



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 11:34 PM UTC

PDB ID : 1U8E / pdb_00001u8e
Title : HUMAN DIPEPTIDYL PEPTIDASE IV/CD26 MUTANT Y547F
Authors : Bjelke, J.R.; Christensen, J.; Branner, S.; Wagtmann, N.; Olsen, C.; Kanstrup, A.B.; Rasmussen, H.B.
Deposited on : 2004-08-05
Resolution : 2.20 Å(reported)

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

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

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

with specific help available everywhere you see the ⓘ 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

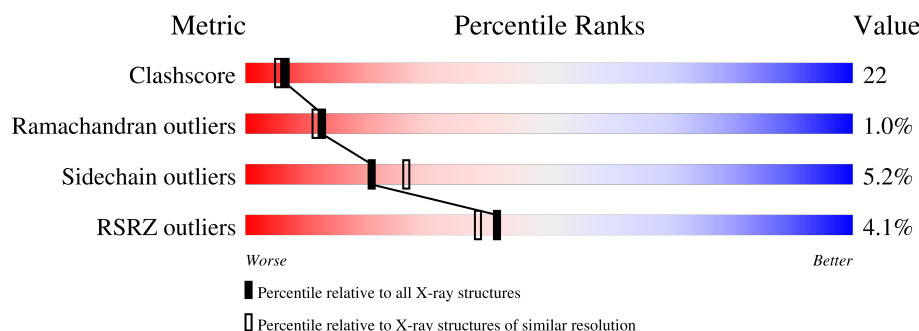
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



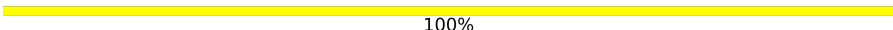
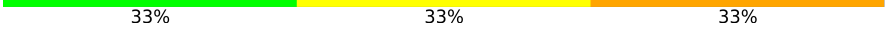
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. 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	728	3% 64% 31% .
1	B	728	5% 61% 34% 5%
2	C	3	33% 67%
2	D	3	33% 67%
2	I	3	33% 67%
3	E	2	100%

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

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
2	NAG	D	1	X	-	-	-
3	NAG	E	1	X	-	-	-
3	NAG	J	1	X	-	-	-
5	NAG	K	1	X	-	-	-
7	NAG	M	1	X	-	-	-
8	NAG	B	781	-	-	X	-

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 13182 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.

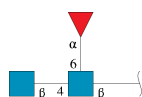
- Molecule 1 is a protein called Dipeptidyl peptidase IV.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	726	Total	C	N	O	S	0	0	0
			5946	3816	980	1124	26			
1	B	728	Total	C	N	O	S	0	0	0
			5961	3827	982	1126	26			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	547	PHE	TYR	engineered mutation	UNP P27487
B	547	PHE	TYR	engineered mutation	UNP P27487

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



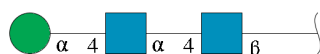
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	D	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	I	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



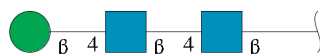
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	J	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



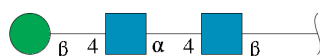
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	F	3	Total	C	N	O	0	0	0
			39	22	2	15			

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



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

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



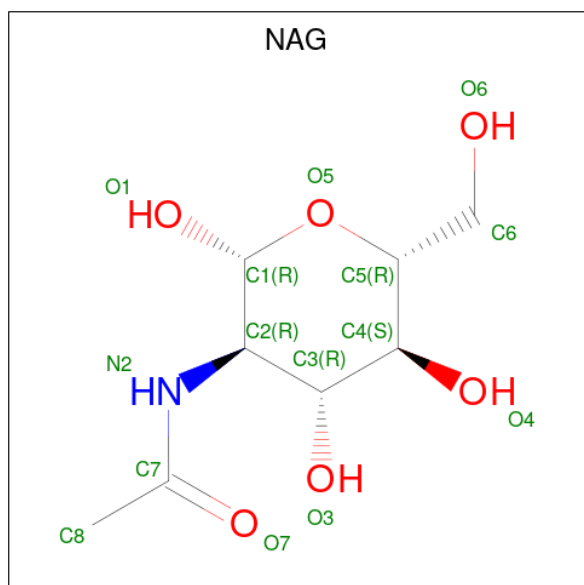
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	L	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 7 is an oligosaccharide called 2-acetamido-2-deoxy- α -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose.



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

- Molecule 8 is 2-acetamido-2-deoxy- β -D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

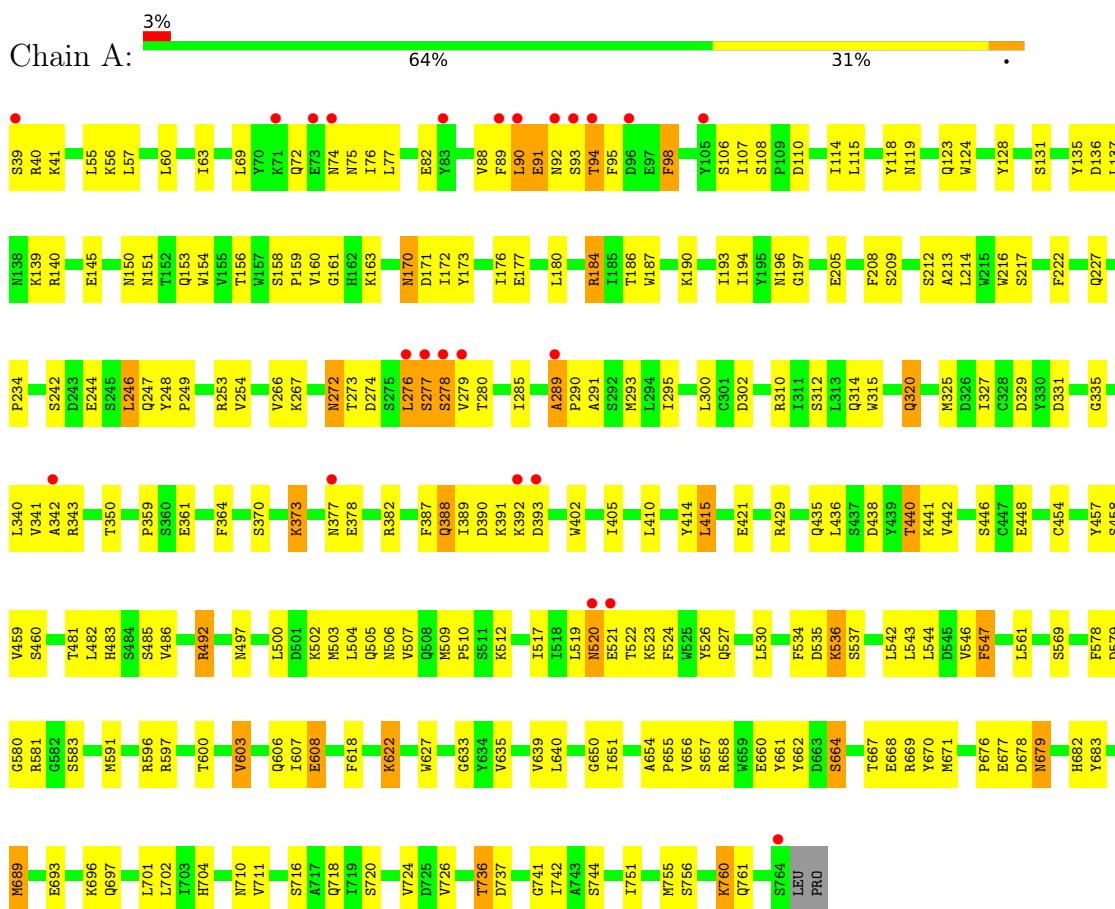
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	448	Total	O	0	0
			448	448		
9	B	414	Total	O	0	0
			414	414		

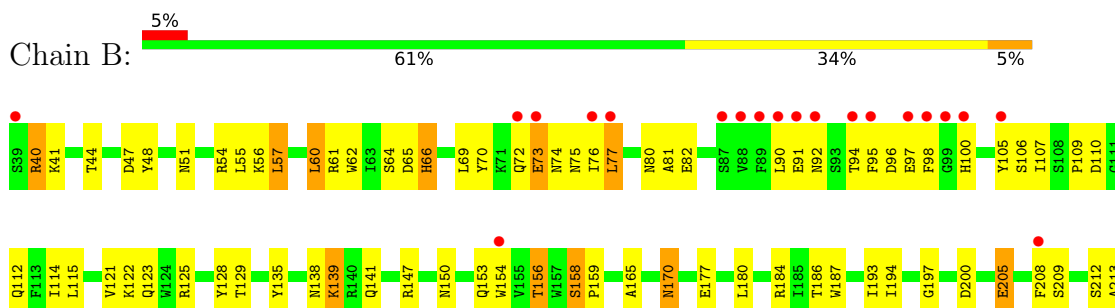
3 Residue-property plots

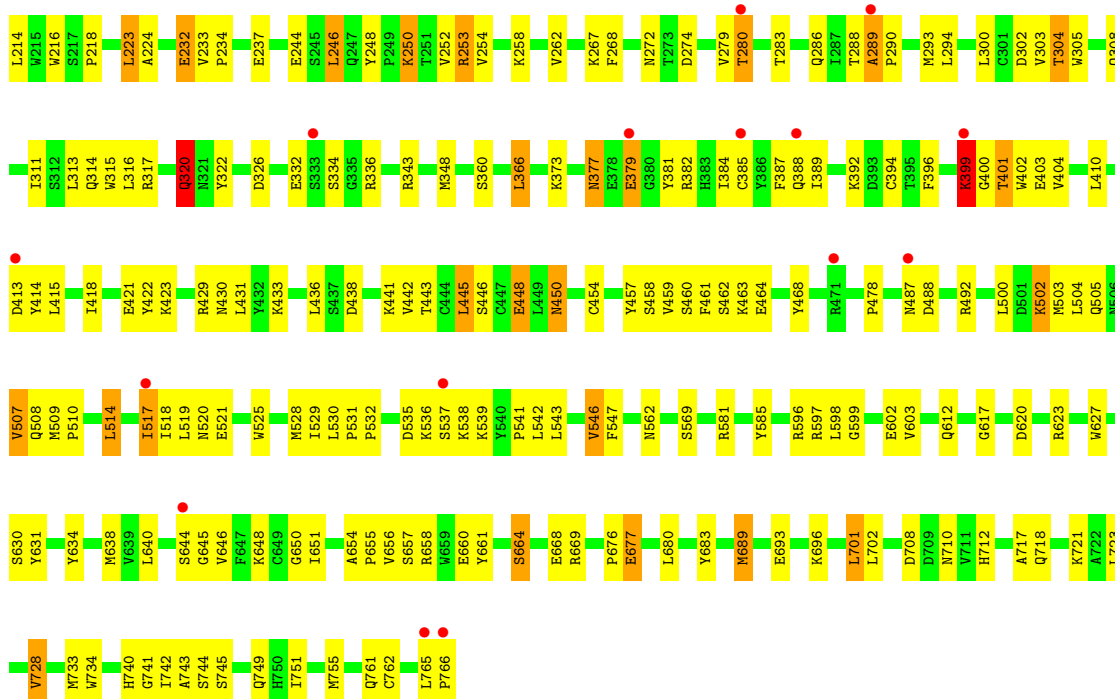
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Dipeptidyl peptidase IV



• Molecule 1: Dipeptidyl peptidase IV





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

Chain C: 33% 67%



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

Chain D: 33% 67%



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

Chain I: 33% 67%



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

Chain E: 100%



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

Chain G:  100%



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

Chain H:  50% 50%



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

Chain J:  50% 50%



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

Chain F:  33% 67%



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

Chain K:  33% 67%



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

Chain L:  33% 33% 33%



- Molecule 7: 2-acetamido-2-deoxy- α -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain M: 

MDG1
MDG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	118.76Å 121.37Å 129.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.84 – 2.20 34.84 – 2.20	Depositor EDS
% Data completeness (in resolution range)	78.5 (34.84-2.20) 78.5 (34.84-2.20)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.52 (at 2.10Å)	Xtriage
Refinement program	CNX 2002	Depositor
R, R_{free}	0.215 , 0.268 (Not available) , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	20.6	Xtriage
Anisotropy	0.459	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.018 for k,h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	13182	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.57% 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, NDG, FUC, MAN, BMA

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	0.39	0/6117	0.92	18/8318 (0.2%)
1	B	0.38	0/6133	0.92	23/8341 (0.3%)
All	All	0.38	0/12250	0.92	41/16659 (0.2%)

There are no bond length outliers.

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	205	GLU	N-CA-C	6.90	118.59	111.14
1	A	454	CYS	N-CA-C	6.84	120.20	108.02
1	B	656	VAL	N-CA-C	-6.77	98.83	108.58
1	A	98	PHE	N-CA-C	-6.55	105.44	113.50
1	B	450	ASN	CA-C-N	6.53	127.05	119.47
1	B	450	ASN	C-N-CA	6.53	127.05	119.47
1	A	458	SER	N-CA-C	-6.37	99.62	109.24
1	B	458	SER	N-CA-C	-6.31	99.74	109.52
1	B	205	GLU	N-CA-C	6.24	118.16	111.36
1	A	91	GLU	N-CA-C	6.08	119.65	112.72
1	B	664	SER	N-CA-C	6.06	117.55	111.07
1	A	300	LEU	N-CA-C	-6.05	97.85	108.20
1	B	388	GLN	N-CA-C	-5.99	99.14	108.90
1	B	300	LEU	N-CA-C	-5.96	97.43	107.99
1	B	617	GLY	N-CA-C	5.87	123.17	115.36
1	B	454	CYS	N-CA-C	5.80	117.85	108.34
1	A	664	SER	N-CA-C	5.77	117.25	111.07
1	A	485	SER	N-CA-C	5.73	118.30	111.71
1	B	546	VAL	N-CA-C	5.67	118.31	108.95
1	B	233	VAL	N-CA-C	5.59	113.40	107.76
1	A	388	GLN	N-CA-C	-5.56	99.45	108.52
1	B	150	ASN	N-CA-C	-5.54	102.28	110.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	399	LYS	N-CA-C	5.54	122.60	110.80
1	A	414	TYR	N-CA-C	5.53	117.59	109.24
1	B	448	GLU	N-CA-C	5.50	120.10	113.17
1	B	158	SER	N-CA-C	-5.45	102.35	110.20
1	B	254	VAL	N-CA-C	5.44	114.08	107.61
1	B	320	GLN	N-CA-C	5.34	122.18	110.80
1	A	678	ASP	N-CA-C	5.32	116.58	108.12
1	A	160	VAL	N-CA-C	5.28	120.32	109.34
1	A	591	MET	N-CA-C	5.28	117.03	111.28
1	B	541	PRO	N-CA-C	-5.28	102.30	111.32
1	B	200	ASP	N-CA-C	-5.25	101.60	109.79
1	A	94	THR	N-CA-C	-5.24	106.39	112.89
1	A	744	SER	N-CA-C	-5.17	103.31	110.35
1	A	547	PHE	N-CA-C	-5.17	99.80	110.80
1	A	320	GLN	N-CA-C	5.06	121.58	110.80
1	A	656	VAL	N-CA-C	-5.06	101.18	108.87
1	B	743	ALA	N-CA-C	5.05	118.48	112.72
1	B	392	LYS	N-CA-C	5.04	117.67	111.82
1	B	744	SER	N-CA-C	-5.00	103.45	110.50

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5946	0	5656	245	0
1	B	5961	0	5673	264	0
2	C	38	0	34	4	0
2	D	38	0	34	3	0
2	I	38	0	34	3	0
3	E	28	0	25	0	0
3	G	28	0	25	2	0
3	H	28	0	25	1	0
3	J	28	0	25	1	0
4	F	39	0	33	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	K	39	0	34	0	0
6	L	39	0	33	1	0
7	M	28	0	24	2	0
8	A	14	0	13	5	0
8	B	28	0	26	10	0
9	A	448	0	0	23	0
9	B	414	0	0	29	0
All	All	13182	0	11694	520	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 22.

All (520) 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:492:ARG:HH11	1:A:492:ARG:HB3	1.25	0.97
1:B:289:ALA:HB1	1:B:290:PRO:HA	1.48	0.93
1:A:289:ALA:HB1	1:A:290:PRO:HA	1.49	0.92
8:B:781:NAG:O7	8:B:781:NAG:H3	1.70	0.89
1:A:253:ARG:HH22	1:B:253:ARG:HH22	1.22	0.87
1:A:172:ILE:H	1:A:186:THR:HG22	1.42	0.85
1:A:194:ILE:HD13	4:F:1:NAG:H82	1.60	0.84
1:B:90:LEU:HD11	1:B:94:THR:HG21	1.57	0.84
1:B:289:ALA:HB1	1:B:290:PRO:CA	2.08	0.84
1:A:289:ALA:HB1	1:A:290:PRO:CA	2.09	0.83
1:B:77:LEU:HD23	1:B:77:LEU:H	1.42	0.82
1:A:676:PRO:HG2	1:A:677:GLU:OE2	1.80	0.80
1:A:176:ILE:HD11	1:A:276:LEU:HD21	1.62	0.79
1:A:272:ASN:ND2	1:A:274:ASP:H	1.79	0.79
1:A:377:ASN:HB3	9:A:783:HOH:O	1.82	0.79
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.64	0.79
1:A:312:SER:HB2	1:A:325:MET:HE3	1.65	0.79
1:B:72:GLN:HE21	1:B:77:LEU:HD21	1.46	0.78
1:B:74:ASN:HB2	9:B:927:HOH:O	1.83	0.78
2:C:1:NAG:H62	2:C:2:NAG:H82	1.65	0.78
1:B:258:LYS:NZ	1:B:712:HIS:HD2	1.82	0.77
1:B:399:LYS:HD3	9:B:843:HOH:O	1.83	0.77
1:B:112:GLN:HB3	1:B:138:ASN:HD21	1.48	0.77
1:A:184:ARG:NH1	1:A:187:TRP:HA	1.99	0.76
1:B:502:LYS:O	1:B:505:GLN:HG2	1.86	0.76
1:B:644:SER:O	1:B:646:VAL:N	2.19	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:399:LYS:HB2	1:B:402:TRP:CZ2	2.22	0.75
2:C:2:NAG:H81	2:C:3:FUC:H61	1.69	0.75
1:A:272:ASN:C	1:A:272:ASN:HD22	1.95	0.74
1:B:651:ILE:HG21	1:B:755:MET:HE3	1.69	0.74
2:D:1:NAG:O3	2:D:2:NAG:H83	1.88	0.74
1:A:492:ARG:HB3	1:A:492:ARG:NH1	2.03	0.73
1:A:177:GLU:HG3	1:A:180:LEU:HD22	1.71	0.73
1:A:581:ARG:CZ	8:A:782:NAG:H62	2.19	0.73
1:B:536:LYS:HD3	9:B:916:HOH:O	1.89	0.73
1:A:388:GLN:HB3	1:A:391:LYS:HD3	1.71	0.72
1:B:258:LYS:HZ1	1:B:712:HIS:HD2	1.35	0.72
1:A:69:LEU:HD13	1:A:107:ILE:HD12	1.72	0.72
1:B:450:ASN:HB2	9:B:907:HOH:O	1.89	0.72
1:B:96:ASP:HB2	9:B:947:HOH:O	1.89	0.72
1:A:736:THR:HG21	1:B:717:ALA:O	1.89	0.72
2:I:1:NAG:H61	2:I:3:FUC:O2	1.90	0.72
1:A:272:ASN:HD22	1:A:274:ASP:H	1.35	0.71
1:B:139:LYS:HG3	1:B:141:GLN:HB2	1.71	0.71
1:A:410:LEU:HD13	1:A:415:LEU:HD23	1.72	0.71
1:B:76:ILE:CG2	1:B:90:LEU:HB3	2.21	0.71
1:A:486:VAL:HG13	9:A:806:HOH:O	1.91	0.71
1:B:184:ARG:HD2	1:B:187:TRP:CE2	2.26	0.70
1:A:253:ARG:NH2	1:B:253:ARG:HH22	1.90	0.70
1:A:756:SER:O	1:A:760:LYS:HG2	1.92	0.70
4:F:2:NDG:H6C2	4:F:3:MAN:H2	1.73	0.70
8:A:782:NAG:O7	8:A:782:NAG:H3	1.91	0.69
1:B:91:GLU:HG2	1:B:94:THR:OG1	1.92	0.69
1:B:509:MET:HG3	1:B:510:PRO:HD2	1.73	0.69
1:A:91:GLU:O	1:A:93:SER:N	2.25	0.69
1:A:377:ASN:CG	1:A:378:GLU:H	2.00	0.69
1:B:308:GLN:HB3	3:J:1:NAG:H62	1.73	0.69
1:B:377:ASN:C	1:B:377:ASN:HD22	2.01	0.69
1:B:360:SER:HB3	1:B:373:LYS:HZ2	1.59	0.68
1:A:658:ARG:HG2	1:A:661:TYR:CE2	2.28	0.68
1:A:172:ILE:H	1:A:186:THR:CG2	2.07	0.68
1:A:502:LYS:O	1:A:505:GLN:HG2	1.93	0.67
1:B:596:ARG:O	1:B:597:ARG:HD2	1.93	0.67
1:A:177:GLU:CG	1:A:180:LEU:HD22	2.24	0.67
1:A:276:LEU:HD23	1:A:276:LEU:O	1.94	0.67
1:B:272:ASN:ND2	1:B:274:ASP:H	1.92	0.67
1:B:377:ASN:ND2	1:B:381:TYR:H	1.92	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:464:GLU:HG3	9:B:995:HOH:O	1.95	0.66
1:A:106:SER:HB3	1:A:115:LEU:HB3	1.77	0.66
1:A:153:GLN:HE22	1:A:170:ASN:ND2	1.92	0.66
1:B:293:MET:HE3	1:B:315:TRP:HB2	1.78	0.66
1:B:723:LEU:HB3	1:B:728:VAL:HG13	1.77	0.66
1:A:76:ILE:HD12	1:A:90:LEU:HD11	1.76	0.66
9:A:1149:HOH:O	3:G:2:NAG:H5	1.94	0.66
1:A:197:GLY:C	1:A:213:ALA:HB3	2.21	0.66
1:B:377:ASN:HD21	1:B:381:TYR:H	1.41	0.66
1:B:69:LEU:HB3	1:B:76:ILE:HD11	1.77	0.66
1:A:377:ASN:CG	1:A:378:GLU:N	2.55	0.65
1:B:74:ASN:HB3	1:B:92:ASN:CG	2.21	0.65
1:B:90:LEU:HD21	1:B:95:PHE:HE2	1.61	0.65
1:A:248:TYR:CZ	1:B:234:PRO:HB2	2.32	0.65
1:A:350:THR:HG22	3:H:1:NAG:H81	1.79	0.65
1:A:75:ASN:HB3	1:A:91:GLU:HA	1.79	0.64
1:B:75:ASN:ND2	1:B:92:ASN:HB2	2.12	0.64
1:B:528:MET:HE2	1:B:530:LEU:HD21	1.79	0.64
1:A:91:GLU:C	1:A:93:SER:H	2.04	0.64
1:B:535:ASP:OD1	1:B:537:SER:HB2	1.98	0.64
1:B:139:LYS:NZ	1:B:139:LYS:HB3	2.12	0.64
1:A:377:ASN:HB3	9:A:831:HOH:O	1.97	0.64
1:B:348:MET:HE2	7:M:1:NAG:H82	1.80	0.64
1:B:520:ASN:HD22	8:B:781:NAG:H4	1.62	0.63
1:B:272:ASN:C	1:B:272:ASN:HD22	2.06	0.63
1:A:291:ALA:O	1:A:295:ILE:HG23	1.99	0.63
1:B:56:LYS:NZ	1:B:56:LYS:HB3	2.14	0.63
1:B:399:LYS:HG3	9:B:825:HOH:O	1.99	0.63
1:B:581:ARG:CZ	8:B:781:NAG:H5	2.28	0.63
1:B:105:TYR:HB2	1:B:114:ILE:HD11	1.80	0.63
1:B:272:ASN:HD22	1:B:274:ASP:H	1.47	0.63
9:B:1194:HOH:O	2:I:2:NAG:H3	1.99	0.63
1:B:542:LEU:C	1:B:542:LEU:HD23	2.24	0.63
1:A:392:LYS:HG2	9:A:907:HOH:O	1.98	0.62
1:A:92:ASN:C	1:A:94:THR:H	2.05	0.62
1:A:438:ASP:OD2	1:A:440:THR:HG22	2.00	0.62
1:B:170:ASN:N	1:B:170:ASN:HD22	1.95	0.62
1:A:622:LYS:HB2	1:A:622:LYS:NZ	2.15	0.62
1:B:95:PHE:O	1:B:98:PHE:HB2	1.98	0.62
2:D:2:NAG:O7	2:D:2:NAG:H3	2.00	0.62
1:B:75:ASN:ND2	1:B:92:ASN:H	1.97	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:156:THR:HG21	9:B:1053:HOH:O	1.99	0.62
1:B:733:MET:HE2	1:B:734:TRP:O	2.00	0.62
1:B:658:ARG:HG2	1:B:661:TYR:CE2	2.34	0.62
1:A:751:ILE:HG12	1:A:755:MET:HE2	1.80	0.61
1:B:184:ARG:HD3	1:B:186:THR:O	1.99	0.61
1:B:399:LYS:HA	9:B:1126:HOH:O	1.98	0.61
1:A:253:ARG:HH22	1:B:253:ARG:NH2	1.95	0.61
1:B:528:MET:CE	1:B:530:LEU:HD21	2.30	0.61
1:A:277:SER:O	1:A:278:SER:HB2	2.00	0.61
1:B:520:ASN:HB2	8:B:781:NAG:H61	1.82	0.61
1:A:173:TYR:CE2	1:A:184:ARG:HG2	2.35	0.60
1:A:500:LEU:HA	1:A:503:MET:HE3	1.82	0.60
1:A:596:ARG:O	1:A:597:ARG:HD2	2.01	0.60
1:B:51:ASN:ND2	1:B:54:ARG:NE	2.48	0.60
1:B:504:LEU:HD22	1:B:509:MET:HE2	1.82	0.60
1:A:171:ASP:OD1	1:A:186:THR:HG23	2.02	0.60
1:A:393:ASP:HB2	9:A:1176:HOH:O	2.00	0.60
1:B:114:ILE:CG2	1:B:135:TYR:HB3	2.32	0.60
1:B:401:THR:HG22	1:B:401:THR:O	2.00	0.60
1:B:272:ASN:HD21	1:B:274:ASP:HB2	1.67	0.60
1:A:341:VAL:O	1:A:342:ALA:HB3	2.02	0.60
1:A:194:ILE:CD1	4:F:1:NAG:H82	2.29	0.60
1:A:40:ARG:HB3	1:A:506:ASN:O	2.02	0.59
1:A:435:GLN:NE2	1:A:441:LYS:HD2	2.17	0.59
1:B:620:ASP:OD2	1:B:623:ARG:HD3	2.01	0.59
1:B:286:GLN:NE2	1:B:288:THR:HG22	2.18	0.59
1:A:341:VAL:C	1:A:343:ARG:H	2.09	0.59
1:A:314:GLN:HE22	1:A:373:LYS:NZ	2.01	0.58
1:A:341:VAL:HG12	9:A:1148:HOH:O	2.03	0.58
1:B:66:HIS:HB2	9:B:930:HOH:O	2.03	0.58
1:B:651:ILE:HG21	1:B:755:MET:CE	2.32	0.58
1:B:40:ARG:NH1	1:B:508:GLN:HG2	2.19	0.58
1:B:153:GLN:HE22	1:B:170:ASN:ND2	2.02	0.58
1:B:599:GLY:H	1:B:602:GLU:CD	2.11	0.57
1:A:119:ASN:HD22	1:A:131:SER:CB	2.16	0.57
1:B:676:PRO:HG2	1:B:677:GLU:OE2	2.04	0.57
1:B:51:ASN:HD21	1:B:54:ARG:NE	2.01	0.57
1:A:273:THR:O	1:A:276:LEU:HD22	2.05	0.57
1:B:433:LYS:HB3	1:B:445:LEU:HD21	1.85	0.57
1:B:546:VAL:HG22	1:B:547:PHE:N	2.19	0.57
1:A:481:THR:OG1	1:A:483:HIS:HE1	1.87	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:THR:HG21	1:A:214:LEU:HD11	1.87	0.57
1:A:388:GLN:CB	1:A:391:LYS:HD3	2.33	0.57
1:A:312:SER:CB	1:A:325:MET:HE3	2.34	0.57
1:B:461:PHE:CD2	1:B:468:TYR:HB3	2.40	0.57
1:A:581:ARG:NH1	8:A:782:NAG:H62	2.20	0.57
1:A:535:ASP:OD2	1:A:537:SER:HB3	2.05	0.56
1:A:696:LYS:HE3	9:A:924:HOH:O	2.04	0.56
1:B:65:ASP:OD1	1:B:464:GLU:HB2	2.05	0.56
1:B:232:GLU:HB3	1:B:262:VAL:HG11	1.87	0.56
1:B:414:TYR:HA	1:B:436:LEU:HD13	1.88	0.56
1:A:741:GLY:O	1:A:742:ILE:C	2.47	0.56
1:B:208:PHE:O	1:B:209:SER:C	2.47	0.56
1:B:289:ALA:HA	1:B:294:LEU:HG	1.87	0.56
1:B:500:LEU:HA	1:B:503:MET:HE3	1.87	0.56
1:B:90:LEU:HD21	1:B:95:PHE:CE2	2.40	0.56
1:A:140:ARG:HG2	1:A:140:ARG:HH11	1.71	0.56
1:A:214:LEU:O	1:A:214:LEU:HD12	2.06	0.56
1:A:293:MET:HE3	1:A:315:TRP:HB2	1.87	0.56
1:B:72:GLN:C	1:B:74:ASN:H	2.13	0.56
1:B:733:MET:HE2	9:B:820:HOH:O	2.05	0.56
1:A:492:ARG:HH11	1:A:492:ARG:CB	2.10	0.56
1:A:92:ASN:C	1:A:94:THR:N	2.64	0.56
1:A:242:SER:HB3	1:A:246:LEU:HD12	1.88	0.56
1:B:504:LEU:HD22	1:B:509:MET:CE	2.36	0.55
1:A:660:GLU:HG3	9:A:911:HOH:O	2.06	0.55
1:B:90:LEU:CD1	1:B:94:THR:HG21	2.34	0.55
2:C:2:NAG:H81	2:C:3:FUC:C6	2.35	0.55
1:A:136:ASP:CG	1:A:139:LYS:HG2	2.32	0.55
1:A:710:ASN:HD22	1:A:710:ASN:C	2.14	0.55
1:B:147:ARG:HH21	8:B:770:NAG:HN2	1.55	0.55
1:B:598:LEU:HA	1:B:602:GLU:OE2	2.06	0.55
1:A:697:GLN:HB3	9:A:1030:HOH:O	2.07	0.55
1:B:268:PHE:CD2	1:B:313:LEU:HD21	2.41	0.55
1:B:384:ILE:HG13	1:B:404:VAL:HG21	1.89	0.55
1:A:310:ARG:HG3	1:A:329:ASP:OD1	2.06	0.55
1:A:196:ASN:OD1	1:A:227:GLN:HG3	2.06	0.55
1:A:546:VAL:HG22	1:A:547:PHE:N	2.21	0.55
1:B:399:LYS:HB2	1:B:402:TRP:HZ2	1.70	0.55
1:B:529:ILE:HD13	9:B:983:HOH:O	2.06	0.55
1:A:658:ARG:HD3	1:A:660:GLU:HB2	1.89	0.55
1:B:56:LYS:HB3	1:B:56:LYS:HZ3	1.72	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:718:GLN:HE21	1:A:718:GLN:HA	1.70	0.54
1:A:89:PHE:HD1	1:A:90:LEU:HD12	1.72	0.54
1:A:234:PRO:HB2	1:B:248:TYR:CZ	2.42	0.54
1:A:535:ASP:C	1:A:537:SER:H	2.15	0.54
1:B:65:ASP:CG	1:B:464:GLU:HB2	2.33	0.54
1:B:289:ALA:CB	1:B:290:PRO:CA	2.82	0.54
1:B:289:ALA:CB	1:B:290:PRO:HA	2.28	0.54
1:A:74:ASN:O	1:A:92:ASN:HB2	2.08	0.54
1:A:75:ASN:CB	1:A:91:GLU:HA	2.37	0.54
1:A:600:THR:O	1:A:603:VAL:HG13	2.07	0.54
1:B:112:GLN:HB3	1:B:138:ASN:ND2	2.21	0.54
1:B:657:SER:OG	1:B:689:MET:HE1	2.07	0.54
1:A:39:SER:O	1:A:40:ARG:HB2	2.07	0.54
1:B:521:GLU:HG3	1:B:521:GLU:O	2.08	0.54
1:B:502:LYS:HD3	1:B:502:LYS:C	2.32	0.54
1:A:41:LYS:HD3	1:A:41:LYS:H	1.72	0.53
1:B:538:LYS:HD3	1:B:539:LYS:N	2.23	0.53
1:B:741:GLY:O	1:B:742:ILE:C	2.51	0.53
1:A:507:VAL:HB	9:A:1113:HOH:O	2.08	0.53
1:A:314:GLN:HE22	1:A:373:LYS:HZ1	1.56	0.53
1:B:75:ASN:CG	1:B:92:ASN:HB2	2.34	0.53
1:B:139:LYS:O	1:B:141:GLN:HG2	2.07	0.53
1:B:399:LYS:HB2	1:B:402:TRP:CE2	2.43	0.53
1:B:510:PRO:HD3	1:B:569:SER:HB2	1.89	0.53
9:B:862:HOH:O	6:L:2:NDG:H6C2	2.08	0.53
1:A:95:PHE:HB3	1:A:98:PHE:HB2	1.88	0.53
1:B:385:CYS:HB3	1:B:396:PHE:CD1	2.44	0.53
1:B:193:ILE:HG22	1:B:194:ILE:HG12	1.89	0.53
1:A:91:GLU:C	1:A:93:SER:N	2.67	0.53
1:A:546:VAL:CG2	1:A:547:PHE:N	2.71	0.53
1:B:627:TRP:CE3	1:B:755:MET:HE1	2.43	0.53
2:D:1:NAG:H4	2:D:2:NAG:N2	2.23	0.53
1:A:402:TRP:CD2	1:A:421:GLU:HB2	2.44	0.53
1:B:57:LEU:N	1:B:57:LEU:HD12	2.23	0.53
1:B:81:ALA:O	1:B:492:ARG:NH2	2.40	0.52
1:B:139:LYS:HB3	1:B:139:LYS:HZ3	1.75	0.52
1:B:74:ASN:O	1:B:92:ASN:HA	2.09	0.52
1:B:147:ARG:NH2	8:B:770:NAG:HN2	2.06	0.52
1:B:644:SER:C	1:B:646:VAL:H	2.14	0.52
1:B:765:LEU:HB3	1:B:766:PRO:HA	1.91	0.52
1:A:542:LEU:HD23	1:A:542:LEU:C	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:538:LYS:HD3	1:B:539:LYS:O	2.10	0.52
1:A:75:ASN:HB2	1:A:90:LEU:O	2.10	0.52
1:B:77:LEU:HD23	1:B:77:LEU:N	2.19	0.52
1:B:290:PRO:HG3	1:B:326:ASP:OD2	2.09	0.52
1:A:186:THR:HG21	1:A:196:ASN:CB	2.39	0.52
1:A:76:ILE:HB	1:A:90:LEU:CD1	2.40	0.52
1:A:435:GLN:HE22	1:A:441:LYS:HD2	1.75	0.52
1:B:320:GLN:OE1	1:B:669:ARG:HD3	2.10	0.52
1:A:662:TYR:OH	1:A:710:ASN:ND2	2.31	0.52
1:B:64:SER:O	1:B:463:LYS:HG2	2.09	0.52
1:B:538:LYS:HD3	1:B:539:LYS:C	2.34	0.52
1:A:150:ASN:O	1:A:151:ASN:HB2	2.09	0.52
1:A:751:ILE:O	1:A:755:MET:HG3	2.10	0.52
1:B:696:LYS:HB2	9:B:1188:HOH:O	2.10	0.52
1:A:289:ALA:CB	1:A:290:PRO:HA	2.32	0.51
1:A:517:ILE:HD11	1:A:578:PHE:CE1	2.45	0.51
1:A:581:ARG:CZ	8:A:782:NAG:C6	2.87	0.51
1:B:348:MET:HE2	7:M:1:NAG:C8	2.39	0.51
1:B:139:LYS:HG3	1:B:141:GLN:CB	2.40	0.51
1:A:76:ILE:HB	1:A:90:LEU:HD13	1.91	0.51
1:A:289:ALA:CB	1:A:290:PRO:CA	2.84	0.51
1:B:316:LEU:HD12	1:B:322:TYR:O	2.11	0.51
1:A:726:VAL:HG12	1:A:726:VAL:O	2.10	0.51
1:B:302:ASP:OD1	1:B:304:THR:HG23	2.11	0.51
1:B:710:ASN:C	1:B:710:ASN:HD22	2.18	0.51
1:B:733:MET:HE3	1:B:734:TRP:H	1.75	0.51
1:A:60:LEU:HD12	1:A:60:LEU:C	2.36	0.51
1:B:431:LEU:HG	1:B:445:LEU:HD23	1.93	0.51
1:A:170:ASN:N	1:A:170:ASN:HD22	2.09	0.51
1:A:272:ASN:HD21	1:A:274:ASP:HB2	1.75	0.51
1:A:325:MET:HE2	1:A:327:ILE:HG12	1.92	0.51
1:A:512:LYS:HD3	9:A:995:HOH:O	2.11	0.51
1:A:718:GLN:HA	1:A:718:GLN:NE2	2.25	0.51
1:A:325:MET:HE2	1:A:327:ILE:CG1	2.42	0.50
1:A:662:TYR:HB3	1:A:667:THR:OG1	2.10	0.50
1:B:76:ILE:HG23	1:B:90:LEU:HB3	1.92	0.50
1:B:651:ILE:CG2	1:B:755:MET:HE3	2.40	0.50
1:B:74:ASN:HB3	1:B:92:ASN:OD1	2.10	0.50
1:B:502:LYS:O	1:B:502:LYS:HD3	2.11	0.50
1:B:546:VAL:CG2	1:B:547:PHE:N	2.74	0.50
1:A:378:GLU:HG3	9:A:1211:HOH:O	2.10	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:704:HIS:HE1	1:A:711:VAL:O	1.95	0.50
1:A:640:LEU:HD11	1:A:650:GLY:HA3	1.93	0.50
1:A:184:ARG:HH12	1:A:187:TRP:HA	1.76	0.50
1:A:302:ASP:HB3	1:A:314:GLN:HB2	1.92	0.50
1:B:55:LEU:HD23	1:B:478:PRO:HG2	1.93	0.50
1:B:60:LEU:HD23	1:B:60:LEU:C	2.36	0.50
1:B:232:GLU:HG2	9:B:1134:HOH:O	2.10	0.50
1:A:69:LEU:CD1	1:A:107:ILE:HD12	2.40	0.50
1:A:140:ARG:HG2	1:A:140:ARG:NH1	2.27	0.50
1:A:635:VAL:O	1:A:639:VAL:HG23	2.11	0.50
1:A:704:HIS:HD2	1:A:716:SER:OG	1.95	0.50
1:B:184:ARG:HH11	1:B:187:TRP:HA	1.77	0.50
1:B:302:ASP:HB3	1:B:314:GLN:HB2	1.94	0.50
1:B:400:GLY:HA2	9:B:1048:HOH:O	2.12	0.50
1:B:72:GLN:O	1:B:74:ASN:N	2.44	0.49
1:A:622:LYS:HB2	1:A:622:LYS:HZ2	1.77	0.49
1:B:72:GLN:HE21	1:B:77:LEU:CD2	2.19	0.49
1:A:658:ARG:HG2	1:A:661:TYR:CD2	2.48	0.49
1:B:293:MET:HE3	1:B:315:TRP:CB	2.42	0.49
1:A:72:GLN:HB3	9:A:1061:HOH:O	2.12	0.49
1:A:156:THR:HG21	1:A:214:LEU:CD1	2.41	0.49
1:B:658:ARG:HD3	1:B:660:GLU:HB2	1.93	0.49
1:B:664:SER:HB2	1:B:668:GLU:OE2	2.12	0.49
1:B:106:SER:HB3	1:B:115:LEU:HB3	1.94	0.49
8:B:781:NAG:O7	8:B:781:NAG:C3	2.53	0.49
1:A:633:GLY:HA3	1:A:655:PRO:HB3	1.95	0.49
1:A:244:GLU:CD	1:B:689:MET:HG3	2.38	0.49
1:B:708:ASP:OD2	1:B:740:HIS:HA	2.13	0.49
2:C:1:NAG:H62	2:C:2:NAG:C8	2.39	0.49
1:A:39:SER:O	1:A:40:ARG:CB	2.60	0.49
1:A:118:TYR:CE2	1:A:119:ASN:ND2	2.80	0.49
1:A:159:PRO:HD3	1:A:216:TRP:CB	2.42	0.49
1:B:403:GLU:OE1	1:B:585:TYR:HA	2.13	0.49
1:B:446:SER:HB2	1:B:457:TYR:CE2	2.48	0.48
1:B:400:GLY:O	1:B:402:TRP:N	2.44	0.48
1:B:96:ASP:O	1:B:97:GLU:HB2	2.13	0.48
1:A:520:ASN:OD1	8:A:782:NAG:O5	2.29	0.48
1:B:114:ILE:HG22	1:B:135:TYR:HB3	1.95	0.48
1:A:277:SER:OG	1:A:280:THR:HB	2.14	0.48
1:B:110:ASP:OD1	1:B:112:GLN:HG3	2.14	0.48
1:B:250:LYS:HB2	1:B:250:LYS:NZ	2.28	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:LYS:HG3	9:A:849:HOH:O	2.13	0.48
1:A:325:MET:HE2	1:A:327:ILE:HD11	1.95	0.48
1:A:410:LEU:HD13	1:A:415:LEU:CD2	2.42	0.48
1:B:517:ILE:CD1	1:B:612:GLN:HG3	2.43	0.48
1:A:108:SER:C	1:A:110:ASP:N	2.70	0.48
1:B:648:LYS:HE2	1:B:762:CYS:O	2.14	0.48
1:A:669:ARG:HD2	1:A:670:TYR:CZ	2.49	0.47
1:B:165:ALA:HB2	1:B:216:TRP:CZ2	2.49	0.47
1:B:630:SER:HB3	9:B:1021:HOH:O	2.12	0.47
1:B:654:ALA:N	1:B:655:PRO:CD	2.77	0.47
1:A:658:ARG:HG3	1:A:658:ARG:O	2.14	0.47
1:B:289:ALA:HB1	1:B:290:PRO:C	2.39	0.47
1:A:77:LEU:CD2	1:A:88:VAL:HA	2.44	0.47
1:A:136:ASP:OD1	1:A:139:LYS:HG2	2.14	0.47
1:B:98:PHE:CD1	1:B:100:HIS:HB2	2.49	0.47
1:A:664:SER:O	1:A:668:GLU:HB2	2.15	0.47
1:A:689:MET:HG3	1:B:244:GLU:CD	2.39	0.47
1:B:402:TRP:CD2	1:B:421:GLU:HB2	2.49	0.47
1:A:504:LEU:HA	1:A:507:VAL:HG12	1.96	0.47
1:B:214:LEU:HD12	1:B:214:LEU:O	2.14	0.47
1:B:445:LEU:H	1:B:445:LEU:HD22	1.80	0.47
8:B:781:NAG:H62	9:B:1099:HOH:O	2.13	0.47
1:A:544:LEU:HD21	1:A:606:GLN:HG3	1.97	0.47
1:A:720:SER:O	1:A:724:VAL:HG23	2.14	0.47
1:B:125:ARG:NH2	1:B:205:GLU:OE2	2.48	0.47
1:B:159:PRO:HD3	1:B:216:TRP:CB	2.45	0.47
1:B:293:MET:HE3	1:B:315:TRP:C	2.39	0.47
1:B:317:ARG:HD2	1:B:322:TYR:HB3	1.97	0.47
1:B:438:ASP:OD2	1:B:441:LYS:HG2	2.15	0.47
1:A:331:ASP:O	1:A:335:GLY:N	2.46	0.47
1:A:438:ASP:CG	1:A:440:THR:HG22	2.39	0.47
1:A:606:GLN:NE2	9:A:799:HOH:O	2.47	0.47
1:A:671:MET:HE1	1:A:682:HIS:CD2	2.48	0.47
1:B:197:GLY:C	1:B:213:ALA:HB3	2.40	0.47
1:B:334:SER:C	1:B:336:ARG:H	2.22	0.47
1:B:293:MET:HG2	1:B:315:TRP:HB3	1.97	0.47
1:B:415:LEU:C	1:B:415:LEU:HD23	2.39	0.47
1:A:446:SER:HB2	1:A:457:TYR:CE1	2.50	0.47
1:B:69:LEU:HD11	1:B:107:ILE:HD12	1.96	0.47
1:B:80:ASN:OD1	1:B:82:GLU:HB3	2.15	0.47
4:F:1:NAG:H62	4:F:2:NDG:O5	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:325:MET:HE1	1:A:364:PHE:HZ	1.79	0.46
1:A:580:GLY:O	1:A:581:ARG:C	2.55	0.46
1:B:279:VAL:O	1:B:280:THR:HG22	2.16	0.46
1:B:77:LEU:H	1:B:77:LEU:CD2	2.20	0.46
1:B:459:VAL:HG22	1:B:460:SER:N	2.31	0.46
1:B:121:VAL:HB	1:B:129:THR:CG2	2.46	0.46
1:A:123:GLN:HG2	1:A:124:TRP:CD2	2.51	0.46
1:B:677:GLU:CD	1:B:677:GLU:H	2.23	0.46
1:A:704:HIS:CE1	1:A:711:VAL:O	2.69	0.46
1:A:177:GLU:HB2	1:A:180:LEU:HD22	1.98	0.46
1:A:340:LEU:O	1:A:343:ARG:HB3	2.16	0.46
1:A:522:THR:HG22	1:A:523:LYS:N	2.31	0.46
9:A:992:HOH:O	4:F:2:NDG:C5	2.64	0.46
1:B:44:THR:O	1:B:47:ASP:HB2	2.16	0.46
1:B:109:PRO:HG2	1:B:158:SER:O	2.16	0.46
1:B:279:VAL:O	1:B:280:THR:CB	2.63	0.46
1:B:596:ARG:C	1:B:597:ARG:HD2	2.41	0.45
1:A:285:ILE:N	1:A:285:ILE:HD12	2.32	0.45
1:B:532:PRO:HD3	1:B:569:SER:HA	1.98	0.45
1:A:677:GLU:H	1:A:677:GLU:CD	2.25	0.45
1:B:634:TYR:O	1:B:638:MET:HG2	2.17	0.45
1:A:392:LYS:HG3	1:A:393:ASP:N	2.31	0.45
1:A:341:VAL:HG22	1:A:342:ALA:H	1.82	0.45
1:A:242:SER:N	1:A:737:ASP:OD2	2.49	0.45
1:A:415:LEU:C	1:A:415:LEU:HD13	2.42	0.45
1:A:128:TYR:CD1	1:A:128:TYR:C	2.95	0.45
1:A:507:VAL:HG13	1:A:509:MET:HG2	1.99	0.45
1:A:534:PHE:HZ	1:A:618:PHE:CD1	2.35	0.45
1:B:343:ARG:HA	1:B:389:ILE:O	2.16	0.45
1:B:399:LYS:HB2	1:B:402:TRP:NE1	2.32	0.45
1:B:751:ILE:O	1:B:755:MET:HG3	2.16	0.45
1:A:77:LEU:HD23	1:A:88:VAL:HA	1.98	0.44
1:A:340:LEU:C	1:A:341:VAL:O	2.57	0.44
1:B:237:GLU:HA	1:B:252:VAL:O	2.17	0.44
1:B:517:ILE:HG12	1:B:518:ILE:N	2.30	0.44
1:B:745:SER:O	1:B:749:GLN:HG3	2.17	0.44
1:A:654:ALA:HA	1:A:704:HIS:CD2	2.51	0.44
1:B:305:TRP:CE2	1:B:311:ILE:HD12	2.53	0.44
1:B:422:TYR:CE2	1:B:423:LYS:HD3	2.52	0.44
1:B:57:LEU:HB3	9:B:1008:HOH:O	2.16	0.44
1:B:366:LEU:HD23	9:B:1122:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:247:GLN:HG2	1:B:258:LYS:HD2	1.99	0.44
1:A:253:ARG:NH2	1:B:253:ARG:HH12	2.16	0.44
1:A:314:GLN:NE2	1:A:373:LYS:NZ	2.64	0.44
1:B:60:LEU:HD21	1:B:62:TRP:NE1	2.32	0.44
1:B:519:LEU:O	1:B:520:ASN:C	2.60	0.44
1:A:119:ASN:HD22	1:A:131:SER:HB2	1.80	0.44
1:A:405:ILE:HG13	1:A:429:ARG:CD	2.46	0.44
1:A:535:ASP:CG	1:A:537:SER:HB3	2.42	0.44
1:B:70:TYR:O	1:B:77:LEU:HD23	2.18	0.44
1:A:55:LEU:HD23	1:A:500:LEU:CD2	2.47	0.44
1:B:223:LEU:HD23	1:B:223:LEU:HA	1.84	0.44
1:B:382:ARG:NH2	9:B:782:HOH:O	2.51	0.44
1:B:413:ASP:O	1:B:436:LEU:HB2	2.17	0.44
1:A:289:ALA:HB1	1:A:290:PRO:C	2.42	0.44
1:A:627:TRP:HB2	1:A:651:ILE:HB	2.00	0.44
1:B:377:ASN:ND2	1:B:379:GLU:H	2.16	0.44
1:B:387:PHE:CE1	1:B:394:CYS:HB3	2.53	0.44
1:A:535:ASP:C	1:A:537:SER:N	2.74	0.43
1:B:258:LYS:NZ	1:B:712:HIS:CD2	2.73	0.43
1:A:56:LYS:HB2	1:A:497:ASN:OD1	2.18	0.43
1:A:114:ILE:HG13	1:A:137:LEU:HD11	2.00	0.43
1:A:718:GLN:HE21	1:A:718:GLN:CA	2.29	0.43
1:B:733:MET:HE3	1:B:734:TRP:N	2.33	0.43
1:A:161:GLY:HA3	9:A:934:HOH:O	2.18	0.43
1:A:377:ASN:CB	9:A:783:HOH:O	2.56	0.43
1:A:459:VAL:HG22	1:A:460:SER:N	2.33	0.43
1:A:693:GLU:HA	1:A:726:VAL:HG11	2.00	0.43
1:A:341:VAL:O	1:A:342:ALA:CB	2.67	0.43
1:A:272:ASN:ND2	1:A:272:ASN:C	2.67	0.43
1:A:314:GLN:NE2	1:A:359:PRO:HB2	2.34	0.43
1:A:370:SER:HB2	1:A:387:PHE:O	2.19	0.43
1:B:648:LYS:HD3	1:B:762:CYS:SG	2.59	0.43
1:A:154:TRP:CE2	1:A:212:SER:HB2	2.54	0.43
1:A:607:ILE:HG12	1:A:639:VAL:HG13	2.01	0.43
1:B:64:SER:C	1:B:463:LYS:HG2	2.44	0.43
1:B:651:ILE:HG23	1:B:701:LEU:HB3	2.00	0.43
1:A:208:PHE:O	1:A:209:SER:C	2.60	0.43
1:A:512:LYS:HE3	1:A:527:GLN:OE1	2.19	0.43
1:B:246:LEU:HD23	1:B:246:LEU:HA	1.84	0.42
1:B:418:ILE:HA	1:B:430:ASN:O	2.19	0.42
1:A:91:GLU:OE1	1:A:91:GLU:N	2.52	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:TYR:CE2	1:A:137:LEU:HD23	2.54	0.42
1:B:69:LEU:CD1	1:B:107:ILE:HD12	2.48	0.42
1:A:392:LYS:HG3	1:A:393:ASP:H	1.85	0.42
1:B:51:ASN:HA	9:B:1149:HOH:O	2.18	0.42
1:B:693:GLU:O	1:B:696:LYS:HG3	2.18	0.42
1:A:158:SER:OG	1:A:163:LYS:HB2	2.19	0.42
1:A:248:TYR:HA	1:A:249:PRO:HD3	1.93	0.42
1:A:82:GLU:HG2	9:A:981:HOH:O	2.18	0.42
1:A:279:VAL:O	1:A:279:VAL:HG12	2.19	0.42
1:B:54:ARG:HE	1:B:54:ARG:HB2	1.58	0.42
1:B:280:THR:HA	9:B:1146:HOH:O	2.18	0.42
1:A:519:LEU:O	1:A:520:ASN:C	2.62	0.42
1:A:658:ARG:HD2	1:A:661:TYR:CE1	2.54	0.42
1:B:218:PRO:HG2	1:B:308:GLN:OE1	2.20	0.42
1:B:289:ALA:HA	1:B:294:LEU:CG	2.49	0.42
1:B:41:LYS:H	1:B:41:LYS:HD3	1.84	0.42
1:B:680:LEU:O	1:B:683:TYR:HB2	2.20	0.42
8:B:781:NAG:C6	9:B:1099:HOH:O	2.67	0.42
1:B:40:ARG:HG2	1:B:508:GLN:HG3	2.02	0.42
1:B:48:TYR:CE1	1:B:562:ASN:HA	2.55	0.42
1:B:128:TYR:CD1	1:B:128:TYR:C	2.98	0.42
1:A:63:ILE:HG21	1:A:69:LEU:HG	2.02	0.42
1:A:187:TRP:HE3	3:G:2:NAG:H83	1.85	0.42
1:A:325:MET:HE2	1:A:327:ILE:CD1	2.50	0.42
1:A:361:GLU:HB3	9:A:1078:HOH:O	2.20	0.42
1:B:125:ARG:HH21	1:B:205:GLU:CD	2.28	0.42
1:B:154:TRP:NE1	1:B:156:THR:HG23	2.35	0.42
1:B:492:ARG:HB3	9:B:960:HOH:O	2.19	0.42
1:A:60:LEU:HD12	1:A:60:LEU:O	2.19	0.42
1:B:442:VAL:HG11	9:B:1077:HOH:O	2.19	0.42
1:B:487:ASN:O	1:B:488:ASP:HB2	2.19	0.42
1:B:602:GLU:HG3	1:B:603:VAL:N	2.35	0.42
4:F:2:NDG:C6	4:F:3:MAN:H2	2.46	0.42
1:A:341:VAL:C	1:A:343:ARG:N	2.74	0.41
1:B:122:LYS:HG2	1:B:123:GLN:N	2.35	0.41
1:B:177:GLU:HB2	1:B:180:LEU:HG	2.02	0.41
1:A:108:SER:C	1:A:110:ASP:H	2.27	0.41
1:A:272:ASN:ND2	1:A:274:ASP:N	2.59	0.41
1:A:519:LEU:HB2	1:A:524:PHE:CE1	2.55	0.41
1:A:760:LYS:HB3	1:A:760:LYS:HE2	1.77	0.41
1:B:76:ILE:HG23	1:B:76:ILE:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:ALA:HB1	1:B:268:PHE:CZ	2.56	0.41
1:B:514:LEU:CD2	1:B:525:TRP:HE3	2.33	0.41
1:A:57:LEU:N	1:A:57:LEU:HD12	2.34	0.41
1:B:718:GLN:HE22	1:B:721:LYS:NZ	2.18	0.41
2:I:1:NAG:C6	2:I:3:FUC:O2	2.66	0.41
1:A:74:ASN:O	1:A:91:GLU:O	2.38	0.41
1:A:266:VAL:HG22	1:A:267:LYS:N	2.35	0.41
1:A:510:PRO:HD3	1:A:569:SER:HB2	2.02	0.41
1:A:622:LYS:NZ	9:A:960:HOH:O	2.54	0.41
1:B:77:LEU:N	1:B:77:LEU:CD2	2.82	0.41
1:B:530:LEU:HA	1:B:531:PRO:HD3	1.83	0.41
1:A:293:MET:HG2	1:A:315:TRP:HB3	2.01	0.41
1:B:154:TRP:CE2	1:B:212:SER:HB2	2.55	0.41
1:A:190:LYS:HE3	1:A:193:ILE:HG13	2.02	0.41
1:A:325:MET:HE1	1:A:364:PHE:CZ	2.55	0.41
1:A:510:PRO:HB2	1:A:530:LEU:O	2.21	0.41
1:A:186:THR:HG21	1:A:196:ASN:HB3	2.02	0.41
1:A:657:SER:HB2	1:A:689:MET:SD	2.61	0.41
1:B:258:LYS:HZ3	1:B:712:HIS:HD2	1.65	0.41
1:B:520:ASN:HD22	8:B:781:NAG:C4	2.26	0.41
1:B:742:ILE:O	1:B:742:ILE:HG22	2.21	0.41
1:A:118:TYR:CD2	1:A:119:ASN:ND2	2.89	0.41
1:A:520:ASN:C	1:A:521:GLU:HG2	2.46	0.41
1:A:608:GLU:OE1	1:A:608:GLU:HA	2.19	0.41
1:B:72:GLN:C	1:B:74:ASN:N	2.79	0.41
1:B:267:LYS:HE2	9:B:1138:HOH:O	2.20	0.41
1:B:289:ALA:HA	1:B:294:LEU:CD1	2.51	0.41
1:B:322:TYR:HD1	1:B:348:MET:HE3	1.85	0.41
1:B:504:LEU:HA	1:B:507:VAL:CG1	2.50	0.41
1:B:62:TRP:CD2	1:B:462:SER:HA	2.56	0.41
1:A:377:ASN:ND2	9:A:783:HOH:O	2.52	0.40
1:A:526:TYR:CD1	1:A:526:TYR:C	2.99	0.40
1:B:121:VAL:HB	1:B:129:THR:HG23	2.03	0.40
1:A:272:ASN:HD22	1:A:273:THR:N	2.17	0.40
1:A:679:ASN:O	1:A:683:TYR:HD2	2.04	0.40
1:B:415:LEU:HB2	1:B:436:LEU:HD11	2.02	0.40
1:B:598:LEU:HG	1:B:631:TYR:OH	2.21	0.40
1:B:658:ARG:HG2	1:B:661:TYR:CD2	2.55	0.40
1:A:579:ASP:HB3	1:A:583:SER:OG	2.21	0.40
1:B:57:LEU:HD11	9:B:1180:HOH:O	2.21	0.40
1:A:217:SER:HB3	1:A:222:PHE:HB2	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:150:ASN:O	1:A:151:ASN:CB	2.70	0.40
1:A:600:THR:O	1:A:603:VAL:CG1	2.70	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	724/728 (100%)	664 (92%)	53 (7%)	7 (1%)	12	11
1	B	726/728 (100%)	669 (92%)	49 (7%)	8 (1%)	11	10
All	All	1450/1456 (100%)	1333 (92%)	102 (7%)	15 (1%)	12	11

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	278	SER
1	A	320	GLN
1	A	520	ASN
1	B	399	LYS
1	B	401	THR
1	B	645	GLY
1	A	289	ALA
1	B	73	GLU
1	B	289	ALA
1	B	320	GLN
1	B	40	ARG
1	A	277	SER
1	A	389	ILE
1	A	536	LYS
1	B	332	GLU

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	650/653 (100%)	619 (95%)	31 (5%)	23	30
1	B	652/653 (100%)	615 (94%)	37 (6%)	18	23
All	All	1302/1306 (100%)	1234 (95%)	68 (5%)	21	26

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	LEU
1	A	145	GLU
1	A	170	ASN
1	A	184	ARG
1	A	246	LEU
1	A	254	VAL
1	A	272	ASN
1	A	276	LEU
1	A	373	LYS
1	A	382	ARG
1	A	390	ASP
1	A	415	LEU
1	A	436	LEU
1	A	440	THR
1	A	442	VAL
1	A	448	GLU
1	A	482	LEU
1	A	492	ARG
1	A	536	LYS
1	A	543	LEU
1	A	561	LEU
1	A	603	VAL
1	A	608	GLU
1	A	622	LYS
1	A	679	ASN
1	A	689	MET
1	A	701	LEU

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Mol	Chain	Res	Type
1	A	702	LEU
1	A	736	THR
1	A	760	LYS
1	A	761	GLN
1	B	57	LEU
1	B	60	LEU
1	B	61	ARG
1	B	66	HIS
1	B	73	GLU
1	B	77	LEU
1	B	139	LYS
1	B	156	THR
1	B	170	ASN
1	B	223	LEU
1	B	232	GLU
1	B	246	LEU
1	B	250	LYS
1	B	253	ARG
1	B	280	THR
1	B	283	THR
1	B	303	VAL
1	B	304	THR
1	B	366	LEU
1	B	377	ASN
1	B	379	GLU
1	B	410	LEU
1	B	429	ARG
1	B	443	THR
1	B	445	LEU
1	B	448	GLU
1	B	502	LYS
1	B	507	VAL
1	B	514	LEU
1	B	517	ILE
1	B	543	LEU
1	B	677	GLU
1	B	689	MET
1	B	701	LEU
1	B	702	LEU
1	B	728	VAL
1	B	761	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (49)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	72	GLN
1	A	92	ASN
1	A	103	ASN
1	A	119	ASN
1	A	169	ASN
1	A	170	ASN
1	A	227	GLN
1	A	247	GLN
1	A	272	ASN
1	A	314	GLN
1	A	338	ASN
1	A	344	GLN
1	A	369	ASN
1	A	430	ASN
1	A	435	GLN
1	A	483	HIS
1	A	612	GLN
1	A	679	ASN
1	A	685	ASN
1	A	694	ASN
1	A	704	HIS
1	A	710	ASN
1	A	718	GLN
1	A	731	GLN
1	A	761	GLN
1	B	51	ASN
1	B	72	GLN
1	B	75	ASN
1	B	112	GLN
1	B	123	GLN
1	B	138	ASN
1	B	169	ASN
1	B	170	ASN
1	B	227	GLN
1	B	272	ASN
1	B	286	GLN
1	B	314	GLN
1	B	344	GLN
1	B	377	ASN
1	B	430	ASN
1	B	572	ASN
1	B	621	ASN

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Mol	Chain	Res	Type
1	B	679	ASN
1	B	685	ASN
1	B	694	ASN
1	B	710	ASN
1	B	712	HIS
1	B	718	GLN
1	B	761	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 ⓘ

28 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	C	1	2,1	14,14,15	0.68	0	17,19,21	1.19	1 (5%)
2	NAG	C	2	2	14,14,15	0.45	0	17,19,21	0.79	1 (5%)
2	FUC	C	3	2	10,10,11	0.58	0	14,14,16	0.61	0
2	NAG	D	1	2,1	14,14,15	1.06	1 (7%)	17,19,21	0.86	0
2	NAG	D	2	2	14,14,15	0.84	1 (7%)	17,19,21	0.65	0
2	FUC	D	3	2	10,10,11	0.61	0	14,14,16	0.40	0
3	NAG	E	1	3,1	14,14,15	0.96	1 (7%)	17,19,21	1.30	2 (11%)
3	NAG	E	2	3	14,14,15	0.61	0	17,19,21	0.79	1 (5%)
4	NAG	F	1	4,1	14,14,15	0.75	0	17,19,21	0.85	1 (5%)
4	NDG	F	2	4	14,14,15	0.89	0	17,19,21	1.47	3 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	MAN	F	3	4	11,11,12	0.82	0	15,15,17	0.53	0
3	NAG	G	1	3,1	14,14,15	0.64	0	17,19,21	0.80	1 (5%)
3	NAG	G	2	3	14,14,15	0.55	0	17,19,21	0.78	0
3	NAG	H	1	3,1	14,14,15	0.63	0	17,19,21	0.80	0
3	NAG	H	2	3	14,14,15	0.65	0	17,19,21	0.93	0
2	NAG	I	1	2,1	14,14,15	0.65	0	17,19,21	1.39	2 (11%)
2	NAG	I	2	2	14,14,15	0.61	0	17,19,21	0.98	1 (5%)
2	FUC	I	3	2	10,10,11	0.67	0	14,14,16	0.68	0
3	NAG	J	1	3,1	14,14,15	0.74	0	17,19,21	0.92	1 (5%)
3	NAG	J	2	3	14,14,15	0.70	0	17,19,21	0.83	1 (5%)
5	NAG	K	1	1,5	14,14,15	1.02	1 (7%)	17,19,21	1.33	3 (17%)
5	NAG	K	2	5	14,14,15	0.77	0	17,19,21	0.82	1 (5%)
5	BMA	K	3	5	11,11,12	0.49	0	15,15,17	0.42	0
6	NAG	L	1	6,1	14,14,15	1.10	1 (7%)	17,19,21	1.22	2 (11%)
6	NDG	L	2	6	14,14,15	1.01	1 (7%)	17,19,21	0.91	1 (5%)
6	BMA	L	3	6	11,11,12	0.63	0	15,15,17	0.58	0
7	NAG	M	1	1,7	14,14,15	0.71	0	17,19,21	0.99	2 (11%)
7	NDG	M	2	7	14,14,15	0.57	0	17,19,21	1.16	2 (11%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	C	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	FUC	C	3	2	-	-	0/1/1/1
2	NAG	D	1	2,1	1/1/5/7	4/6/23/26	0/1/1/1
2	NAG	D	2	2	-	3/6/23/26	0/1/1/1
2	FUC	D	3	2	-	-	0/1/1/1
3	NAG	E	1	3,1	1/1/5/7	4/6/23/26	0/1/1/1
3	NAG	E	2	3	-	4/6/23/26	0/1/1/1
4	NAG	F	1	4,1	-	0/6/23/26	0/1/1/1
4	NDG	F	2	4	-	0/6/23/26	0/1/1/1
4	MAN	F	3	4	-	2/2/19/22	0/1/1/1
3	NAG	G	1	3,1	-	4/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	H	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	H	2	3	-	5/6/23/26	0/1/1/1
2	NAG	I	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	I	2	2	-	0/6/23/26	0/1/1/1
2	FUC	I	3	2	-	-	0/1/1/1
3	NAG	J	1	3,1	1/1/5/7	1/6/23/26	0/1/1/1
3	NAG	J	2	3	-	4/6/23/26	0/1/1/1
5	NAG	K	1	1,5	1/1/5/7	0/6/23/26	0/1/1/1
5	NAG	K	2	5	-	2/6/23/26	0/1/1/1
5	BMA	K	3	5	-	0/2/19/22	0/1/1/1
6	NAG	L	1	6,1	-	2/6/23/26	0/1/1/1
6	NDG	L	2	6	-	2/6/23/26	0/1/1/1
6	BMA	L	3	6	-	0/2/19/22	0/1/1/1
7	NAG	M	1	1,7	1/1/5/7	4/6/23/26	0/1/1/1
7	NDG	M	2	7	-	2/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	1	NAG	C1-C2	2.81	1.56	1.52
2	D	1	NAG	C1-C2	2.63	1.55	1.52
2	D	2	NAG	C1-C2	2.52	1.55	1.52
5	K	1	NAG	C1-C2	2.51	1.55	1.52
3	E	1	NAG	C1-C2	2.12	1.55	1.52
6	L	2	NDG	O5-C5	2.02	1.47	1.43

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	2	NDG	C2-N2-C7	-3.54	118.15	122.90
5	K	1	NAG	C1-O5-C5	3.29	116.60	112.19
2	C	1	NAG	C1-O5-C5	3.18	116.45	112.19
6	L	1	NAG	C2-N2-C7	-3.10	118.75	122.90
3	E	1	NAG	C1-O5-C5	3.07	116.30	112.19
2	I	1	NAG	C1-C2-N2	2.94	115.07	110.43
2	I	2	NAG	C2-N2-C7	-2.90	119.02	122.90
2	I	1	NAG	O5-C1-C2	-2.80	106.95	111.29
4	F	1	NAG	C2-N2-C7	-2.72	119.25	122.90
7	M	1	NAG	C2-N2-C7	-2.68	119.31	122.90
7	M	2	NDG	C1-O5-C5	2.63	115.71	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	F	2	NDG	C1-O5-C5	2.63	115.71	112.19
4	F	2	NDG	C6-C5-C4	-2.63	106.57	113.02
3	E	1	NAG	C2-N2-C7	-2.61	119.40	122.90
5	K	1	NAG	C2-N2-C7	-2.45	119.62	122.90
3	E	2	NAG	C2-N2-C7	-2.45	119.62	122.90
7	M	2	NDG	C2-N2-C7	-2.42	119.66	122.90
5	K	2	NAG	C2-N2-C7	-2.42	119.66	122.90
3	J	2	NAG	C2-N2-C7	-2.35	119.75	122.90
6	L	1	NAG	C3-C4-C5	-2.32	106.03	110.23
3	G	1	NAG	C2-N2-C7	-2.28	119.84	122.90
3	J	1	NAG	C2-N2-C7	-2.23	119.91	122.90
5	K	1	NAG	O5-C1-C2	2.13	114.59	111.29
2	C	2	NAG	C2-N2-C7	-2.10	120.08	122.90
7	M	1	NAG	O5-C1-C2	-2.09	108.05	111.29
6	L	2	NDG	C2-N2-C7	-2.07	120.13	122.90

All (5) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	D	1	NAG	C1
3	E	1	NAG	C1
3	J	1	NAG	C1
5	K	1	NAG	C1
7	M	1	NAG	C1

All (52) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	H	2	NAG	C1-C2-N2-C7
3	H	2	NAG	O7-C7-N2-C2
3	H	2	NAG	C8-C7-N2-C2
3	J	2	NAG	O5-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
7	M	2	NDG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
7	M	2	NDG	C4-C5-C6-O6
3	J	2	NAG	C4-C5-C6-O6
6	L	2	NDG	C4-C5-C6-O6
7	M	1	NAG	O5-C5-C6-O6
2	D	1	NAG	C8-C7-N2-C2
2	D	1	NAG	O7-C7-N2-C2
3	J	2	NAG	C8-C7-N2-C2

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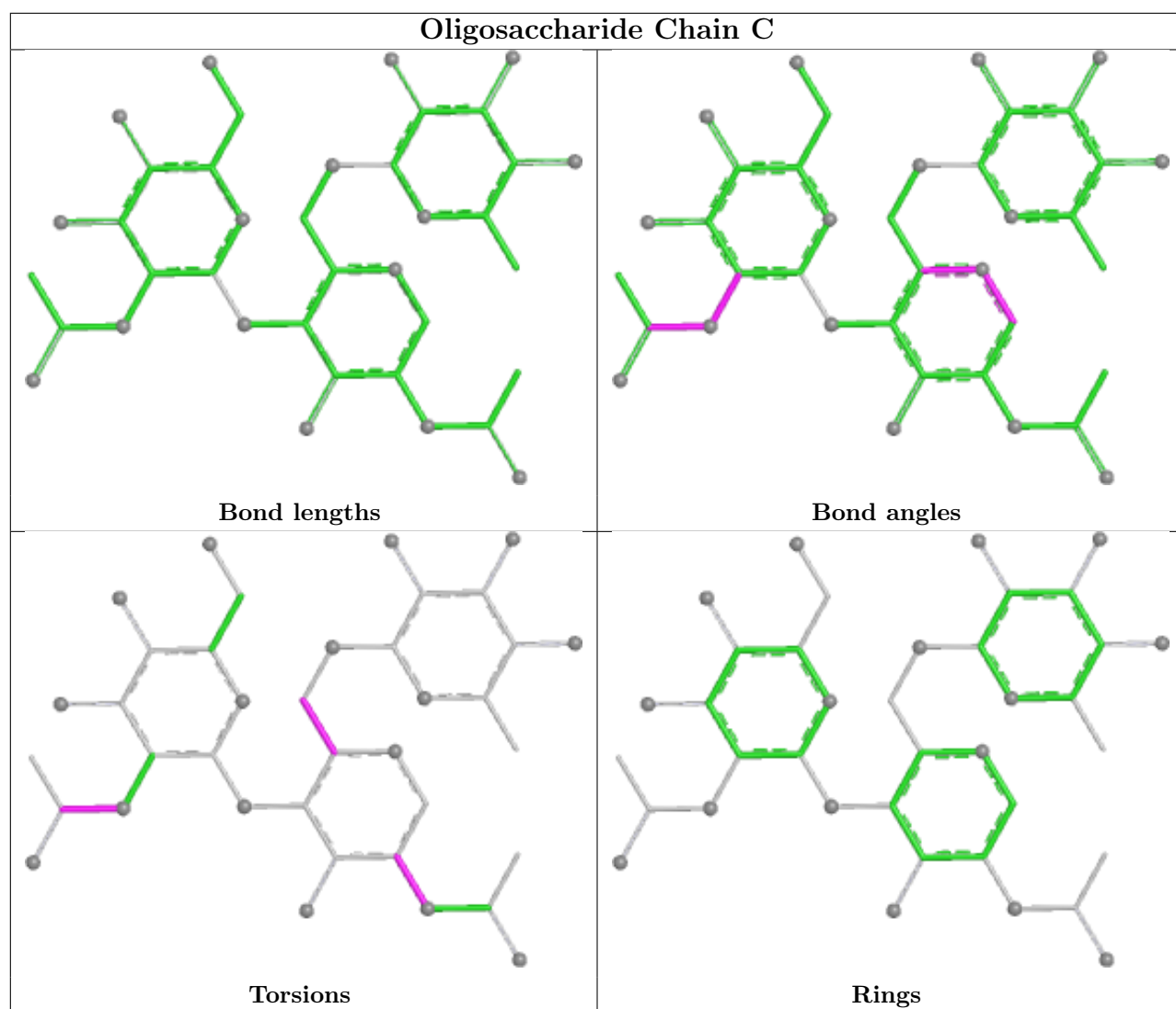
Mol	Chain	Res	Type	Atoms
5	K	2	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
5	K	2	NAG	C4-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	J	2	NAG	O7-C7-N2-C2
4	F	3	MAN	O5-C5-C6-O6
6	L	2	NDG	O5-C5-C6-O6
2	C	2	NAG	C8-C7-N2-C2
4	F	3	MAN	C4-C5-C6-O6
3	E	2	NAG	O5-C5-C6-O6
3	E	1	NAG	C8-C7-N2-C2
6	L	1	NAG	C4-C5-C6-O6
2	C	2	NAG	O7-C7-N2-C2
6	L	1	NAG	O5-C5-C6-O6
7	M	1	NAG	C4-C5-C6-O6
3	E	1	NAG	O7-C7-N2-C2
3	J	1	NAG	O5-C5-C6-O6
3	E	1	NAG	C4-C5-C6-O6
3	G	1	NAG	C8-C7-N2-C2
2	D	2	NAG	C3-C2-N2-C7
3	E	1	NAG	O5-C5-C6-O6
2	D	1	NAG	C1-C2-N2-C7
3	G	2	NAG	C1-C2-N2-C7
3	G	1	NAG	O7-C7-N2-C2
2	C	1	NAG	O5-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
2	D	2	NAG	C8-C7-N2-C2
2	D	1	NAG	C3-C2-N2-C7
3	G	2	NAG	C3-C2-N2-C7
3	E	2	NAG	C8-C7-N2-C2
2	D	2	NAG	O7-C7-N2-C2
3	G	1	NAG	O5-C5-C6-O6
2	C	1	NAG	C1-C2-N2-C7
7	M	1	NAG	C8-C7-N2-C2
3	E	2	NAG	O7-C7-N2-C2
7	M	1	NAG	O7-C7-N2-C2
3	H	1	NAG	C4-C5-C6-O6

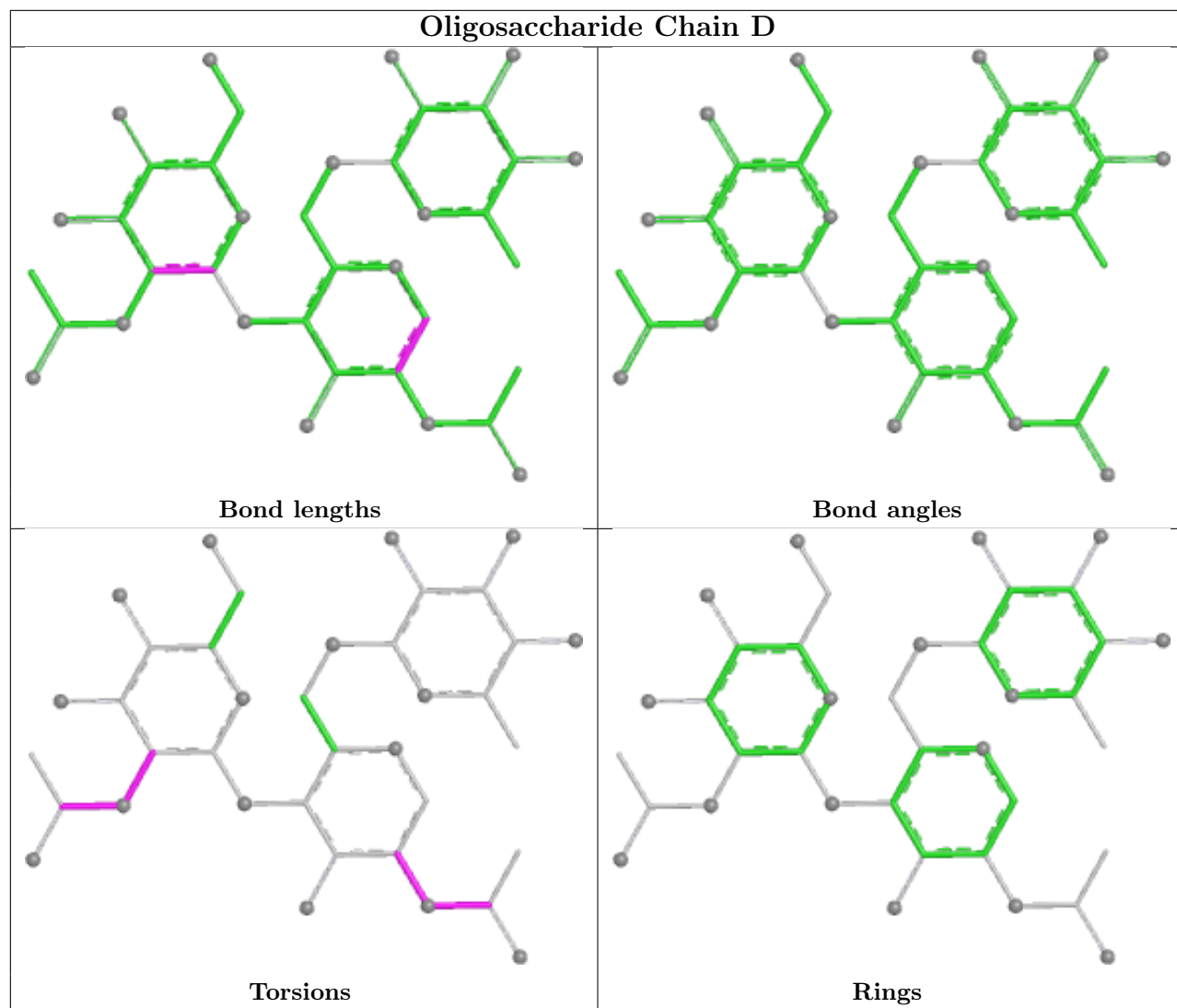
There are no ring outliers.

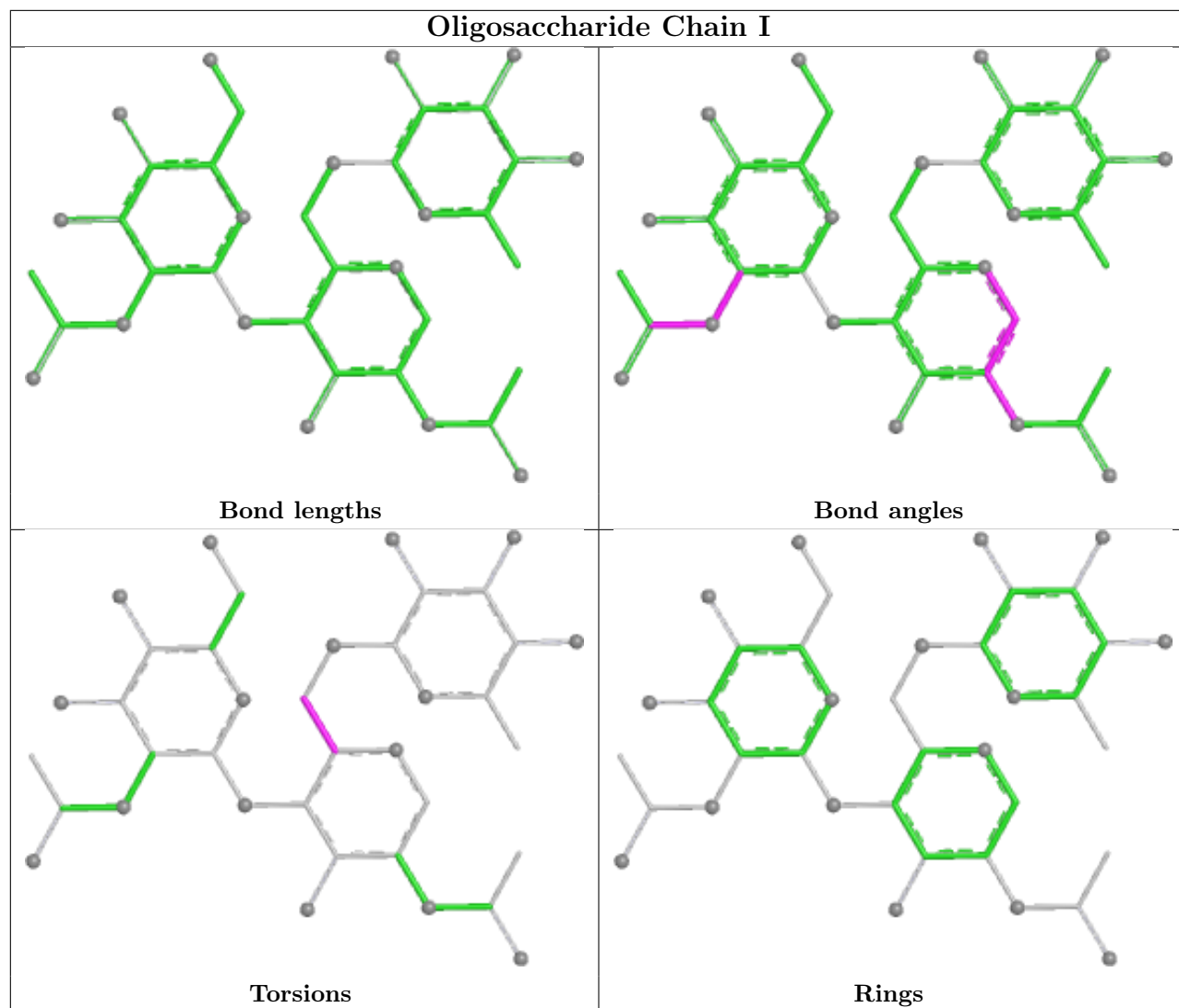
16 monomers are involved in 23 short contacts:

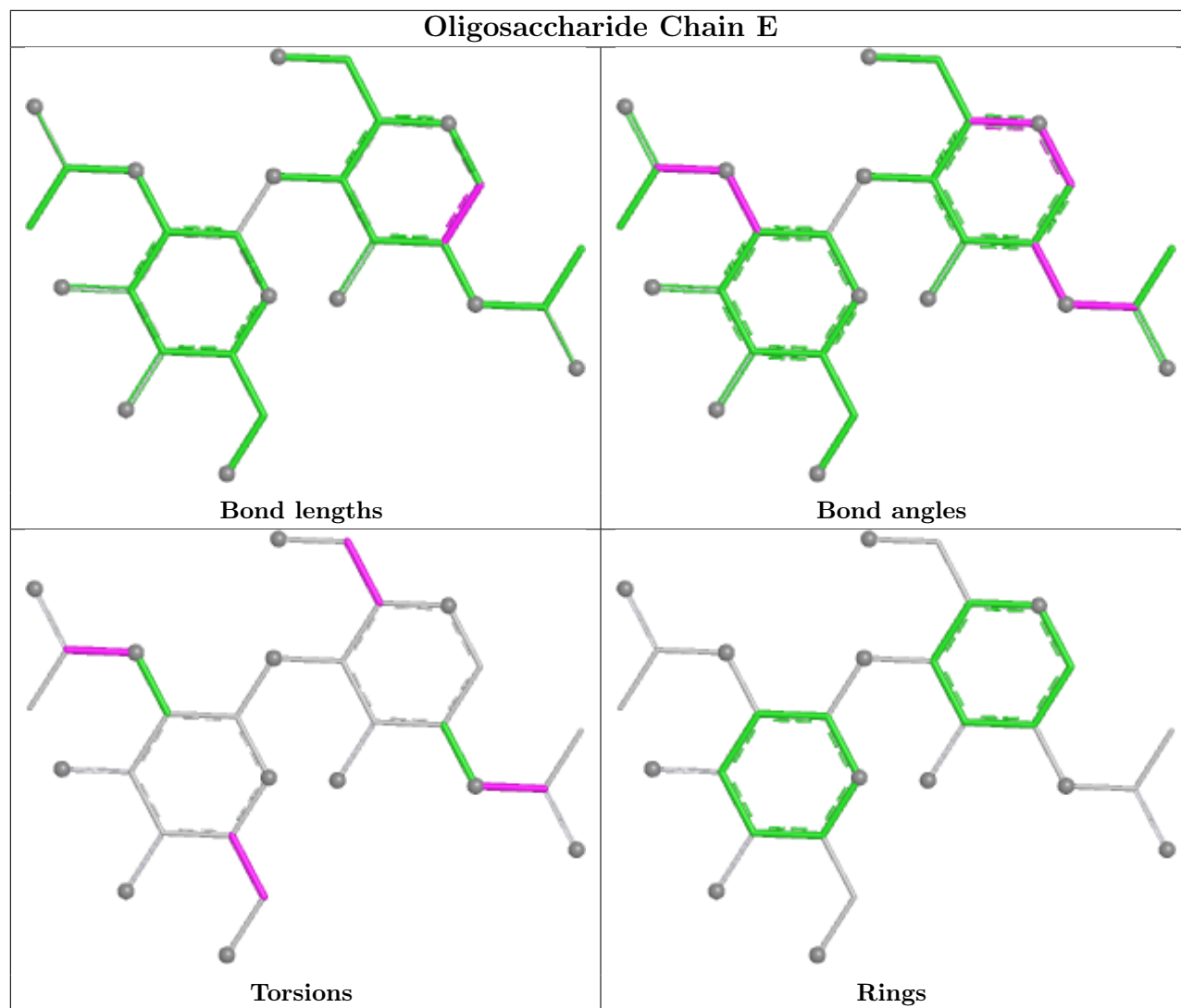
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1	NAG	1	0
7	M	1	NAG	2	0
2	D	1	NAG	2	0
2	C	1	NAG	2	0
4	F	1	NAG	3	0
3	G	2	NAG	2	0
2	I	2	NAG	1	0
2	I	3	FUC	2	0
4	F	2	NDG	4	0
2	C	2	NAG	4	0
6	L	2	NDG	1	0
2	C	3	FUC	2	0
3	J	1	NAG	1	0
4	F	3	MAN	2	0
2	D	2	NAG	3	0
2	I	1	NAG	2	0

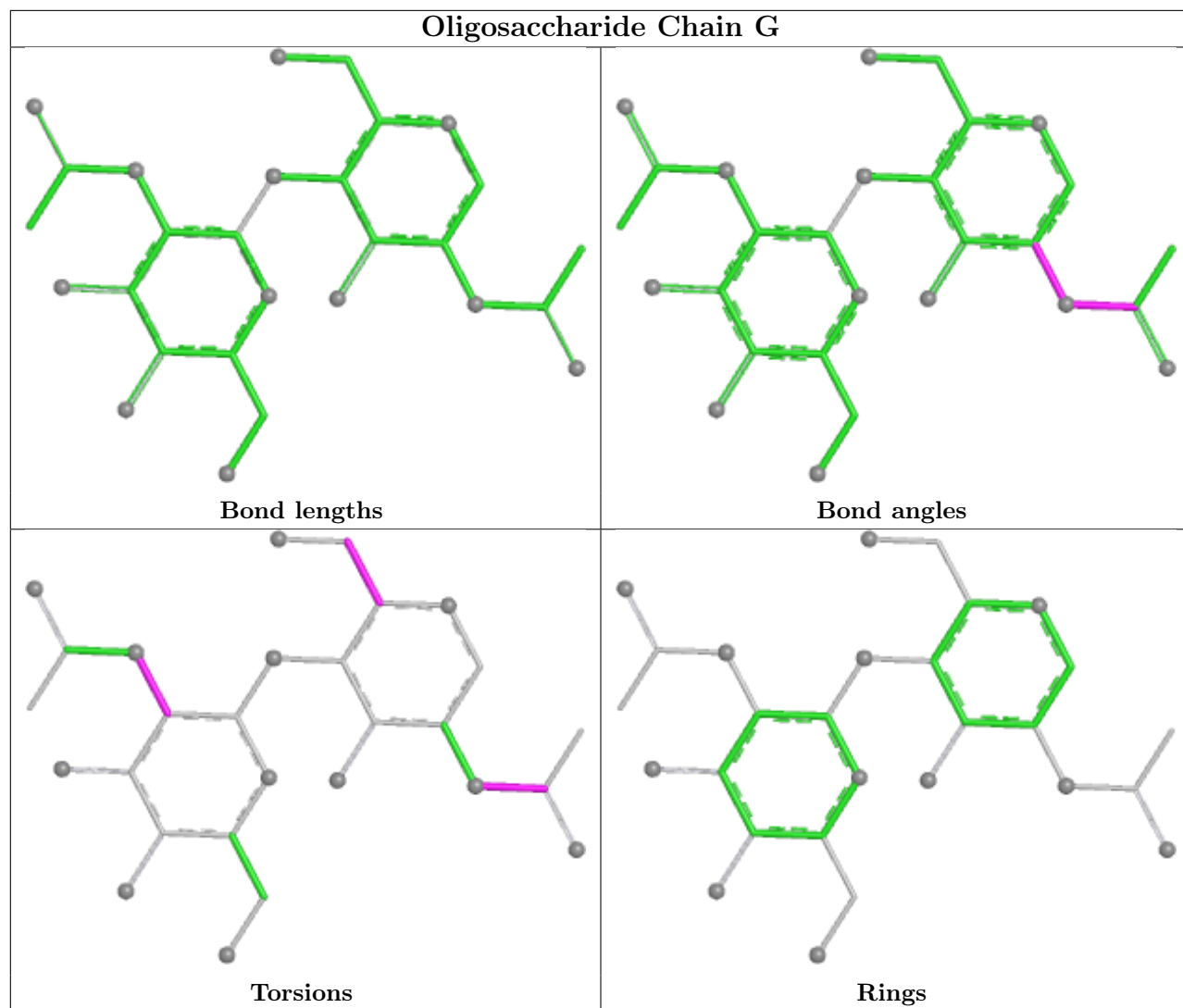
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

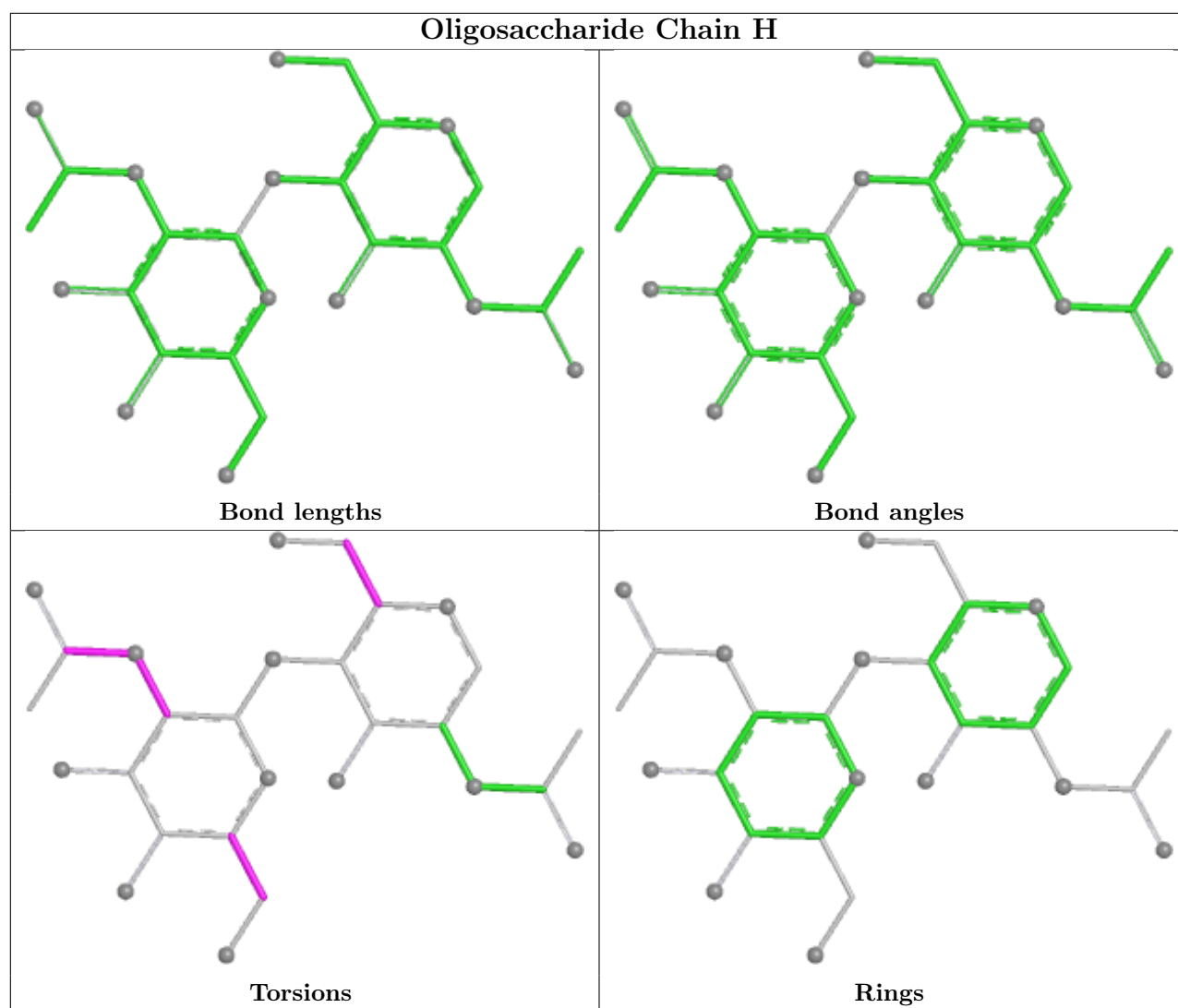


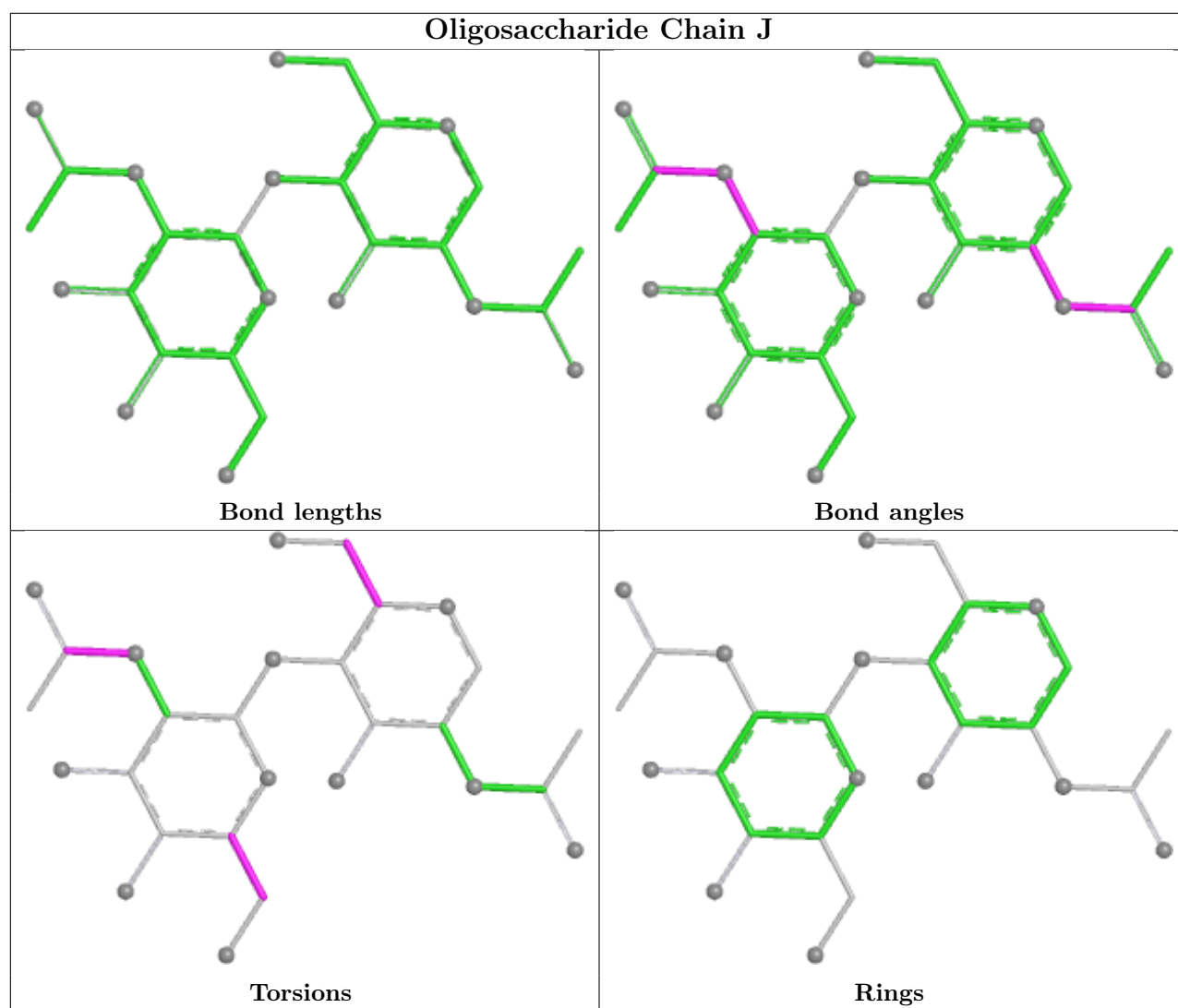


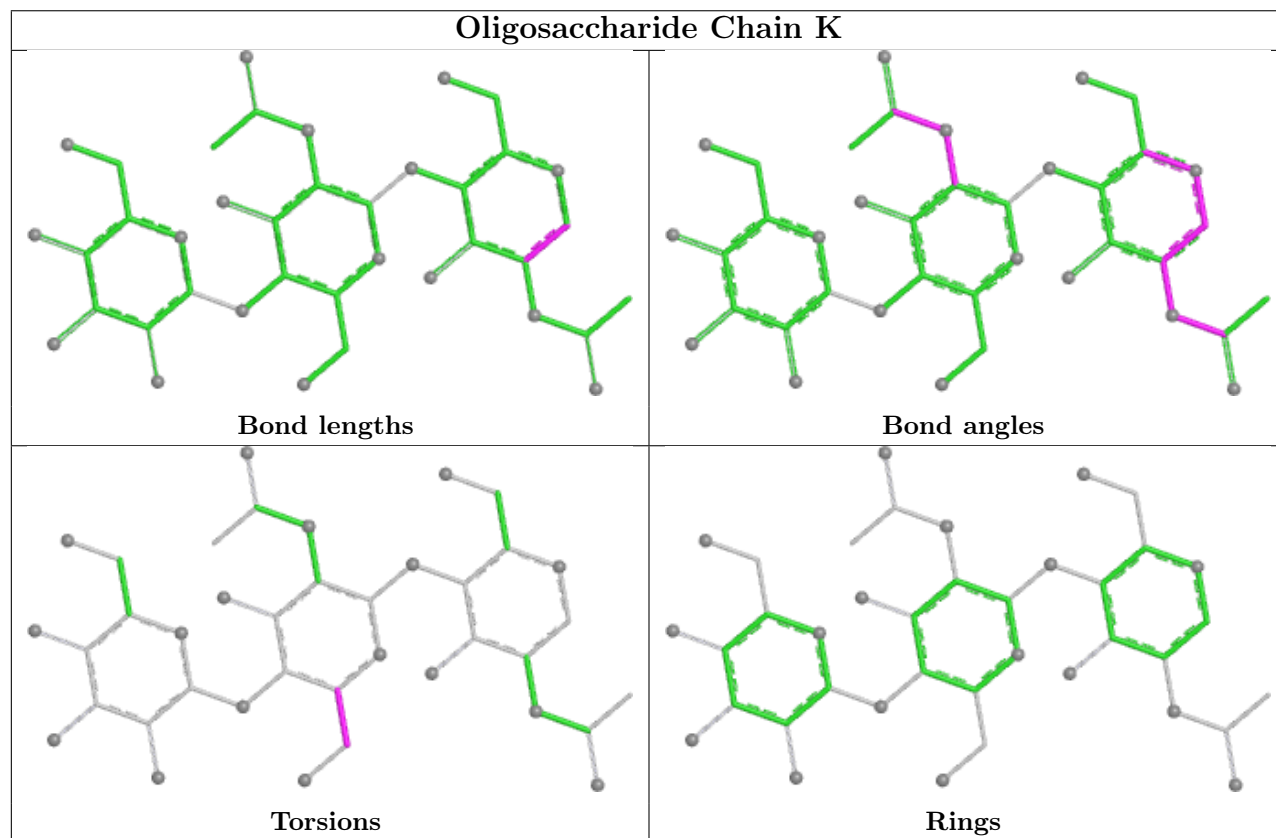
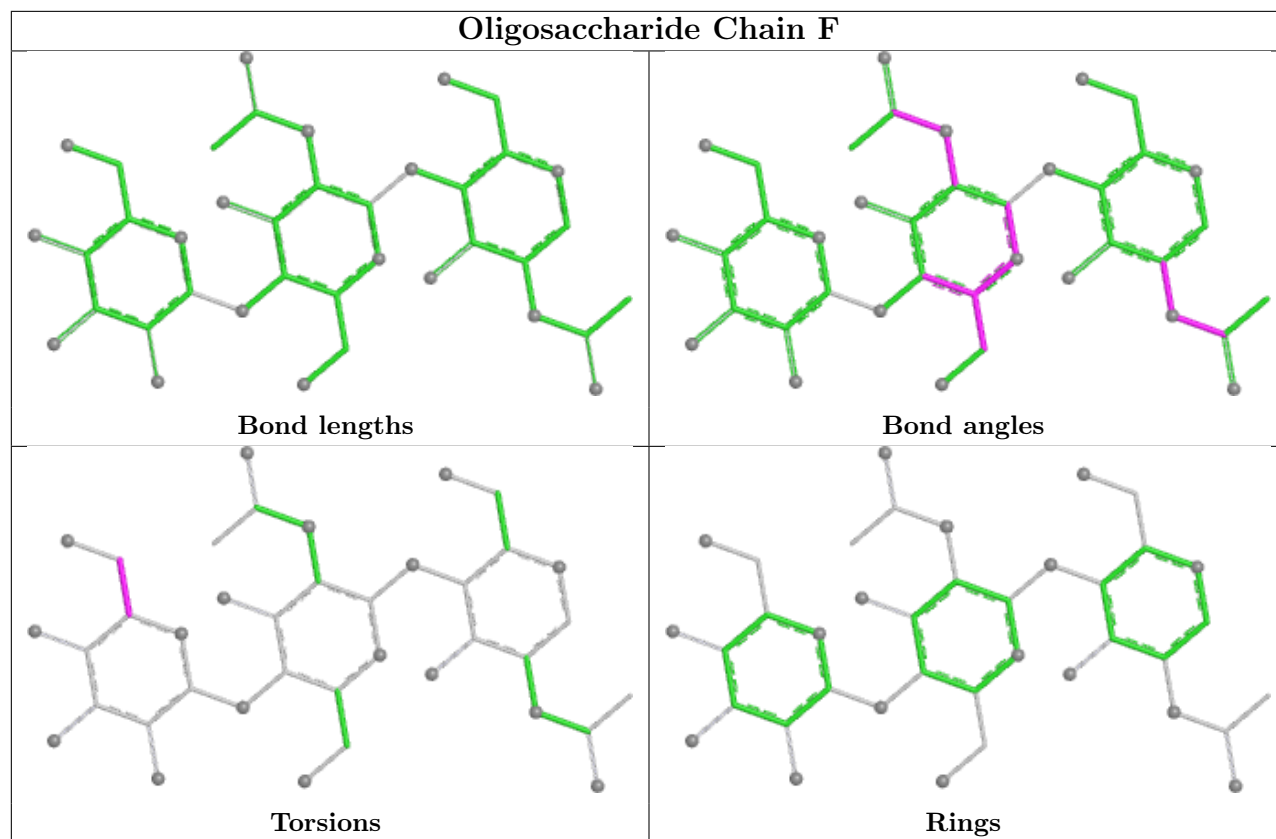


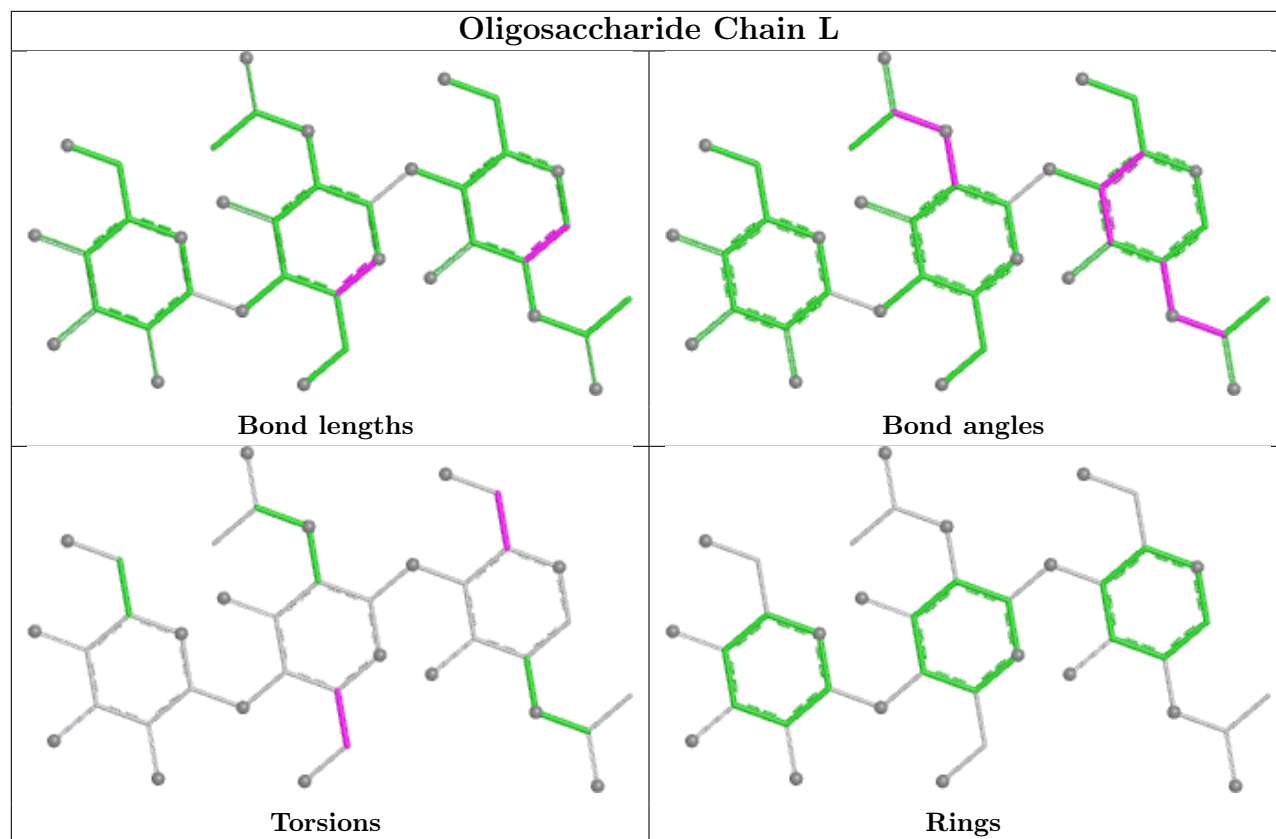


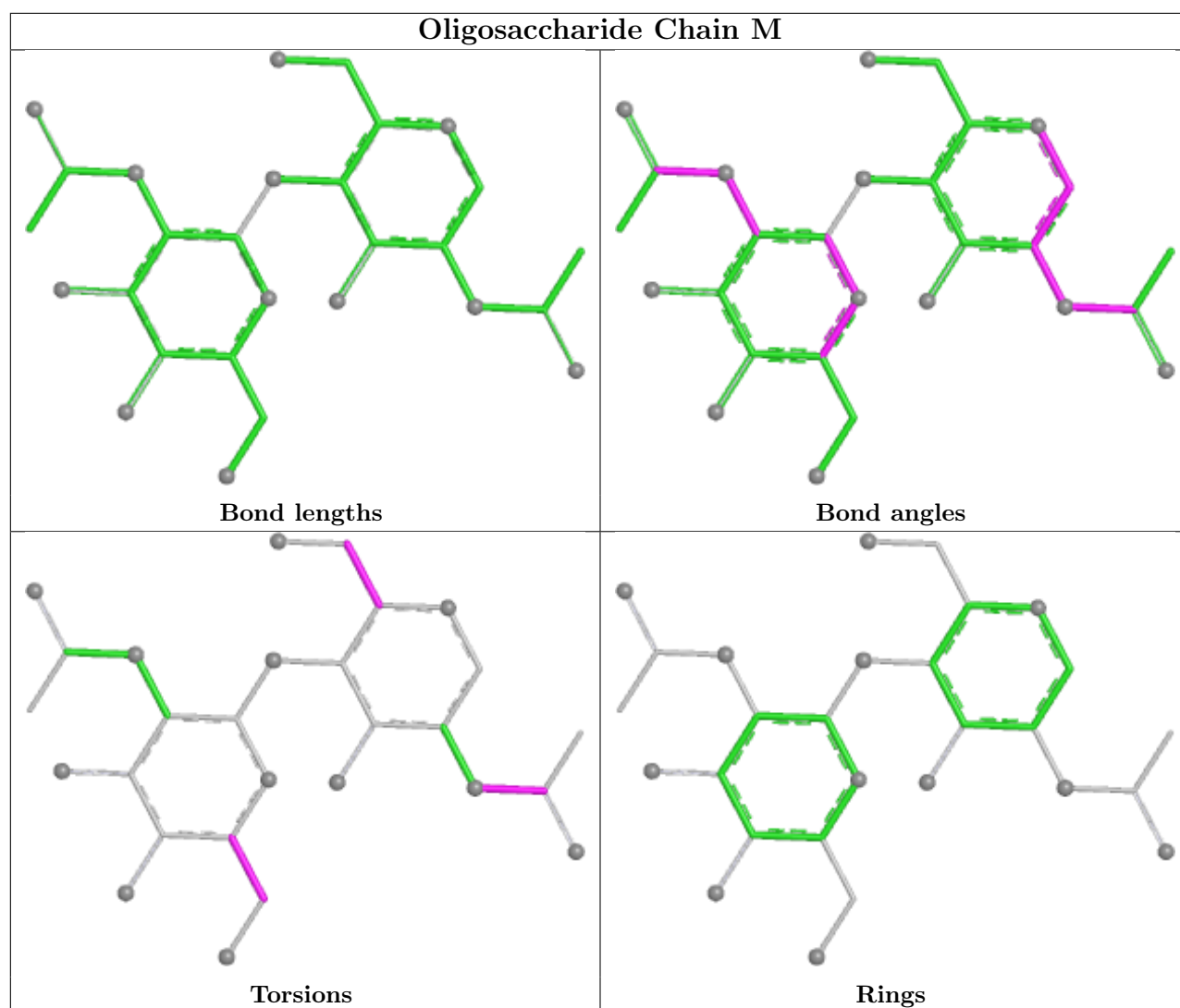












5.6 Ligand geometry [i](#)

3 ligands 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	$\# Z > 2$
8	NAG	B	770	1	14,14,15	0.62	0	17,19,21	0.77	1 (5%)
8	NAG	B	781	1	14,14,15	0.96	1 (7%)	17,19,21	0.77	0
8	NAG	A	782	1	14,14,15	0.78	1 (7%)	17,19,21	0.69	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
8	NAG	B	770	1	-	1/6/23/26	0/1/1/1
8	NAG	B	781	1	-	3/6/23/26	0/1/1/1
8	NAG	A	782	1	-	3/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	781	NAG	C1-C2	2.79	1.56	1.52
8	A	782	NAG	C1-C2	2.38	1.55	1.52

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	770	NAG	C2-N2-C7	-2.23	119.91	122.90

There are no chirality outliers.

All (7) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	782	NAG	C3-C2-N2-C7
8	B	781	NAG	C3-C2-N2-C7
8	A	782	NAG	O5-C5-C6-O6
8	A	782	NAG	C4-C5-C6-O6
8	B	781	NAG	C8-C7-N2-C2
8	B	781	NAG	O7-C7-N2-C2
8	B	770	NAG	C4-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 15 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	B	770	NAG	2	0
8	B	781	NAG	8	0
8	A	782	NAG	5	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	726/728 (99%)	0.11	24 (3%)	49 46	9, 24, 49, 68	0
1	B	728/728 (100%)	0.15	35 (4%)	35 32	9, 24, 53, 74	0
All	All	1454/1456 (99%)	0.13	59 (4%)	41 38	9, 24, 51, 74	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	280	THR	4.1
1	A	377	ASN	4.0
1	B	76	ILE	4.0
1	A	289	ALA	3.9
1	A	39	SER	3.7
1	B	289	ALA	3.6
1	B	73	GLU	3.6
1	A	520	ASN	3.4
1	B	94	THR	3.4
1	B	98	PHE	3.4
1	A	393	ASP	3.3
1	A	276	LEU	3.3
1	A	92	ASN	3.3
1	A	342	ALA	3.2
1	B	39	SER	3.1
1	B	379	GLU	3.0
1	A	392	LYS	3.0
1	B	399	LYS	3.0
1	A	279	VAL	2.9
1	A	105	TYR	2.9
1	B	385	CYS	2.8
1	B	388	GLN	2.8
1	A	90	LEU	2.8
1	B	95	PHE	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	208	PHE	2.7
1	A	94	THR	2.7
1	A	277	SER	2.6
1	B	88	VAL	2.6
1	B	644	SER	2.6
1	B	91	GLU	2.5
1	B	99	GLY	2.5
1	B	105	TYR	2.5
1	A	96	ASP	2.4
1	B	537	SER	2.4
1	B	90	LEU	2.4
1	A	83	TYR	2.4
1	B	72	GLN	2.3
1	A	278	SER	2.3
1	A	71	LYS	2.3
1	B	77	LEU	2.2
1	B	765	LEU	2.2
1	B	333	SER	2.2
1	A	73	GLU	2.2
1	B	92	ASN	2.2
1	B	413	ASP	2.1
1	A	764	SER	2.1
1	B	471	ARG	2.1
1	A	93	SER	2.1
1	B	87	SER	2.1
1	A	521	GLU	2.1
1	A	74	ASN	2.1
1	B	89	PHE	2.1
1	B	154	TRP	2.1
1	B	487	ASN	2.1
1	A	89	PHE	2.0
1	B	97	GLU	2.0
1	B	100	HIS	2.0
1	B	517	ILE	2.0
1	B	766	PRO	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 ⓘ

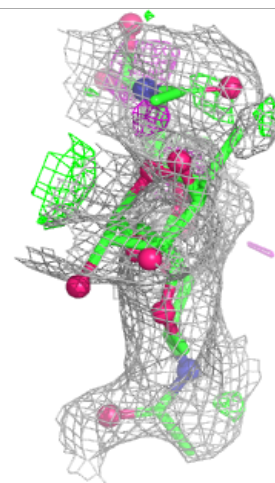
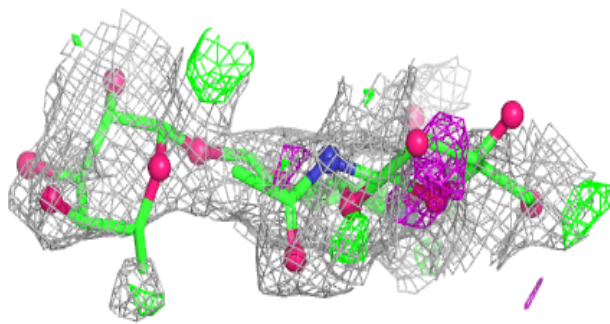
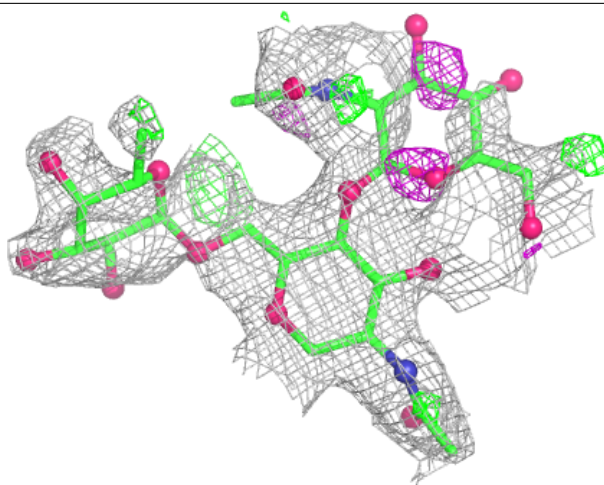
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	D	2	14/15	0.19	0.21	80,81,83,84	0
2	NAG	C	2	14/15	0.32	0.25	78,79,81,81	0
2	FUC	D	3	10/11	0.35	0.20	78,79,79,79	0
2	NAG	I	2	14/15	0.38	0.19	80,83,84,85	0
4	MAN	F	3	11/12	0.42	0.20	72,73,74,74	0
2	FUC	I	3	10/11	0.45	0.22	77,79,79,79	0
5	BMA	K	3	11/12	0.45	0.18	73,75,77,77	0
6	BMA	L	3	11/12	0.45	0.20	80,82,83,84	0
3	NAG	H	2	14/15	0.47	0.21	63,65,66,66	0
2	FUC	C	3	10/11	0.50	0.22	77,78,78,78	0
3	NAG	J	2	14/15	0.51	0.20	66,68,68,68	0
6	NDG	L	2	14/15	0.51	0.25	64,68,72,76	0
3	NAG	G	2	14/15	0.51	0.20	73,76,77,77	0
7	NDG	M	2	14/15	0.52	0.19	73,76,79,79	0
2	NAG	D	1	14/15	0.54	0.19	71,76,78,78	0
4	NDG	F	2	14/15	0.59	0.18	62,65,67,70	0
7	NAG	M	1	14/15	0.68	0.20	56,60,63,69	0
3	NAG	E	2	14/15	0.69	0.17	63,66,66,67	0
3	NAG	J	1	14/15	0.70	0.17	54,59,62,63	0
2	NAG	I	1	14/15	0.72	0.14	67,69,76,76	0
5	NAG	K	2	14/15	0.73	0.16	58,61,66,70	0
3	NAG	G	1	14/15	0.74	0.15	63,66,67,70	0
3	NAG	E	1	14/15	0.77	0.14	47,51,54,59	0
2	NAG	C	1	14/15	0.81	0.14	63,66,74,74	0
3	NAG	H	1	14/15	0.82	0.13	47,51,55,59	0
6	NAG	L	1	14/15	0.85	0.14	37,41,48,56	0
5	NAG	K	1	14/15	0.86	0.12	30,39,45,51	0
4	NAG	F	1	14/15	0.87	0.11	43,45,49,56	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.

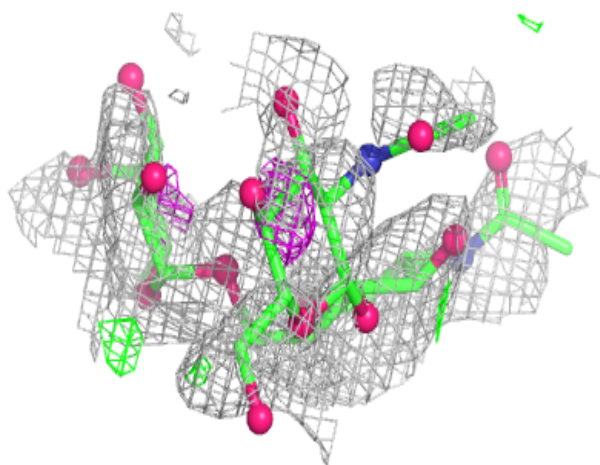
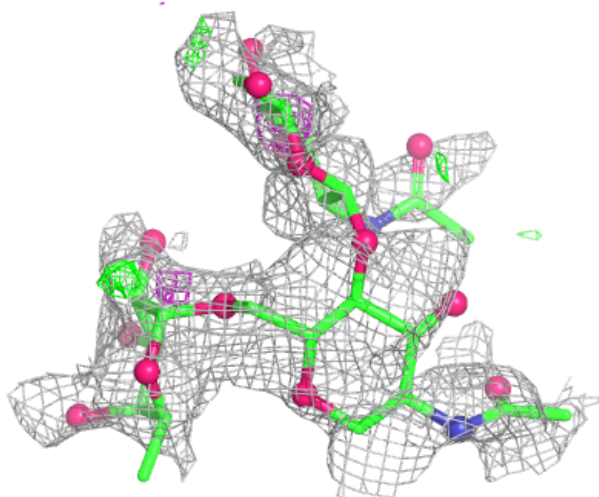
Electron density around Chain C:

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



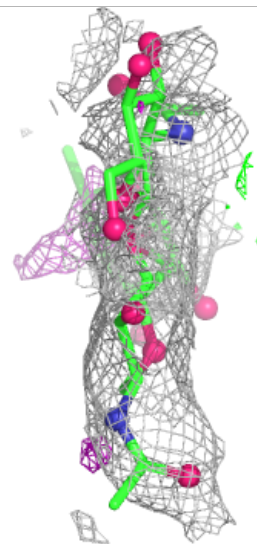
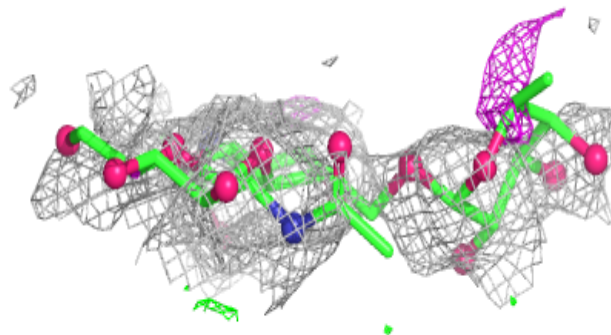
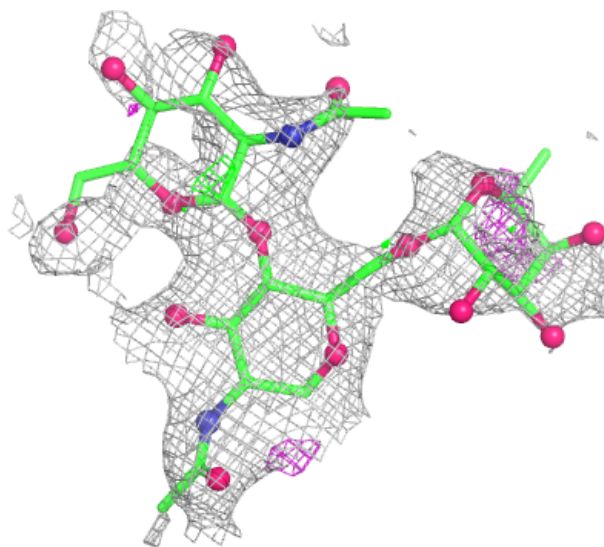
Electron density around Chain D:

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



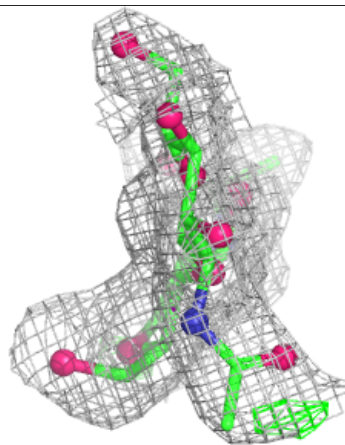
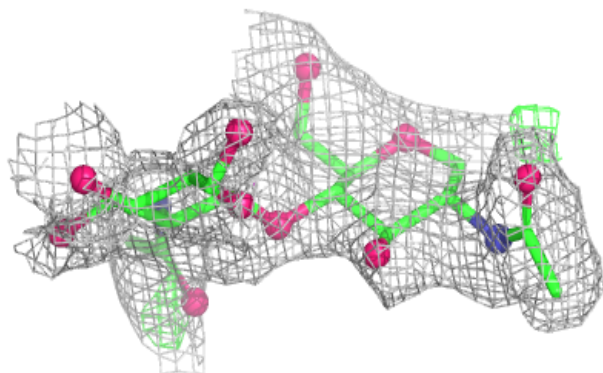
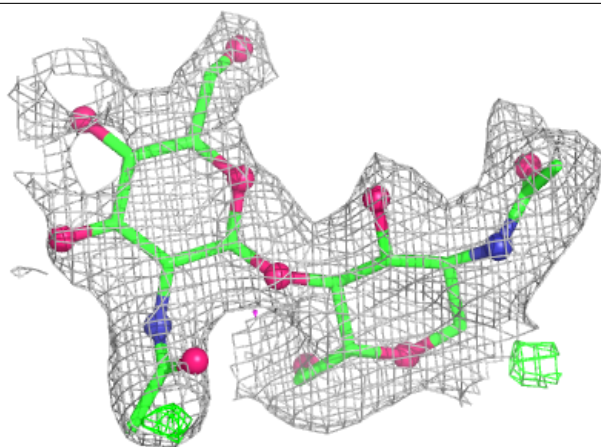
Electron density around Chain I:

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



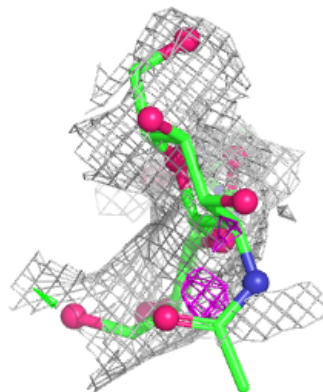
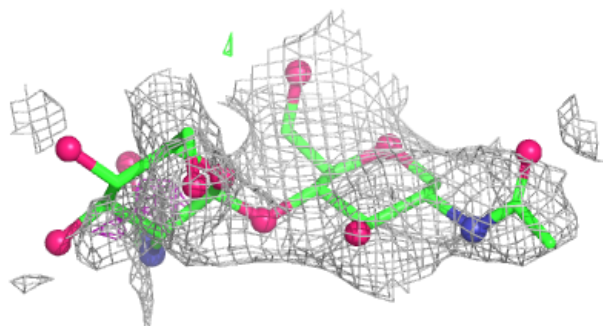
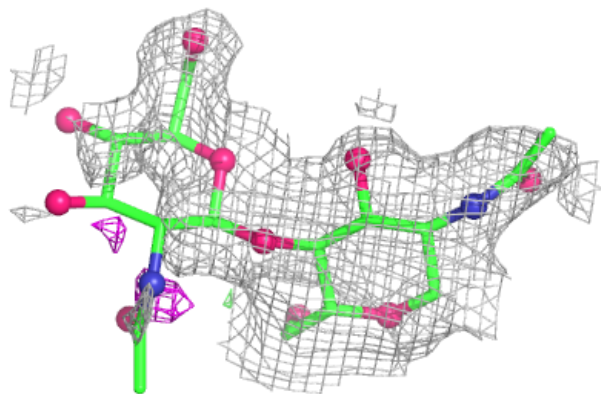
Electron density around Chain E:

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



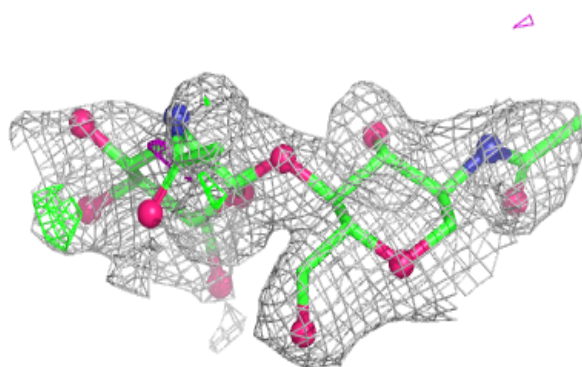
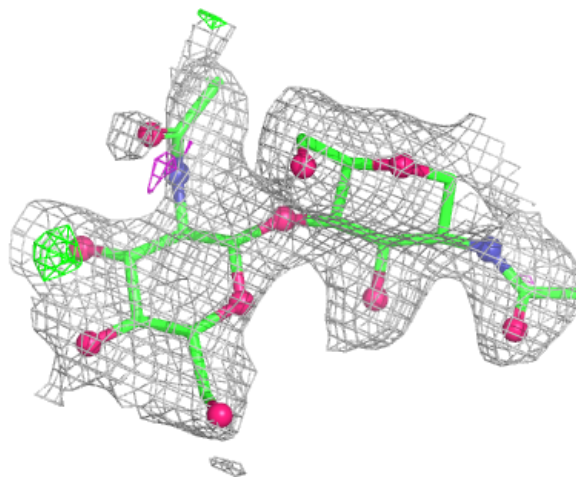
Electron density around Chain G:

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



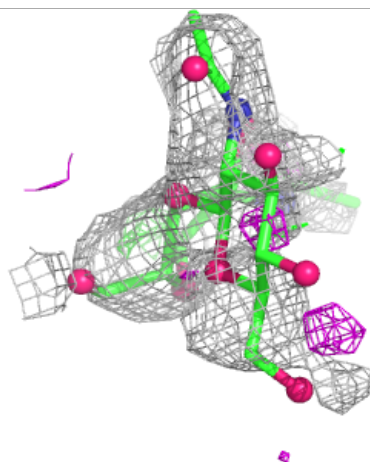
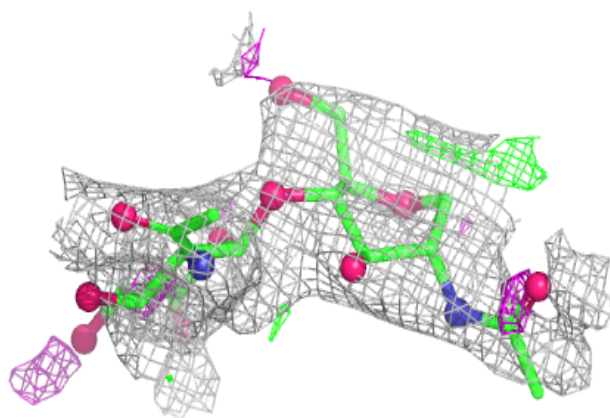
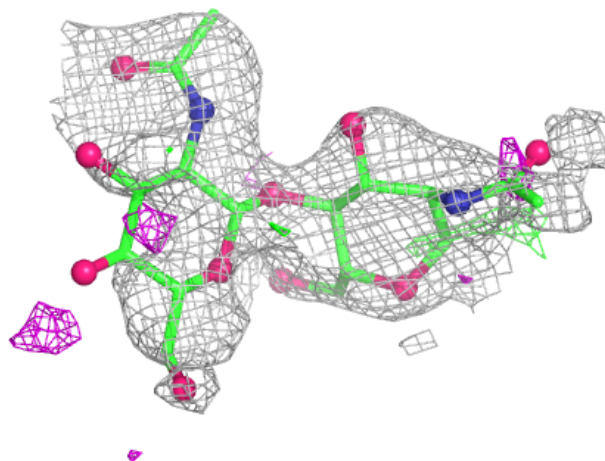
Electron density around Chain H:

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



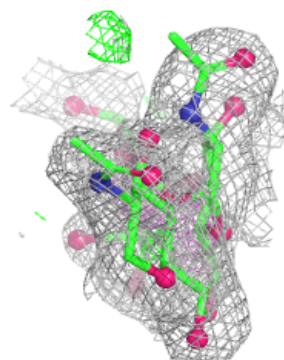
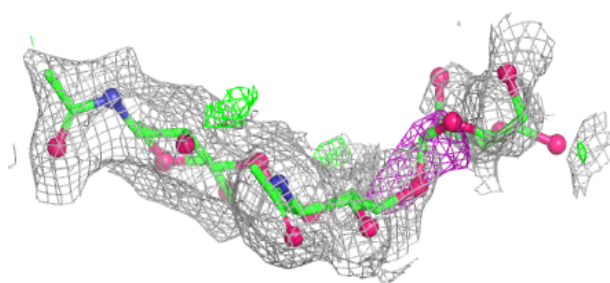
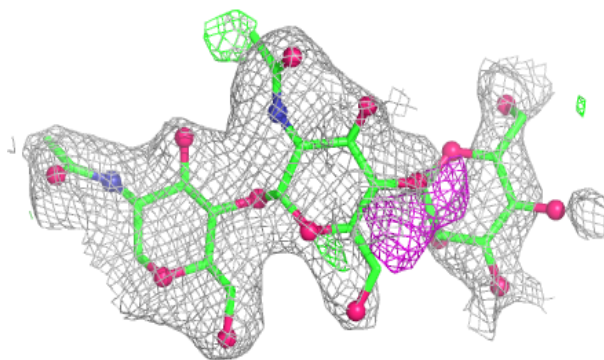
Electron density around Chain J:

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

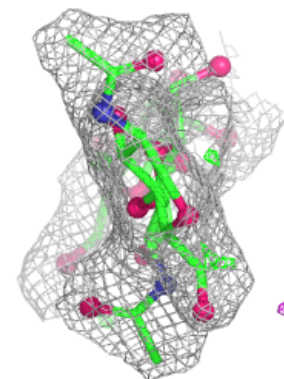
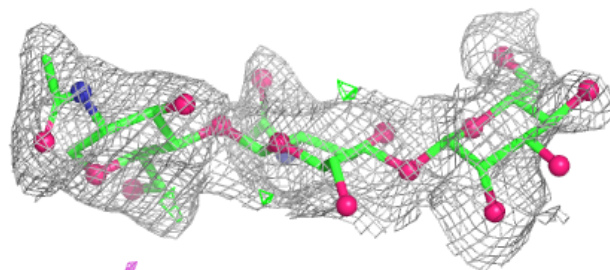
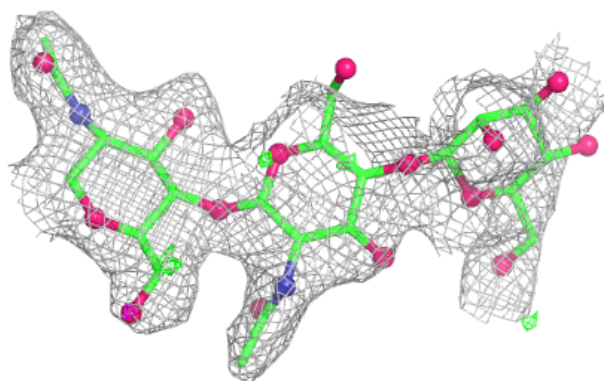


Electron density around Chain F:

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

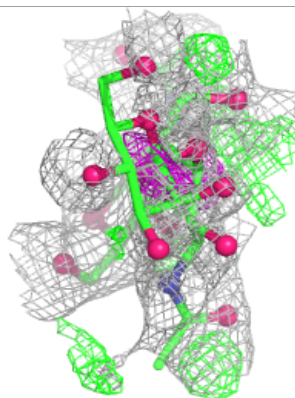
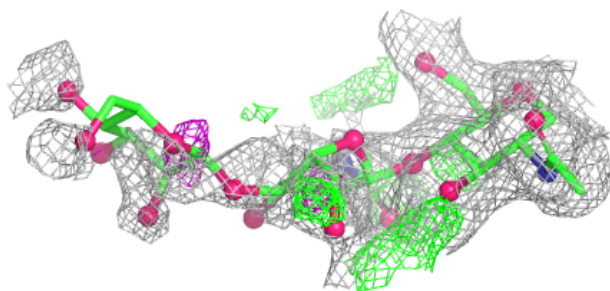
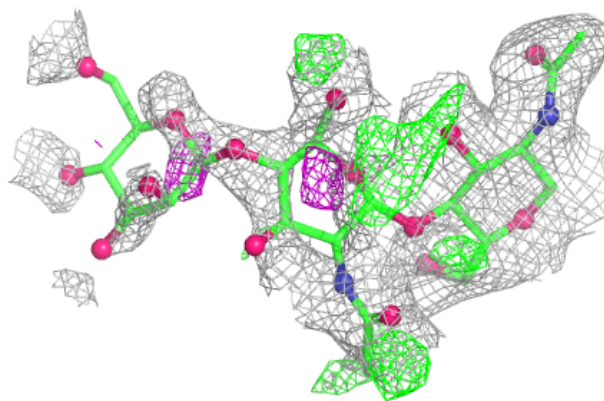
**Electron density around Chain K:**

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

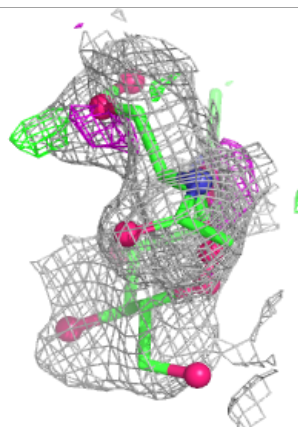
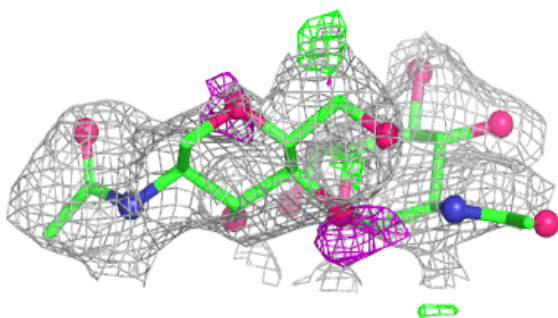
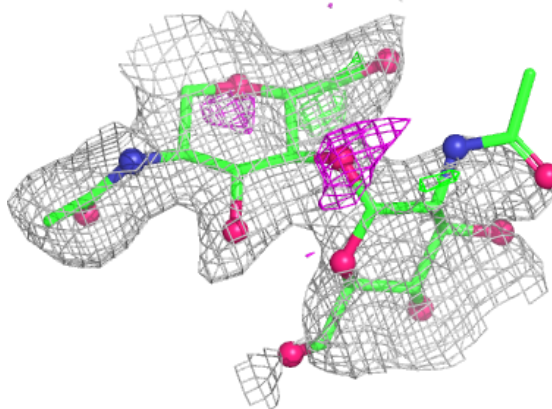


Electron density around Chain L:

$2mF_o-DF_c$ (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)



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
8	NAG	A	782	14/15	0.41	0.23	73,76,77,77	0
8	NAG	B	781	14/15	0.46	0.21	66,68,69,69	0
8	NAG	B	770	14/15	0.56	0.18	60,64,65,66	0

6.5 Other polymers [i](#)

There are no such residues in this entry.