



# Full wwPDB X-ray Structure Validation Report ⓘ

Jul 2, 2020 – 04:48 AM BST

PDB ID : 1O7A  
Title : Human beta-Hexosaminidase B  
Authors : Maier, T.; Strater, N.; Schuette, C.; Klingenstein, R.; Sandhoff, K.; Saenger, W.  
Deposited on : 2002-10-29  
Resolution : 2.25 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](mailto: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.

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

MolProbity : 4.02b-467  
Mogul : 1.8.5 (274361), CSD as541be (2020)  
Xtriage (Phenix) : 1.13  
EDS : 2.13.dev2  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
Refmac : 5.8.0158  
CCP4 : 7.0.044 (Gargrove)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.13.dev2

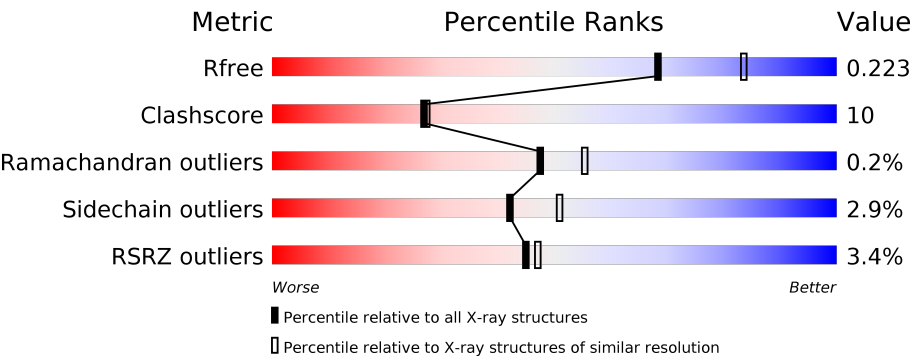
# 1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.25 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))
$R_{free}$	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	515	<div><div>3%</div><div>72%20%6%</div></div>
1	B	515	<div><div>2%</div><div>77%16%6%</div></div>
1	C	515	<div><div>4%</div><div>77%16%6%</div></div>
1	D	515	<div><div>3%</div><div>74%18%6%</div></div>
1	E	515	<div><div>3%</div><div>74%18%6%</div></div>
1	F	515	<div><div>5%</div><div>73%20%6%</div></div>

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Mol	Chain	Length	Quality of chain
2	G	2	 100%
2	H	2	 100%
2	I	2	 50% 50%
2	J	2	 50% 50%
2	K	2	 50% 50%
2	L	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	G	2	-	-	-	X
2	NAG	H	2	-	-	-	X
2	NAG	I	2	-	-	-	X
2	NAG	J	2	-	-	-	X
2	NAG	K	2	-	-	-	X
4	NAG	C	703	-	-	-	X
4	NAG	E	702	-	-	-	X
4	NAG	F	702	-	-	-	X
4	NAG	F	703	-	-	-	X
5	EDO	B	801	-	-	X	-

## 2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 26351 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 BETA-HEXOSAMINIDASE BETA CHAIN.

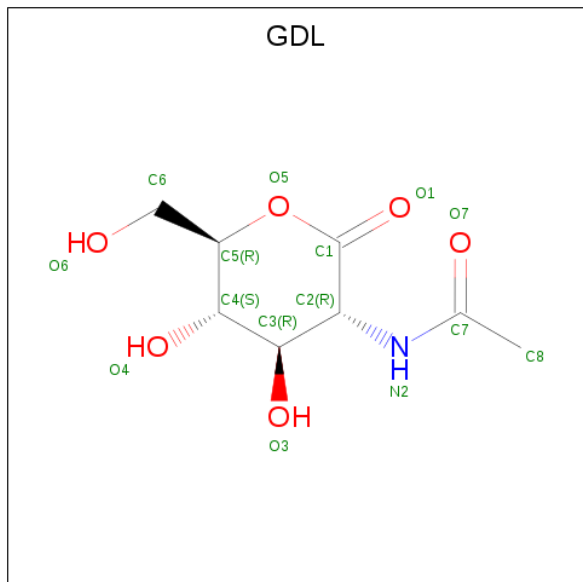
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	484	Total	C	N	O	S	0	4	1
			3933	2538	654	725	16			
1	B	484	Total	C	N	O	S	0	3	1
			3927	2535	653	724	15			
1	C	484	Total	C	N	O	S	0	4	1
			3935	2538	656	727	14			
1	D	483	Total	C	N	O	S	0	6	1
			3939	2542	655	726	16			
1	E	484	Total	C	N	O	S	0	5	1
			3941	2542	655	728	16			
1	F	484	Total	C	N	O	S	0	6	1
			3948	2546	656	730	16			

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



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

- Molecule 3 is 2-(acetylamido)-2-deoxy-D-glucono-1,5-lactone (three-letter code: GDL) (formula:  $C_8H_{13}NO_6$ ).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			15	8	1	6		
3	B	1	Total	C	N	O	0	0
			15	8	1	6		
3	C	1	Total	C	N	O	0	0
			15	8	1	6		
3	D	1	Total	C	N	O	0	0
			15	8	1	6		
3	E	1	Total	C	N	O	0	0
			15	8	1	6		
3	F	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula:  $C_8H_{15}NO_6$ ).



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

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	C	1	Total	C	O	0	0
			4	2	2		
5	E	1	Total	C	O	0	0
			4	2	2		
5	E	1	Total	C	O	0	0
			4	2	2		
5	E	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		
5	F	1	Total	C	O	0	0
			4	2	2		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	F	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

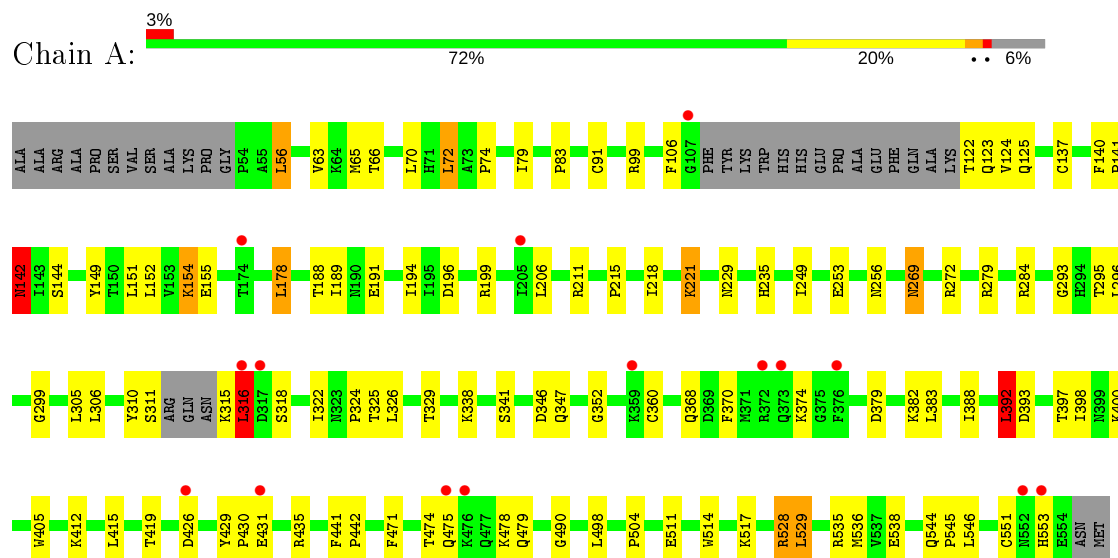
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	385	Total	O	0	0
			385	385		
6	B	399	Total	O	0	0
			399	399		
6	C	385	Total	O	0	0
			385	385		
6	D	346	Total	O	0	0
			346	346		
6	E	448	Total	O	0	0
			448	448		
6	F	363	Total	O	0	0
			363	363		



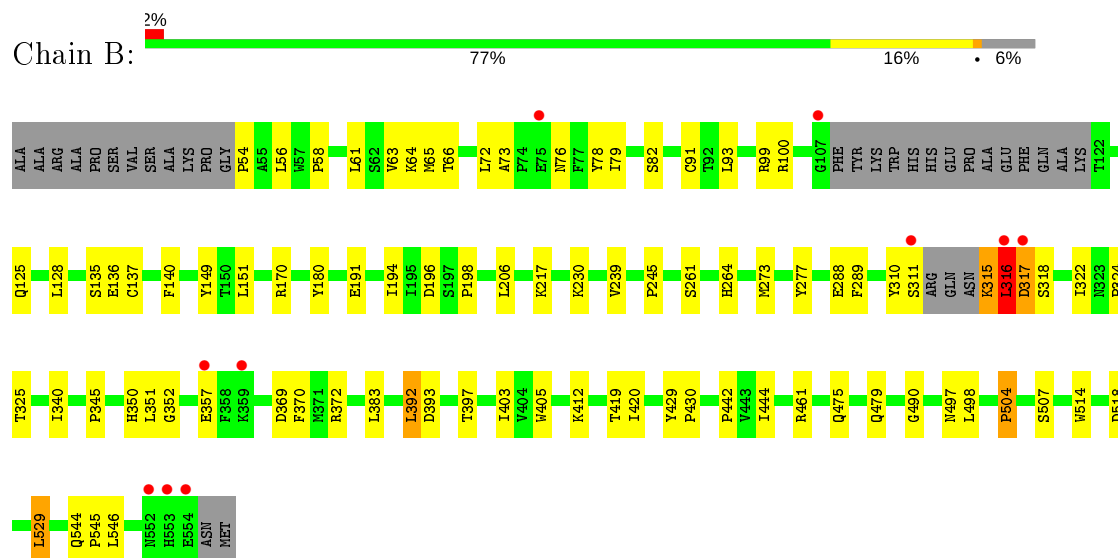
### 3 Residue-property plots [i](#)

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.

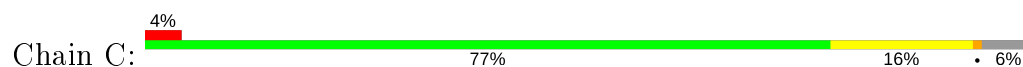
#### • Molecule 1: BETA-HEXOSAMINIDASE BETA CHAIN



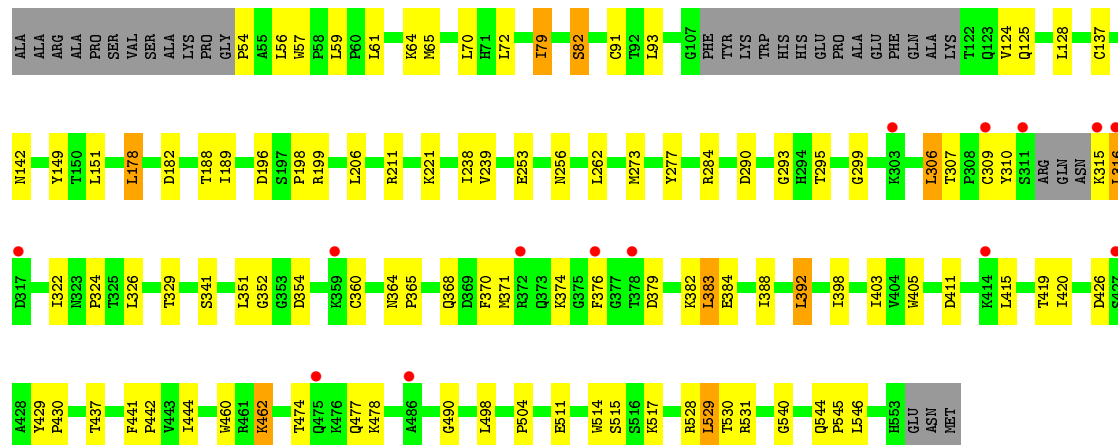
#### • Molecule 1: BETA-HEXOSAMINIDASE BETA CHAIN



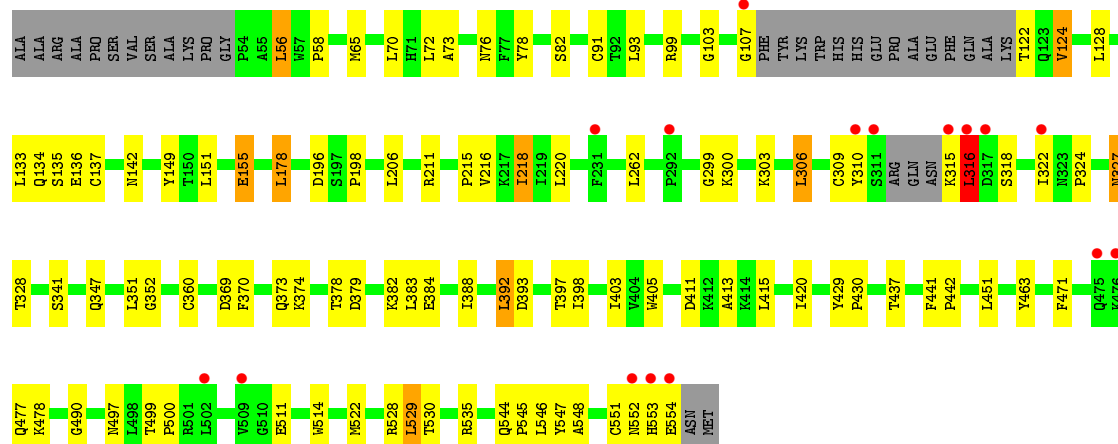
#### • Molecule 1: BETA-HEXOSAMINIDASE BETA CHAIN



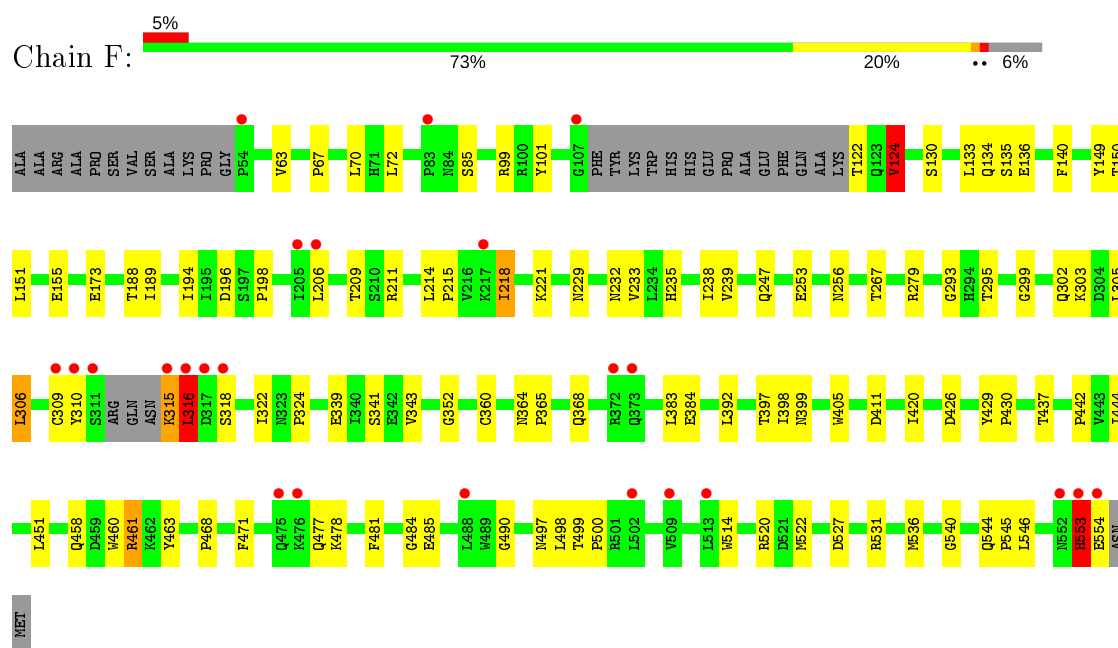
● Molecule 1: BETA-HEXOSAMINIDASE BETA CHAIN



● Molecule 1: BETA-HEXOSAMINIDASE BETA CHAIN



- Molecule 1: BETA-HEXOSAMINIDASE BETA CHAIN



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



MAG1  
MAG2

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



MAG1  
MAG2

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



MAG1  
MAG2

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




MAG1  
MAG2

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

Chain K:  50% 50%

MAG1  
MAG2

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

Chain L:  50% 50%

MAG1  
MAG2

## 4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	163.93Å 163.93Å 244.72Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.12 – 2.25 28.12 – 2.25	Depositor EDS
% Data completeness (in resolution range)	97.4 (28.12-2.25) 97.4 (28.12-2.25)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	4.84 (at 2.24Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.196 , 0.236 0.187 , 0.223	Depositor DCC
$R_{free}$ test set	1715 reflections (0.98%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	35.3	Xtriage
Anisotropy	0.622	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.33 , 47.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.53$ , $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	0.004 for -h,-k,l	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	26351	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	41.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.00% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical 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: GDL, NAG, EDO

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	0.54	1/4045 (0.0%)	0.73	6/5499 (0.1%)
1	B	0.50	0/4039	0.71	1/5491 (0.0%)
1	C	0.49	0/4047	0.68	2/5503 (0.0%)
1	D	0.46	0/4050	0.69	2/5506 (0.0%)
1	E	0.51	0/4053	0.72	3/5510 (0.1%)
1	F	0.47	0/4060	0.70	3/5520 (0.1%)
All	All	0.50	1/24294 (0.0%)	0.70	17/33029 (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

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	ASN	CG-ND2	12.73	1.64	1.32

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	142	ASN	OD1-CG-ND2	-11.23	96.07	121.90
1	E	155	GLU	C-N-CD	-7.06	105.07	120.60
1	A	142	ASN	CB-CG-ND2	6.98	133.45	116.70
1	F	316	LEU	CA-CB-CG	-6.56	100.20	115.30
1	E	490	GLY	N-CA-C	5.88	127.81	113.10
1	A	316	LEU	N-CA-C	-5.81	95.33	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	490	GLY	N-CA-C	5.63	127.17	113.10
1	F	124	VAL	N-CA-C	-5.63	95.80	111.00
1	C	123	GLN	N-CA-C	5.47	125.78	111.00
1	A	528	ARG	NE-CZ-NH2	-5.46	117.57	120.30
1	A	392	LEU	CA-CB-CG	5.32	127.54	115.30
1	B	490	GLY	N-CA-C	5.25	126.22	113.10
1	A	490	GLY	N-CA-C	5.23	126.17	113.10
1	C	490	GLY	N-CA-C	5.18	126.04	113.10
1	D	490	GLY	N-CA-C	5.11	125.87	113.10
1	D	79	ILE	N-CA-C	-5.04	97.40	111.00
1	E	535	ARG	NE-CZ-NH2	-5.01	117.80	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	142	ASN	Sidechain

## 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	3933	0	3832	91	0
1	B	3927	0	3828	70	0
1	C	3935	0	3830	72	0
1	D	3939	0	3843	86	0
1	E	3941	0	3835	79	0
1	F	3948	0	3840	75	0
2	G	28	0	25	2	0
2	H	28	0	25	4	0
2	I	28	0	25	4	0
2	J	28	0	25	2	0
2	K	28	0	25	2	0
2	L	28	0	25	2	0
3	A	15	0	8	0	0
3	B	15	0	8	0	0
3	C	15	0	8	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	15	0	8	0	0
3	E	15	0	8	0	0
3	F	15	0	8	0	0
4	A	14	0	13	1	0
4	B	14	0	13	0	0
4	C	14	0	13	0	0
4	E	14	0	13	0	0
4	F	28	0	26	0	0
5	A	12	0	18	4	0
5	B	12	0	18	4	0
5	C	12	0	18	2	0
5	E	12	0	18	1	0
5	F	12	0	18	1	0
6	A	385	0	0	17	0
6	B	399	0	0	11	0
6	C	385	0	0	11	0
6	D	346	0	0	8	0
6	E	448	0	0	14	0
6	F	363	0	0	11	0
All	All	26351	0	23374	479	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 10.

All (479) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:56:LEU:HD21	1:D:178:LEU:HD13	1.51	0.93
1:C:488:LEU:HD23	1:C:502:LEU:HD13	1.51	0.92
1:C:393:ASP:O	1:C:397:THR:HG23	1.70	0.92
1:C:488:LEU:HD12	1:C:488:LEU:O	1.74	0.87
1:F:553:HIS:ND1	1:F:554:GLU:N	2.22	0.86
1:A:63:VAL:HG22	1:A:194:ILE:HD12	1.58	0.85
1:E:315:LYS:HG2	1:E:316:LEU:H	1.40	0.84
1:D:56:LEU:HD21	1:D:178:LEU:CD1	2.06	0.84
1:E:56:LEU:HD11	1:E:178:LEU:HD13	1.60	0.84
1:D:371:MET:HE3	1:D:383:LEU:HG	1.59	0.84
1:F:303:LYS:HA	1:F:303:LYS:HE2	1.60	0.83
1:E:315:LYS:N	6:E:2243:HOH:O	2.11	0.83
1:C:125:GLN:NE2	1:C:156:PRO:HB2	1.93	0.83
1:A:310:TYR:CE1	1:A:315:LYS:HG2	2.15	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:316:LEU:HD13	1:B:318:SER:HB3	1.62	0.82
1:E:553:HIS:ND1	1:E:554:GLU:N	2.28	0.81
1:C:56:LEU:HD11	1:C:178:LEU:HD13	1.60	0.81
1:C:125:GLN:HE21	1:C:156:PRO:HB2	1.46	0.81
1:D:315:LYS:HG2	6:D:2201:HOH:O	1.81	0.79
1:E:122:THR:N	6:E:2057:HOH:O	2.15	0.78
1:E:347:GLN:HB2	6:E:2082:HOH:O	1.83	0.78
1:E:300:LYS:HA	1:E:300:LYS:HE2	1.66	0.77
6:F:2360:HOH:O	2:L:2:NAG:H82	1.85	0.77
1:C:371:MET:HE2	1:C:376:PHE:HD1	1.49	0.76
1:F:99:ARG:NH1	6:F:2044:HOH:O	2.17	0.76
1:F:315:LYS:O	1:F:316:LEU:HB2	1.86	0.75
1:B:78:TYR:CE2	1:D:262:LEU:HD11	2.22	0.74
2:G:1:NAG:H62	2:G:2:NAG:HN2	1.51	0.74
1:D:309:CYS:O	1:D:316:LEU:HB2	1.87	0.74
1:F:458:GLN:O	1:F:461:ARG:HG2	1.89	0.73
1:C:488:LEU:CD2	1:C:502:LEU:HD13	2.18	0.73
1:B:170:ARG:HD2	1:B:230:LYS:HD2	1.71	0.72
1:E:56:LEU:HD11	1:E:178:LEU:CD1	2.19	0.72
1:A:310:TYR:HE1	1:A:315:LYS:HG2	1.53	0.72
1:A:310:TYR:CD1	1:A:315:LYS:HA	2.25	0.71
1:C:91:CYS:SG	1:C:137[B]:CYS:HB2	2.31	0.71
1:A:431:GLU:HG2	1:A:435:ARG:NH1	2.05	0.70
1:A:346:ASP:O	1:A:400:LYS:HE3	1.91	0.70
1:D:299:GLY:HA3	1:D:306:LEU:HD22	1.73	0.70
1:A:91:CYS:SG	1:A:137[B]:CYS:HB2	2.31	0.69
1:F:437:THR:HB	1:F:477:GLN:HG2	1.74	0.69
1:E:99:ARG:NH2	6:E:2038:HOH:O	2.23	0.69
1:E:315:LYS:HG2	1:E:316:LEU:N	2.07	0.69
1:A:341:SER:HB3	1:A:398:ILE:HD12	1.74	0.69
1:A:56:LEU:HD11	1:A:178:LEU:HD13	1.75	0.68
1:E:299:GLY:HA3	1:E:306:LEU:HD22	1.75	0.68
1:B:315:LYS:O	1:B:316:LEU:HB2	1.93	0.68
1:C:371:MET:HE3	1:C:383:LEU:HB2	1.75	0.67
1:F:67:PRO:O	2:L:1:NAG:H61	1.94	0.67
1:E:310:TYR:CE2	1:E:315:LYS:HA	2.30	0.67
1:C:371:MET:HE2	1:C:376:PHE:CD1	2.29	0.66
1:E:478:LYS:HE2	6:E:2380:HOH:O	1.95	0.66
1:C:317:ASP:O	1:C:318:SER:HB2	1.95	0.66
1:D:379:ASP:OD2	1:D:382:LYS:HG3	1.94	0.65
1:C:102:HIS:ND1	6:C:2048:HOH:O	2.29	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:310:TYR:HD2	1:D:315:LYS:HA	1.62	0.64
1:E:91:CYS:SG	1:E:137[B]:CYS:HB2	2.37	0.64
1:B:79:ILE:H	5:B:801:EDO:C1	2.11	0.64
1:A:99:ARG:HD3	6:C:2181:HOH:O	1.98	0.64
1:C:67:PRO:O	2:I:1:NAG:H61	1.98	0.64
1:F:316:LEU:HD23	1:F:318:SER:OG	1.97	0.64
1:F:315:LYS:HD2	6:F:2111:HOH:O	1.98	0.63
1:E:478:LYS:HE3	6:E:2372:HOH:O	1.98	0.63
1:F:520:ARG:O	1:F:522:MET:HE3	1.98	0.63
1:E:529:LEU:HB3	1:E:546:LEU:HD11	1.80	0.63
1:D:529:LEU:HB3	1:D:546[A]:LEU:HD11	1.79	0.63
1:F:229:ASN:HD21	1:F:536:MET:CE	2.12	0.63
2:H:1:NAG:H61	2:H:2:NAG:HN2	1.64	0.63
1:D:221:LYS:HE2	6:D:2122:HOH:O	2.00	0.62
1:C:317:ASP:O	1:C:318:SER:CB	2.46	0.62
1:B:91:CYS:SG	1:B:137[B]:CYS:HB2	2.38	0.62
1:A:72:LEU:N	1:A:72:LEU:HD12	2.14	0.62
1:D:310:TYR:CD2	1:D:315:LYS:HA	2.34	0.62
5:C:801:EDO:H12	1:D:79:ILE:H	1.64	0.61
1:A:360[B]:CYS:HB3	6:A:2261:HOH:O	1.99	0.61
1:C:206:LEU:O	1:C:488:LEU:HG	2.00	0.61
1:C:529:LEU:HB3	1:C:546:LEU:HD11	1.83	0.61
1:A:551:CYS:HB3	1:A:553:HIS:HD2	1.66	0.61
1:D:315:LYS:N	6:D:2201:HOH:O	2.34	0.61
1:A:152:LEU:CD2	1:A:154:LYS:HE2	2.31	0.60
1:F:299:GLY:HA3	1:F:306:LEU:HD22	1.83	0.60
1:F:497:ASN:HB2	6:F:2322:HOH:O	2.01	0.60
1:F:341:SER:HB3	1:F:398:ILE:HD12	1.84	0.60
1:D:368:GLN:NE2	1:D:368:GLN:HA	2.16	0.60
1:D:65:MET:HG2	2:J:1:NAG:C7	2.31	0.60
1:D:511:GLU:OE2	1:D:528:ARG:NH2	2.34	0.60
1:D:529:LEU:HB3	1:D:546[B]:LEU:HD21	1.83	0.60
1:D:273[A]:MET:HE2	1:D:277:TYR:HB2	1.84	0.60
1:C:125:GLN:HB2	6:C:2064:HOH:O	2.01	0.60
1:A:347:GLN:HB2	6:A:2077:HOH:O	2.01	0.60
1:A:475:GLN:O	1:A:479:GLN:HG3	2.02	0.60
1:A:511:GLU:OE2	1:A:528:ARG:NH2	2.35	0.60
1:A:65:MET:HE2	2:G:1:NAG:C7	2.32	0.59
1:D:371:MET:CE	1:D:376:PHE:HD1	2.14	0.59
1:A:529:LEU:HB3	1:A:546:LEU:HD11	1.84	0.59
1:C:125:GLN:HE21	1:C:156:PRO:CB	2.13	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:303:LYS:CA	1:F:303:LYS:HE2	2.33	0.59
1:B:79:ILE:H	5:B:801:EDO:H12	1.66	0.59
1:A:66:THR:OG1	1:A:191:GLU:HB3	2.03	0.59
1:D:142:ASN:HB3	6:D:2062:HOH:O	2.02	0.59
1:A:400:LYS:HE2	6:A:2249:HOH:O	2.03	0.59
1:E:215:PRO:O	1:E:218:ILE:HG22	2.03	0.59
1:D:91:CYS:SG	1:D:137[B]:CYS:HB2	2.43	0.58
1:E:316:LEU:C	1:E:318:SER:H	2.06	0.58
1:F:215:PRO:O	1:F:218:ILE:HG22	2.03	0.58
1:A:310:TYR:C	1:A:316:LEU:HD12	2.23	0.58
1:A:125:GLN:HB2	6:A:2051:HOH:O	2.04	0.58
1:B:100:ARG:HD3	6:B:2048:HOH:O	2.03	0.58
1:A:311:SER:N	1:A:316:LEU:HD12	2.19	0.57
1:F:310:TYR:CD1	1:F:315:LYS:N	2.72	0.57
1:B:412:LYS:HA	1:B:412:LYS:HE2	1.85	0.57
1:A:529:LEU:HD13	1:A:546:LEU:HD21	1.86	0.57
1:F:63:VAL:HG22	1:F:194:ILE:HD12	1.86	0.57
1:A:122:THR:HG22	1:A:123:GLN:H	1.69	0.57
1:B:322:ILE:O	1:B:324:PRO:HD3	2.05	0.57
1:B:351:LEU:HD12	1:B:392:LEU:HD13	1.85	0.57
5:B:801:EDO:O1	1:D:262:LEU:HD13	2.04	0.57
1:E:198:PRO:HB3	1:E:514:TRP:CE3	2.40	0.57
1:B:206:LEU:C	1:B:206:LEU:HD23	2.25	0.57
1:F:527[A]:ASP:OD2	1:F:531:ARG:NH1	2.37	0.57
1:F:498:LEU:HD23	1:F:498:LEU:C	2.26	0.57
1:A:56:LEU:HD11	1:A:178:LEU:CD1	2.34	0.56
1:C:125:GLN:HG2	1:C:156:PRO:O	2.05	0.56
1:C:310:TYR:HD2	1:C:315:LYS:HB3	1.69	0.56
1:A:310:TYR:HD1	1:A:315:LYS:HA	1.70	0.56
1:B:72:LEU:HD12	1:B:72:LEU:N	2.20	0.56
1:E:370:PHE:CZ	1:E:374:LYS:HE3	2.39	0.56
2:I:1:NAG:H62	2:I:2:NAG:HN2	1.69	0.56
1:D:420:ILE:HG21	1:D:444:ILE:HD12	1.86	0.56
1:D:474:THR:O	1:D:478:LYS:HG2	2.06	0.56
1:F:444:ILE:HG23	1:F:484:GLY:HA2	1.88	0.56
1:B:529:LEU:HB3	1:B:546:LEU:HD11	1.87	0.56
1:F:399:ASN:ND2	6:F:2248:HOH:O	2.38	0.56
1:F:352:GLY:HA2	1:F:405:TRP:CD1	2.41	0.56
1:F:451:LEU:HD22	1:F:463:TYR:CE2	2.41	0.56
1:A:379:ASP:OD2	1:A:382:LYS:HG3	2.06	0.56
1:D:125:GLN:HB2	6:D:2046:HOH:O	2.06	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:315:LYS:CG	1:E:316:LEU:H	2.14	0.56
1:C:488:LEU:CD1	6:C:2134:HOH:O	2.52	0.56
1:A:299:GLY:HA3	1:A:306:LEU:HD22	1.89	0.55
1:D:307:THR:HG22	1:D:360[B]:CYS:SG	2.47	0.55
1:D:415:LEU:HB2	1:D:441:PHE:CZ	2.40	0.55
1:A:316:LEU:HD22	1:A:318:SER:HB3	1.87	0.55
1:C:488:LEU:HD23	1:C:502:LEU:CD1	2.30	0.55
1:B:125:GLN:HB3	6:B:2061:HOH:O	2.06	0.55
1:D:61:LEU:HD23	1:D:61:LEU:C	2.27	0.55
1:A:284:ARG:HD3	6:A:2202:HOH:O	2.05	0.55
1:B:310:TYR:O	1:B:311:SER:HB2	2.05	0.55
1:E:72:LEU:HG	1:E:124:VAL:HG13	1.88	0.55
1:F:238:ILE:HG23	1:F:239:VAL:HG13	1.87	0.54
1:C:371:MET:CE	1:C:376:PHE:HD1	2.20	0.54
1:F:499:THR:HB	1:F:500:PRO:HD3	1.89	0.54
1:E:70:LEU:HD22	1:E:155:GLU:HB3	1.90	0.54
1:B:273[B]:MET:HE2	1:B:277:TYR:HB2	1.90	0.54
1:D:419:THR:O	1:D:442:PRO:HD2	2.07	0.54
1:B:429:TYR:N	1:B:430:PRO:CD	2.70	0.54
1:B:461:ARG:NH2	6:B:2329:HOH:O	2.40	0.54
1:C:488:LEU:HD12	6:C:2134:HOH:O	2.07	0.54
1:E:499:THR:HB	1:E:500:PRO:HD3	1.90	0.54
1:D:72:LEU:HD12	1:D:72:LEU:N	2.24	0.53
1:E:216:VAL:O	1:E:220:LEU:HG	2.09	0.53
1:F:364:ASN:O	1:F:368:GLN:HG2	2.08	0.53
1:F:72:LEU:HD12	1:F:72:LEU:N	2.23	0.53
1:C:370:PHE:CZ	1:C:374:LYS:HE3	2.44	0.53
1:E:378:THR:HA	6:E:2297:HOH:O	2.08	0.53
1:F:471:PHE:O	1:F:478:LYS:HE2	2.09	0.53
1:B:544:GLN:HB2	1:B:545:PRO:HD2	1.91	0.53
1:B:99:ARG:HD3	6:B:2043:HOH:O	2.09	0.53
1:B:315:LYS:O	1:B:316:LEU:CB	2.56	0.52
1:C:341:SER:HB3	1:C:398:ILE:HD12	1.91	0.52
1:C:530:THR:HG22	1:C:546:LEU:HD12	1.92	0.52
1:C:529:LEU:HD13	1:C:546:LEU:HD21	1.90	0.52
1:C:311:SER:N	1:C:315:LYS:O	2.37	0.52
1:E:388:ILE:O	1:E:392:LEU:HD22	2.10	0.52
1:D:206:LEU:HD23	1:D:206:LEU:C	2.29	0.52
1:A:249:ILE:HD12	6:A:2092:HOH:O	2.09	0.52
1:D:59:LEU:HB2	1:D:528:ARG:NH2	2.25	0.52
1:B:56:LEU:HD12	1:B:56:LEU:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:497:ASN:HB2	6:B:2352:HOH:O	2.10	0.52
1:D:82:SER:HA	1:D:128:LEU:HD22	1.91	0.52
1:D:64:LYS:HD2	6:D:2011:HOH:O	2.10	0.52
1:E:65:MET:HG2	2:K:1:NAG:C7	2.40	0.51
1:A:293:GLY:O	1:A:295:THR:HG23	2.09	0.51
1:B:54:PRO:HD3	1:B:180:TYR:CD2	2.45	0.51
1:B:63:VAL:HG13	1:B:194:ILE:CD1	2.40	0.51
1:B:78:TYR:CD2	1:D:262:LEU:HD11	2.45	0.51
1:A:142:ASN:ND2	6:A:2079:HOH:O	2.40	0.51
1:E:206:LEU:HD23	1:E:206:LEU:C	2.31	0.51
1:E:300:LYS:HE2	1:E:300:LYS:CA	2.39	0.51
1:E:352:GLY:HA2	1:E:405:TRP:CD1	2.46	0.51
1:A:74:PRO:HG3	1:A:106:PHE:CD2	2.45	0.51
1:A:388:ILE:HG13	1:A:392:LEU:HD22	1.93	0.51
1:E:309:CYS:SG	1:E:360[B]:CYS:SG	3.06	0.51
1:E:437:THR:OG1	1:E:477:GLN:HG2	2.11	0.51
1:A:498:LEU:HD23	1:A:498:LEU:C	2.30	0.50
1:B:66:THR:OG1	1:B:191:GLU:HB3	2.11	0.50
1:F:221:LYS:HD2	6:F:2138:HOH:O	2.11	0.50
1:C:125:GLN:NE2	1:C:156:PRO:CB	2.71	0.50
1:C:142:ASN:HB3	6:C:2083:HOH:O	2.09	0.50
1:D:54:PRO:HD2	1:D:531:ARG:NH2	2.26	0.50
1:E:511:GLU:OE2	1:E:528:ARG:NH2	2.39	0.50
1:C:530:THR:CG2	1:C:546:LEU:HD12	2.41	0.50
1:E:529:LEU:HD13	1:E:546:LEU:HD11	1.93	0.50
1:A:431:GLU:HG2	1:A:435:ARG:HH11	1.77	0.50
1:E:522:MET:HE3	6:E:2409:HOH:O	2.11	0.50
1:A:370:PHE:CZ	1:A:374:LYS:HE3	2.46	0.50
1:F:309:CYS:O	1:F:316:LEU:HB3	2.11	0.50
2:I:1:NAG:H62	2:I:2:NAG:C1	2.42	0.50
1:C:504:PRO:O	1:C:507:SER:HB2	2.12	0.50
1:C:322:ILE:O	1:C:324:PRO:HD3	2.12	0.50
6:B:2014:HOH:O	1:D:262:LEU:HD12	2.10	0.50
1:A:141:PRO:HB2	6:A:2199:HOH:O	2.11	0.50
1:A:83:PRO:HD2	4:A:702:NAG:H82	1.93	0.50
1:C:70:LEU:HD22	1:C:155:GLU:HB3	1.93	0.50
1:D:326:LEU:O	1:D:329:THR:HB	2.12	0.50
1:D:370:PHE:CZ	1:D:374:LYS:HE3	2.46	0.50
1:D:70:LEU:HB3	1:D:124:VAL:HG23	1.94	0.49
1:B:58:PRO:HD2	1:B:230:LYS:HG3	1.94	0.49
1:D:352:GLY:HA2	1:D:405:TRP:CD1	2.47	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:70:LEU:HD22	1:A:155:GLU:HB3	1.94	0.49
1:F:293:GLY:O	1:F:295:THR:HG23	2.11	0.49
1:B:529:LEU:HD13	1:B:546:LEU:HD21	1.95	0.49
1:E:327:ASN:HD22	1:E:328:THR:N	2.10	0.49
1:A:352:GLY:HA2	1:A:405:TRP:CD1	2.47	0.49
1:A:535:ARG:HA	1:A:538:GLU:HG2	1.95	0.49
1:D:322:ILE:O	1:D:324:PRO:HD3	2.13	0.49
1:D:429:TYR:N	1:D:430:PRO:CD	2.75	0.49
1:D:198:PRO:HB3	1:D:514:TRP:CE3	2.48	0.49
1:A:426:ASP:HB2	5:A:800:EDO:H12	1.94	0.49
1:A:429:TYR:N	1:A:430:PRO:CD	2.76	0.49
1:B:351:LEU:CD1	1:B:392:LEU:HD13	2.43	0.49
1:F:397:THR:HB	6:F:2245:HOH:O	2.11	0.49
1:A:412:LYS:HE2	1:A:412:LYS:HA	1.94	0.48
1:B:261:SER:OG	1:B:264:HIS:ND1	2.39	0.48
1:C:288:GLU:HG3	1:C:350:HIS:HB3	1.95	0.48
1:E:262:LEU:HA	6:E:2202:HOH:O	2.13	0.48
1:F:426:ASP:HB2	5:F:800:EDO:H12	1.95	0.48
1:A:152:LEU:HG	1:A:154:LYS:HG2	1.95	0.48
1:A:322:ILE:O	1:A:324:PRO:HD3	2.14	0.48
1:C:56:LEU:HD11	1:C:178:LEU:CD1	2.39	0.48
1:D:253:GLU:HA	1:D:256:ASN:HB2	1.96	0.48
1:E:149:TYR:CZ	1:E:196:ASP:HB3	2.48	0.48
1:E:471:PHE:O	1:E:478:LYS:NZ	2.45	0.48
1:F:198:PRO:HB3	1:F:514:TRP:CE3	2.47	0.48
1:C:488:LEU:CD2	1:C:502:LEU:HD22	2.43	0.48
1:D:284:ARG:HD3	6:D:2181:HOH:O	2.13	0.48
1:A:311:SER:HB2	6:A:2228:HOH:O	2.13	0.48
1:B:64:LYS:HD2	6:B:2010:HOH:O	2.14	0.48
1:C:91:CYS:SG	1:C:137[B]:CYS:CB	2.99	0.48
1:B:289:PHE:CE2	1:B:340:ILE:HD12	2.49	0.48
1:B:393:ASP:O	1:B:397:THR:HG23	2.14	0.48
1:D:514:TRP:CD1	1:D:514:TRP:C	2.87	0.48
1:E:310:TYR:CD2	1:E:315:LYS:HA	2.48	0.48
1:F:310:TYR:HD1	1:F:315:LYS:N	2.10	0.48
1:A:142:ASN:OD1	1:A:144:SER:HB2	2.13	0.48
1:F:322:ILE:O	1:F:324:PRO:HD3	2.13	0.48
1:D:498:LEU:C	1:D:498:LEU:HD23	2.34	0.48
1:E:73:ALA:HB3	1:E:76:ASN:HB3	1.95	0.48
1:E:341:SER:HB3	1:E:398:ILE:HD12	1.95	0.47
2:H:1:NAG:H61	2:H:2:NAG:N2	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:498:LEU:C	1:C:498:LEU:HD23	2.34	0.47
1:F:429:TYR:N	1:F:430:PRO:CD	2.77	0.47
1:A:253:GLU:HA	1:A:256:ASN:HB2	1.97	0.47
1:A:368:GLN:HG3	6:A:2265:HOH:O	2.14	0.47
1:B:65:MET:HG2	2:H:1:NAG:C7	2.44	0.47
1:D:544:GLN:HB2	1:D:545:PRO:HD2	1.96	0.47
1:B:91:CYS:SG	1:B:137[B]:CYS:CB	3.00	0.47
1:C:526:TYR:O	1:C:530:THR:HG23	2.14	0.47
1:D:290:ASP:OD2	1:D:354:ASP:OD1	2.32	0.47
1:F:140:PHE:CE2	1:F:279:ARG:HD3	2.50	0.47
1:E:553:HIS:CG	1:E:554:GLU:N	2.82	0.47
1:B:317:ASP:N	1:B:317:ASP:OD1	2.45	0.47
1:B:419:THR:O	1:B:442:PRO:HD2	2.15	0.47
1:C:135:SER:O	1:C:136:GLU:HB2	2.15	0.47
1:D:315:LYS:HB2	1:D:316:LEU:H	1.25	0.47
1:A:63:VAL:HG13	1:A:194:ILE:HD11	1.97	0.47
1:C:284:ARG:HD3	6:C:2199:HOH:O	2.13	0.47
1:D:315:LYS:O	1:D:316:LEU:HG	2.15	0.47
1:E:91:CYS:SG	1:E:137[B]:CYS:CB	3.02	0.47
1:C:80:SER:HB3	6:C:2018:HOH:O	2.15	0.46
1:A:229:ASN:HD21	1:A:536:MET:CE	2.28	0.46
1:A:325:THR:HG22	1:A:370:PHE:CD2	2.51	0.46
1:A:415:LEU:HB2	1:A:441:PHE:CZ	2.50	0.46
1:E:429:TYR:N	1:E:430:PRO:CD	2.77	0.46
1:A:419:THR:O	1:A:442:PRO:HD2	2.16	0.46
1:E:351:LEU:HD12	1:E:392:LEU:HD13	1.96	0.46
1:A:63:VAL:HG22	1:A:194:ILE:CD1	2.39	0.46
1:F:221:LYS:HE3	1:F:540:GLY:O	2.16	0.46
1:D:182:ASP:C	1:D:182:ASP:OD1	2.54	0.46
1:E:379:ASP:HB3	1:E:382:LYS:HD2	1.98	0.46
1:A:99:ARG:HB2	6:A:2036:HOH:O	2.16	0.46
1:A:99:ARG:HD3	6:A:2036:HOH:O	2.15	0.46
1:D:511:GLU:CD	1:D:528:ARG:HH22	2.18	0.46
1:C:206:LEU:HD23	1:C:206:LEU:C	2.36	0.45
1:F:514:TRP:CD1	1:F:514:TRP:C	2.89	0.45
1:A:140:PHE:CE2	1:A:279:ARG:HD3	2.50	0.45
1:B:54:PRO:HD3	1:B:180:TYR:HB3	1.98	0.45
1:C:72:LEU:HD12	1:C:72:LEU:N	2.31	0.45
1:D:61:LEU:HD23	1:D:61:LEU:O	2.16	0.45
1:C:415:LEU:HB2	1:C:441:PHE:CZ	2.51	0.45
1:D:199:ARG:CZ	1:D:517:LYS:HB2	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:451:LEU:HD22	1:E:463:TYR:CE2	2.51	0.45
1:F:364:ASN:HA	1:F:365:PRO:HD3	1.80	0.45
1:F:553:HIS:CG	1:F:554:GLU:N	2.84	0.45
1:B:79:ILE:H	5:B:801:EDO:H11	1.78	0.45
1:C:368:GLN:HA	1:C:368:GLN:OE1	2.15	0.45
1:C:544:GLN:HB2	1:C:545:PRO:HD2	1.97	0.45
1:D:368:GLN:HE21	1:D:368:GLN:HA	1.82	0.45
1:D:65:MET:HG2	2:J:1:NAG:O7	2.16	0.45
1:A:63:VAL:HG13	1:A:194:ILE:CD1	2.47	0.45
1:B:316:LEU:HD13	1:B:318:SER:CB	2.41	0.45
1:B:65:MET:HG2	2:H:1:NAG:O7	2.17	0.45
1:A:206:LEU:HD12	1:A:235:HIS:CE1	2.52	0.45
1:D:403:ILE:HG12	1:D:420:ILE:HB	1.99	0.45
1:E:351:LEU:CD1	1:E:392:LEU:HD13	2.47	0.45
1:C:161:LYS:HB2	1:C:161:LYS:HE3	1.88	0.45
1:D:460:TRP:CD1	1:D:546[A]:LEU:HD22	2.51	0.45
1:F:135:SER:O	1:F:136:GLU:HB2	2.16	0.45
1:B:412:LYS:HE2	1:B:412:LYS:CA	2.46	0.45
1:C:352:GLY:HA2	1:C:405:TRP:CD1	2.52	0.45
1:C:429:TYR:N	1:C:430:PRO:CD	2.80	0.45
1:E:149:TYR:CE2	1:E:196:ASP:HB3	2.51	0.45
1:E:544:GLN:HB2	1:E:545:PRO:HD2	1.99	0.45
2:K:1:NAG:H62	2:K:2:NAG:C1	2.47	0.45
1:B:420:ILE:HG21	1:B:444:ILE:HD12	1.99	0.45
1:C:497:ASN:HB2	6:C:2341:HOH:O	2.17	0.45
1:F:360[B]:CYS:HB3	6:F:2233:HOH:O	2.17	0.45
1:A:152:LEU:HD23	1:A:154:LYS:HE2	1.97	0.45
1:D:93:LEU:C	1:D:93:LEU:HD23	2.38	0.45
1:E:78:TYR:HA	5:E:801:EDO:O1	2.17	0.44
1:C:370:PHE:CE2	1:C:374:LYS:HG3	2.52	0.44
1:E:178:LEU:HG	6:E:2127:HOH:O	2.17	0.44
1:E:393:ASP:O	1:E:397:THR:HG23	2.17	0.44
1:F:70:LEU:HD22	1:F:155:GLU:HB3	1.99	0.44
1:A:188:THR:HG22	1:A:189:ILE:N	2.32	0.44
1:B:420:ILE:HG12	1:B:442:PRO:HB2	1.99	0.44
1:D:462:LYS:HD2	6:D:2274:HOH:O	2.17	0.44
1:E:82:SER:HA	1:E:128:LEU:HD22	2.00	0.44
1:A:326:LEU:O	1:A:329:THR:HB	2.17	0.44
1:B:239:VAL:HG12	1:B:245:PRO:HD2	1.98	0.44
1:A:272:ARG:NH2	6:A:2191:HOH:O	2.50	0.44
1:A:471:PHE:O	1:A:478:LYS:HE2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:315:LYS:NZ	6:B:2219:HOH:O	2.51	0.44
1:C:198:PRO:HB3	1:C:514:TRP:CE3	2.52	0.44
1:F:339:GLU:O	1:F:343:VAL:HG23	2.18	0.44
1:A:310:TYR:CE1	1:A:315:LYS:HA	2.52	0.44
1:F:149:TYR:CE2	1:F:196:ASP:HB3	2.53	0.44
1:B:518[A]:ASP:N	1:B:518[A]:ASP:OD1	2.50	0.44
1:C:428:ALA:HB1	1:C:431:GLU:OE1	2.18	0.44
1:E:262:LEU:HD12	6:E:2202:HOH:O	2.16	0.44
1:D:371:MET:HE2	1:D:376:PHE:HD1	1.83	0.43
1:F:247:GLN:OE1	1:F:267:THR:HG22	2.18	0.43
1:B:498:LEU:HD23	1:B:498:LEU:C	2.38	0.43
1:E:93:LEU:HD23	1:E:93:LEU:C	2.39	0.43
1:F:63:VAL:HG13	1:F:194:ILE:CD1	2.48	0.43
1:F:209:THR:HG22	1:F:214:LEU:HD12	1.99	0.43
1:B:352:GLY:HA2	1:B:405:TRP:CD1	2.54	0.43
1:B:403:ILE:HG12	1:B:420:ILE:HB	2.00	0.43
1:D:351:LEU:HD12	1:D:392:LEU:HD13	2.01	0.43
1:F:229:ASN:HD21	1:F:536:MET:HE3	1.79	0.43
1:A:393:ASP:O	1:A:397:THR:HG23	2.18	0.43
1:A:544:GLN:HB2	1:A:545:PRO:HD2	2.00	0.43
1:E:72:LEU:N	1:E:72:LEU:HD12	2.33	0.43
1:A:284:ARG:HG3	6:A:2199:HOH:O	2.18	0.43
1:F:232:ASN:OD1	1:F:233:VAL:HG23	2.18	0.43
1:C:529:LEU:HD13	1:C:546:LEU:HD11	1.99	0.43
1:F:188:THR:HG22	1:F:189:ILE:N	2.33	0.43
1:C:497:ASN:O	1:C:500:PRO:HD2	2.18	0.43
5:A:800:EDO:H22	6:A:2303:HOH:O	2.18	0.43
1:D:91:CYS:SG	1:D:137[B]:CYS:CB	3.04	0.43
1:D:188:THR:HG22	1:D:189:ILE:N	2.33	0.43
1:E:303:LYS:HA	1:E:303:LYS:HD3	1.75	0.43
1:F:72:LEU:HG	1:F:124:VAL:HG13	2.00	0.43
1:A:206:LEU:HD23	1:A:206:LEU:C	2.39	0.43
1:B:73:ALA:HB3	1:B:76:ASN:HB3	2.01	0.43
1:D:57:TRP:CH2	1:D:529:LEU:HG	2.54	0.43
1:F:149:TYR:CZ	1:F:196:ASP:HB3	2.54	0.43
1:B:504:PRO:O	1:B:507:SER:HB2	2.19	0.43
1:D:371:MET:HE2	1:D:376:PHE:CD1	2.54	0.43
1:D:341:SER:HB3	1:D:398:ILE:HD12	2.01	0.43
1:D:437:THR:HB	1:D:477:GLN:HG2	2.01	0.42
1:E:514:TRP:CD1	1:E:514:TRP:C	2.92	0.42
1:F:206:LEU:C	1:F:206:LEU:HD23	2.39	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:325:THR:HB	1:B:370:PHE:CD1	2.54	0.42
1:C:431:GLU:HB3	1:C:435:ARG:HH12	1.84	0.42
1:F:85:SER:HB2	1:F:130:SER:HA	2.01	0.42
1:D:384:GLU:OE1	1:D:411:ASP:OD2	2.37	0.42
1:F:460:TRP:CD1	1:F:546:LEU:HD22	2.54	0.42
1:B:419:THR:HA	6:B:2293:HOH:O	2.18	0.42
1:D:364:ASN:HA	1:D:365:PRO:HD3	1.93	0.42
1:D:383:LEU:HA	1:D:383:LEU:HD23	1.83	0.42
1:A:72:LEU:HG	1:A:124:VAL:HG13	2.02	0.42
1:B:140:PHE:CE2	1:B:345:PRO:HG3	2.55	0.42
1:E:420:ILE:HA	1:E:442:PRO:HG2	2.01	0.42
1:E:529:LEU:HD13	1:E:546:LEU:HD21	2.02	0.42
1:E:551:CYS:O	1:E:553:HIS:N	2.41	0.42
1:F:253:GLU:HA	1:F:256:ASN:HB2	2.01	0.42
2:I:1:NAG:C6	2:I:2:NAG:HN2	2.32	0.42
1:B:135:SER:O	1:B:136:GLU:HB2	2.19	0.42
1:C:253:GLU:HA	1:C:256:ASN:HB2	2.02	0.42
1:E:142:ASN:HB2	6:E:2080:HOH:O	2.20	0.42
1:E:415:LEU:HB2	1:E:441:PHE:CZ	2.54	0.42
1:F:302:GLN:HG3	1:F:305:LEU:HB2	2.01	0.42
1:C:149:TYR:CE2	1:C:196:ASP:HB3	2.54	0.42
1:A:529:LEU:HD13	1:A:546:LEU:HD11	2.02	0.42
1:A:79:ILE:H	5:A:801:EDO:C1	2.33	0.42
1:B:149:TYR:CZ	1:B:196:ASP:HB3	2.54	0.42
1:D:149:TYR:CZ	1:D:196:ASP:HB3	2.55	0.42
1:D:371:MET:CE	1:D:376:PHE:CD1	3.01	0.42
1:D:511:GLU:O	1:D:515:SER:HB2	2.20	0.42
1:E:547:TYR:CG	1:E:548:ALA:N	2.88	0.42
1:A:269:ASN:HB2	6:C:2038:HOH:O	2.20	0.41
1:A:305:LEU:HG	1:A:322:ILE:HG12	2.02	0.41
1:B:99:ARG:HG3	6:B:2045:HOH:O	2.18	0.41
1:E:413:ALA:HB2	6:E:2302:HOH:O	2.20	0.41
1:F:150[A]:THR:HG23	6:F:2093:HOH:O	2.20	0.41
1:A:514:TRP:CD1	1:A:514:TRP:C	2.93	0.41
1:B:514:TRP:CD1	1:B:514:TRP:C	2.93	0.41
1:C:476:LYS:O	1:C:479:GLN:HB2	2.20	0.41
1:E:103:GLY:O	1:E:107:GLY:N	2.50	0.41
1:E:369:ASP:O	1:E:373:GLN:HG3	2.20	0.41
1:F:384:GLU:OE1	1:F:411:ASP:OD2	2.37	0.41
1:F:461:ARG:NH2	6:F:2296:HOH:O	2.47	0.41
1:F:527[B]:ASP:O	1:F:531:ARG:HG3	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:TYR:CE2	1:A:196:ASP:HB3	2.55	0.41
1:A:221:LYS:NZ	6:A:2145:HOH:O	2.52	0.41
1:B:475:GLN:O	1:B:479:GLN:HG3	2.20	0.41
1:F:544:GLN:HB2	1:F:545:PRO:HD2	2.01	0.41
1:E:135:SER:O	1:E:136:GLU:HB2	2.21	0.41
1:C:427:SER:HB2	6:C:2289:HOH:O	2.19	0.41
1:D:238:ILE:HG23	1:D:239:VAL:HG13	2.02	0.41
1:E:58:PRO:HB3	1:E:511:GLU:HA	2.01	0.41
1:F:235:HIS:ND1	1:F:485:GLU:OE1	2.53	0.41
1:F:309:CYS:O	1:F:316:LEU:CB	2.68	0.41
1:C:76:ASN:ND2	1:C:78:TYR:HE1	2.19	0.41
1:D:307:THR:CG2	1:D:360[B]:CYS:SG	3.07	0.41
1:B:288:GLU:HG3	1:B:350:HIS:HB3	2.03	0.41
1:C:290:ASP:OD2	1:C:354:ASP:OD1	2.38	0.41
1:A:79:ILE:HB	5:A:801:EDO:H12	2.02	0.41
1:E:497:ASN:HB2	6:E:2398:HOH:O	2.20	0.41
1:F:310:TYR:CE1	1:F:315:LYS:N	2.89	0.41
1:D:293:GLY:O	1:D:295:THR:HG23	2.21	0.41
1:E:384:GLU:OE1	1:E:411:ASP:OD2	2.39	0.41
1:B:217:LYS:HG3	6:B:2134:HOH:O	2.21	0.41
1:B:93:LEU:C	1:B:93:LEU:HD23	2.41	0.41
1:D:371:MET:HE2	1:D:376:PHE:HB2	2.02	0.41
1:E:403:ILE:HG12	1:E:420:ILE:HB	2.02	0.41
1:F:437:THR:CB	1:F:477:GLN:HG2	2.48	0.41
1:D:149:TYR:CE2	1:D:196:ASP:HB3	2.55	0.41
1:F:461:ARG:NH2	6:F:2295:HOH:O	2.49	0.41
1:B:369:ASP:O	1:B:372:ARG:HB3	2.21	0.40
1:C:140:PHE:CE2	1:C:279:ARG:HD3	2.56	0.40
1:E:322:ILE:O	1:E:324:PRO:HD3	2.20	0.40
1:F:133:LEU:O	1:F:134:GLN:C	2.60	0.40
1:F:420:ILE:HG12	1:F:442:PRO:HB2	2.04	0.40
1:F:468:PRO:HB3	1:F:481:PHE:CZ	2.56	0.40
1:A:199:ARG:CZ	1:A:517:LYS:HB2	2.51	0.40
1:A:221:LYS:HE2	6:A:2147:HOH:O	2.21	0.40
1:B:198:PRO:HB3	1:B:514:TRP:CE3	2.56	0.40
5:C:801:EDO:H12	1:D:79:ILE:N	2.35	0.40
1:E:309:CYS:SG	1:E:360[B]:CYS:HB2	2.62	0.40
1:A:474:THR:O	1:A:478:LYS:HG2	2.21	0.40
1:C:93:LEU:C	1:C:93:LEU:HD23	2.41	0.40
1:D:221:LYS:NZ	1:D:540:GLY:O	2.43	0.40
1:E:316:LEU:C	1:E:318:SER:N	2.73	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:82:SER:HA	1:B:128:LEU:HD22	2.02	0.40
1:B:54:PRO:HD3	1:B:180:TYR:CG	2.56	0.40
1:C:326:LEU:O	1:C:329:THR:HB	2.21	0.40
1:C:488:LEU:O	1:C:488:LEU:CD1	2.58	0.40
1:B:78:TYR:HE2	1:D:262:LEU:HD11	1.80	0.40
1:D:530:THR:CG2	1:D:546[A]:LEU:HD12	2.51	0.40
1:F:101:TYR:CE2	1:F:173:GLU:HA	2.57	0.40
1:A:149:TYR:CZ	1:A:196:ASP:HB3	2.56	0.40
1:A:215:PRO:O	1:A:218:ILE:HG22	2.21	0.40
1:A:338:LYS:HD2	1:A:338:LYS:HA	1.87	0.40
1:A:431:GLU:CG	1:A:435:ARG:NH1	2.79	0.40
1:C:149:TYR:CZ	1:C:196:ASP:HB3	2.55	0.40
1:C:474:THR:O	1:C:478:LYS:HG2	2.21	0.40
1:E:133:LEU:O	1:E:134:GLN:C	2.60	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	482/515 (94%)	466 (97%)	16 (3%)	0	100	100
1	B	481/515 (93%)	462 (96%)	18 (4%)	1 (0%)	47	55
1	C	482/515 (94%)	467 (97%)	14 (3%)	1 (0%)	47	55
1	D	483/515 (94%)	462 (96%)	20 (4%)	1 (0%)	47	55
1	E	483/515 (94%)	463 (96%)	18 (4%)	2 (0%)	34	37
1	F	484/515 (94%)	463 (96%)	19 (4%)	2 (0%)	34	37
All	All	2895/3090 (94%)	2783 (96%)	105 (4%)	7 (0%)	47	55

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	316	LEU
1	C	318	SER
1	D	316	LEU
1	E	552	ASN
1	F	316	LEU
1	E	316	LEU
1	F	553	HIS

### 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	433/454 (95%)	418 (96%)	15 (4%)	36	43
1	B	432/454 (95%)	422 (98%)	10 (2%)	50	59
1	C	433/454 (95%)	420 (97%)	13 (3%)	41	50
1	D	434/454 (96%)	422 (97%)	12 (3%)	43	52
1	E	434/454 (96%)	421 (97%)	13 (3%)	41	50
1	F	435/454 (96%)	423 (97%)	12 (3%)	43	52
All	All	2601/2724 (96%)	2526 (97%)	75 (3%)	42	51

All (75) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	56	LEU
1	A	72	LEU
1	A	142	ASN
1	A	151	LEU
1	A	154	LYS
1	A	178	LEU
1	A	211	ARG
1	A	221	LYS
1	A	269	ASN
1	A	296	LEU
1	A	316	LEU
1	A	383	LEU

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Mol	Chain	Res	Type
1	A	392	LEU
1	A	504	PRO
1	A	529	LEU
1	B	61	LEU
1	B	151	LEU
1	B	315	LYS
1	B	316	LEU
1	B	317	ASP
1	B	357	GLU
1	B	383	LEU
1	B	392	LEU
1	B	504	PRO
1	B	529	LEU
1	C	61	LEU
1	C	80	SER
1	C	99	ARG
1	C	124	VAL
1	C	151	LEU
1	C	155	GLU
1	C	178	LEU
1	C	315	LYS
1	C	373	GLN
1	C	383	LEU
1	C	388	ILE
1	C	392	LEU
1	C	529	LEU
1	D	82	SER
1	D	151	LEU
1	D	178	LEU
1	D	211	ARG
1	D	306	LEU
1	D	383	LEU
1	D	388	ILE
1	D	392	LEU
1	D	426	ASP
1	D	462	LYS
1	D	504	PRO
1	D	529	LEU
1	E	56	LEU
1	E	124	VAL
1	E	151	LEU
1	E	178	LEU

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Mol	Chain	Res	Type
1	E	211	ARG
1	E	218	ILE
1	E	306	LEU
1	E	316	LEU
1	E	327	ASN
1	E	383	LEU
1	E	392	LEU
1	E	529	LEU
1	E	530	THR
1	F	122	THR
1	F	124	VAL
1	F	151	LEU
1	F	211	ARG
1	F	218	ILE
1	F	306	LEU
1	F	315	LYS
1	F	316	LEU
1	F	383	LEU
1	F	392	LEU
1	F	461	ARG
1	F	553	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	76	ASN
1	A	553	HIS
1	B	76	ASN
1	B	134	GLN
1	B	373	GLN
1	C	71	HIS
1	C	76	ASN
1	C	125	GLN
1	D	68	ASN
1	D	76	ASN
1	D	125	GLN
1	D	126	GLN
1	D	368	GLN
1	D	373	GLN
1	D	479	GLN
1	E	71	HIS
1	E	76	ASN

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Mol	Chain	Res	Type
1	E	327	ASN
1	F	102	HIS
1	F	125	GLN
1	F	142	ASN

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates ⓘ

12 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	G	1	1,2	14,14,15	0.51	0	17,19,21	0.92	0
2	NAG	G	2	2	14,14,15	0.54	0	17,19,21	0.71	0
2	NAG	H	1	1,2	14,14,15	0.42	0	17,19,21	0.68	0
2	NAG	H	2	2	14,14,15	0.52	0	17,19,21	0.66	0
2	NAG	I	1	1,2	14,14,15	0.43	0	17,19,21	0.88	1 (5%)
2	NAG	I	2	2	14,14,15	0.56	0	17,19,21	0.67	0
2	NAG	J	1	1,2	14,14,15	0.49	0	17,19,21	0.92	1 (5%)
2	NAG	J	2	2	14,14,15	0.54	0	17,19,21	0.78	1 (5%)
2	NAG	K	1	1,2	14,14,15	0.47	0	17,19,21	1.11	2 (11%)
2	NAG	K	2	2	14,14,15	0.52	0	17,19,21	0.67	0
2	NAG	L	1	1,2	14,14,15	0.52	0	17,19,21	0.80	1 (5%)
2	NAG	L	2	2	14,14,15	0.63	0	17,19,21	0.62	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
2	NAG	G	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	G	2	2	-	6/6/23/26	0/1/1/1
2	NAG	H	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	H	2	2	-	4/6/23/26	0/1/1/1
2	NAG	I	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	I	2	2	-	4/6/23/26	0/1/1/1
2	NAG	J	1	1,2	-	3/6/23/26	0/1/1/1
2	NAG	J	2	2	-	4/6/23/26	0/1/1/1
2	NAG	K	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	K	2	2	-	5/6/23/26	0/1/1/1
2	NAG	L	1	1,2	-	4/6/23/26	0/1/1/1
2	NAG	L	2	2	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	K	1	NAG	C3-C4-C5	2.60	114.88	110.24
2	J	1	NAG	C2-N2-C7	-2.58	119.23	122.90
2	L	1	NAG	C2-N2-C7	-2.23	119.72	122.90
2	J	2	NAG	C2-N2-C7	-2.20	119.78	122.90
2	K	1	NAG	C2-N2-C7	-2.10	119.91	122.90
2	I	1	NAG	C2-N2-C7	-2.02	120.03	122.90

There are no chirality outliers.

All (44) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	G	2	NAG	C8-C7-N2-C2
2	G	2	NAG	O7-C7-N2-C2
2	H	2	NAG	C8-C7-N2-C2
2	H	2	NAG	O7-C7-N2-C2
2	K	2	NAG	C8-C7-N2-C2
2	K	2	NAG	O7-C7-N2-C2
2	I	2	NAG	C8-C7-N2-C2

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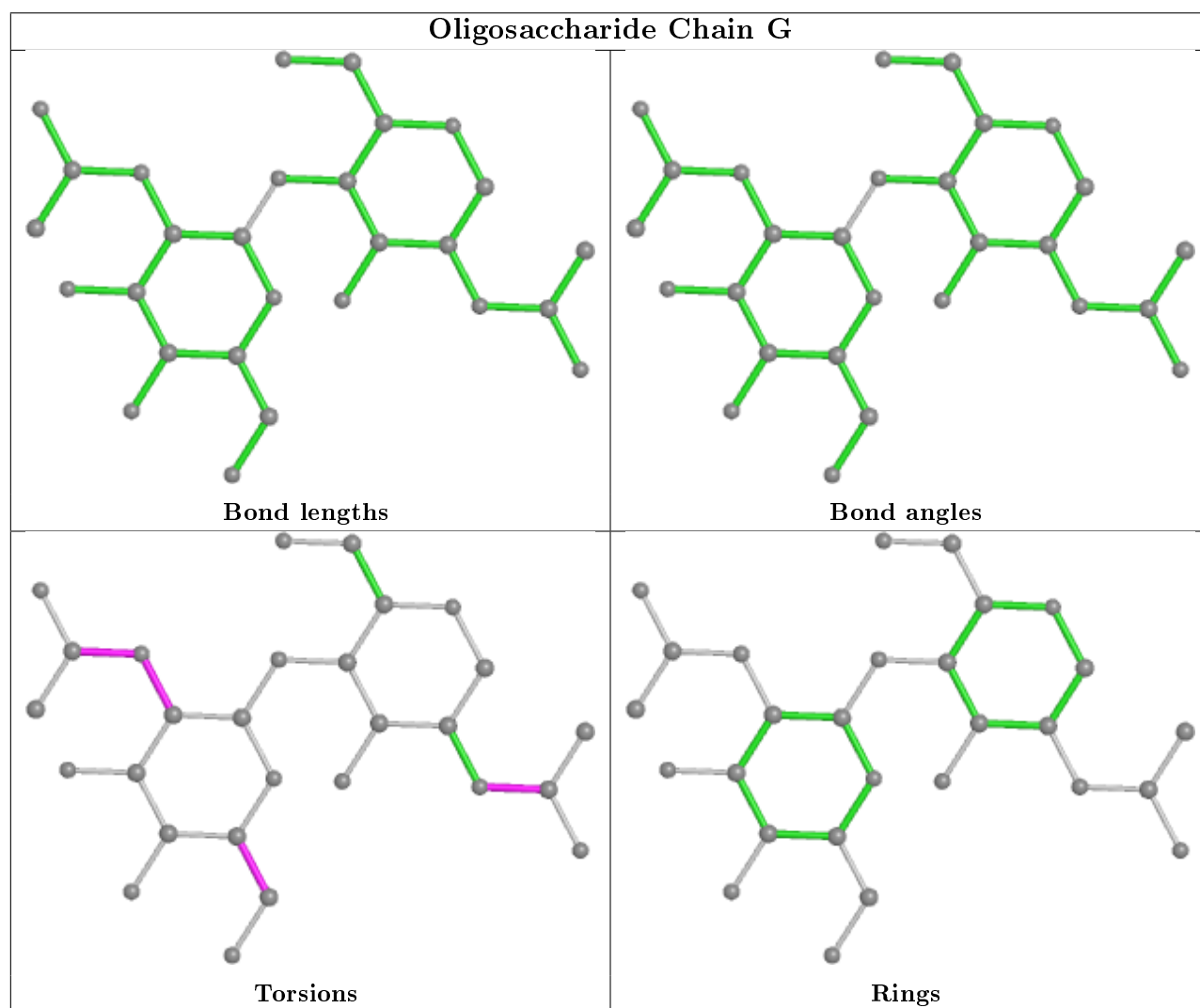
Mol	Chain	Res	Type	Atoms
2	I	2	NAG	O7-C7-N2-C2
2	L	2	NAG	C8-C7-N2-C2
2	L	2	NAG	O7-C7-N2-C2
2	J	2	NAG	C8-C7-N2-C2
2	J	2	NAG	O7-C7-N2-C2
2	G	1	NAG	C8-C7-N2-C2
2	G	1	NAG	O7-C7-N2-C2
2	L	1	NAG	O5-C5-C6-O6
2	I	2	NAG	O5-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
2	L	1	NAG	C4-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6
2	G	2	NAG	C1-C2-N2-C7
2	H	2	NAG	C4-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	I	2	NAG	C4-C5-C6-O6
2	J	1	NAG	C8-C7-N2-C2
2	K	1	NAG	C8-C7-N2-C2
2	K	2	NAG	C4-C5-C6-O6
2	J	2	NAG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
2	J	2	NAG	C4-C5-C6-O6
2	J	1	NAG	O7-C7-N2-C2
2	K	1	NAG	O7-C7-N2-C2
2	L	2	NAG	C1-C2-N2-C7
2	I	1	NAG	C8-C7-N2-C2
2	I	1	NAG	O5-C5-C6-O6
2	L	1	NAG	C8-C7-N2-C2
2	H	1	NAG	C8-C7-N2-C2
2	I	1	NAG	O7-C7-N2-C2
2	H	1	NAG	O7-C7-N2-C2
2	L	1	NAG	O7-C7-N2-C2
2	G	2	NAG	C3-C2-N2-C7
2	L	2	NAG	C3-C2-N2-C7
2	J	1	NAG	C4-C5-C6-O6
2	K	2	NAG	C1-C2-N2-C7

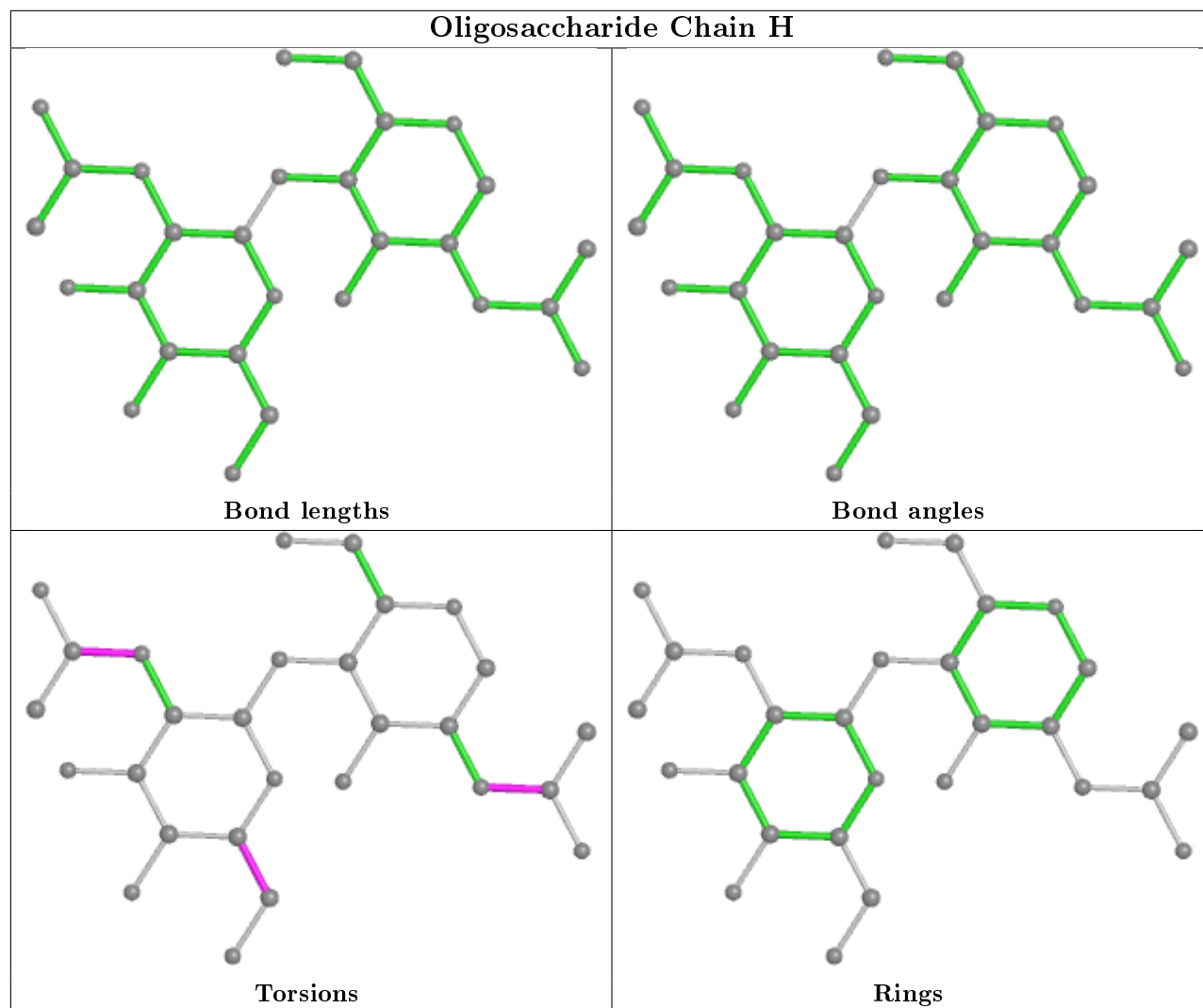
There are no ring outliers.

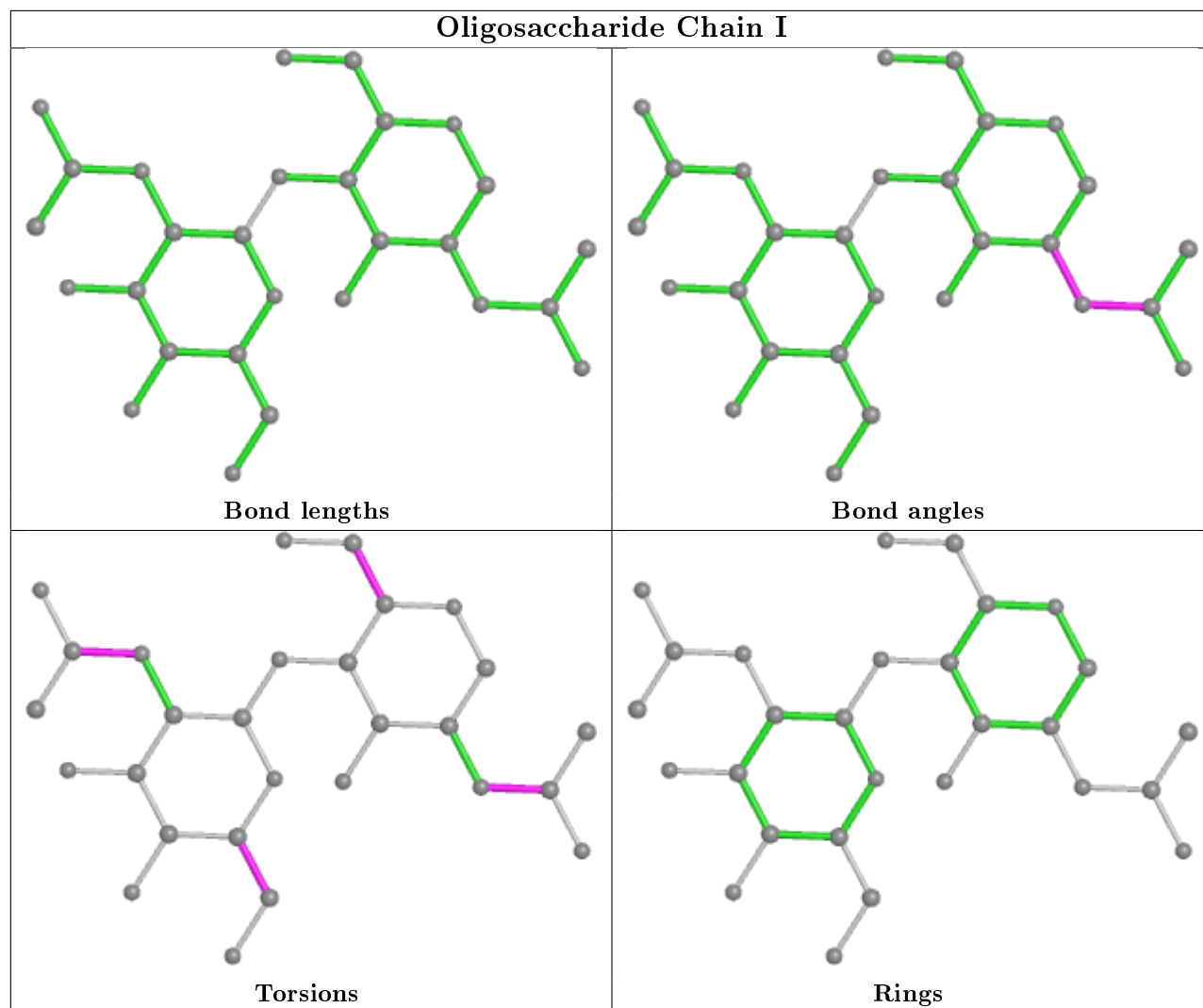
11 monomers are involved in 16 short contacts:

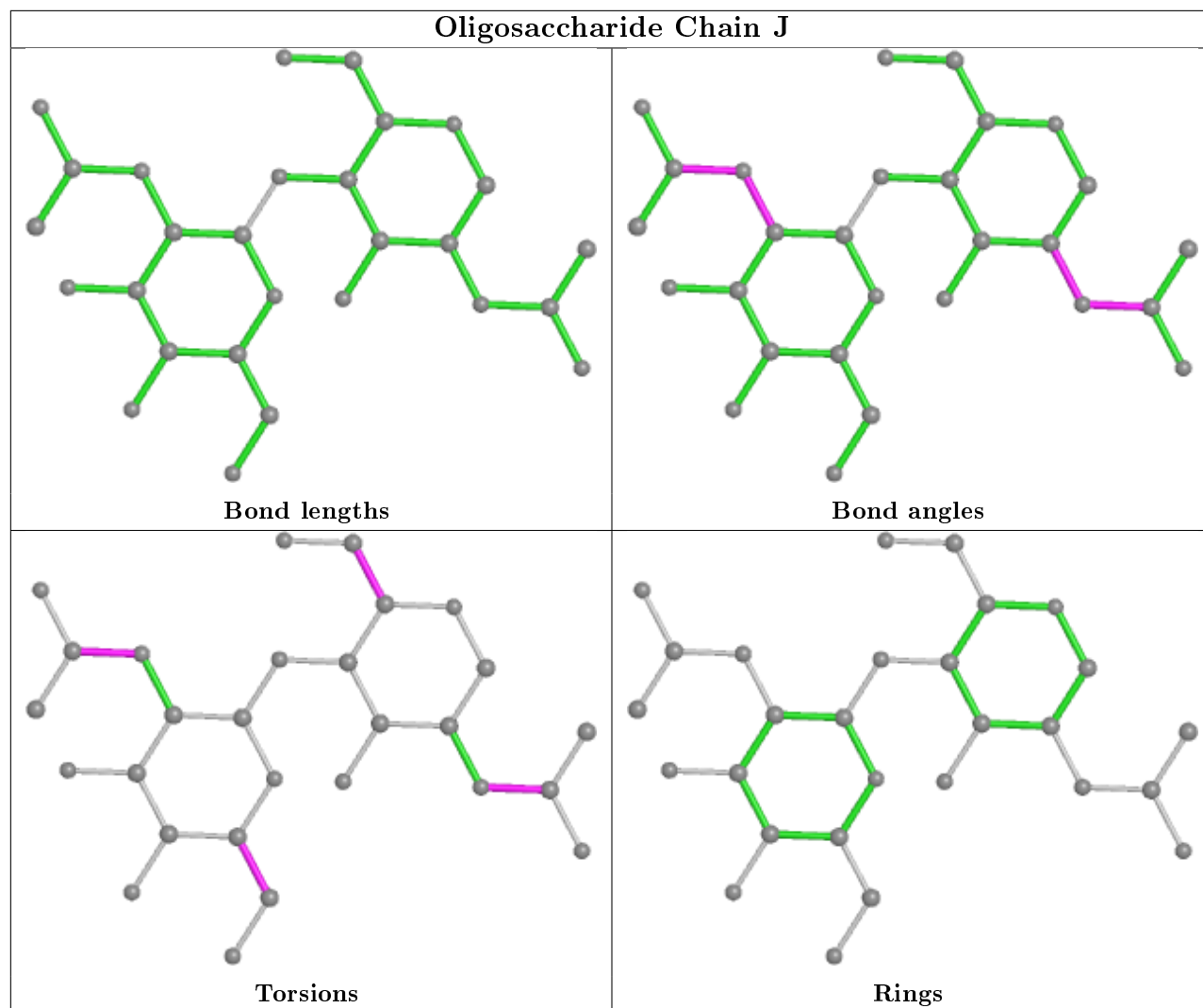
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	J	1	NAG	2	0
2	I	1	NAG	4	0
2	G	2	NAG	1	0
2	H	2	NAG	2	0
2	K	2	NAG	1	0
2	I	2	NAG	3	0
2	K	1	NAG	2	0
2	L	1	NAG	1	0
2	L	2	NAG	1	0
2	G	1	NAG	2	0
2	H	1	NAG	4	0

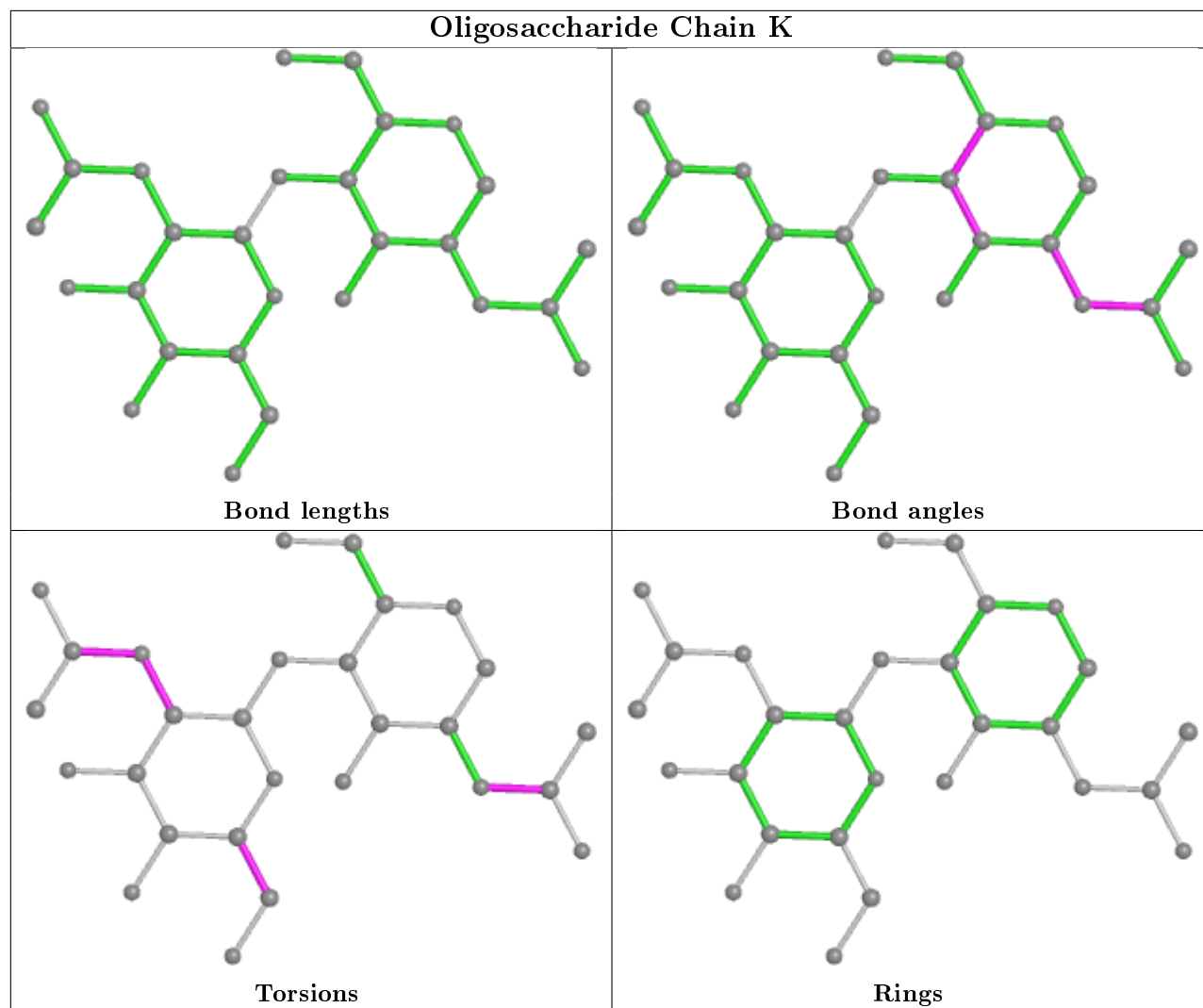
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

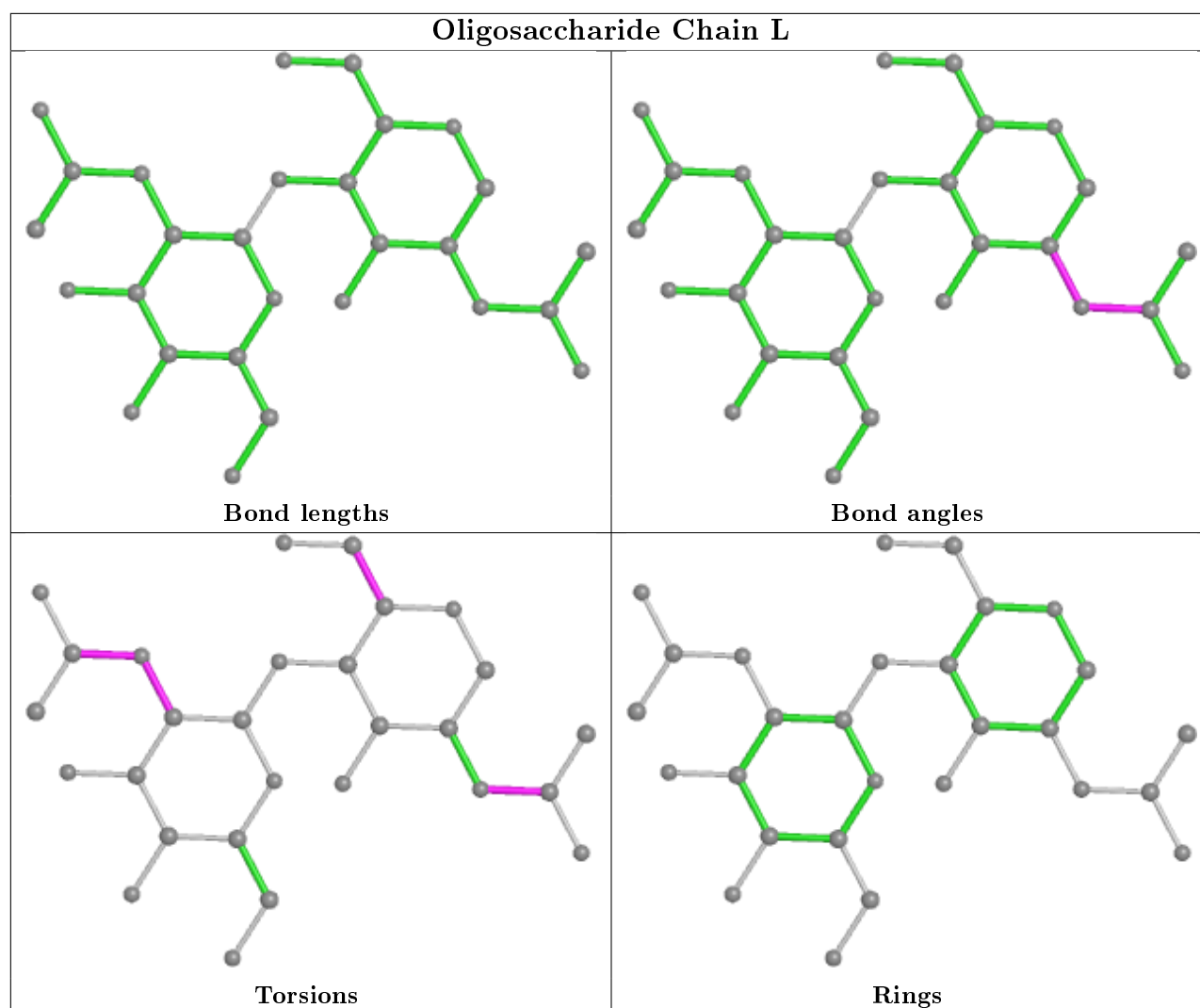












## 5.6 Ligand geometry ⓘ

27 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	$\# Z  > 2$
3	GDL	F	600	-	15,15,15	3.50	3 (20%)	17,21,21	1.66	4 (23%)
5	EDO	C	802	-	3,3,3	0.58	0	2,2,2	0.44	0
4	NAG	F	702	1	14,14,15	0.60	0	17,19,21	0.76	1 (5%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
4	NAG	A	702	1	14,14,15	0.69	0	17,19,21	0.67	0
3	GDL	A	600	-	15,15,15	3.69	4 (26%)	17,21,21	1.69	4 (23%)
5	EDO	B	801	-	3,3,3	0.52	0	2,2,2	0.38	0
5	EDO	E	802	-	3,3,3	0.69	0	2,2,2	0.40	0
5	EDO	A	802	-	3,3,3	0.49	0	2,2,2	0.40	0
5	EDO	B	800	-	3,3,3	0.58	0	2,2,2	0.47	0
4	NAG	B	702	1	14,14,15	0.62	0	17,19,21	0.65	0
5	EDO	F	800	-	3,3,3	0.53	0	2,2,2	0.41	0
5	EDO	F	802	-	3,3,3	0.50	0	2,2,2	0.40	0
3	GDL	E	600	-	15,15,15	3.96	3 (20%)	17,21,21	1.58	4 (23%)
5	EDO	A	800	-	3,3,3	0.57	0	2,2,2	0.52	0
3	GDL	C	600	-	15,15,15	3.76	4 (26%)	17,21,21	1.58	4 (23%)
5	EDO	F	801	-	3,3,3	0.59	0	2,2,2	0.45	0
3	GDL	D	600	-	15,15,15	3.81	5 (33%)	17,21,21	1.57	4 (23%)
5	EDO	A	801	-	3,3,3	0.60	0	2,2,2	0.43	0
4	NAG	F	703	1	14,14,15	0.59	0	17,19,21	0.60	0
5	EDO	E	800	-	3,3,3	0.60	0	2,2,2	0.44	0
5	EDO	E	801	-	3,3,3	0.58	0	2,2,2	0.38	0
5	EDO	C	801	-	3,3,3	0.59	0	2,2,2	0.43	0
4	NAG	E	702	1	14,14,15	0.59	0	17,19,21	0.67	0
3	GDL	B	600	-	15,15,15	3.97	3 (20%)	17,21,21	1.65	3 (17%)
4	NAG	C	703	1	14,14,15	0.60	0	17,19,21	0.70	0
5	EDO	B	802	-	3,3,3	0.57	0	2,2,2	0.44	0
5	EDO	C	800	-	3,3,3	0.56	0	2,2,2	0.46	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	GDL	F	600	-	-	1/6/26/26	0/1/1/1
5	EDO	C	802	-	-	0/1/1/1	-
4	NAG	F	702	1	-	1/6/23/26	0/1/1/1
4	NAG	A	702	1	-	0/6/23/26	0/1/1/1
3	GDL	A	600	-	-	1/6/26/26	0/1/1/1
5	EDO	B	801	-	-	0/1/1/1	-
5	EDO	E	802	-	-	0/1/1/1	-
5	EDO	A	802	-	-	0/1/1/1	-
5	EDO	B	800	-	-	0/1/1/1	-
4	NAG	B	702	1	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	EDO	F	800	-	-	1/1/1/1	-
5	EDO	F	802	-	-	1/1/1/1	-
3	GDL	E	600	-	-	1/6/26/26	0/1/1/1
5	EDO	A	800	-	-	1/1/1/1	-
3	GDL	C	600	-	-	1/6/26/26	0/1/1/1
5	EDO	F	801	-	-	1/1/1/1	-
3	GDL	D	600	-	-	1/6/26/26	0/1/1/1
5	EDO	A	801	-	-	0/1/1/1	-
4	NAG	F	703	1	-	0/6/23/26	0/1/1/1
5	EDO	E	800	-	-	1/1/1/1	-
5	EDO	E	801	-	-	1/1/1/1	-
5	EDO	C	801	-	-	1/1/1/1	-
4	NAG	E	702	1	-	2/6/23/26	0/1/1/1
3	GDL	B	600	-	-	1/6/26/26	0/1/1/1
4	NAG	C	703	1	-	1/6/23/26	0/1/1/1
5	EDO	B	802	-	-	0/1/1/1	-
5	EDO	C	800	-	-	1/1/1/1	-

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	600	GDL	C2-C1	10.20	1.66	1.51
3	B	600	GDL	C3-C2	9.90	1.66	1.52
3	B	600	GDL	C2-C1	9.67	1.65	1.51
3	E	600	GDL	C3-C2	9.56	1.65	1.52
3	C	600	GDL	C2-C1	9.54	1.65	1.51
3	D	600	GDL	C2-C1	9.43	1.65	1.51
3	D	600	GDL	C3-C2	9.29	1.65	1.52
3	A	600	GDL	C3-C2	9.17	1.65	1.52
3	A	600	GDL	C2-C1	9.12	1.65	1.51
3	F	600	GDL	C2-C1	8.97	1.64	1.51
3	C	600	GDL	C3-C2	8.87	1.65	1.52
3	F	600	GDL	C3-C2	7.91	1.63	1.52
3	B	600	GDL	O5-C1	-5.07	1.27	1.34
3	E	600	GDL	O5-C1	-5.04	1.27	1.34
3	F	600	GDL	O5-C1	-4.75	1.27	1.34
3	C	600	GDL	O5-C1	-4.75	1.27	1.34
3	D	600	GDL	O5-C1	-4.47	1.28	1.34
3	A	600	GDL	O5-C1	-4.25	1.28	1.34
3	C	600	GDL	C2-N2	2.30	1.50	1.45
3	D	600	GDL	C2-N2	2.29	1.50	1.45
3	D	600	GDL	C8-C7	2.22	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	600	GDL	C8-C7	2.21	1.55	1.50

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	600	GDL	O5-C1-O1	3.96	124.26	118.47
3	B	600	GDL	O5-C1-O1	3.90	124.16	118.47
3	F	600	GDL	O5-C1-O1	3.83	124.06	118.47
3	C	600	GDL	O5-C1-O1	3.78	123.98	118.47
3	D	600	GDL	O5-C1-O1	3.78	123.98	118.47
3	E	600	GDL	O5-C1-O1	3.48	123.55	118.47
3	A	600	GDL	O1-C1-C2	-3.07	113.78	123.56
3	D	600	GDL	O1-C1-C2	-3.04	113.89	123.56
3	E	600	GDL	O5-C5-C4	3.01	115.08	109.73
3	C	600	GDL	O1-C1-C2	-2.97	114.10	123.56
3	B	600	GDL	O5-C5-C4	2.91	114.90	109.73
3	F	600	GDL	O1-C1-C2	-2.82	114.56	123.56
3	B	600	GDL	O1-C1-C2	-2.79	114.68	123.56
3	A	600	GDL	O5-C5-C4	2.79	114.68	109.73
3	F	600	GDL	C3-C2-N2	-2.66	108.11	112.28
3	E	600	GDL	O1-C1-C2	-2.64	115.14	123.56
3	D	600	GDL	O5-C5-C4	2.23	113.69	109.73
3	A	600	GDL	C3-C2-N2	-2.23	108.79	112.28
3	C	600	GDL	O5-C5-C4	2.23	113.69	109.73
4	F	702	NAG	C2-N2-C7	-2.23	119.73	122.90
3	E	600	GDL	C3-C2-N2	-2.18	108.87	112.28
3	D	600	GDL	C3-C2-N2	-2.14	108.93	112.28
3	F	600	GDL	O5-C5-C4	2.09	113.45	109.73
3	C	600	GDL	C3-C2-N2	-2.09	109.01	112.28

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	600	GDL	C1-C2-N2-C7
4	E	702	NAG	O5-C5-C6-O6
4	E	702	NAG	C4-C5-C6-O6
5	A	800	EDO	O1-C1-C2-O2
5	C	801	EDO	O1-C1-C2-O2
5	F	802	EDO	O1-C1-C2-O2
3	F	600	GDL	C1-C2-N2-C7
3	A	600	GDL	C1-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
3	E	600	GDL	C1-C2-N2-C7
3	C	600	GDL	C1-C2-N2-C7
3	B	600	GDL	C1-C2-N2-C7
5	C	800	EDO	O1-C1-C2-O2
5	F	800	EDO	O1-C1-C2-O2
5	F	801	EDO	O1-C1-C2-O2
5	E	800	EDO	O1-C1-C2-O2
4	C	703	NAG	C4-C5-C6-O6
4	F	702	NAG	C8-C7-N2-C2
5	E	801	EDO	O1-C1-C2-O2

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	702	NAG	1	0
5	B	801	EDO	4	0
5	F	800	EDO	1	0
5	A	800	EDO	2	0
5	A	801	EDO	2	0
5	E	801	EDO	1	0
5	C	801	EDO	2	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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	484/515 (93%)	-0.04	15 (3%)	49	52	27, 38, 58, 85	0
1	B	484/515 (93%)	-0.15	10 (2%)	63	66	24, 37, 57, 85	0
1	C	484/515 (93%)	-0.05	19 (3%)	39	42	27, 38, 67, 88	0
1	D	483/515 (93%)	-0.03	14 (2%)	51	55	27, 39, 64, 93	0
1	E	484/515 (93%)	-0.11	16 (3%)	46	48	25, 36, 54, 86	0
1	F	484/515 (93%)	-0.05	24 (4%)	28	31	28, 39, 66, 90	0
All	All	2903/3090 (93%)	-0.07	98 (3%)	45	47	24, 38, 62, 93	0

All (98) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	554	GLU	13.6
1	C	554	GLU	10.2
1	B	554	GLU	9.1
1	D	317	ASP	7.8
1	E	315	LYS	7.5
1	F	317	ASP	7.3
1	A	317	ASP	6.2
1	E	317	ASP	6.2
1	A	553	HIS	5.7
1	B	317	ASP	5.7
1	D	315	LYS	4.8
1	E	316	LEU	4.7
1	C	553	HIS	4.6
1	A	316	LEU	4.4
1	F	554	GLU	4.4
1	F	553	HIS	4.3
1	C	317	ASP	4.0
1	B	316	LEU	3.9
1	E	553	HIS	3.9

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Mol	Chain	Res	Type	RSRZ
1	B	553	HIS	3.8
1	D	378	THR	3.7
1	A	107	GLY	3.7
1	F	372	ARG	3.6
1	E	311	SER	3.6
1	D	376	PHE	3.6
1	B	359	LYS	3.6
1	A	372	ARG	3.6
1	F	316	LEU	3.4
1	D	309	CYS	3.2
1	C	318	SER	3.2
1	A	431	GLU	3.2
1	C	488	LEU	3.2
1	F	502	LEU	3.2
1	E	107	GLY	3.2
1	C	316	LEU	3.0
1	D	316	LEU	3.0
1	E	476	LYS	3.0
1	A	552	ASN	2.9
1	E	502	LEU	2.9
1	C	310	TYR	2.9
1	B	357	GLU	2.8
1	D	427	SER	2.8
1	F	509	VAL	2.8
1	A	359	LYS	2.8
1	A	376	PHE	2.8
1	B	311	SER	2.8
1	D	359	LYS	2.8
1	D	372	ARG	2.8
1	C	378	THR	2.7
1	F	310	TYR	2.7
1	F	373	GLN	2.6
1	E	509	VAL	2.6
1	F	488	LEU	2.6
1	A	373	GLN	2.6
1	C	172	LEU	2.6
1	F	309	CYS	2.5
1	B	552	ASN	2.5
1	C	359	LYS	2.5
1	F	217	LYS	2.5
1	E	552	ASN	2.5
1	F	475	GLN	2.4

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Mol	Chain	Res	Type	RSRZ
1	F	318	SER	2.4
1	C	369	ASP	2.4
1	C	233	VAL	2.4
1	A	426	ASP	2.4
1	D	303	LYS	2.3
1	C	427	SER	2.3
1	F	205	ILE	2.3
1	F	311	SER	2.3
1	F	476	LYS	2.3
1	D	486	ALA	2.3
1	F	315	LYS	2.3
1	B	75	GLU	2.3
1	E	475	GLN	2.3
1	A	205	ILE	2.3
1	B	107	GLY	2.2
1	C	372	ARG	2.2
1	D	311	SER	2.2
1	F	107	GLY	2.2
1	C	205	ILE	2.2
1	C	309	CYS	2.2
1	F	54	PRO	2.2
1	E	310	TYR	2.2
1	A	174	THR	2.2
1	F	206	LEU	2.2
1	C	552	ASN	2.1
1	E	292	PRO	2.1
1	A	475	GLN	2.1
1	D	475	GLN	2.1
1	E	322	ILE	2.1
1	C	373	GLN	2.1
1	A	476	LYS	2.0
1	D	414	LYS	2.0
1	F	513	LEU	2.0
1	C	377	GLY	2.0
1	F	83	PRO	2.0
1	E	231	PHE	2.0
1	F	552	ASN	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, 95<sup>th</sup> 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.

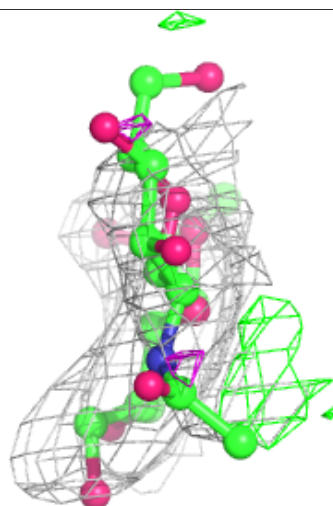
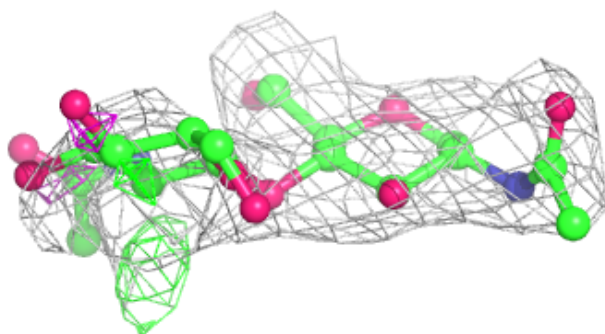
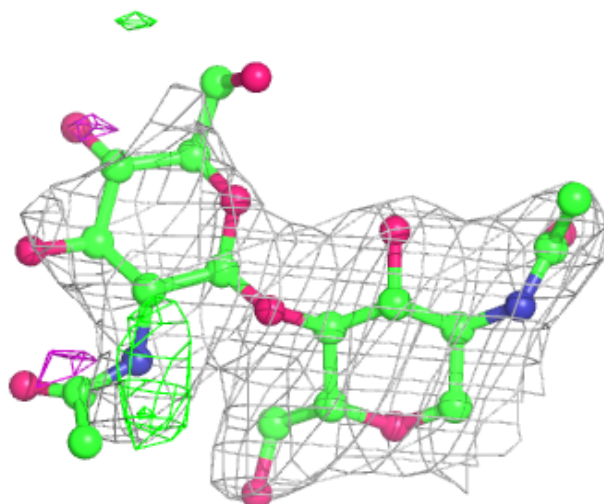
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å <sup>2</sup> )	Q<0.9
2	NAG	H	2	14/15	0.60	0.47	80,84,87,90	0
2	NAG	G	2	14/15	0.71	0.56	89,93,95,96	0
2	NAG	I	2	14/15	0.72	0.56	85,88,90,91	0
2	NAG	J	2	14/15	0.76	0.42	81,83,86,88	0
2	NAG	K	2	14/15	0.78	0.46	81,84,86,87	0
2	NAG	L	2	14/15	0.82	0.54	80,85,88,90	0
2	NAG	L	1	14/15	0.84	0.25	57,60,66,74	0
2	NAG	K	1	14/15	0.85	0.26	64,67,70,76	0
2	NAG	J	1	14/15	0.85	0.21	64,68,72,77	0
2	NAG	H	1	14/15	0.86	0.18	62,65,69,75	0
2	NAG	G	1	14/15	0.88	0.17	66,70,74,82	0
2	NAG	I	1	14/15	0.89	0.24	66,69,74,80	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.



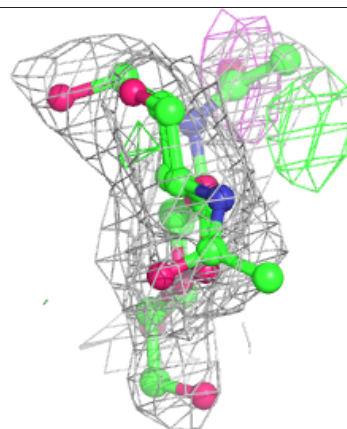
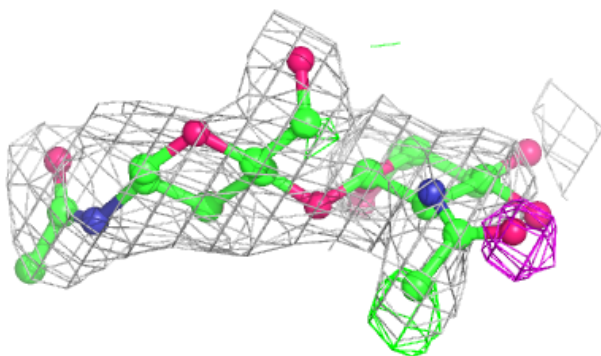
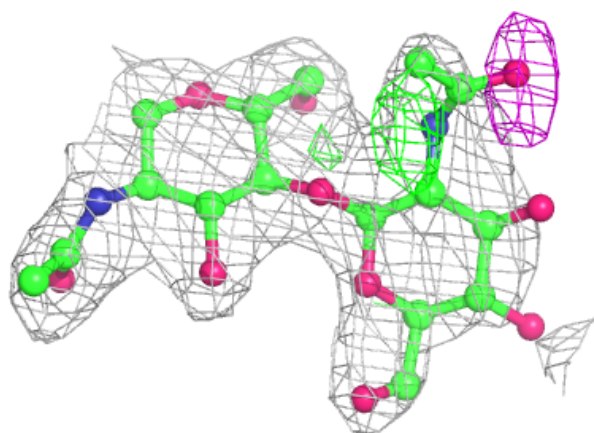
**Electron density around Chain G:**

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

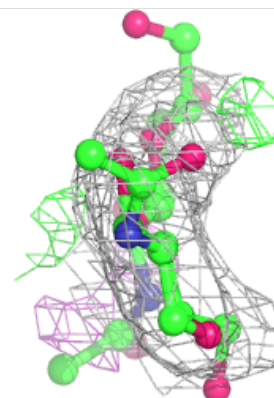
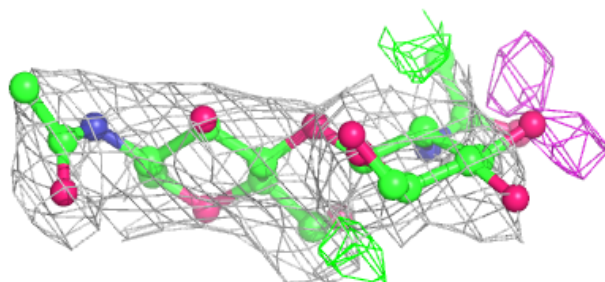
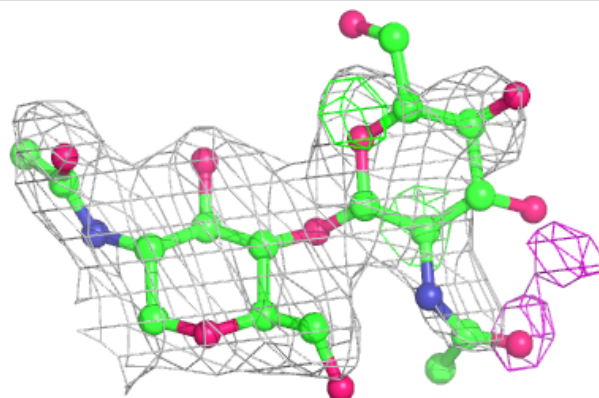


**Electron density around Chain H:**

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

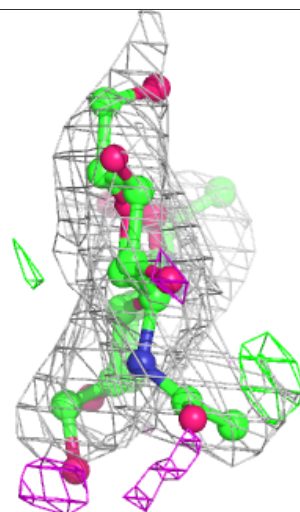
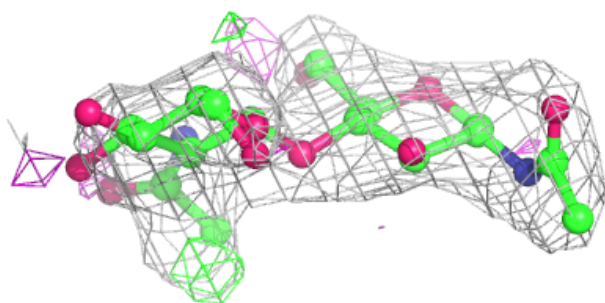
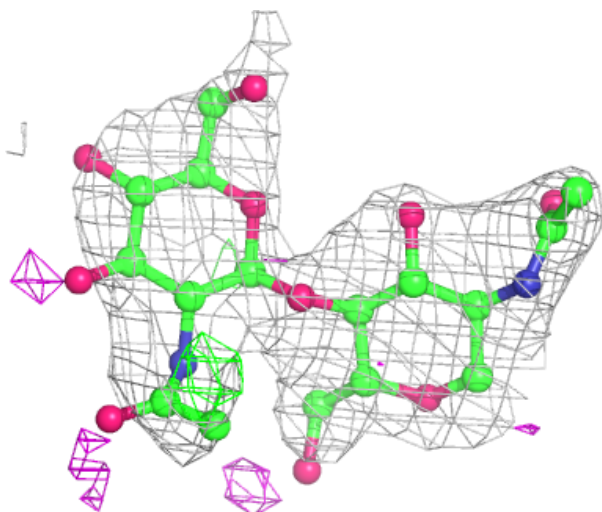
**Electron density around Chain I:**

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



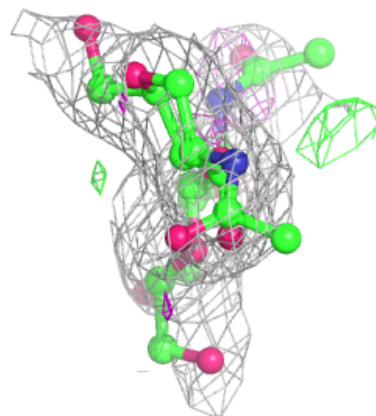
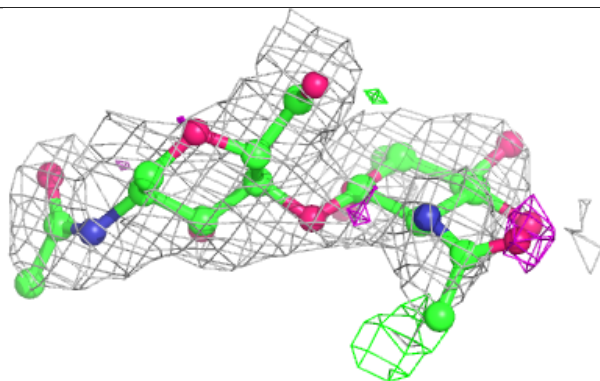
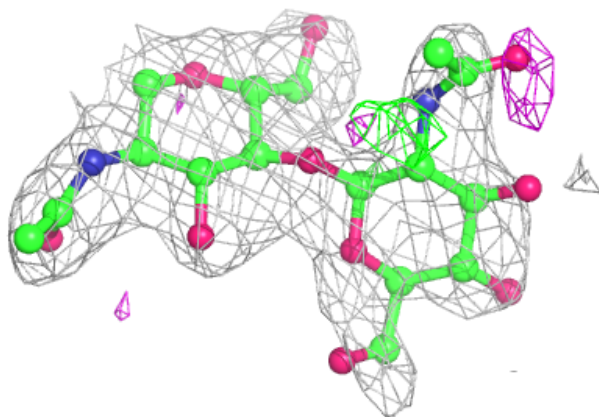
**Electron density around Chain J:**

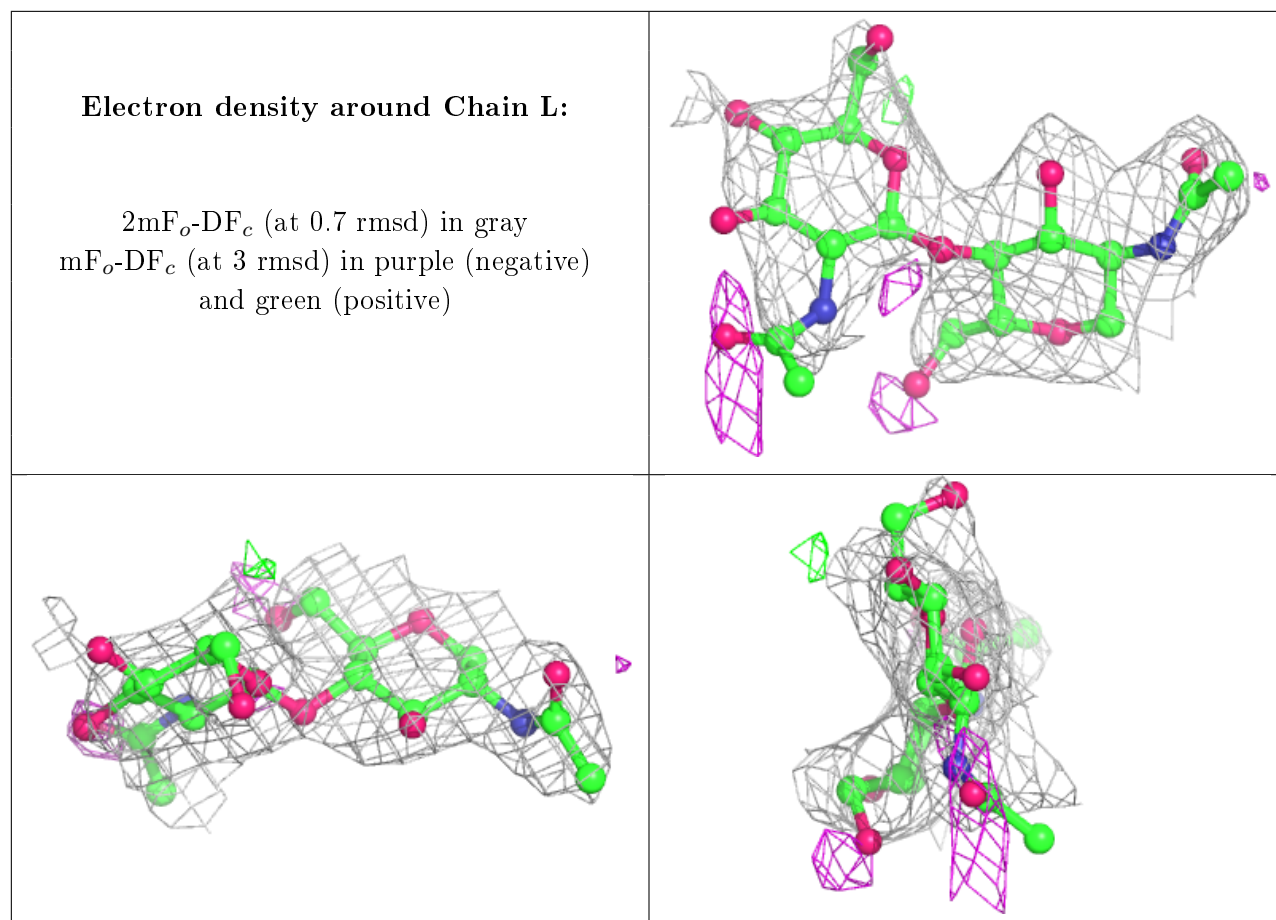
$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



**Electron density around Chain K:**

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





## 6.4 Ligands ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
4	NAG	B	702	14/15	0.67	0.40	70,72,75,76	0
4	NAG	E	702	14/15	0.68	0.48	77,81,82,83	0
4	NAG	C	703	14/15	0.69	0.54	84,88,90,90	0
4	NAG	A	702	14/15	0.70	0.34	64,71,75,78	0
4	NAG	F	702	14/15	0.70	0.52	83,88,90,91	0
5	EDO	C	801	4/4	0.77	0.30	65,66,67,67	0
4	NAG	F	703	14/15	0.80	0.45	79,82,84,84	0
5	EDO	B	801	4/4	0.84	0.31	61,62,63,65	0
5	EDO	A	801	4/4	0.84	0.35	76,76,76,77	0
5	EDO	E	801	4/4	0.89	0.27	65,66,66,67	0
3	GDL	F	600	15/15	0.90	0.10	33,38,40,41	0
5	EDO	E	802	4/4	0.91	0.16	30,31,34,38	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
5	EDO	F	800	4/4	0.91	0.17	48,51,52,56	0
5	EDO	F	801	4/4	0.91	0.31	66,67,67,69	0
5	EDO	C	800	4/4	0.91	0.19	59,59,59,61	0
5	EDO	E	800	4/4	0.92	0.13	49,49,49,51	0
3	GDL	C	600	15/15	0.93	0.10	30,34,38,38	0
5	EDO	B	802	4/4	0.93	0.17	30,35,36,38	0
3	GDL	E	600	15/15	0.93	0.12	29,33,34,36	0
3	GDL	D	600	15/15	0.94	0.11	32,34,37,40	0
3	GDL	A	600	15/15	0.94	0.10	30,32,34,35	0
3	GDL	B	600	15/15	0.95	0.10	30,33,34,35	0
5	EDO	F	802	4/4	0.95	0.17	37,38,39,39	0
5	EDO	A	802	4/4	0.96	0.17	31,38,40,40	0
5	EDO	A	800	4/4	0.97	0.11	56,57,57,61	0
5	EDO	B	800	4/4	0.97	0.12	38,38,38,41	0
5	EDO	C	802	4/4	0.98	0.08	35,36,37,38	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.