



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:40 PM UTC

PDB ID : 1JMA / pdb_00001jma
Title : CRYSTAL STRUCTURE OF THE HERPES SIMPLEX VIRUS GLYCOPROTEIN D BOUND TO THE CELLULAR RECEPTOR HVEA/HVEM
Authors : Carfi, A.; Willis, S.H.; Whitbeck, J.C.; Krummenacker, C.; Cohen, G.H.; Eisenberg, R.J.; Wiley, D.C.
Deposited on : 2001-07-17
Resolution : 2.65 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

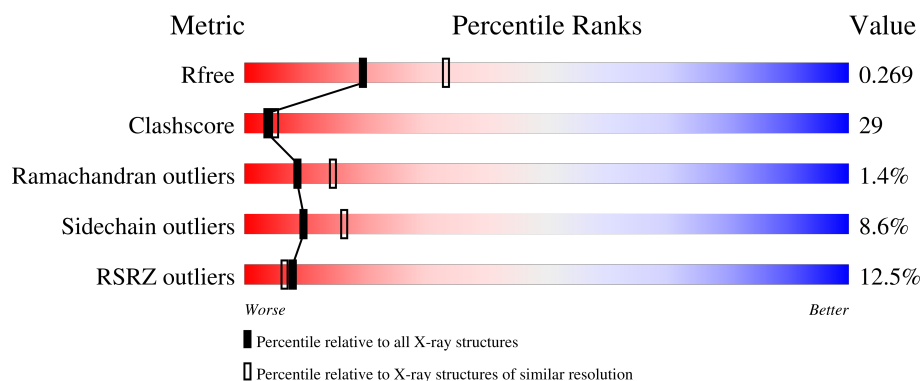
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

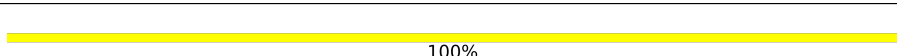
The reported resolution of this entry is 2.65 Å.

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	1110 (2.66-2.66)
Clashscore	190562	1141 (2.66-2.66)
Ramachandran outliers	187476	1126 (2.66-2.66)
Sidechain outliers	187428	1126 (2.66-2.66)
RSRZ outliers	180081	1110 (2.66-2.66)

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	B	167	
2	A	290	
3	C	2	

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 2885 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 HERPESVIRUS ENTRY MEDIATOR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	100	Total	C	N	O	S	0	0	0
			729	439	131	141	18			

- Molecule 2 is a protein called GLYCOPROTEIN D.

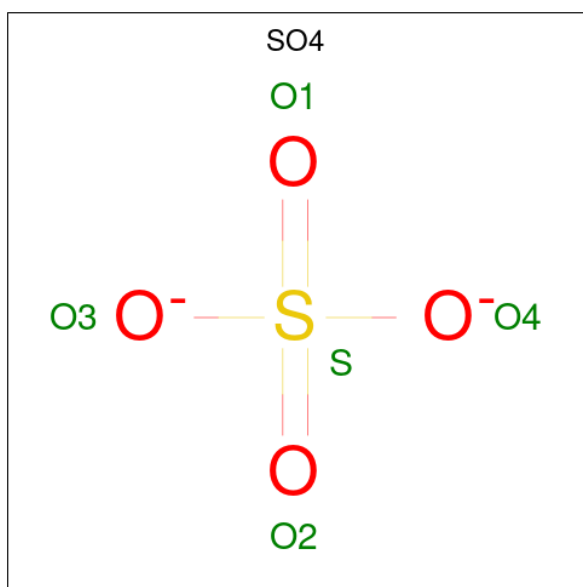
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	259	Total	C	N	O	S	0	2	0
			2054	1316	352	375	11			

- Molecule 3 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
3	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	O	S	0	0
			5	4	1		
4	B	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		
4	A	1	Total	O	S	0	0
			5	4	1		

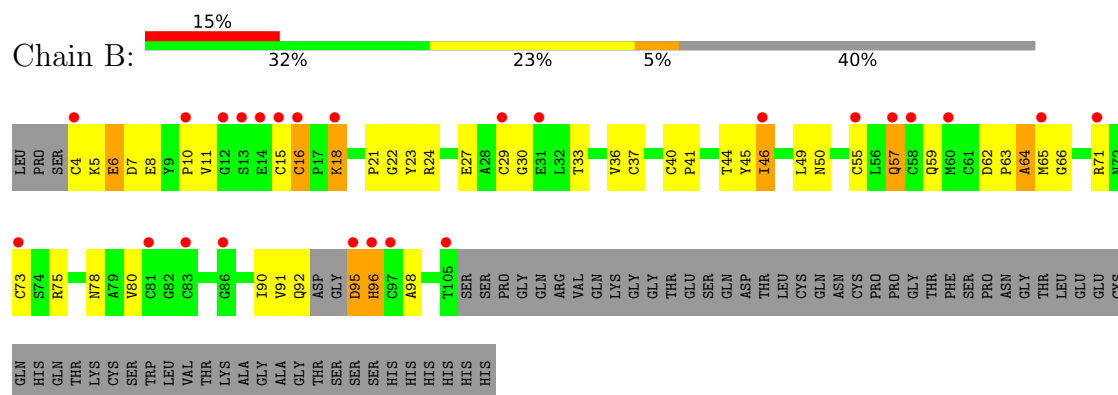
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	4	Total	O	0	0
			4	4		
5	A	45	Total	O	0	0
			45	45		

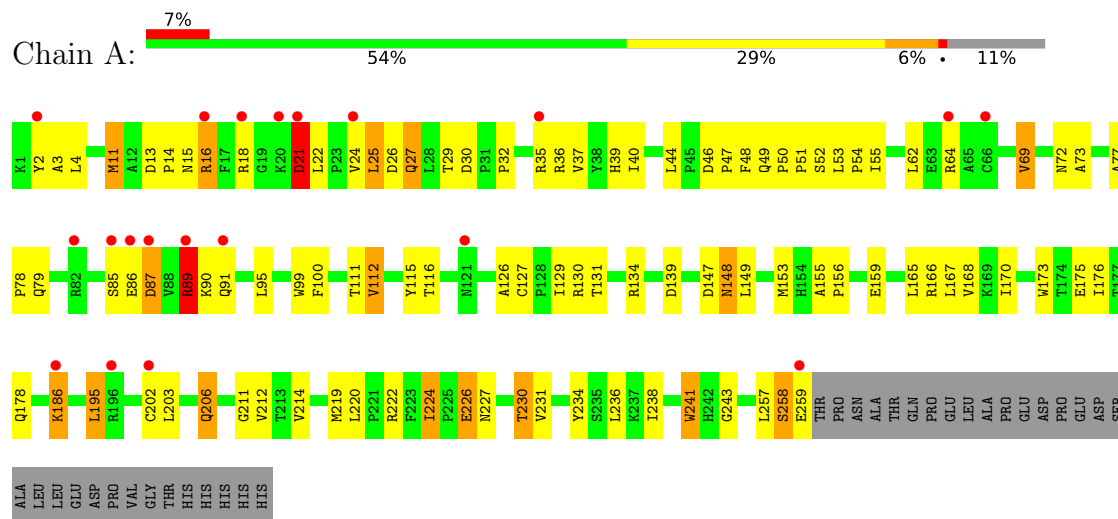
3 Residue-property plots

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: HERPESVIRUS ENTRY MEDIATOR



• Molecule 2: GLYCOPROTEIN D



• Molecule 3: 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 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	129.15Å 129.15Å 80.94Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	20.00 – 2.65 20.00 – 2.66	Depositor EDS
% Data completeness (in resolution range)	99.7 (20.00-2.65) 99.4 (20.00-2.66)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.41 (at 2.65Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.236 , 0.259 0.249 , 0.269	Depositor DCC
R_{free} test set	1101 reflections (4.86%)	wwPDB-VP
Wilson B-factor (Å ²)	46.9	Xtriage
Anisotropy	0.665	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 38.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	2885	wwPDB-VP
Average B, all atoms (Å ²)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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: SO4, 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	B	1.25	4/743 (0.5%)	1.28	11/1004 (1.1%)
2	A	0.76	5/2113 (0.2%)	1.18	15/2886 (0.5%)
All	All	0.91	9/2856 (0.3%)	1.20	26/3890 (0.7%)

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
2	A	0	2

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	6	GLU	CG-CD	27.17	2.19	1.52
2	A	89	ARG	CD-NE	-14.76	1.25	1.46
2	A	86	GLU	CA-CB	12.01	1.72	1.53
1	B	71	ARG	CB-CG	-11.57	1.17	1.52
2	A	258	SER	C-N	-9.86	1.19	1.33
2	A	87	ASP	CB-CG	-9.55	1.28	1.52
1	B	6	GLU	CB-CG	9.09	1.79	1.52
2	A	241	TRP	C-N	9.02	1.46	1.33
1	B	96	HIS	CB-CG	6.09	1.58	1.50

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	89	ARG	CG-CD-NE	16.73	148.82	112.00
2	A	91	GLN	CG-CD-NE2	13.86	137.19	116.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	86	GLU	N-CA-CB	12.61	128.82	110.16
2	A	258	SER	O-C-N	-11.78	106.93	122.59
1	B	92	GLN	OE1-CD-NE2	-10.21	112.39	122.60
2	A	91	GLN	OE1-CD-NE2	-9.82	112.78	122.60
1	B	6	GLU	CB-CG-CD	8.79	127.54	112.60
1	B	6	GLU	CA-CB-CG	-8.77	96.56	114.10
1	B	6	GLU	CG-CD-OE2	-8.60	98.62	118.40
1	B	6	GLU	CG-CD-OE1	8.57	138.11	118.40
2	A	87	ASP	CA-CB-CG	8.20	120.80	112.60
1	B	96	HIS	CB-CA-C	8.20	124.95	110.36
1	B	96	HIS	CA-CB-CG	-7.47	106.33	113.80
1	B	92	GLN	CG-CD-NE2	7.30	127.35	116.40
2	A	86	GLU	CA-CB-CG	7.23	128.57	114.10
2	A	139	ASP	N-CA-C	7.03	120.95	112.38
2	A	21	ASP	CB-CA-C	6.70	122.05	110.79
1	B	95	ASP	CA-C-N	-6.31	111.76	123.05
1	B	95	ASP	C-N-CA	-6.31	111.76	123.05
2	A	127	CYS	CA-C-N	6.23	125.64	118.85
2	A	127	CYS	C-N-CA	6.23	125.64	118.85
1	B	16	CYS	N-CA-C	5.67	117.41	108.23
2	A	91	GLN	CG-CD-OE1	-5.57	109.67	120.80
2	A	52	SER	N-CA-C	-5.40	106.72	113.15
2	A	29	THR	N-CA-C	-5.23	102.42	109.95
2	A	203	LEU	N-CA-C	5.00	117.68	110.28

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	A	258	SER	Mainchain
2	A	89	ARG	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	729	0	674	62	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	2054	0	2035	105	0
3	C	28	0	25	0	0
4	A	15	0	0	0	0
4	B	10	0	0	0	0
5	A	45	0	0	0	0
5	B	4	0	0	0	0
All	All	2885	0	2734	160	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 29.

All (160) 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:B:6:GLU:CB	1:B:6:GLU:CG	1.79	1.51
2:A:35:ARG:HD3	2:A:129:ILE:CD1	1.75	1.16
1:B:6:GLU:CG	1:B:6:GLU:CD	2.20	1.15
2:A:186:LYS:H	2:A:186:LYS:CD	1.60	1.10
2:A:11:MET:HB2	2:A:27:GLN:HG2	1.34	1.08
2:A:186:LYS:HD3	2:A:186:LYS:N	1.61	1.05
1:B:18:LYS:HD3	1:B:49:LEU:HB3	1.39	1.04
1:B:7:ASP:HB2	1:B:18:LYS:HE2	1.40	1.03
1:B:80:VAL:CG1	1:B:96:HIS:ND1	2.23	1.00
1:B:80:VAL:HG13	1:B:96:HIS:ND1	1.78	0.97
2:A:35:ARG:HG2	2:A:129:ILE:HD13	1.46	0.97
2:A:35:ARG:HD3	2:A:129:ILE:HD11	1.48	0.94
2:A:214:VAL:HG13	2:A:219:MET:HB2	1.52	0.91
2:A:186:LYS:H	2:A:186:LYS:HD3	0.75	0.91
1:B:6:GLU:CG	1:B:6:GLU:CA	2.49	0.90
2:A:35:ARG:CD	2:A:129:ILE:CD1	2.49	0.89
1:B:7:ASP:O	1:B:18:LYS:HG3	1.75	0.86
1:B:80:VAL:HG13	1:B:96:HIS:CG	2.10	0.86
2:A:147:ASP:O	2:A:148:ASN:HB2	1.77	0.83
2:A:11:MET:CB	2:A:27:GLN:HG2	2.09	0.83
2:A:35:ARG:HD3	2:A:129:ILE:HD12	1.56	0.82
2:A:30:ASP:CG	2:A:224:ILE:HD11	2.06	0.81
2:A:35:ARG:HG2	2:A:129:ILE:CD1	2.12	0.79
1:B:18:LYS:HD2	1:B:18:LYS:C	2.07	0.79
1:B:57:GLN:H	1:B:57:GLN:CD	1.92	0.78
2:A:2:TYR:HB3	2:A:37:VAL:HG11	1.64	0.78
2:A:35:ARG:CD	2:A:129:ILE:HD12	2.13	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:80:VAL:HG11	1:B:96:HIS:ND1	1.99	0.77
2:A:35:ARG:CG	2:A:129:ILE:CD1	2.63	0.77
1:B:18:LYS:HD3	1:B:49:LEU:CB	2.15	0.77
1:B:80:VAL:HG13	1:B:96:HIS:CB	2.14	0.76
1:B:30:GLY:HA3	1:B:33:THR:HG22	1.67	0.75
1:B:95:ASP:C	1:B:96:HIS:CD2	2.66	0.74
1:B:95:ASP:C	1:B:96:HIS:HD2	1.96	0.73
2:A:87:ASP:O	2:A:90:LYS:HG2	1.90	0.71
2:A:153:MET:HB3	2:A:156:PRO:HG3	1.73	0.70
1:B:80:VAL:CG1	1:B:96:HIS:CG	2.73	0.70
1:B:18:LYS:HD2	1:B:18:LYS:O	1.92	0.69
1:B:23:TYR:CD1	2:A:11:MET:HE2	2.28	0.68
2:A:30:ASP:OD1	2:A:224:ILE:HD11	1.94	0.67
2:A:35:ARG:CG	2:A:129:ILE:HD13	2.24	0.67
1:B:4:CYS:SG	1:B:10:PRO:HG3	2.36	0.65
2:A:227:ASN:O	2:A:231:VAL:HG23	1.96	0.65
2:A:47:PRO:HG3	2:A:100:PHE:CZ	2.31	0.65
1:B:90:ILE:HG13	1:B:91:VAL:HG23	1.80	0.64
2:A:167:LEU:C	2:A:167:LEU:HD23	2.22	0.64
2:A:165:LEU:HD21	2:A:176:ILE:HG23	1.80	0.63
2:A:35:ARG:CG	2:A:129:ILE:HD12	2.28	0.63
2:A:85:SER:O	2:A:89:ARG:HB2	1.99	0.63
2:A:25:LEU:HD13	2:A:27:GLN:HE22	1.62	0.62
2:A:72:ASN:HD22	2:A:73:ALA:H	1.48	0.62
2:A:18:ARG:O	2:A:22:LEU:HD22	1.99	0.61
1:B:30:GLY:HA3	1:B:33:THR:CG2	2.29	0.61
2:A:2:TYR:CB	2:A:37:VAL:HG11	2.32	0.60
2:A:27:GLN:HG3	2:A:227:ASN:OD1	2.02	0.60
1:B:18:LYS:CD	1:B:49:LEU:HB3	2.24	0.59
1:B:95:ASP:O	1:B:96:HIS:HD2	1.84	0.59
2:A:234:TYR:O	2:A:238:ILE:HG12	2.03	0.59
2:A:186:LYS:CD	2:A:186:LYS:N	2.37	0.57
2:A:25:LEU:HD21	2:A:230:THR:CG2	2.34	0.57
2:A:112:VAL:HG13	2:A:134:ARG:HB2	1.86	0.57
2:A:111:THR:HA	2:A:134:ARG:O	2.05	0.56
2:A:44:LEU:HD12	2:A:212:VAL:HG13	1.87	0.56
2:A:134:ARG:CD	2:A:214:VAL:HG11	2.36	0.56
2:A:220:LEU:CD1	2:A:222:ARG:HH12	2.18	0.56
1:B:24:ARG:HA	1:B:50:ASN:ND2	2.20	0.56
1:B:64:ALA:O	1:B:66:GLY:N	2.40	0.55
1:B:75:ARG:HD2	2:A:14:PRO:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:62:LEU:HD13	2:A:69:VAL:HG22	1.88	0.54
1:B:23:TYR:CG	2:A:11:MET:HE2	2.42	0.54
1:B:80:VAL:HG13	1:B:96:HIS:HB3	1.88	0.54
2:A:21:ASP:N	2:A:21:ASP:OD1	2.39	0.54
2:A:25:LEU:HD22	2:A:26:ASP:O	2.08	0.54
2:A:134:ARG:HD2	2:A:214:VAL:HG11	1.89	0.54
1:B:24:ARG:HA	1:B:50:ASN:HD22	1.71	0.53
2:A:220:LEU:HD12	2:A:222:ARG:HH12	1.73	0.53
1:B:15:CYS:O	1:B:16:CYS:SG	2.65	0.53
2:A:27:GLN:HG3	2:A:227:ASN:CG	2.33	0.53
1:B:27:GLU:HB3	1:B:36:VAL:HB	1.90	0.53
1:B:6:GLU:CG	1:B:6:GLU:C	2.81	0.53
2:A:155:ALA:N	2:A:156:PRO:HD3	2.23	0.53
1:B:22:GLY:HA2	1:B:75:ARG:O	2.09	0.53
1:B:64:ALA:C	1:B:66:GLY:H	2.16	0.53
1:B:45:TYR:O	1:B:55:CYS:HB3	2.09	0.52
2:A:72:ASN:ND2	2:A:73:ALA:H	2.07	0.52
2:A:95:LEU:HD13	2:A:170:ILE:HD13	1.91	0.52
1:B:80:VAL:CG2	1:B:96:HIS:ND1	2.73	0.51
2:A:224:ILE:HG23	2:A:226[A]:GLU:HG2	1.90	0.51
1:B:5:LYS:HD3	1:B:8:GLU:OE2	2.10	0.51
1:B:96:HIS:CD2	1:B:96:HIS:N	2.76	0.51
2:A:159:GLU:CD	2:A:159:GLU:H	2.18	0.51
2:A:116:THR:CG2	2:A:129:ILE:HB	2.41	0.50
1:B:18:LYS:HD3	1:B:49:LEU:CG	2.42	0.50
1:B:18:LYS:HB2	1:B:50:ASN:O	2.12	0.50
2:A:165:LEU:HD12	2:A:178:GLN:HE21	1.78	0.49
2:A:77:ALA:HB3	2:A:78:PRO:HD3	1.95	0.48
2:A:165:LEU:HD12	2:A:178:GLN:NE2	2.28	0.48
2:A:234:TYR:CZ	2:A:238:ILE:HD11	2.49	0.48
1:B:24:ARG:NH2	1:B:41:PRO:HG3	2.29	0.48
2:A:37:VAL:HG12	2:A:39:HIS:H	1.78	0.48
1:B:62:ASP:OD1	1:B:64:ALA:HB3	2.14	0.47
1:B:37:CYS:HB2	2:A:11:MET:HE1	1.96	0.47
2:A:165:LEU:HD21	2:A:176:ILE:HD12	1.96	0.47
2:A:85:SER:O	2:A:89:ARG:CB	2.61	0.47
2:A:51:PRO:HG3	2:A:176:ILE:HG13	1.96	0.47
1:B:40:CYS:HB3	1:B:44:THR:OG1	2.15	0.47
1:B:6:GLU:C	1:B:6:GLU:HG3	2.40	0.47
2:A:126:ALA:HA	2:A:130:ARG:NH2	2.30	0.47
2:A:47:PRO:HG3	2:A:100:PHE:CE2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:23:TYR:CE1	2:A:11:MET:HE2	2.50	0.46
2:A:173:TRP:CH2	2:A:175:GLU:HB2	2.50	0.46
1:B:49:LEU:N	1:B:49:LEU:CD1	2.78	0.46
2:A:53:LEU:HD23	2:A:54:PRO:HD2	1.96	0.46
1:B:21:PRO:HA	1:B:46:ILE:HG22	1.97	0.46
1:B:44:THR:HB	1:B:55:CYS:HB2	1.96	0.46
2:A:36:ARG:HA	2:A:130:ARG:O	2.15	0.46
2:A:95:LEU:HD13	2:A:170:ILE:CD1	2.45	0.46
2:A:46:ASP:O	2:A:49:GLN:HG2	2.16	0.45
2:A:79:GLN:NE2	2:A:79:GLN:HA	2.31	0.45
1:B:45:TYR:CZ	1:B:78:ASN:HB2	2.50	0.45
2:A:50:PRO:HA	2:A:51:PRO:HD3	1.74	0.45
2:A:64:ARG:HB3	2:A:64:ARG:NH1	2.32	0.45
2:A:51:PRO:C	2:A:53:LEU:H	2.24	0.45
2:A:13:ASP:HB3	2:A:16:ARG:HB2	1.97	0.45
2:A:206:GLN:H	2:A:206:GLN:CD	2.23	0.45
1:B:37:CYS:CB	2:A:11:MET:HE1	2.47	0.44
2:A:77:ALA:HB2	2:A:168:VAL:HG11	1.98	0.44
2:A:37:VAL:O	2:A:131:THR:HA	2.17	0.44
2:A:11:MET:HE3	2:A:27:GLN:HB3	1.98	0.44
2:A:195:LEU:HD23	2:A:195:LEU:HA	1.84	0.44
2:A:51:PRO:HG2	2:A:55:ILE:HD11	2.00	0.43
2:A:15:ASN:O	2:A:18:ARG:HD2	2.18	0.43
2:A:53:LEU:HD23	2:A:54:PRO:N	2.33	0.43
1:B:18:LYS:HE2	1:B:49:LEU:HD23	2.00	0.43
1:B:95:ASP:O	1:B:96:HIS:CD2	2.68	0.43
1:B:45:TYR:O	1:B:46:ILE:HG12	2.18	0.43
2:A:72:ASN:HD22	2:A:73:ALA:N	2.16	0.43
2:A:241:TRP:CH2	2:A:243:GLY:HA3	2.54	0.43
1:B:33:THR:O	2:A:32:PRO:HD3	2.19	0.43
2:A:18:ARG:HD3	2:A:18:ARG:H	1.84	0.42
2:A:220:LEU:HD12	2:A:222:ARG:NH1	2.33	0.42
2:A:112:VAL:CG1	2:A:134:ARG:HB2	2.49	0.42
2:A:95:LEU:HB3	2:A:115:TYR:HB2	2.02	0.42
2:A:167:LEU:C	2:A:167:LEU:CD2	2.92	0.42
1:B:59:GLN:HB2	1:B:78:ASN:HD21	1.83	0.42
2:A:148:ASN:HD22	2:A:148:ASN:HA	1.64	0.42
2:A:62:LEU:HD13	2:A:69:VAL:CG2	2.49	0.42
2:A:55:ILE:HG12	2:A:176:ILE:HB	2.02	0.42
2:A:134:ARG:HD2	2:A:214:VAL:CG1	2.49	0.42
1:B:91:VAL:HG12	1:B:98:ALA:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:2:TYR:HB3	2:A:37:VAL:CG1	2.42	0.41
1:B:7:ASP:HB2	1:B:18:LYS:CE	2.29	0.41
2:A:46:ASP:OD1	2:A:48:PHE:N	2.50	0.41
2:A:53:LEU:HD23	2:A:54:PRO:CD	2.50	0.41
2:A:116:THR:HG23	2:A:129:ILE:HB	2.02	0.41
1:B:80:VAL:CG1	1:B:96:HIS:HB3	2.49	0.41
2:A:99:TRP:CH2	2:A:166:ARG:HB2	2.56	0.41
1:B:62:ASP:HA	1:B:63:PRO:HD3	1.96	0.40
2:A:3:ALA:O	2:A:37:VAL:HG13	2.21	0.40
2:A:25:LEU:CD1	2:A:27:GLN:HE22	2.32	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	B	96/167 (58%)	86 (90%)	7 (7%)	3 (3%)	3	5
2	A	259/290 (89%)	236 (91%)	21 (8%)	2 (1%)	16	27
All	All	355/457 (78%)	322 (91%)	28 (8%)	5 (1%)	9	14

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	29	CYS
1	B	65	MET
2	A	211	GLY
1	B	64	ALA
2	A	40	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	B	82/139 (59%)	77 (94%)	5 (6%)	17	30
2	A	223/248 (90%)	201 (90%)	22 (10%)	7	11
All	All	305/387 (79%)	278 (91%)	27 (9%)	10	15

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	11	VAL
1	B	18	LYS
1	B	46	ILE
1	B	57	GLN
1	B	73	CYS
2	A	4	LEU
2	A	11	MET
2	A	16	ARG
2	A	21	ASP
2	A	24	VAL
2	A	25	LEU
2	A	27	GLN
2	A	69	VAL
2	A	112	VAL
2	A	148	ASN
2	A	149	LEU
2	A	186	LYS
2	A	195	LEU
2	A	202	CYS
2	A	206	GLN
2	A	224	ILE
2	A	226[A]	GLU
2	A	226[B]	GLU
2	A	230	THR
2	A	236	LEU
2	A	257	LEU
2	A	259	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	27	GLN
2	A	72	ASN
2	A	79	GLN
2	A	148	ASN
2	A	178	GLN
2	A	206	GLN
2	A	209	GLN

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 ⓘ

2 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	$\# Z > 2$
3	NAG	C	1	2,3	14,14,15	0.51	0	17,19,21	0.91	1 (5%)
3	NAG	C	2	3	14,14,15	0.49	0	17,19,21	0.78	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
3	NAG	C	1	2,3	-	0/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	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($^{\circ}$)	Ideal($^{\circ}$)
3	C	1	NAG	C2-N2-C7	-2.93	118.98	122.90
3	C	2	NAG	C2-N2-C7	-2.18	119.97	122.90

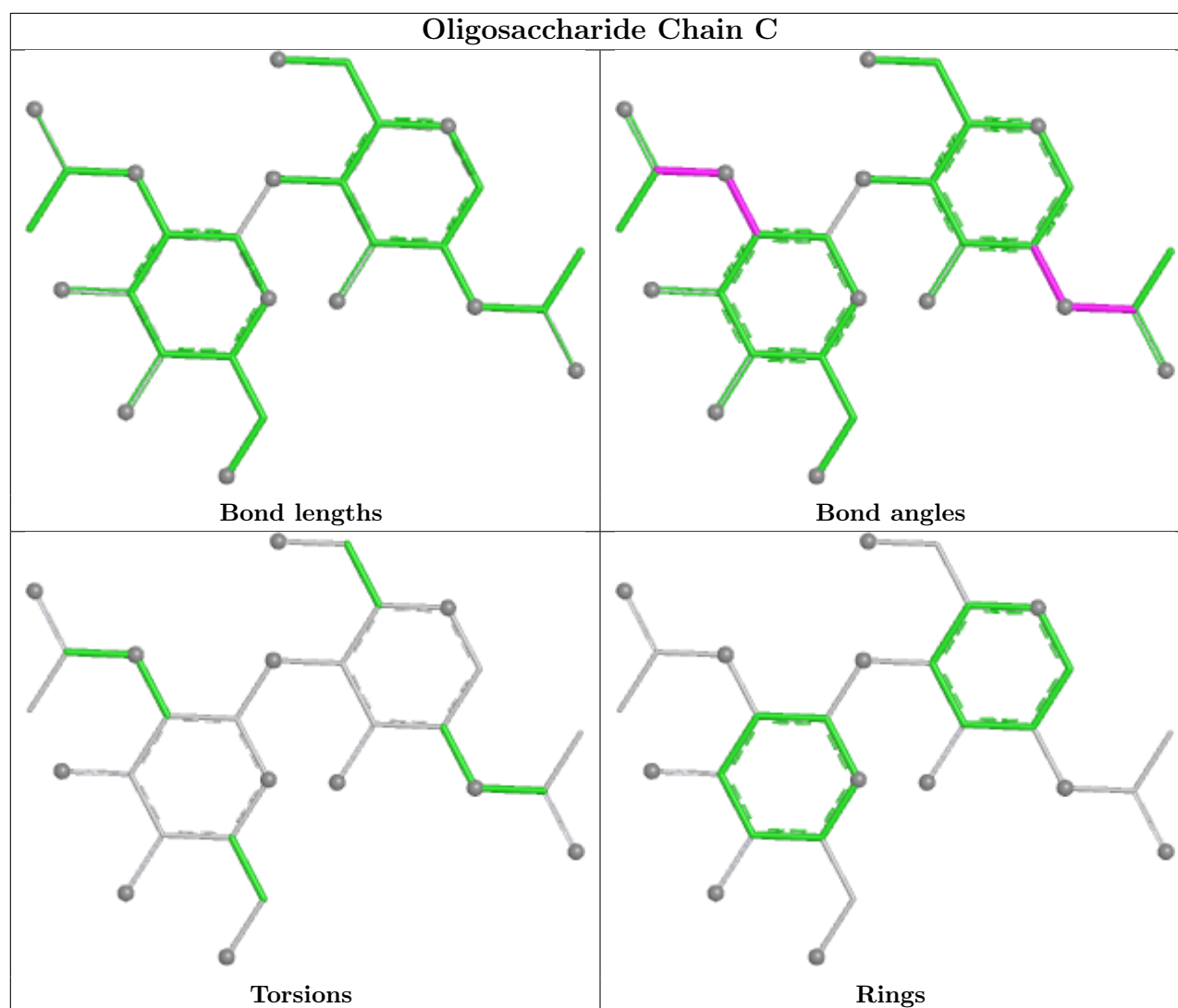
There are no chirality outliers.

There are no torsion outliers.

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](#)

5 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	$\# Z > 2$
4	SO4	A	528	-	4,4,4	0.34	0	6,6,6	0.19	0
4	SO4	B	526	-	4,4,4	0.36	0	6,6,6	0.21	0
4	SO4	B	529	-	4,4,4	0.36	0	6,6,6	0.22	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	SO4	A	527	-	4,4,4	0.35	0	6,6,6	0.14	0
4	SO4	A	525	-	4,4,4	0.38	0	6,6,6	0.19	0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
2	A	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	258:SER	C	259:GLU	N	1.19

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	B	100/167 (59%)	1.39	25 (25%) 2 1	39, 58, 72, 79	0
2	A	259/290 (89%)	0.26	20 (7%) 19 14	16, 41, 61, 70	2 (0%)
All	All	359/457 (78%)	0.58	45 (12%) 8 6	16, 46, 68, 79	2 (0%)

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	96	HIS	7.2
1	B	95	ASP	6.5
1	B	18	LYS	6.3
2	A	86	GLU	6.2
2	A	87	ASP	5.9
2	A	91	GLN	4.5
2	A	18	ARG	4.5
2	A	21	ASP	4.5
1	B	31	GLU	4.3
2	A	35	ARG	4.1
1	B	105	THR	3.9
2	A	196	ARG	3.5
2	A	89	ARG	3.4
1	B	71	ARG	3.1
2	A	20	LYS	3.1
1	B	4	CYS	3.0
1	B	15	CYS	3.0
1	B	46	ILE	3.0
1	B	81	CYS	2.8
1	B	16	CYS	2.7
1	B	60	MET	2.7
1	B	65	MET	2.7
2	A	16	ARG	2.7
2	A	202	CYS	2.6

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Mol	Chain	Res	Type	RSRZ
2	A	66	CYS	2.6
2	A	24	VAL	2.6
2	A	186	LYS	2.5
1	B	73	CYS	2.5
1	B	55	CYS	2.5
1	B	13	SER	2.5
2	A	2	TYR	2.5
1	B	14	GLU	2.4
2	A	259	GLU	2.4
1	B	86	GLY	2.4
2	A	64	ARG	2.3
1	B	83	CYS	2.3
1	B	97	CYS	2.3
1	B	57	GLN	2.3
1	B	29	CYS	2.2
2	A	85	SER	2.2
2	A	121	ASN	2.1
1	B	12	GLY	2.1
1	B	10	PRO	2.1
2	A	82	ARG	2.1
1	B	58	CYS	2.0

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

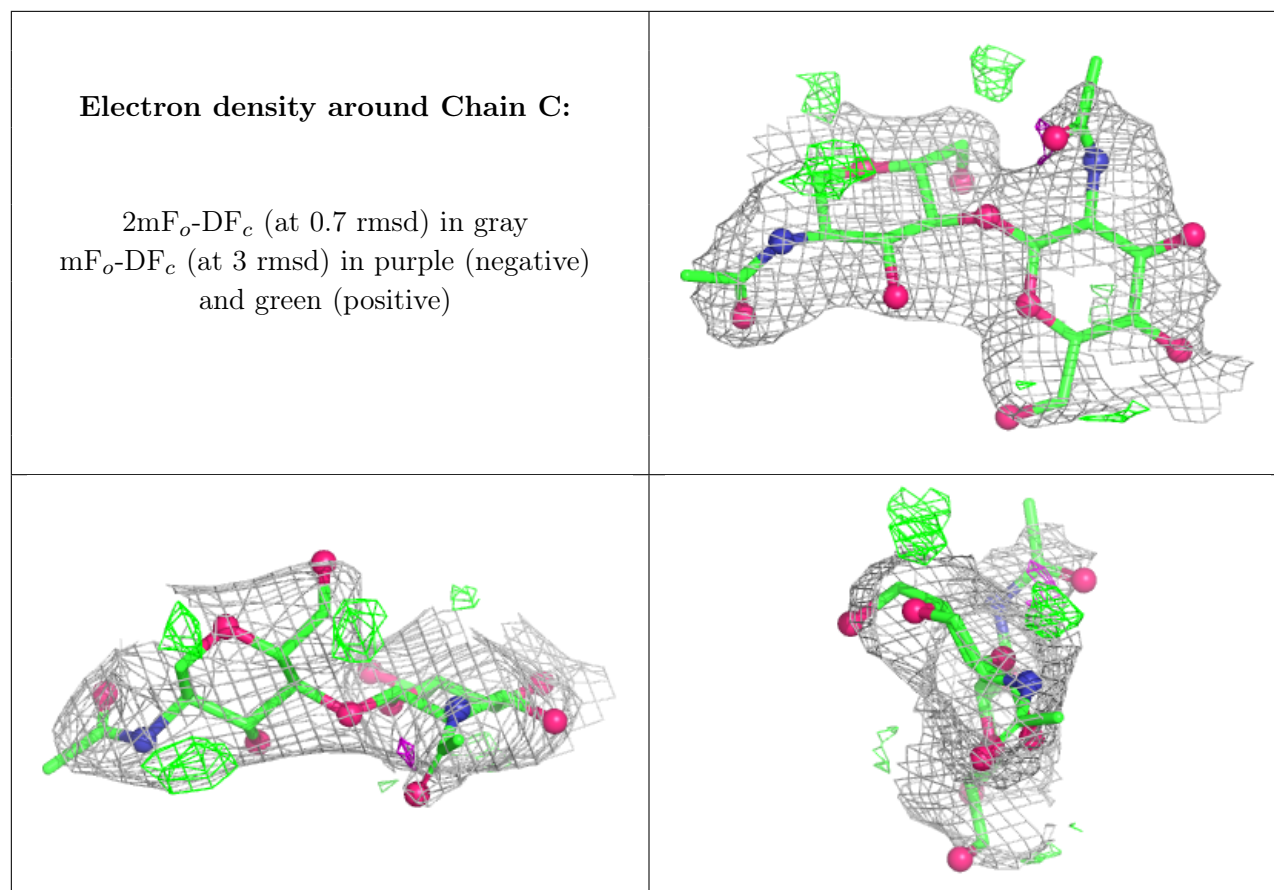
There are no non-standard protein/DNA/RNA residues in this entry.

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
3	NAG	C	2	14/15	0.76	0.15	69,71,75,76	0
3	NAG	C	1	14/15	0.82	0.13	59,61,66,68	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 ⓘ

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	SO4	A	527	5/5	0.69	0.23	102,102,103,103	0
4	SO4	B	526	5/5	0.71	0.34	105,105,106,106	0
4	SO4	B	529	5/5	0.79	0.30	109,110,111,111	0
4	SO4	A	525	5/5	0.80	0.16	100,101,102,102	0
4	SO4	A	528	5/5	0.93	0.25	96,98,98,99	0

6.5 Other polymers ⓘ

There are no such residues in this entry.