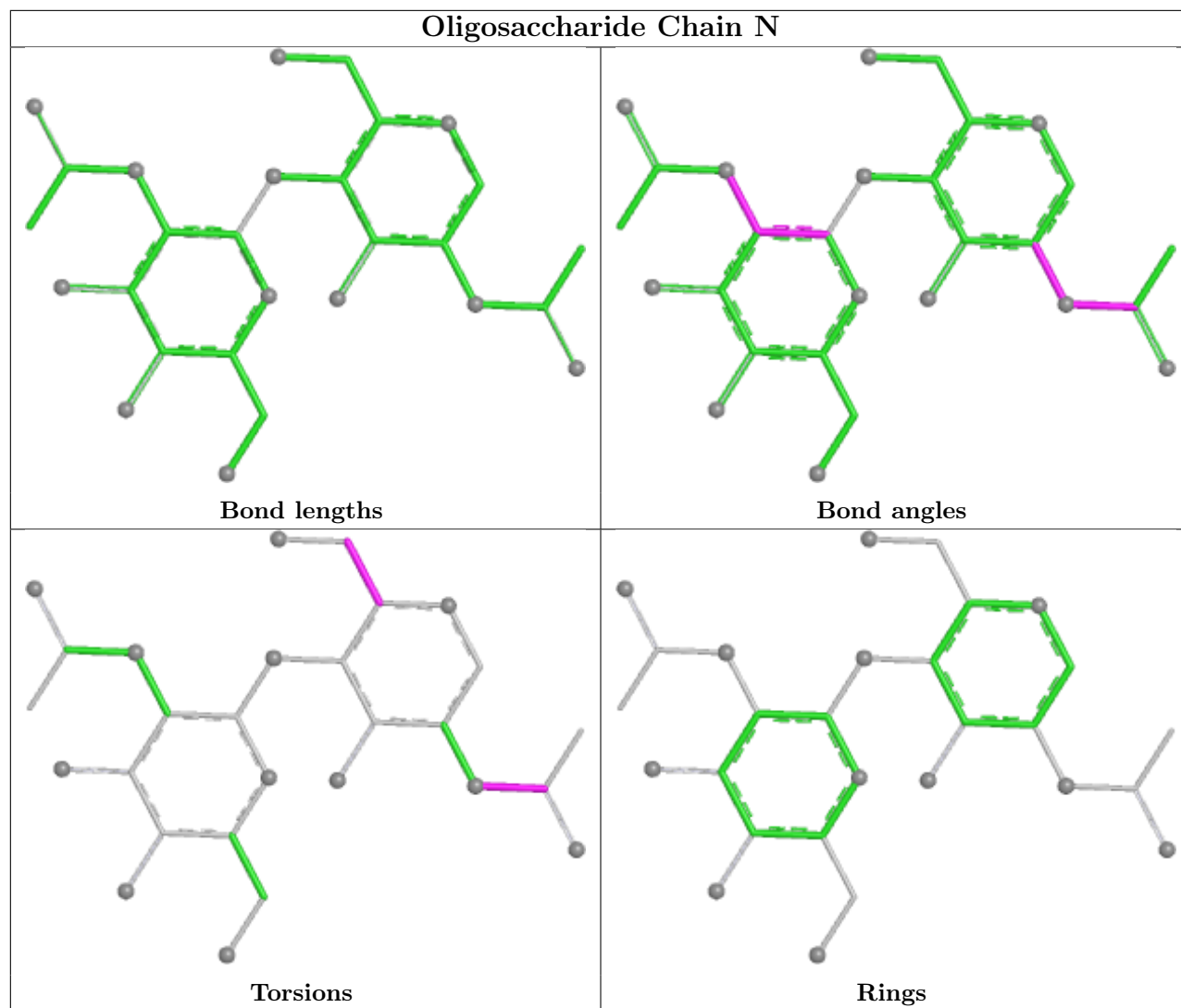


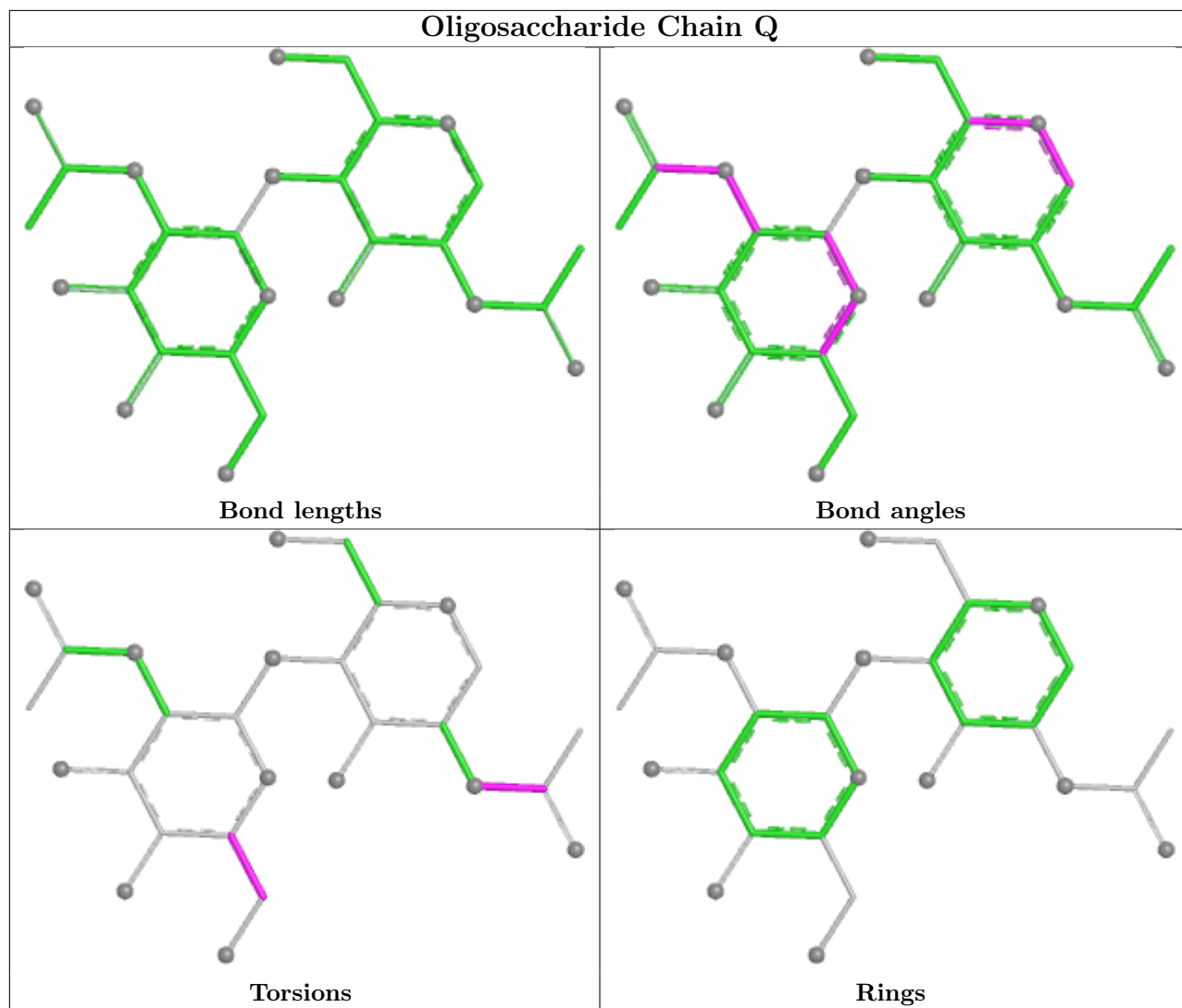


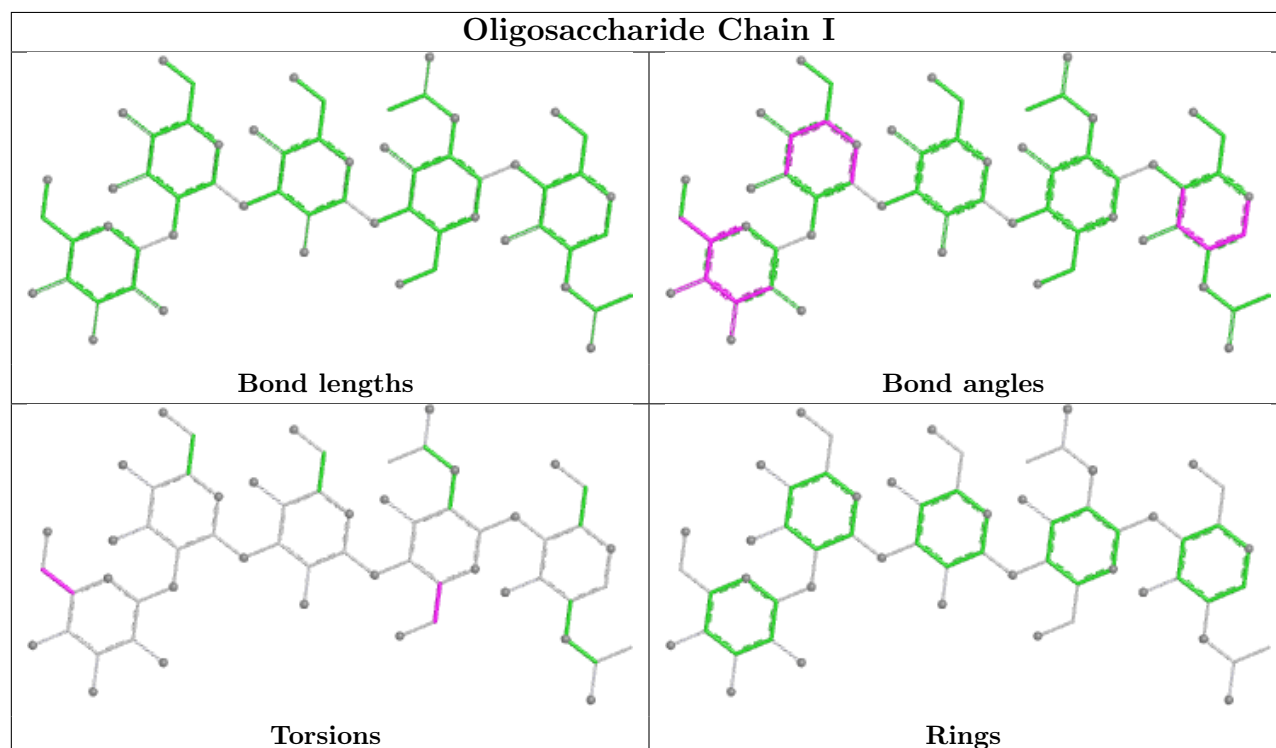
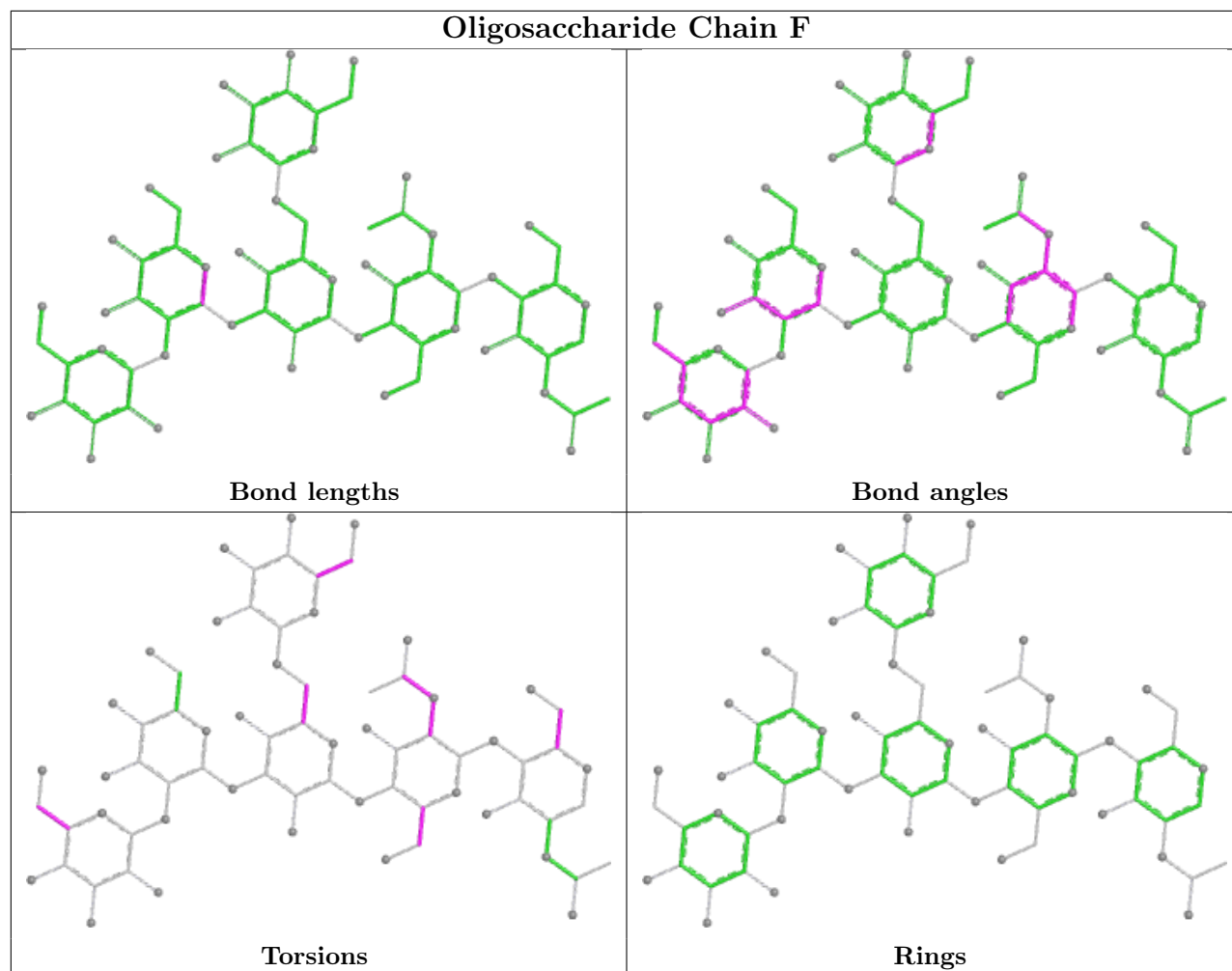


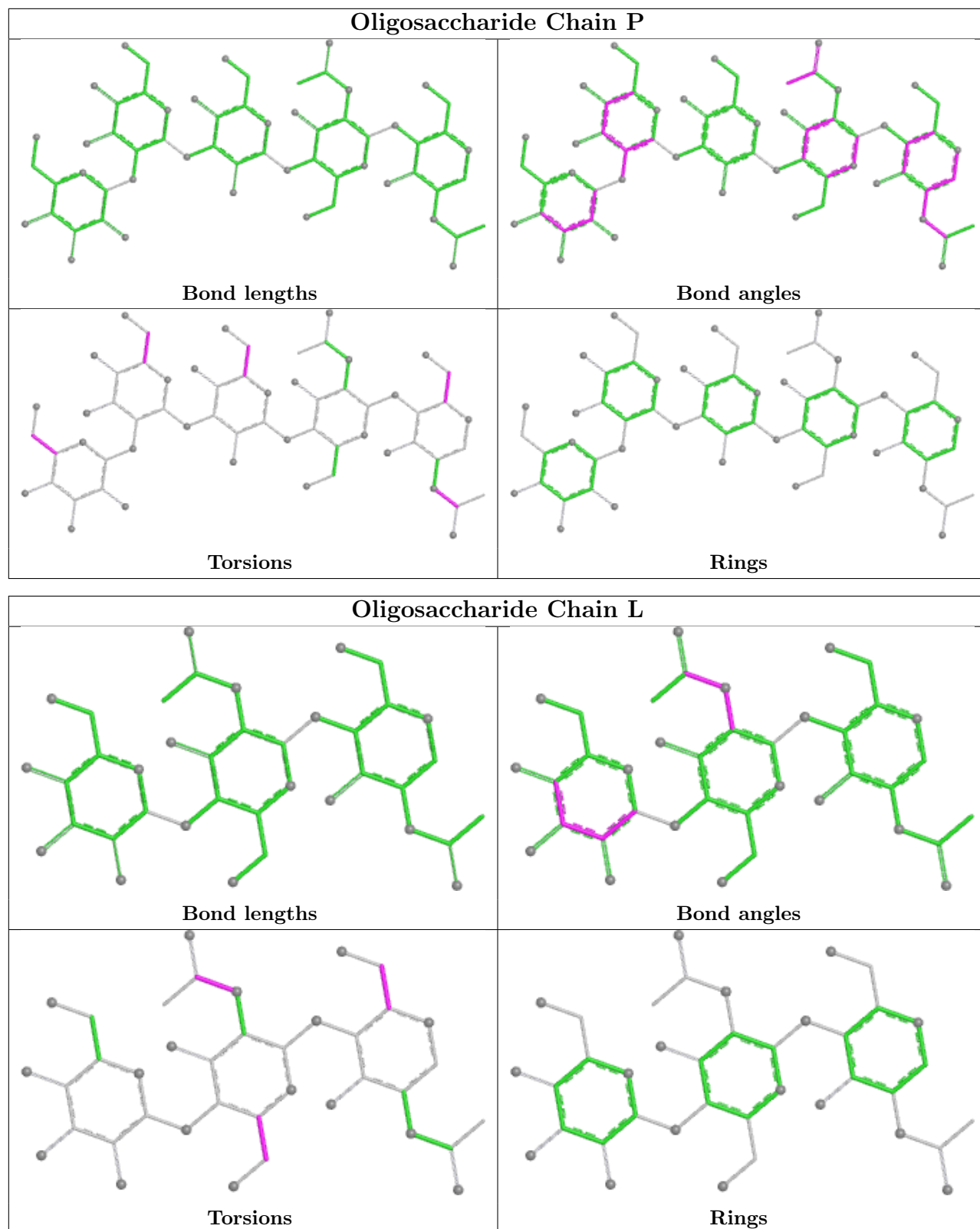


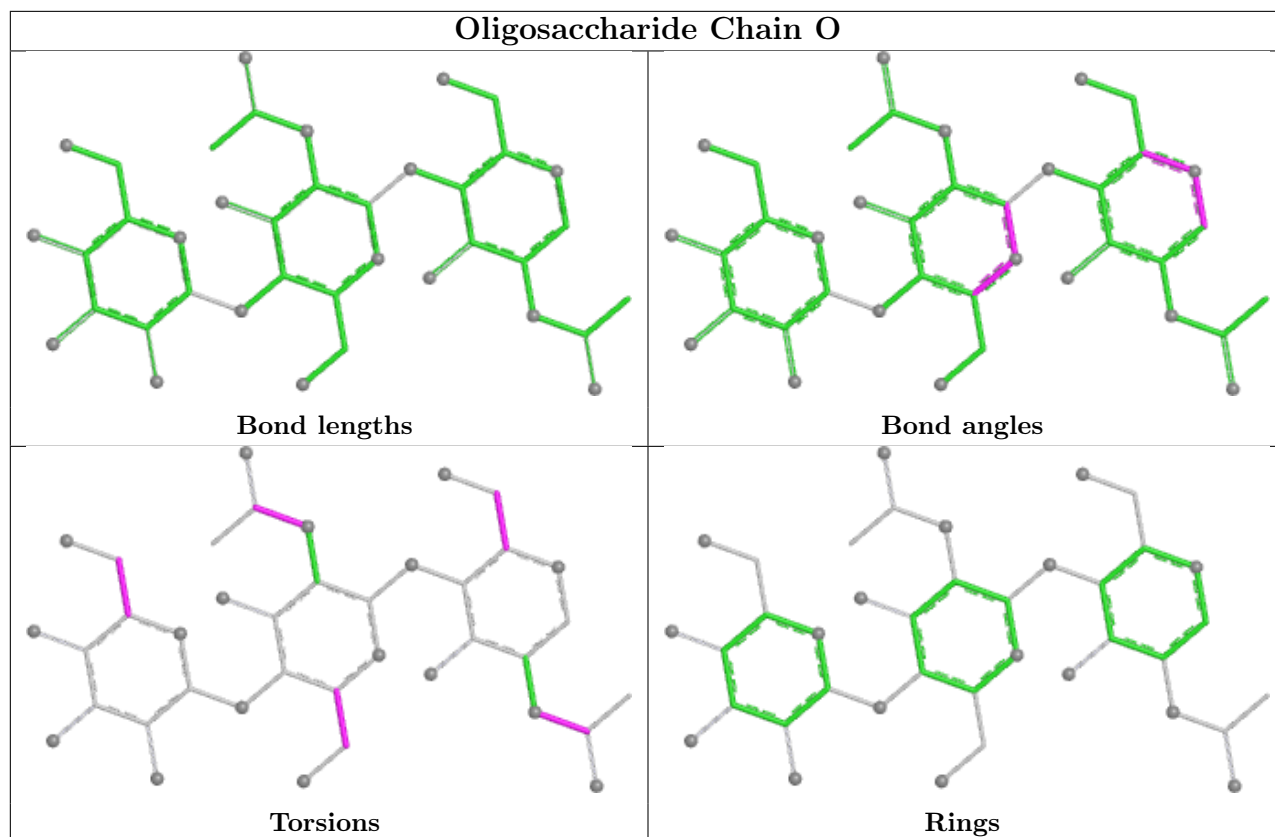












5.6 Ligand geometry [i](#)

Of 25 ligands modelled in this entry, 16 are monoatomic - leaving 9 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
7	OXY	A	620	6	1,1,1	0.17	0	-		
8	NAG	A	720	1	14,14,15	0.58	0	17,19,21	1.03	1 (5%)
8	NAG	C	730	1	14,14,15	0.47	0	17,19,21	1.64	2 (11%)
8	NAG	A	760	1	14,14,15	0.64	0	17,19,21	1.37	1 (5%)
8	NAG	B	770	1	14,14,15	0.58	0	17,19,21	1.34	3 (17%)
8	NAG	D	720	1	14,14,15	0.54	0	17,19,21	1.52	3 (17%)
8	NAG	B	760	1	14,14,15	0.57	0	17,19,21	1.57	3 (17%)
8	NAG	B	720	1	14,14,15	0.53	0	17,19,21	1.06	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	NAG	A	730	1	14,14,15	0.60	0	17,19,21	1.53	2 (11%)

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	NAG	A	720	1	-	4/6/23/26	0/1/1/1
8	NAG	C	730	1	-	4/6/23/26	0/1/1/1
8	NAG	A	760	1	-	2/6/23/26	0/1/1/1
8	NAG	B	770	1	-	2/6/23/26	0/1/1/1
8	NAG	D	720	1	-	2/6/23/26	0/1/1/1
8	NAG	B	760	1	-	2/6/23/26	0/1/1/1
8	NAG	B	720	1	-	2/6/23/26	0/1/1/1
8	NAG	A	730	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	A	760	NAG	C2-N2-C7	-4.27	117.17	122.90
8	B	760	NAG	C2-N2-C7	-4.27	117.18	122.90
8	A	730	NAG	C1-O5-C5	3.72	117.17	112.19
8	C	730	NAG	C1-O5-C5	3.69	117.14	112.19
8	D	720	NAG	C1-O5-C5	3.64	117.06	112.19
8	C	730	NAG	C4-C3-C2	-3.42	106.01	111.02
8	A	730	NAG	O5-C1-C2	-3.09	106.51	111.29
8	B	770	NAG	C1-O5-C5	-2.87	108.34	112.19
8	D	720	NAG	O5-C1-C2	-2.78	106.98	111.29
8	B	720	NAG	C1-O5-C5	2.55	115.60	112.19
8	B	770	NAG	O5-C5-C6	2.48	112.49	107.66
8	D	720	NAG	C3-C4-C5	2.42	114.63	110.23
8	A	720	NAG	O5-C1-C2	-2.36	107.64	111.29
8	B	760	NAG	O3-C3-C2	2.35	114.29	109.40
8	B	770	NAG	C4-C3-C2	2.32	114.42	111.02
8	B	760	NAG	C8-C7-N2	2.16	119.70	116.12

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	B	770	NAG	C8-C7-N2-C2
8	B	770	NAG	O7-C7-N2-C2
8	A	720	NAG	O5-C5-C6-O6
8	A	720	NAG	C4-C5-C6-O6
8	B	720	NAG	O5-C5-C6-O6
8	A	730	NAG	O5-C5-C6-O6
8	B	760	NAG	O5-C5-C6-O6
8	B	720	NAG	C4-C5-C6-O6
8	B	760	NAG	C4-C5-C6-O6
8	A	720	NAG	C8-C7-N2-C2
8	A	720	NAG	O7-C7-N2-C2
8	C	730	NAG	C8-C7-N2-C2
8	C	730	NAG	O7-C7-N2-C2
8	A	730	NAG	C4-C5-C6-O6
8	A	760	NAG	O5-C5-C6-O6
8	C	730	NAG	O5-C5-C6-O6
8	A	760	NAG	C4-C5-C6-O6
8	D	720	NAG	O5-C5-C6-O6
8	D	720	NAG	C4-C5-C6-O6
8	C	730	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	760	NAG	1	0
8	D	720	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	564/604 (93%)	-1.55	0 100 100	11, 19, 31, 50	0
1	B	564/604 (93%)	-1.55	0 100 100	9, 19, 34, 47	0
1	C	564/604 (93%)	-1.52	0 100 100	10, 20, 34, 48	0
1	D	564/604 (93%)	-1.46	0 100 100	11, 20, 32, 51	0
All	All	2256/2416 (93%)	-1.52	0 100 100	9, 19, 33, 51	0

There are no RSRZ outliers to report.

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	BMA	O	3	11/12	0.96	0.05	38,51,61,65	0
4	BMA	P	3	11/12	0.97	0.04	32,48,57,68	0
3	MAN	F	6	11/12	0.98	0.05	16,32,37,42	0
2	NAG	G	2	14/15	0.98	0.04	19,25,31,35	0
4	MAN	P	5	11/12	0.98	0.05	32,50,60,63	0
2	NAG	N	2	14/15	0.98	0.04	18,33,42,48	0
2	NAG	J	1	14/15	0.99	0.03	8,24,32,34	0
2	NAG	J	2	14/15	0.99	0.03	17,28,33,34	0
2	NAG	K	1	14/15	0.99	0.03	12,20,28,29	0

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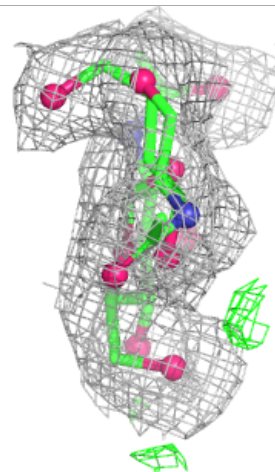
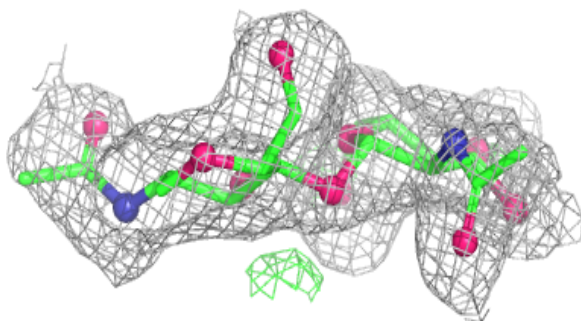
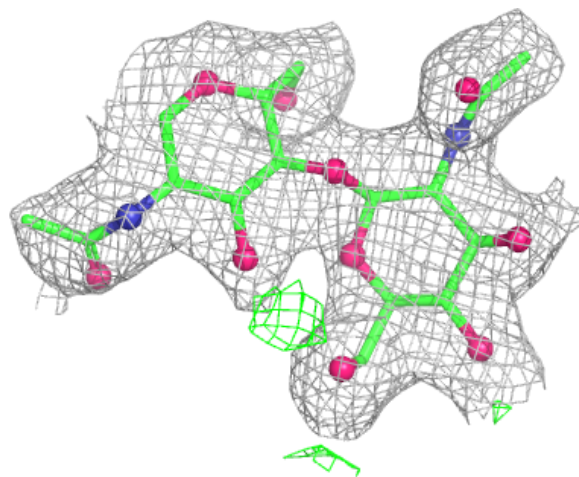
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	NAG	K	2	14/15	0.99	0.03	14,28,40,40	0
2	NAG	M	1	14/15	0.99	0.03	13,23,30,36	0
2	NAG	M	2	14/15	0.99	0.03	12,24,29,34	0
2	NAG	N	1	14/15	0.99	0.04	8,27,36,45	0
2	NAG	E	2	14/15	0.99	0.03	12,19,29,30	0
2	NAG	Q	1	14/15	0.99	0.03	12,23,34,35	0
2	NAG	Q	2	14/15	0.99	0.03	16,36,42,44	0
3	NAG	F	1	14/15	0.99	0.03	11,15,24,25	0
3	NAG	F	2	14/15	0.99	0.03	7,18,25,28	0
3	BMA	F	3	11/12	0.99	0.03	12,25,33,34	0
3	MAN	F	4	11/12	0.99	0.03	17,24,33,39	0
3	MAN	F	5	11/12	0.99	0.05	18,26,37,41	0
2	NAG	G	1	14/15	0.99	0.03	11,16,25,28	0
4	NAG	I	1	14/15	0.99	0.03	12,19,27,34	0
4	NAG	I	2	14/15	0.99	0.03	8,22,28,30	0
4	BMA	I	3	11/12	0.99	0.04	9,24,28,40	0
4	MAN	I	4	11/12	0.99	0.03	16,30,35,37	0
4	MAN	I	5	11/12	0.99	0.04	18,29,36,38	0
4	NAG	P	1	14/15	0.99	0.03	17,28,32,33	0
4	NAG	P	2	14/15	0.99	0.04	21,28,39,44	0
2	NAG	E	1	14/15	0.99	0.03	9,16,22,30	0
4	MAN	P	4	11/12	0.99	0.04	24,41,53,54	0
2	NAG	H	1	14/15	0.99	0.03	8,16,23,26	0
5	NAG	L	1	14/15	0.99	0.03	12,20,29,29	0
5	NAG	L	2	14/15	0.99	0.04	19,25,33,38	0
5	BMA	L	3	11/12	0.99	0.04	12,28,36,37	0
5	NAG	O	1	14/15	0.99	0.03	13,18,25,25	0
5	NAG	O	2	14/15	0.99	0.04	15,26,42,54	0
2	NAG	H	2	14/15	0.99	0.03	10,14,25,26	0

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

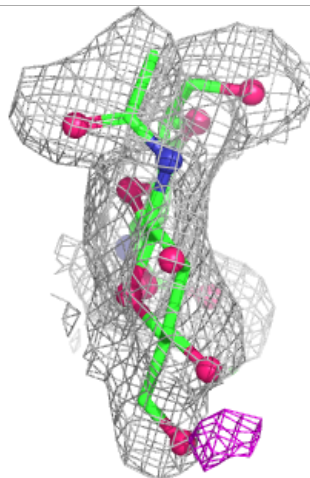
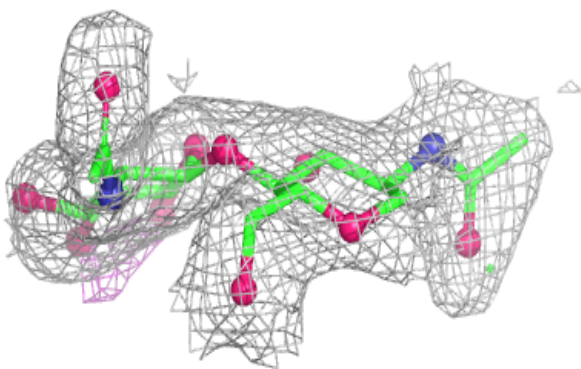
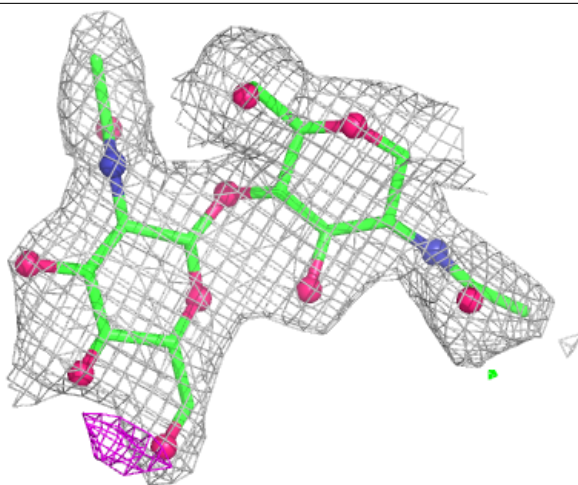
Electron density around Chain E:

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and green (positive)



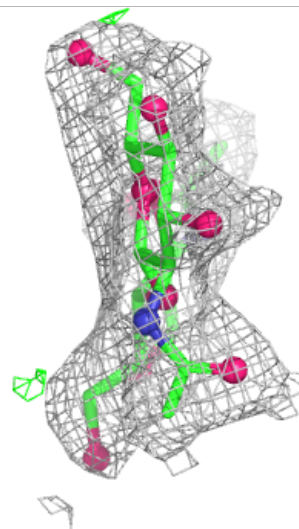
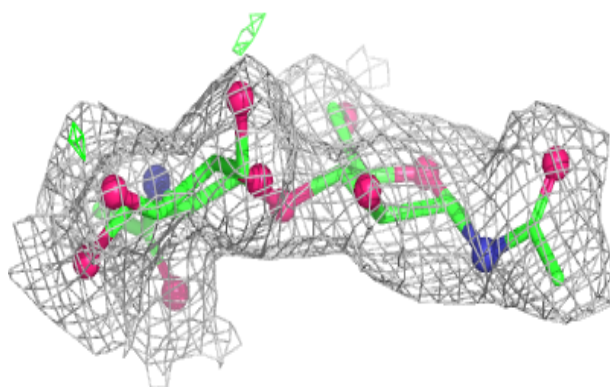
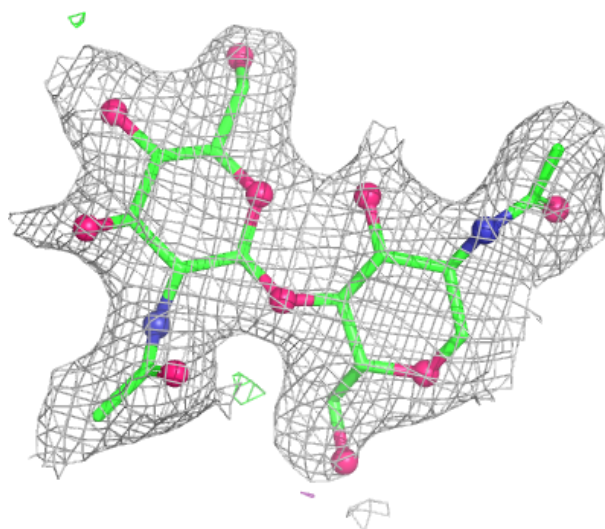
Electron density around Chain G:

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and green (positive)



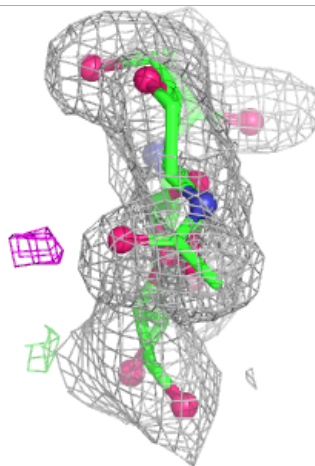
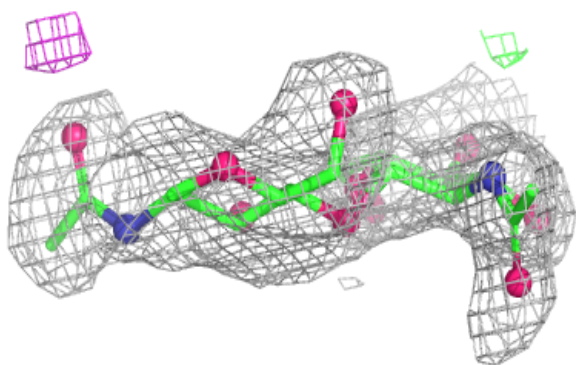
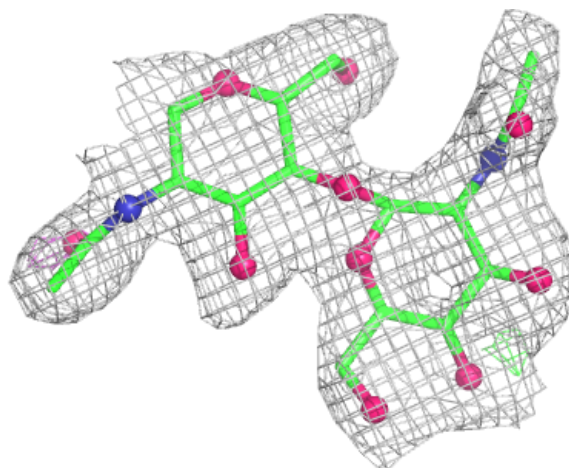
Electron density around Chain H:

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and green (positive)



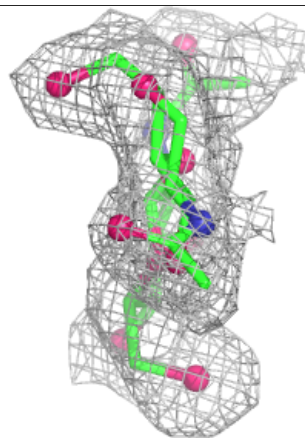
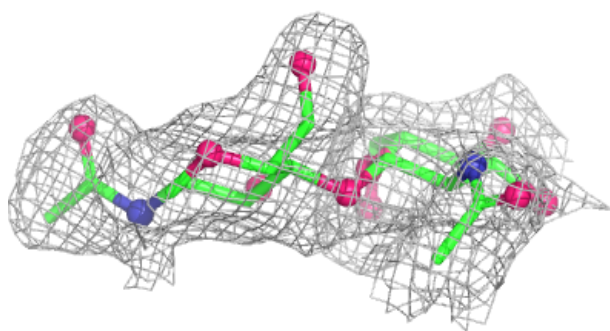
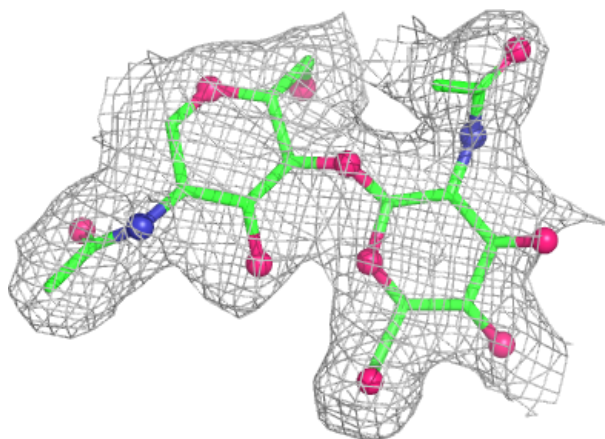
Electron density around Chain J:

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and green (positive)



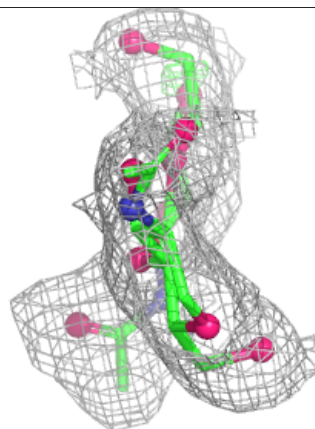
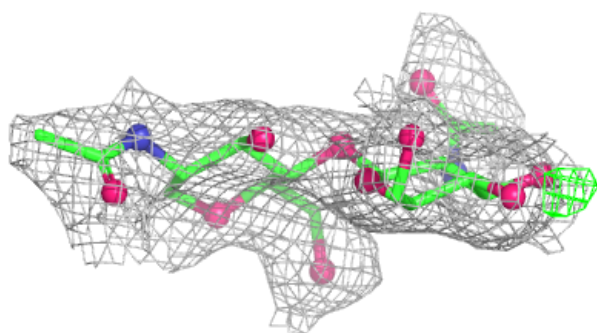
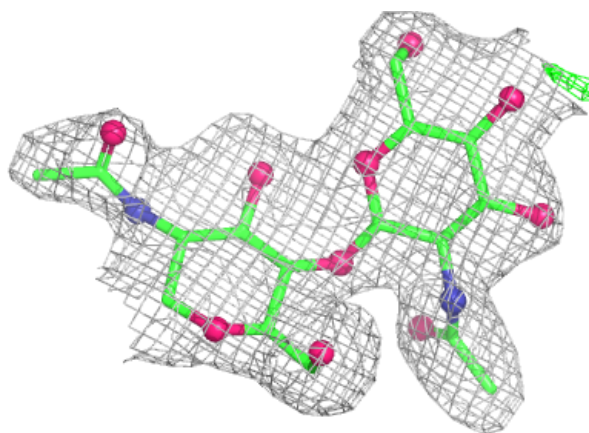
Electron density around Chain K:

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and green (positive)



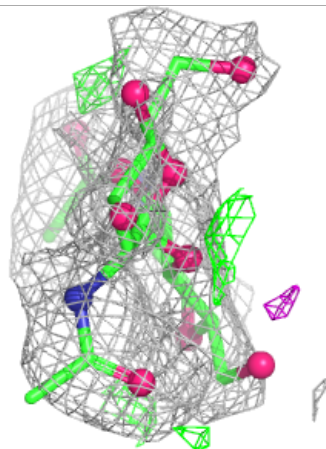
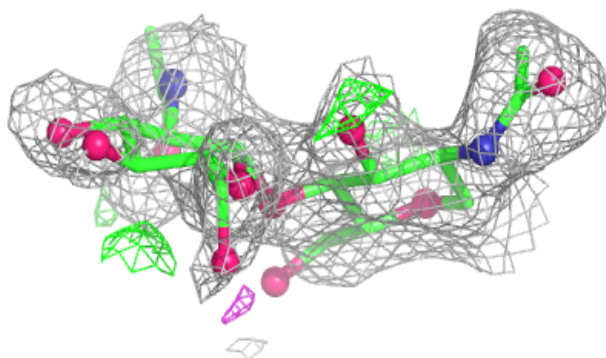
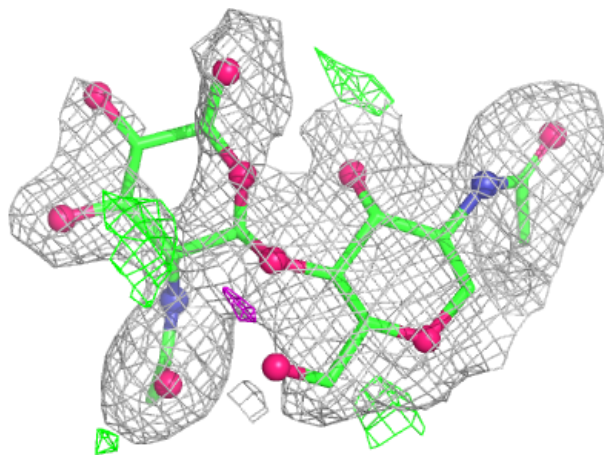
Electron density around Chain M:

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



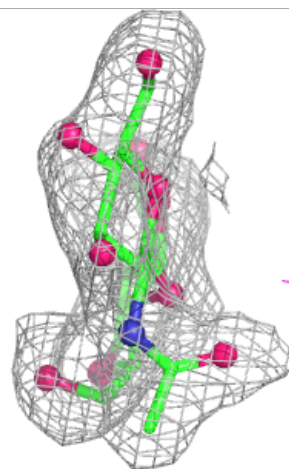
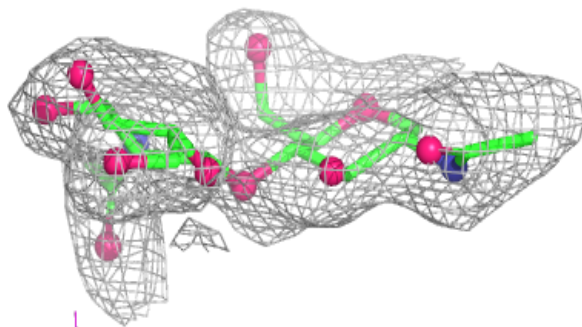
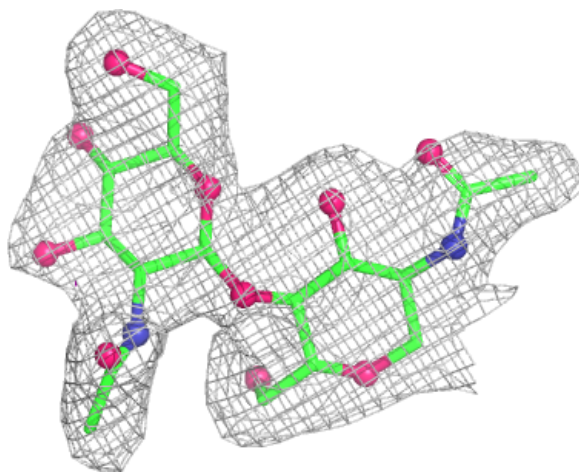
Electron density around Chain N:

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



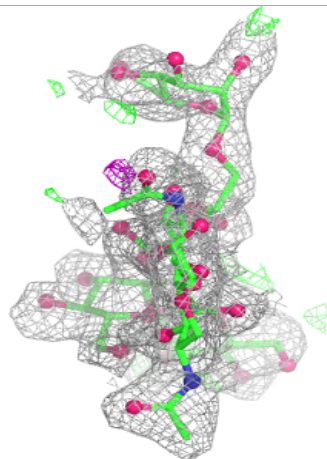
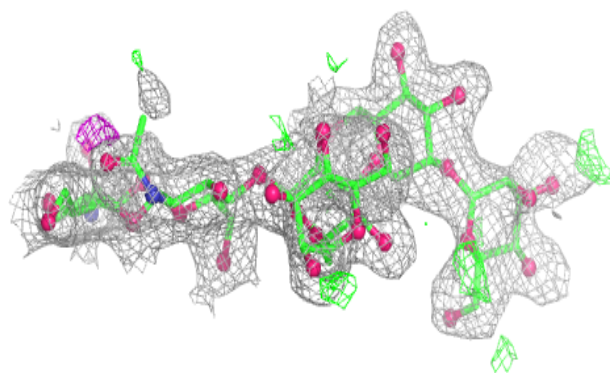
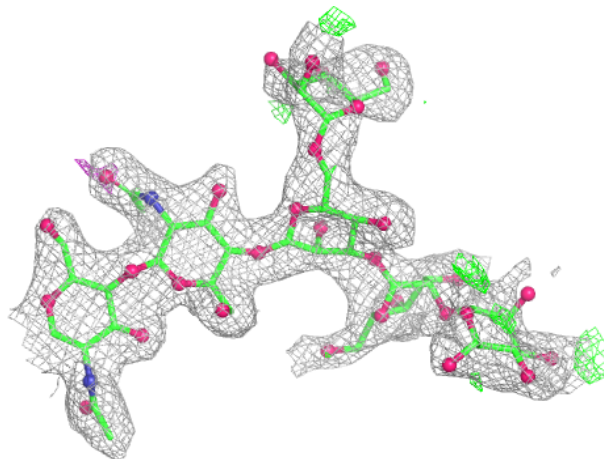
Electron density around Chain Q:

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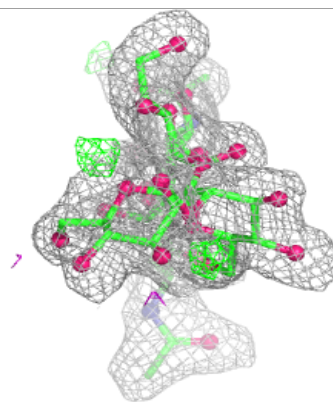
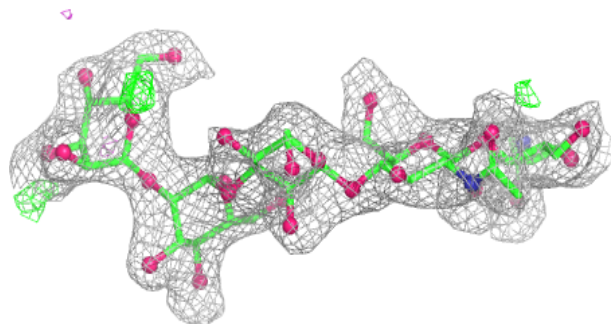
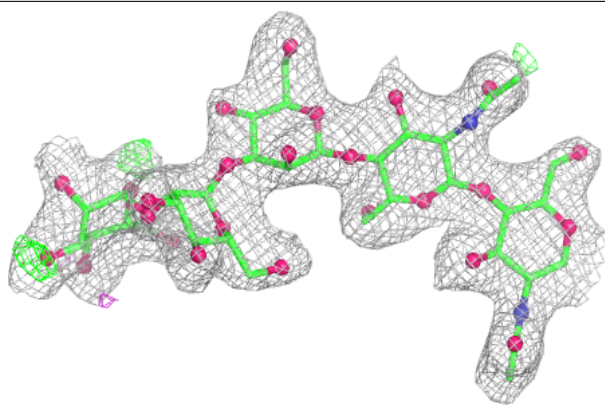
Electron density around Chain F:

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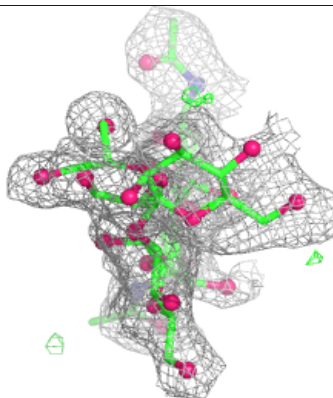
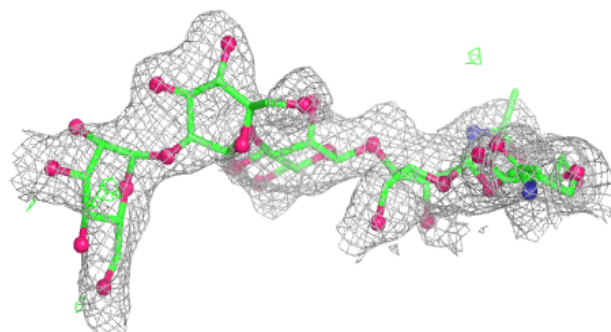
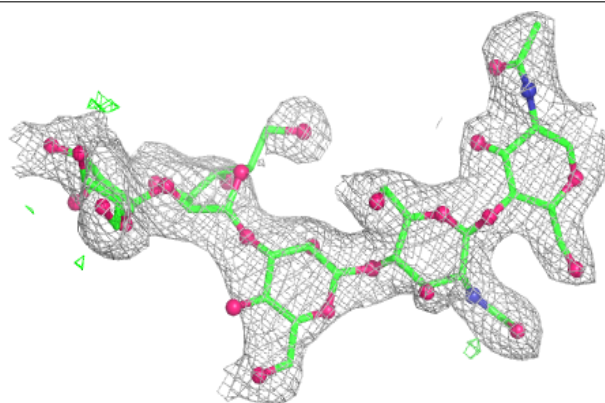


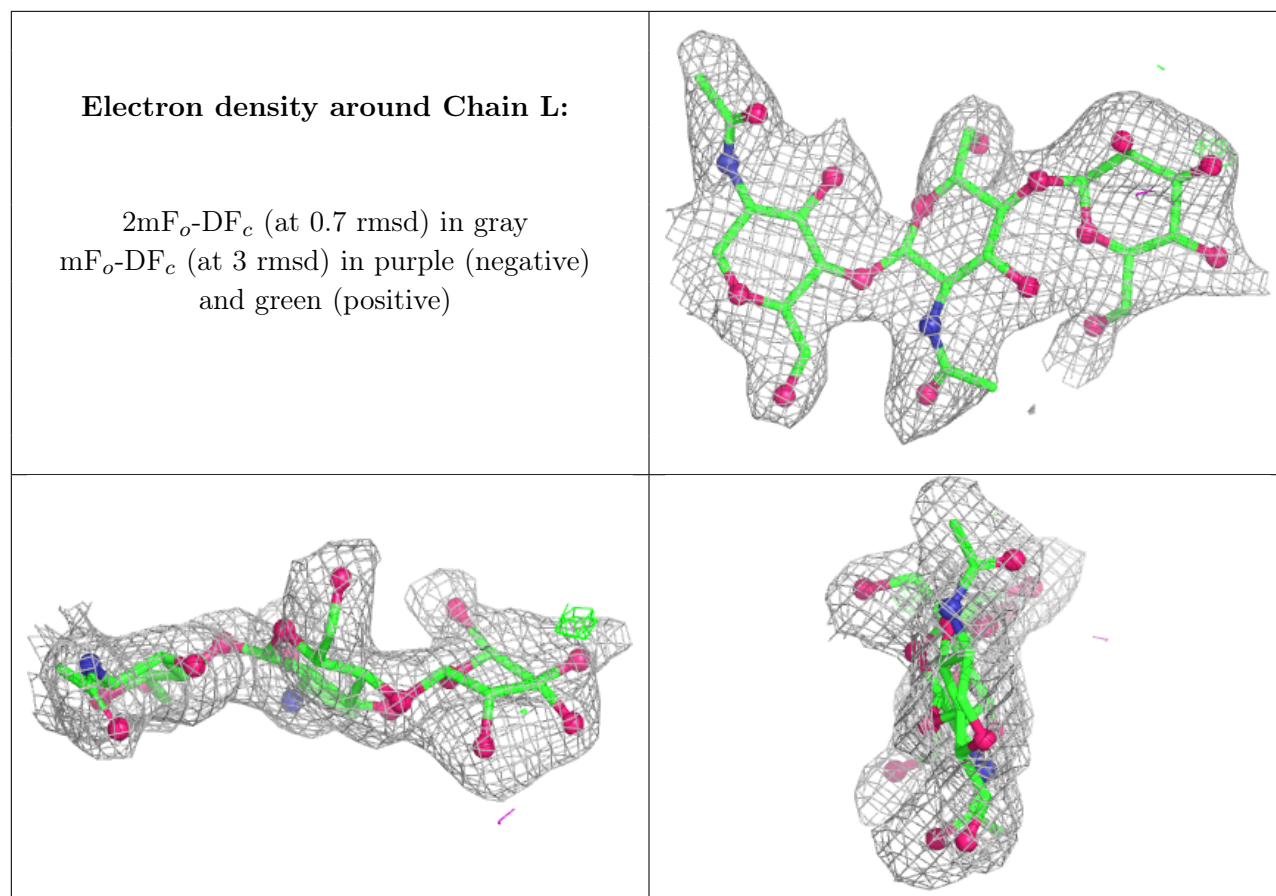
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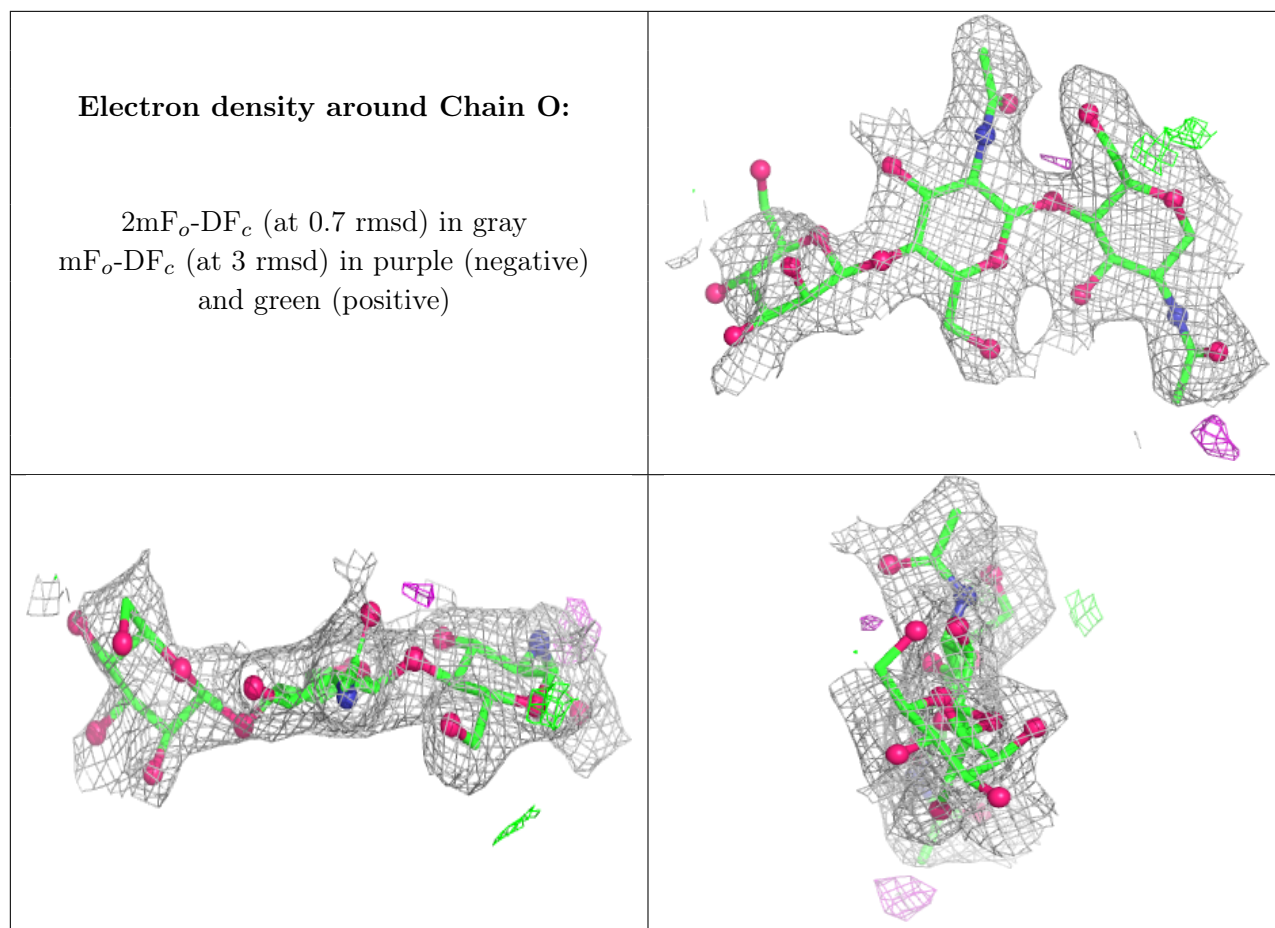
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around Chain P:**

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







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(Å ²)	Q<0.9
8	NAG	C	730	14/15	0.98	0.04	19,29,44,58	0
8	NAG	A	730	14/15	0.99	0.04	13,31,41,56	0
8	NAG	A	760	14/15	0.99	0.04	12,21,25,30	0
8	NAG	B	720	14/15	0.99	0.03	13,23,35,36	0
8	NAG	B	760	14/15	0.99	0.03	6,15,25,28	0
8	NAG	B	770	14/15	0.99	0.03	17,27,34,34	0
8	NAG	A	720	14/15	0.99	0.03	15,20,29,44	0
8	NAG	D	720	14/15	0.99	0.04	14,20,28,32	0
6	CU	C	601	1/1	1.00	0.01	21,21,21,21	0
6	CU	C	602	1/1	1.00	0.01	21,21,21,21	0
6	CU	C	603	1/1	1.00	0.01	18,18,18,18	0
6	CU	C	604	1/1	1.00	0.01	30,30,30,30	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
6	CU	D	601	1/1	1.00	0.01	19,19,19,19	0
6	CU	D	602	1/1	1.00	0.01	26,26,26,26	0
6	CU	D	603	1/1	1.00	0.01	20,20,20,20	0
6	CU	D	604	1/1	1.00	0.01	33,33,33,33	0
7	OXY	A	620	2/2	1.00	0.04	5,5,5,8	0
6	CU	A	601	1/1	1.00	0.01	31,31,31,31	0
6	CU	A	602	1/1	1.00	0.00	11,11,11,11	0
6	CU	A	603	1/1	1.00	0.01	14,14,14,14	0
6	CU	A	604	1/1	1.00	0.01	27,27,27,27	0
6	CU	B	601	1/1	1.00	0.01	21,21,21,21	0
6	CU	B	602	1/1	1.00	0.01	17,17,17,17	0
6	CU	B	603	1/1	1.00	0.01	13,13,13,13	0
6	CU	B	604	1/1	1.00	0.02	26,26,26,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.