



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:50 AM UTC

PDB ID : 4P59 / pdb_00004p59
Title : HER3 extracellular domain in complex with Fab fragment of MOR09825
Authors : Sprague, E.R.
Deposited on : 2014-03-15
Resolution : 3.40 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

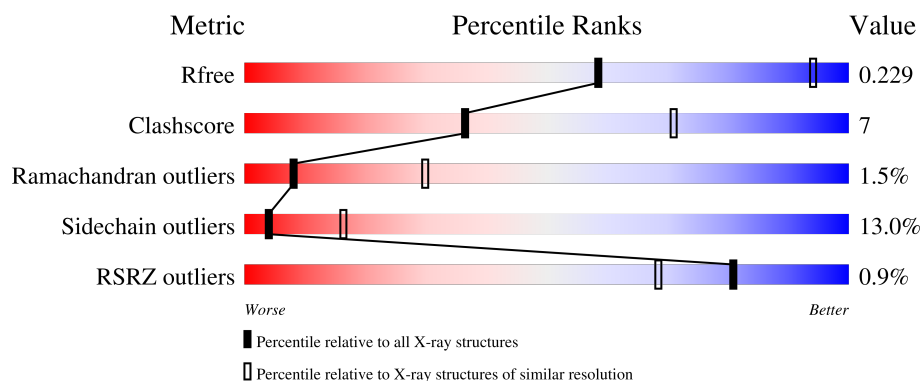
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	627	71% 21% 6%
2	H	238	58% 24% 6% 12%
3	L	214	69% 24% 6%
4	B	2	100%
4	C	2	100%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 7834 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 Receptor tyrosine-protein kinase erbB-3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	592	Total	C	N	O	S	0	0	0
			4547	2813	823	852	59			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	641	GLU	HIS	cloning artifact	UNP P21860
A	642	PHE	LEU	cloning artifact	UNP P21860
A	643	ARG	THR	cloning artifact	UNP P21860
A	644	HIS	MET	cloning artifact	UNP P21860
A	645	ASP	ALA	cloning artifact	UNP P21860
A	646	SER	LEU	cloning artifact	UNP P21860

- Molecule 2 is a protein called MOR09825 Fab fragment heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	210	Total	C	N	O	S	0	0	0
			1568	990	263	309	6			

- Molecule 3 is a protein called MOR09825 Fab fragment light chain.

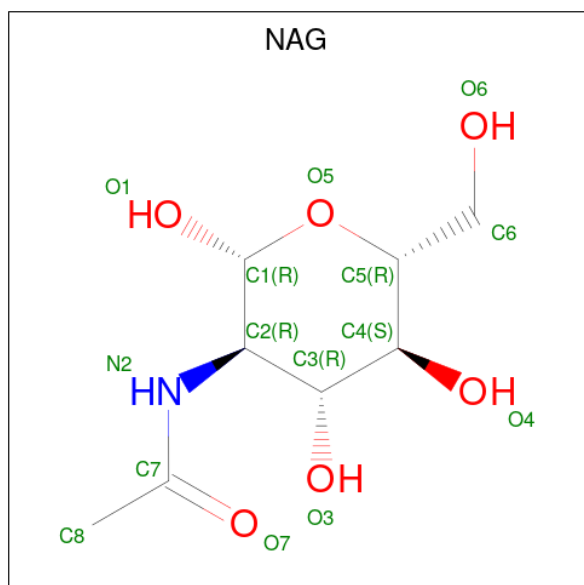
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1618	1015	271	327	5			

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



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

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

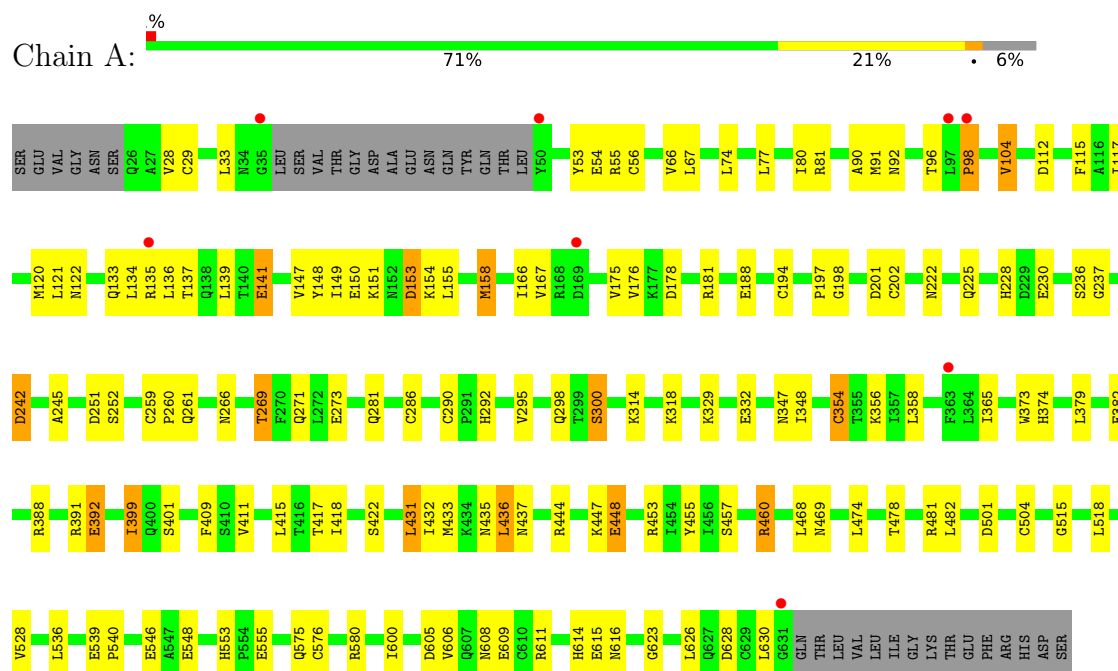
- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	H	2	Total O 2 2	0	0
6	L	1	Total O 1 1	0	0

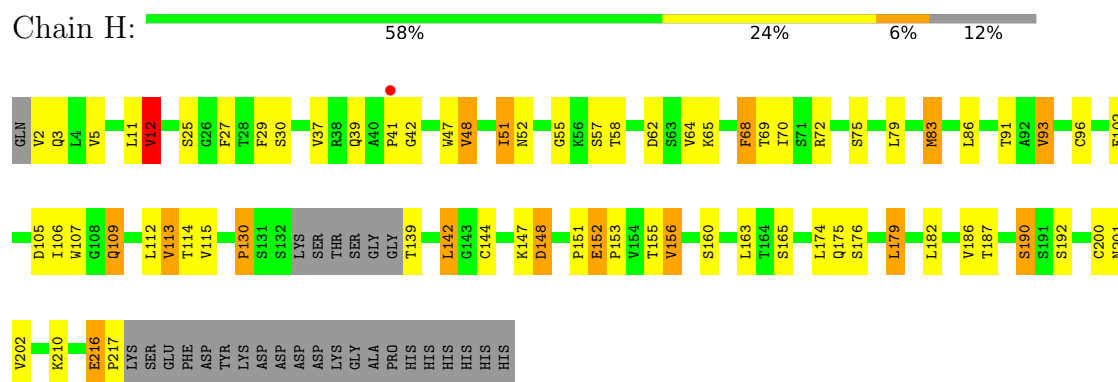
3 Residue-property plots [i](#)

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: Receptor tyrosine-protein kinase erbB-3

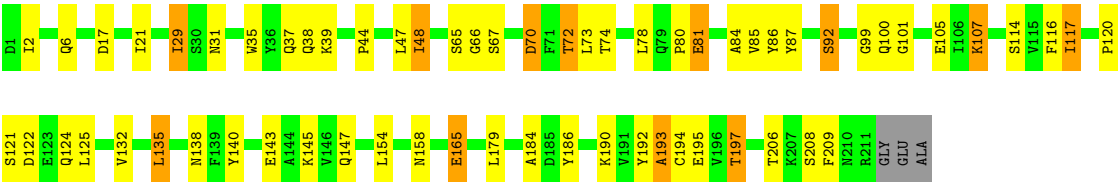


- Molecule 2: MOR09825 Fab fragment heavy chain



- Molecule 3: MOR09825 Fab fragment light chain





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



• Molecule 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	I 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	124.23Å 140.94Å 180.25Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	111.03 – 3.40 111.03 – 3.40	Depositor EDS
% Data completeness (in resolution range)	99.8 (111.03-3.40) 99.9 (111.03-3.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 3.41Å)	Xtriage
Refinement program	BUSTER-TNT BUSTER 2.11.4, BUSTER 2.11.4	Depositor
R, R_{free}	0.184 , 0.250 (Not available) , 0.229	Depositor DCC
R_{free} test set	1132 reflections (5.12%)	wwPDB-VP
Wilson B-factor (Å ²)	76.5	Xtriage
Anisotropy	0.778	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 112.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7834	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.79% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NAG

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	0.79	0/4657	1.32	24/6317 (0.4%)
2	H	0.83	1/1605 (0.1%)	1.30	15/2186 (0.7%)
3	L	0.76	0/1654	1.26	8/2246 (0.4%)
All	All	0.79	1/7916 (0.0%)	1.30	47/10749 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	H	12	VAL	CA-C	-5.71	1.45	1.52

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	236	SER	N-CA-C	-8.98	102.24	113.55
1	A	437	ASN	CA-CB-CG	7.18	119.78	112.60
3	L	66	GLY	N-CA-C	6.73	118.88	112.08
2	H	130	PRO	CA-C-N	6.17	132.81	121.70
2	H	130	PRO	C-N-CA	6.17	132.81	121.70
1	A	55	ARG	CA-C-N	6.00	131.01	120.87
1	A	55	ARG	C-N-CA	6.00	131.01	120.87
1	A	260	PRO	CA-C-N	5.98	129.18	120.51
1	A	260	PRO	C-N-CA	5.98	129.18	120.51
1	A	141	GLU	CA-C-N	5.96	131.11	123.13
1	A	141	GLU	C-N-CA	5.96	131.11	123.13
3	L	80	PRO	CA-C-N	5.74	128.25	120.38
3	L	80	PRO	C-N-CA	5.74	128.25	120.38
2	H	147	LYS	CA-C-N	5.67	132.37	121.54
2	H	147	LYS	C-N-CA	5.67	132.37	121.54
2	H	175	GLN	CA-C-N	5.63	127.83	120.28
2	H	175	GLN	C-N-CA	5.63	127.83	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	409	PHE	CA-C-N	5.62	128.58	120.38
1	A	409	PHE	C-N-CA	5.62	128.58	120.38
1	A	435	ASN	CA-C-N	5.60	128.35	120.28
1	A	435	ASN	C-N-CA	5.60	128.35	120.28
1	A	347	ASN	CA-CB-CG	5.50	118.10	112.60
3	L	70	ASP	CA-CB-CG	5.43	118.03	112.60
2	H	29	PHE	CA-C-N	5.42	128.30	120.38
2	H	29	PHE	C-N-CA	5.42	128.30	120.38
2	H	155	THR	CA-C-N	-5.37	115.95	123.10
2	H	155	THR	C-N-CA	-5.37	115.95	123.10
3	L	17	ASP	CA-CB-CG	5.32	117.92	112.60
1	A	606	VAL	N-CA-C	-5.32	104.45	111.09
2	H	93	VAL	N-CA-C	-5.26	100.83	108.36
2	H	179	LEU	N-CA-C	5.26	118.64	110.17
1	A	197	PRO	CA-C-N	5.21	125.05	120.10
1	A	197	PRO	C-N-CA	5.21	125.05	120.10
2	H	187	THR	CB-CA-C	5.21	118.63	110.19
2	H	105	ASP	CA-CB-CG	5.18	117.78	112.60
1	A	53	TYR	CA-C-N	5.18	130.22	122.86
1	A	53	TYR	C-N-CA	5.18	130.22	122.86
1	A	198	GLY	N-CA-C	5.18	117.88	111.93
2	H	113	VAL	N-CA-CB	5.15	117.67	110.56
1	A	115	PHE	N-CA-C	5.08	117.84	109.72
1	A	136	LEU	CA-C-N	5.07	127.07	120.28
1	A	136	LEU	C-N-CA	5.07	127.07	120.28
3	L	107	LYS	N-CA-C	-5.04	100.94	108.96
3	L	193	ALA	N-CA-C	5.04	116.84	109.24
3	L	85	VAL	N-CA-CB	5.03	117.02	110.99
1	A	374	HIS	CA-C-N	5.01	131.11	121.54
1	A	374	HIS	C-N-CA	5.01	131.11	121.54

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4547	0	4313	53	0
2	H	1568	0	1528	34	0
3	L	1618	0	1575	25	0
4	B	28	0	25	2	0
4	C	28	0	25	0	0
5	A	42	0	39	2	0
6	H	2	0	0	0	0
6	L	1	0	0	0	0
All	All	7834	0	7505	110	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 7.

All (110) 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:H:91:THR:HG22	2:H:115:VAL:H	1.38	0.89
2:H:48:VAL:HG22	2:H:64:VAL:HG21	1.61	0.81
3:L:6:GLN:HE22	3:L:87:TYR:HA	1.50	0.77
1:A:269:THR:HG23	1:A:271:GLN:HG2	1.66	0.76
3:L:145:LYS:HB2	3:L:197:THR:HG23	1.69	0.74
1:A:266:ASN:HB3	1:A:269:THR:HG22	1.73	0.70
2:H:216:GLU:HB3	2:H:217:PRO:HD3	1.73	0.70
5:A:702:NAG:O3	5:A:702:NAG:H82	1.94	0.68
1:A:623:GLY:H	1:A:628:ASP:HB3	1.57	0.67
2:H:51:ILE:HD12	2:H:58:THR:HG22	1.76	0.67
1:A:614:HIS:HD2	1:A:616:ASN:H	1.43	0.66
3:L:37:GLN:HB2	3:L:47:LEU:HD11	1.77	0.66
1:A:251:ASP:HA	1:A:286:CYS:HB2	1.79	0.64
2:H:5:VAL:HA	2:H:109:GLN:HE22	1.61	0.64
2:H:64:VAL:HG13	2:H:68:PHE:CG	2.34	0.62
3:L:147:GLN:HB2	3:L:195:GLU:HB3	1.81	0.61
2:H:37:VAL:HG22	2:H:47:TRP:HA	1.81	0.60
1:A:453:ARG:HG2	1:A:481:ARG:HB2	1.84	0.59
2:H:139:THR:N	2:H:190:SER:HG	2.02	0.58
3:L:116:PHE:HB2	3:L:135:LEU:HB3	1.86	0.58
1:A:158:MET:HB3	1:A:176:VAL:HG11	1.86	0.57
3:L:194:CYS:O	3:L:206:THR:HA	2.03	0.57
2:H:148:ASP:HB3	2:H:179:LEU:HD13	1.86	0.56
1:A:605:ASP:HB2	1:A:609:GLU:H	1.71	0.55
1:A:539:GLU:HB2	1:A:540:PRO:HD3	1.88	0.55
1:A:96:THR:HG22	1:A:133:GLN:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:86:TYR:O	3:L:101:GLY:HA2	2.07	0.54
1:A:365:ILE:H	1:A:365:ILE:HD12	1.73	0.54
2:H:91:THR:CG2	2:H:115:VAL:H	2.14	0.54
2:H:2:VAL:HG12	2:H:27:PHE:HB3	1.90	0.53
1:A:269:THR:CG2	1:A:271:GLN:HG2	2.37	0.53
1:A:431:LEU:HG	1:A:455:TYR:HB3	1.91	0.53
2:H:130:PRO:HG3	2:H:142:LEU:HB3	1.90	0.52
1:A:66:VAL:HG12	1:A:90:ALA:HB3	1.90	0.52
2:H:64:VAL:CG1	2:H:68:PHE:CG	2.93	0.52
2:H:12:VAL:HG21	2:H:86:LEU:HD13	1.92	0.52
2:H:93:VAL:HG22	2:H:112:LEU:HD12	1.92	0.51
1:A:237:GLY:H	1:A:242:ASP:HB3	1.76	0.51
1:A:104:VAL:HG11	1:A:245:ALA:HB2	1.92	0.50
1:A:417:THR:HG23	1:A:448:GLU:HB3	1.93	0.50
1:A:141:GLU:HA	1:A:166:ILE:HG23	1.94	0.50
3:L:117:ILE:HG12	3:L:209:PHE:HD2	1.77	0.49
1:A:281:GLN:HB2	1:A:300:SER:HB3	1.93	0.49
2:H:107:TRP:CE3	3:L:44:PRO:HD2	2.48	0.48
3:L:120:PRO:HD3	3:L:132:VAL:HG22	1.95	0.48
1:A:474:LEU:HD23	1:A:478:THR:HG22	1.95	0.48
1:A:222:ASN:HB2	1:A:225:GLN:HE21	1.77	0.48
1:A:98:PRO:HA	1:A:135:ARG:HB2	1.94	0.48
2:H:52:ASN:HB2	2:H:57:SER:O	2.14	0.47
1:A:134:LEU:HG	1:A:155:LEU:HD11	1.96	0.47
3:L:29:ILE:HG13	3:L:92:SER:HB3	1.95	0.47
1:A:614:HIS:CD2	1:A:616:ASN:H	2.28	0.47
3:L:81:GLU:H	3:L:81:GLU:HG3	1.54	0.47
3:L:39:LYS:HG2	3:L:84:ALA:HB2	1.97	0.46
2:H:64:VAL:HG13	2:H:68:PHE:CD2	2.51	0.46
2:H:139:THR:HA	2:H:190:SER:H	1.80	0.46
1:A:92:ASN:H	1:A:122:ASN:HA	1.81	0.46
2:H:152:GLU:OE1	2:H:153:PRO:HA	2.16	0.46
3:L:165:GLU:CD	3:L:165:GLU:H	2.24	0.46
1:A:273:GLU:OE1	2:H:52:ASN:OD1	2.32	0.46
1:A:295:VAL:HG12	1:A:318:LYS:HB3	1.97	0.46
3:L:21:ILE:HD12	3:L:73:LEU:HD23	1.97	0.46
4:B:1:NAG:H61	4:B:2:NAG:C7	2.46	0.46
1:A:399:ILE:HD12	1:A:432:ILE:HG12	1.99	0.45
1:A:515:GLY:HA2	1:A:528:VAL:HA	1.98	0.45
1:A:149:ILE:HG21	1:A:158:MET:HG2	1.97	0.45
1:A:269:THR:HG21	1:A:273:GLU:CG	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:LYS:HG3	1:A:392:GLU:HB3	1.99	0.44
2:H:91:THR:HG22	2:H:115:VAL:N	2.19	0.44
3:L:107:LYS:HA	3:L:140:TYR:OH	2.18	0.44
1:A:433:MET:HA	1:A:457:SER:O	2.18	0.44
3:L:186:TYR:HA	3:L:192:TYR:OH	2.17	0.44
3:L:2:ILE:HD12	3:L:29:ILE:CD1	2.48	0.44
1:A:77:LEU:HA	1:A:80:ILE:HD12	2.00	0.43
1:A:614:HIS:CD2	1:A:626:LEU:HD13	2.53	0.43
3:L:65:SER:HB3	3:L:72:THR:HG22	1.98	0.43
1:A:120:MET:HG2	1:A:121:LEU:HG	2.00	0.43
1:A:415:LEU:HD21	1:A:418:ILE:HD11	2.00	0.43
1:A:388:ARG:HG2	1:A:411:VAL:HA	1.99	0.43
2:H:51:ILE:HG13	2:H:52:ASN:N	2.32	0.43
1:A:120:MET:HE3	1:A:121:LEU:HD12	2.00	0.43
2:H:39:GLN:HE22	3:L:38:GLN:HE22	1.67	0.43
2:H:68:PHE:N	2:H:68:PHE:CD1	2.87	0.43
1:A:615:GLU:HB3	3:L:31:ASN:ND2	2.34	0.42
3:L:193:ALA:HB2	3:L:208:SER:HB2	2.01	0.42
1:A:553:HIS:HD2	1:A:555:GLU:H	1.67	0.42
3:L:6:GLN:HE21	3:L:99:GLY:HA3	1.84	0.42
3:L:35:TRP:HB2	3:L:48:ILE:HB	2.01	0.42
3:L:122:ASP:HA	3:L:125:LEU:HD12	2.01	0.42
2:H:64:VAL:HG12	2:H:68:PHE:HB2	2.02	0.42
1:A:553:HIS:CD2	1:A:555:GLU:H	2.38	0.42
1:A:576:CYS:SG	1:A:580:ARG:HB2	2.59	0.42
1:A:348:ILE:HD13	1:A:379:LEU:HD11	2.01	0.41
1:A:436:LEU:HD21	5:A:702:NAG:H83	2.03	0.41
2:H:64:VAL:CG1	2:H:68:PHE:HB2	2.49	0.41
2:H:11:LEU:HB2	2:H:151:PRO:HG3	2.03	0.41
2:H:52:ASN:HB3	2:H:55:GLY:H	1.85	0.41
2:H:83:MET:HB2	2:H:86:LEU:HD21	2.02	0.41
1:A:148:TYR:CZ	1:A:150:GLU:HB2	2.55	0.41
1:A:194:CYS:HA	1:A:202:CYS:HA	2.02	0.41
2:H:11:LEU:HD12	2:H:114:THR:O	2.20	0.41
2:H:156:VAL:HB	2:H:202:VAL:HG22	2.02	0.41
1:A:228:HIS:CD2	1:A:230:GLU:H	2.38	0.41
1:A:329:LYS:HB3	1:A:354:CYS:HA	2.03	0.41
1:A:81:ARG:CZ	4:B:1:NAG:H2	2.51	0.41
1:A:74:LEU:HB3	1:A:77:LEU:HD12	2.03	0.41
1:A:222:ASN:H	1:A:225:GLN:HE21	1.67	0.41
2:H:41:PRO:HA	2:H:42:GLY:HA2	1.78	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:ILE:HG13	1:A:401:SER:HB2	2.01	0.40
2:H:64:VAL:CG1	2:H:68:PHE:CD2	3.05	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	588/627 (94%)	520 (88%)	58 (10%)	10 (2%)	7	28
2	H	206/238 (87%)	190 (92%)	13 (6%)	3 (2%)	8	30
3	L	209/214 (98%)	190 (91%)	17 (8%)	2 (1%)	12	40
All	All	1003/1079 (93%)	900 (90%)	88 (9%)	15 (2%)	8	30

All (15) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	H	216	GLU
1	A	444	ARG
2	H	148	ASP
2	H	192	SER
3	L	138	ASN
1	A	153	ASP
1	A	298	GLN
1	A	139	LEU
1	A	167	VAL
1	A	188	GLU
1	A	242	ASP
3	L	184	ALA
1	A	460	ARG
1	A	608	ASN
1	A	98	PRO

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	512/543 (94%)	457 (89%)	55 (11%)	6	23
2	H	175/199 (88%)	140 (80%)	35 (20%)	1	4
3	L	185/186 (100%)	162 (88%)	23 (12%)	4	18
All	All	872/928 (94%)	759 (87%)	113 (13%)	4	16

All (113) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	VAL
1	A	29	CYS
1	A	33	LEU
1	A	54	GLU
1	A	56	CYS
1	A	67	LEU
1	A	91	MET
1	A	104	VAL
1	A	112	ASP
1	A	117	ILE
1	A	137	THR
1	A	147	VAL
1	A	151	LYS
1	A	153	ASP
1	A	154	LYS
1	A	158	MET
1	A	175	VAL
1	A	178	ASP
1	A	181	ARG
1	A	201	ASP
1	A	252	SER
1	A	259	CYS
1	A	261	GLN
1	A	269	THR
1	A	290	CYS
1	A	292	HIS

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Mol	Chain	Res	Type
1	A	300	SER
1	A	314	LYS
1	A	332	GLU
1	A	354	CYS
1	A	358	LEU
1	A	373	TRP
1	A	382	GLU
1	A	391	ARG
1	A	392	GLU
1	A	399	ILE
1	A	422	SER
1	A	431	LEU
1	A	436	LEU
1	A	447	LYS
1	A	448	GLU
1	A	460	ARG
1	A	468	LEU
1	A	469	ASN
1	A	482	LEU
1	A	501	ASP
1	A	504	CYS
1	A	518	LEU
1	A	536	LEU
1	A	546	GLU
1	A	548	GLU
1	A	575	GLN
1	A	600	ILE
1	A	611	ARG
1	A	630	LEU
2	H	3	GLN
2	H	12	VAL
2	H	25	SER
2	H	30	SER
2	H	48	VAL
2	H	51	ILE
2	H	62	ASP
2	H	65	LYS
2	H	68	PHE
2	H	69	THR
2	H	70	ILE
2	H	72	ARG
2	H	75	SER

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Mol	Chain	Res	Type
2	H	79	LEU
2	H	83	MET
2	H	96	CYS
2	H	102	GLU
2	H	106	ILE
2	H	109	GLN
2	H	113	VAL
2	H	142	LEU
2	H	144	CYS
2	H	152	GLU
2	H	156	VAL
2	H	160	SER
2	H	163	LEU
2	H	165	SER
2	H	174	LEU
2	H	176	SER
2	H	182	LEU
2	H	186	VAL
2	H	190	SER
2	H	200	CYS
2	H	201	ASN
2	H	210	LYS
3	L	29	ILE
3	L	48	ILE
3	L	67	SER
3	L	70	ASP
3	L	72	THR
3	L	74	THR
3	L	78	LEU
3	L	81	GLU
3	L	92	SER
3	L	100	GLN
3	L	105	GLU
3	L	114	SER
3	L	117	ILE
3	L	121	SER
3	L	124	GLN
3	L	135	LEU
3	L	143	GLU
3	L	154	LEU
3	L	158	ASN
3	L	165	GLU

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Mol	Chain	Res	Type
3	L	179	LEU
3	L	190	LYS
3	L	197	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	71	ASN
1	A	126	ASN
1	A	133	GLN
1	A	138	GLN
1	A	225	GLN
1	A	228	HIS
1	A	248	HIS
1	A	315	ASN
1	A	487	ASN
1	A	534	ASN
1	A	553	HIS
1	A	566	ASN
1	A	578	HIS
2	H	109	GLN
2	H	168	HIS
2	H	196	GLN
3	L	3	GLN
3	L	6	GLN
3	L	31	ASN
3	L	38	GLN
3	L	79	GLN
3	L	138	ASN
3	L	199	GLN
3	L	210	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

4 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	B	1	4,1	14,14,15	0.40	0	17,19,21	2.57	4 (23%)
4	NAG	B	2	4	14,14,15	0.40	0	17,19,21	2.57	4 (23%)
4	NAG	C	1	4,1	14,14,15	0.47	0	17,19,21	2.57	4 (23%)
4	NAG	C	2	4	14,14,15	0.46	0	17,19,21	0.95	1 (5%)

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	B	1	4,1	-	2/6/23/26	0/1/1/1
4	NAG	B	2	4	-	0/6/23/26	0/1/1/1
4	NAG	C	1	4,1	-	4/6/23/26	0/1/1/1
4	NAG	C	2	4	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	2	NAG	C1-O5-C5	8.33	123.34	112.19
4	C	1	NAG	C1-O5-C5	7.21	121.85	112.19
4	C	1	NAG	C1-C2-N2	5.93	119.78	110.43
4	B	1	NAG	O5-C1-C2	-5.87	102.20	111.29
4	B	1	NAG	C1-O5-C5	5.65	119.75	112.19
4	B	1	NAG	C1-C2-N2	5.18	118.59	110.43
4	B	2	NAG	O5-C1-C2	4.46	118.19	111.29
4	B	2	NAG	C3-C4-C5	3.41	116.41	110.23
4	B	1	NAG	C2-N2-C7	3.05	126.98	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	1	NAG	C2-N2-C7	3.03	126.97	122.90
4	C	2	NAG	C1-C2-N2	-2.62	106.30	110.43
4	B	2	NAG	O5-C5-C4	2.07	115.87	110.83
4	C	1	NAG	C4-C3-C2	-2.04	108.03	111.02

There are no chirality outliers.

All (8) torsion outliers are listed below:

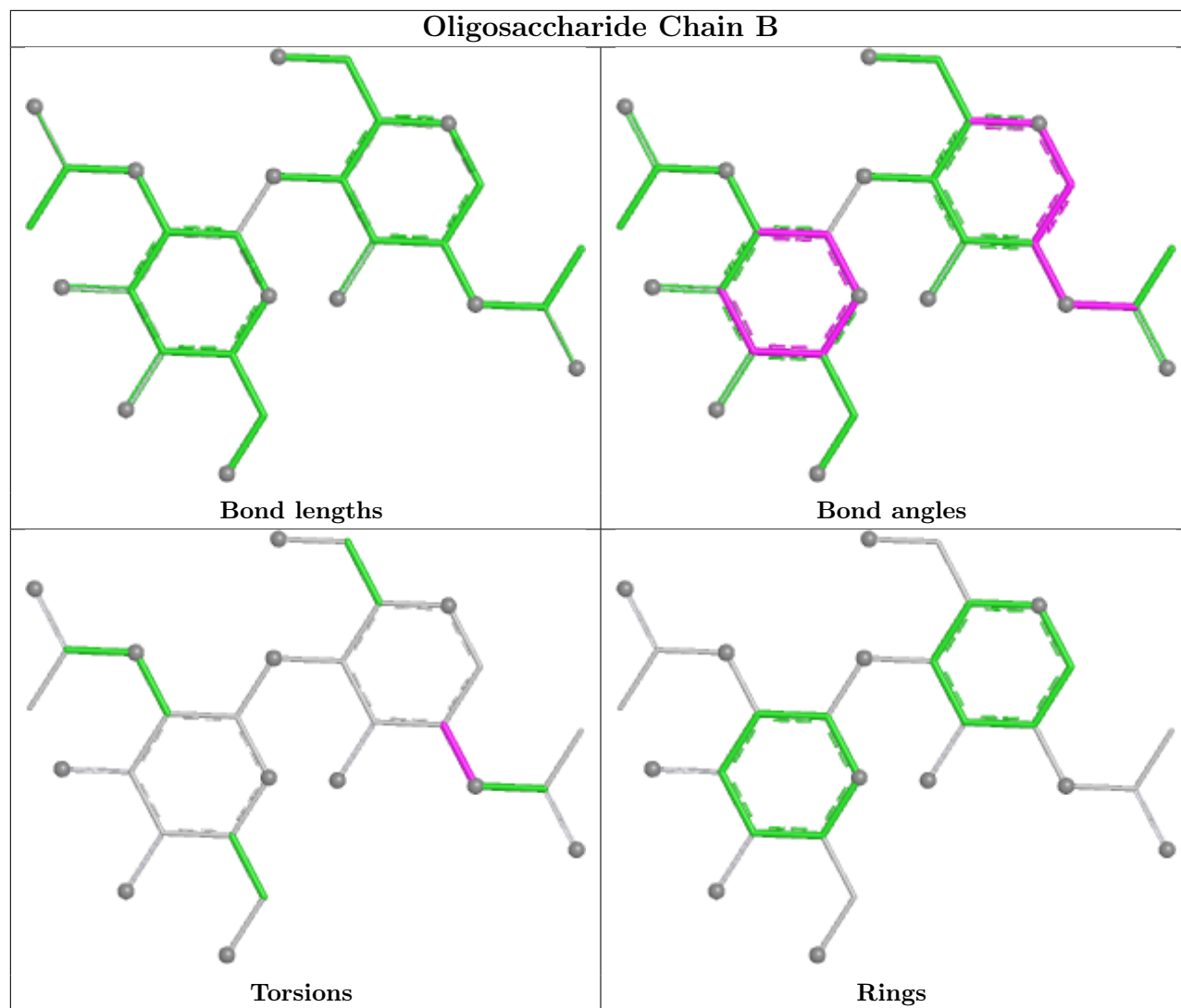
Mol	Chain	Res	Type	Atoms
4	B	1	NAG	C1-C2-N2-C7
4	C	1	NAG	C1-C2-N2-C7
4	C	1	NAG	O5-C5-C6-O6
4	C	1	NAG	C4-C5-C6-O6
4	C	2	NAG	O5-C5-C6-O6
4	B	1	NAG	C3-C2-N2-C7
4	C	1	NAG	C3-C2-N2-C7
4	C	2	NAG	C4-C5-C6-O6

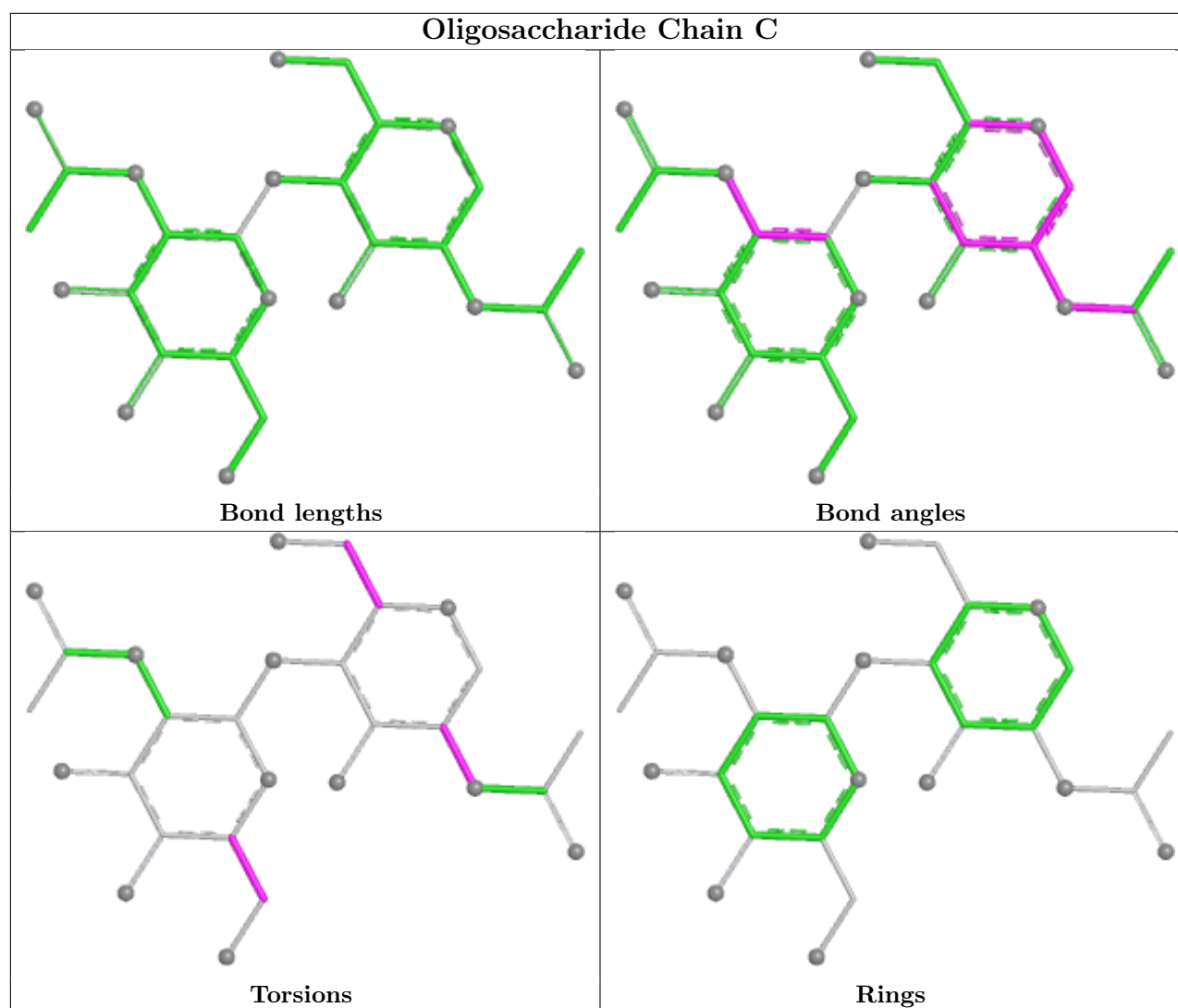
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	2	NAG	1	0
4	B	1	NAG	2	0

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





5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	A	701	1	14,14,15	0.39	0	17,19,21	1.04	1 (5%)
5	NAG	A	702	1	14,14,15	0.78	0	17,19,21	2.73	4 (23%)
5	NAG	A	707	1	14,14,15	0.25	0	17,19,21	0.59	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
5	NAG	A	701	1	-	5/6/23/26	0/1/1/1
5	NAG	A	702	1	-	3/6/23/26	0/1/1/1
5	NAG	A	707	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	702	NAG	C1-O5-C5	-8.74	100.47	112.19
5	A	702	NAG	C1-C2-N2	-4.43	103.45	110.43
5	A	701	NAG	O5-C1-C2	-3.17	106.38	111.29
5	A	702	NAG	C6-C5-C4	3.16	120.77	113.02
5	A	702	NAG	C2-N2-C7	2.72	126.55	122.90

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	702	NAG	C8-C7-N2-C2
5	A	701	NAG	C8-C7-N2-C2
5	A	701	NAG	O7-C7-N2-C2
5	A	702	NAG	O7-C7-N2-C2
5	A	701	NAG	O5-C5-C6-O6
5	A	701	NAG	C4-C5-C6-O6
5	A	701	NAG	C1-C2-N2-C7
5	A	702	NAG	C3-C2-N2-C7

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	702	NAG	2	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	592/627 (94%)	0.06	8 (1%) 73 59	42, 116, 172, 228	0
2	H	210/238 (88%)	-0.44	1 (0%) 87 78	37, 60, 86, 119	0
3	L	211/214 (98%)	-0.40	0 100 100	42, 72, 113, 139	0
All	All	1013/1079 (93%)	-0.14	9 (0%) 81 68	37, 88, 159, 228	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	631	GLY	3.7
1	A	35	GLY	3.4
1	A	363	PHE	2.8
2	H	41	PRO	2.5
1	A	169	ASP	2.4
1	A	50	TYR	2.3
1	A	135	ARG	2.2
1	A	98	PRO	2.0
1	A	97	LEU	2.0

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

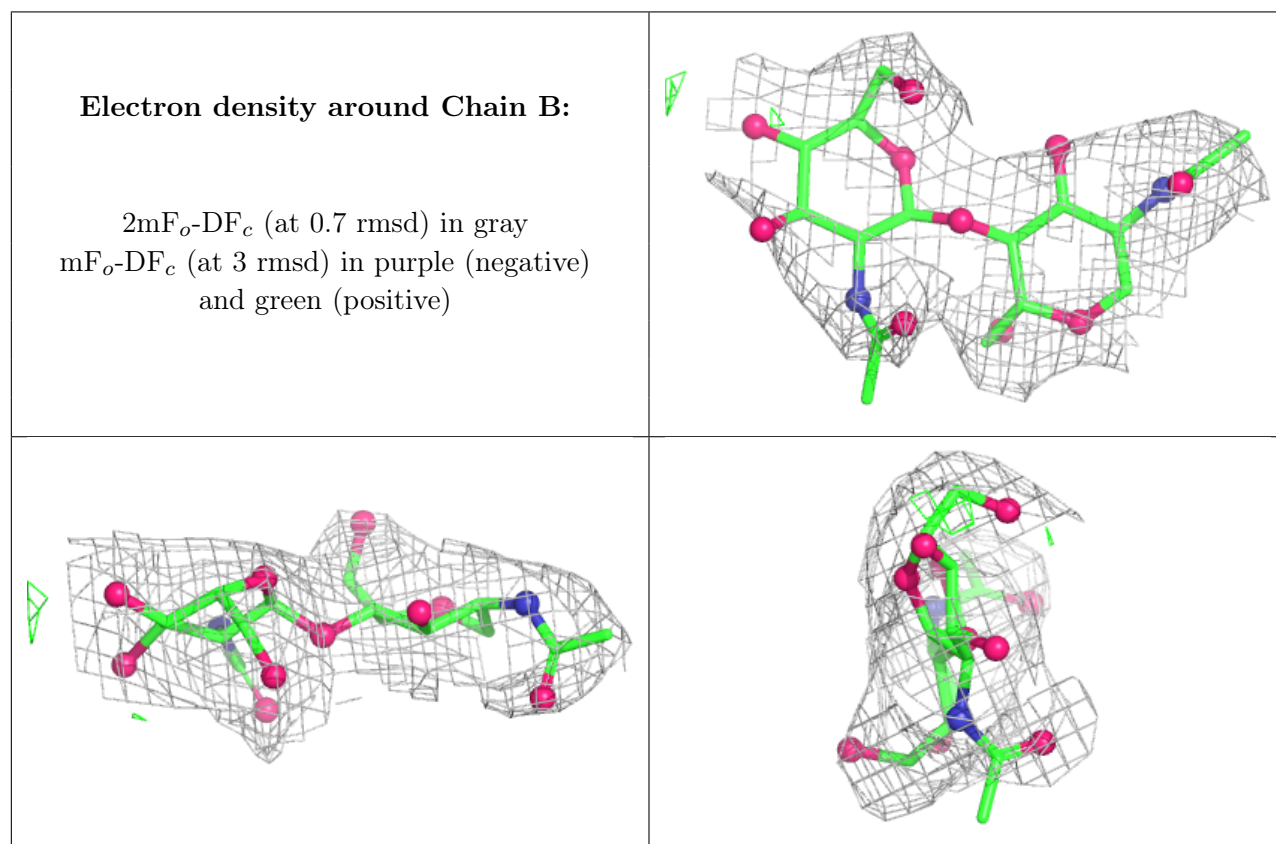
There are no non-standard protein/DNA/RNA residues in this entry.

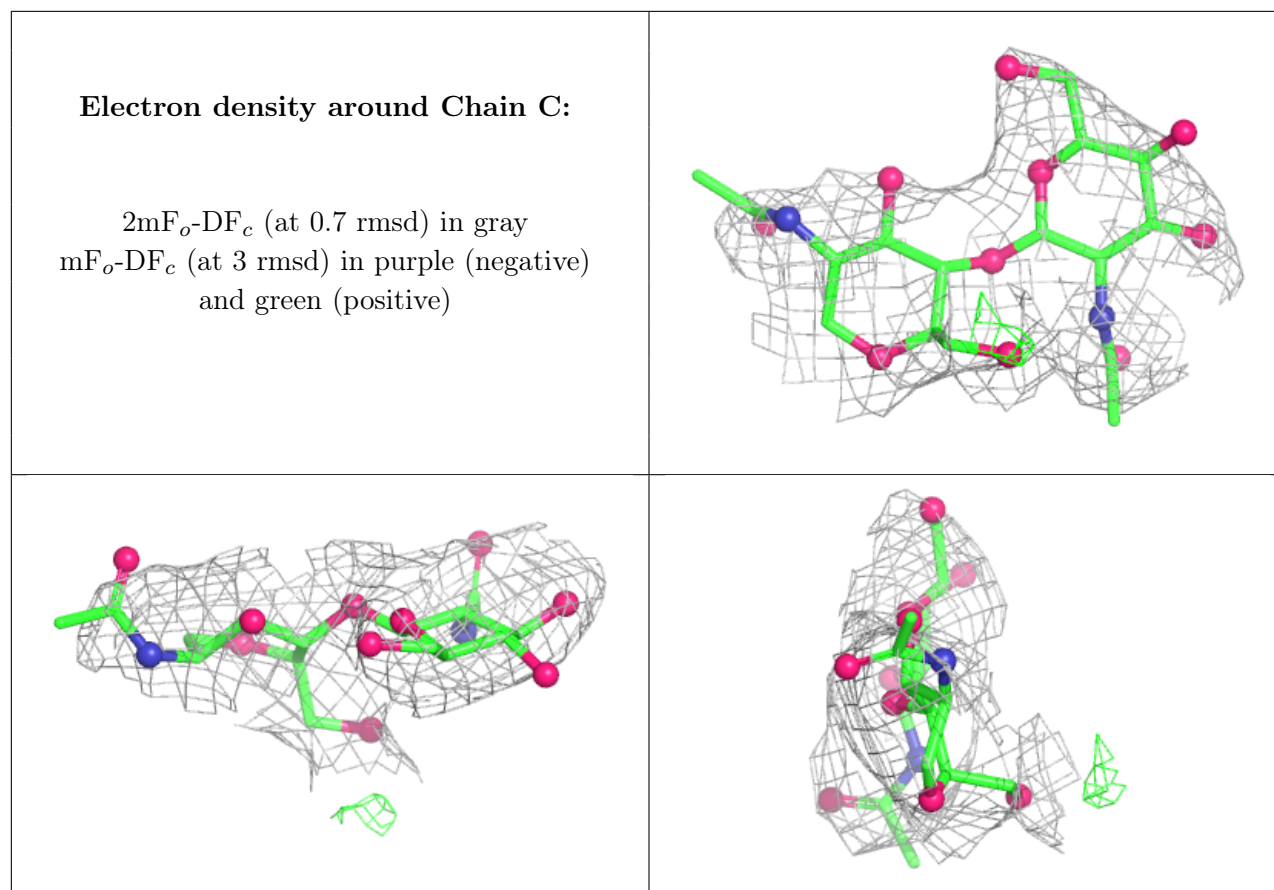
6.3 Carbohydrates [i](#)

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
4	NAG	B	1	14/15	-	-	97,100,104,108	0
4	NAG	B	2	14/15	-	-	107,111,112,113	0
4	NAG	C	1	14/15	-	-	147,149,149,149	0
4	NAG	C	2	14/15	-	-	148,151,151,151	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.





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	A	707	14/15	0.48	0.14	168,171,174,174	0
5	NAG	A	702	14/15	0.71	0.15	106,111,120,120	0
5	NAG	A	701	14/15	0.71	0.13	114,116,120,121	0

6.5 Other polymers [i](#)

There are no such residues in this entry.