



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 02:55 PM UTC

PDB ID : 8SIQ / pdb_00008siq
Title : Crystal structure of SARS-CoV-2 spike receptor-binding domain in complex with broadly neutralizing antibodies CC25.36 and CV38-142 Fab
Authors : Liu, H.; Wilson, I.A.
Deposited on : 2023-04-16
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:

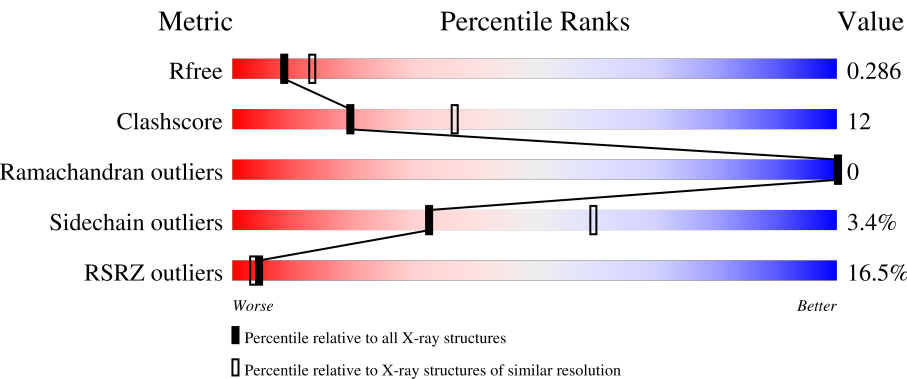
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

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))
R_{free}	180053	5829 (2.50-2.50)
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	205	12% 66% 29% .
2	H	230	26% 69% 26% ..
3	L	217	36% 63% 31% ..
4	M	226	4% 82% 16% .
5	N	217	3% 73% 26%

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Mol	Chain	Length	Quality of chain
6	B	4	50% 50%

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 7963 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 Spike protein S1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	196	Total	C	N	O	S	0	0	0
			1530	979	255	288	8			

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	531	GLY	-	expression tag	UNP P0DTC2
A	532	HIS	-	expression tag	UNP P0DTC2
A	533	HIS	-	expression tag	UNP P0DTC2
A	534	HIS	-	expression tag	UNP P0DTC2
A	535	HIS	-	expression tag	UNP P0DTC2
A	536	HIS	-	expression tag	UNP P0DTC2
A	537	HIS	-	expression tag	UNP P0DTC2

- Molecule 2 is a protein called CC25.36 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	H	220	Total	C	N	O	S	0	0	0
			1591	1003	272	309	7			

- Molecule 3 is a protein called CC25.36 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	210	Total	C	N	O	S	0	0	0
			1474	921	246	303	4			

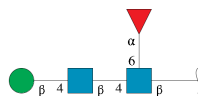
- Molecule 4 is a protein called CV38-142 Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	M	222	Total	C	N	O	S	0	0	0
			1665	1058	273	326	8			

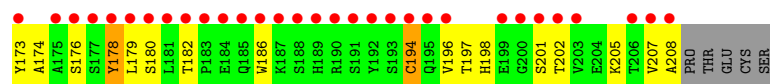
- Molecule 5 is a protein called CV38-142 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	N	216	Total	C	N	O	S	0	0	0
			1654	1031	278	339	6			

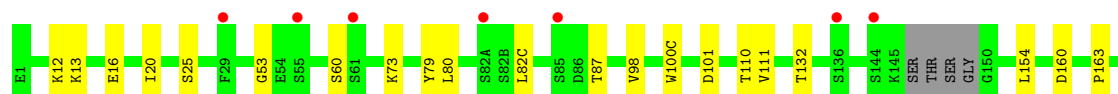
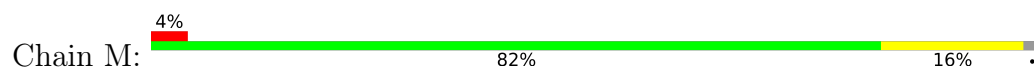
- Molecule 6 is an oligosaccharide called beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



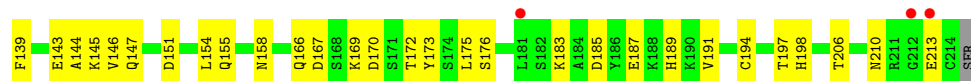
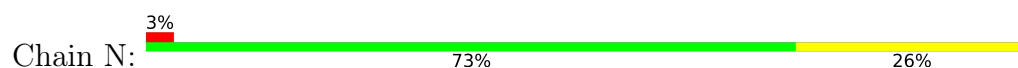
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
6	B	4	Total	C	N	O	0	0	0
			49	28	2	19			



• Molecule 4: CV38-142 Fab heavy chain



• Molecule 5: CV38-142 Fab light chain



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	60.40Å 77.03Å 266.44Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.52 – 2.50 38.52 – 2.50	Depositor EDS
% Data completeness (in resolution range)	77.8 (38.52-2.50) 77.7 (38.52-2.50)	Depositor EDS
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.86 (at 2.51Å)	Xtriage
Refinement program	PHENIX (1.19.2_4158: ???)	Depositor
R, R_{free}	0.247 , 0.287 0.247 , 0.286	Depositor DCC
R_{free} test set	1635 reflections (3.70%)	wwPDB-VP
Wilson B-factor (Å ²)	41.3	Xtriage
Anisotropy	0.050	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 40.0	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	0.89	EDS
Total number of atoms	7963	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.14	0/1573	0.40	0/2142
2	H	0.14	0/1631	0.39	0/2228
3	L	0.37	1/1510 (0.1%)	0.57	2/2074 (0.1%)
4	M	0.12	0/1709	0.37	0/2326
5	N	0.13	0/1690	0.39	0/2297
All	All	0.20	1/8113 (0.0%)	0.43	2/11067 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	L	142	PRO	CG-CD	-10.69	1.22	1.51

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	L	142	PRO	N-CD-CG	-9.33	92.60	103.80
3	L	142	PRO	CA-CB-CG	-5.76	93.06	104.00

There are no chirality outliers.

There are no planarity outliers.

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	1530	0	1422	38	0
2	H	1591	0	1448	44	0
3	L	1474	0	1337	51	0
4	M	1665	0	1621	25	0
5	N	1654	0	1590	32	1
6	B	49	0	43	1	0
All	All	7963	0	7461	179	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (179) 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)
3:L:80:ALA:HA	3:L:106:VAL:HG21	1.49	0.94
3:L:14:ALA:N	3:L:17:GLN:OE1	2.16	0.77
1:A:439:ASN:HD21	1:A:499:PRO:HA	1.47	0.77
1:A:455:LEU:HD11	1:A:493:GLN:HB2	1.65	0.77
5:N:155:GLN:HB3	5:N:158:ASN:HD21	1.49	0.76
1:A:360:ASN:H	1:A:523:THR:HB	1.51	0.75
2:H:18:LEU:HD22	2:H:82(C):LEU:HD11	1.66	0.75
2:H:200:HIS:HD2	2:H:202:PRO:HD2	1.53	0.73
1:A:406:GLU:HA	3:L:53:ASN:HD21	1.54	0.72
4:M:211:ILE:HG12	4:M:226:LYS:HG2	1.73	0.70
4:M:60:SER:HA	5:N:95(A):ARG:HA	1.75	0.68
5:N:147:GLN:HG3	5:N:154:LEU:HD21	1.74	0.68
2:H:87:THR:HG23	2:H:110:THR:HA	1.78	0.66
2:H:48:VAL:HG13	2:H:63:VAL:HG11	1.77	0.66
5:N:183:LYS:HG3	5:N:187:GLU:OE2	1.95	0.66
3:L:132:THR:HA	3:L:180:SER:CB	2.28	0.63
3:L:135:CYS:HB2	3:L:149:TRP:CZ2	2.35	0.62
3:L:83:GLU:OE2	3:L:105:THR:HG22	2.00	0.62
2:H:18:LEU:HD23	2:H:82:MET:HB2	1.80	0.62
2:H:121:VAL:HG22	2:H:142:VAL:HG12	1.82	0.62
2:H:200:HIS:CD2	2:H:202:PRO:HD2	2.35	0.61
3:L:186:TRP:CH2	3:L:208:ALA:HA	2.35	0.61
2:H:144:ASP:HA	2:H:175:LEU:HB3	1.83	0.61
1:A:366:SER:HA	1:A:369:TYR:CE1	2.36	0.60
1:A:357:ARG:HG3	1:A:396:TYR:CE1	2.37	0.58
2:H:5:VAL:O	2:H:22:CYS:HA	2.03	0.58
1:A:420:ASP:HB3	1:A:460:ASN:OD1	2.03	0.58
5:N:170:ASP:OD2	5:N:172:THR:N	2.36	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:N:35:TRP:HB2	5:N:48:ILE:HB	1.85	0.58
2:H:119:PRO:HD3	2:H:200:HIS:ND1	2.19	0.58
1:A:433:VAL:HG22	1:A:512:VAL:HG22	1.86	0.57
5:N:170:ASP:OD2	5:N:170:ASP:C	2.47	0.57
5:N:166:GLN:HG3	5:N:173:TYR:CZ	2.40	0.56
2:H:82:MET:HB3	2:H:82(C):LEU:HD21	1.87	0.56
1:A:363:ALA:O	1:A:527:PRO:HD3	2.05	0.56
2:H:35:ASN:CG	2:H:100(I):ASN:HD21	2.13	0.56
1:A:388:ASN:O	1:A:526:GLY:HA3	2.05	0.55
1:A:342:PHE:HB2	6:B:1:NAG:H82	1.89	0.55
1:A:376:THR:O	1:A:434:ILE:HA	2.07	0.55
2:H:123:PRO:HD3	2:H:209:LYS:HE2	1.89	0.54
3:L:12:SER:HB2	3:L:107:LEU:HD21	1.89	0.54
4:M:179:VAL:HG22	4:M:198:VAL:HB	1.90	0.54
3:L:35:TRP:HB2	3:L:48:ILE:HB	1.90	0.53
5:N:136:LEU:HD13	5:N:175:LEU:HD22	1.91	0.53
5:N:144:ALA:C	5:N:145:LYS:HD2	2.34	0.53
1:A:394:ASN:OD1	1:A:516:GLU:HB2	2.09	0.52
4:M:132:THR:HG22	4:M:163:PRO:HD3	1.91	0.52
1:A:462:LYS:HE3	1:A:465:GLU:CD	2.35	0.51
3:L:106:VAL:O	3:L:141:TYR:OH	2.28	0.51
1:A:376:THR:HG23	1:A:378:LYS:HE2	1.93	0.51
4:M:168:VAL:HG22	4:M:214:VAL:HG22	1.92	0.51
5:N:120:PRO:HG3	5:N:132:VAL:HG22	1.92	0.51
5:N:155:GLN:HB3	5:N:158:ASN:ND2	2.22	0.51
3:L:179:LEU:HD13	3:L:180:SER:N	2.26	0.51
1:A:490:PHE:CD1	1:A:491:PRO:HD2	2.46	0.50
3:L:163:THR:HG22	3:L:164:THR:O	2.11	0.50
4:M:53:GLY:HA2	4:M:73:LYS:HD2	1.93	0.50
3:L:140:PHE:HB2	3:L:198:HIS:CD2	2.46	0.50
2:H:66:ARG:HB3	2:H:82(A):ASN:O	2.12	0.50
3:L:168:GLN:HB2	3:L:170:ASN:OD1	2.12	0.50
1:A:350:VAL:HG22	1:A:422:ASN:HB3	1.92	0.50
4:M:12:LYS:O	4:M:111:VAL:HA	2.11	0.50
3:L:14:ALA:HA	3:L:107:LEU:HB2	1.94	0.49
4:M:170:TRP:CH2	4:M:212:CYS:HB3	2.48	0.49
4:M:20:ILE:HD11	4:M:80:LEU:HD23	1.96	0.48
2:H:163:VAL:HA	2:H:182:VAL:HA	1.95	0.48
4:M:175:LEU:HD21	4:M:198:VAL:HG21	1.95	0.48
2:H:138:LEU:HD11	2:H:194:TYR:CD2	2.49	0.48
3:L:13:GLY:O	3:L:106:VAL:HA	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:37:VAL:HG22	2:H:47:TRP:HA	1.96	0.47
2:H:35:ASN:OD1	2:H:100(I):ASN:ND2	2.42	0.47
3:L:37:GLN:HB2	3:L:86:TYR:CE2	2.49	0.47
3:L:83:GLU:HG3	3:L:104:LEU:O	2.14	0.47
4:M:154:LEU:HG	4:M:198:VAL:HG13	1.95	0.47
4:M:87:THR:HG23	4:M:110:THR:HA	1.96	0.47
1:A:354:ASN:O	1:A:398:ASP:HA	2.13	0.47
2:H:50:TYR:CD1	2:H:58:TYR:HD2	2.32	0.47
4:M:100(C):TRP:CG	5:N:96:TRP:HZ2	2.33	0.47
4:M:178:GLY:O	4:M:198:VAL:HA	2.14	0.47
5:N:35:TRP:CE2	5:N:73:LEU:HB2	2.49	0.47
5:N:108:ARG:HD2	5:N:170:ASP:O	2.15	0.47
5:N:151:ASP:OD1	5:N:189:HIS:HB3	2.15	0.47
3:L:18:ARG:HG3	3:L:76:THR:HG22	1.95	0.47
4:M:215:ASN:ND2	4:M:222:LYS:HG3	2.29	0.47
2:H:67:PHE:CE2	2:H:82:MET:HG2	2.49	0.47
2:H:94:ARG:HB3	2:H:102:ALA:HB3	1.97	0.47
4:M:20:ILE:O	4:M:79:TYR:HA	2.14	0.47
4:M:101:ASP:OD1	4:M:101:ASP:N	2.48	0.47
1:A:462:LYS:HE3	1:A:465:GLU:OE2	2.14	0.47
2:H:24:ALA:HB1	2:H:27:PHE:CE1	2.50	0.47
3:L:123:SER:HA	3:L:126:LEU:HB2	1.97	0.47
2:H:12:VAL:O	2:H:111:VAL:HA	2.15	0.46
4:M:100(C):TRP:HB2	5:N:91:SER:HB2	1.97	0.46
5:N:8:PRO:O	5:N:102:THR:OG1	2.27	0.46
5:N:167:ASP:OD1	5:N:169:LYS:N	2.43	0.46
3:L:186:TRP:HH2	3:L:208:ALA:HA	1.78	0.46
5:N:143:GLU:O	5:N:198:HIS:HD2	1.98	0.46
3:L:196:VAL:O	3:L:202:THR:HA	2.16	0.46
5:N:105:GLU:OE1	5:N:173:TYR:OH	2.33	0.46
1:A:393:THR:HA	1:A:522:ALA:HA	1.98	0.46
3:L:107:LEU:HD23	3:L:107:LEU:HA	1.73	0.46
5:N:210:ASN:HB2	5:N:213:GLU:HB2	1.98	0.46
3:L:176:SER:HB2	3:L:178:TYR:CE1	2.50	0.46
4:M:216:HIS:CD2	4:M:218:PRO:HD2	2.51	0.46
2:H:70:SER:OG	2:H:79:TYR:HB2	2.16	0.45
2:H:39:GLN:HB2	2:H:45:LEU:HD23	1.97	0.45
2:H:67:PHE:CZ	2:H:82:MET:HG2	2.52	0.45
1:A:401:VAL:HG22	1:A:509:ARG:HG2	1.97	0.45
2:H:32:ARG:HG3	2:H:94:ARG:HD2	1.98	0.45
2:H:122:PHE:O	2:H:141:LEU:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:160:VAL:HA	3:L:179:LEU:CD2	2.47	0.45
5:N:37:GLN:HB2	5:N:47:LEU:HD11	1.99	0.45
1:A:358:ILE:HB	1:A:395:VAL:HB	1.98	0.45
3:L:116:VAL:O	3:L:205:LYS:HG3	2.16	0.45
2:H:13:GLN:HE21	2:H:13:GLN:HB3	1.59	0.45
3:L:37:GLN:HB3	3:L:47:LEU:HG	1.98	0.45
5:N:147:GLN:CG	5:N:154:LEU:HD21	2.46	0.45
3:L:163:THR:HB	3:L:176:SER:O	2.17	0.44
3:L:5:THR:HB	3:L:24:THR:OG1	2.17	0.44
3:L:141:TYR:HA	3:L:142:PRO:C	2.42	0.44
1:A:365:TYR:CD2	1:A:365:TYR:N	2.85	0.44
1:A:425:LEU:HD21	1:A:512:VAL:HG11	2.00	0.44
2:H:119:PRO:HA	2:H:145:TYR:HB3	1.98	0.44
2:H:167:PRO:HB2	3:L:163:THR:HG21	1.99	0.44
1:A:380:TYR:CE2	2:H:99:TYR:HB3	2.52	0.44
1:A:438:SER:HB3	1:A:509:ARG:HG3	2.00	0.44
2:H:67:PHE:HA	2:H:81:GLN:O	2.18	0.44
3:L:140:PHE:CE1	3:L:174:ALA:HA	2.53	0.44
2:H:169:VAL:O	2:H:176:TYR:HA	2.18	0.43
2:H:94:ARG:NH2	2:H:101:ASP:OD2	2.42	0.43
3:L:79:GLN:O	3:L:106:VAL:HG11	2.18	0.43
3:L:143:GLY:HA3	3:L:173:TYR:CG	2.54	0.43
2:H:201:LYS:HB3	2:H:202:PRO:HD3	2.00	0.43
1:A:334:ASN:N	1:A:334:ASN:OD1	2.52	0.43
3:L:116:VAL:HA	3:L:136:LEU:O	2.18	0.43
3:L:3:VAL:HG12	3:L:27(B):ASN:ND2	2.33	0.43
5:N:118:PHE:HA	5:N:119:PRO:HD3	1.90	0.43
5:N:194:CYS:O	5:N:206:THR:HA	2.19	0.42
1:A:439:ASN:HA	1:A:507:PRO:HG2	2.01	0.42
1:A:340:GLU:O	1:A:344:ALA:HB2	2.19	0.42
3:L:54:ARG:NE	3:L:62:PHE:O	2.46	0.42
4:M:13:LYS:HB2	4:M:16:GLU:OE1	2.20	0.42
3:L:4:LEU:HD11	3:L:90:SER:HB3	2.01	0.42
3:L:146:THR:O	3:L:196:VAL:HA	2.19	0.42
4:M:160:ASP:HA	4:M:191:LEU:HB3	2.01	0.42
2:H:18:LEU:CD2	2:H:82(C):LEU:HD11	2.44	0.42
3:L:123:SER:O	3:L:127:GLN:N	2.37	0.42
1:A:406:GLU:HA	3:L:53:ASN:ND2	2.28	0.42
2:H:50:TYR:HD1	2:H:58:TYR:HD2	1.67	0.42
2:H:154:TRP:CZ3	2:H:196:CYS:HB3	2.55	0.42
3:L:62:PHE:CD1	3:L:75:ILE:HG12	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:118:LEU:HD12	3:L:194:CYS:HB3	2.02	0.42
5:N:191:VAL:HG22	5:N:210:ASN:OD1	2.19	0.42
2:H:166:PHE:CE2	3:L:136:LEU:HD13	2.55	0.41
1:A:408:ARG:HG3	2:H:100(J):TRP:CH2	2.55	0.41
4:M:82(C):LEU:HD23	4:M:82(C):LEU:HA	1.91	0.41
5:N:151:ASP:OD1	5:N:189:HIS:CD2	2.74	0.41
3:L:197:THR:HA	3:L:201:SER:O	2.21	0.41
1:A:335:LEU:HA	1:A:362:VAL:O	2.21	0.41
1:A:398:ASP:OD2	1:A:423:TYR:OH	2.31	0.41
1:A:452:LEU:HD23	1:A:492:LEU:HB3	2.01	0.41
2:H:148:GLU:OE1	2:H:150:VAL:HG22	2.21	0.41
2:H:154:TRP:CH2	2:H:196:CYS:HB3	2.56	0.41
3:L:179:LEU:HD22	3:L:179:LEU:HA	1.53	0.41
3:L:39:LEU:HD12	3:L:84:ALA:HB2	2.02	0.41
3:L:131:ALA:HB3	3:L:182:THR:HA	2.02	0.41
3:L:34:HIS:HD2	3:L:50:GLY:H	1.69	0.41
3:L:207:VAL:O	3:L:208:ALA:HB2	2.20	0.41
4:M:170:TRP:CZ3	4:M:212:CYS:HB3	2.56	0.41
1:A:345:THR:HA	4:M:98:VAL:HG22	2.03	0.41
1:A:502:GLY:O	1:A:506:GLN:HG3	2.20	0.41
2:H:121:VAL:HA	2:H:141:LEU:O	2.21	0.41
4:M:182:PHE:CE2	5:N:176:SER:HB3	2.56	0.41
3:L:34:HIS:CD2	3:L:50:GLY:H	2.40	0.40
5:N:89:GLN:HB2	5:N:98:PHE:CD2	2.56	0.40
5:N:95:PRO:HA	5:N:96:TRP:CD2	2.56	0.40
5:N:113:PRO:HA	5:N:139:PHE:HB3	2.03	0.40
1:A:336:CYS:SG	1:A:363:ALA:HB2	2.61	0.40
1:A:425:LEU:HD23	1:A:425:LEU:HA	1.91	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:N:60:SER:OG	5:N:185:ASP:OD2[3_455]	2.19	0.01

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	194/205 (95%)	191 (98%)	3 (2%)	0	100	100
2	H	216/230 (94%)	216 (100%)	0	0	100	100
3	L	206/217 (95%)	200 (97%)	6 (3%)	0	100	100
4	M	218/226 (96%)	217 (100%)	1 (0%)	0	100	100
5	N	214/217 (99%)	210 (98%)	4 (2%)	0	100	100
All	All	1048/1095 (96%)	1034 (99%)	14 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	162/177 (92%)	157 (97%)	5 (3%)	35	62
2	H	159/194 (82%)	154 (97%)	5 (3%)	35	62
3	L	150/180 (83%)	142 (95%)	8 (5%)	20	42
4	M	186/190 (98%)	184 (99%)	2 (1%)	65	84
5	N	189/192 (98%)	180 (95%)	9 (5%)	23	46
All	All	846/933 (91%)	817 (97%)	29 (3%)	32	60

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	399	SER
1	A	455	LEU
1	A	469	SER
1	A	477	SER
1	A	525	CYS
2	H	70	SER
2	H	107	THR
2	H	115	SER
2	H	142	VAL
2	H	153	SER
3	L	24	THR
3	L	63	SER
3	L	70	SER
3	L	125	GLU
3	L	146	THR
3	L	162	THR
3	L	178	TYR
3	L	194	CYS
4	M	25	SER
4	M	198	VAL
5	N	7	SER
5	N	10	SER
5	N	12	SER
5	N	30	SER
5	N	48	ILE
5	N	67	SER
5	N	121	SER
5	N	146	VAL
5	N	197	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	334	ASN
1	A	360	ASN
1	A	450	ASN
2	H	13	GLN
2	H	81	GLN
3	L	53	ASN
3	L	168	GLN
4	M	3	GLN
4	M	215	ASN
4	M	220	ASN

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Mol	Chain	Res	Type
5	N	34	ASN
5	N	155	GLN
5	N	158	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	$\# Z > 2$
6	NAG	B	1	1,6	14,14,15	0.33	0	17,19,21	0.61	0
6	NAG	B	2	6	14,14,15	0.31	0	17,19,21	0.45	0
6	BMA	B	3	6	11,11,12	0.64	0	15,15,17	0.65	0
6	FUC	B	4	6	10,10,11	0.76	0	14,14,16	0.98	2 (14%)

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	B	1	1,6	-	0/6/23/26	0/1/1/1
6	NAG	B	2	6	-	0/6/23/26	0/1/1/1
6	BMA	B	3	6	-	0/2/19/22	0/1/1/1
6	FUC	B	4	6	-	-	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	4	FUC	O5-C5-C4	2.32	113.72	109.55
6	B	4	FUC	C1-O5-C5	2.00	117.69	112.97

There are no chirality outliers.

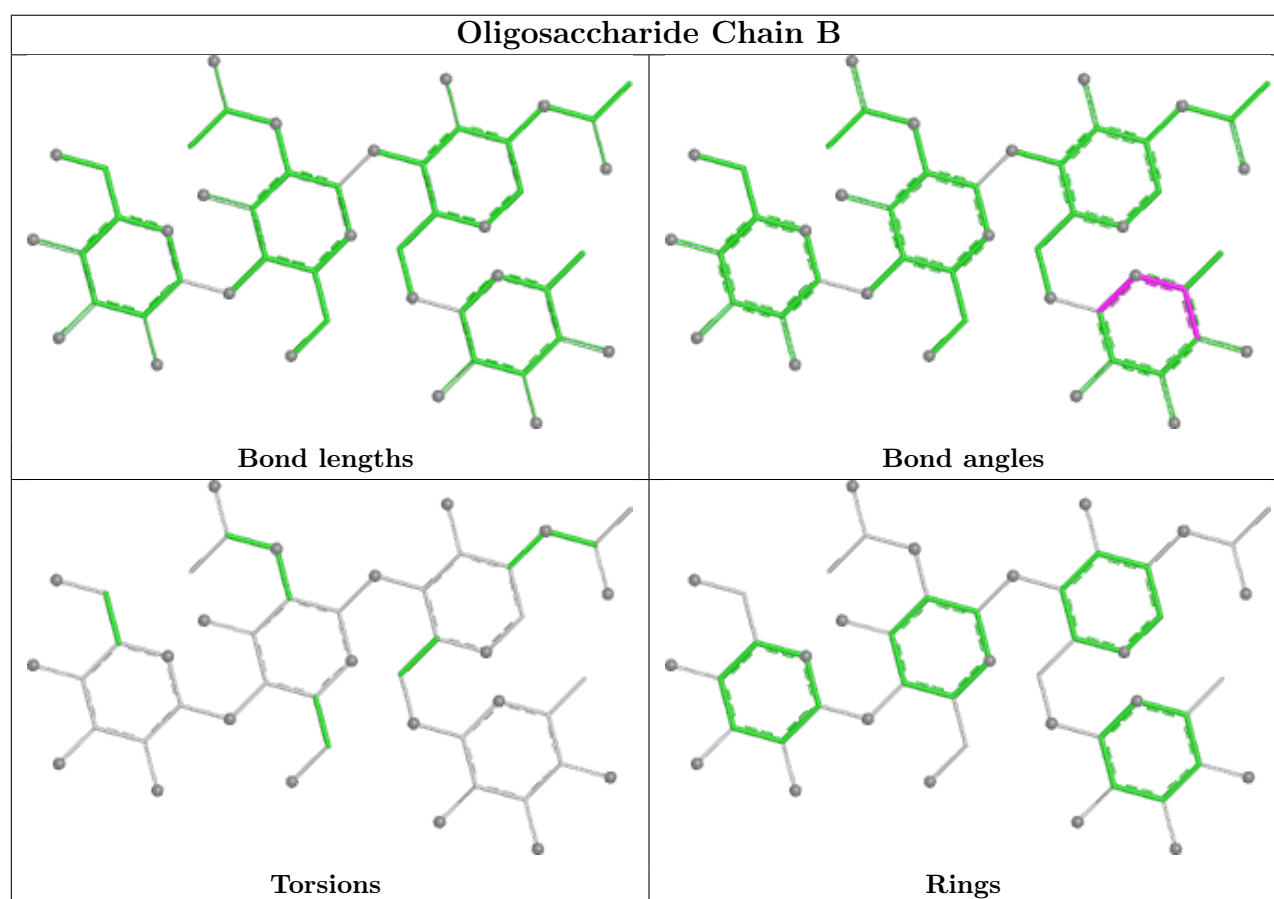
There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	1	NAG	1	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

There are no ligands in this entry.

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	196/205 (95%)	0.82	24 (12%) 8 7	25, 34, 99, 189	0
2	H	220/230 (95%)	1.31	59 (26%) 1 1	24, 39, 108, 177	0
3	L	210/217 (96%)	1.74	78 (37%) 1 1	24, 60, 118, 182	0
4	M	222/226 (98%)	0.58	8 (3%) 46 41	27, 43, 61, 90	0
5	N	216/217 (99%)	0.55	7 (3%) 50 46	28, 44, 62, 206	0
All	All	1064/1095 (97%)	1.00	176 (16%) 4 3	24, 44, 103, 206	0

All (176) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	L	194	CYS	7.8
5	N	212	GLY	6.4
2	H	125	ALA	6.0
3	L	108	GLY	6.0
3	L	134	VAL	6.0
2	H	141	LEU	5.9
2	H	120	SER	5.9
2	H	202	PRO	5.7
3	L	119	PHE	5.4
3	L	120	PRO	5.3
1	A	384	PRO	5.3
3	L	118	LEU	5.2
1	A	373	SER	5.2
2	H	127	SER	5.0
3	L	207	VAL	5.0
2	H	194	TYR	5.0
1	A	527	PRO	5.0
2	H	195	ILE	4.9
3	L	130	LYS	4.8
3	L	193	SER	4.8

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Mol	Chain	Res	Type	RSRZ
3	L	179	LEU	4.8
1	A	517	LEU	4.8
3	L	137	ILE	4.7
2	H	188	SER	4.5
3	L	175	ALA	4.5
3	L	196	VAL	4.5
2	H	138	LEU	4.5
3	L	181	LEU	4.5
3	L	133	LEU	4.4
2	H	154	TRP	4.4
2	H	140	CYS	4.4
3	L	149	TRP	4.3
2	H	207	VAL	4.3
3	L	147	VAL	4.3
3	L	148	ALA	4.3
3	L	183	PRO	4.2
3	L	180	SER	4.2
2	H	193	THR	4.2
2	H	208	ASP	4.1
2	H	198	VAL	4.1
3	L	185	GLN	4.1
2	H	121	VAL	4.1
3	L	208	ALA	4.0
1	A	368	LEU	3.9
2	H	135	THR	3.8
2	H	150	VAL	3.8
3	L	117	THR	3.8
2	H	192	GLN	3.8
2	H	205	THR	3.7
3	L	177	SER	3.7
2	H	126	PRO	3.7
3	L	131	ALA	3.7
3	L	156	VAL	3.6
3	L	186	TRP	3.6
2	H	186	SER	3.6
3	L	141	TYR	3.5
2	H	158	ALA	3.5
3	L	136	LEU	3.5
2	H	201	LYS	3.4
2	H	213	PRO	3.4
2	H	151	THR	3.4
2	H	124	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
2	H	190	GLY	3.4
3	L	123	SER	3.4
2	H	185	PRO	3.3
3	L	135	CYS	3.3
1	A	518	LEU	3.3
3	L	184	GLU	3.3
3	L	178	TYR	3.2
1	A	525	CYS	3.2
2	H	152	VAL	3.2
3	L	160	VAL	3.2
3	L	154	SER	3.2
1	A	519	HIS	3.2
2	H	159	LEU	3.2
3	L	132	THR	3.1
3	L	124	GLU	3.1
3	L	195	GLN	3.1
4	M	61	SER	3.1
2	H	191	THR	3.1
3	L	112	ALA	3.1
1	A	521	PRO	3.1
4	M	29	PHE	3.0
2	H	179	SER	3.0
3	L	153	SER	3.0
3	L	128	ALA	3.0
2	H	183	THR	2.9
1	A	522	ALA	2.9
2	H	157	GLY	2.9
2	H	199	ASN	2.9
2	H	189	LEU	2.9
3	L	189	HIS	2.9
1	A	364	ASP	2.9
3	L	192	TYR	2.9
3	L	201	SER	2.9
3	L	93	SER	2.8
1	A	372	ALA	2.8
3	L	199	GLU	2.8
3	L	176	SER	2.8
3	L	116	VAL	2.8
3	L	187	LYS	2.8
3	L	182	THR	2.8
3	L	81	GLU	2.8
3	L	191	SER	2.8

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Mol	Chain	Res	Type	RSRZ
2	H	181	VAL	2.8
2	H	122	PHE	2.7
2	H	156	SER	2.7
2	H	139	GLY	2.7
3	L	155	PRO	2.7
3	L	129	ASN	2.7
2	H	187	SER	2.7
3	L	115	SER	2.6
2	H	168	ALA	2.6
1	A	394	ASN	2.6
1	A	516	GLU	2.6
2	H	137	ALA	2.6
5	N	45	LYS	2.6
2	H	161	SER	2.6
3	L	200	GLY	2.6
2	H	136	ALA	2.5
3	L	110	PRO	2.5
1	A	392	PHE	2.5
3	L	206	THR	2.5
3	L	127	GLN	2.5
2	H	155	ASN	2.5
3	L	27(C)	ILE	2.4
1	A	367	VAL	2.4
2	H	142	VAL	2.4
4	M	82(A)	SER	2.4
3	L	122	SER	2.4
3	L	168	GLN	2.4
1	A	524	VAL	2.4
3	L	114	PRO	2.4
3	L	190	ARG	2.4
4	M	85	SER	2.3
1	A	391	CYS	2.3
3	L	203	VAL	2.3
5	N	63	SER	2.3
2	H	160	THR	2.2
3	L	140	PHE	2.2
1	A	369	TYR	2.2
4	M	194	LEU	2.2
2	H	203	SER	2.2
3	L	188	SER	2.2
3	L	142	PRO	2.2
1	A	388	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
3	L	113	ALA	2.2
2	H	148	GLU	2.2
2	H	184	VAL	2.2
5	N	181	LEU	2.2
2	H	204	ASN	2.1
3	L	1	GLN	2.1
3	L	126	LEU	2.1
3	L	173	TYR	2.1
2	H	119	PRO	2.1
3	L	163	THR	2.1
3	L	202	THR	2.1
2	H	167	PRO	2.1
1	A	390	LEU	2.1
1	A	393	THR	2.1
5	N	41	GLY	2.1
3	L	166	SER	2.1
2	H	166	PHE	2.1
3	L	152	ASP	2.1
1	A	395	VAL	2.0
2	H	100	TYR	2.0
4	M	136	SER	2.0
2	H	211	VAL	2.0
5	N	116	PHE	2.0
3	L	111	LYS	2.0
5	N	213	GLU	2.0
3	L	171	ASN	2.0
1	A	383	SER	2.0
2	H	200	HIS	2.0
4	M	55	SER	2.0
4	M	144	SER	2.0

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

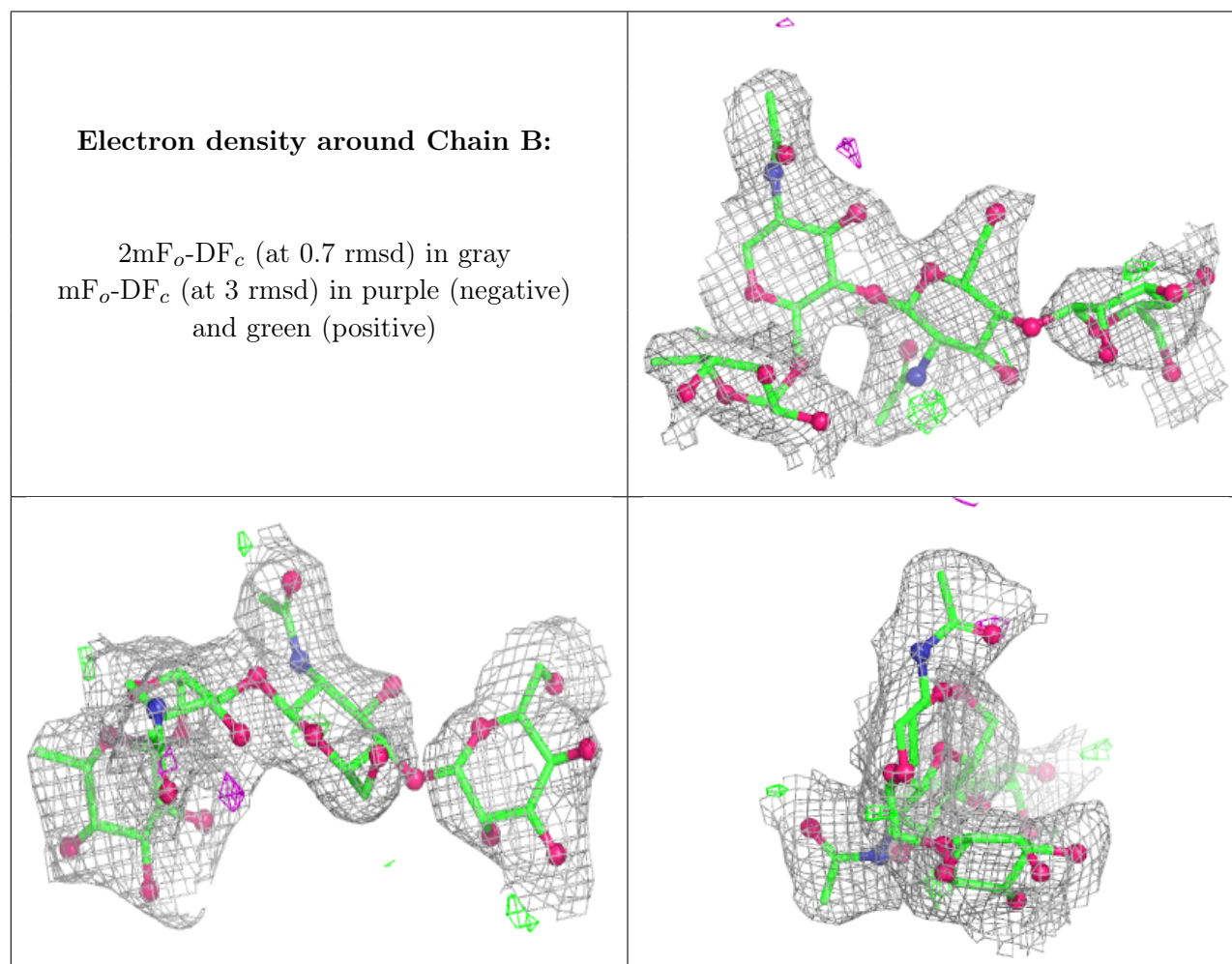
There are no non-standard protein/DNA/RNA residues in this entry.

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
6	NAG	B	1	14/15	-	-	24,36,47,50	0
6	NAG	B	2	14/15	-	-	34,44,57,76	0
6	BMA	B	3	11/12	-	-	59,68,76,77	0
6	FUC	B	4	10/11	-	-	31,40,42,42	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.