



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 07:35 AM EDT

PDB ID : 3IJE / pdb_00003ije
Title : Crystal structure of the complete integrin α V β 3 ectodomain plus an Alpha/beta transmembrane fragment
Authors : Xiong, J.-P.; Mahalingham, B.; Rui, X.; Hyman, B.T.; Goodman, S.L.; Arnaout, M.A.
Deposited on : 2009-08-04
Resolution : 2.90 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

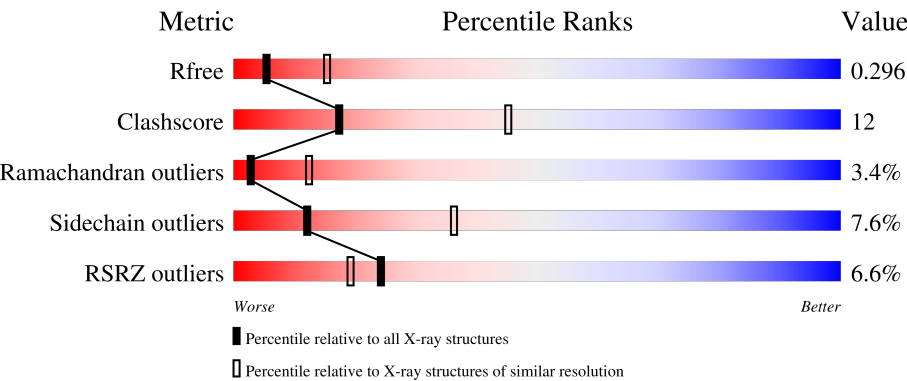
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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.



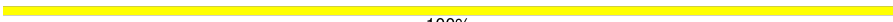
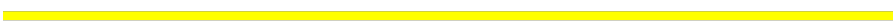
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	967	
2	B	695	
3	C	5	
4	D	2	
4	F	2	

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Mol	Chain	Length	Quality of chain
4	G	2	 100%
4	H	2	 50% 50%
4	I	2	 100%
4	K	2	 100%
5	E	6	 50% 50%
6	J	3	 33% 67%
6	L	3	 33% 67%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NAG	C	2	-	-	X	-

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 13087 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

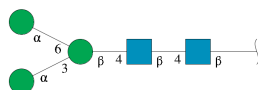
- Molecule 1 is a protein called Integrin alpha-V.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	938	Total	C	N	O	S	0	0	0
			7300	4628	1236	1400	36			

- Molecule 2 is a protein called Integrin beta-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	695	Total	C	N	O	S	0	0	0
			5332	3276	909	1077	70			

- Molecule 3 is an oligosaccharide called 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
3	C	5	Total	C	N	O	0	0	0
			61	34	2	25			

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



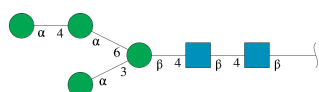
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	2	Total	C	N	O	0	0	0
			28	16	2	10			

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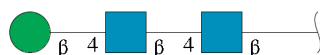
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
4	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-4)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



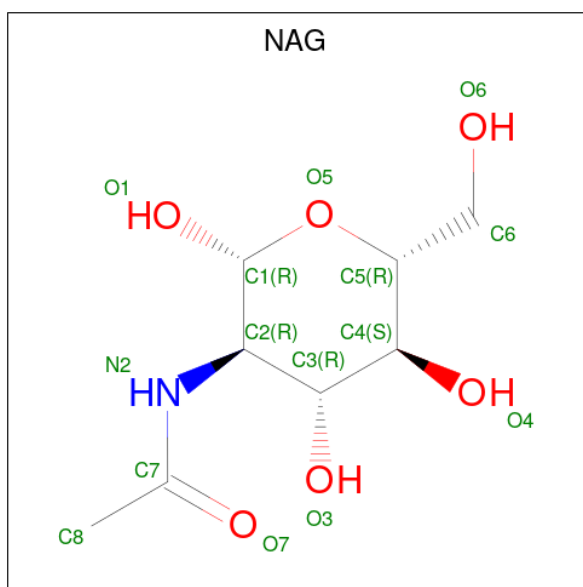
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	E	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 6 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
6	J	3	Total	C	N	O	0	0	0
			39	22	2	15			
6	L	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

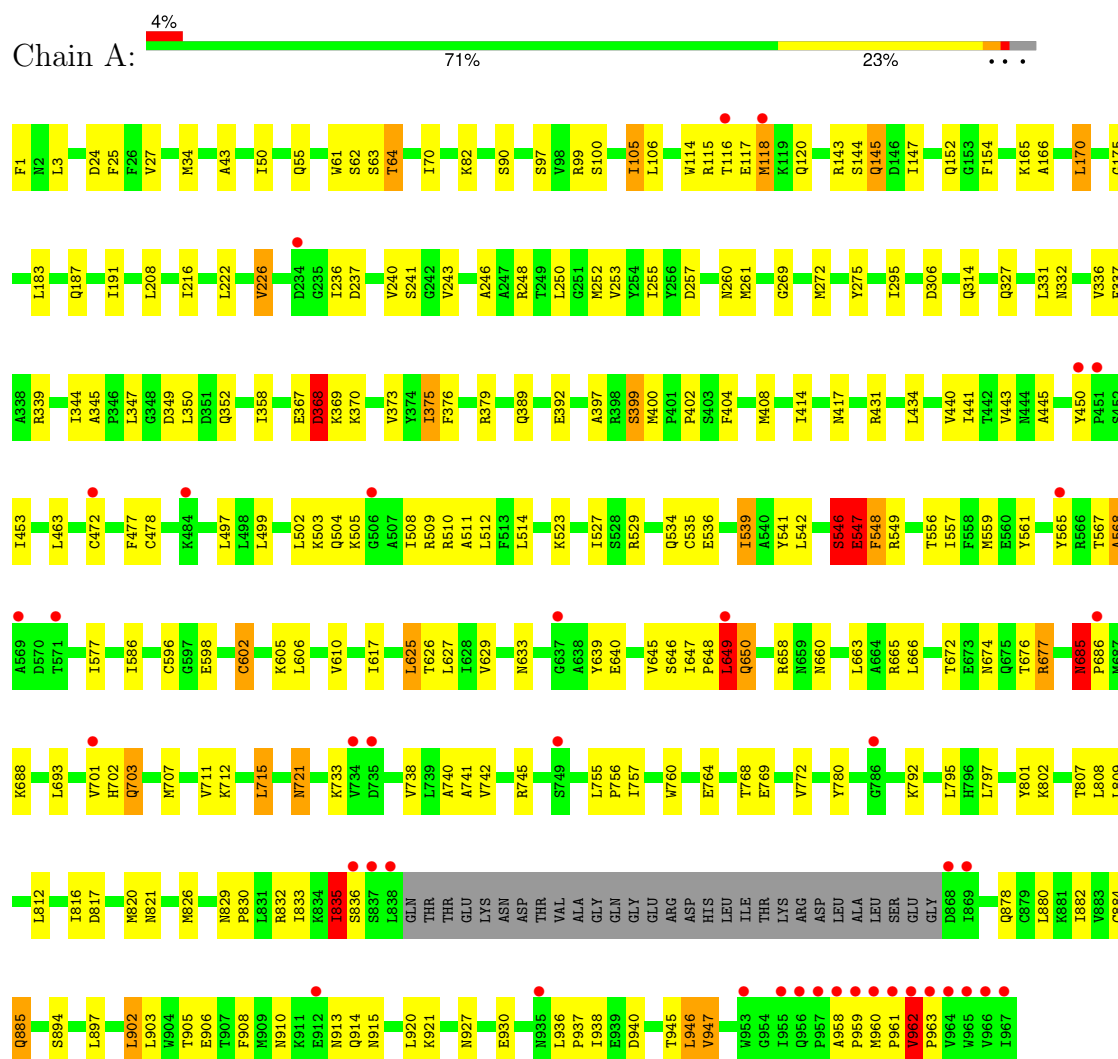
- Molecule 8 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	5	Total	Ca	0	0
			5	5		
8	B	1	Total	Ca	0	0
			1	1		

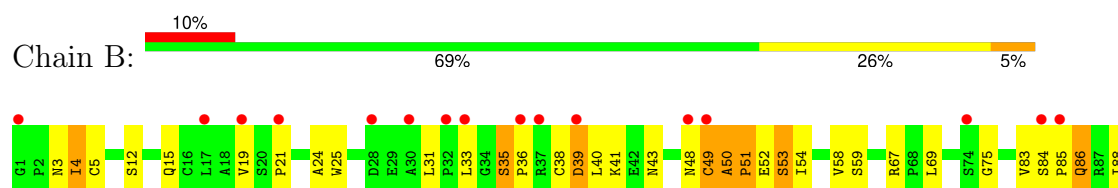
3 Residue-property plots

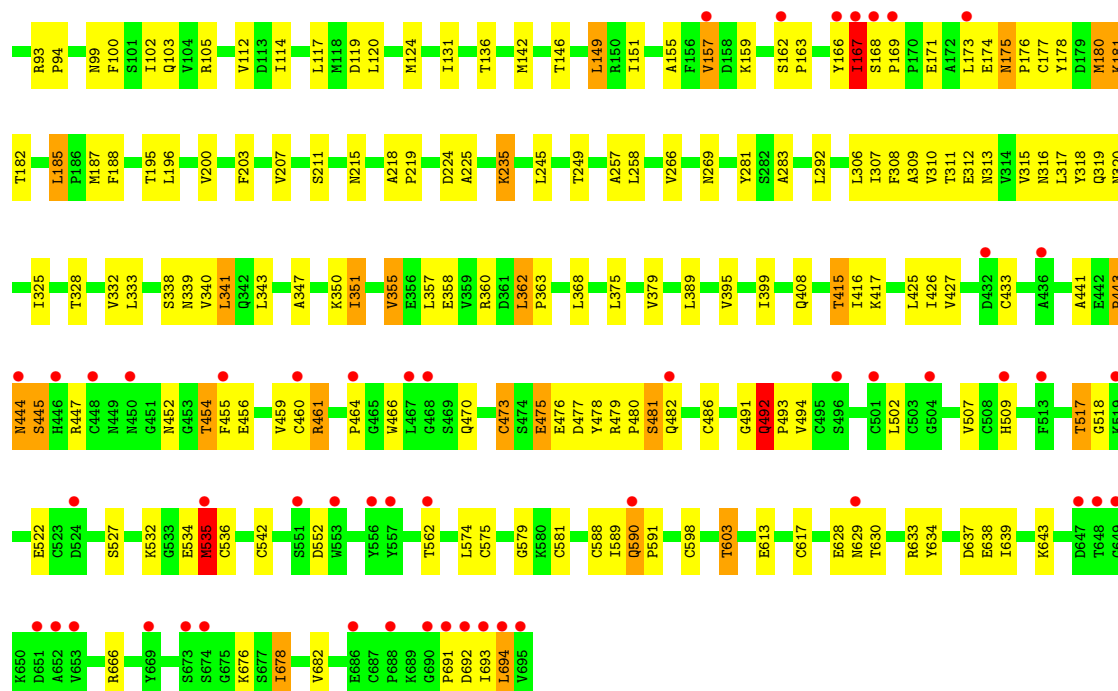
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: Integrin alpha-V



• Molecule 2: Integrin beta-3





- Molecule 3: 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 C: 40% 60%



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

Chain D: 50% 50%



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

Chain F: 50% 50%



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

Chain G: 100%

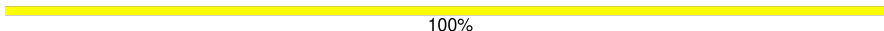
MAG1
MAG2

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

Chain H:  50% 50%

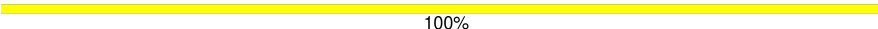
MAG1
MAG2

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

Chain I:  100%

MAG1
MAG2

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

Chain K:  100%

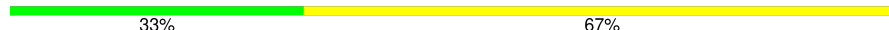
MAG1
MAG2

- Molecule 5: alpha-D-mannopyranose-(1-4)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:  50% 50%

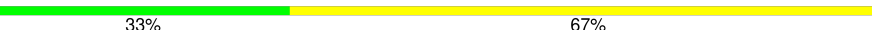
MAG1
MAG2
BMA3
MAN4
MAN5
MAN6

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

Chain J:  33% 67%

MAG1
MAG2
BMA3

- Molecule 6: 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

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.87Å 129.87Å 305.90Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.90 20.00 – 2.90	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.90) 92.3 (20.00-2.90)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.16 (at 2.91Å)	Xtriage
Refinement program	REFMAC 5.2.0019, CNS	Depositor
R, R_{free}	0.244 , 0.285 0.244 , 0.296	Depositor DCC
R_{free} test set	2924 reflections (4.70%)	wwPDB-VP
Wilson B-factor (Å ²)	82.4	Xtriage
Anisotropy	0.191	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 51.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.020 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	13087	wwPDB-VP
Average B, all atoms (Å ²)	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.04% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: BMA, NAG, CA, 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	0.44	0/7462	0.71	6/10123 (0.1%)
2	B	0.46	0/5429	0.77	3/7344 (0.0%)
All	All	0.45	0/12891	0.74	9/17467 (0.1%)

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	A	0	1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	548	PHE	N-CA-C	-6.87	100.14	110.24
1	A	547	GLU	CA-C-N	-6.87	111.23	121.05
1	A	547	GLU	C-N-CA	-6.87	111.23	121.05
1	A	915	ASN	N-CA-C	5.70	116.80	108.14
2	B	492	GLN	CA-C-N	-5.64	113.95	119.76
2	B	492	GLN	C-N-CA	-5.64	113.95	119.76
1	A	685	ASN	CA-C-N	-5.51	112.95	119.84
1	A	685	ASN	C-N-CA	-5.51	112.95	119.84
2	B	590	GLN	N-CA-C	5.14	121.17	109.81

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	546	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	7300	0	7126	171	0
2	B	5332	0	5071	147	0
3	C	61	0	52	7	0
4	D	28	0	25	0	0
4	F	28	0	25	0	0
4	G	28	0	25	0	0
4	H	28	0	25	2	0
4	I	28	0	25	0	0
4	K	28	0	25	0	0
5	E	72	0	61	0	0
6	J	39	0	34	0	0
6	L	39	0	34	0	0
7	A	28	0	26	0	0
7	B	42	0	39	5	0
8	A	5	0	0	0	0
8	B	1	0	0	0	0
All	All	13087	0	12593	319	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 12.

All (319) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:27:VAL:HG13	1:A:34:MET:HE3	1.53	0.90
1:A:3:LEU:HD11	1:A:350:LEU:HD11	1.58	0.85
1:A:445:ALA:HB1	1:A:559:MET:HE1	1.60	0.84
1:A:559:MET:HE2	1:A:586:ILE:HG22	1.61	0.82
2:B:249:THR:HG22	2:B:309:ALA:HB3	1.59	0.82
2:B:281:TYR:CE1	2:B:283:ALA:HB3	2.14	0.82
1:A:639:TYR:H	1:A:721:ASN:HD21	1.30	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:596:CYS:HG	1:A:602:CYS:HG	0.81	0.78
1:A:801:TYR:CD2	1:A:826:MET:HE1	2.18	0.78
1:A:546:SER:O	1:A:547:GLU:HB2	1.82	0.77
1:A:958:ALA:HB1	1:A:959:PRO:HD2	1.65	0.77
2:B:169:PRO:HG2	2:B:173:LEU:HD23	1.65	0.77
1:A:821:ASN:HD21	4:H:1:NAG:C1	1.98	0.76
1:A:3:LEU:CD1	1:A:350:LEU:HD11	2.15	0.76
1:A:772:VAL:HG21	1:A:833:ILE:HD13	1.70	0.74
2:B:579:GLY:HA2	2:B:589:ILE:HD13	1.71	0.73
2:B:375:LEU:HD13	2:B:633:ARG:HD2	1.70	0.73
1:A:443:VAL:HG21	1:A:561:TYR:CE1	2.23	0.72
1:A:617:ILE:HD13	1:A:701:VAL:HG11	1.70	0.71
2:B:454:THR:HG23	2:B:461:ARG:O	1.91	0.71
2:B:589:ILE:HG22	2:B:591:PRO:HD3	1.71	0.71
2:B:574:LEU:HD11	2:B:581:CYS:HB2	1.73	0.70
2:B:5:CYS:HB3	2:B:38:CYS:SG	2.33	0.69
1:A:400:MET:HE1	2:B:266:VAL:HG13	1.73	0.68
2:B:292:LEU:HD22	2:B:325:ILE:HD11	1.75	0.68
1:A:499:LEU:HD23	1:A:557:ILE:HD12	1.77	0.67
2:B:333:LEU:HD11	2:B:340:VAL:HG22	1.77	0.67
2:B:479:ARG:O	2:B:481:SER:N	2.28	0.66
1:A:116:THR:HG21	1:A:147:ILE:HD13	1.78	0.66
2:B:320:ASN:OD1	7:B:3320:NAG:H82	1.96	0.66
2:B:355:VAL:HG12	2:B:389:LEU:CD1	2.26	0.66
2:B:454:THR:O	2:B:460:CYS:HB2	1.96	0.65
1:A:885:GLN:HE21	1:A:885:GLN:N	1.95	0.65
2:B:362:LEU:HD23	2:B:363:PRO:HD2	1.79	0.65
2:B:589:ILE:HG23	2:B:590:GLN:H	1.62	0.64
1:A:368:ASP:O	1:A:370:LYS:N	2.31	0.63
1:A:527:ILE:HG22	1:A:534:GLN:HB3	1.79	0.63
1:A:147:ILE:HD12	1:A:147:ILE:O	1.97	0.63
1:A:617:ILE:HD11	1:A:625:LEU:HD22	1.81	0.63
1:A:314:GLN:HE21	1:A:332:ASN:HD21	1.45	0.63
1:A:443:VAL:HG21	1:A:561:TYR:HE1	1.63	0.62
1:A:116:THR:CG2	1:A:118:MET:SD	2.87	0.62
2:B:588:CYS:HG	2:B:598:CYS:HG	0.70	0.62
1:A:246:ALA:HB1	1:A:252:MET:HE3	1.82	0.62
3:C:2:NAG:H82	3:C:4:MAN:H62	1.82	0.62
1:A:236:ILE:HD12	1:A:237:ASP:H	1.65	0.61
1:A:807:THR:HG22	1:A:807:THR:O	2.00	0.61
2:B:149:LEU:HD21	2:B:151:ILE:HG23	1.83	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:50:ALA:O	2:B:54:ILE:HG23	2.00	0.61
1:A:768:THR:HB	1:A:835:ILE:HD12	1.82	0.61
2:B:613:GLU:O	2:B:617:CYS:HB2	2.00	0.61
1:A:769:GLU:HG2	1:A:812:LEU:HD11	1.82	0.60
1:A:248:ARG:HD2	7:B:3320:NAG:H81	1.83	0.60
2:B:281:TYR:CZ	2:B:283:ALA:HB3	2.36	0.60
2:B:589:ILE:CG2	2:B:590:GLN:H	2.14	0.60
2:B:180:MET:O	2:B:182:THR:HG22	2.00	0.60
2:B:149:LEU:C	2:B:149:LEU:HD23	2.27	0.60
2:B:332:VAL:O	2:B:343:LEU:HD21	2.02	0.60
2:B:316:ASN:HB3	7:B:3320:NAG:H83	1.84	0.60
2:B:589:ILE:CG2	2:B:590:GLN:N	2.65	0.60
1:A:610:VAL:HG11	1:A:715:LEU:HD21	1.85	0.59
2:B:218:ALA:HB3	2:B:219:PRO:HD3	1.84	0.59
1:A:546:SER:O	1:A:547:GLU:CB	2.51	0.59
1:A:508:ILE:HD11	2:B:475:GLU:O	2.02	0.59
2:B:502:LEU:HD13	2:B:507:VAL:CG2	2.33	0.59
1:A:745:ARG:HG3	2:B:603:THR:HG21	1.84	0.59
2:B:694:LEU:HD23	2:B:694:LEU:O	2.03	0.59
1:A:187:GLN:O	1:A:191:ILE:HD12	2.03	0.58
1:A:672:THR:HG23	1:A:676:THR:O	2.03	0.58
1:A:547:GLU:CD	2:B:477:ASP:HA	2.28	0.58
1:A:627:LEU:O	1:A:627:LEU:HD12	2.03	0.58
2:B:355:VAL:HG12	2:B:389:LEU:HD11	1.85	0.58
1:A:755:LEU:HD11	1:A:908:PHE:HB3	1.86	0.58
1:A:70:ILE:HD13	1:A:105:ILE:HD11	1.86	0.58
1:A:547:GLU:O	1:A:548:PHE:C	2.43	0.58
1:A:165:LYS:O	1:A:166:ALA:HB3	2.04	0.57
1:A:666:LEU:HD21	1:A:693:LEU:HD23	1.86	0.57
2:B:51:PRO:O	2:B:53:SER:N	2.37	0.57
1:A:721:ASN:N	1:A:721:ASN:HD22	2.02	0.57
1:A:936:LEU:HD22	1:A:937:PRO:HD2	1.85	0.57
1:A:567:THR:HG23	1:A:567:THR:O	2.05	0.56
2:B:24:ALA:O	2:B:38:CYS:HA	2.05	0.56
2:B:169:PRO:CG	2:B:173:LEU:HD23	2.35	0.56
1:A:685:ASN:HB3	1:A:686:PRO:CD	2.36	0.56
1:A:116:THR:HG22	1:A:118:MET:H	1.68	0.56
1:A:170:LEU:HD13	1:A:226:VAL:HG22	1.87	0.56
2:B:441:ALA:HB1	2:B:456:GLU:CB	2.35	0.56
1:A:24:ASP:OD2	1:A:25:PHE:N	2.38	0.56
2:B:39:ASP:O	2:B:40:LEU:HD22	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:119:ASP:HB3	2:B:124:MET:HE2	1.88	0.56
1:A:880:LEU:C	1:A:880:LEU:HD23	2.30	0.56
1:A:903:LEU:C	1:A:903:LEU:HD23	2.31	0.56
1:A:61:TRP:O	1:A:63:SER:N	2.39	0.56
2:B:159:LYS:NZ	2:B:224:ASP:OD1	2.39	0.56
1:A:502:LEU:HD12	2:B:509:HIS:CD2	2.41	0.56
1:A:3:LEU:H	1:A:389:GLN:HE22	1.53	0.55
1:A:757:ILE:O	1:A:757:ILE:HG23	2.04	0.55
1:A:936:LEU:HD13	1:A:937:PRO:HD2	1.89	0.55
1:A:3:LEU:HD13	1:A:350:LEU:HD21	1.88	0.55
1:A:208:LEU:HB3	1:A:261:MET:HE2	1.89	0.55
1:A:332:ASN:N	1:A:332:ASN:HD22	2.02	0.55
2:B:112:VAL:HG11	2:B:142:MET:CE	2.38	0.54
2:B:50:ALA:HB3	2:B:51:PRO:HD3	1.89	0.54
2:B:173:LEU:HD12	2:B:173:LEU:O	2.06	0.54
1:A:946:LEU:H	1:A:946:LEU:HD23	1.73	0.54
2:B:83:VAL:O	2:B:86:GLN:NE2	2.36	0.54
1:A:816:ILE:HG23	1:A:820:MET:CE	2.37	0.54
2:B:15:GLN:O	2:B:19:VAL:HG23	2.07	0.54
1:A:936:LEU:HD13	1:A:937:PRO:CD	2.38	0.54
1:A:27:VAL:CG1	1:A:34:MET:HE3	2.33	0.53
1:A:792:LYS:HB2	1:A:930:GLU:HB2	1.90	0.53
2:B:441:ALA:CB	2:B:456:GLU:HB2	2.39	0.53
2:B:491:GLY:O	2:B:492:GLN:HB2	2.09	0.53
1:A:648:PRO:HG3	1:A:711:VAL:HG22	1.90	0.53
1:A:795:LEU:HB3	1:A:884:CYS:HB2	1.90	0.53
1:A:154:PHE:O	1:A:175:GLY:HA3	2.08	0.53
1:A:414:ILE:HG21	1:A:434:LEU:HD21	1.91	0.53
2:B:157:VAL:HG12	2:B:157:VAL:O	2.09	0.53
2:B:308:PHE:CE2	2:B:328:THR:HG21	2.44	0.53
1:A:617:ILE:HD11	1:A:625:LEU:CD2	2.39	0.52
2:B:131:ILE:HG22	2:B:207:VAL:HG21	1.91	0.52
1:A:248:ARG:O	1:A:272:MET:HE3	2.09	0.52
1:A:314:GLN:HE21	1:A:332:ASN:ND2	2.06	0.52
1:A:755:LEU:O	1:A:757:ILE:N	2.43	0.52
2:B:637:ASP:O	2:B:639:ILE:HD12	2.09	0.52
1:A:347:LEU:CD1	1:A:375:ILE:HD13	2.37	0.52
1:A:510:ARG:O	1:A:542:LEU:HD12	2.10	0.52
1:A:347:LEU:HD13	1:A:350:LEU:HD22	1.91	0.52
2:B:375:LEU:HD21	2:B:630:THR:HG22	1.91	0.52
2:B:459:VAL:HG22	2:B:460:CYS:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:LYS:O	1:A:236:ILE:HD11	2.09	0.52
1:A:453:ILE:HD11	1:A:639:TYR:CE2	2.45	0.52
2:B:441:ALA:HB1	2:B:456:GLU:HB3	1.92	0.52
1:A:27:VAL:HG22	1:A:34:MET:HE2	1.92	0.52
2:B:693:ILE:O	2:B:693:ILE:HG23	2.10	0.52
1:A:222:LEU:HA	1:A:243:VAL:HG22	1.92	0.51
2:B:166:TYR:CD2	2:B:173:LEU:HD21	2.45	0.51
1:A:100:SER:HB2	1:A:105:ILE:HG22	1.91	0.51
1:A:903:LEU:HD21	1:A:908:PHE:CD1	2.45	0.51
1:A:246:ALA:HB1	1:A:252:MET:CE	2.41	0.51
2:B:112:VAL:HG12	2:B:146:THR:HG21	1.93	0.51
1:A:769:GLU:HG2	1:A:902:LEU:HD21	1.92	0.51
1:A:685:ASN:HD22	1:A:685:ASN:C	2.19	0.51
1:A:43:ALA:HB3	1:A:55:GLN:HG2	1.92	0.51
2:B:112:VAL:CG1	2:B:146:THR:HG21	2.41	0.51
2:B:185:LEU:HD11	2:B:211:SER:O	2.11	0.51
1:A:960:MET:HE3	1:A:963:PRO:HG2	1.91	0.50
2:B:3:ASN:O	2:B:5:CYS:N	2.43	0.50
1:A:116:THR:HG22	1:A:117:GLU:N	2.25	0.50
1:A:946:LEU:HD23	1:A:946:LEU:N	2.25	0.50
2:B:310:VAL:HG11	2:B:318:TYR:CD1	2.46	0.50
2:B:39:ASP:C	2:B:40:LEU:HD22	2.37	0.50
2:B:33:LEU:HD21	2:B:470:GLN:HE21	1.76	0.50
1:A:625:LEU:HD12	1:A:626:THR:N	2.28	0.49
2:B:333:LEU:CD1	2:B:340:VAL:HG22	2.40	0.49
1:A:243:VAL:HG12	1:A:246:ALA:HB2	1.94	0.49
2:B:3:ASN:HB3	2:B:4:ILE:HD12	1.94	0.49
1:A:253:VAL:HG21	1:A:295:ILE:HD13	1.93	0.49
2:B:40:LEU:HD23	2:B:43:ASN:HD22	1.78	0.49
2:B:39:ASP:OD1	2:B:40:LEU:N	2.43	0.49
1:A:514:LEU:HD23	1:A:539:ILE:HD13	1.95	0.49
2:B:99:ASN:HD22	7:B:3099:NAG:H61	1.78	0.49
2:B:166:TYR:O	2:B:167:ILE:C	2.56	0.48
2:B:535:MET:SD	2:B:536:CYS:N	2.78	0.48
1:A:472:CYS:HA	1:A:541:TYR:HA	1.94	0.48
2:B:177:CYS:HB3	2:B:182:THR:HG23	1.94	0.48
1:A:183:LEU:HD13	1:A:222:LEU:HG	1.95	0.48
1:A:809:LEU:HG	1:A:920:LEU:HD13	1.95	0.48
1:A:936:LEU:HD13	1:A:937:PRO:N	2.29	0.48
2:B:40:LEU:HD23	2:B:43:ASN:ND2	2.27	0.48
2:B:117:LEU:HD21	2:B:225:ALA:HB1	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:816:ILE:HG23	1:A:820:MET:HE2	1.94	0.48
2:B:456:GLU:O	2:B:459:VAL:O	2.32	0.48
2:B:574:LEU:HD13	2:B:575:CYS:N	2.29	0.48
1:A:645:VAL:HG13	1:A:645:VAL:O	2.13	0.48
2:B:67:ARG:N	2:B:86:GLN:OE1	2.47	0.48
1:A:1:PHE:HA	1:A:389:GLN:HB2	1.94	0.47
1:A:647:ILE:C	1:A:647:ILE:HD12	2.39	0.47
1:A:567:THR:O	1:A:568:ALA:HB2	2.14	0.47
1:A:70:ILE:HD13	1:A:105:ILE:CD1	2.45	0.47
2:B:21:PRO:HA	2:B:93:ARG:HD2	1.96	0.47
1:A:499:LEU:CD2	1:A:557:ILE:HD12	2.45	0.47
1:A:514:LEU:CD2	1:A:539:ILE:HD13	2.44	0.47
2:B:175:ASN:N	2:B:175:ASN:OD1	2.47	0.47
2:B:307:ILE:HD13	2:B:347:ALA:CB	2.45	0.47
2:B:375:LEU:HD21	2:B:630:THR:CG2	2.43	0.47
2:B:455:PHE:O	2:B:456:GLU:HB3	2.13	0.47
2:B:166:TYR:HD2	2:B:173:LEU:HD21	1.78	0.47
1:A:702:HIS:O	1:A:703:GLN:C	2.58	0.47
2:B:88:ILE:CG2	2:B:427:VAL:HG22	2.44	0.47
2:B:195:THR:HG22	2:B:235:LYS:O	2.15	0.47
3:C:2:NAG:H82	3:C:4:MAN:C6	2.46	0.47
1:A:397:ALA:HB2	1:A:402:PRO:HD3	1.97	0.46
2:B:48:ASN:O	2:B:49:CYS:C	2.58	0.46
2:B:517:THR:HG22	2:B:518:GLY:H	1.79	0.46
1:A:596:CYS:CB	1:A:602:CYS:HG	2.28	0.46
1:A:905:THR:O	1:A:906:GLU:C	2.58	0.46
2:B:332:VAL:O	2:B:332:VAL:HG13	2.15	0.46
1:A:257:ASP:O	1:A:260:ASN:O	2.34	0.46
1:A:927:ASN:HD21	1:A:940:ASP:HB3	1.81	0.46
1:A:962:VAL:N	1:A:963:PRO:CD	2.78	0.46
2:B:93:ARG:HD3	2:B:94:PRO:HD2	1.97	0.46
2:B:362:LEU:HD11	2:B:368:LEU:HD13	1.97	0.46
3:C:2:NAG:O3	3:C:3:BMA:C2	2.63	0.46
1:A:145:GLN:C	1:A:147:ILE:HG23	2.40	0.46
1:A:649:LEU:HD12	1:A:677:ARG:HB3	1.97	0.46
2:B:249:THR:HG22	2:B:309:ALA:CB	2.38	0.46
2:B:312:GLU:O	2:B:315:VAL:HG12	2.15	0.46
2:B:360:ARG:HB2	2:B:415:THR:HG23	1.98	0.46
1:A:738:VAL:HG23	1:A:936:LEU:HD12	1.98	0.46
2:B:389:LEU:HD23	2:B:633:ARG:HH12	1.81	0.45
2:B:426:ILE:HD12	2:B:426:ILE:N	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:114:ILE:HB	2:B:151:ILE:HG22	1.98	0.45
2:B:340:VAL:O	2:B:341:LEU:C	2.59	0.45
1:A:63:SER:O	1:A:64:THR:C	2.60	0.45
1:A:170:LEU:HD13	1:A:226:VAL:CG2	2.47	0.45
1:A:821:ASN:ND2	4:H:1:NAG:C1	2.74	0.45
2:B:25:TRP:HZ2	2:B:459:VAL:HG11	1.80	0.45
2:B:84:SER:HB3	2:B:85:PRO:HD3	1.98	0.45
2:B:339:ASN:O	2:B:343:LEU:HD13	2.17	0.45
2:B:445:SER:OG	7:B:3452:NAG:H82	2.16	0.45
1:A:144:SER:OG	1:A:152:GLN:NE2	2.50	0.45
1:A:511:ALA:O	1:A:512:LEU:HD23	2.17	0.45
1:A:100:SER:CB	1:A:105:ILE:HG22	2.47	0.45
2:B:441:ALA:HB1	2:B:456:GLU:HB2	1.99	0.45
1:A:253:VAL:HG21	1:A:295:ILE:CD1	2.47	0.44
1:A:610:VAL:HG12	1:A:629:VAL:HG22	2.00	0.44
2:B:355:VAL:HG11	2:B:395:VAL:HG21	1.98	0.44
2:B:103:GLN:OE1	2:B:103:GLN:N	2.51	0.44
1:A:797:LEU:HD23	1:A:882:ILE:HD12	1.99	0.44
1:A:504:GLN:HE22	1:A:509:ARG:HB3	1.81	0.44
2:B:306:LEU:O	2:B:328:THR:HG23	2.17	0.44
2:B:666:ARG:O	2:B:682:VAL:HG12	2.18	0.44
1:A:272:MET:HE1	2:B:317:LEU:HD12	1.98	0.44
2:B:180:MET:O	2:B:182:THR:N	2.50	0.44
1:A:250:LEU:HD22	1:A:269:GLY:O	2.18	0.44
1:A:477:PHE:HD2	1:A:497:LEU:HD21	1.82	0.44
1:A:740:ALA:O	1:A:742:VAL:HG23	2.17	0.44
1:A:757:ILE:HG23	1:A:760:TRP:CB	2.48	0.44
1:A:116:THR:HG22	1:A:118:MET:SD	2.57	0.44
1:A:757:ILE:HG23	1:A:760:TRP:HB3	2.00	0.44
1:A:962:VAL:HG12	1:A:962:VAL:O	2.18	0.44
2:B:441:ALA:HB2	2:B:456:GLU:HB2	2.00	0.44
1:A:885:GLN:HE21	1:A:885:GLN:H	1.63	0.43
2:B:69:LEU:HB3	2:B:105:ARG:HE	1.83	0.43
1:A:183:LEU:HD21	1:A:241:SER:HB2	2.00	0.43
1:A:165:LYS:O	1:A:166:ALA:CB	2.66	0.43
1:A:243:VAL:CG1	1:A:246:ALA:HB2	2.49	0.43
1:A:253:VAL:CG2	1:A:295:ILE:HD13	2.48	0.43
1:A:640:GLU:N	1:A:685:ASN:O	2.52	0.43
1:A:660:ASN:ND2	1:A:663:LEU:HD13	2.32	0.43
2:B:638:GLU:HB2	2:B:678:ILE:HG22	2.00	0.43
1:A:375:ILE:HD12	1:A:376:PHE:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:PHE:CD2	1:A:497:LEU:HD21	2.54	0.43
2:B:3:ASN:C	2:B:4:ILE:HD12	2.44	0.43
1:A:50:ILE:HG23	1:A:90:SER:HA	2.01	0.43
2:B:88:ILE:HG23	2:B:427:VAL:HG22	2.00	0.43
2:B:100:PHE:CE1	2:B:399:ILE:HD12	2.54	0.43
1:A:502:LEU:HD21	1:A:556:THR:OG1	2.19	0.43
1:A:633:ASN:C	1:A:633:ASN:HD22	2.27	0.43
2:B:58:VAL:HG12	2:B:59:SER:O	2.18	0.43
2:B:83:VAL:HG13	2:B:102:ILE:HD11	2.00	0.42
1:A:945:THR:HG22	1:A:946:LEU:N	2.35	0.42
2:B:136:THR:HA	2:B:200:VAL:HG11	1.99	0.42
2:B:311:THR:OG1	2:B:313:ASN:ND2	2.52	0.42
2:B:443:PRO:O	2:B:444:ASN:C	2.62	0.42
1:A:236:ILE:HD12	1:A:237:ASP:N	2.32	0.42
1:A:345:ALA:CB	1:A:408:MET:HE3	2.48	0.42
1:A:826:MET:HE2	1:A:880:LEU:HB2	2.01	0.42
3:C:2:NAG:C3	3:C:3:BMA:H2	2.50	0.42
1:A:757:ILE:O	1:A:757:ILE:CG2	2.66	0.42
2:B:157:VAL:HG23	2:B:188:PHE:CZ	2.54	0.42
1:A:50:ILE:HG23	1:A:90:SER:CA	2.49	0.42
1:A:336:VAL:HG12	1:A:337:PHE:CD2	2.55	0.42
1:A:648:PRO:O	1:A:650:GLN:N	2.52	0.42
2:B:162:SER:HB3	2:B:167:ILE:HG23	2.02	0.42
1:A:344:ILE:HG22	1:A:358:ILE:HD11	2.00	0.42
1:A:801:TYR:CD2	1:A:802:LYS:HG3	2.55	0.42
2:B:332:VAL:C	2:B:343:LEU:HD21	2.45	0.42
3:C:2:NAG:O3	3:C:3:BMA:H2	2.20	0.42
2:B:358:GLU:N	2:B:417:LYS:O	2.45	0.42
1:A:114:TRP:CE3	1:A:143:ARG:HD2	2.55	0.42
1:A:240:VAL:HG22	1:A:255:ILE:HG12	2.01	0.42
2:B:180:MET:O	2:B:181:LYS:C	2.62	0.42
2:B:257:ALA:O	2:B:258:LEU:HB2	2.19	0.42
1:A:672:THR:O	1:A:672:THR:HG22	2.20	0.41
1:A:835:ILE:HG22	1:A:836:SER:N	2.34	0.41
2:B:310:VAL:HG11	2:B:318:TYR:CG	2.55	0.41
2:B:102:ILE:HD12	2:B:425:LEU:HB2	2.02	0.41
1:A:936:LEU:O	1:A:938:ILE:HG23	2.20	0.41
2:B:492:GLN:N	2:B:493:PRO:HD2	2.36	0.41
1:A:780:TYR:CZ	1:A:947:VAL:HG13	2.55	0.41
1:A:829:ASN:N	1:A:830:PRO:HD3	2.36	0.41
2:B:50:ALA:O	2:B:51:PRO:O	2.38	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:203:PHE:O	2:B:207:VAL:HG12	2.21	0.41
2:B:630:THR:HB	2:B:634:TYR:CD1	2.56	0.41
1:A:685:ASN:HB3	1:A:686:PRO:HD3	2.02	0.41
2:B:245:LEU:CD2	2:B:351:ILE:HD11	2.50	0.41
2:B:589:ILE:HG22	2:B:590:GLN:N	2.36	0.41
2:B:48:ASN:O	2:B:50:ALA:HB2	2.20	0.41
2:B:552:ASP:HA	2:B:562:THR:OG1	2.20	0.41
2:B:94:PRO:HD3	2:B:433:CYS:HB3	2.02	0.41
1:A:349:ASP:O	1:A:352:GLN:NE2	2.47	0.41
2:B:12:SER:HB3	2:B:461:ARG:HH21	1.86	0.41
2:B:399:ILE:HD13	2:B:416:ILE:HD13	2.02	0.41
2:B:682:VAL:O	2:B:682:VAL:HG13	2.20	0.41
1:A:373:VAL:HG23	1:A:404:PHE:CE2	2.56	0.41
1:A:440:VAL:HG22	1:A:577:ILE:CG2	2.50	0.41
2:B:187:MET:HE1	2:B:215:ASN:HA	2.02	0.41
1:A:306:ASP:OD1	1:A:306:ASP:N	2.48	0.40
1:A:332:ASN:N	1:A:332:ASN:ND2	2.69	0.40
1:A:625:LEU:HD12	1:A:625:LEU:C	2.46	0.40
2:B:35:SER:HB3	2:B:36:PRO:HD3	2.03	0.40
2:B:502:LEU:HD13	2:B:507:VAL:HG23	2.03	0.40
2:B:173:LEU:HD13	2:B:178:TYR:CE1	2.56	0.40
3:C:2:NAG:HO3	3:C:3:BMA:C2	2.33	0.40
1:A:463:LEU:HD23	1:A:463:LEU:C	2.46	0.40
3:C:2:NAG:C3	3:C:3:BMA:C2	3.00	0.40
2:B:120:LEU:HG	2:B:155:ALA:HB1	2.03	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	934/967 (97%)	843 (90%)	68 (7%)	23 (2%)	4 17

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	693/695 (100%)	579 (84%)	82 (12%)	32 (5%)	2	7
All	All	1627/1662 (98%)	1422 (87%)	150 (9%)	55 (3%)	3	12

All (55) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	62	SER
1	A	505	LYS
1	A	547	GLU
1	A	568	ALA
1	A	703	GLN
1	A	741	ALA
1	A	961	PRO
2	B	39	ASP
2	B	51	PRO
2	B	52	GLU
2	B	167	ILE
2	B	168	SER
2	B	181	LYS
2	B	473	CYS
2	B	478	TYR
2	B	480	PRO
2	B	481	SER
2	B	492	GLN
1	A	64	THR
1	A	368	ASP
1	A	369	LYS
1	A	546	SER
1	A	707	MET
1	A	835	ILE
1	A	914	GLN
2	B	35	SER
2	B	49	CYS
2	B	53	SER
2	B	196	LEU
2	B	444	ASN
2	B	535	MET
1	A	503	LYS
1	A	674	ASN
1	A	808	LEU
2	B	50	ALA
2	B	482	GLN

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Mol	Chain	Res	Type
2	B	527	SER
1	A	399	SER
1	A	529	ARG
1	A	649	LEU
2	B	4	ILE
2	B	443	PRO
2	B	532	LYS
1	A	910	ASN
2	B	41	LYS
2	B	157	VAL
2	B	445	SER
1	A	756	PRO
2	B	464	PRO
2	B	691	PRO
2	B	163	PRO
2	B	176	PRO
1	A	962	VAL
2	B	75	GLY
2	B	494	VAL

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	797/821 (97%)	733 (92%)	64 (8%)	11	34
2	B	617/617 (100%)	573 (93%)	44 (7%)	13	40
All	All	1414/1438 (98%)	1306 (92%)	108 (8%)	12	36

All (108) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	82	LYS
1	A	97	SER
1	A	99	ARG
1	A	105	ILE

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Mol	Chain	Res	Type
1	A	106	LEU
1	A	115	ARG
1	A	118	MET
1	A	120	GLN
1	A	145	GLN
1	A	170	LEU
1	A	216	ILE
1	A	226	VAL
1	A	275	TYR
1	A	327	GLN
1	A	331	LEU
1	A	339	ARG
1	A	367	GLU
1	A	368	ASP
1	A	375	ILE
1	A	379	ARG
1	A	392	GLU
1	A	399	SER
1	A	417	ASN
1	A	431	ARG
1	A	441	ILE
1	A	450	TYR
1	A	478	CYS
1	A	523	LYS
1	A	535	CYS
1	A	536	GLU
1	A	539	ILE
1	A	549	ARG
1	A	565	TYR
1	A	598	GLU
1	A	602	CYS
1	A	605	LYS
1	A	606	LEU
1	A	625	LEU
1	A	646	SER
1	A	649	LEU
1	A	650	GLN
1	A	658	ARG
1	A	665	ARG
1	A	677	ARG
1	A	685	ASN
1	A	688	LYS

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Mol	Chain	Res	Type
1	A	712	LYS
1	A	715	LEU
1	A	721	ASN
1	A	733	LYS
1	A	764	GLU
1	A	817	ASP
1	A	832	ARG
1	A	835	ILE
1	A	878	GLN
1	A	885	GLN
1	A	894	SER
1	A	897	LEU
1	A	902	LEU
1	A	913	ASN
1	A	921	LYS
1	A	946	LEU
1	A	947	VAL
1	A	962	VAL
2	B	31	LEU
2	B	86	GLN
2	B	149	LEU
2	B	167	ILE
2	B	171	GLU
2	B	174	GLU
2	B	175	ASN
2	B	180	MET
2	B	185	LEU
2	B	235	LYS
2	B	269	ASN
2	B	319	GLN
2	B	338	SER
2	B	341	LEU
2	B	350	LYS
2	B	351	ILE
2	B	355	VAL
2	B	357	LEU
2	B	362	LEU
2	B	379	VAL
2	B	408	GLN
2	B	415	THR
2	B	447	ARG
2	B	452	ASN

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Mol	Chain	Res	Type
2	B	454	THR
2	B	461	ARG
2	B	466	TRP
2	B	473	CYS
2	B	475	GLU
2	B	476	GLU
2	B	486	CYS
2	B	517	THR
2	B	522	GLU
2	B	534	GLU
2	B	535	MET
2	B	542	CYS
2	B	603	THR
2	B	628	GLU
2	B	629	ASN
2	B	643	LYS
2	B	676	LYS
2	B	678	ILE
2	B	692	ASP
2	B	694	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (46) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	91	HIS
1	A	120	GLN
1	A	131	GLN
1	A	152	GLN
1	A	156	GLN
1	A	187	GLN
1	A	205	ASN
1	A	206	ASN
1	A	207	GLN
1	A	332	ASN
1	A	389	GLN
1	A	394	GLN
1	A	456	GLN
1	A	474	ASN
1	A	492	ASN
1	A	504	GLN
1	A	580	GLN
1	A	614	GLN

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Mol	Chain	Res	Type
1	A	623	ASN
1	A	633	ASN
1	A	650	GLN
1	A	674	ASN
1	A	685	ASN
1	A	704	GLN
1	A	721	ASN
1	A	732	HIS
1	A	796	HIS
1	A	885	GLN
1	A	927	ASN
2	B	3	ASN
2	B	43	ASN
2	B	210	GLN
2	B	240	ASN
2	B	244	HIS
2	B	269	ASN
2	B	301	GLN
2	B	303	ASN
2	B	313	ASN
2	B	376	ASN
2	B	408	GLN
2	B	438	GLN
2	B	446	HIS
2	B	450	ASN
2	B	470	GLN
2	B	483	GLN
2	B	541	GLN

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 ⓘ

29 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	$\# Z > 2$
3	NAG	C	1	1,3	14,14,15	0.61	0	17,19,21	1.25	2 (11%)
3	NAG	C	2	3	14,14,15	0.41	0	17,19,21	1.50	3 (17%)
3	BMA	C	3	3	11,11,12	0.51	0	15,15,17	2.44	2 (13%)
3	MAN	C	4	3	11,11,12	0.55	0	15,15,17	0.97	1 (6%)
3	MAN	C	5	3	11,11,12	0.58	0	15,15,17	0.93	1 (6%)
4	NAG	D	1	4,1	14,14,15	0.51	0	17,19,21	0.72	0
4	NAG	D	2	4	14,14,15	0.46	0	17,19,21	0.82	1 (5%)
5	NAG	E	1	5,1	14,14,15	0.54	0	17,19,21	1.14	1 (5%)
5	NAG	E	2	5	14,14,15	0.53	0	17,19,21	0.66	0
5	BMA	E	3	5	11,11,12	0.25	0	15,15,17	0.70	0
5	MAN	E	4	5	11,11,12	0.61	0	15,15,17	1.15	1 (6%)
5	MAN	E	5	5	11,11,12	0.63	0	15,15,17	0.80	1 (6%)
5	MAN	E	6	5	11,11,12	0.61	0	15,15,17	0.66	0
4	NAG	F	1	4,1	14,14,15	0.59	0	17,19,21	1.25	1 (5%)
4	NAG	F	2	4	14,14,15	0.50	0	17,19,21	0.77	0
4	NAG	G	1	4,1	14,14,15	0.65	0	17,19,21	1.42	2 (11%)
4	NAG	G	2	4	14,14,15	0.71	0	17,19,21	0.93	1 (5%)
4	NAG	H	1	4	14,14,15	0.66	0	17,19,21	1.65	2 (11%)
4	NAG	H	2	4	14,14,15	0.70	0	17,19,21	0.97	1 (5%)
4	NAG	I	1	4,1	14,14,15	0.68	0	17,19,21	1.25	1 (5%)
4	NAG	I	2	4	14,14,15	0.53	0	17,19,21	1.20	1 (5%)
6	NAG	J	1	1,6	14,14,15	0.48	0	17,19,21	0.96	2 (11%)
6	NAG	J	2	6	14,14,15	0.67	0	17,19,21	1.18	1 (5%)
6	BMA	J	3	6	11,11,12	0.31	0	15,15,17	0.61	0
4	NAG	K	1	2,4	14,14,15	0.47	0	17,19,21	1.93	2 (11%)
4	NAG	K	2	4	14,14,15	0.50	0	17,19,21	1.34	1 (5%)
6	NAG	L	1	2,6	14,14,15	0.51	0	17,19,21	1.10	2 (11%)
6	NAG	L	2	6	14,14,15	0.58	0	17,19,21	1.18	1 (5%)
6	BMA	L	3	6	11,11,12	0.44	0	15,15,17	0.75	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	C	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	C	2	3	-	2/6/23/26	0/1/1/1
3	BMA	C	3	3	-	0/2/19/22	0/1/1/1
3	MAN	C	4	3	-	1/2/19/22	0/1/1/1
3	MAN	C	5	3	-	2/2/19/22	0/1/1/1
4	NAG	D	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1
5	NAG	E	1	5,1	-	4/6/23/26	0/1/1/1
5	NAG	E	2	5	-	0/6/23/26	0/1/1/1
5	BMA	E	3	5	-	2/2/19/22	0/1/1/1
5	MAN	E	4	5	-	0/2/19/22	0/1/1/1
5	MAN	E	5	5	-	2/2/19/22	0/1/1/1
5	MAN	E	6	5	-	2/2/19/22	0/1/1/1
4	NAG	F	1	4,1	-	3/6/23/26	0/1/1/1
4	NAG	F	2	4	-	2/6/23/26	0/1/1/1
4	NAG	G	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	G	2	4	-	2/6/23/26	0/1/1/1
4	NAG	H	1	4	-	2/6/23/26	0/1/1/1
4	NAG	H	2	4	-	3/6/23/26	0/1/1/1
4	NAG	I	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	I	2	4	-	2/6/23/26	0/1/1/1
6	NAG	J	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	J	2	6	-	5/6/23/26	0/1/1/1
6	BMA	J	3	6	-	2/2/19/22	0/1/1/1
4	NAG	K	1	2,4	-	4/6/23/26	0/1/1/1
4	NAG	K	2	4	-	4/6/23/26	0/1/1/1
6	NAG	L	1	2,6	-	4/6/23/26	0/1/1/1
6	NAG	L	2	6	-	2/6/23/26	0/1/1/1
6	BMA	L	3	6	-	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(°)
3	C	3	BMA	C1-O5-C5	7.86	122.72	112.19
4	K	1	NAG	C1-O5-C5	5.65	119.76	112.19
4	G	1	NAG	C4-C3-C2	4.49	117.59	111.02
4	H	1	NAG	C3-C4-C5	4.33	118.07	110.23
4	K	1	NAG	O4-C4-C5	4.27	119.85	109.32
4	I	2	NAG	C1-O5-C5	4.19	117.81	112.19
3	C	3	BMA	C1-C2-C3	4.10	115.61	109.64
4	I	1	NAG	C4-C3-C2	4.08	117.00	111.02
4	K	2	NAG	C1-O5-C5	3.89	117.40	112.19
6	J	2	NAG	C4-C3-C2	3.79	116.57	111.02
4	H	1	NAG	O5-C1-C2	-3.63	105.67	111.29
6	L	1	NAG	O5-C1-C2	-3.48	105.91	111.29
4	F	1	NAG	C4-C3-C2	3.45	116.07	111.02
6	L	2	NAG	C4-C3-C2	3.31	115.87	111.02
3	C	2	NAG	O4-C4-C3	3.07	117.61	110.38
3	C	2	NAG	C1-O5-C5	3.04	116.26	112.19
3	C	4	MAN	C1-O5-C5	2.92	116.11	112.19
3	C	1	NAG	C1-O5-C5	2.86	116.02	112.19
3	C	2	NAG	O3-C3-C2	-2.79	103.62	109.40
4	H	2	NAG	C4-C3-C2	2.75	115.05	111.02
5	E	4	MAN	O4-C4-C3	-2.67	104.09	110.38
3	C	5	MAN	C1-C2-C3	2.62	113.46	109.64
5	E	1	NAG	O5-C1-C2	-2.59	107.28	111.29
4	G	2	NAG	C4-C3-C2	2.58	114.81	111.02
4	G	1	NAG	C3-C4-C5	2.49	114.74	110.23
6	L	1	NAG	C1-O5-C5	2.30	115.27	112.19
6	J	1	NAG	O5-C1-C2	-2.27	107.78	111.29
4	D	2	NAG	C1-O5-C5	2.14	115.05	112.19
5	E	5	MAN	C1-C2-C3	2.12	112.73	109.64
3	C	1	NAG	O4-C4-C5	-2.08	104.20	109.32
6	J	1	NAG	C4-C3-C2	-2.01	108.08	111.02

There are no chirality outliers.

All (68) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	2	NAG	C8-C7-N2-C2
3	C	2	NAG	O7-C7-N2-C2
4	D	2	NAG	C8-C7-N2-C2
4	D	2	NAG	O7-C7-N2-C2
4	F	1	NAG	C3-C2-N2-C7
4	F	2	NAG	C8-C7-N2-C2
4	F	2	NAG	O7-C7-N2-C2

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

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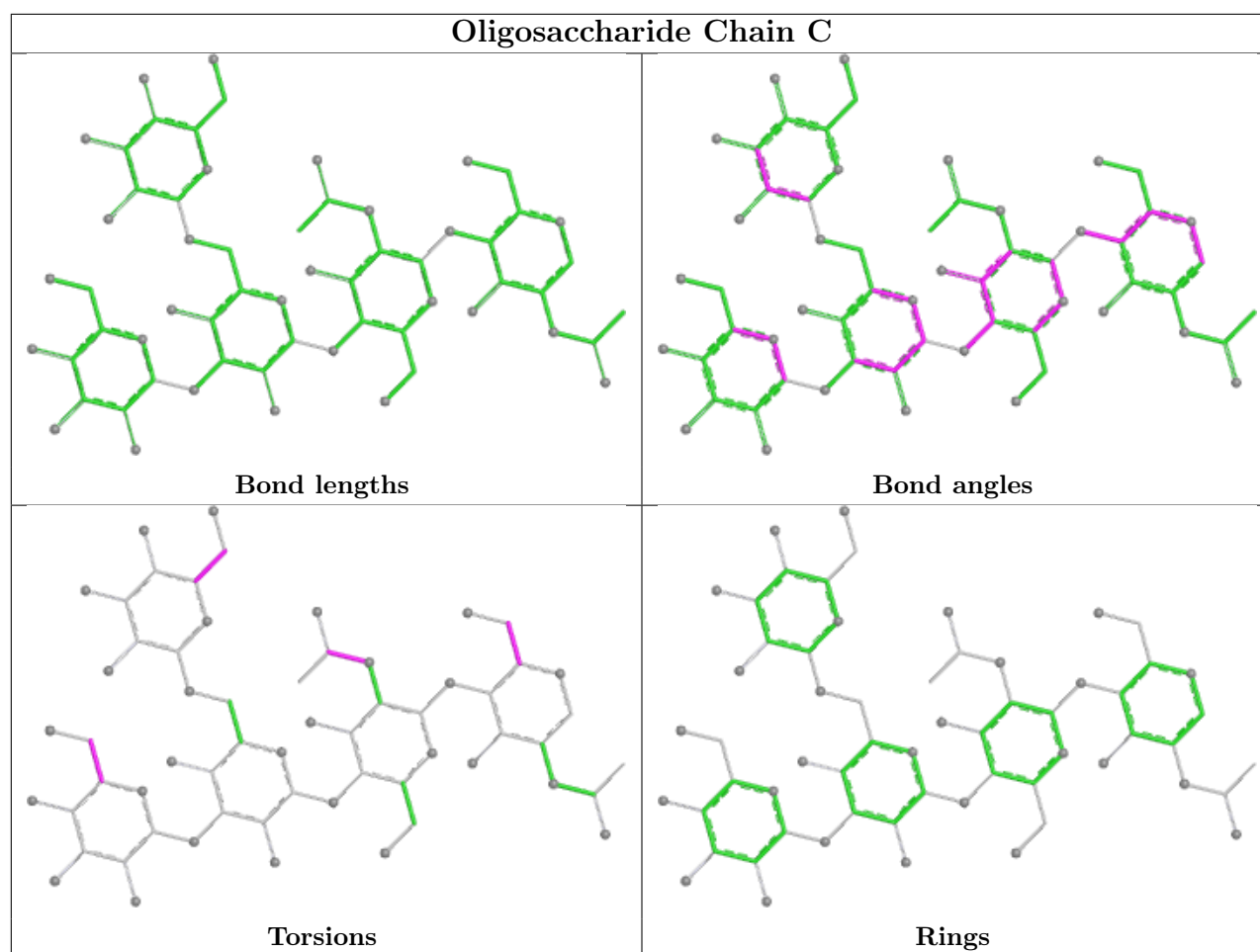
Mol	Chain	Res	Type	Atoms
4	H	1	NAG	O7-C7-N2-C2
4	I	1	NAG	C8-C7-N2-C2
4	I	1	NAG	O7-C7-N2-C2
4	K	2	NAG	C4-C5-C6-O6
5	E	6	MAN	O5-C5-C6-O6
3	C	5	MAN	O5-C5-C6-O6
4	K	1	NAG	C8-C7-N2-C2
5	E	3	BMA	C4-C5-C6-O6
4	D	1	NAG	C4-C5-C6-O6
3	C	4	MAN	O5-C5-C6-O6
6	J	2	NAG	C3-C2-N2-C7
5	E	3	BMA	O5-C5-C6-O6
4	K	1	NAG	O7-C7-N2-C2
5	E	1	NAG	C8-C7-N2-C2
5	E	6	MAN	C4-C5-C6-O6
5	E	1	NAG	O7-C7-N2-C2
6	J	2	NAG	C4-C5-C6-O6
4	H	2	NAG	O5-C5-C6-O6
6	J	2	NAG	O5-C5-C6-O6

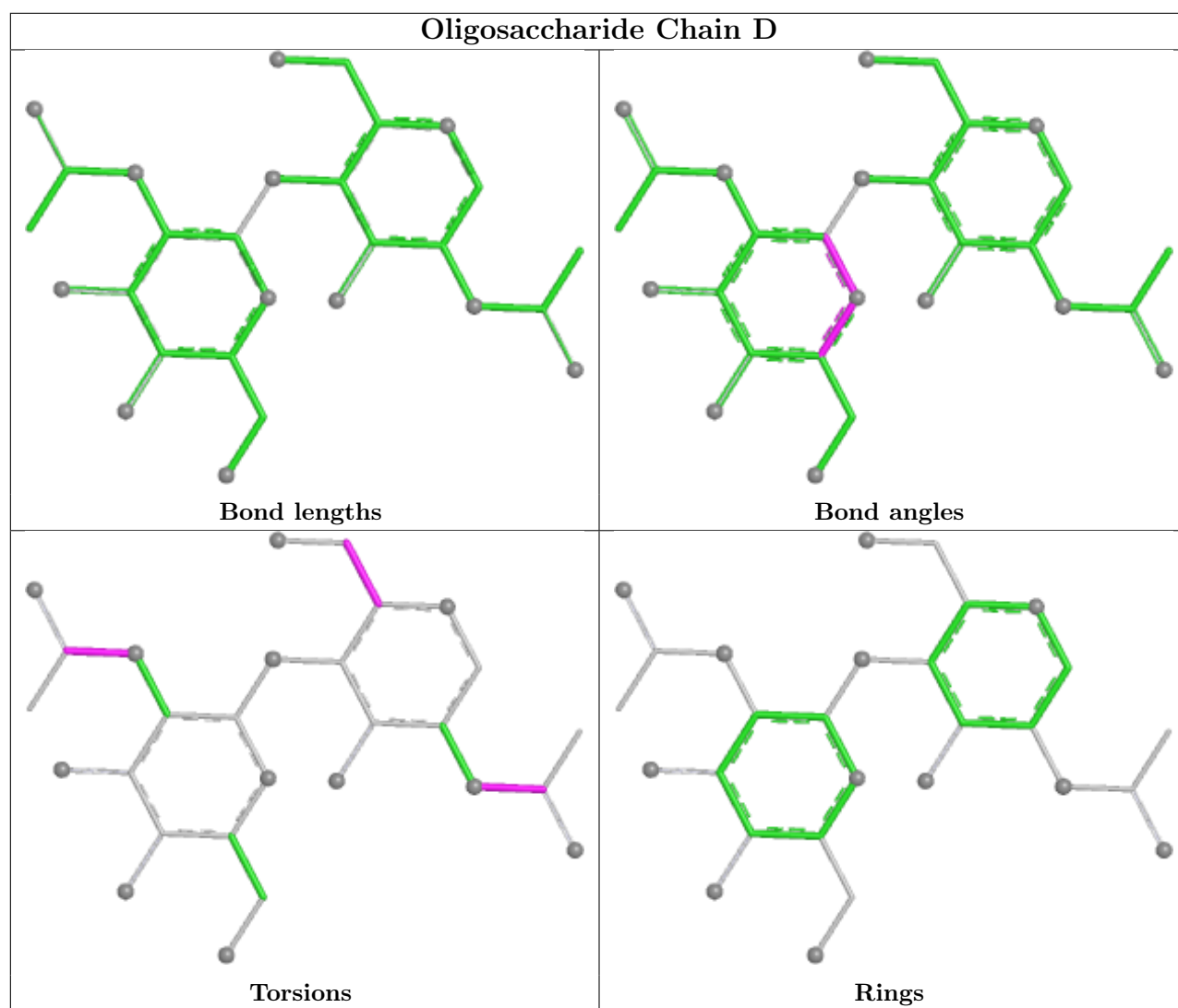
There are no ring outliers.

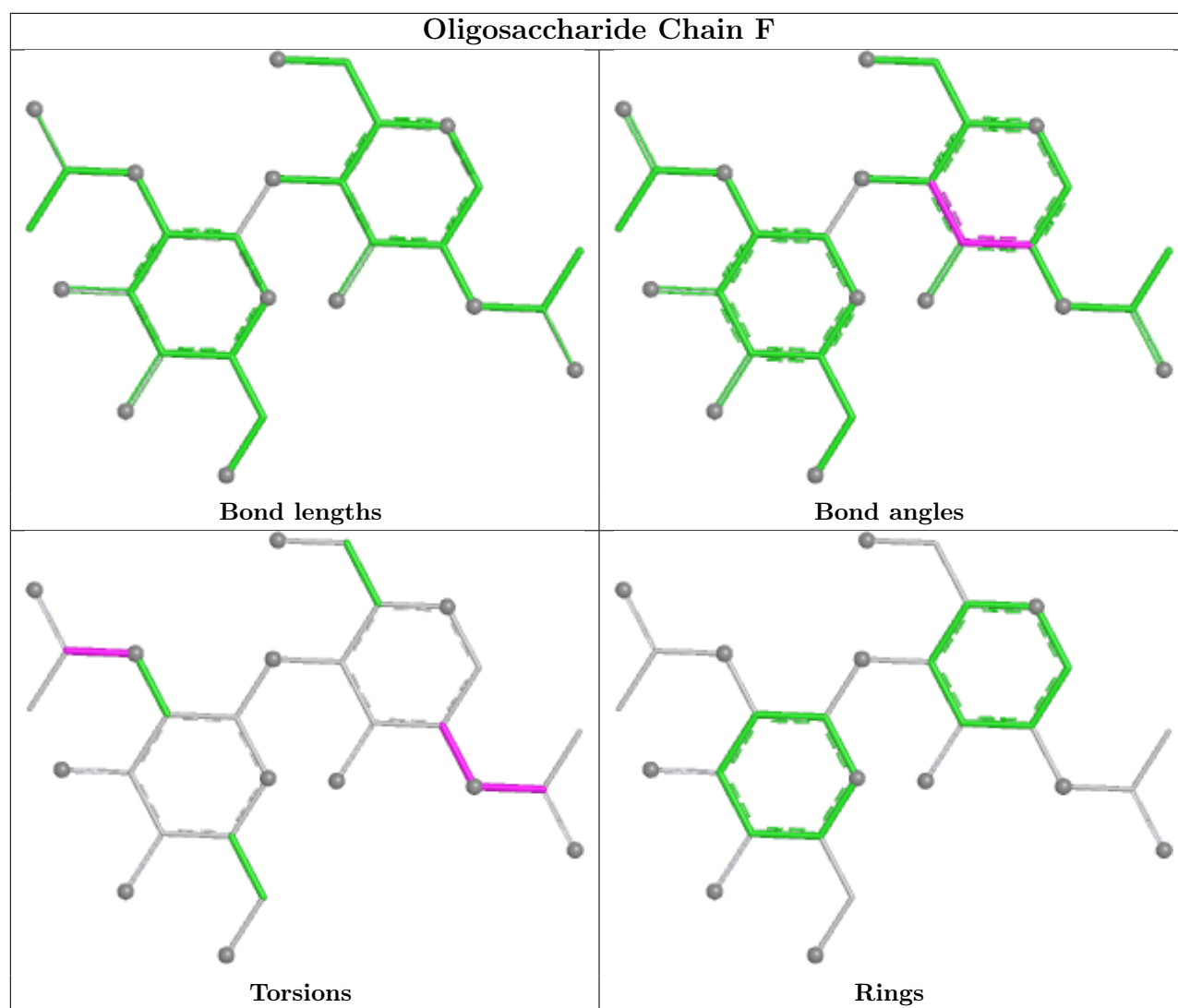
4 monomers are involved in 9 short contacts:

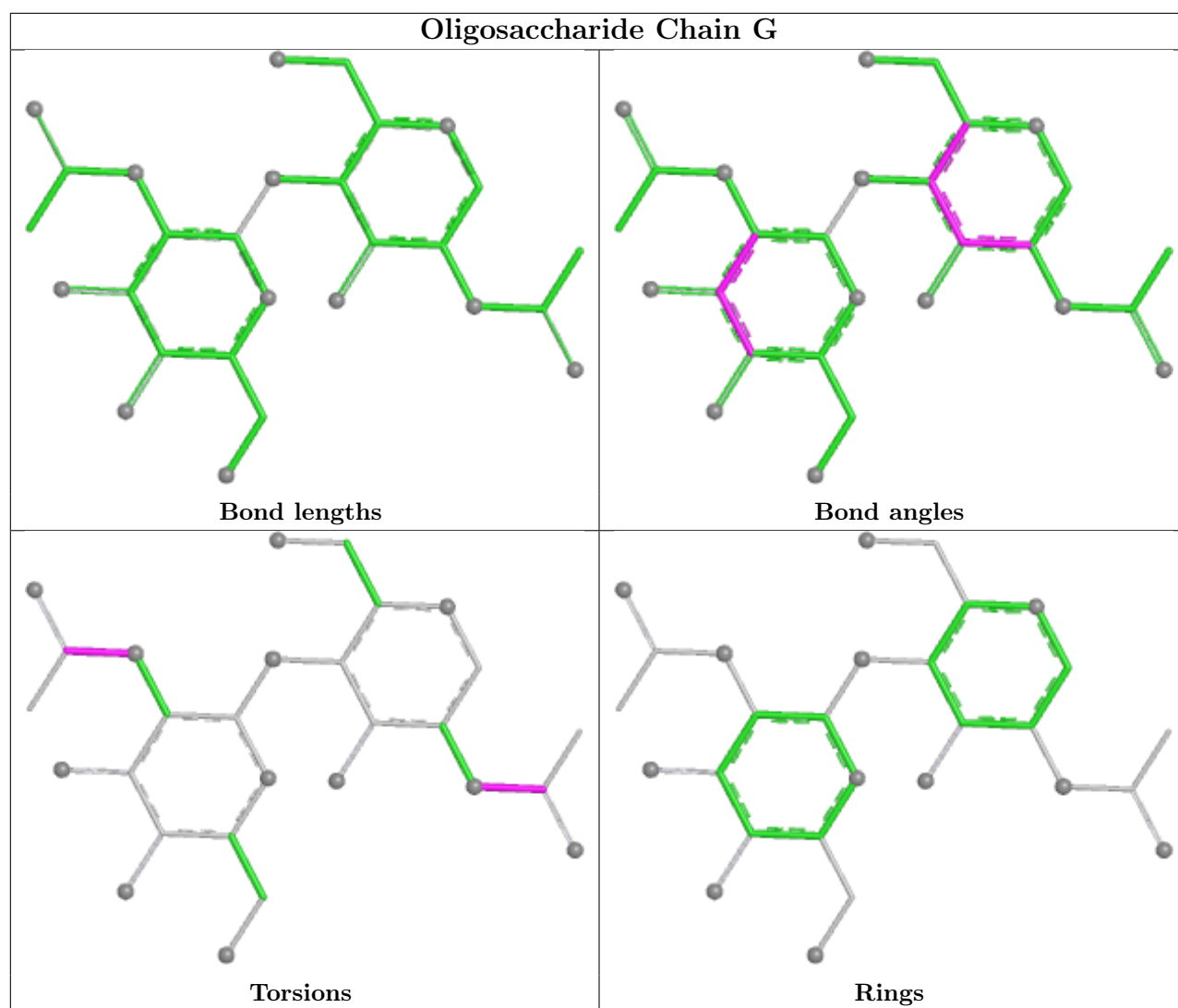
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	3	BMA	5	0
3	C	4	MAN	2	0
3	C	2	NAG	7	0
4	H	1	NAG	2	0

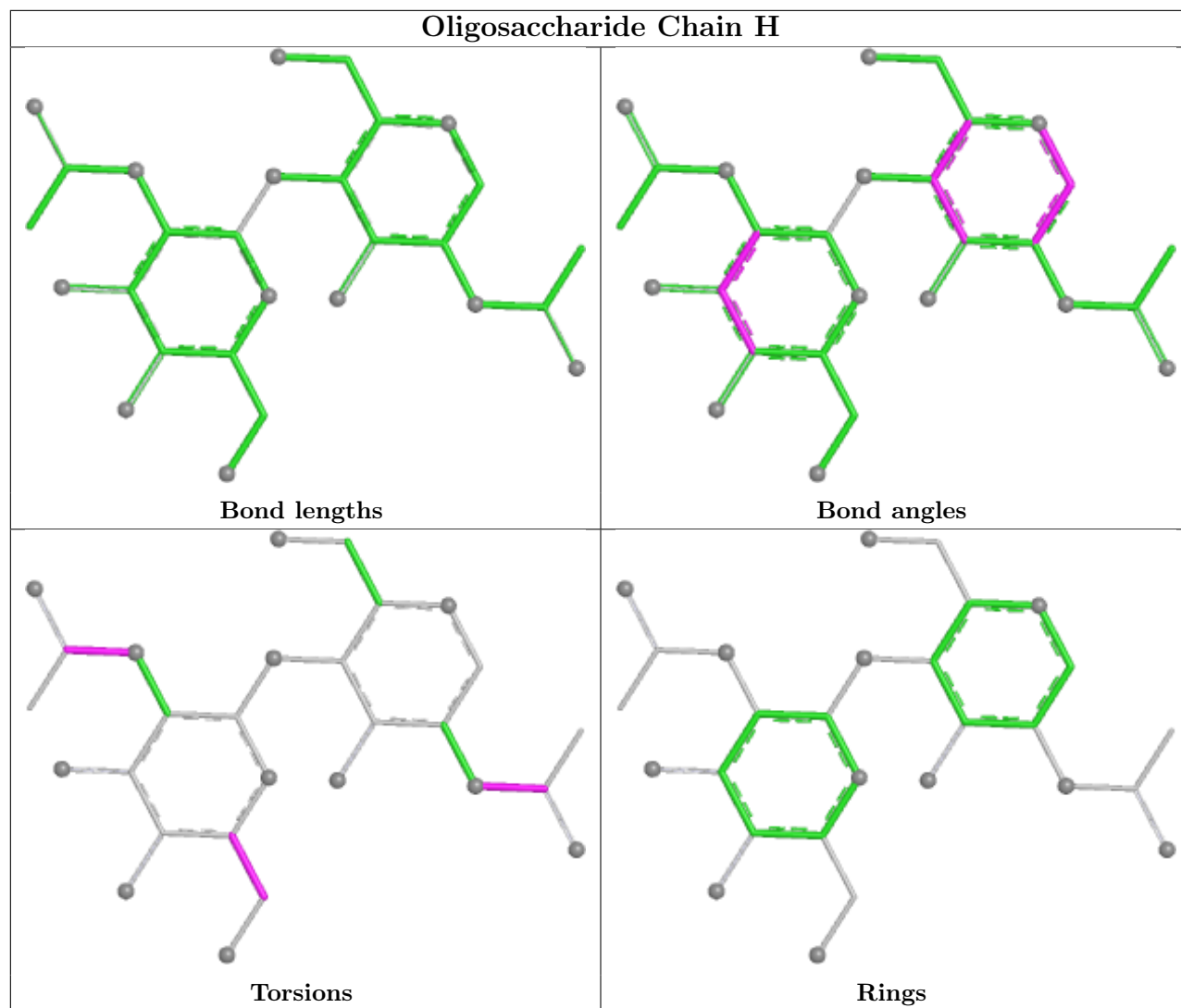
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

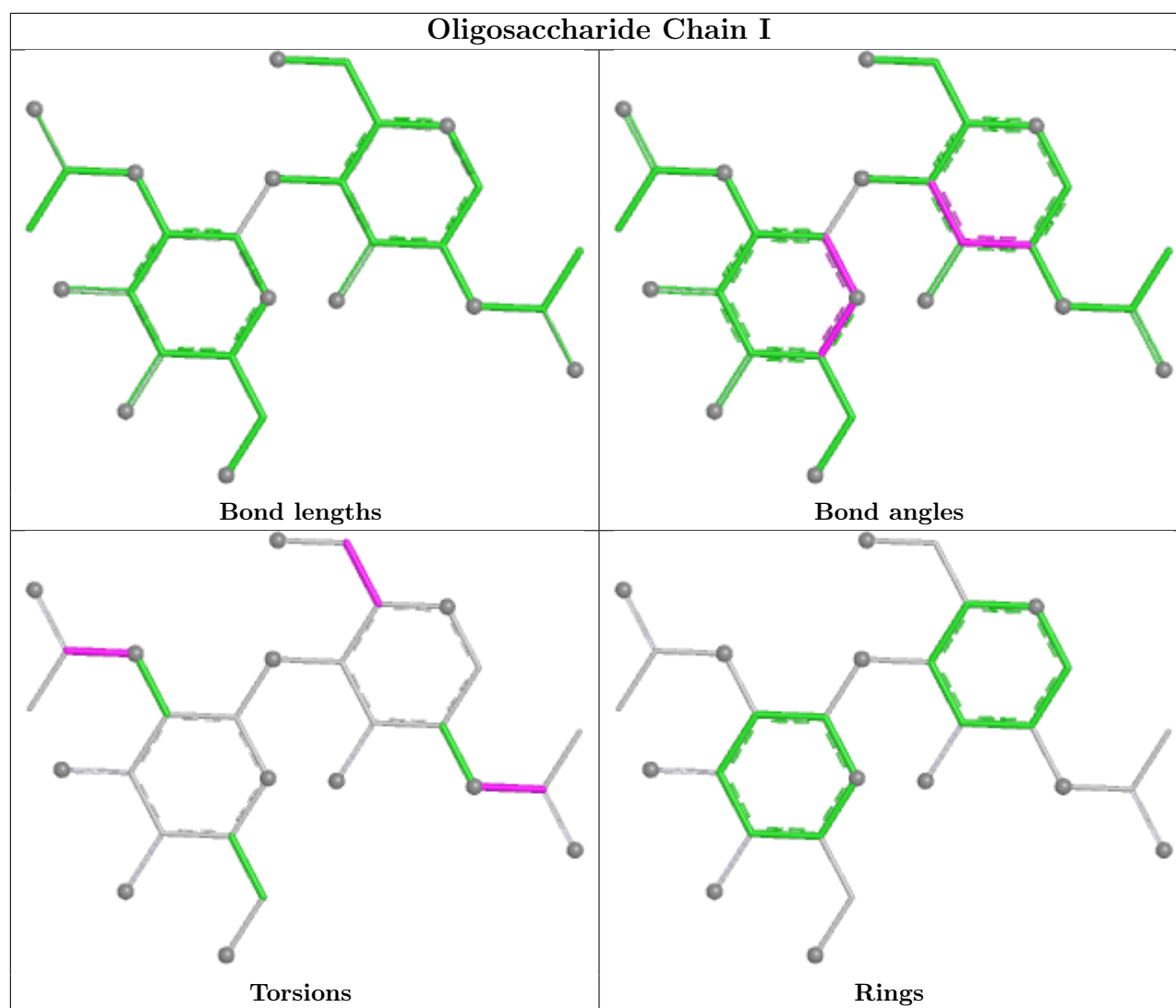


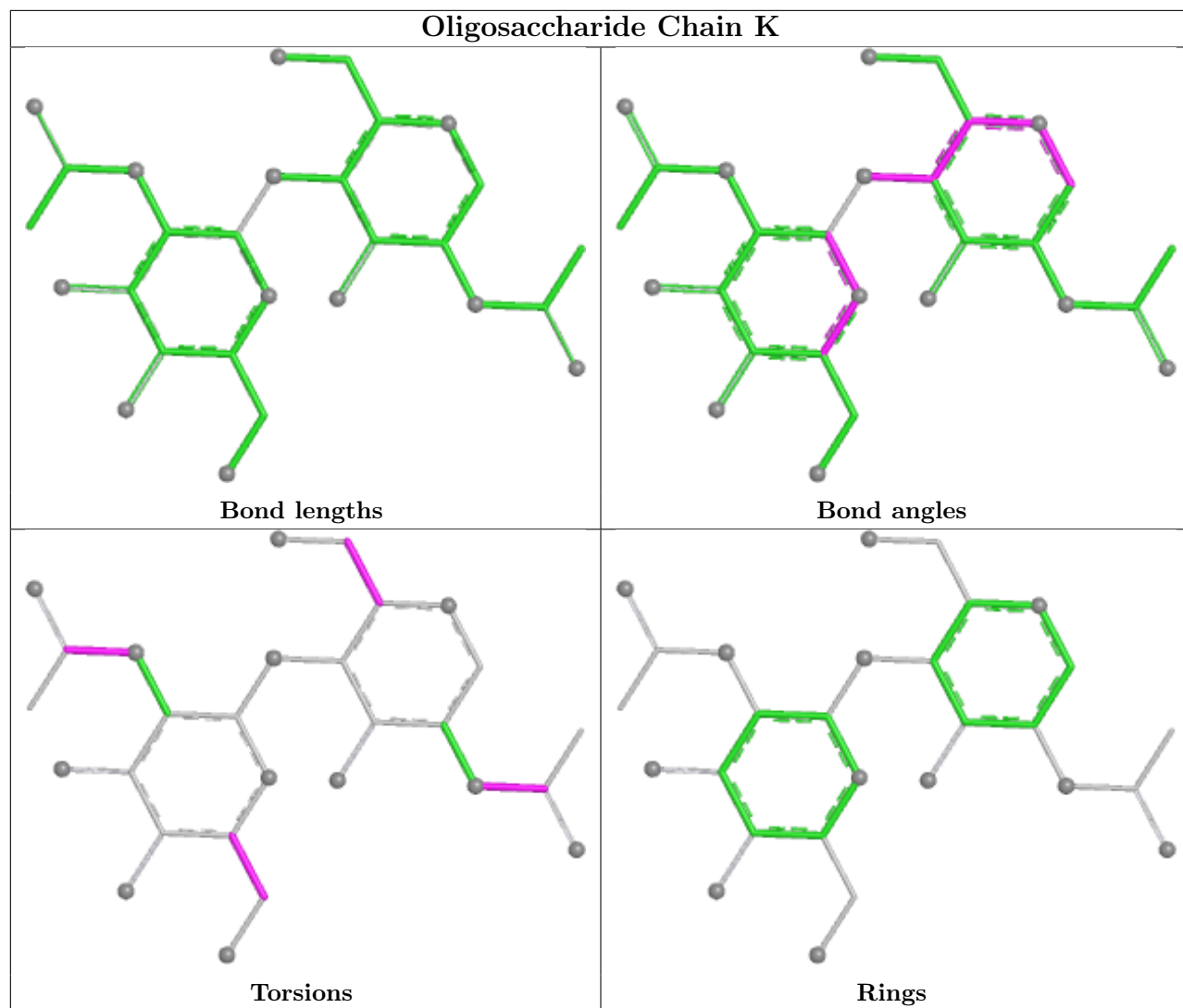




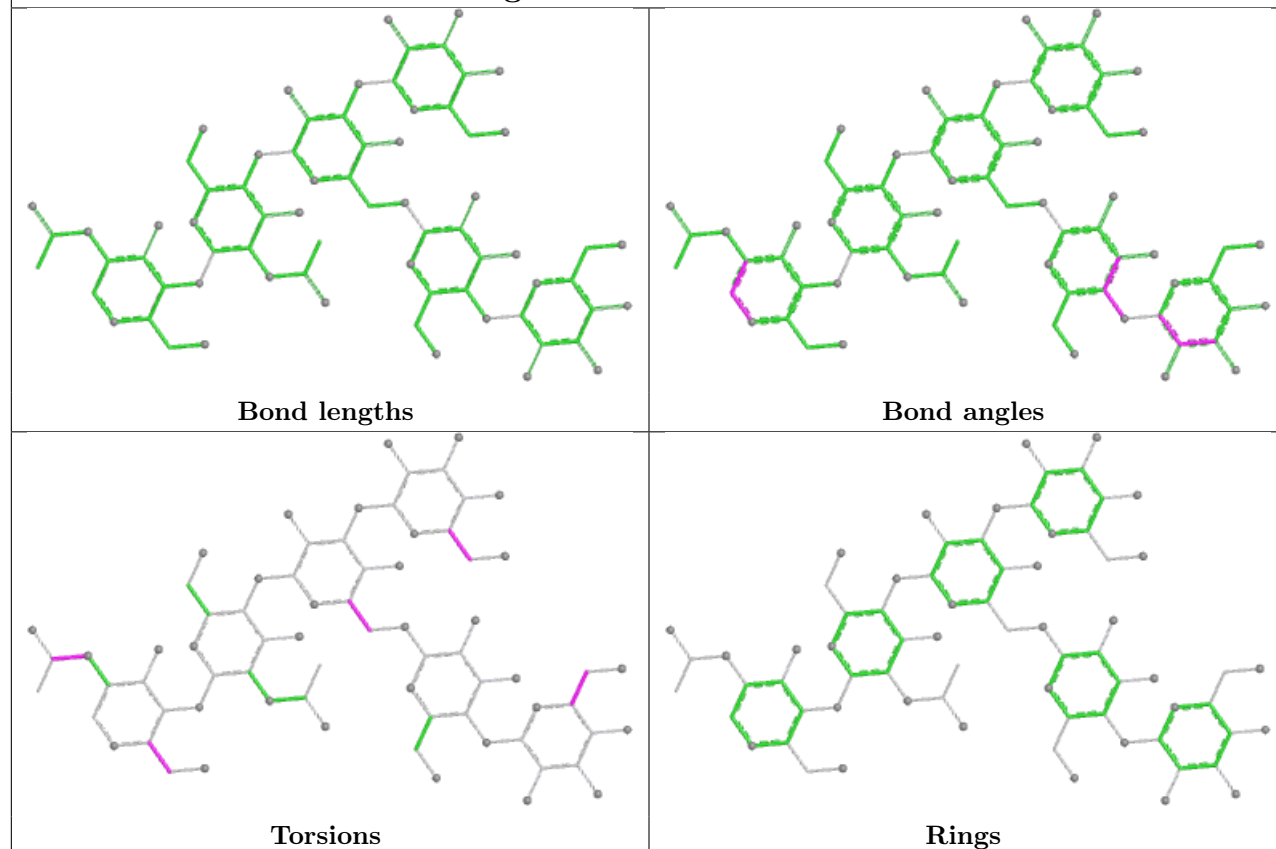




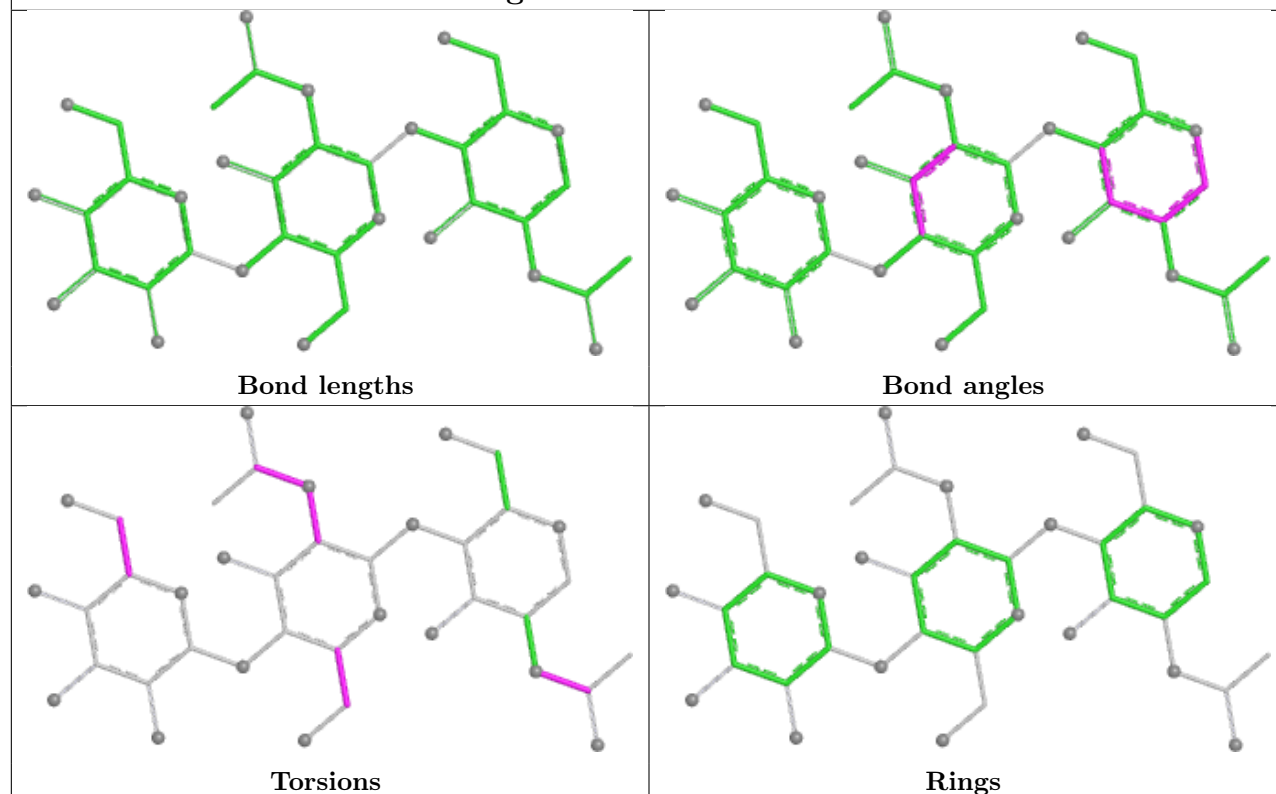


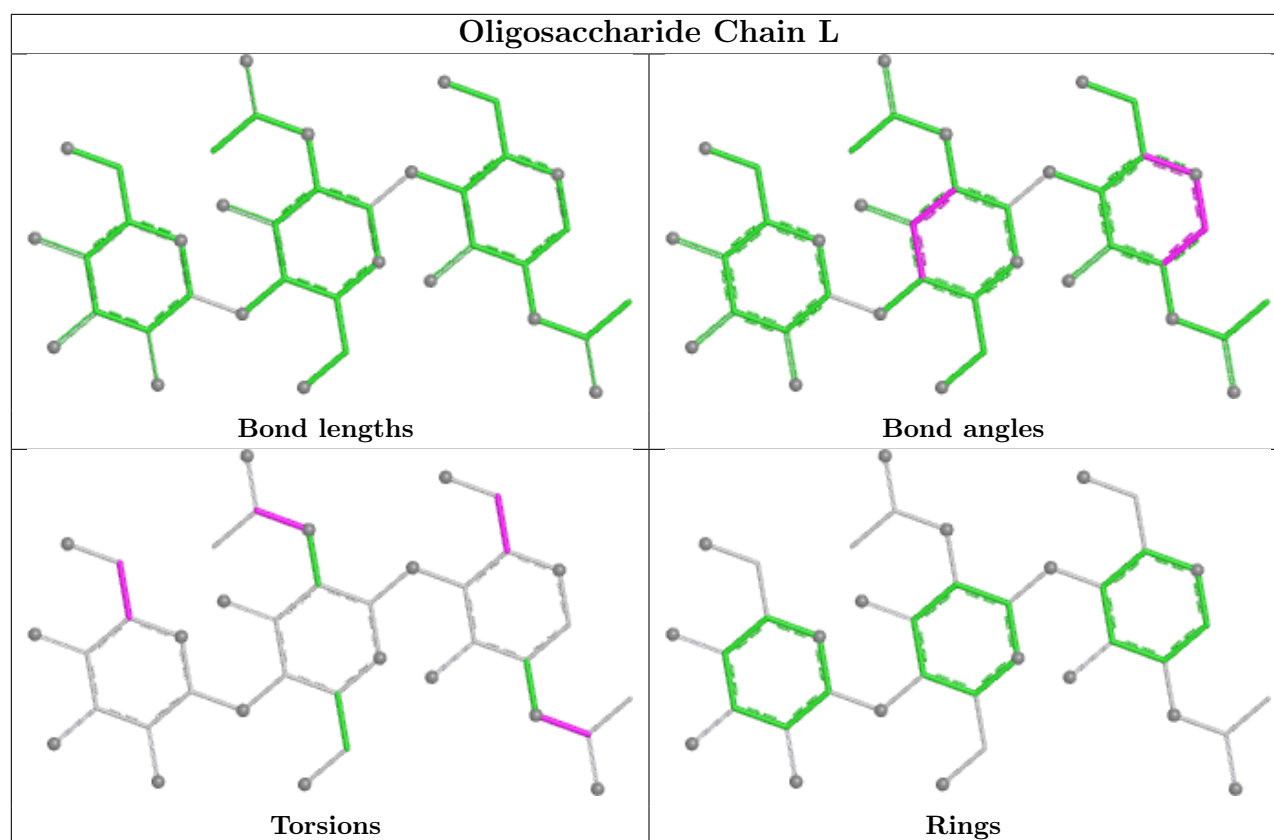


Oligosaccharide Chain E



Oligosaccharide Chain J





5.6 Ligand geometry [i](#)

Of 11 ligands modelled in this entry, 6 are monoatomic - leaving 5 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	NAG	B	3099	2	14,14,15	0.49	0	17,19,21	1.27	1 (5%)
7	NAG	B	3452	2	14,14,15	0.62	0	17,19,21	0.84	1 (5%)
7	NAG	B	3320	2	14,14,15	0.70	0	17,19,21	1.60	4 (23%)
7	NAG	A	2524	1	14,14,15	0.44	0	17,19,21	1.00	2 (11%)
7	NAG	A	2805	1	14,14,15	0.59	0	17,19,21	0.91	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the

Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns.
 '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
7	NAG	B	3099	2	-	4/6/23/26	0/1/1/1
7	NAG	B	3452	2	-	4/6/23/26	0/1/1/1
7	NAG	B	3320	2	-	4/6/23/26	0/1/1/1
7	NAG	A	2524	1	-	3/6/23/26	0/1/1/1
7	NAG	A	2805	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	B	3320	NAG	C4-C3-C2	4.26	117.25	111.02
7	B	3099	NAG	C1-O5-C5	3.78	117.26	112.19
7	B	3320	NAG	O5-C5-C6	2.80	113.11	107.66
7	A	2524	NAG	C1-O5-C5	2.78	115.91	112.19
7	B	3452	NAG	C1-O5-C5	2.58	115.64	112.19
7	B	3320	NAG	C1-O5-C5	2.37	115.36	112.19
7	B	3320	NAG	C3-C4-C5	2.37	114.52	110.23
7	A	2524	NAG	C3-C4-C5	2.01	113.88	110.23

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	2524	NAG	C8-C7-N2-C2
7	A	2524	NAG	O7-C7-N2-C2
7	B	3099	NAG	C8-C7-N2-C2
7	B	3099	NAG	O7-C7-N2-C2
7	B	3320	NAG	C8-C7-N2-C2
7	B	3320	NAG	O7-C7-N2-C2
7	B	3452	NAG	C8-C7-N2-C2
7	B	3452	NAG	O7-C7-N2-C2
7	B	3452	NAG	O5-C5-C6-O6
7	B	3099	NAG	O5-C5-C6-O6
7	A	2805	NAG	C8-C7-N2-C2
7	B	3452	NAG	C4-C5-C6-O6
7	B	3099	NAG	C4-C5-C6-O6
7	A	2805	NAG	O7-C7-N2-C2
7	B	3320	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
7	A	2524	NAG	C4-C5-C6-O6
7	B	3320	NAG	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	B	3099	NAG	1	0
7	B	3452	NAG	1	0
7	B	3320	NAG	3	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	A	938/967 (97%)	0.28	40 (4%)	40	31	42, 51, 60, 123	0
2	B	695/695 (100%)	0.65	67 (9%)	13	11	44, 51, 61, 94	0
All	All	1633/1662 (98%)	0.44	107 (6%)	24	19	42, 51, 60, 123	0

All (107) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	957	PRO	6.4
2	B	39	ASP	5.5
1	A	868	ASP	5.2
2	B	21	PRO	5.1
1	A	963	PRO	4.8
2	B	694	LEU	4.7
2	B	693	ILE	4.5
2	B	691	PRO	4.5
2	B	695	VAL	4.5
1	A	965	TRP	4.4
1	A	958	ALA	4.3
1	A	869	ILE	4.3
2	B	692	ASP	4.3
2	B	648	THR	4.3
2	B	649	GLY	4.1
1	A	959	PRO	4.1
2	B	450	ASN	4.0
2	B	19	VAL	4.0
1	A	234	ASP	3.9
2	B	167	ILE	3.8
1	A	964	VAL	3.7
2	B	436	ALA	3.7
2	B	653	VAL	3.7
1	A	962	VAL	3.7

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Mol	Chain	Res	Type	RSRZ
1	A	836	SER	3.7
2	B	84	SER	3.6
1	A	961	PRO	3.6
2	B	509	HIS	3.5
1	A	960	MET	3.5
2	B	36	PRO	3.5
2	B	504	GLY	3.5
2	B	669	TYR	3.4
2	B	173	LEU	3.3
2	B	524	ASP	3.3
2	B	652	ALA	3.2
1	A	967	ILE	3.2
1	A	649	LEU	3.2
2	B	513	PHE	3.2
2	B	690	GLY	3.2
2	B	28	ASP	3.2
1	A	956	GLN	3.2
1	A	955	ILE	3.1
1	A	565	TYR	3.1
1	A	735	ASP	3.1
2	B	556	TYR	2.9
2	B	535	MET	2.9
1	A	838	LEU	2.9
2	B	448	CYS	2.8
1	A	116	THR	2.8
2	B	562	THR	2.8
2	B	553	TRP	2.7
2	B	157	VAL	2.7
2	B	168	SER	2.7
2	B	460	CYS	2.6
2	B	501	CYS	2.6
2	B	37	ARG	2.6
1	A	451	PRO	2.6
2	B	162	SER	2.6
1	A	786	GLY	2.6
2	B	467	LEU	2.6
2	B	33	LEU	2.5
1	A	734	VAL	2.5
1	A	749	SER	2.5
2	B	557	TYR	2.5
1	A	484	LYS	2.5
1	A	686	PRO	2.5

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Mol	Chain	Res	Type	RSRZ
2	B	1	GLY	2.5
2	B	446	HIS	2.5
2	B	74	SER	2.5
1	A	966	VAL	2.5
2	B	432	ASP	2.4
1	A	118	MET	2.4
1	A	837	SER	2.4
1	A	701	VAL	2.3
2	B	49	CYS	2.3
2	B	551	SER	2.3
2	B	85	PRO	2.3
1	A	506	GLY	2.3
2	B	686	GLU	2.3
1	A	953	TRP	2.3
2	B	17	LEU	2.3
1	A	450	TYR	2.3
1	A	637	GLY	2.3
1	A	569	ALA	2.2
2	B	444	ASN	2.2
2	B	674	SER	2.2
2	B	688	PRO	2.2
2	B	519	LYS	2.2
1	A	472	CYS	2.2
2	B	455	PHE	2.2
2	B	169	PRO	2.2
2	B	166	TYR	2.2
1	A	571	THR	2.1
2	B	30	ALA	2.1
1	A	912	GLU	2.1
2	B	629	ASN	2.1
2	B	496	SER	2.1
2	B	647	ASP	2.1
2	B	468	GLY	2.1
2	B	673	SER	2.0
2	B	482	GLN	2.0
2	B	32	PRO	2.0
2	B	464	PRO	2.0
1	A	935	ASN	2.0
2	B	590	GLN	2.0
2	B	48	ASN	2.0
2	B	651	ASP	2.0

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 ⓘ

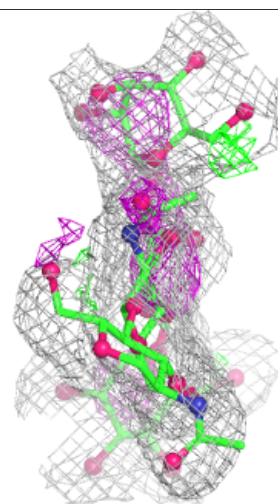
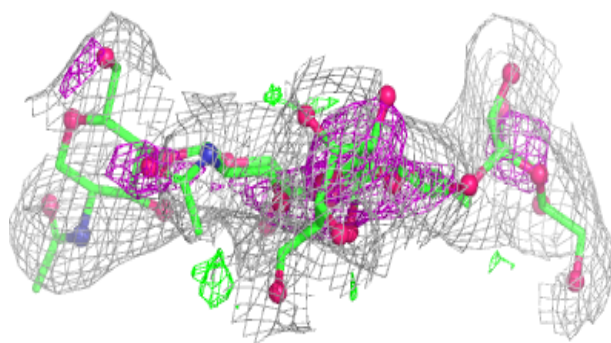
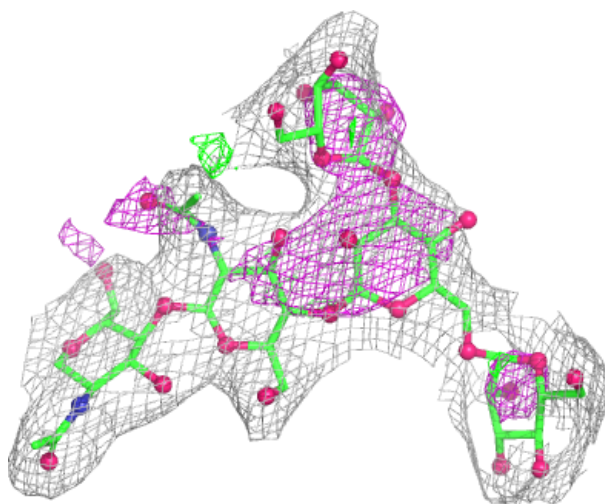
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	I	2	14/15	0.22	0.16	92,93,94,94	0
4	NAG	H	2	14/15	0.24	0.17	110,111,111,111	0
6	BMA	J	3	11/12	0.27	0.15	105,106,106,107	0
3	BMA	C	3	11/12	0.30	0.15	97,100,102,103	0
3	MAN	C	5	11/12	0.33	0.15	104,104,104,104	0
3	MAN	C	4	11/12	0.33	0.15	104,105,105,105	0
4	NAG	G	2	14/15	0.35	0.17	91,93,93,94	0
4	NAG	K	2	14/15	0.36	0.17	83,85,86,86	0
6	NAG	J	2	14/15	0.44	0.15	98,101,102,104	0
5	MAN	E	6	11/12	0.44	0.13	91,92,93,93	0
4	NAG	I	1	14/15	0.56	0.13	80,83,86,89	0
4	NAG	D	2	14/15	0.57	0.15	85,87,88,88	0
3	NAG	C	2	14/15	0.58	0.15	81,85,88,93	0
4	NAG	G	1	14/15	0.58	0.15	79,84,86,89	0
4	NAG	F	2	14/15	0.60	0.15	84,86,86,86	0
4	NAG	F	1	14/15	0.60	0.15	72,77,79,81	0
5	MAN	E	5	11/12	0.61	0.13	97,98,98,98	0
4	NAG	K	1	14/15	0.66	0.12	71,76,78,80	0
6	BMA	L	3	11/12	0.66	0.11	88,90,90,90	0
5	MAN	E	4	11/12	0.67	0.14	93,95,95,97	0
5	BMA	E	3	11/12	0.67	0.12	83,87,89,91	0
4	NAG	H	1	14/15	0.72	0.12	106,107,108,110	0
4	NAG	D	1	14/15	0.73	0.13	74,78,80,82	0
6	NAG	J	1	14/15	0.79	0.11	85,89,91,95	0
6	NAG	L	2	14/15	0.80	0.12	82,84,86,87	0
3	NAG	C	1	14/15	0.87	0.09	61,64,69,75	0
6	NAG	L	1	14/15	0.87	0.11	69,73,75,78	0
5	NAG	E	1	14/15	0.91	0.07	58,59,62,65	0
5	NAG	E	2	14/15	0.91	0.09	65,69,74,79	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.

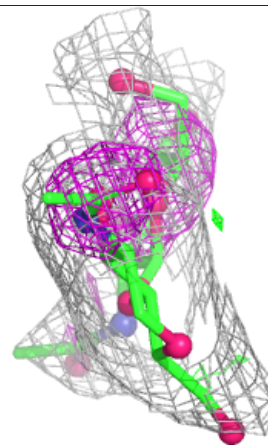
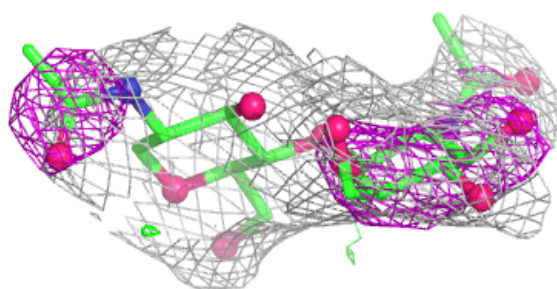
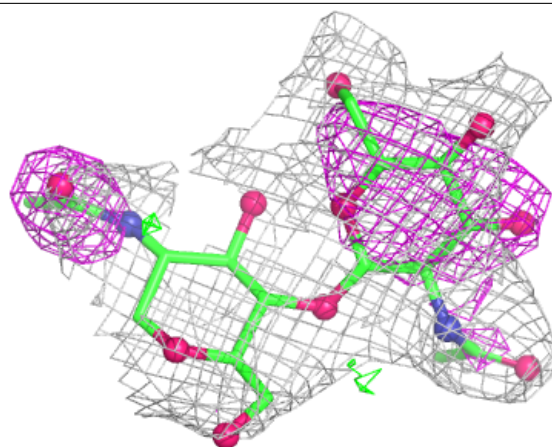
Electron density around Chain C:

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



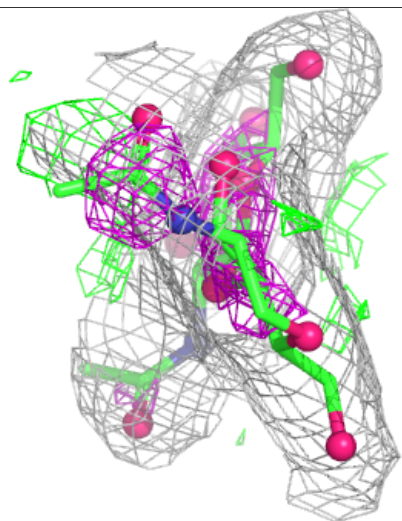
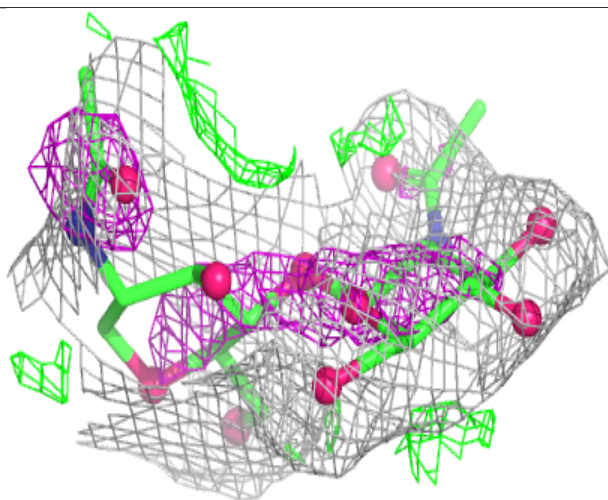
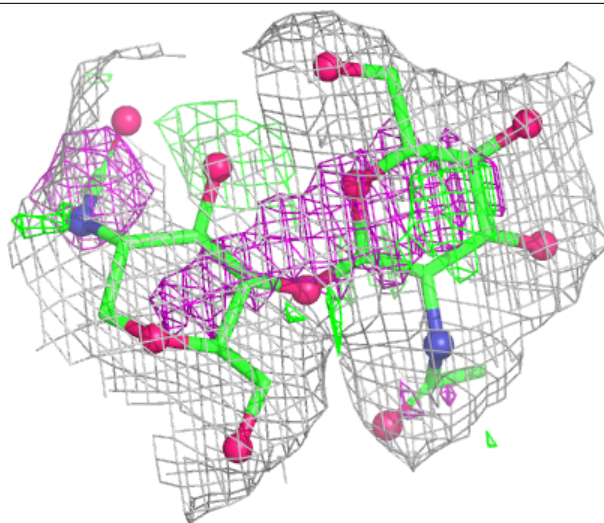
Electron density around Chain D:

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



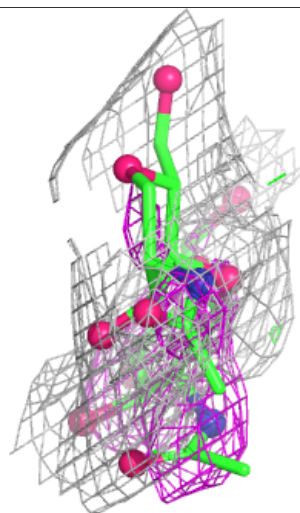
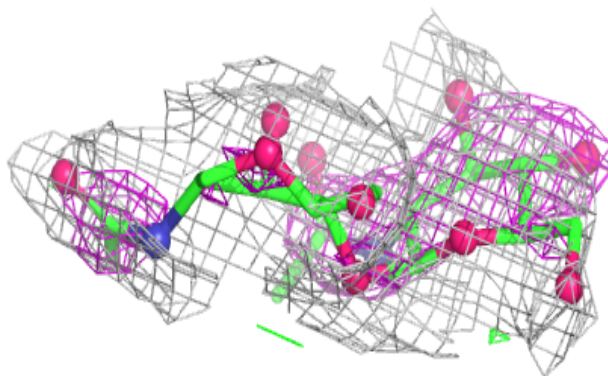
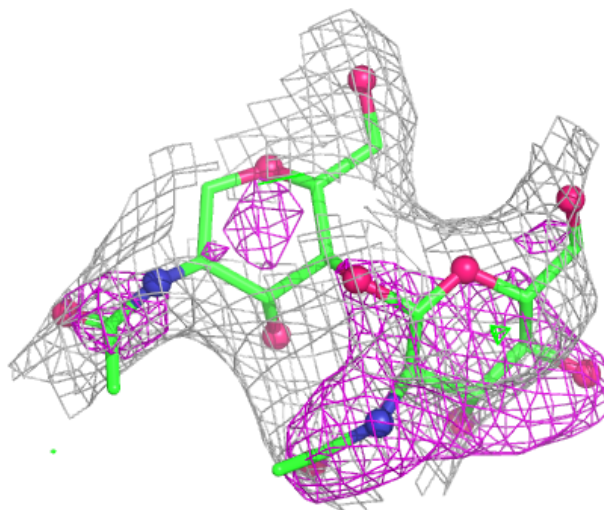
Electron density around Chain F:

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



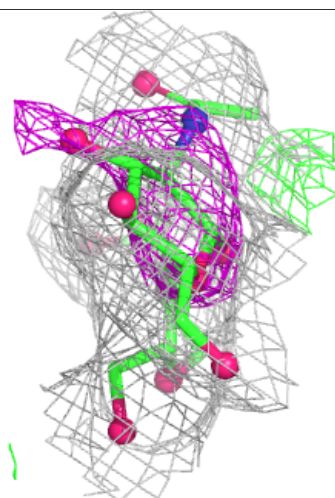
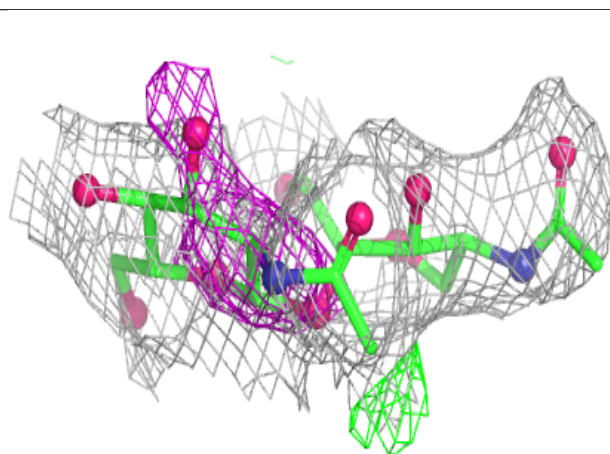
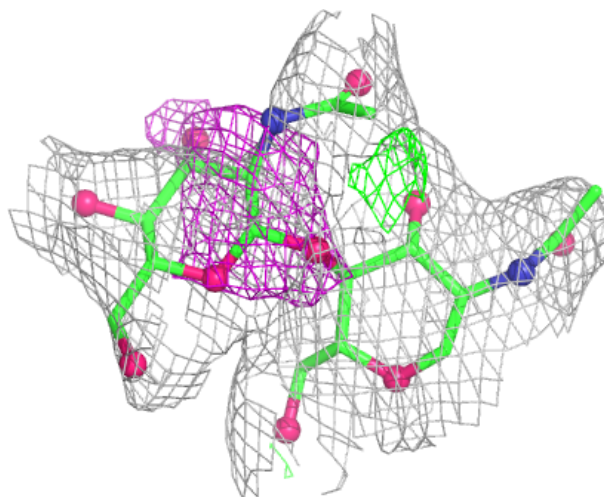
Electron density around Chain G:

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



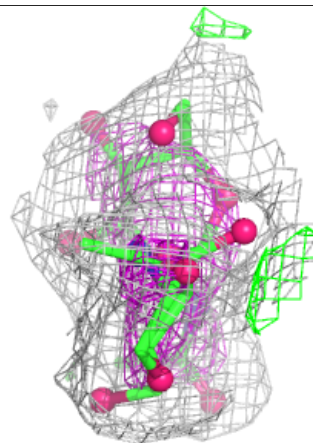
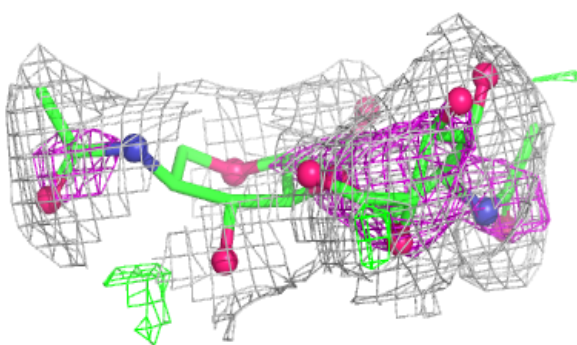
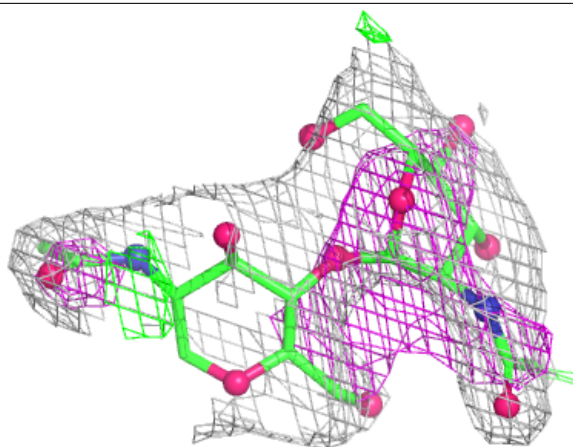
Electron density around Chain H:

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



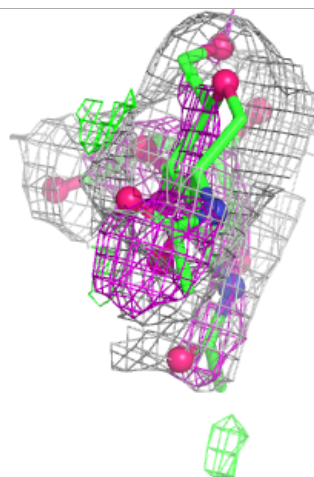
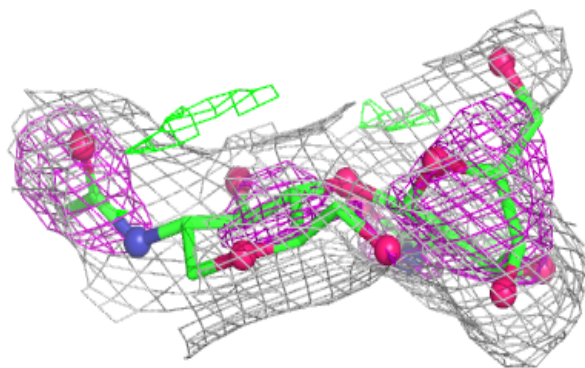
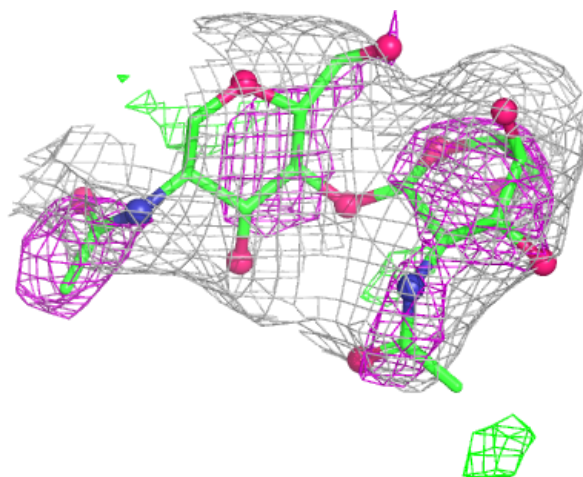
Electron density around Chain I:

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



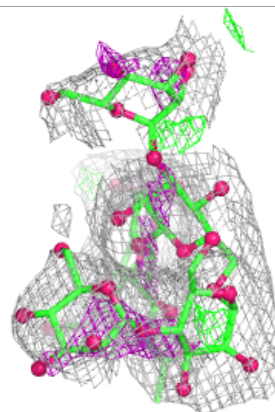
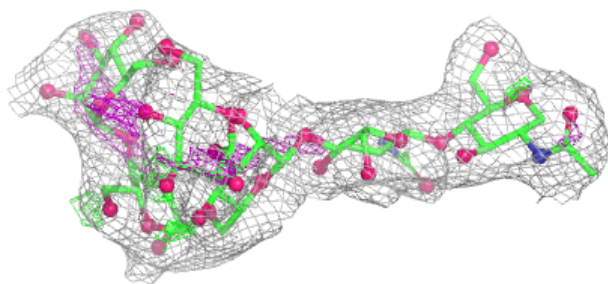
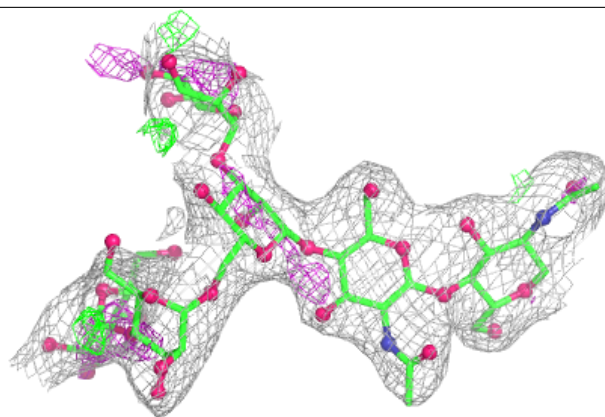
Electron density around Chain K:

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

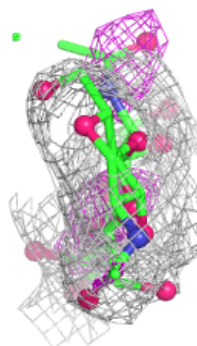
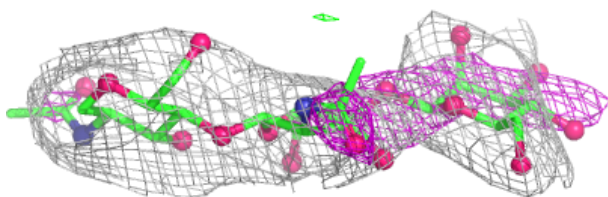
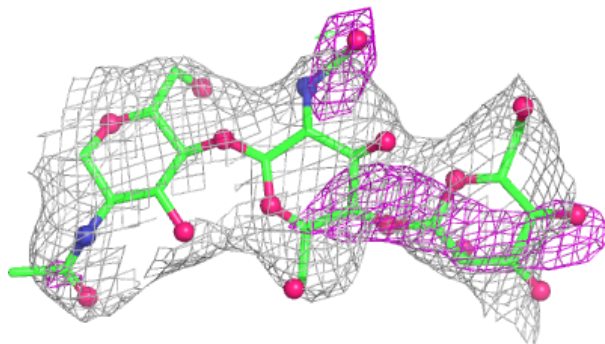


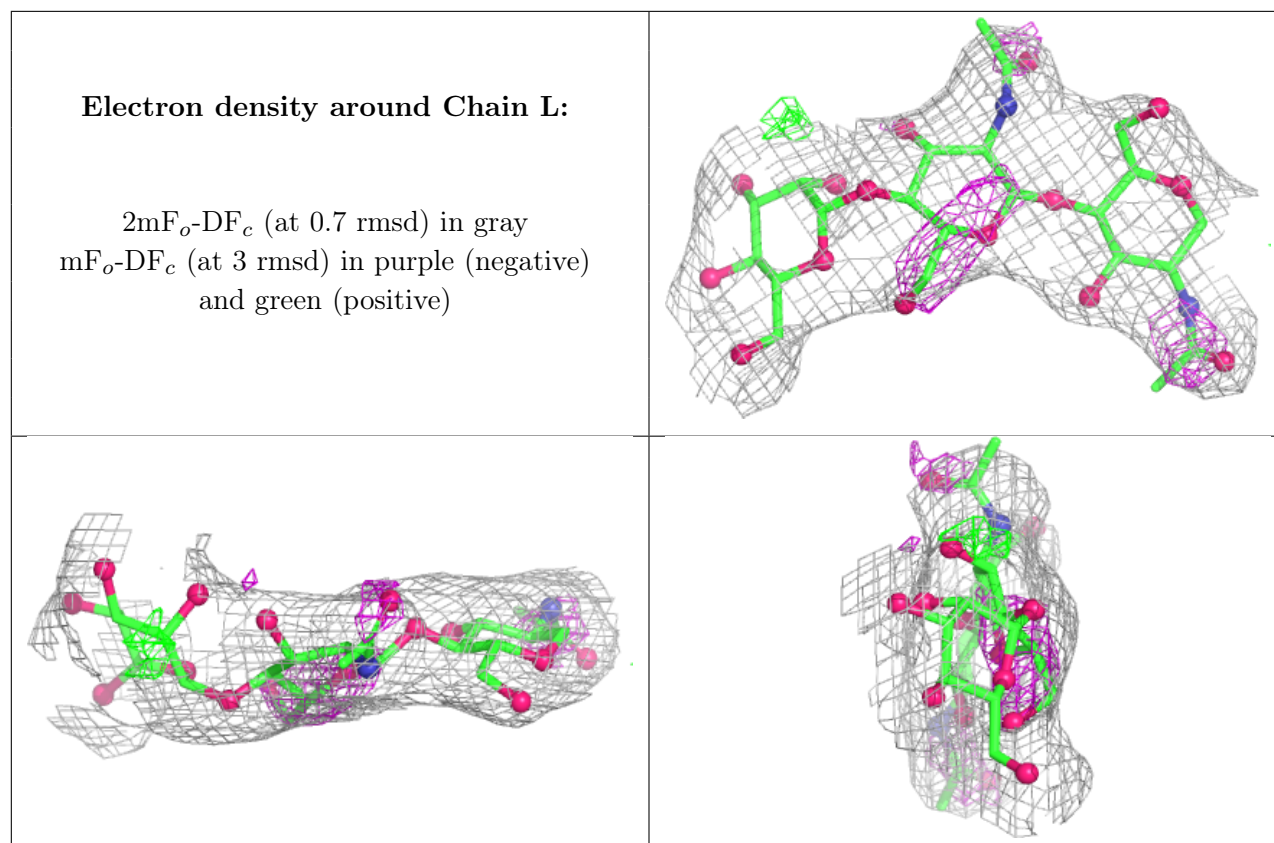
Electron density around Chain E:

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

**Electron density around Chain J:**

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	NAG	A	2524	14/15	0.22	0.19	81,86,87,87	0
7	NAG	B	3452	14/15	0.23	0.18	86,90,91,91	0
7	NAG	A	2805	14/15	0.30	0.17	74,79,79,80	0
7	NAG	B	3320	14/15	0.45	0.17	68,71,72,73	0
7	NAG	B	3099	14/15	0.45	0.16	74,79,80,81	0
8	CA	A	4008	1/1	0.93	0.11	32,32,32,32	0
8	CA	A	4007	1/1	0.95	0.04	34,34,34,34	0
8	CA	A	4006	1/1	0.96	0.06	59,59,59,59	0
8	CA	A	4004	1/1	0.97	0.04	78,78,78,78	0
8	CA	B	4002	1/1	0.97	0.03	55,55,55,55	0
8	CA	A	4005	1/1	0.98	0.02	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.