



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:31 AM UTC

PDB ID : 7MP4 / pdb_00007mp4
Title : Crystal structure of Epiphyas postvittana antennal carboxylesterase 24
Authors : Hamiaux, C.; Carraher, C.
Deposited on : 2021-05-04
Resolution : 2.43 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)
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

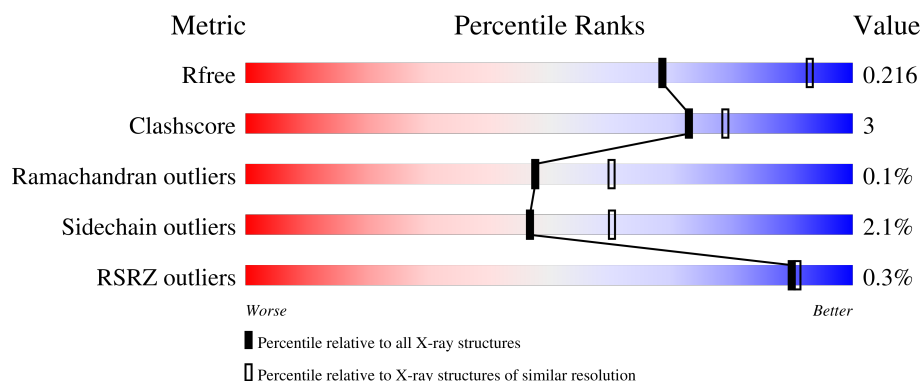
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.43 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	2340 (2.46-2.42)
Clashscore	190562	2400 (2.46-2.42)
Ramachandran outliers	187476	2379 (2.46-2.42)
Sidechain outliers	187428	2379 (2.46-2.42)
RSRZ outliers	180081	2340 (2.46-2.42)

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	556	 88% 8% .
1	B	556	 89% 7% .
1	C	556	 86% 9% . .
1	D	556	 88% 8% . .
1	E	556	 87% 8% . .

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Mol	Chain	Length	Quality of chain
1	F	556	86% 10%
1	G	556	85% 10%
1	H	556	87% 8%
2	I	5	20% 80%
2	J	5	20% 80%
2	K	5	60% 40%
2	L	5	60% 40%
2	M	5	60% 40%
2	N	5	40% 60%
2	O	5	100%
2	P	5	40% 60%

2 Entry composition [i](#)

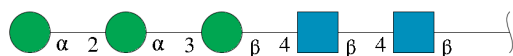
There are 3 unique types of molecules in this entry. The entry contains 34862 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Carboxylesterase-24.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			
1	B	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			
1	C	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			
1	D	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			
1	E	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			
1	F	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			
1	G	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			
1	H	534	Total	C	N	O	S	0	0	0
			4147	2645	708	776	18			

- Molecule 2 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
2	I	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	J	5	Total	C	N	O	0	0	0
			61	34	2	25			
2	K	5	Total	C	N	O	0	0	0
			61	34	2	25			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	L	5	Total 61	C 34	N 2	O 25	0	0	0
2	M	5	Total 61	C 34	N 2	O 25	0	0	0
2	N	5	Total 61	C 34	N 2	O 25	0	0	0
2	O	5	Total 61	C 34	N 2	O 25	0	0	0
2	P	5	Total 61	C 34	N 2	O 25	0	0	0

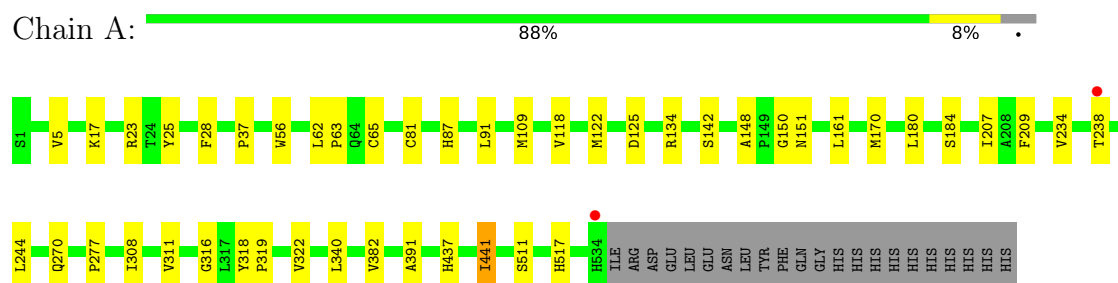
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	214	Total 214	O 214	0	0
3	B	92	Total 92	O 92	0	0
3	C	129	Total 129	O 129	0	0
3	D	149	Total 149	O 149	0	0
3	E	190	Total 190	O 190	0	0
3	F	104	Total 104	O 104	0	0
3	G	166	Total 166	O 166	0	0
3	H	154	Total 154	O 154	0	0

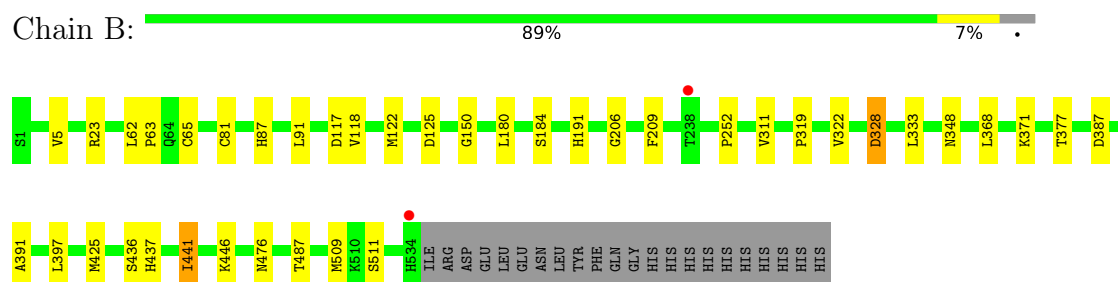
3 Residue-property plots

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.

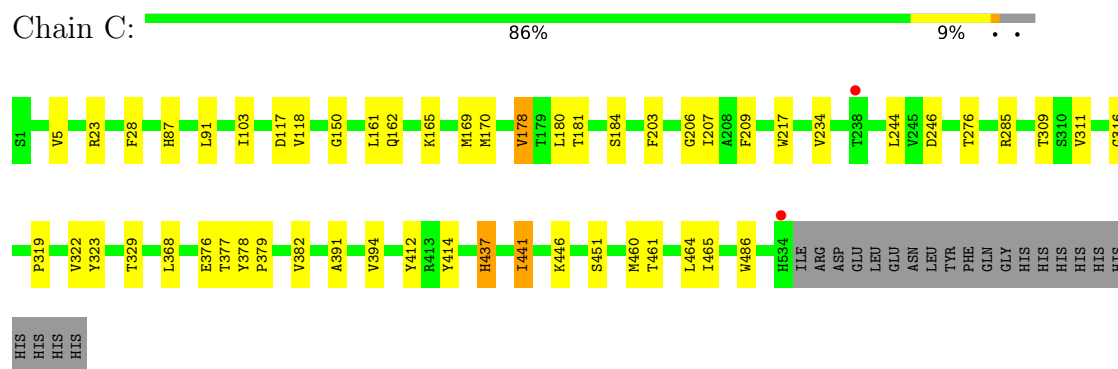
• Molecule 1: Carboxylesterase-24



• Molecule 1: Carboxylesterase-24

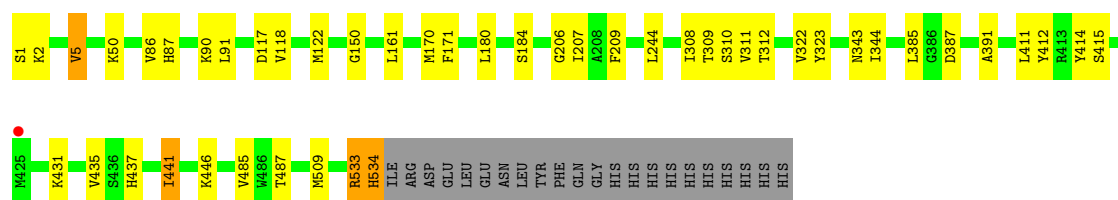


• Molecule 1: Carboxylesterase-24



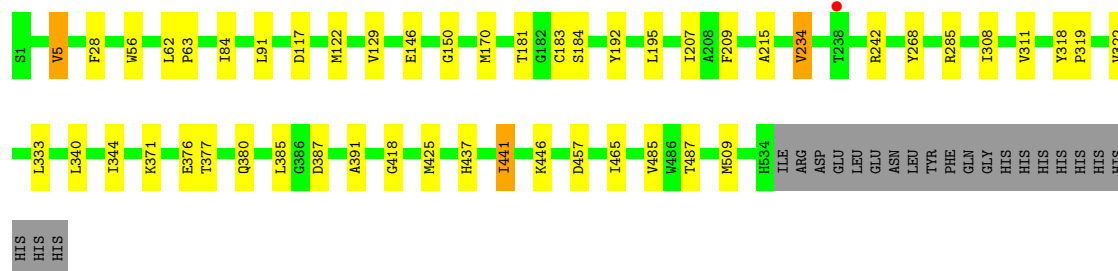
• Molecule 1: Carboxylesterase-24





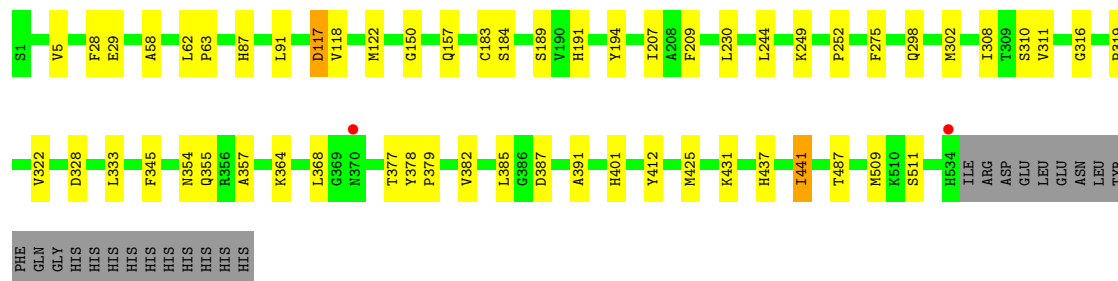
• Molecule 1: Carboxylesterase-24

Chain E: 87% 8% . .



• Molecule 1: Carboxylesterase-24

Chain F: 86% 10% .



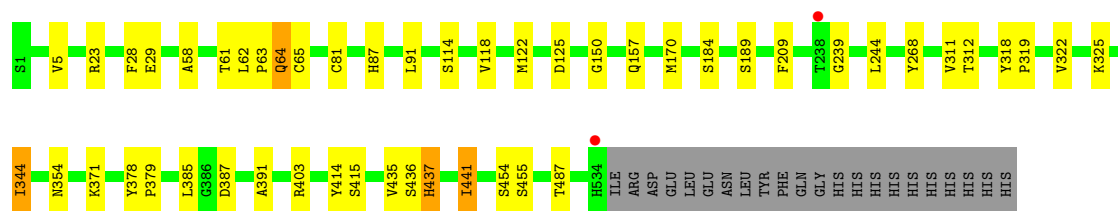
• Molecule 1: Carboxylesterase-24

Chain G: 85% 10% .

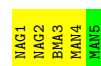


• Molecule 1: Carboxylesterase-24

Chain H: 87% 8% . .



- Molecule 2: 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



- Molecule 2: 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



- Molecule 2: 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



- Molecule 2: 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



- Molecule 2: 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

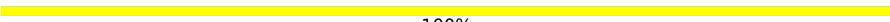


- Molecule 2: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  40% 60%

MAG1	MAG2	BMA3	MAN4	MAN5
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- Molecule 2: 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 O:  100%

MAG1	MAG2	BMA3	MAN4	MAN5
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- Molecule 2: 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:  40% 60%

MAG1	MAG2	BMA3	MAN4	MAN5
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4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	107.47Å 113.97Å 125.77Å 78.31° 70.18° 65.88°	Depositor
Resolution (Å)	48.84 – 2.43 48.84 – 2.43	Depositor EDS
% Data completeness (in resolution range)	98.2 (48.84-2.43) 98.2 (48.84-2.43)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.42Å)	Xtriage
Refinement program	REFMAC 5.8.0267	Depositor
R, R_{free}	0.180 , 0.215 0.186 , 0.216	Depositor DCC
R_{free} test set	9515 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	36.8	Xtriage
Anisotropy	0.215	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 35.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	34862	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.48% 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: NAG, BMA, MAN

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	0.99	0/4266	1.34	3/5808 (0.1%)
1	B	1.00	0/4266	1.39	1/5808 (0.0%)
1	C	0.99	0/4266	1.37	1/5808 (0.0%)
1	D	0.99	0/4266	1.37	3/5808 (0.1%)
1	E	0.98	0/4266	1.35	3/5808 (0.1%)
1	F	1.00	0/4266	1.37	3/5808 (0.1%)
1	G	0.99	0/4266	1.37	4/5808 (0.1%)
1	H	1.00	0/4266	1.38	2/5808 (0.0%)
All	All	0.99	0/34128	1.37	20/46464 (0.0%)

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

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

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	150	GLY	CA-C-O	-5.99	117.99	122.37
1	F	150	GLY	CA-C-O	-5.97	118.01	122.37
1	A	150	GLY	CA-C-O	-5.81	118.27	122.45
1	C	150	GLY	CA-C-O	-5.61	118.41	122.45
1	D	343	ASN	CA-C-N	5.35	127.50	120.60
1	D	343	ASN	C-N-CA	5.35	127.50	120.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	215	ALA	CA-C-N	5.30	127.39	120.28
1	E	215	ALA	C-N-CA	5.30	127.39	120.28
1	H	354	ASN	CB-CA-C	-5.28	99.92	110.42
1	F	357	ALA	CA-C-N	5.20	125.61	119.94
1	F	357	ALA	C-N-CA	5.20	125.61	119.94
1	B	150	GLY	CA-C-O	-5.06	118.67	122.37
1	H	150	GLY	CA-C-O	-5.06	118.68	122.37
1	G	215	ALA	CA-C-N	5.05	127.31	120.38
1	G	215	ALA	C-N-CA	5.05	127.31	120.38
1	A	517	HIS	CA-C-N	5.05	127.35	120.54
1	A	517	HIS	C-N-CA	5.05	127.35	120.54
1	G	145	ASP	CA-C-N	5.04	127.28	120.38
1	G	145	ASP	C-N-CA	5.04	127.28	120.38
1	E	150	GLY	CA-C-O	-5.01	118.71	122.37

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	436	SER	Peptide
1	H	436	SER	Peptide

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	4147	0	4011	26	0
1	B	4147	0	4011	21	0
1	C	4147	0	4011	32	0
1	D	4147	0	4011	29	0
1	E	4147	0	4011	27	0
1	F	4147	0	4011	29	0
1	G	4147	0	4011	25	0
1	H	4147	0	4011	25	0
2	I	61	0	52	1	0
2	J	61	0	52	0	0
2	K	61	0	52	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	L	61	0	52	0	0
2	M	61	0	52	0	0
2	N	61	0	52	0	0
2	O	61	0	52	0	0
2	P	61	0	52	0	0
3	A	214	0	0	0	0
3	B	92	0	0	0	0
3	C	129	0	0	0	0
3	D	149	0	0	1	0
3	E	190	0	0	0	0
3	F	104	0	0	0	0
3	G	166	0	0	0	0
3	H	154	0	0	1	0
All	All	34862	0	32504	212	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 3.

All (212) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:344:ILE:HG22	1:D:385:LEU:HD21	1.62	0.81
1:C:368:LEU:HD23	1:C:377:THR:HB	1.63	0.79
1:E:344:ILE:HG23	1:E:385:LEU:HD21	1.66	0.78
1:C:414:TYR:HE2	1:C:460:MET:HE2	1.50	0.76
1:H:344:ILE:HG23	1:H:385:LEU:HD21	1.68	0.74
1:H:209:PHE:CG	1:H:441:ILE:HD11	2.22	0.73
1:C:209:PHE:CG	1:C:441:ILE:HD11	2.27	0.70
1:B:319:PRO:O	1:B:322:VAL:HG12	1.90	0.69
1:C:184:SER:HB2	1:C:437:HIS:NE2	2.08	0.69
1:E:344:ILE:CG2	1:E:385:LEU:HD21	2.23	0.69
1:D:184:SER:HB2	1:D:437:HIS:NE2	2.09	0.68
1:D:344:ILE:CG2	1:D:385:LEU:HD21	2.23	0.68
1:C:5:VAL:HG23	1:C:28:PHE:CZ	2.28	0.68
1:H:311:VAL:HG11	1:H:391:ALA:HA	1.74	0.68
1:B:209:PHE:CG	1:B:441:ILE:HD11	2.28	0.67
1:F:319:PRO:O	1:F:322:VAL:HG12	1.94	0.67
1:C:460:MET:HE3	1:C:464:LEU:HG	1.77	0.67
1:D:209:PHE:CG	1:D:441:ILE:HD11	2.29	0.66
1:F:209:PHE:CG	1:F:441:ILE:HD11	2.31	0.66
1:H:209:PHE:CD2	1:H:441:ILE:HD11	2.31	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:178:VAL:HG22	1:C:203:PHE:HA	1.78	0.65
1:B:5:VAL:CG2	1:B:91:LEU:HD11	2.28	0.64
1:C:311:VAL:HG11	1:C:391:ALA:HA	1.80	0.64
1:C:414:TYR:CE2	1:C:460:MET:HE2	2.31	0.63
1:B:368:LEU:HD23	1:B:377:THR:HB	1.81	0.62
1:E:5:VAL:HG22	1:E:28:PHE:CE2	2.34	0.62
1:D:487:THR:HB	1:D:509:MET:HE2	1.80	0.62
1:F:311:VAL:HG11	1:F:391:ALA:HA	1.81	0.61
1:H:122:MET:HA	1:H:122:MET:HE2	1.81	0.61
1:A:209:PHE:CD2	1:A:441:ILE:HD11	2.36	0.60
1:A:184:SER:HB2	1:A:437:HIS:NE2	2.16	0.60
1:G:311:VAL:HG11	1:G:391:ALA:HA	1.83	0.60
1:A:311:VAL:HG11	1:A:391:ALA:HA	1.82	0.59
1:E:209:PHE:CD2	1:E:441:ILE:HD11	2.37	0.59
1:A:319:PRO:O	1:A:322:VAL:HG12	2.03	0.59
1:H:170:MET:HA	1:H:170:MET:HE2	1.84	0.59
1:A:238:THR:HG21	2:I:4:MAN:O4	2.02	0.58
1:F:184:SER:HB2	1:F:437:HIS:NE2	2.18	0.58
1:C:169:MET:HB3	1:C:170:MET:HE2	1.86	0.58
1:F:487:THR:HB	1:F:509:MET:HE2	1.85	0.58
1:D:170:MET:HE3	1:D:170:MET:HA	1.86	0.58
1:F:5:VAL:CG2	1:F:91:LEU:HD11	2.34	0.57
1:B:387:ASP:HA	1:B:391:ALA:HB3	1.86	0.56
1:G:209:PHE:CG	1:G:441:ILE:HD11	2.39	0.56
1:E:5:VAL:CG1	1:E:91:LEU:HD11	2.35	0.56
1:D:209:PHE:CD2	1:D:441:ILE:HD11	2.40	0.56
1:B:368:LEU:CD2	1:B:377:THR:HB	2.35	0.56
1:H:344:ILE:CG2	1:H:385:LEU:HD21	2.36	0.55
1:H:62:LEU:HB3	1:H:63:PRO:CD	2.37	0.55
1:D:117:ASP:HB3	1:D:446:LYS:HB2	1.89	0.54
1:C:460:MET:HE3	1:C:464:LEU:CG	2.38	0.54
1:D:534:HIS:ND1	1:D:534:HIS:C	2.63	0.54
1:C:209:PHE:CD2	1:C:441:ILE:HD11	2.42	0.54
1:G:87:HIS:CG	1:G:118:VAL:HG21	2.43	0.54
1:H:268:TYR:CD2	1:H:344:ILE:HD11	2.42	0.54
1:E:5:VAL:HG13	1:E:91:LEU:HD11	1.89	0.54
1:G:184:SER:HB2	1:G:437:HIS:NE2	2.23	0.54
1:B:23:ARG:NH2	1:B:125:ASP:OD1	2.34	0.53
1:D:323:TYR:OH	1:D:344:ILE:HD12	2.08	0.53
1:G:191:HIS:CE1	1:G:397:LEU:HD13	2.44	0.53
1:F:209:PHE:CD2	1:F:441:ILE:HD11	2.43	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:209:PHE:CD2	1:B:441:ILE:HD11	2.44	0.53
1:B:122:MET:HE2	1:B:122:MET:HA	1.91	0.52
1:F:87:HIS:CG	1:F:118:VAL:HG21	2.44	0.52
1:H:184:SER:HB2	1:H:437:HIS:NE2	2.25	0.52
1:E:311:VAL:HG11	1:E:391:ALA:HA	1.90	0.52
1:H:5:VAL:CG2	1:H:91:LEU:HD11	2.40	0.52
1:D:87:HIS:CG	1:D:118:VAL:HG21	2.45	0.52
1:D:5:VAL:CG2	1:D:91:LEU:HD11	2.39	0.52
1:G:161:LEU:HD21	1:G:180:LEU:HD13	1.92	0.52
1:B:487:THR:HB	1:B:509:MET:HE2	1.92	0.51
1:G:418:GLY:HA3	1:G:457:ASP:OD1	2.11	0.51
1:C:368:LEU:CD2	1:C:377:THR:HB	2.36	0.51
1:G:62:LEU:HB3	1:G:63:PRO:CD	2.40	0.51
1:C:319:PRO:O	1:C:322:VAL:HG12	2.10	0.51
1:G:117:ASP:HB3	1:G:446:LYS:HB2	1.93	0.51
1:G:209:PHE:CD2	1:G:441:ILE:HD11	2.45	0.51
1:D:1:SER:N	3:D:602:HOH:O	2.43	0.51
1:E:170:MET:HE3	1:E:170:MET:HA	1.93	0.50
1:E:122:MET:HA	1:E:122:MET:HE2	1.94	0.50
1:A:207:ILE:HA	1:A:308:ILE:O	2.12	0.49
1:G:322:VAL:HG13	1:G:426:MET:HE2	1.94	0.49
1:A:170:MET:HE2	1:A:170:MET:HA	1.93	0.49
1:D:534:HIS:ND1	1:D:534:HIS:O	2.45	0.49
1:F:302:MET:HE2	1:F:401:HIS:ND1	2.27	0.49
1:A:62:LEU:HB3	1:A:63:PRO:CD	2.42	0.49
1:E:377:THR:HA	1:E:380:GLN:OE1	2.12	0.49
1:A:161:LEU:HD21	1:A:180:LEU:HD13	1.95	0.49
1:F:122:MET:HE2	1:F:122:MET:HA	1.93	0.49
1:C:5:VAL:HG22	1:C:91:LEU:HD11	1.94	0.49
1:F:183:CYS:HB2	1:F:441:ILE:HG12	1.95	0.49
1:D:533:ARG:HH11	1:D:533:ARG:CG	2.26	0.49
1:D:161:LEU:HD21	1:D:180:LEU:HD13	1.93	0.48
1:A:5:VAL:HG22	1:A:91:LEU:HD11	1.94	0.48
1:A:5:VAL:CG2	1:A:91:LEU:HD11	2.43	0.48
1:E:209:PHE:CG	1:E:441:ILE:HD11	2.48	0.48
1:B:62:LEU:HB3	1:B:63:PRO:CD	2.44	0.48
1:B:311:VAL:HG11	1:B:391:ALA:HA	1.96	0.48
1:G:5:VAL:HG11	1:G:171:PHE:CE2	2.49	0.48
1:F:378:TYR:HB3	1:F:379:PRO:HD3	1.95	0.47
1:F:191:HIS:O	1:F:194:TYR:HB2	2.14	0.47
1:A:87:HIS:CG	1:A:118:VAL:HG21	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:62:LEU:HB3	1:E:63:PRO:CD	2.45	0.47
1:C:244:LEU:C	1:C:244:LEU:HD23	2.40	0.47
1:D:312:THR:HA	1:D:414:TYR:O	2.15	0.47
1:E:184:SER:HB2	1:E:437:HIS:NE2	2.30	0.47
1:H:23:ARG:NH2	1:H:125:ASP:OD1	2.45	0.47
1:C:87:HIS:CG	1:C:118:VAL:HG21	2.50	0.47
1:H:5:VAL:HG23	1:H:28:PHE:CZ	2.50	0.47
1:E:117:ASP:HB3	1:E:446:LYS:HB2	1.96	0.46
1:F:368:LEU:CD2	1:F:377:THR:HB	2.46	0.46
1:G:65:CYS:HA	1:G:81:CYS:CB	2.46	0.46
1:C:162:GLN:HE22	1:C:165:LYS:HE2	1.79	0.46
1:E:268:TYR:CD2	1:E:344:ILE:HD11	2.51	0.46
1:B:117:ASP:HB3	1:B:446:LYS:HB2	1.97	0.46
1:E:487:THR:HB	1:E:509:MET:HE2	1.96	0.46
1:H:387:ASP:HA	1:H:391:ALA:HB3	1.98	0.46
1:C:309:THR:OG1	1:C:394:VAL:HG13	2.16	0.46
1:G:217:TRP:CZ3	1:G:218:THR:HG22	2.51	0.46
1:B:184:SER:HB2	1:B:437:HIS:NE2	2.31	0.46
1:E:234:VAL:O	1:E:234:VAL:CG1	2.64	0.46
1:A:5:VAL:HG23	1:A:28:PHE:CZ	2.51	0.45
1:B:87:HIS:CG	1:B:118:VAL:HG21	2.51	0.45
1:C:180:LEU:O	1:C:206:GLY:HA2	2.17	0.45
1:D:533:ARG:HH11	1:D:533:ARG:HG3	1.81	0.45
1:E:387:ASP:HA	1:E:391:ALA:HB3	1.98	0.45
1:G:142:SER:HA	1:G:148:ALA:O	2.17	0.45
1:H:87:HIS:CG	1:H:118:VAL:HG21	2.52	0.45
1:E:56:TRP:CH2	1:F:252:PRO:HD3	2.51	0.45
1:E:207:ILE:HA	1:E:308:ILE:O	2.17	0.45
1:D:180:LEU:O	1:D:206:GLY:HA2	2.16	0.45
1:C:378:TYR:HB3	1:C:379:PRO:HD3	1.99	0.45
1:C:117:ASP:HB3	1:C:446:LYS:HB2	1.99	0.45
1:H:157:GLN:OE1	1:H:189:SER:HB3	2.16	0.45
1:A:17:LYS:HB2	1:A:25:TYR:CZ	2.51	0.44
1:B:191:HIS:CD2	1:B:397:LEU:HD13	2.52	0.44
1:G:29:GLU:O	1:G:58:ALA:HA	2.16	0.44
1:C:5:VAL:CG2	1:C:28:PHE:CZ	2.99	0.44
1:H:29:GLU:O	1:H:58:ALA:HA	2.17	0.44
1:H:244:LEU:C	1:H:244:LEU:HD23	2.43	0.44
1:A:122:MET:HE2	1:A:122:MET:HA	1.99	0.44
1:A:65:CYS:HA	1:A:81:CYS:CB	2.48	0.44
1:G:170:MET:HE2	1:G:170:MET:HA	1.98	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:378:TYR:HB3	1:H:379:PRO:HD3	2.00	0.44
1:G:441:ILE:HG22	1:G:442:PHE:CD1	2.53	0.43
1:F:157:GLN:OE1	1:F:189:SER:HB3	2.18	0.43
1:G:302:MET:HE2	1:G:401:HIS:ND1	2.33	0.43
1:A:209:PHE:CG	1:A:441:ILE:HD11	2.53	0.43
1:C:316:GLY:HA3	1:C:382:VAL:O	2.18	0.43
1:E:84:ILE:HD11	1:E:129:VAL:HG11	2.00	0.43
1:A:23:ARG:NH2	1:A:125:ASP:OD1	2.34	0.43
1:A:244:LEU:C	1:A:244:LEU:HD23	2.44	0.43
1:A:318:TYR:HA	1:A:319:PRO:HA	1.85	0.43
1:D:310:SER:HB2	1:D:412:TYR:CE1	2.54	0.43
1:E:318:TYR:HA	1:E:319:PRO:HA	1.84	0.43
1:H:65:CYS:HA	1:H:81:CYS:CB	2.48	0.43
1:C:323:TYR:HA	1:C:329:THR:HG21	2.01	0.43
1:C:412:TYR:HB3	1:C:486:TRP:CZ2	2.54	0.43
1:G:318:TYR:HA	1:G:319:PRO:HA	1.83	0.43
1:E:340:LEU:HB3	1:E:344:ILE:HD13	2.01	0.42
1:F:387:ASP:HA	1:F:391:ALA:HB3	2.00	0.42
1:A:142:SER:HA	1:A:148:ALA:O	2.19	0.42
1:C:161:LEU:HD23	1:C:161:LEU:HA	1.94	0.42
1:D:311:VAL:HG11	1:D:391:ALA:HA	2.01	0.42
1:F:316:GLY:HA3	1:F:382:VAL:O	2.19	0.42
1:D:387:ASP:HA	1:D:391:ALA:HB3	2.01	0.42
1:H:61:THR:HG21	1:H:114:SER:HB2	2.01	0.42
1:H:312:THR:HA	1:H:414:TYR:O	2.20	0.42
1:A:109:MET:HE1	1:A:270:GLN:O	2.19	0.42
1:F:345:PHE:CD1	1:F:385:LEU:HD23	2.54	0.42
1:G:207:ILE:HA	1:G:308:ILE:O	2.19	0.42
1:H:318:TYR:HA	1:H:319:PRO:HA	1.86	0.42
1:D:122:MET:HE2	1:D:122:MET:HA	2.02	0.42
1:F:310:SER:HB2	1:F:412:TYR:CE1	2.54	0.42
1:C:217:TRP:O	1:C:276:THR:HG21	2.20	0.42
1:D:5:VAL:HG23	1:D:91:LEU:HD11	2.02	0.42
1:G:183:CYS:HB2	1:G:441:ILE:HG12	2.01	0.42
1:A:316:GLY:HA3	1:A:382:VAL:O	2.20	0.41
1:F:5:VAL:HG22	1:F:91:LEU:HD11	2.02	0.41
1:F:244:LEU:C	1:F:244:LEU:HD23	2.45	0.41
1:F:5:VAL:HG23	1:F:28:PHE:CZ	2.55	0.41
1:H:415:SER:C	1:H:435:VAL:HG21	2.45	0.41
1:F:207:ILE:HA	1:F:308:ILE:O	2.20	0.41
1:C:181:THR:HG22	1:C:207:ILE:HB	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:162:GLN:NE2	1:G:202:ASN:HD21	2.18	0.41
1:G:390:PHE:O	1:G:394:VAL:HG23	2.20	0.41
1:D:207:ILE:HA	1:D:308:ILE:O	2.21	0.41
1:G:35:ARG:HB3	1:G:47:GLN:HB2	2.01	0.41
1:C:103:ILE:HD11	1:C:180:LEU:HD11	2.02	0.41
1:B:191:HIS:NE2	1:B:397:LEU:HD13	2.36	0.41
1:D:244:LEU:C	1:D:244:LEU:HD23	2.46	0.41
1:H:64:GLN:NE2	3:H:614:HOH:O	2.53	0.41
1:B:180:LEU:O	1:B:206:GLY:HA2	2.21	0.41
1:B:328:ASP:OD1	1:B:328:ASP:N	2.53	0.41
1:E:181:THR:HA	1:E:207:ILE:O	2.21	0.41
1:E:183:CYS:HB2	1:E:441:ILE:HG13	2.03	0.41
1:F:29:GLU:O	1:F:58:ALA:HA	2.20	0.41
1:A:37:PRO:CG	1:A:134:ARG:HD3	2.51	0.40
1:A:151:ASN:OD1	1:A:277:PRO:HA	2.21	0.40
1:C:234:VAL:O	1:C:234:VAL:CG1	2.68	0.40
1:D:415:SER:C	1:D:435:VAL:HG21	2.47	0.40
1:E:192:TYR:HA	1:E:195:LEU:HD12	2.03	0.40
1:E:418:GLY:HA3	1:E:457:ASP:OD1	2.21	0.40
1:F:62:LEU:HB3	1:F:63:PRO:HD2	2.03	0.40
1:B:65:CYS:HA	1:B:81:CYS:CB	2.51	0.40
1:C:461:THR:O	1:C:465:ILE:HG12	2.21	0.40
1:F:117:ASP:OD1	1:F:117:ASP:C	2.63	0.40
1:F:364:LYS:HA	1:F:368:LEU:HD12	2.03	0.40
1:D:86:VAL:HG11	1:D:171:PHE:CG	2.57	0.40
1:F:230:LEU:HD22	1:F:275:PHE:CZ	2.56	0.40
1:A:56:TRP:CH2	1:B:252:PRO:HD3	2.57	0.40
1:D:309:THR:O	1:D:411:LEU:HA	2.22	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	532/556 (96%)	514 (97%)	18 (3%)	0	100	100
1	B	532/556 (96%)	511 (96%)	21 (4%)	0	100	100
1	C	532/556 (96%)	509 (96%)	22 (4%)	1 (0%)	43	53
1	D	532/556 (96%)	514 (97%)	18 (3%)	0	100	100
1	E	532/556 (96%)	510 (96%)	22 (4%)	0	100	100
1	F	532/556 (96%)	511 (96%)	21 (4%)	0	100	100
1	G	532/556 (96%)	512 (96%)	19 (4%)	1 (0%)	43	53
1	H	532/556 (96%)	513 (96%)	17 (3%)	2 (0%)	30	36
All	All	4256/4448 (96%)	4094 (96%)	158 (4%)	4 (0%)	48	60

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	437	HIS
1	H	239	GLY
1	H	437	HIS
1	G	239	GLY

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	444/465 (96%)	440 (99%)	4 (1%)	70	80
1	B	444/465 (96%)	436 (98%)	8 (2%)	51	64
1	C	444/465 (96%)	437 (98%)	7 (2%)	55	68
1	D	444/465 (96%)	434 (98%)	10 (2%)	44	57
1	E	444/465 (96%)	431 (97%)	13 (3%)	37	49
1	F	444/465 (96%)	433 (98%)	11 (2%)	42	54
1	G	444/465 (96%)	431 (97%)	13 (3%)	37	49
1	H	444/465 (96%)	434 (98%)	10 (2%)	44	57
All	All	3552/3720 (96%)	3476 (98%)	76 (2%)	47	60

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	234	VAL
1	A	340	LEU
1	A	441	ILE
1	A	511	SER
1	B	328	ASP
1	B	333	LEU
1	B	348	ASN
1	B	371	LYS
1	B	425	MET
1	B	441	ILE
1	B	476	ASN
1	B	511	SER
1	C	23	ARG
1	C	178	VAL
1	C	246	ASP
1	C	285	ARG
1	C	376	GLU
1	C	441	ILE
1	C	451	SER
1	D	2	LYS
1	D	5	VAL
1	D	50	LYS
1	D	90	LYS
1	D	322	VAL
1	D	431	LYS
1	D	441	ILE
1	D	485	VAL
1	D	533	ARG
1	D	534	HIS
1	E	5	VAL
1	E	146	GLU
1	E	234	VAL
1	E	242	ARG
1	E	285	ARG
1	E	322	VAL
1	E	333	LEU
1	E	371	LYS
1	E	376	GLU
1	E	425	MET
1	E	441	ILE
1	E	465	ILE
1	E	485	VAL

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Mol	Chain	Res	Type
1	F	117	ASP
1	F	249	LYS
1	F	298	GLN
1	F	328	ASP
1	F	333	LEU
1	F	354	ASN
1	F	355	GLN
1	F	425	MET
1	F	431	LYS
1	F	441	ILE
1	F	511	SER
1	G	4	VAL
1	G	169	MET
1	G	322	VAL
1	G	333	LEU
1	G	354	ASN
1	G	403	ARG
1	G	428	SER
1	G	431	LYS
1	G	441	ILE
1	G	451	SER
1	G	459	ARG
1	G	531	ASN
1	G	534	HIS
1	H	64	GLN
1	H	322	VAL
1	H	325	LYS
1	H	344	ILE
1	H	371	LYS
1	H	403	ARG
1	H	441	ILE
1	H	454	SER
1	H	455	SER
1	H	487	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (53) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	162	GLN
1	A	226	ASN
1	A	336	ASN

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Mol	Chain	Res	Type
1	A	365	GLN
1	A	404	HIS
1	A	407	GLN
1	B	64	GLN
1	B	73	GLN
1	B	162	GLN
1	B	303	HIS
1	B	336	ASN
1	B	355	GLN
1	B	407	GLN
1	C	48	HIS
1	C	64	GLN
1	C	73	GLN
1	C	80	ASN
1	C	162	GLN
1	C	191	HIS
1	C	269	GLN
1	C	355	GLN
1	C	383	GLN
1	C	531	ASN
1	D	219	HIS
1	D	298	GLN
1	D	336	ASN
1	D	339	GLN
1	D	481	ASN
1	E	13	GLN
1	E	21	HIS
1	E	47	GLN
1	E	48	HIS
1	F	48	HIS
1	F	73	GLN
1	F	219	HIS
1	F	336	ASN
1	F	365	GLN
1	F	529	ASN
1	G	191	HIS
1	G	202	ASN
1	G	226	ASN
1	G	282	GLN
1	G	355	GLN
1	G	404	HIS
1	G	407	GLN

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Mol	Chain	Res	Type
1	H	191	HIS
1	H	336	ASN
1	H	365	GLN
1	H	404	HIS
1	H	407	GLN
1	H	476	ASN
1	H	517	HIS

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 ⓘ

40 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	I	1	2,1	14,14,15	0.47	0	17,19,21	0.82	1 (5%)
2	NAG	I	2	2	14,14,15	0.42	0	17,19,21	0.93	1 (5%)
2	BMA	I	3	2	11,11,12	0.79	0	15,15,17	1.11	2 (13%)
2	MAN	I	4	2	11,11,12	0.51	0	15,15,17	0.93	0
2	MAN	I	5	2	11,11,12	0.44	0	15,15,17	0.77	0
2	NAG	J	1	2,1	14,14,15	0.51	0	17,19,21	1.48	3 (17%)
2	NAG	J	2	2	14,14,15	0.43	0	17,19,21	0.94	1 (5%)
2	BMA	J	3	2	11,11,12	0.68	0	15,15,17	1.07	1 (6%)
2	MAN	J	4	2	11,11,12	0.37	0	15,15,17	0.95	1 (6%)
2	MAN	J	5	2	11,11,12	0.47	0	15,15,17	0.62	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	K	1	2,1	14,14,15	0.40	0	17,19,21	0.90	1 (5%)
2	NAG	K	2	2	14,14,15	0.39	0	17,19,21	0.98	2 (11%)
2	BMA	K	3	2	11,11,12	0.69	0	15,15,17	1.01	0
2	MAN	K	4	2	11,11,12	0.35	0	15,15,17	0.74	0
2	MAN	K	5	2	11,11,12	0.48	0	15,15,17	0.76	0
2	NAG	L	1	2,1	14,14,15	0.42	0	17,19,21	0.84	1 (5%)
2	NAG	L	2	2	14,14,15	0.36	0	17,19,21	0.83	0
2	BMA	L	3	2	11,11,12	0.40	0	15,15,17	0.83	0
2	MAN	L	4	2	11,11,12	0.43	0	15,15,17	0.91	1 (6%)
2	MAN	L	5	2	11,11,12	0.33	0	15,15,17	0.45	0
2	NAG	M	1	2,1	14,14,15	0.42	0	17,19,21	0.76	0
2	NAG	M	2	2	14,14,15	0.39	0	17,19,21	1.07	2 (11%)
2	BMA	M	3	2	11,11,12	0.71	0	15,15,17	0.90	1 (6%)
2	MAN	M	4	2	11,11,12	0.44	0	15,15,17	0.63	0
2	MAN	M	5	2	11,11,12	0.45	0	15,15,17	0.87	0
2	NAG	N	1	2,1	14,14,15	0.40	0	17,19,21	0.85	1 (5%)
2	NAG	N	2	2	14,14,15	0.39	0	17,19,21	1.05	2 (11%)
2	BMA	N	3	2	11,11,12	0.52	0	15,15,17	0.86	1 (6%)
2	MAN	N	4	2	11,11,12	0.44	0	15,15,17	0.77	0
2	MAN	N	5	2	11,11,12	0.40	0	15,15,17	0.57	0
2	NAG	O	1	2,1	14,14,15	0.40	0	17,19,21	0.92	1 (5%)
2	NAG	O	2	2	14,14,15	0.35	0	17,19,21	0.94	2 (11%)
2	BMA	O	3	2	11,11,12	0.60	0	15,15,17	1.13	1 (6%)
2	MAN	O	4	2	11,11,12	0.43	0	15,15,17	0.86	1 (6%)
2	MAN	O	5	2	11,11,12	0.29	0	15,15,17	0.96	1 (6%)
2	NAG	P	1	2,1	14,14,15	0.43	0	17,19,21	0.61	0
2	NAG	P	2	2	14,14,15	0.40	0	17,19,21	0.76	1 (5%)
2	BMA	P	3	2	11,11,12	0.45	0	15,15,17	0.70	0
2	MAN	P	4	2	11,11,12	0.38	0	15,15,17	0.91	1 (6%)
2	MAN	P	5	2	11,11,12	0.52	0	15,15,17	0.93	1 (6%)

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	I	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1

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

There are no bond length outliers.

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	J	1	NAG	O5-C1-C2	-4.05	105.02	111.29
2	O	3	BMA	C1-O5-C5	3.36	116.69	112.19
2	J	2	NAG	C1-C2-N2	2.91	115.02	110.43
2	M	3	BMA	C1-O5-C5	2.89	116.06	112.19
2	M	2	NAG	C1-O5-C5	2.82	115.97	112.19
2	N	2	NAG	C1-C2-N2	2.79	114.82	110.43
2	N	2	NAG	C1-O5-C5	2.69	115.79	112.19
2	M	2	NAG	C1-C2-N2	2.68	114.66	110.43
2	N	3	BMA	C1-O5-C5	2.58	115.64	112.19
2	K	2	NAG	C1-C2-N2	2.56	114.47	110.43
2	L	4	MAN	C1-C2-C3	2.55	113.36	109.64
2	J	1	NAG	C1-C2-N2	2.47	114.33	110.43
2	O	2	NAG	C1-C2-N2	2.45	114.29	110.43
2	I	3	BMA	C1-O5-C5	2.42	115.42	112.19
2	I	3	BMA	C3-C4-C5	2.40	114.59	110.23
2	I	1	NAG	O5-C1-C2	-2.40	107.58	111.29
2	J	4	MAN	C1-O5-C5	2.38	115.38	112.19
2	O	2	NAG	C1-O5-C5	2.36	115.35	112.19
2	O	4	MAN	C1-C2-C3	2.35	113.06	109.64
2	J	1	NAG	C2-N2-C7	2.34	126.04	122.90
2	J	3	BMA	C3-C4-C5	2.30	114.40	110.23
2	O	5	MAN	O2-C2-C3	2.29	114.90	110.15
2	K	2	NAG	C1-O5-C5	2.22	115.16	112.19
2	P	2	NAG	C1-O5-C5	2.20	115.13	112.19
2	I	2	NAG	C1-O5-C5	2.19	115.12	112.19
2	P	4	MAN	C1-O5-C5	2.18	115.11	112.19
2	K	1	NAG	O5-C1-C2	-2.18	107.92	111.29
2	P	5	MAN	C3-C4-C5	-2.13	106.38	110.23
2	L	1	NAG	C1-C2-N2	2.09	113.72	110.43
2	O	1	NAG	O5-C1-C2	-2.07	108.09	111.29
2	N	1	NAG	C2-N2-C7	2.07	125.68	122.90

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	I	3	BMA	O5-C5-C6-O6
2	L	3	BMA	O5-C5-C6-O6
2	L	3	BMA	C4-C5-C6-O6
2	L	2	NAG	O5-C5-C6-O6
2	M	3	BMA	C4-C5-C6-O6
2	P	5	MAN	C4-C5-C6-O6
2	I	3	BMA	C4-C5-C6-O6

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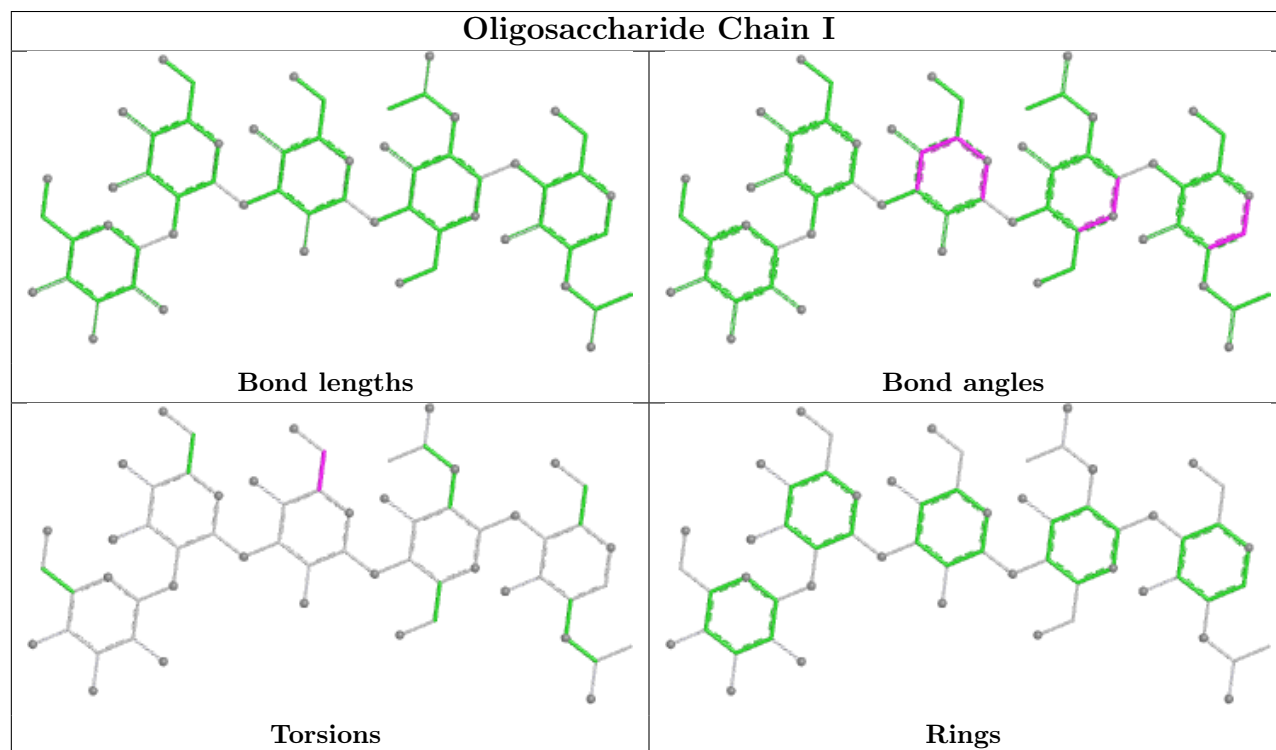
Mol	Chain	Res	Type	Atoms
2	P	5	MAN	O5-C5-C6-O6
2	M	3	BMA	O5-C5-C6-O6
2	N	3	BMA	O5-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	J	3	BMA	O5-C5-C6-O6
2	K	3	BMA	O5-C5-C6-O6
2	L	2	NAG	C4-C5-C6-O6
2	L	2	NAG	C8-C7-N2-C2
2	P	4	MAN	O5-C5-C6-O6
2	L	2	NAG	O7-C7-N2-C2
2	K	1	NAG	C1-C2-N2-C7
2	P	1	NAG	C1-C2-N2-C7

There are no ring outliers.

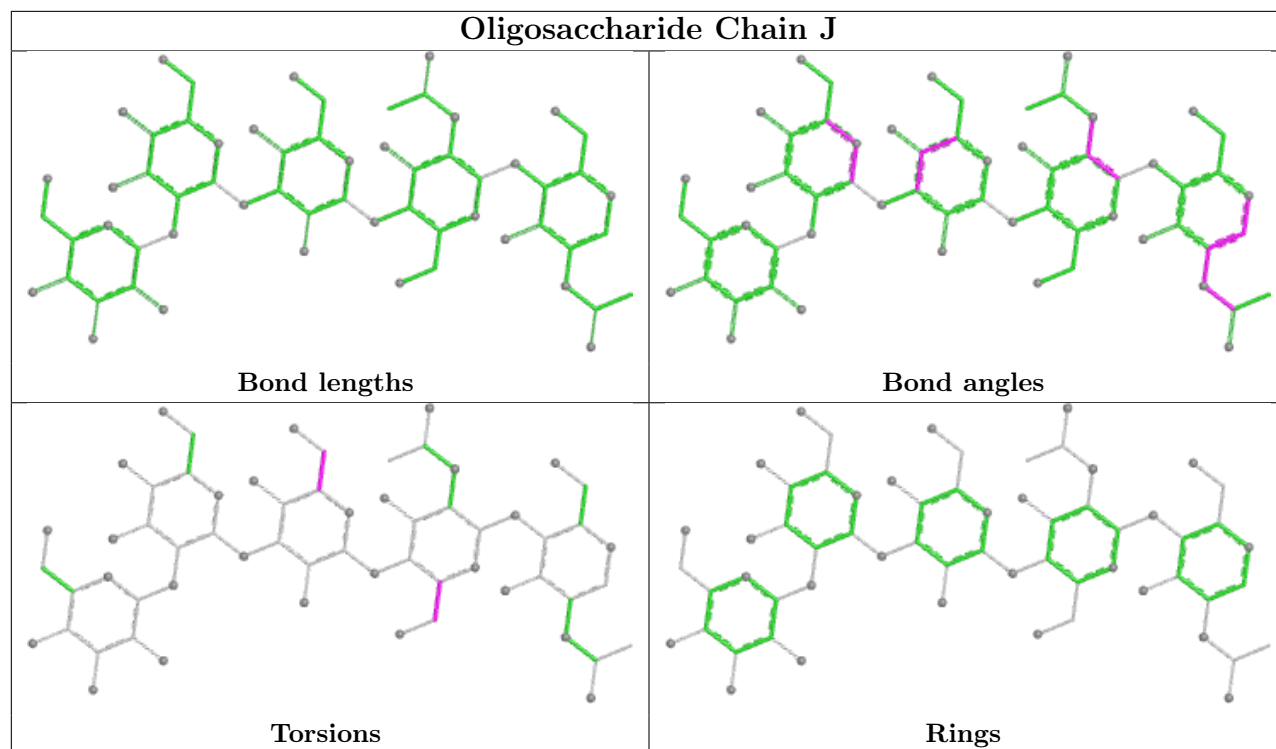
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	4	MAN	1	0

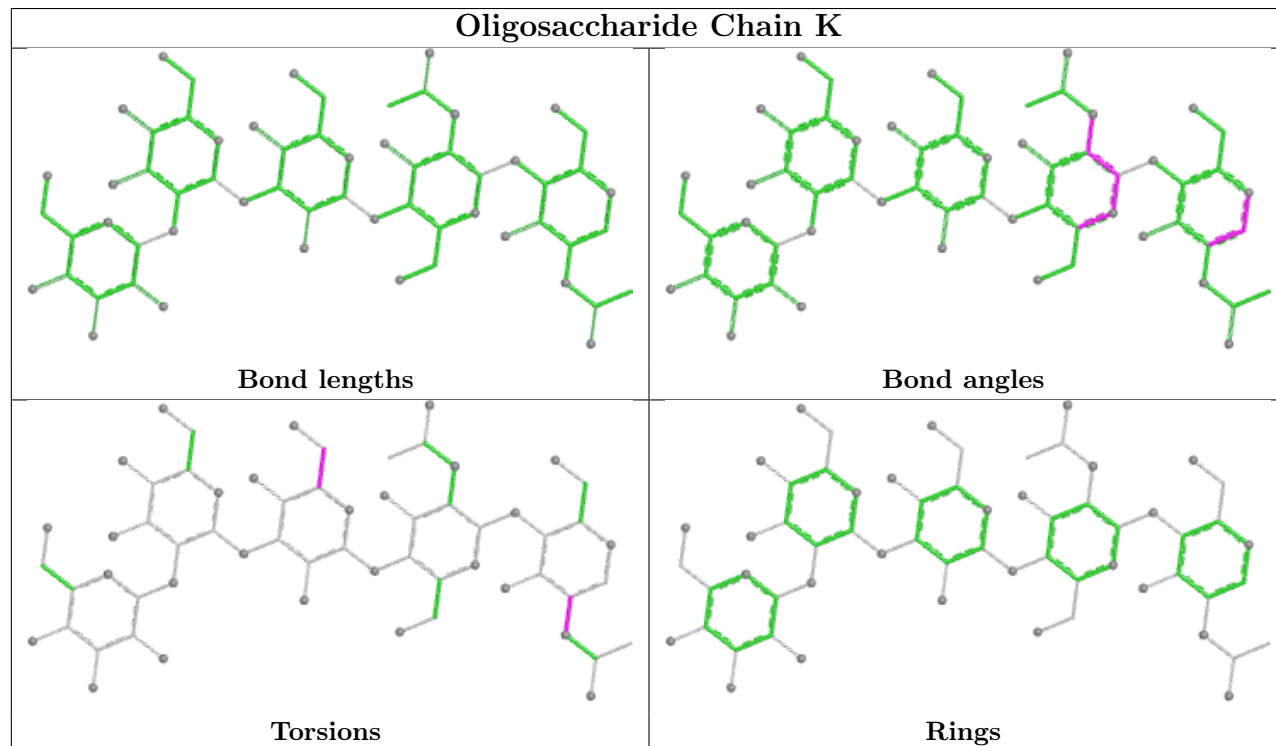
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

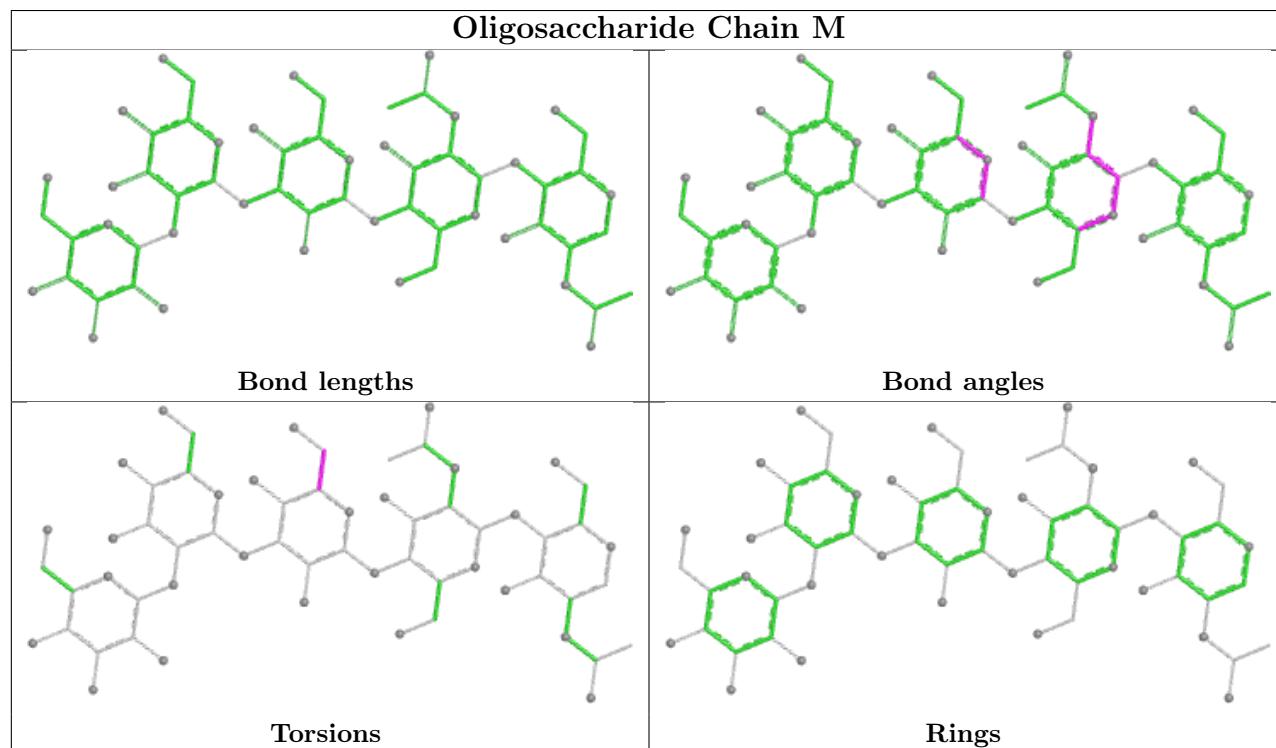
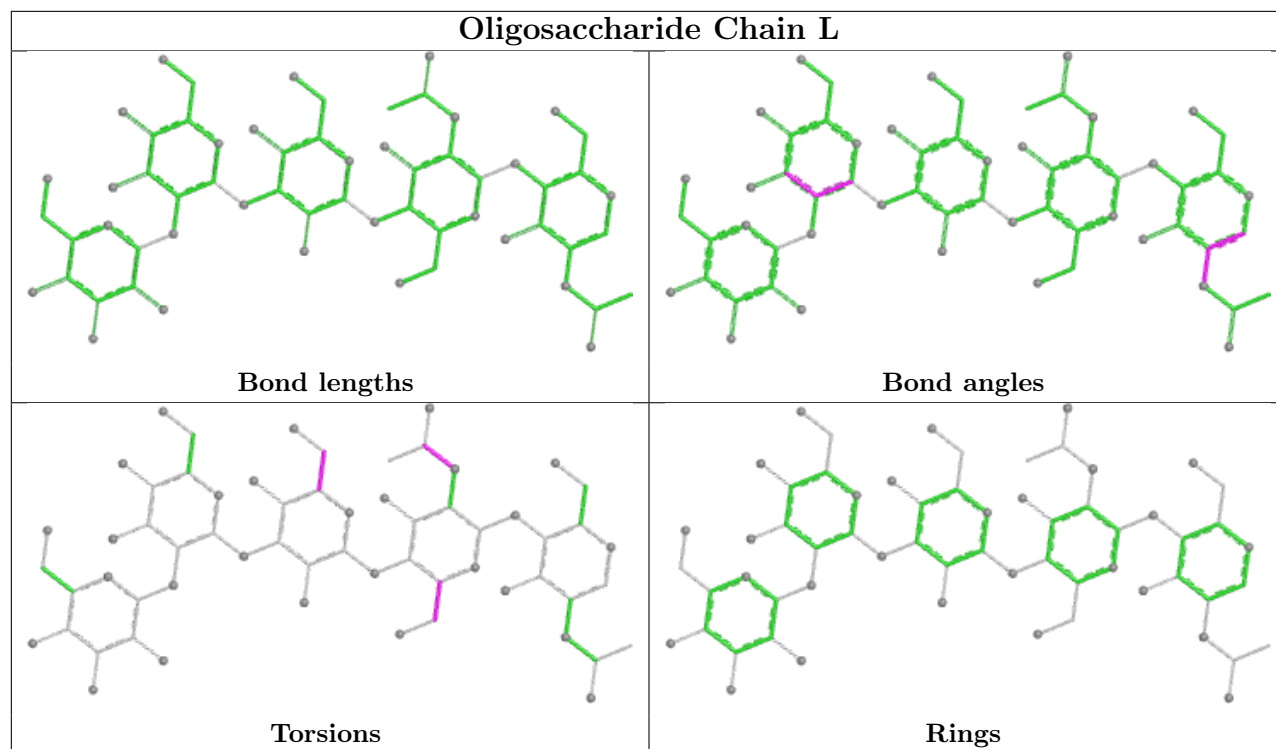


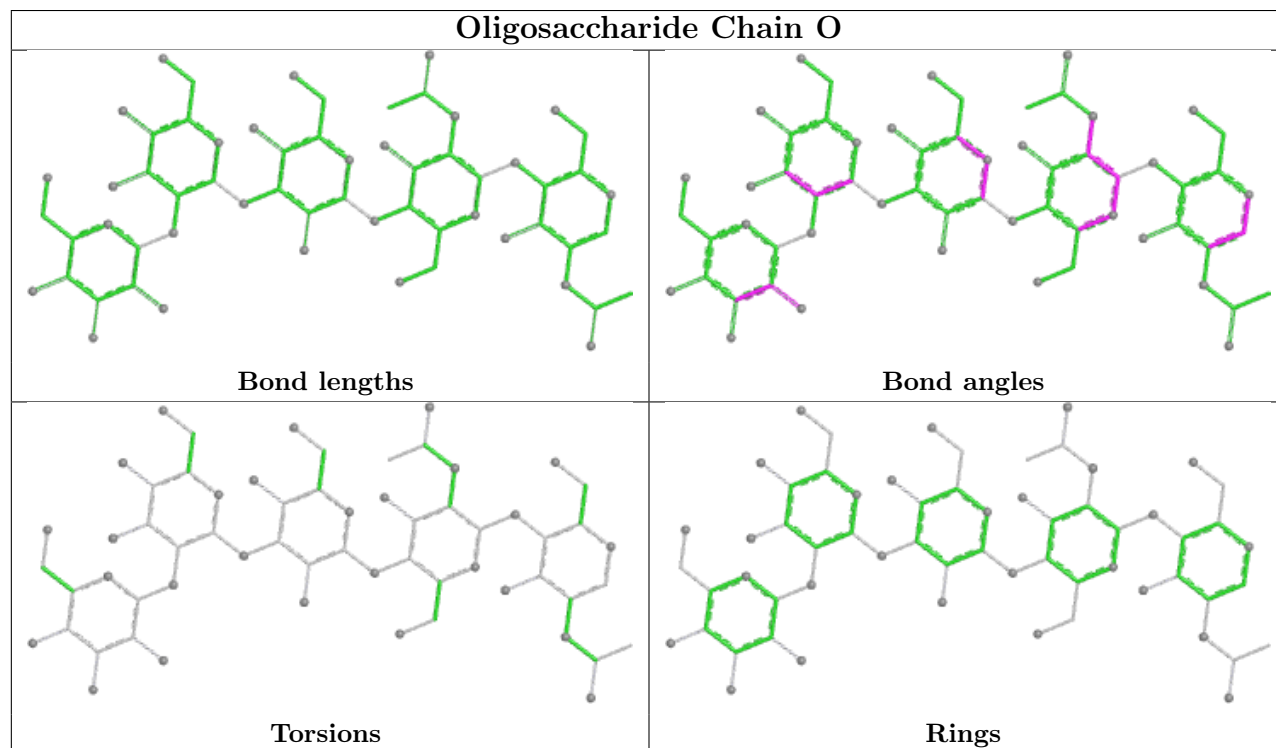
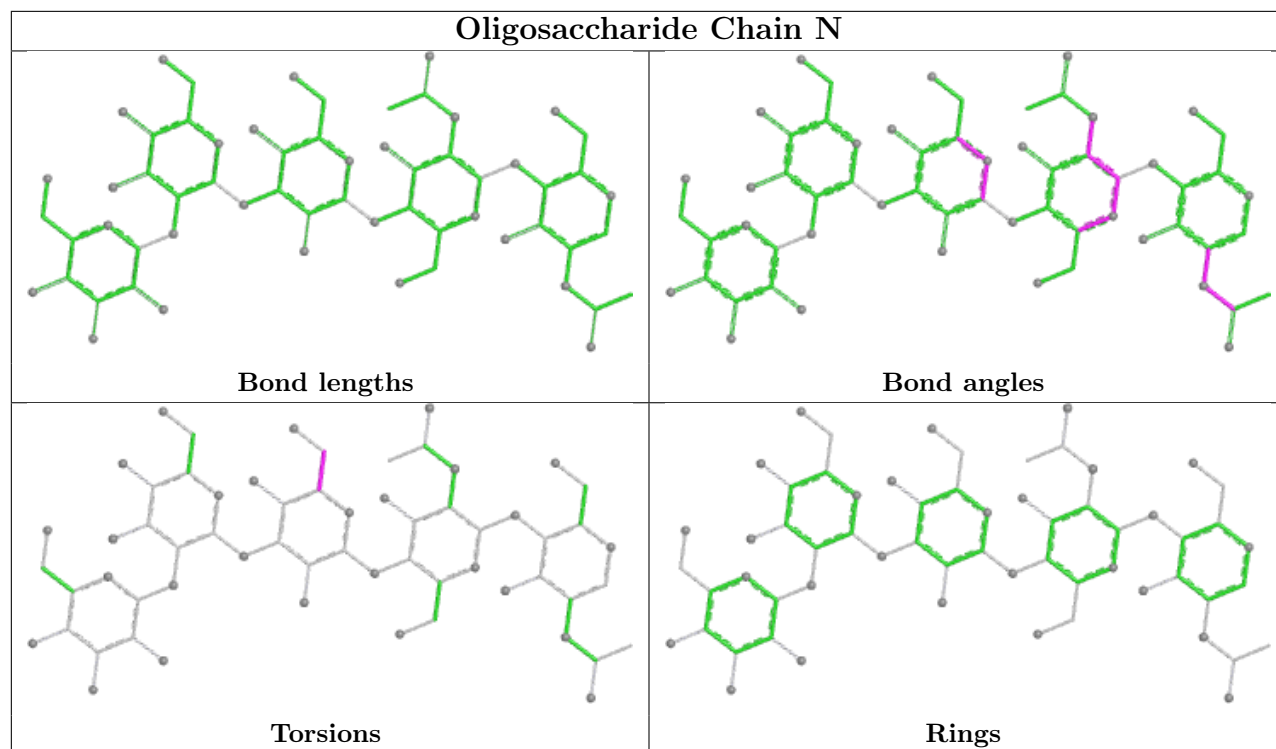
Oligosaccharide Chain J

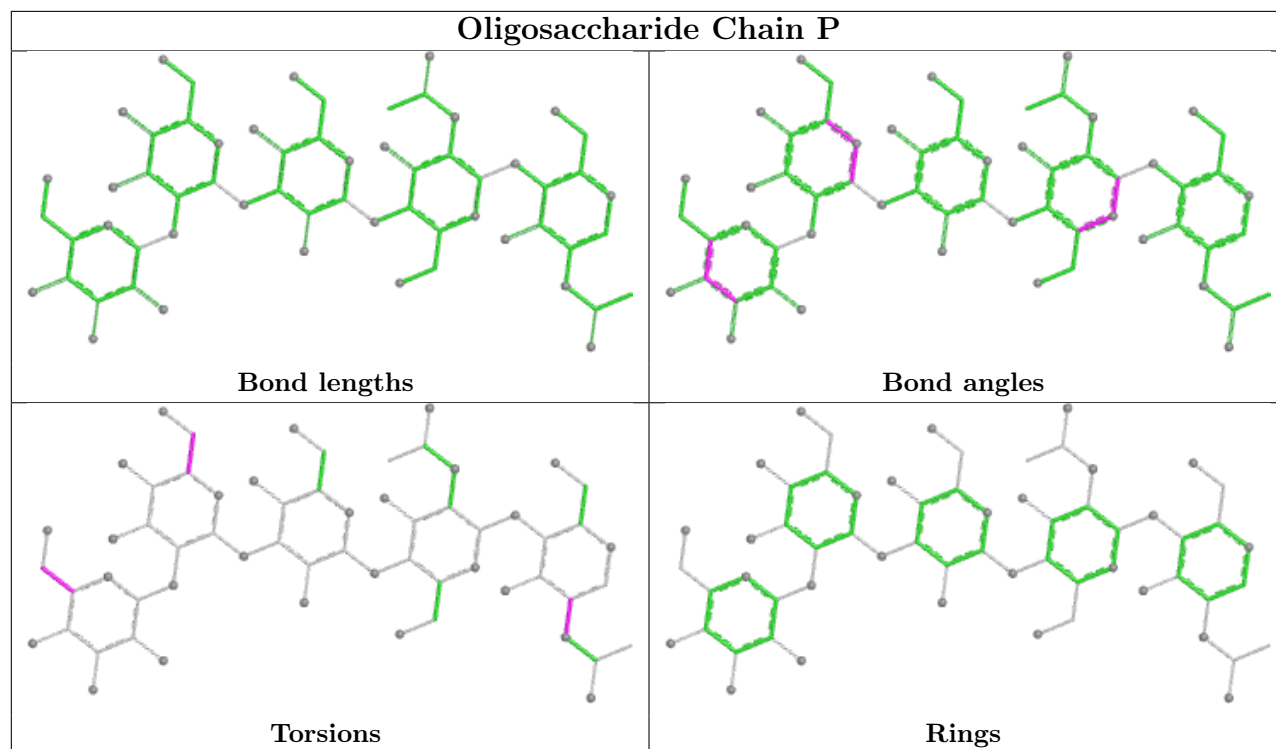


Oligosaccharide Chain K









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	534/556 (96%)	-0.59	2 (0%) 88 90	28, 39, 53, 91	0
1	B	534/556 (96%)	-0.30	2 (0%) 88 90	35, 53, 76, 100	0
1	C	534/556 (96%)	-0.42	2 (0%) 88 90	33, 47, 69, 104	0
1	D	534/556 (96%)	-0.46	1 (0%) 91 92	32, 44, 61, 89	0
1	E	534/556 (96%)	-0.53	1 (0%) 91 92	30, 41, 58, 90	0
1	F	534/556 (96%)	-0.32	2 (0%) 88 90	34, 50, 76, 104	0
1	G	534/556 (96%)	-0.47	2 (0%) 88 90	31, 43, 63, 98	0
1	H	534/556 (96%)	-0.47	2 (0%) 88 90	33, 43, 60, 94	0
All	All	4272/4448 (96%)	-0.44	14 (0%) 90 91	28, 44, 69, 104	0

All (14) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	534	HIS	3.6
1	C	534	HIS	3.5
1	G	238	THR	3.3
1	E	238	THR	3.1
1	C	238	THR	3.1
1	A	534	HIS	3.0
1	A	238	THR	2.8
1	B	534	HIS	2.8
1	D	425	MET	2.8
1	G	534	HIS	2.6
1	F	534	HIS	2.6
1	H	238	THR	2.6
1	B	238	THR	2.5
1	F	370	ASN	2.3

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

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

6.3 Carbohydrates ⓘ

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

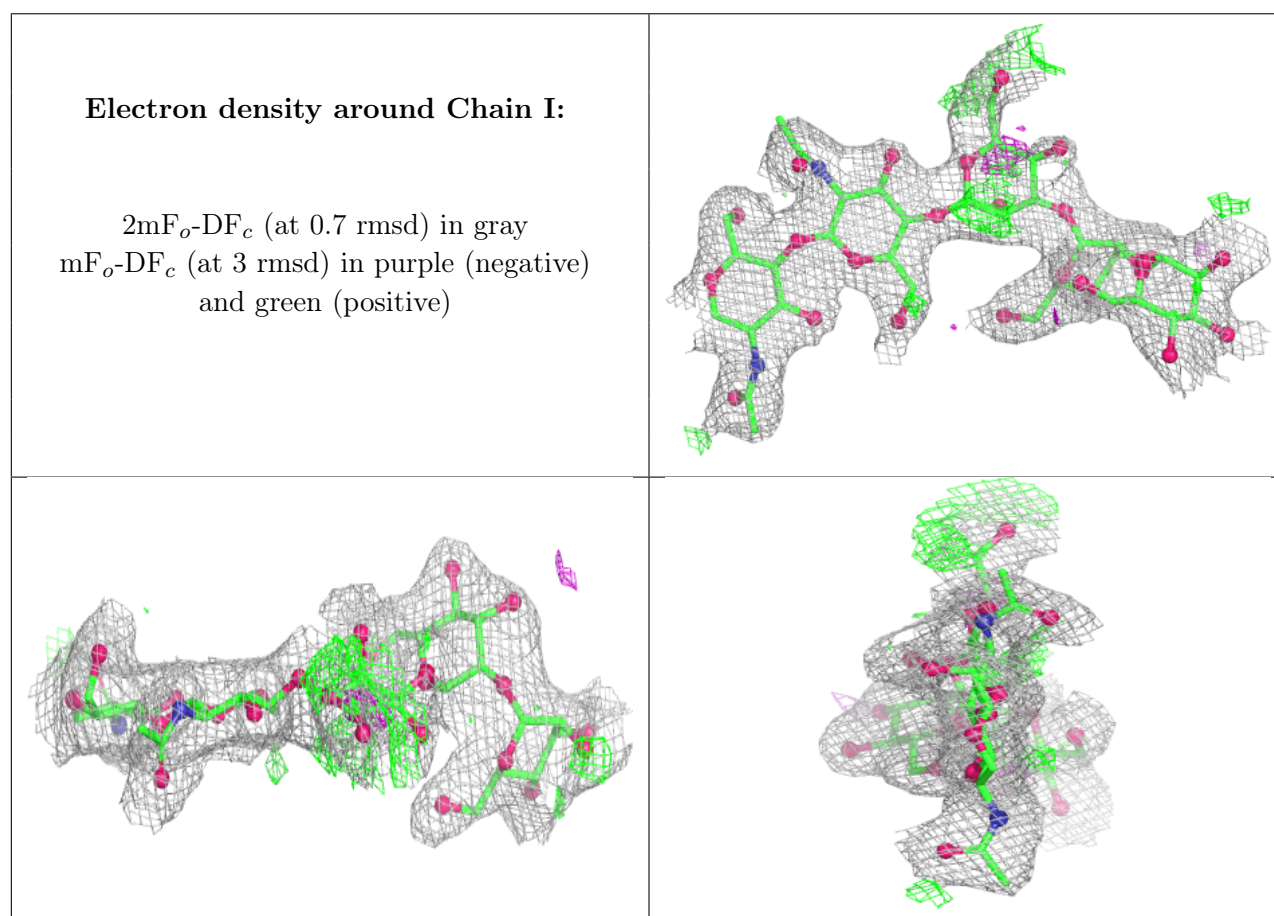
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	BMA	I	3	11/12	0.71	0.13	55,58,61,63	0
2	BMA	K	3	11/12	0.71	0.14	65,67,71,73	0
2	BMA	J	3	11/12	0.74	0.12	70,72,77,79	0
2	MAN	K	5	11/12	0.77	0.12	70,72,74,75	0
2	BMA	N	3	11/12	0.79	0.12	63,65,69,70	0
2	BMA	M	3	11/12	0.80	0.11	58,61,65,65	0
2	MAN	J	5	11/12	0.80	0.12	71,73,76,77	0
2	MAN	P	5	11/12	0.80	0.12	71,75,77,79	0
2	MAN	N	5	11/12	0.81	0.11	69,71,73,74	0
2	MAN	M	5	11/12	0.81	0.11	70,72,73,73	0
2	BMA	O	3	11/12	0.82	0.11	61,64,68,68	0
2	MAN	I	5	11/12	0.84	0.10	65,66,67,67	0
2	BMA	L	3	11/12	0.85	0.12	64,67,72,73	0
2	MAN	O	5	11/12	0.85	0.10	64,66,67,67	0
2	MAN	L	5	11/12	0.85	0.09	72,73,75,76	0
2	BMA	P	3	11/12	0.86	0.11	62,65,68,70	0
2	NAG	K	2	14/15	0.88	0.10	58,62,66,66	0
2	NAG	L	2	14/15	0.89	0.09	55,59,61,62	0
2	NAG	J	1	14/15	0.91	0.09	54,57,60,61	0
2	NAG	J	2	14/15	0.91	0.09	63,66,67,70	0
2	NAG	K	1	14/15	0.92	0.08	50,55,58,58	0
2	MAN	I	4	11/12	0.93	0.08	52,56,59,62	0
2	NAG	O	2	14/15	0.93	0.08	50,52,54,58	0
2	MAN	M	4	11/12	0.93	0.08	56,61,64,66	0
2	MAN	O	4	11/12	0.93	0.08	55,59,63,64	0
2	MAN	K	4	11/12	0.93	0.10	58,63,66,68	0
2	NAG	N	2	14/15	0.93	0.09	57,61,64,65	0
2	NAG	M	2	14/15	0.93	0.07	46,48,50,54	0
2	MAN	L	4	11/12	0.94	0.09	59,62,66,68	0
2	NAG	I	2	14/15	0.94	0.07	45,47,49,52	0
2	MAN	N	4	11/12	0.94	0.07	59,62,65,66	0
2	MAN	J	4	11/12	0.94	0.08	63,68,70,72	0

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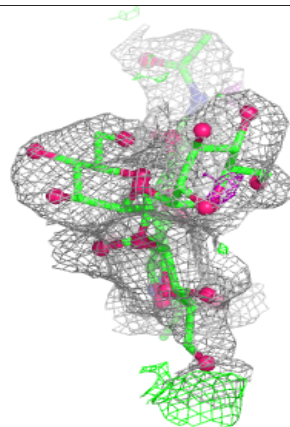
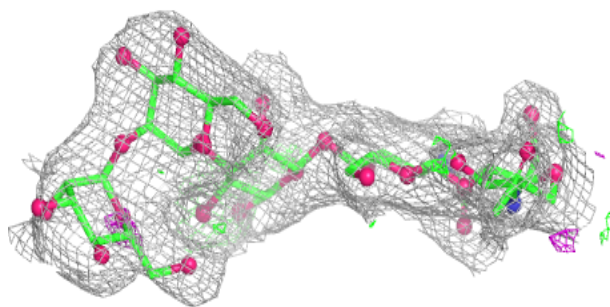
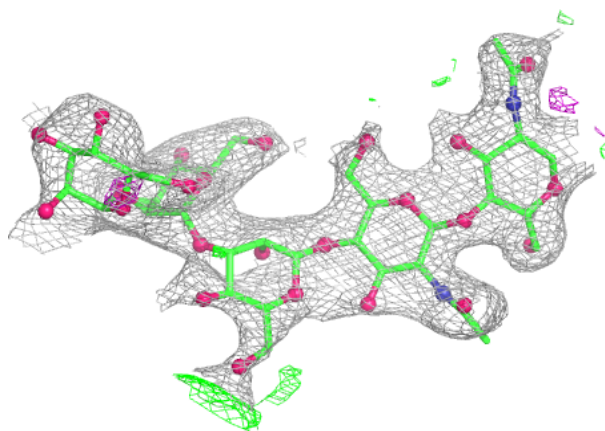
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	N	1	14/15	0.94	0.07	49,53,57,57	0
2	NAG	L	1	14/15	0.95	0.07	44,48,50,53	0
2	MAN	P	4	11/12	0.95	0.08	56,62,66,70	0
2	NAG	P	2	14/15	0.95	0.07	50,53,55,59	0
2	NAG	O	1	14/15	0.96	0.06	43,44,46,47	0
2	NAG	P	1	14/15	0.96	0.05	39,42,44,47	0
2	NAG	M	1	14/15	0.96	0.06	40,41,43,44	0
2	NAG	I	1	14/15	0.98	0.05	37,39,41,42	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.



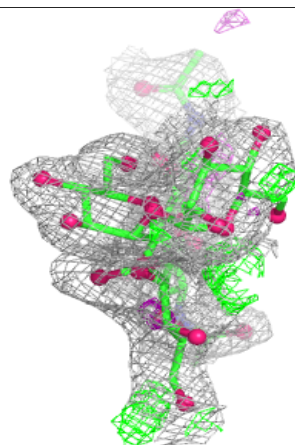
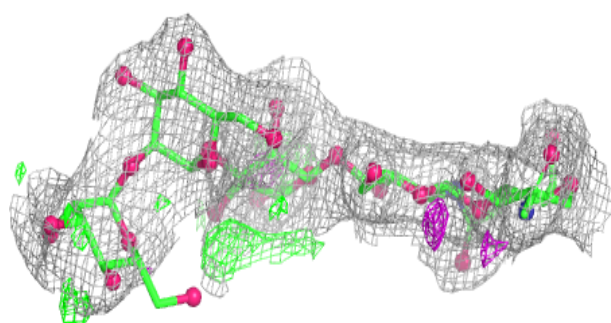
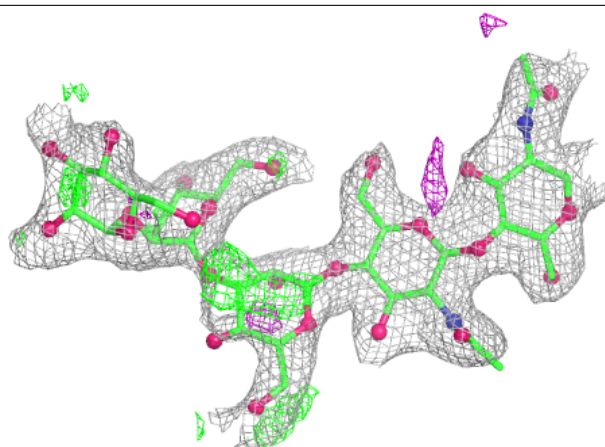
Electron density around Chain J:

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

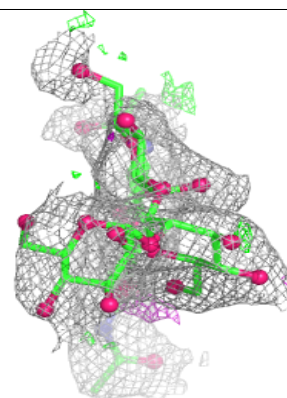
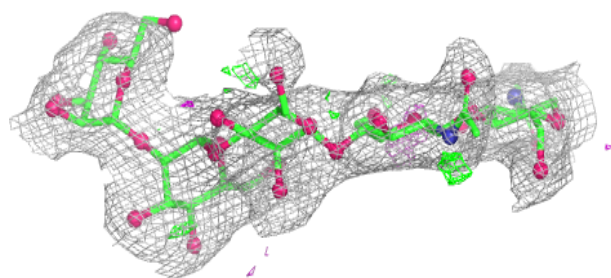
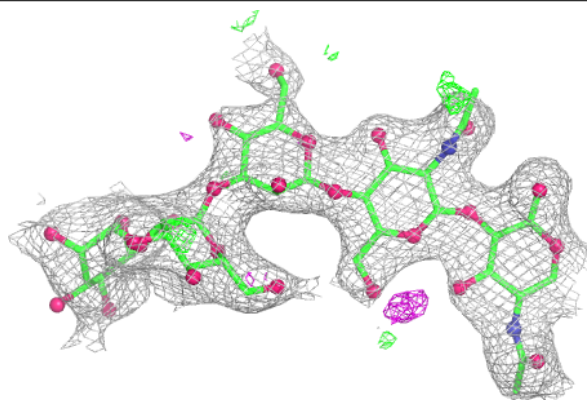


Electron density around Chain K:

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

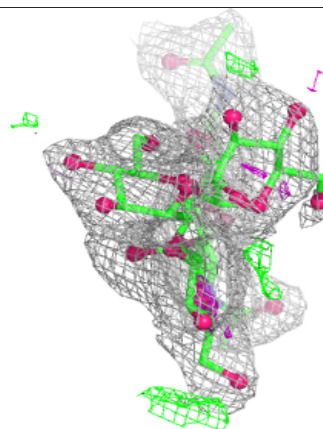
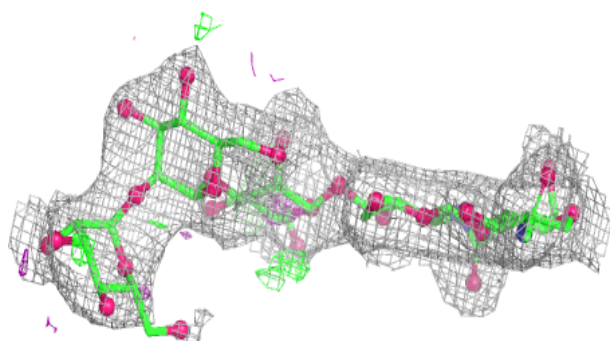
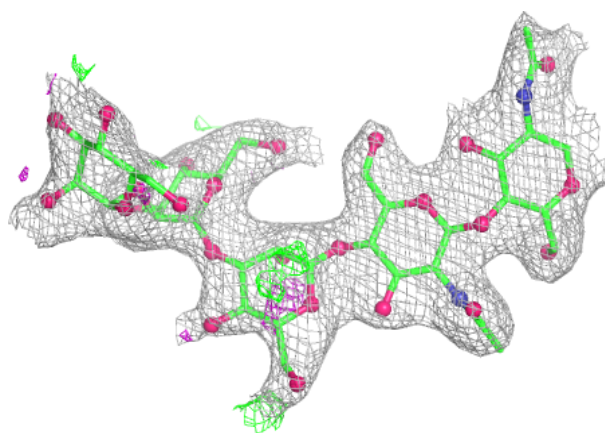
**Electron density around Chain L:**

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

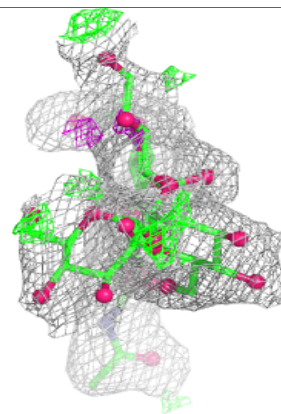
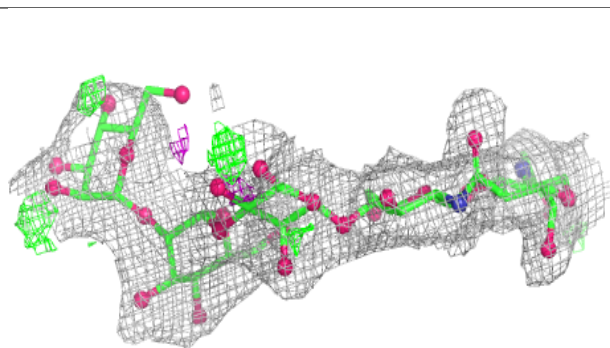
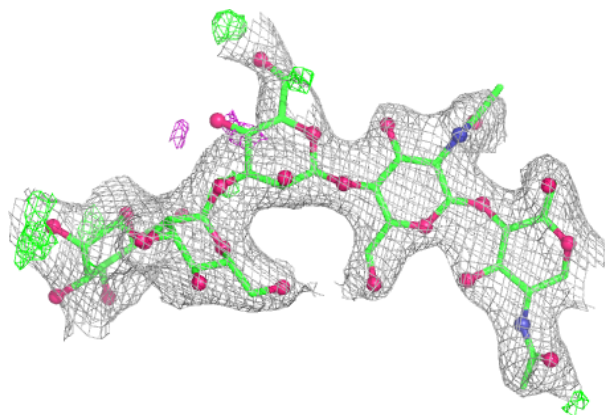


Electron density around Chain M:

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

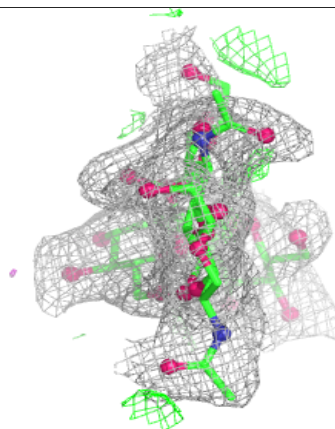
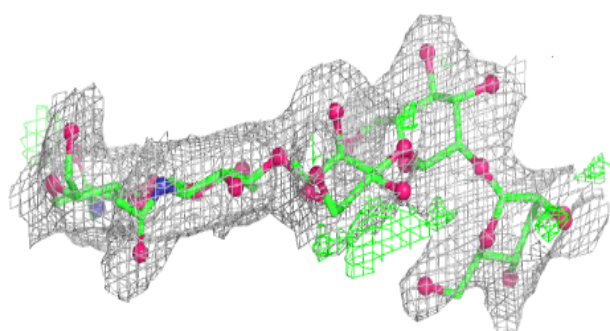
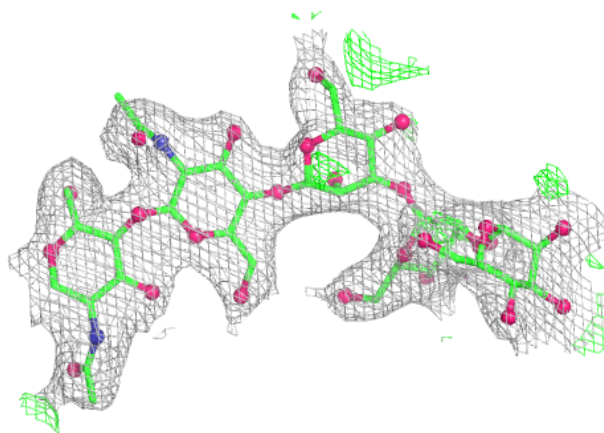
**Electron density around Chain N:**

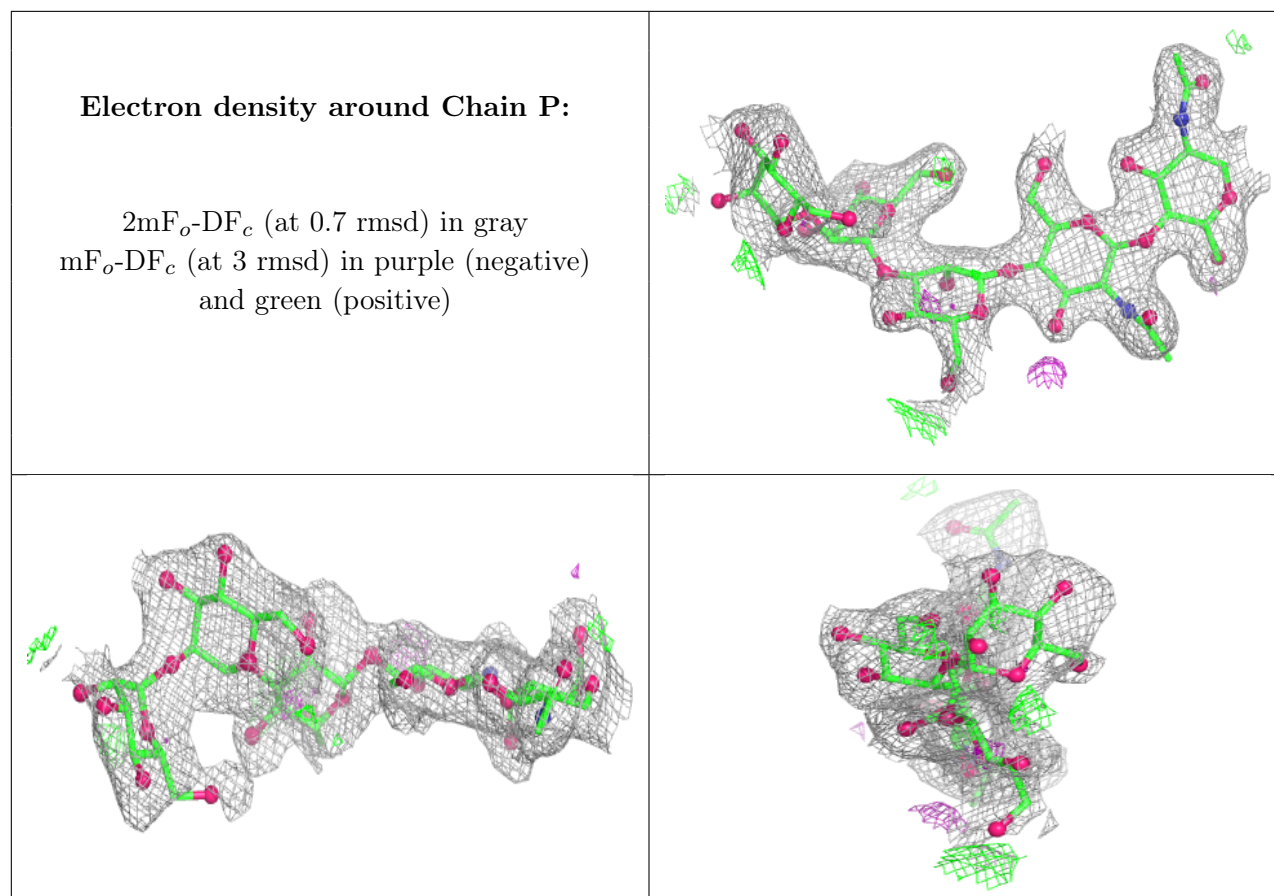
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain O:

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





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.