



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:16 PM UTC

PDB ID : 1NMC / pdb_00001nmc
Title : COMPLEX BETWEEN NC10 ANTI-INFLUENZA VIRUS NEURAMINIDASE SINGLE CHAIN ANTIBODY WITH A 15 RESIDUE LINKER AND INFLUENZA VIRUS NEURAMINIDASE
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Deposited on : 1997-12-21
Resolution : 2.50 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
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

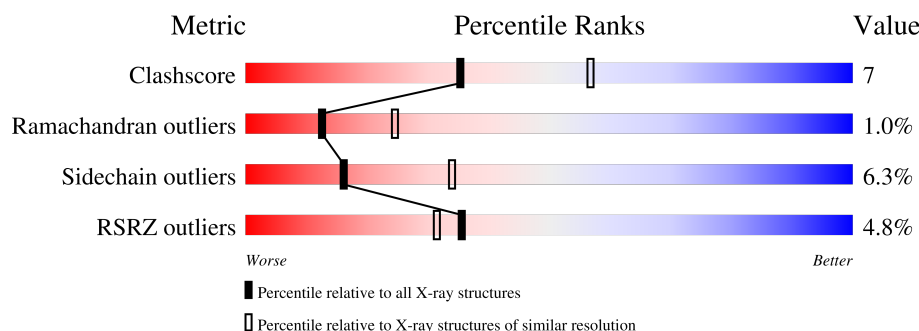
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



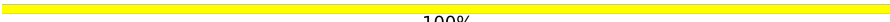
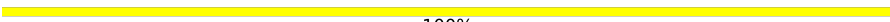
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	388	4% 81% 15% .
1	N	388	2% 83% 14% .
2	B	122	6% 76% 17% 7%
2	H	122	2% 76% 17% 7%
3	C	109	16% 72% 23% 6%
3	L	109	11% 70% 25% 6%

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Mol	Chain	Length	Quality of chain
4	D	6	 100%
4	E	6	 100%

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 10138 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 NEURAMINIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	N	388	Total	C	N	O	S	0	0	0
			3067	1914	538	592	23			
1	A	388	Total	C	N	O	S	0	0	0
			3067	1914	538	592	23			

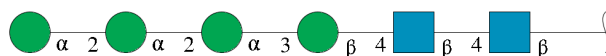
- Molecule 2 is a protein called SINGLE CHAIN ANTIBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	122	Total	C	N	O	S	0	0	0
			943	591	155	192	5			
2	B	122	Total	C	N	O	S	0	0	0
			943	591	155	192	5			

- Molecule 3 is a protein called SINGLE CHAIN ANTIBODY.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	109	Total	C	N	O	S	0	0	0
			857	535	138	182	2			
3	C	109	Total	C	N	O	S	0	0	0
			857	535	138	182	2			

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



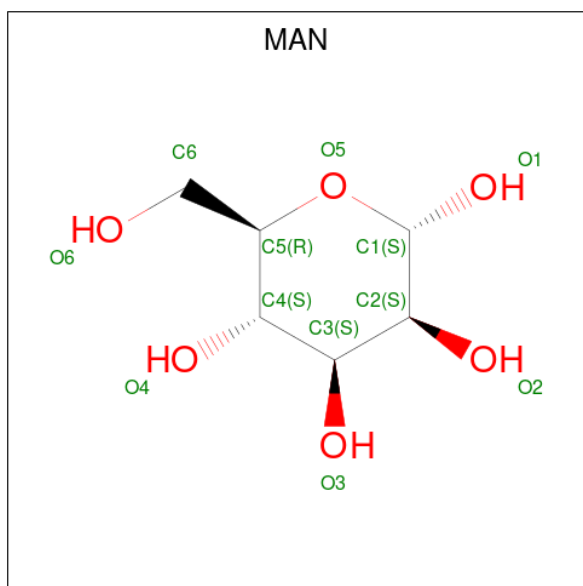
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	6	Total	C	N	O	0	0	0
			72	40	2	30			

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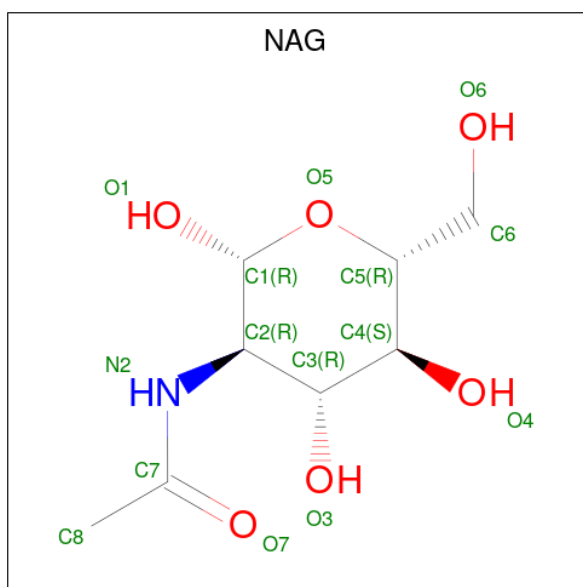
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	E	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 5 is alpha-D-mannopyranose (CCD ID: MAN) (formula: C₆H₁₂O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	N	1	Total	C	O	0	0
			11	6	5		
5	A	1	Total	C	O	0	0
			11	6	5		

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 7 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	N	1	Total	Ca	0	0
			1	1		
7	A	1	Total	Ca	0	0
			1	1		

- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	N	81	Total	O	0	0
			81	81		
8	H	6	Total	O	0	0
			6	6		
8	L	3	Total	O	0	0
			3	3		

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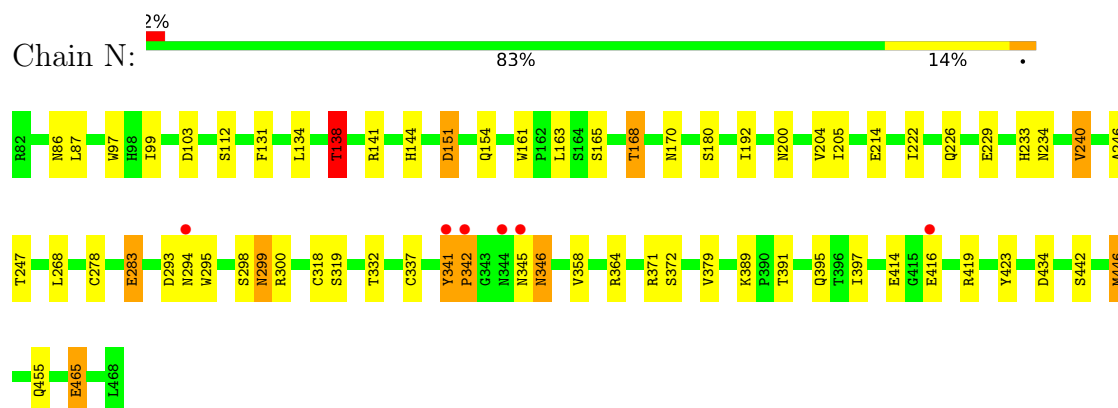
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	81	Total 81	O 81	0	0
8	B	6	Total 6	O 6	0	0
8	C	3	Total 3	O 3	0	0

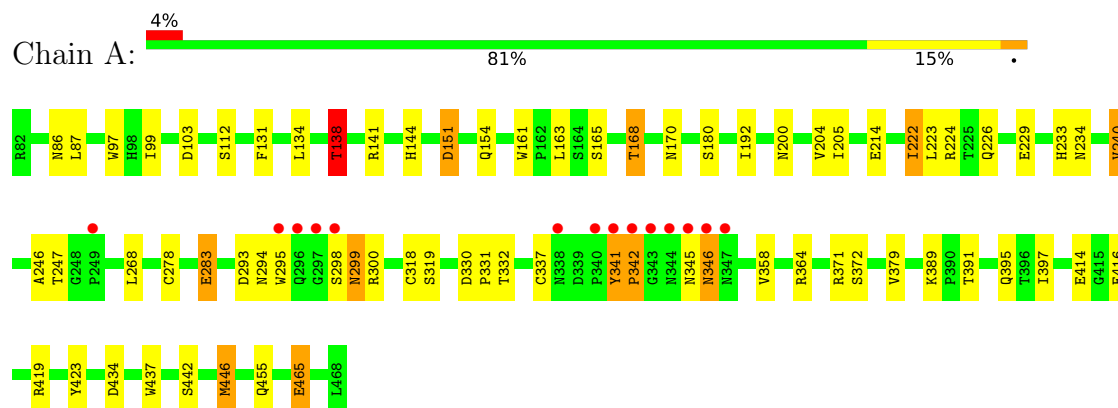
3 Residue-property plots

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

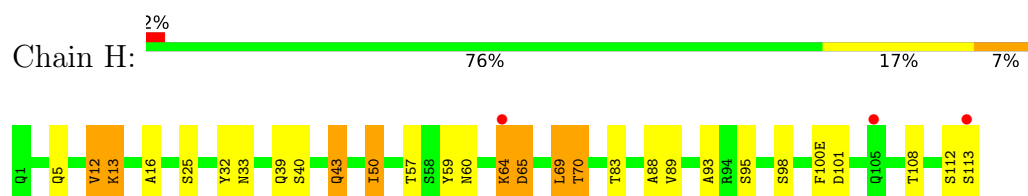
• Molecule 1: NEURAMINIDASE



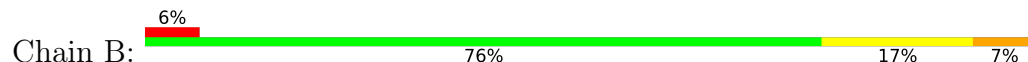
• Molecule 1: NEURAMINIDASE

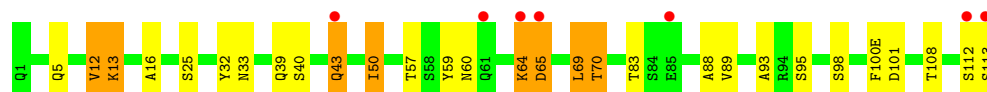


• Molecule 2: SINGLE CHAIN ANTIBODY



• Molecule 2: SINGLE CHAIN ANTIBODY





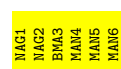
• Molecule 3: SINGLE CHAIN ANTIBODY



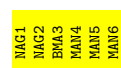
• Molecule 3: SINGLE CHAIN ANTIBODY



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	144.40Å 144.40Å 227.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	7.00 – 2.50 7.00 – 2.50	Depositor EDS
% Data completeness (in resolution range)	80.0 (7.00-2.50) 76.7 (7.00-2.50)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	185.30 (at 2.51Å)	Xtrriage
Refinement program	X-PLOR 3.1	Depositor
R, R_{free}	0.220 , 0.260 0.233 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	24.7	Xtrriage
Anisotropy	0.470	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 69.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	10138	wwPDB-VP
Average B, all atoms (Å ²)	23.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 37.20 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.4472e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: BMA, CA, NAG, MAN

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.73	3/3150 (0.1%)	1.14	26/4290 (0.6%)
1	N	0.73	3/3150 (0.1%)	1.14	25/4290 (0.6%)
2	B	0.55	0/966	0.95	2/1306 (0.2%)
2	H	0.54	0/966	0.95	2/1306 (0.2%)
3	C	0.54	0/874	0.92	3/1187 (0.3%)
3	L	0.54	0/874	0.92	3/1187 (0.3%)
All	All	0.67	6/9980 (0.1%)	1.07	61/13566 (0.4%)

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
1	N	0	1
All	All	0	2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	200	ASN	CG-OD1	5.61	1.34	1.23
1	N	200	ASN	CG-OD1	5.60	1.34	1.23
1	N	161	TRP	NE1-CE2	-5.44	1.31	1.37
1	A	161	TRP	NE1-CE2	-5.43	1.31	1.37
1	A	240	VAL	CA-CB	5.29	1.60	1.54
1	N	240	VAL	CA-CB	5.23	1.60	1.54

All (61) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	N	226	GLN	N-CA-C	11.21	123.06	111.07
1	A	226	GLN	N-CA-C	11.20	123.05	111.07
1	A	300	ARG	N-CA-C	8.74	122.35	110.29
1	N	300	ARG	N-CA-C	8.73	122.34	110.29
1	N	138	THR	CB-CA-C	-7.45	96.76	110.62
1	A	138	THR	CB-CA-C	-7.45	96.77	110.62
1	A	299	ASN	N-CA-C	-7.09	98.75	110.17
1	N	299	ASN	N-CA-C	-7.08	98.76	110.17
1	N	180	SER	N-CA-C	6.88	117.66	108.38
2	H	101	ASP	N-CA-C	6.87	121.41	112.89
2	B	101	ASP	N-CA-C	6.86	121.39	112.89
1	A	180	SER	N-CA-C	6.84	117.61	108.38
1	N	229	GLU	N-CA-C	6.63	119.85	110.50
1	A	229	GLU	N-CA-C	6.62	119.83	110.50
3	L	76	SER	N-CA-C	6.35	118.77	111.02
3	C	76	SER	N-CA-C	6.34	118.75	111.02
2	H	60	ASN	N-CA-C	-6.25	99.83	109.76
2	B	60	ASN	N-CA-C	-6.23	99.85	109.76
1	A	389	LYS	CA-C-N	6.07	126.09	119.90
1	A	389	LYS	C-N-CA	6.07	126.09	119.90
1	N	389	LYS	CA-C-N	6.05	126.07	119.90
1	N	389	LYS	C-N-CA	6.05	126.07	119.90
1	A	319	SER	N-CA-C	6.02	116.52	109.60
1	N	319	SER	N-CA-C	6.01	116.51	109.60
1	A	318	CYS	N-CA-C	5.89	117.78	111.36
1	N	318	CYS	N-CA-C	5.87	117.75	111.36
1	A	346	ASN	N-CA-C	5.77	119.65	111.92
1	N	346	ASN	N-CA-C	5.77	119.65	111.92
1	A	337	CYS	N-CA-C	5.68	120.34	112.90
1	N	337	CYS	N-CA-C	5.68	120.34	112.90
1	N	298	SER	N-CA-C	-5.49	106.58	113.28
1	A	298	SER	N-CA-C	-5.49	106.59	113.28
1	A	151	ASP	N-CA-C	5.46	120.22	113.50
1	N	151	ASP	N-CA-C	5.46	120.21	113.50
1	A	341	TYR	CA-C-N	5.44	126.64	119.84
1	A	341	TYR	C-N-CA	5.44	126.64	119.84
1	N	341	TYR	CA-C-N	5.43	126.63	119.84
1	N	341	TYR	C-N-CA	5.43	126.63	119.84
1	N	165	SER	N-CA-C	-5.36	103.97	110.13
1	A	165	SER	N-CA-C	-5.35	103.98	110.13
1	N	379	VAL	CA-C-N	5.31	125.61	119.93
1	N	379	VAL	C-N-CA	5.31	125.61	119.93
1	A	434	ASP	N-CA-C	5.31	119.40	113.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	N	434	ASP	N-CA-C	5.31	119.39	113.02
1	A	278	CYS	N-CA-C	5.31	118.14	110.28
1	A	379	VAL	CA-C-N	5.31	125.61	119.93
1	A	379	VAL	C-N-CA	5.31	125.61	119.93
1	N	278	CYS	N-CA-C	5.30	118.13	110.28
1	N	455	GLN	N-CA-C	5.29	118.62	110.42
1	A	455	GLN	N-CA-C	5.29	118.61	110.42
1	N	246	ALA	N-CA-C	-5.24	104.48	111.24
1	A	246	ALA	N-CA-C	-5.21	104.52	111.24
1	N	240	VAL	N-CA-C	5.16	115.40	108.17
1	A	240	VAL	N-CA-C	5.16	115.39	108.17
3	L	17	ASP	N-CA-C	5.15	117.48	109.60
3	C	17	ASP	N-CA-C	5.13	117.45	109.60
1	A	103	ASP	N-CA-C	5.09	119.76	113.50
1	N	103	ASP	N-CA-C	5.07	119.74	113.50
3	C	78	LEU	N-CA-C	5.05	117.42	110.35
3	L	78	LEU	N-CA-C	5.03	117.40	110.35
1	A	223	LEU	N-CA-C	-5.00	102.97	110.23

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	423	TYR	Sidechain
1	N	423	TYR	Sidechain

5.2 Too-close contacts [i](#)

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	3067	0	2893	30	1
1	N	3067	0	2893	27	1
2	B	943	0	880	20	0
2	H	943	0	880	20	0
3	C	857	0	805	22	0
3	L	857	0	805	23	0
4	D	72	0	61	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	72	0	61	0	0
5	A	11	0	10	0	0
5	N	11	0	10	0	0
6	A	28	0	26	1	0
6	N	28	0	26	0	0
7	A	1	0	0	0	0
7	N	1	0	0	0	0
8	A	81	0	0	3	0
8	B	6	0	0	1	0
8	C	3	0	0	0	0
8	H	6	0	0	1	0
8	L	3	0	0	0	0
8	N	81	0	0	3	0
All	All	10138	0	9350	142	2

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 7.

All (142) 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)
2:B:89:VAL:HG22	2:B:108:THR:HG22	1.57	0.87
2:H:89:VAL:HG22	2:H:108:THR:HG22	1.57	0.87
1:N:168:THR:HB	1:N:170:ASN:OD1	1.85	0.77
1:A:168:THR:HB	1:A:170:ASN:OD1	1.85	0.77
1:A:168:THR:HG22	1:A:170:ASN:H	1.49	0.77
1:N:168:THR:HG22	1:N:170:ASN:H	1.49	0.76
1:N:87:LEU:H	1:N:233:HIS:HD2	1.33	0.75
1:A:87:LEU:H	1:A:233:HIS:HD2	1.33	0.75
2:H:64:LYS:HE3	2:H:65:ASP:H	1.51	0.74
2:B:64:LYS:HE3	2:B:65:ASP:H	1.51	0.74
3:C:59:PRO:HG2	3:C:62:PHE:HE2	1.53	0.73
3:L:59:PRO:HG2	3:L:62:PHE:HE2	1.53	0.73
1:A:138:THR:HG23	1:A:144:HIS:HB2	1.73	0.69
3:C:59:PRO:HG2	3:C:62:PHE:CE2	2.27	0.69
1:N:138:THR:HG23	1:N:144:HIS:HB2	1.73	0.68
3:L:59:PRO:HG2	3:L:62:PHE:CE2	2.27	0.68
1:N:416:GLU:H	1:N:416:GLU:CD	2.09	0.61
1:N:345:ASN:HA	8:N:543:HOH:O	2.00	0.61
1:A:416:GLU:CD	1:A:416:GLU:H	2.09	0.61
1:A:345:ASN:HA	8:A:543:HOH:O	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:THR:HG22	1:A:170:ASN:N	2.16	0.60
1:N:168:THR:HG22	1:N:170:ASN:N	2.16	0.60
2:H:59:TYR:CE2	2:H:69:LEU:HD22	2.37	0.60
3:L:24:ARG:HG3	3:L:70:ASP:OD1	2.02	0.60
3:C:24:ARG:HG3	3:C:70:ASP:OD1	2.02	0.59
2:B:59:TYR:CE2	2:B:69:LEU:HD22	2.37	0.59
2:H:32:TYR:CD2	2:H:98:SER:HA	2.38	0.59
2:B:32:TYR:CD2	2:B:98:SER:HA	2.38	0.59
3:C:12:SER:HA	3:C:105:GLU:O	2.03	0.59
3:L:12:SER:HA	3:L:105:GLU:O	2.03	0.59
2:B:64:LYS:HE3	2:B:65:ASP:N	2.18	0.58
1:A:247:THR:HA	1:A:346:ASN:ND2	2.19	0.58
1:A:138:THR:CG2	1:A:144:HIS:HB2	2.34	0.58
1:N:138:THR:CG2	1:N:144:HIS:HB2	2.34	0.58
1:N:295:TRP:O	1:N:345:ASN:HB2	2.05	0.57
2:H:64:LYS:HE3	2:H:65:ASP:N	2.18	0.57
1:A:295:TRP:O	1:A:345:ASN:HB2	2.05	0.57
2:B:64:LYS:HE3	2:B:64:LYS:HA	1.86	0.57
1:N:247:THR:HA	1:N:346:ASN:ND2	2.19	0.56
2:H:64:LYS:HE3	2:H:64:LYS:HA	1.86	0.56
2:B:64:LYS:O	2:B:65:ASP:HB2	2.05	0.56
2:H:64:LYS:O	2:H:65:ASP:HB2	2.05	0.55
1:N:247:THR:HG22	1:N:346:ASN:HB3	1.88	0.55
1:N:397:ILE:HD12	1:N:446:MET:HE2	1.88	0.55
1:A:247:THR:HG22	1:A:346:ASN:HB3	1.88	0.55
1:A:397:ILE:HD12	1:A:446:MET:HE2	1.88	0.54
3:L:18:ARG:HB2	3:L:18:ARG:NH1	2.23	0.54
1:N:295:TRP:HA	1:N:346:ASN:ND2	2.23	0.54
1:A:295:TRP:HA	1:A:346:ASN:ND2	2.23	0.54
3:L:21:ILE:HD12	3:L:22:SER:H	1.73	0.53
3:C:21:ILE:HD12	3:C:22:SER:H	1.73	0.53
1:A:295:TRP:HD1	1:A:346:ASN:HD21	1.54	0.53
3:C:13:ALA:O	3:C:106:ILE:HA	2.08	0.53
1:N:295:TRP:HD1	1:N:346:ASN:HD21	1.55	0.53
3:C:18:ARG:HB2	3:C:18:ARG:NH1	2.22	0.53
1:N:86:ASN:OD1	1:N:234:ASN:HB2	2.09	0.53
3:L:13:ALA:O	3:L:106:ILE:HA	2.08	0.52
1:N:293:ASP:O	1:N:295:TRP:N	2.42	0.52
1:A:293:ASP:O	1:A:295:TRP:N	2.42	0.52
3:C:19:VAL:HG22	3:C:78:LEU:HD11	1.91	0.52
1:A:86:ASN:OD1	1:A:234:ASN:HB2	2.09	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:19:VAL:HG22	3:L:78:LEU:HD11	1.91	0.52
2:B:43:GLN:NE2	2:B:43:GLN:HA	2.24	0.52
2:H:43:GLN:HA	2:H:43:GLN:NE2	2.24	0.51
3:C:48:ILE:CD1	3:C:54:LEU:HD12	2.40	0.51
3:L:48:ILE:CD1	3:L:54:LEU:HD12	2.40	0.51
1:N:87:LEU:H	1:N:233:HIS:CD2	2.22	0.51
3:L:80:GLN:OE1	3:L:106:ILE:HD12	2.11	0.51
1:A:87:LEU:H	1:A:233:HIS:CD2	2.23	0.50
3:C:80:GLN:OE1	3:C:106:ILE:HD12	2.11	0.50
2:H:12:VAL:CG1	2:H:16:ALA:HB3	2.42	0.50
1:A:299:ASN:OD1	1:A:341:TYR:N	2.45	0.50
1:N:299:ASN:OD1	1:N:341:TYR:N	2.45	0.50
3:L:38:GLN:HG3	3:L:44:VAL:HG22	1.94	0.50
3:C:80:GLN:OE1	3:C:80:GLN:HA	2.12	0.49
2:H:40:SER:OG	2:H:43:GLN:HB3	2.12	0.49
3:L:11:LEU:O	3:L:104:LEU:HD12	2.12	0.49
2:B:40:SER:OG	2:B:43:GLN:HB3	2.12	0.49
3:C:38:GLN:HG3	3:C:44:VAL:HG22	1.94	0.49
2:B:12:VAL:CG1	2:B:16:ALA:HB3	2.42	0.49
1:A:295:TRP:HA	1:A:346:ASN:HD22	1.78	0.49
1:A:295:TRP:CD1	1:A:346:ASN:HD21	2.30	0.49
2:B:93:ALA:HB1	2:B:100(E):PHE:HB3	1.95	0.49
1:N:295:TRP:HA	1:N:346:ASN:HD22	1.78	0.49
3:L:80:GLN:OE1	3:L:80:GLN:HA	2.12	0.49
1:N:295:TRP:CD1	1:N:346:ASN:HD21	2.30	0.49
3:C:11:LEU:O	3:C:104:LEU:HD12	2.12	0.49
2:H:93:ALA:HB1	2:H:100(E):PHE:HB3	1.95	0.48
1:A:442:SER:HB2	8:A:519:HOH:O	2.14	0.48
2:H:13:LYS:HD2	2:H:113:SER:HA	1.95	0.48
2:B:33:ASN:HB2	2:B:95:SER:HB2	1.95	0.48
2:H:33:ASN:HB2	2:H:95:SER:HB2	1.95	0.48
1:A:131:PHE:CE1	1:A:163:LEU:HD12	2.49	0.48
1:N:192:ILE:HG12	1:N:205:ILE:HG13	1.95	0.47
1:N:442:SER:HB2	8:N:519:HOH:O	2.14	0.47
2:H:64:LYS:O	2:H:65:ASP:CB	2.62	0.47
2:B:13:LYS:HD2	2:B:113:SER:HA	1.95	0.47
2:B:64:LYS:O	2:B:65:ASP:CB	2.62	0.47
3:L:49:TYR:O	3:L:53:ASN:HB2	2.14	0.47
3:C:49:TYR:O	3:C:53:ASN:HB2	2.14	0.47
1:N:131:PHE:CE1	1:N:163:LEU:HD12	2.49	0.47
1:A:192:ILE:HG12	1:A:205:ILE:HG13	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:341:TYR:HA	1:A:342:PRO:HD3	1.76	0.47
3:C:12:SER:HB3	3:C:107:ARG:HB2	1.97	0.47
3:L:12:SER:HB3	3:L:107:ARG:HB2	1.97	0.46
2:H:12:VAL:HG13	2:H:16:ALA:HB3	1.97	0.46
2:B:12:VAL:HG13	2:B:16:ALA:HB3	1.97	0.46
2:H:100(E):PHE:CD1	2:H:100(E):PHE:N	2.84	0.45
3:C:48:ILE:HA	3:C:53:ASN:O	2.17	0.45
3:L:48:ILE:HA	3:L:53:ASN:O	2.17	0.45
2:B:100(E):PHE:N	2:B:100(E):PHE:CD1	2.84	0.45
1:N:97:TRP:HB2	1:N:395:GLN:NE2	2.32	0.44
1:A:97:TRP:HB2	1:A:395:GLN:NE2	2.32	0.44
2:H:50:ILE:O	2:H:57:THR:HA	2.18	0.43
2:B:50:ILE:O	2:B:57:THR:HA	2.18	0.43
1:N:341:TYR:HA	1:N:342:PRO:HD3	1.76	0.43
3:L:86:TYR:O	3:L:101:GLY:HA2	2.18	0.43
3:C:86:TYR:O	3:C:101:GLY:HA2	2.18	0.42
1:N:465:GLU:CD	1:N:465:GLU:H	2.27	0.42
1:A:330:ASP:HA	1:A:331:PRO:HD2	1.95	0.42
3:L:21:ILE:HD12	3:L:22:SER:N	2.35	0.41
1:A:465:GLU:H	1:A:465:GLU:CD	2.27	0.41
3:C:21:ILE:HD12	3:C:22:SER:N	2.35	0.41
2:H:39:GLN:O	2:H:88:ALA:HB1	2.21	0.41
1:A:222:ILE:O	1:A:224:ARG:HG2	2.21	0.41
2:B:39:GLN:O	2:B:88:ALA:HB1	2.21	0.41
3:C:7:THR:O	3:C:8:THR:HG22	2.21	0.41
1:A:283:GLU:HG3	8:A:535:HOH:O	2.21	0.41
3:C:15:LEU:HD23	3:C:16:GLY:N	2.36	0.41
3:L:11:LEU:HB3	3:L:104:LEU:HD12	2.02	0.41
1:N:283:GLU:HG3	8:N:535:HOH:O	2.21	0.40
2:H:12:VAL:CG1	2:H:13:LYS:N	2.84	0.40
2:H:70:THR:HB	8:H:116:HOH:O	2.21	0.40
3:L:7:THR:O	3:L:8:THR:HG22	2.21	0.40
1:A:437:TRP:CD1	6:A:476(A):NAG:H82	2.56	0.40
3:C:11:LEU:HB3	3:C:104:LEU:HD12	2.02	0.40
3:L:15:LEU:HD23	3:L:16:GLY:N	2.36	0.40
3:L:18:ARG:HG3	3:L:76:SER:HA	2.03	0.40
3:C:18:ARG:HG3	3:C:76:SER:HA	2.03	0.40
3:L:30:SER:O	3:L:31:ASN:HB2	2.21	0.40
2:B:12:VAL:CG1	2:B:13:LYS:N	2.84	0.40
2:B:70:THR:HB	8:B:149:HOH:O	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the sym-

metry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:N:214:GLU:OE2	1:N:419:ARG:NH2[4_555]	2.09	0.11
1:A:214:GLU:OE2	1:A:419:ARG:NH2[4_555]	2.09	0.11

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	386/388 (100%)	357 (92%)	26 (7%)	3 (1%)	16	31
1	N	386/388 (100%)	357 (92%)	26 (7%)	3 (1%)	16	31
2	B	120/122 (98%)	114 (95%)	4 (3%)	2 (2%)	7	13
2	H	120/122 (98%)	114 (95%)	4 (3%)	2 (2%)	7	13
3	C	107/109 (98%)	95 (89%)	11 (10%)	1 (1%)	14	27
3	L	107/109 (98%)	95 (89%)	11 (10%)	1 (1%)	14	27
All	All	1226/1238 (99%)	1132 (92%)	82 (7%)	12 (1%)	12	24

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	N	294	ASN
3	L	8	THR
1	A	294	ASN
3	C	8	THR
2	H	43	GLN
2	H	65	ASP
2	B	43	GLN
2	B	65	ASP
1	N	342	PRO
1	A	342	PRO
1	N	222	ILE
1	A	222	ILE

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	341/341 (100%)	320 (94%)	21 (6%)	16	34
1	N	341/341 (100%)	320 (94%)	21 (6%)	16	34
2	B	100/100 (100%)	90 (90%)	10 (10%)	7	15
2	H	100/100 (100%)	90 (90%)	10 (10%)	7	15
3	C	98/98 (100%)	95 (97%)	3 (3%)	35	62
3	L	98/98 (100%)	95 (97%)	3 (3%)	35	62
All	All	1078/1078 (100%)	1010 (94%)	68 (6%)	16	34

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

Mol	Chain	Res	Type
1	N	99	ILE
1	N	112	SER
1	N	134	LEU
1	N	138	THR
1	N	141	ARG
1	N	151	ASP
1	N	154	GLN
1	N	168	THR
1	N	204	VAL
1	N	240	VAL
1	N	268	LEU
1	N	283	GLU
1	N	332	THR
1	N	358	VAL
1	N	364	ARG
1	N	371	ARG
1	N	372	SER
1	N	391	THR
1	N	414	GLU
1	N	446	MET
1	N	465	GLU
2	H	5	GLN

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Mol	Chain	Res	Type
2	H	12	VAL
2	H	13	LYS
2	H	25	SER
2	H	50	ILE
2	H	64	LYS
2	H	69	LEU
2	H	70	THR
2	H	83	THR
2	H	112	SER
3	L	54	LEU
3	L	105	GLU
3	L	107	ARG
1	A	99	ILE
1	A	112	SER
1	A	134	LEU
1	A	138	THR
1	A	141	ARG
1	A	151	ASP
1	A	154	GLN
1	A	168	THR
1	A	204	VAL
1	A	240	VAL
1	A	268	LEU
1	A	283	GLU
1	A	332	THR
1	A	358	VAL
1	A	364	ARG
1	A	371	ARG
1	A	372	SER
1	A	391	THR
1	A	414	GLU
1	A	446	MET
1	A	465	GLU
2	B	5	GLN
2	B	12	VAL
2	B	13	LYS
2	B	25	SER
2	B	50	ILE
2	B	64	LYS
2	B	69	LEU
2	B	70	THR
2	B	83	THR

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Mol	Chain	Res	Type
2	B	112	SER
3	C	54	LEU
3	C	105	GLU
3	C	107	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	N	154	GLN
1	N	198	ASN
1	N	233	HIS
1	N	345	ASN
1	N	346	ASN
1	N	381	ASN
2	H	3	GLN
2	H	5	GLN
2	H	76	ASN
1	A	198	ASN
1	A	233	HIS
1	A	345	ASN
1	A	346	ASN
1	A	381	ASN
2	B	76	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

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
4	NAG	D	1	4,1	14,14,15	1.47	3 (21%)	17,19,21	4.00	7 (41%)
4	NAG	D	2	4	14,14,15	1.28	1 (7%)	17,19,21	1.43	3 (17%)
4	BMA	D	3	4	11,11,12	1.11	0	15,15,17	1.24	1 (6%)
4	MAN	D	4	4	11,11,12	0.95	1 (9%)	15,15,17	1.33	2 (13%)
4	MAN	D	5	4	11,11,12	1.36	1 (9%)	15,15,17	1.42	3 (20%)
4	MAN	D	6	4	11,11,12	1.91	2 (18%)	15,15,17	1.55	1 (6%)
4	NAG	E	1	4,1	14,14,15	1.46	3 (21%)	17,19,21	4.00	7 (41%)
4	NAG	E	2	4	14,14,15	1.27	1 (7%)	17,19,21	1.43	3 (17%)
4	BMA	E	3	4	11,11,12	1.12	1 (9%)	15,15,17	1.24	1 (6%)
4	MAN	E	4	4	11,11,12	0.94	0	15,15,17	1.34	2 (13%)
4	MAN	E	5	4	11,11,12	1.35	1 (9%)	15,15,17	1.43	3 (20%)
4	MAN	E	6	4	11,11,12	1.90	2 (18%)	15,15,17	1.54	1 (6%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	D	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	D	2	4	-	0/6/23/26	0/1/1/1
4	BMA	D	3	4	-	1/2/19/22	0/1/1/1
4	MAN	D	4	4	-	0/2/19/22	0/1/1/1
4	MAN	D	5	4	-	0/2/19/22	0/1/1/1
4	MAN	D	6	4	-	2/2/19/22	0/1/1/1
4	NAG	E	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	E	2	4	-	0/6/23/26	0/1/1/1
4	BMA	E	3	4	-	1/2/19/22	0/1/1/1
4	MAN	E	4	4	-	0/2/19/22	0/1/1/1
4	MAN	E	5	4	-	0/2/19/22	0/1/1/1
4	MAN	E	6	4	-	2/2/19/22	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	D	6	MAN	O5-C5	4.91	1.53	1.43
4	E	6	MAN	O5-C5	4.90	1.53	1.43
4	D	5	MAN	C2-C3	-2.78	1.48	1.52
4	E	5	MAN	C2-C3	-2.77	1.48	1.52
4	D	2	NAG	C1-C2	-2.69	1.48	1.52
4	E	1	NAG	O5-C5	2.65	1.48	1.43
4	D	1	NAG	O5-C5	2.65	1.48	1.43
4	E	2	NAG	C1-C2	-2.62	1.48	1.52
4	D	1	NAG	C6-C5	2.61	1.60	1.51
4	E	1	NAG	C6-C5	2.60	1.60	1.51
4	E	6	MAN	C6-C5	2.37	1.59	1.51
4	D	6	MAN	C6-C5	2.37	1.59	1.51
4	D	1	NAG	C1-C2	2.30	1.55	1.52
4	E	1	NAG	C1-C2	2.29	1.55	1.52
4	D	4	MAN	O5-C5	2.03	1.47	1.43
4	E	3	BMA	C2-C3	2.01	1.55	1.52

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	E	1	NAG	C1-O5-C5	11.88	128.11	112.19
4	D	1	NAG	C1-O5-C5	11.86	128.09	112.19
4	E	1	NAG	C2-N2-C7	7.13	132.45	122.90
4	D	1	NAG	C2-N2-C7	7.11	132.43	122.90
4	D	1	NAG	C1-C2-N2	-6.08	100.86	110.43
4	E	1	NAG	C1-C2-N2	-6.07	100.87	110.43
4	D	6	MAN	C1-O5-C5	4.49	118.21	112.19
4	E	6	MAN	C1-O5-C5	4.48	118.19	112.19
4	E	4	MAN	C6-C5-C4	-3.33	104.85	113.02
4	D	4	MAN	C6-C5-C4	-3.31	104.88	113.02
4	D	1	NAG	C3-C4-C5	2.99	115.65	110.23
4	E	1	NAG	C3-C4-C5	2.98	115.64	110.23
4	E	1	NAG	O5-C1-C2	2.97	115.89	111.29
4	D	1	NAG	O5-C1-C2	2.96	115.87	111.29
4	E	2	NAG	C6-C5-C4	2.96	120.29	113.02
4	D	2	NAG	C6-C5-C4	2.95	120.25	113.02
4	E	5	MAN	C3-C4-C5	2.75	115.22	110.23
4	D	5	MAN	C3-C4-C5	2.74	115.20	110.23
4	E	1	NAG	C6-C5-C4	-2.49	106.90	113.02
4	D	3	BMA	O5-C5-C6	2.49	112.51	107.66
4	D	1	NAG	C6-C5-C4	-2.49	106.92	113.02
4	E	3	BMA	O5-C5-C6	2.47	112.48	107.66
4	E	4	MAN	C1-O5-C5	2.44	115.45	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	D	4	MAN	C1-O5-C5	2.42	115.42	112.19
4	D	2	NAG	O7-C7-C8	-2.37	117.84	122.05
4	E	2	NAG	O7-C7-C8	-2.36	117.86	122.05
4	E	5	MAN	O5-C5-C4	-2.35	105.12	110.83
4	D	5	MAN	O5-C5-C4	-2.34	105.14	110.83
4	D	1	NAG	O3-C3-C2	2.31	114.19	109.40
4	E	1	NAG	O3-C3-C2	2.30	114.17	109.40
4	D	2	NAG	C8-C7-N2	2.22	119.79	116.12
4	E	2	NAG	C8-C7-N2	2.20	119.77	116.12
4	E	5	MAN	O5-C5-C6	2.17	111.89	107.66
4	D	5	MAN	O5-C5-C6	2.14	111.84	107.66

There are no chirality outliers.

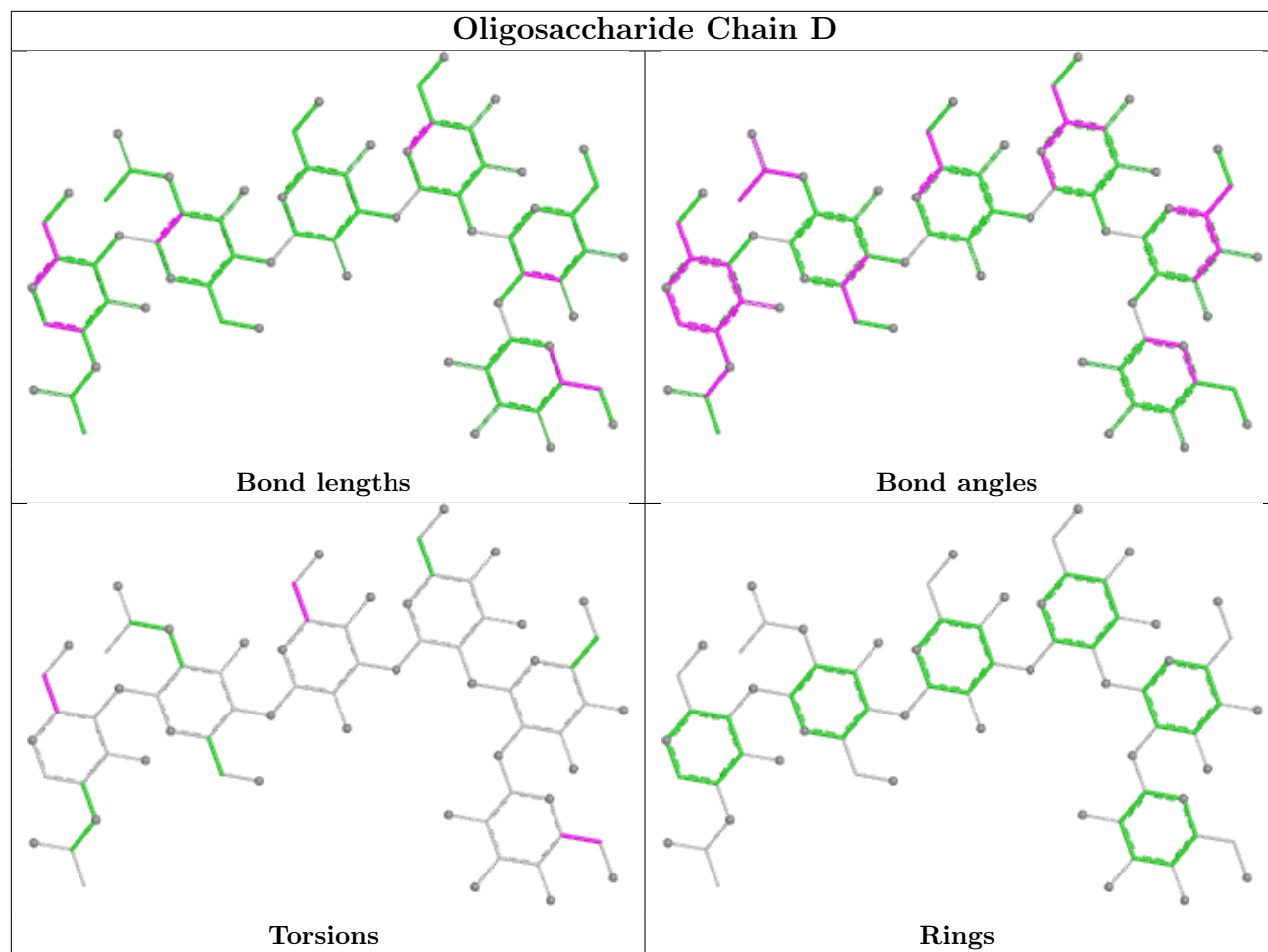
All (10) torsion outliers are listed below:

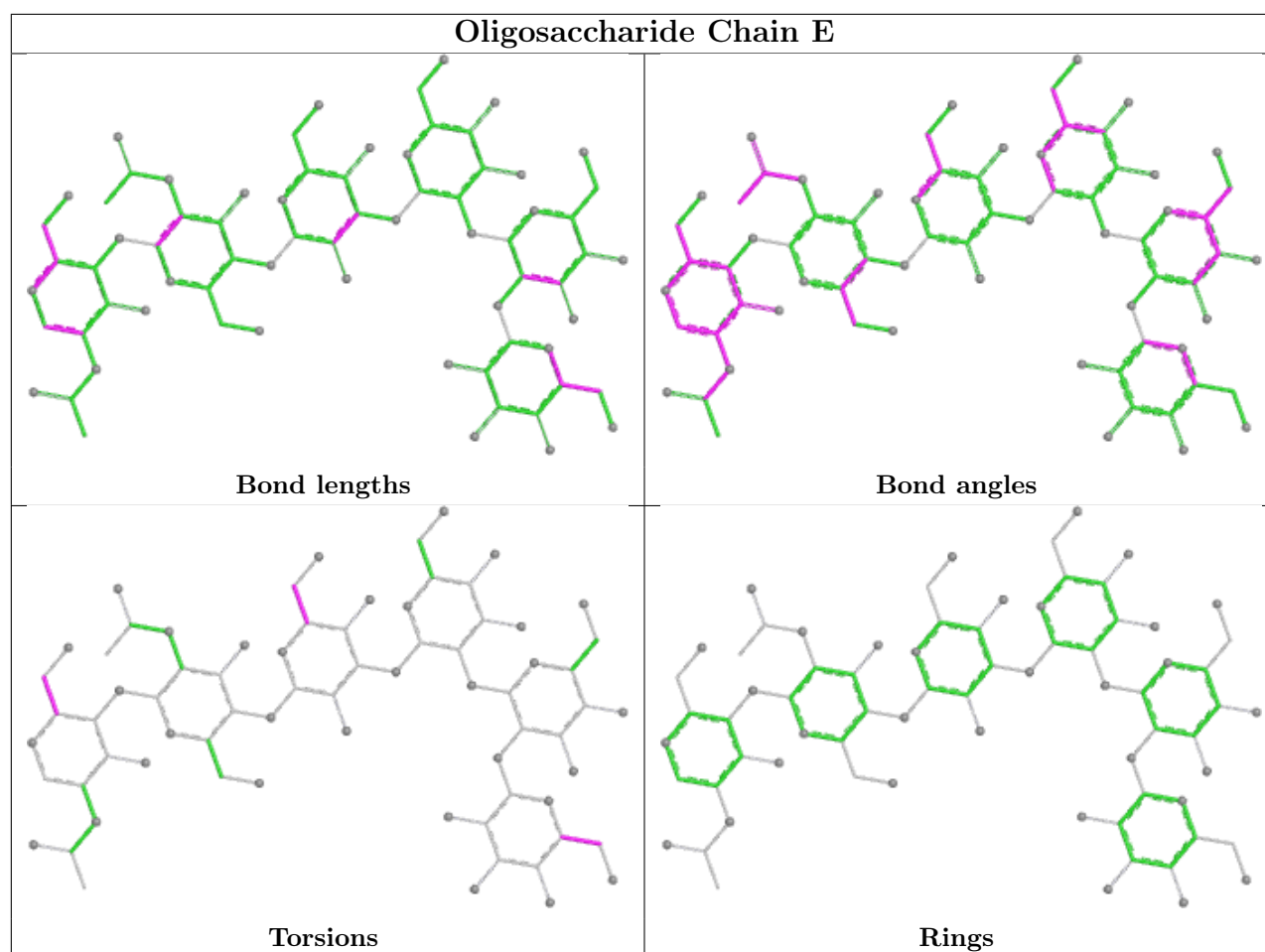
Mol	Chain	Res	Type	Atoms
4	D	6	MAN	O5-C5-C6-O6
4	E	6	MAN	O5-C5-C6-O6
4	D	1	NAG	C4-C5-C6-O6
4	E	1	NAG	C4-C5-C6-O6
4	D	1	NAG	O5-C5-C6-O6
4	E	1	NAG	O5-C5-C6-O6
4	D	6	MAN	C4-C5-C6-O6
4	E	6	MAN	C4-C5-C6-O6
4	D	3	BMA	O5-C5-C6-O6
4	E	3	BMA	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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





5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 2 are monoatomic - leaving 6 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
6	NAG	A	477(A)	1	14,14,15	1.16	3 (21%)	17,19,21	1.59	5 (29%)
5	MAN	N	475(G)	-	11,11,12	1.80	3 (27%)	15,15,17	2.50	7 (46%)
6	NAG	N	477(A)	1	14,14,15	1.16	3 (21%)	17,19,21	1.58	5 (29%)
6	NAG	N	476(A)	1	14,14,15	1.42	3 (21%)	17,19,21	0.76	0
5	MAN	A	475(G)	-	11,11,12	1.80	3 (27%)	15,15,17	2.51	7 (46%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	476(A)	1	14,14,15	1.41	3 (21%)	17,19,21	0.76	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
6	NAG	A	477(A)	1	-	1/6/23/26	0/1/1/1
5	MAN	N	475(G)	-	-	0/2/19/22	0/1/1/1
6	NAG	N	477(A)	1	-	1/6/23/26	0/1/1/1
6	NAG	N	476(A)	1	-	0/6/23/26	0/1/1/1
5	MAN	A	475(G)	-	-	0/2/19/22	0/1/1/1
6	NAG	A	476(A)	1	-	0/6/23/26	0/1/1/1

All (18) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	475(G)	MAN	C4-C5	-4.27	1.43	1.53
5	N	475(G)	MAN	C4-C5	-4.27	1.43	1.53
6	N	476(A)	NAG	C1-C2	-3.05	1.48	1.52
6	A	476(A)	NAG	C1-C2	-3.04	1.48	1.52
6	N	476(A)	NAG	O5-C5	2.78	1.48	1.43
6	A	476(A)	NAG	O5-C5	2.76	1.48	1.43
5	N	475(G)	MAN	O4-C4	-2.29	1.37	1.43
5	A	475(G)	MAN	O4-C4	-2.27	1.37	1.43
6	N	477(A)	NAG	C1-C2	2.22	1.55	1.52
6	A	477(A)	NAG	C1-C2	2.20	1.55	1.52
6	A	477(A)	NAG	C4-C5	2.12	1.57	1.53
6	N	476(A)	NAG	O3-C3	-2.10	1.37	1.43
6	A	476(A)	NAG	O3-C3	-2.10	1.37	1.43
6	N	477(A)	NAG	C4-C5	2.08	1.57	1.53
6	A	477(A)	NAG	C8-C7	2.01	1.54	1.50
5	N	475(G)	MAN	C1-C2	2.01	1.57	1.52
5	A	475(G)	MAN	C1-C2	2.01	1.57	1.52
6	N	477(A)	NAG	C8-C7	2.00	1.54	1.50

All (24) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	475(G)	MAN	C2-C3-C4	4.48	118.74	110.86
5	N	475(G)	MAN	C2-C3-C4	4.47	118.72	110.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	N	475(G)	MAN	O2-C2-C1	4.46	119.44	109.22
5	A	475(G)	MAN	O2-C2-C1	4.46	119.43	109.22
5	N	475(G)	MAN	O3-C3-C4	-4.06	100.80	110.38
5	A	475(G)	MAN	O3-C3-C4	-4.05	100.82	110.38
5	N	475(G)	MAN	C1-O5-C5	3.23	116.52	112.19
5	A	475(G)	MAN	C1-O5-C5	3.22	116.50	112.19
6	A	477(A)	NAG	O5-C1-C2	-3.19	106.36	111.29
6	N	477(A)	NAG	O5-C1-C2	-3.19	106.36	111.29
5	A	475(G)	MAN	O5-C5-C4	-3.01	103.50	110.83
5	N	475(G)	MAN	O5-C5-C4	-3.01	103.51	110.83
5	A	475(G)	MAN	C1-C2-C3	-2.91	105.41	109.64
5	N	475(G)	MAN	C1-C2-C3	-2.88	105.45	109.64
5	A	475(G)	MAN	O5-C5-C6	2.74	113.00	107.66
5	N	475(G)	MAN	O5-C5-C6	2.73	112.98	107.66
6	N	477(A)	NAG	C8-C7-N2	2.45	120.18	116.12
6	A	477(A)	NAG	C8-C7-N2	2.43	120.15	116.12
6	N	477(A)	NAG	O7-C7-C8	-2.42	117.74	122.05
6	A	477(A)	NAG	O7-C7-C8	-2.42	117.74	122.05
6	A	477(A)	NAG	C1-O5-C5	2.21	115.15	112.19
6	N	477(A)	NAG	C1-O5-C5	2.19	115.12	112.19
6	A	477(A)	NAG	O3-C3-C2	2.13	113.83	109.40
6	N	477(A)	NAG	O3-C3-C2	2.12	113.81	109.40

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	N	477(A)	NAG	O5-C5-C6-O6
6	A	477(A)	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	476(A)	NAG	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	388/388 (100%)	-0.46	14 (3%) 46 41	3, 12, 43, 78	0
1	N	388/388 (100%)	-0.50	6 (1%) 72 68	3, 12, 43, 78	0
2	B	122/122 (100%)	0.05	7 (5%) 29 26	13, 27, 50, 80	0
2	H	122/122 (100%)	0.03	3 (2%) 58 54	13, 27, 50, 80	0
3	C	109/109 (100%)	0.62	17 (15%) 5 4	13, 38, 72, 102	0
3	L	109/109 (100%)	0.47	12 (11%) 10 8	13, 38, 72, 102	0
All	All	1238/1238 (100%)	-0.20	59 (4%) 35 31	3, 18, 64, 102	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	342	PRO	6.6
3	C	108	ASP	6.2
1	A	340	PRO	5.9
3	L	109	TYR	5.5
3	C	7	THR	5.4
2	H	113	SER	5.0
3	C	109	TYR	4.7
3	C	76	SER	4.6
1	A	343	GLY	4.6
1	N	345	ASN	4.2
3	L	7	THR	4.1
3	C	1	ASP	4.1
3	L	108	ASP	4.1
2	B	113	SER	3.5
1	A	296	GLN	3.4
3	C	107	ARG	3.4
1	A	249	PRO	3.3
1	A	345	ASN	3.2
3	L	8	THR	3.1

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Mol	Chain	Res	Type	RSRZ
3	C	8	THR	3.1
2	H	105	GLN	3.0
3	C	79	GLU	3.0
3	C	17	ASP	3.0
3	L	12	SER	2.9
3	C	12	SER	2.8
3	L	18	ARG	2.8
1	A	295	TRP	2.8
2	B	64	LYS	2.8
3	L	14	SER	2.7
1	N	344	ASN	2.7
1	A	344	ASN	2.7
3	C	81	GLU	2.7
3	C	14	SER	2.7
3	L	17	ASP	2.6
1	A	297	GLY	2.6
3	L	77	ASN	2.6
1	A	347	ASN	2.6
2	B	65	ASP	2.6
1	A	338	ASN	2.5
1	A	346	ASN	2.5
1	N	341	TYR	2.5
3	L	105	GLU	2.5
2	B	85	GLU	2.5
1	N	416	GLU	2.4
3	C	105	GLU	2.4
2	B	112	SER	2.3
3	L	11	LEU	2.3
2	H	64	LYS	2.2
3	C	39	ASN	2.2
2	B	43	GLN	2.2
3	C	10	SER	2.2
1	N	342	PRO	2.2
1	N	294	ASN	2.2
1	A	298	SER	2.2
3	C	82	ASP	2.1
2	B	61	GLN	2.1
3	L	15	LEU	2.1
3	C	41	ASP	2.1
1	A	341	TYR	2.0

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

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

6.3 Carbohydrates ⓘ

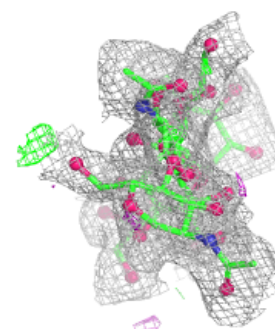
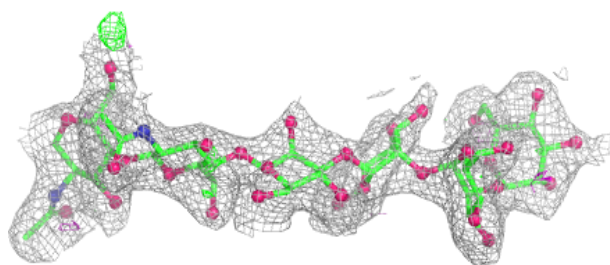
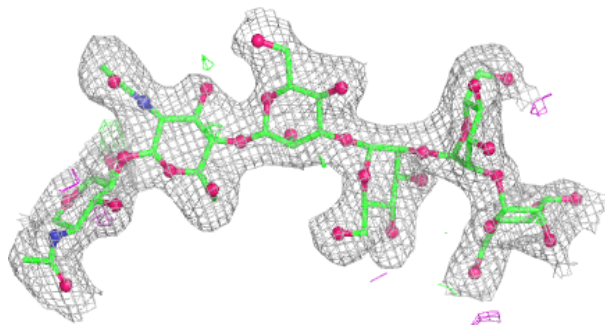
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	D	1	14/15	-	-	5,15,25,29	0
4	NAG	D	2	14/15	-	-	7,12,19,28	0
4	BMA	D	3	11/12	-	-	11,13,17,23	0
4	MAN	D	4	11/12	-	-	12,14,17,24	0
4	MAN	D	5	11/12	-	-	9,16,20,27	0
4	MAN	D	6	11/12	-	-	10,12,15,17	0
4	NAG	E	1	14/15	-	-	5,15,25,29	0
4	NAG	E	2	14/15	-	-	7,12,19,28	0
4	BMA	E	3	11/12	-	-	11,13,17,23	0
4	MAN	E	4	11/12	-	-	12,14,17,24	0
4	MAN	E	5	11/12	-	-	9,16,20,27	0
4	MAN	E	6	11/12	-	-	10,12,15,17	0

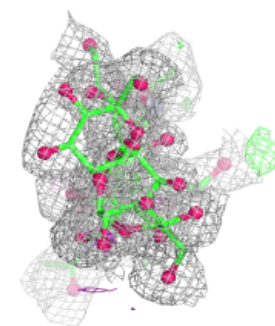
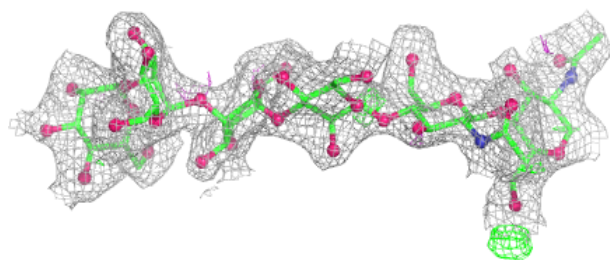
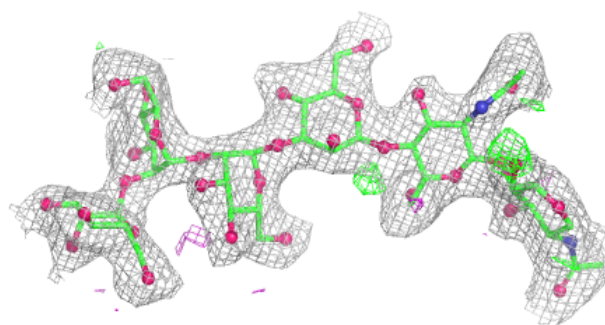
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain 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 E:**

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



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	CA	N	478	1/1	0.59	0.12	65,65,65,65	0
6	NAG	A	477(A)	14/15	0.73	0.12	44,48,51,56	0
6	NAG	N	477(A)	14/15	0.81	0.12	44,48,51,56	0
6	NAG	A	476(A)	14/15	0.81	0.14	40,53,58,64	0
5	MAN	N	475(G)	11/12	0.84	0.11	34,36,40,42	0
6	NAG	N	476(A)	14/15	0.84	0.11	40,53,58,64	0
5	MAN	A	475(G)	11/12	0.88	0.12	34,36,40,42	0
7	CA	A	478	1/1	0.93	0.09	65,65,65,65	0

6.5 Other polymers [i](#)

There are no such residues in this entry.