



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:33 AM UTC

PDB ID : 3S88 / pdb_00003s88
Title : Crystal structure of Sudan Ebolavirus Glycoprotein (strain Gulu) bound to 16F6
Authors : Sapphire, E.O.; Dias, J.M.; Bale, S.
Deposited on : 2011-05-27
Resolution : 3.35 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:

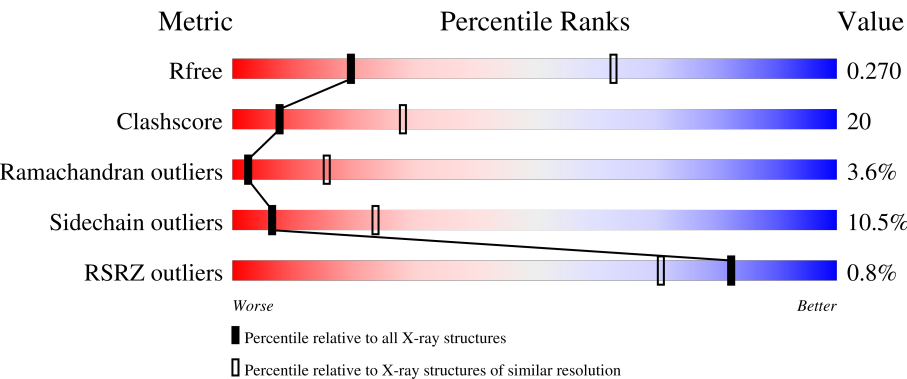
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (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 3.35 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	1099 (3.40-3.32)
Clashscore	190562	1116 (3.40-3.32)
Ramachandran outliers	187476	1101 (3.40-3.32)
Sidechain outliers	187428	1101 (3.40-3.32)
RSRZ outliers	180081	1099 (3.40-3.32)

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	H	220	62%29%9%
2	I	298	% 51%27%6%16%
3	J	167	% 33%27%•37%
4	L	212	62%34%••
5	A	2	100%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
5	B	2	 100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 6040 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 16F6 - Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	H	220	Total	C	N	O	S	0	0	0
			1660	1055	273	324	8			

- Molecule 2 is a protein called Envelope glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	I	249	Total	C	N	O	S	0	0	0
			1859	1180	318	355	6			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
I	16	TYR	-	expression tag	UNP Q7T9D9
I	17	PRO	-	expression tag	UNP Q7T9D9
I	18	TYR	-	expression tag	UNP Q7T9D9
I	19	ASP	-	expression tag	UNP Q7T9D9
I	20	VAL	-	expression tag	UNP Q7T9D9
I	21	PRO	-	expression tag	UNP Q7T9D9
I	22	ASP	-	expression tag	UNP Q7T9D9
I	23	TYR	-	expression tag	UNP Q7T9D9
I	24	ALA	-	expression tag	UNP Q7T9D9
I	25	ILE	-	expression tag	UNP Q7T9D9
I	26	GLU	-	expression tag	UNP Q7T9D9
I	27	GLY	-	expression tag	UNP Q7T9D9
I	28	ARG	-	expression tag	UNP Q7T9D9
I	29	GLY	-	expression tag	UNP Q7T9D9
I	30	ALA	-	expression tag	UNP Q7T9D9
I	31	ARG	-	expression tag	UNP Q7T9D9

- Molecule 3 is a protein called Envelope glycoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	J	106	Total	C	N	O	S	0	0	0
			817	514	149	148	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
J	631	VAL	ILE	engineered mutation	UNP Q7T9D9
J	638	VAL	GLN	engineered mutation	UNP Q7T9D9

- Molecule 4 is a protein called 16F6 - Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	L	212	Total	C	N	O	S	0	0	0
			1648	1030	281	330	7			

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

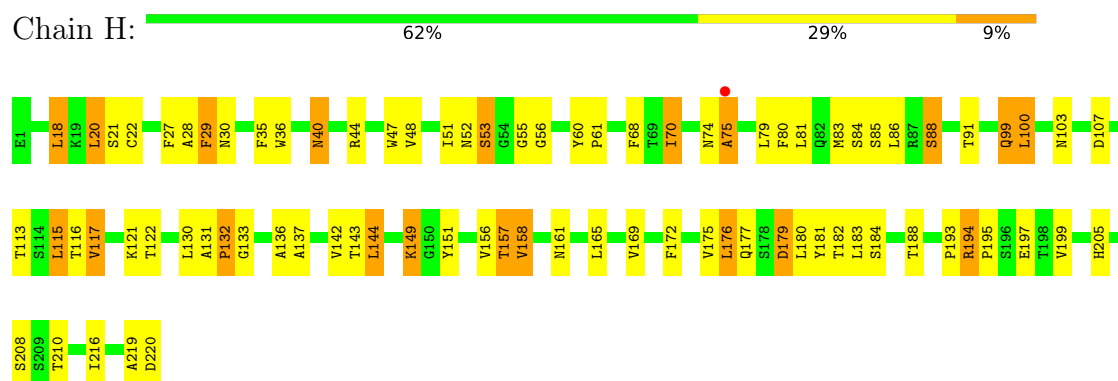


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	A	2	Total	C	N	O	0	0	0
			28	16	2	10			
5	B	2	Total	C	N	O	0	0	0
			28	16	2	10			

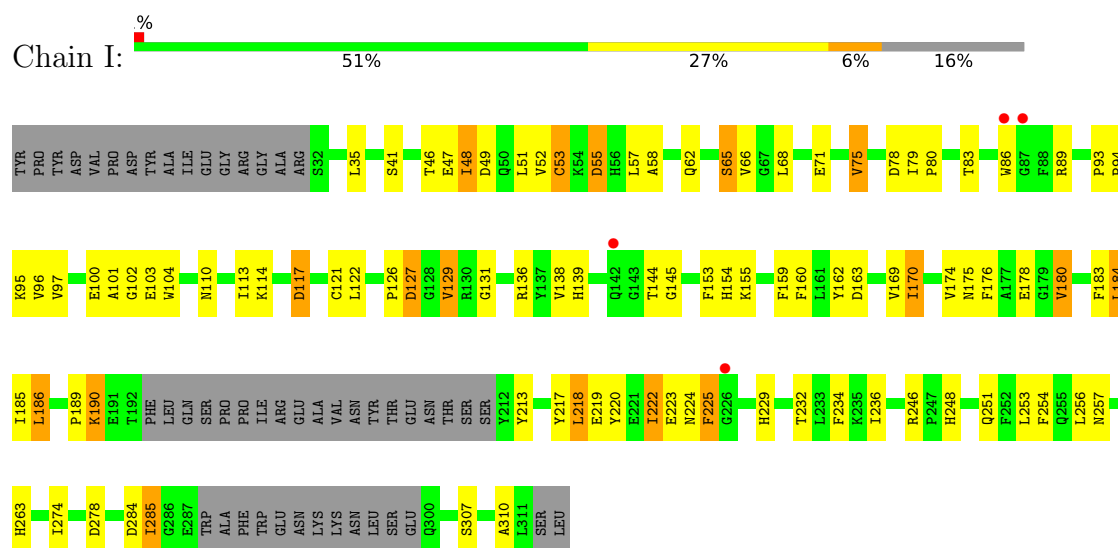
3 Residue-property plots [i](#)

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.

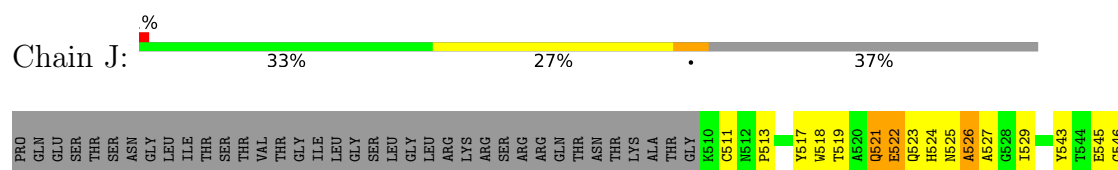
• Molecule 1: 16F6 - Heavy chain



• Molecule 2: Envelope glycoprotein



• Molecule 3: Envelope glycoprotein





• Molecule 4: 16F6 - Light chain



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



• Molecule 5: 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 3	Depositor
Cell constants a, b, c, α , β , γ	193.59Å 193.59Å 193.59Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.63 – 3.35 45.63 – 3.35	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.63-3.35) 97.2 (45.63-3.35)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.56 (at 3.32Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.222 , 0.278 0.216 , 0.270	Depositor DCC
R_{free} test set	864 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	104.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 91.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.021 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6040	wwPDB-VP
Average B, all atoms (Å ²)	121.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% 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	H	0.89	0/1704	1.01	3/2326 (0.1%)
2	I	0.74	0/1904	0.93	2/2593 (0.1%)
3	J	0.80	0/838	0.91	1/1144 (0.1%)
4	L	0.81	0/1687	0.97	3/2295 (0.1%)
All	All	0.81	0/6133	0.96	9/8358 (0.1%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L	80	ALA	N-CA-C	-7.10	102.73	111.40
4	L	178	THR	N-CA-C	6.05	118.27	107.80
3	J	555	VAL	CB-CA-C	-5.96	104.35	111.97
1	H	149	LYS	N-CA-C	5.71	118.78	109.59
2	I	222	ILE	N-CA-C	5.37	116.22	108.48
4	L	108	ARG	N-CA-C	5.30	115.53	108.38
2	I	263	HIS	N-CA-C	5.25	113.09	108.78
1	H	176	LEU	CA-C-N	-5.04	117.74	123.13
1	H	176	LEU	C-N-CA	-5.04	117.74	123.13

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	H	1660	0	1615	69	0
2	I	1859	0	1709	73	0
3	J	817	0	756	42	0
4	L	1648	0	1588	66	0
5	A	28	0	25	0	0
5	B	28	0	25	2	0
All	All	6040	0	5718	234	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 20.

All (234) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:88:SER:O	1:H:91:THR:HG22	1.47	1.13
3:J:521:GLN:HA	3:J:522:GLU:HB3	1.20	1.13
3:J:521:GLN:HA	3:J:522:GLU:CB	1.89	1.01
4:L:163:TRP:HE1	4:L:175:MET:HE2	1.32	0.92
4:L:163:TRP:NE1	4:L:175:MET:HE2	1.86	0.90
4:L:193:THR:HG22	4:L:208:SER:HB2	1.53	0.89
4:L:4:MET:HE3	4:L:23:CYS:SG	2.13	0.87
4:L:20:THR:CG2	4:L:74:THR:HG23	2.09	0.81
2:I:185:ILE:C	2:I:186:LEU:HD23	2.05	0.81
4:L:20:THR:HG23	4:L:74:THR:HG23	1.64	0.78
2:I:66:VAL:HG12	2:I:184:LEU:HD11	1.65	0.78
2:I:180:VAL:HG22	3:J:562:ALA:HB1	1.65	0.77
1:H:179:ASP:O	1:H:180:LEU:HD12	1.85	0.77
2:I:126:PRO:O	2:I:129:VAL:HG12	1.87	0.75
1:H:74:ASN:O	1:H:75:ALA:HB3	1.85	0.74
4:L:195:GLU:HG2	4:L:206:VAL:HG22	1.70	0.73
1:H:40:ASN:HD21	1:H:44:ARG:H	1.37	0.73
4:L:189:HIS:O	4:L:211:ARG:HD3	1.90	0.72
1:H:133:GLY:HA2	1:H:219:ALA:HB2	1.72	0.71
3:J:521:GLN:CA	3:J:522:GLU:HB3	2.11	0.70
1:H:133:GLY:HA2	1:H:219:ALA:CB	2.21	0.70
4:L:193:THR:HG22	4:L:208:SER:CB	2.23	0.68
3:J:573:LEU:O	3:J:576:THR:HG22	1.92	0.68
2:I:222:ILE:HG23	2:I:232:THR:HG22	1.76	0.67
1:H:74:ASN:O	1:H:75:ALA:CB	2.43	0.67
1:H:40:ASN:C	1:H:40:ASN:HD22	2.03	0.66
2:I:256:LEU:HD13	2:I:256:LEU:C	2.21	0.66
4:L:190:ASN:ND2	4:L:210:ASN:HB3	2.11	0.66

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:159:PHE:CZ	3:J:569:LEU:HD12	2.32	0.65
2:I:101:ALA:HB2	3:J:519:THR:HG22	1.78	0.65
3:J:563:ASN:O	3:J:566:THR:HG22	1.98	0.64
1:H:35:PHE:HE2	1:H:99:GLN:HB2	1.61	0.64
1:H:83:MET:HB3	1:H:86:LEU:HD21	1.80	0.64
2:I:71:GLU:HA	2:I:75:VAL:HG22	1.82	0.62
4:L:29:VAL:O	4:L:29:VAL:HG12	1.98	0.62
2:I:217:TYR:O	2:I:218:LEU:HD13	1.99	0.62
2:I:234:PHE:CD1	2:I:253:LEU:HD23	2.34	0.61
2:I:236:ILE:HD12	2:I:257:ASN:HA	1.81	0.61
1:H:51:ILE:HB	1:H:70:ILE:HD13	1.82	0.61
2:I:66:VAL:CG1	2:I:184:LEU:HD11	2.31	0.61
3:J:549:HIS:O	3:J:550:ASN:ND2	2.34	0.60
2:I:96:VAL:HG13	3:J:580:ARG:HA	1.82	0.60
4:L:62:PHE:CE1	4:L:75:LEU:HD13	2.36	0.59
2:I:103:GLU:O	2:I:104:TRP:C	2.44	0.59
2:I:138:VAL:CG2	2:I:217:TYR:CD1	2.85	0.59
1:H:131:ALA:HB2	1:H:216:ILE:CG2	2.34	0.58
2:I:52:VAL:HG13	2:I:53:CYS:O	2.03	0.58
2:I:68:LEU:HD11	3:J:558:LEU:HD12	1.86	0.58
4:L:24:LYS:NZ	4:L:70:ALA:HB2	2.19	0.58
2:I:101:ALA:CB	3:J:519:THR:HG22	2.33	0.58
3:J:558:LEU:HD12	3:J:558:LEU:C	2.29	0.58
4:L:161:ASN:HD22	4:L:177:SER:HA	1.68	0.58
1:H:99:GLN:HG2	1:H:100:LEU:O	2.04	0.58
4:L:142:LYS:HB2	4:L:173:TYR:CE2	2.38	0.57
2:I:236:ILE:HG23	2:I:257:ASN:OD1	2.04	0.57
2:I:68:LEU:HD11	3:J:558:LEU:CD1	2.34	0.57
4:L:135:PHE:C	4:L:136:LEU:HD12	2.29	0.57
4:L:29:VAL:HG13	4:L:92:TYR:HB2	1.85	0.57
4:L:136:LEU:HD12	4:L:136:LEU:N	2.20	0.57
2:I:162:TYR:CE1	2:I:176:PHE:HB3	2.39	0.57
1:H:40:ASN:ND2	1:H:44:ARG:H	2.02	0.57
1:H:177:GLN:NE2	4:L:160:LEU:HD11	2.19	0.57
4:L:75:LEU:CD2	4:L:78:VAL:HG22	2.35	0.56
3:J:549:HIS:O	3:J:550:ASN:CG	2.48	0.56
1:H:142:VAL:HG22	1:H:143:THR:N	2.20	0.56
1:H:194:ARG:HD2	1:H:195:PRO:HA	1.88	0.56
2:I:52:VAL:O	2:I:55:ASP:OD1	2.23	0.56
2:I:79:ILE:HG21	2:I:220:TYR:CE1	2.41	0.56
2:I:184:LEU:N	2:I:184:LEU:HD12	2.21	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:219:GLU:C	2:I:220:TYR:CD2	2.83	0.56
4:L:82:ASP:O	4:L:86:TYR:OH	2.24	0.56
1:H:176:LEU:HD13	1:H:181:TYR:CD1	2.40	0.55
4:L:75:LEU:HD23	4:L:78:VAL:HG22	1.87	0.55
3:J:583:THR:HG22	3:J:583:THR:O	2.06	0.55
2:I:138:VAL:HG21	2:I:217:TYR:CD1	2.41	0.55
5:B:1:NAG:O4	5:B:2:NAG:H83	2.07	0.55
1:H:20:LEU:HD13	1:H:83:MET:HE2	1.87	0.55
2:I:97:VAL:HG12	3:J:581:THR:HB	1.88	0.55
3:J:561:LEU:O	3:J:565:THR:HG23	2.05	0.55
2:I:131:GLY:HA2	2:I:174:VAL:HG12	1.89	0.54
4:L:15:VAL:O	4:L:15:VAL:HG13	2.06	0.54
2:I:234:PHE:CG	2:I:253:LEU:HD23	2.42	0.54
2:I:154:HIS:NE2	2:I:178:GLU:OE1	2.30	0.54
4:L:30:THR:OG1	4:L:31:THR:N	2.39	0.54
4:L:181:LEU:N	4:L:181:LEU:HD12	2.22	0.54
1:H:40:ASN:C	1:H:40:ASN:ND2	2.66	0.54
4:L:20:THR:HG22	4:L:74:THR:HG23	1.87	0.54
1:H:131:ALA:HB2	1:H:216:ILE:HG22	1.90	0.54
2:I:79:ILE:O	2:I:83:THR:OG1	2.25	0.54
2:I:144:THR:O	2:I:224:ASN:OD1	2.25	0.54
1:H:161:ASN:ND2	1:H:165:LEU:HD12	2.23	0.53
4:L:115:VAL:HG22	4:L:136:LEU:HG	1.90	0.53
3:J:550:ASN:O	3:J:551:GLN:C	2.51	0.53
4:L:16:GLY:O	4:L:17:ASP:O	2.26	0.53
2:I:127:ASP:O	3:J:580:ARG:NH2	2.42	0.53
4:L:96:LEU:H	4:L:96:LEU:HD22	1.74	0.53
1:H:35:PHE:CE2	1:H:99:GLN:HB2	2.44	0.53
1:H:208:SER:HB3	1:H:210:THR:HG23	1.91	0.52
1:H:179:ASP:C	1:H:180:LEU:HD12	2.34	0.52
1:H:143:THR:C	1:H:144:LEU:HD13	2.34	0.52
2:I:110:ASN:HD22	2:I:175:ASN:HD22	1.55	0.52
3:J:529:ILE:HG22	3:J:529:ILE:O	2.10	0.52
3:J:583:THR:HA	3:J:586:ASN:ND2	2.25	0.52
2:I:46:THR:OG1	2:I:47:GLU:N	2.43	0.51
4:L:197:THR:HG22	4:L:204:PRO:HG3	1.92	0.51
1:H:36:TRP:O	1:H:48:VAL:HG22	2.10	0.51
2:I:68:LEU:CD1	3:J:559:ARG:HA	2.40	0.51
2:I:186:LEU:HD23	2:I:186:LEU:N	2.25	0.51
3:J:550:ASN:OD1	3:J:551:GLN:N	2.44	0.51
1:H:52:ASN:O	1:H:55:GLY:N	2.40	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:124:GLN:HE22	4:L:131:SER:CB	2.25	0.50
4:L:2:ILE:HD12	4:L:2:ILE:H	1.76	0.50
4:L:117:ILE:HD13	4:L:208:SER:HA	1.93	0.50
4:L:85:LEU:HD11	4:L:103:LYS:HG2	1.93	0.50
2:I:48:ILE:HD12	2:I:49:ASP:H	1.77	0.50
2:I:126:PRO:O	2:I:129:VAL:CG1	2.60	0.49
1:H:60:TYR:OH	1:H:70:ILE:HG22	2.12	0.49
4:L:36:TYR:OH	4:L:89:GLN:NE2	2.46	0.49
4:L:181:LEU:HD23	4:L:185:GLU:HG2	1.94	0.49
1:H:158:VAL:HG21	1:H:183:LEU:HD11	1.95	0.49
2:I:126:PRO:HD2	2:I:129:VAL:HG11	1.95	0.49
4:L:2:ILE:HD12	4:L:2:ILE:N	2.28	0.49
1:H:18:LEU:CD2	1:H:115:LEU:HD23	2.42	0.48
2:I:86:TRP:HB2	2:I:153:PHE:O	2.13	0.48
2:I:139:HIS:ND1	2:I:139:HIS:N	2.61	0.48
1:H:193:PRO:O	1:H:194:ARG:CB	2.61	0.48
4:L:163:TRP:HE1	4:L:175:MET:CE	2.13	0.48
1:H:151:TYR:OH	1:H:183:LEU:HD23	2.14	0.48
4:L:29:VAL:O	4:L:29:VAL:CG1	2.62	0.48
4:L:24:LYS:HZ3	4:L:70:ALA:HB2	1.78	0.47
4:L:170:ASP:O	4:L:172:THR:HG23	2.14	0.47
1:H:132:PRO:O	1:H:219:ALA:HB3	2.13	0.47
2:I:65:SER:OG	2:I:100:GLU:HG3	2.14	0.47
2:I:57:LEU:HD13	2:I:57:LEU:C	2.39	0.47
3:J:590:ILE:O	3:J:591:ASP:C	2.57	0.47
2:I:35:LEU:HD12	2:I:183:PHE:O	2.14	0.47
1:H:172:PHE:CD1	4:L:164:THR:HG23	2.50	0.47
3:J:554:LEU:C	3:J:554:LEU:HD23	2.40	0.47
1:H:103:ASN:HD21	3:J:553:ALA:HA	1.79	0.47
1:H:144:LEU:HB3	1:H:216:ILE:HG21	1.97	0.47
4:L:137:ASN:HB3	4:L:174:SER:HB3	1.97	0.47
1:H:161:ASN:HD22	1:H:165:LEU:HD12	1.80	0.46
1:H:176:LEU:HD13	1:H:181:TYR:CE1	2.50	0.46
1:H:149:LYS:HA	1:H:182:THR:HG23	1.95	0.46
4:L:134:CYS:SG	4:L:136:LEU:HD11	2.55	0.46
1:H:205:HIS:CE1	1:H:208:SER:HG	2.31	0.46
2:I:138:VAL:CG2	2:I:217:TYR:HD1	2.28	0.46
2:I:47:GLU:O	2:I:48:ILE:C	2.59	0.46
1:H:22:CYS:HB3	1:H:79:LEU:HB3	1.98	0.46
3:J:603:ILE:O	3:J:605:GLY:N	2.49	0.46
1:H:130:LEU:HB3	4:L:118:PHE:CD2	2.52	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:177:GLN:HA	1:H:177:GLN:OE1	2.17	0.45
4:L:95:PRO:O	4:L:97:THR:HG23	2.16	0.45
2:I:256:LEU:HD13	2:I:256:LEU:O	2.16	0.45
3:J:566:THR:HG21	5:B:1:NAG:H82	1.98	0.45
1:H:179:ASP:OD1	1:H:179:ASP:N	2.49	0.45
3:J:517:TYR:CZ	3:J:546:GLY:HA3	2.52	0.45
1:H:183:LEU:HD12	1:H:184:SER:N	2.31	0.45
4:L:16:GLY:O	4:L:17:ASP:C	2.60	0.45
1:H:133:GLY:CA	1:H:219:ALA:CB	2.93	0.44
2:I:138:VAL:HG21	2:I:217:TYR:HD1	1.80	0.44
4:L:181:LEU:HD23	4:L:185:GLU:CD	2.42	0.44
2:I:223:GLU:HA	2:I:224:ASN:OD1	2.16	0.44
4:L:86:TYR:CE2	4:L:104:LEU:HD22	2.52	0.44
2:I:113:ILE:HG22	2:I:121:CYS:SG	2.57	0.44
4:L:2:ILE:CG2	4:L:3:VAL:N	2.80	0.44
4:L:29:VAL:CG1	4:L:32:ALA:HB3	2.47	0.44
1:H:161:ASN:HD21	1:H:199:VAL:HA	1.82	0.44
2:I:52:VAL:HG22	2:I:53:CYS:N	2.33	0.44
2:I:68:LEU:HD13	3:J:559:ARG:HG2	2.00	0.44
3:J:554:LEU:HD23	3:J:554:LEU:O	2.17	0.44
1:H:107:ASP:OD1	1:H:107:ASP:N	2.44	0.44
1:H:121:LYS:O	1:H:122:THR:C	2.58	0.44
4:L:15:VAL:O	4:L:15:VAL:CG1	2.65	0.44
1:H:52:ASN:O	1:H:53:SER:C	2.61	0.44
2:I:58:ALA:HB3	2:I:62:GLN:OE1	2.17	0.43
2:I:138:VAL:HG22	2:I:217:TYR:CD1	2.53	0.43
4:L:31:THR:O	4:L:31:THR:OG1	2.33	0.43
1:H:91:THR:HB	1:H:117:VAL:H	1.82	0.43
2:I:110:ASN:HD22	2:I:175:ASN:ND2	2.16	0.43
4:L:150:ILE:HG21	4:L:189:HIS:CD2	2.53	0.43
1:H:132:PRO:O	1:H:219:ALA:CB	2.67	0.43
4:L:37:GLN:HG3	4:L:86:TYR:CE1	2.53	0.43
4:L:122:SER:HA	4:L:125:LEU:HD23	2.01	0.43
3:J:579:LEU:HD12	3:J:579:LEU:H	1.83	0.43
1:H:113:THR:O	1:H:113:THR:HG23	2.18	0.43
2:I:160:PHE:CE1	2:I:170:ILE:HD11	2.54	0.43
4:L:170:ASP:OD1	4:L:172:THR:HG23	2.19	0.43
1:H:136:ALA:O	1:H:137:ALA:HB3	2.18	0.42
1:H:177:GLN:O	1:H:179:ASP:N	2.48	0.42
2:I:102:GLY:HA3	3:J:518:TRP:CZ2	2.53	0.42
4:L:36:TYR:HE1	4:L:89:GLN:HE21	1.66	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:L:133:VAL:HG22	4:L:178:THR:HB	2.01	0.42
2:I:251:GLN:OE1	2:I:251:GLN:N	2.47	0.42
4:L:77:SER:O	4:L:78:VAL:C	2.62	0.42
1:H:156:VAL:HG12	1:H:157:THR:N	2.32	0.42
2:I:176:PHE:CD1	2:I:176:PHE:C	2.97	0.42
2:I:284:ASP:O	2:I:285:ILE:CB	2.67	0.42
2:I:117:ASP:OD2	2:I:117:ASP:N	2.50	0.42
3:J:601:CYS:SG	3:J:609:CYS:O	2.78	0.42
1:H:208:SER:O	1:H:210:THR:HG23	2.19	0.42
2:I:93:PRO:HA	2:I:94:PRO:HD3	1.92	0.42
1:H:142:VAL:CG2	1:H:143:THR:N	2.83	0.42
2:I:113:ILE:HD13	2:I:225:PHE:CG	2.55	0.42
3:J:513:PRO:HB2	3:J:552:ASN:OD1	2.19	0.42
3:J:523:GLN:CD	3:J:523:GLN:H	2.27	0.42
3:J:525:ASN:O	3:J:526:ALA:HB3	2.20	0.42
1:H:28:ALA:O	1:H:29:PHE:C	2.62	0.42
1:H:47:TRP:O	1:H:61:PRO:HG3	2.19	0.42
3:J:569:LEU:HD22	3:J:569:LEU:HA	1.85	0.42
4:L:2:ILE:HG22	4:L:3:VAL:N	2.35	0.42
2:I:189:PRO:CB	2:I:190:LYS:CB	2.98	0.42
2:I:78:ASP:C	2:I:80:PRO:HD2	2.45	0.41
1:H:131:ALA:O	1:H:133:GLY:N	2.53	0.41
2:I:217:TYR:CD2	2:I:217:TYR:N	2.88	0.41
2:I:223:GLU:HA	2:I:224:ASN:HA	1.81	0.41
3:J:569:LEU:O	3:J:573:LEU:HD13	2.19	0.41
4:L:190:ASN:HD21	4:L:210:ASN:HB3	1.84	0.41
4:L:96:LEU:HD22	4:L:96:LEU:N	2.35	0.41
1:H:133:GLY:O	1:H:219:ALA:CB	2.69	0.41
2:I:113:ILE:CG2	2:I:114:LYS:N	2.83	0.41
4:L:85:LEU:HD13	4:L:85:LEU:HA	1.91	0.41
4:L:85:LEU:HD21	4:L:103:LYS:HE2	2.03	0.41
3:J:583:THR:O	3:J:583:THR:CG2	2.67	0.41
1:H:52:ASN:HD22	1:H:56:GLY:H	1.68	0.41
1:H:84:SER:O	1:H:85:SER:C	2.61	0.41
2:I:169:VAL:HG12	2:I:170:ILE:N	2.34	0.41
4:L:161:ASN:HB2	4:L:163:TRP:CH2	2.56	0.41
1:H:113:THR:O	1:H:113:THR:CG2	2.69	0.41
2:I:163:ASP:OD1	3:J:543:TYR:OH	2.38	0.41
2:I:145:GLY:N	2:I:225:PHE:HZ	2.18	0.40
1:H:79:LEU:HD12	1:H:80:PHE:H	1.86	0.40
1:H:68:PHE:N	1:H:68:PHE:CD1	2.89	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:27:PHE:CE1	1:H:29:PHE:HA	2.56	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	H	218/220 (99%)	200 (92%)	11 (5%)	7 (3%)	3	17
2	I	243/298 (82%)	196 (81%)	37 (15%)	10 (4%)	2	14
3	J	104/167 (62%)	88 (85%)	7 (7%)	9 (9%)	0	4
4	L	210/212 (99%)	191 (91%)	17 (8%)	2 (1%)	12	38
All	All	775/897 (86%)	675 (87%)	72 (9%)	28 (4%)	2	15

All (28) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	194	ARG
2	I	225	PHE
2	I	274	ILE
3	J	522	GLU
3	J	604	LEU
3	J	605	GLY
4	L	17	ASP
1	H	75	ALA
2	I	190	LYS
4	L	78	VAL
1	H	18	LEU
1	H	29	PHE
2	I	278	ASP
2	I	285	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	J	608	CYS
3	J	610	ILE
1	H	53	SER
2	I	41	SER
2	I	53	CYS
2	I	229	HIS
3	J	526	ALA
1	H	30	ASN
2	I	310	ALA
3	J	511	CYS
3	J	527	ALA
1	H	132	PRO
2	I	307	SER
3	J	606	PRO

5.3.2 Protein sidechains [i](#)

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	H	183/183 (100%)	163 (89%)	20 (11%)	6	22
2	I	186/256 (73%)	164 (88%)	22 (12%)	5	19
3	J	80/142 (56%)	72 (90%)	8 (10%)	7	26
4	L	187/187 (100%)	170 (91%)	17 (9%)	9	30
All	All	636/768 (83%)	569 (90%)	67 (10%)	6	24

All (67) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	20	LEU
1	H	21	SER
1	H	40	ASN
1	H	70	ILE
1	H	81	LEU
1	H	88	SER
1	H	99	GLN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	H	100	LEU
1	H	115	LEU
1	H	116	THR
1	H	117	VAL
1	H	144	LEU
1	H	157	THR
1	H	158	VAL
1	H	169	VAL
1	H	175	VAL
1	H	179	ASP
1	H	188	THR
1	H	197	GLU
1	H	220	ASP
2	I	48	ILE
2	I	51	LEU
2	I	55	ASP
2	I	65	SER
2	I	75	VAL
2	I	89	ARG
2	I	95	LYS
2	I	117	ASP
2	I	122	LEU
2	I	127	ASP
2	I	129	VAL
2	I	136	ARG
2	I	155	LYS
2	I	170	ILE
2	I	180	VAL
2	I	184	LEU
2	I	186	LEU
2	I	213	TYR
2	I	218	LEU
2	I	246	ARG
2	I	248	HIS
2	I	254	PHE
3	J	521	GLN
3	J	524	HIS
3	J	545	GLU
3	J	548	MET
3	J	560	GLN
3	J	569	LEU
3	J	584	ILE

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
3	J	601	CYS
4	L	5	THR
4	L	12	SER
4	L	17	ASP
4	L	31	THR
4	L	61	ARG
4	L	67	SER
4	L	85	LEU
4	L	89	GLN
4	L	105	GLU
4	L	116	SER
4	L	153	SER
4	L	167	ASP
4	L	168	SER
4	L	175	MET
4	L	176	SER
4	L	178	THR
4	L	181	LEU

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

Mol	Chain	Res	Type
1	H	39	GLN
1	H	40	ASN
1	H	52	ASN
1	H	82	GLN
1	H	99	GLN
1	H	103	ASN
1	H	161	ASN
2	I	56	HIS
2	I	175	ASN
3	J	524	HIS
3	J	551	GLN
3	J	570	GLN
4	L	6	GLN
4	L	8	HIS
4	L	38	GLN
4	L	89	GLN
4	L	124	GLN
4	L	145	ASN
4	L	161	ASN
4	L	189	HIS

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
4	L	190	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	$\# Z > 2$
5	NAG	A	1	2,5	14,14,15	0.51	0	17,19,21	0.86	0
5	NAG	A	2	5	14,14,15	0.57	0	17,19,21	0.87	0
5	NAG	B	1	3,5	14,14,15	0.56	0	17,19,21	1.83	4 (23%)
5	NAG	B	2	5	14,14,15	0.46	0	17,19,21	1.83	3 (17%)

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	1	2,5	-	0/6/23/26	0/1/1/1
5	NAG	A	2	5	-	1/6/23/26	0/1/1/1
5	NAG	B	1	3,5	-	0/6/23/26	0/1/1/1
5	NAG	B	2	5	-	4/6/23/26	0/1/1/1

There are no bond length outliers.

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	1	NAG	O5-C1-C2	-4.65	104.09	111.29
5	B	2	NAG	C4-C3-C2	-3.58	105.77	111.02
5	B	2	NAG	O5-C1-C2	-3.50	105.88	111.29
5	B	2	NAG	C1-O5-C5	3.37	116.70	112.19
5	B	1	NAG	C3-C4-C5	-3.22	104.39	110.23
5	B	1	NAG	O6-C6-C5	-2.72	102.07	111.33
5	B	1	NAG	C2-N2-C7	-2.55	119.48	122.90

There are no chirality outliers.

All (5) torsion outliers are listed below:

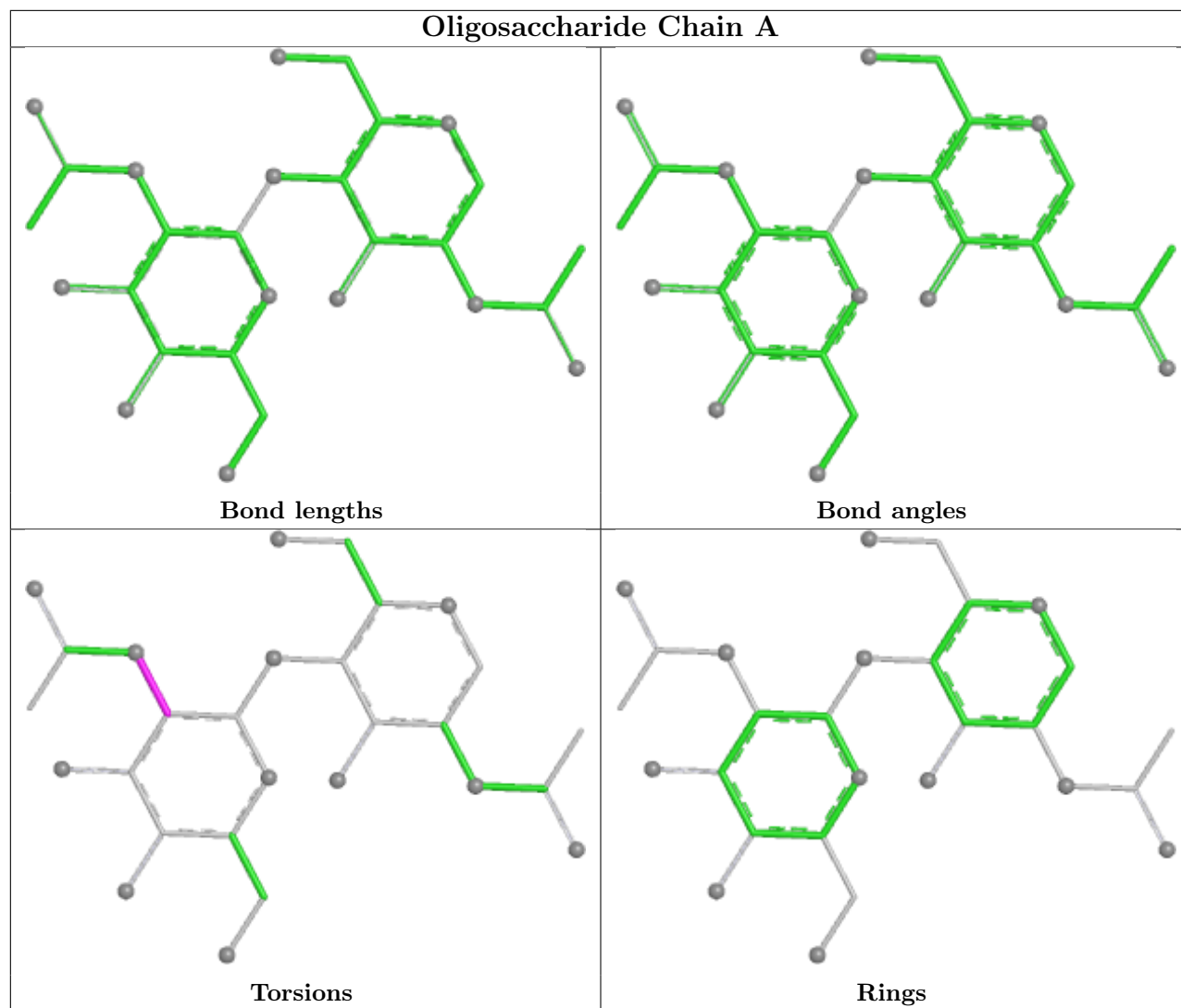
Mol	Chain	Res	Type	Atoms
5	B	2	NAG	C4-C5-C6-O6
5	B	2	NAG	O5-C5-C6-O6
5	B	2	NAG	C8-C7-N2-C2
5	B	2	NAG	O7-C7-N2-C2
5	A	2	NAG	C1-C2-N2-C7

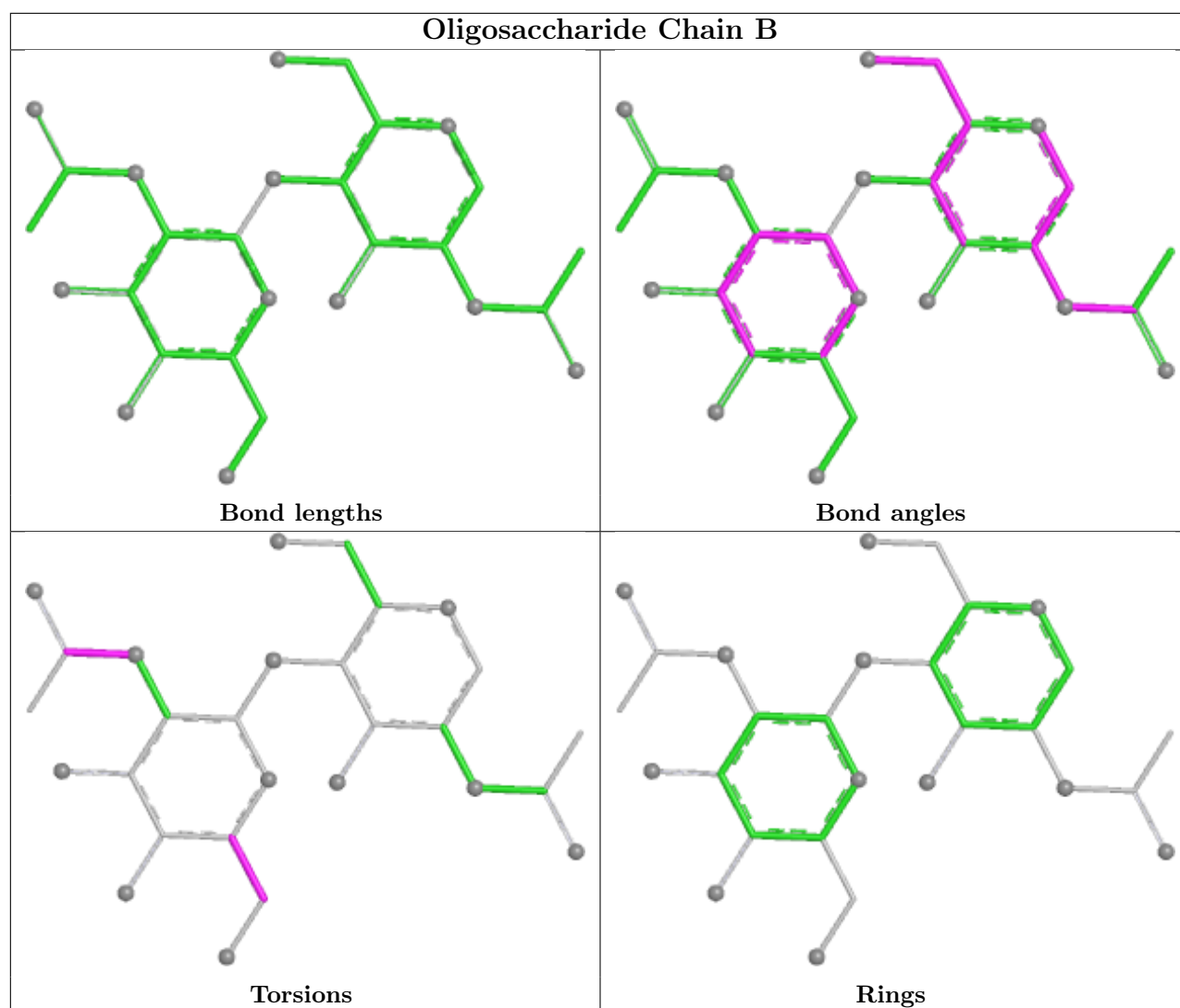
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	2	NAG	1	0
5	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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	H	220/220 (100%)	-0.45	1 (0%) 87 78	62, 93, 130, 162	0
2	I	249/298 (83%)	0.03	4 (1%) 70 55	81, 144, 248, 271	0
3	J	106/167 (63%)	-0.11	1 (0%) 81 69	78, 121, 229, 260	0
4	L	212/212 (100%)	-0.52	0 100 100	67, 104, 145, 165	0
All	All	787/897 (87%)	-0.27	6 (0%) 82 71	62, 111, 229, 271	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	I	86	TRP	4.0
2	I	87	GLY	3.6
1	H	75	ALA	2.9
3	J	597	TRP	2.6
2	I	142	GLN	2.5
2	I	226	GLY	2.1

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.

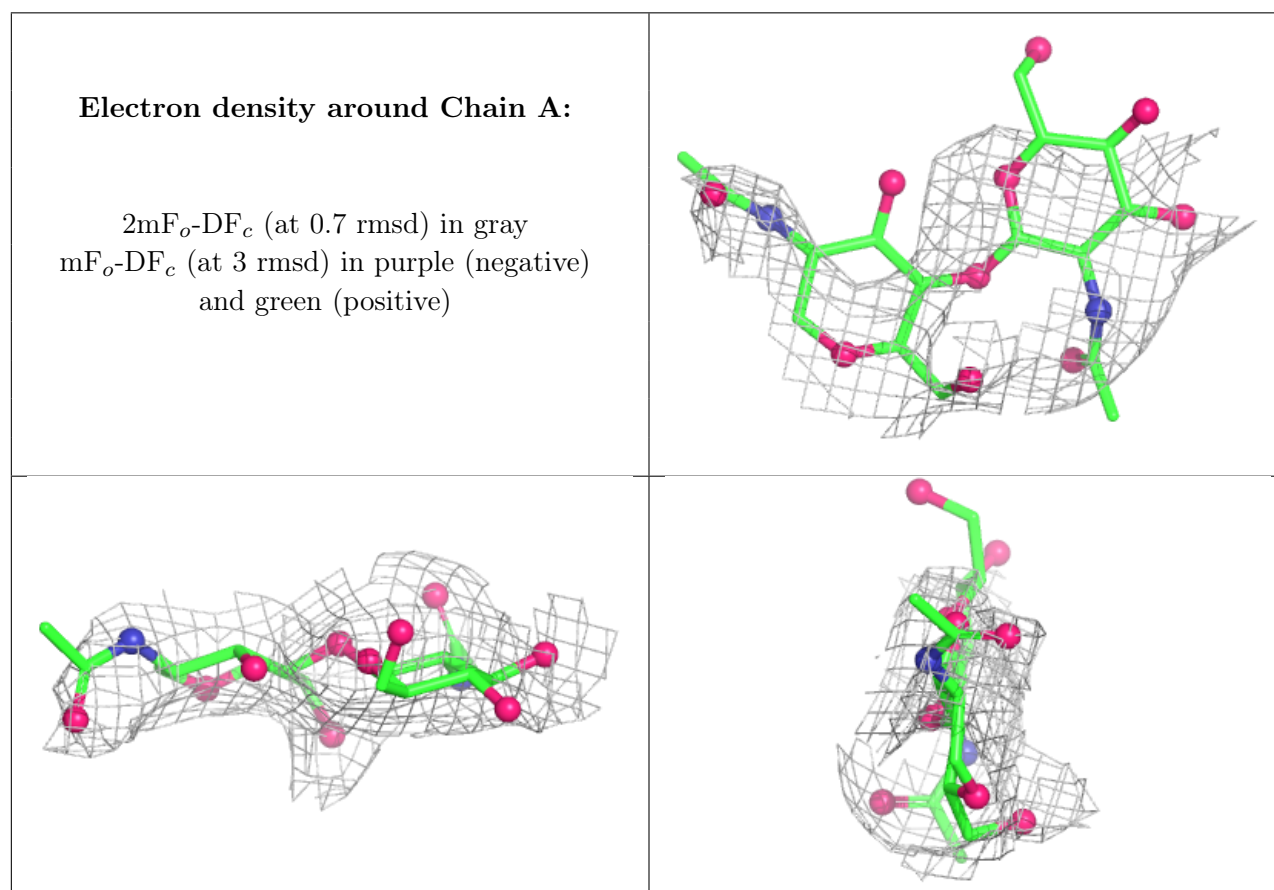
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	A	1	14/15	-	-	190,198,203,203	0

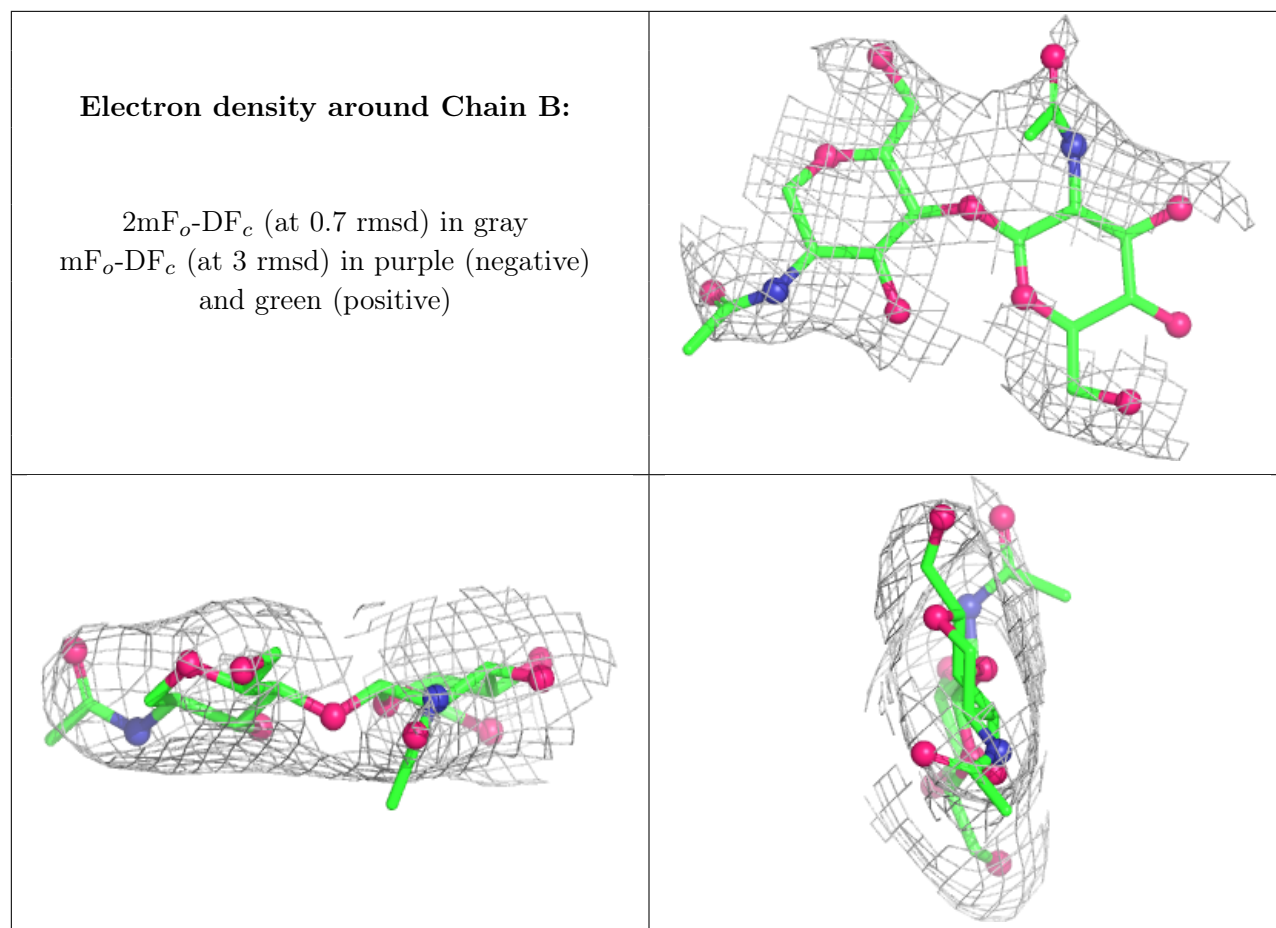
Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	NAG	A	2	14/15	-	-	202,206,212,213	0
5	NAG	B	1	14/15	-	-	129,136,143,150	0
5	NAG	B	2	14/15	-	-	156,162,165,165	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

There are no ligands in this entry.

6.5 Other polymers i

There are no such residues in this entry.