



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:20 AM UTC

PDB ID : 2CWG / pdb_00002cwg
Title : CRYSTALLOGRAPHIC REFINEMENT AND STRUCTURE ANALYSIS
OF THE COMPLEX OF WHEAT GERM AGGLUTININ WITH A BIVA-
LENT SIALOGLYCOPEPTIDE FROM GLYCOPHORIN A
Authors : Wright, C.S.
Deposited on : 1993-01-08
Resolution : 2.00 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

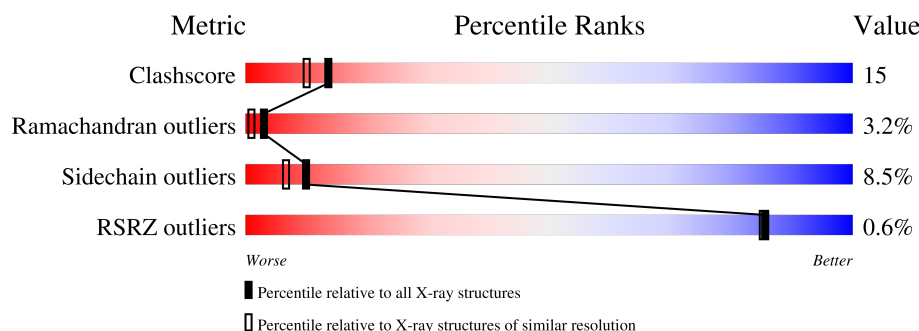
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	171	
1	B	171	
2	D	8	
2	E	8	
3	C	4	
3	F	4	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	NDG	C	1	X	-	-	-
3	NDG	F	1	X	-	-	-

2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2935 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.

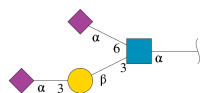
- Molecule 1 is a protein called AGGLUTININ ISOLECTIN 1 (WGA1).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	171	Total	C	N	O	S	0	0	0
			1167	681	214	238	34			
1	B	171	Total	C	N	O	S	0	0	0
			1167	681	214	238	34			

- Molecule 2 is a protein called T5 SIALOGLYCOPEPTIDE OF GLYCOPHORIN A.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	8	Total	C	N	O	0	0	0
			63	38	11	14			
2	E	3	Total	C	N	O	0	0	0
			19	12	3	4			

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)]2-acetamido-2-deoxy-alpha-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	4	Total	C	N	O	0	0	0
			65	36	3	26			
3	F	4	Total	C	N	O	0	0	0
			65	36	3	26			

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	187	Total	O	0	0
			187	187		

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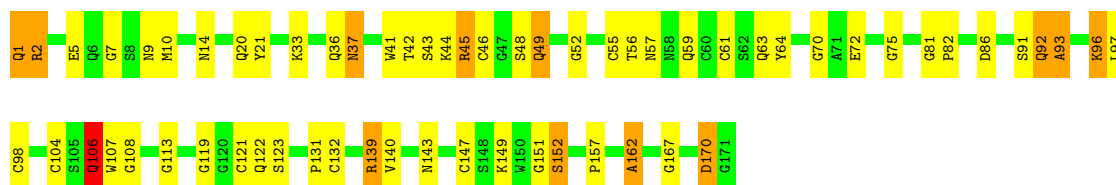
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	D	24	Total 24	O 24	0	0
4	B	176	Total 176	O 176	0	0
4	E	2	Total 2	O 2	0	0

3 Residue-property plots

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

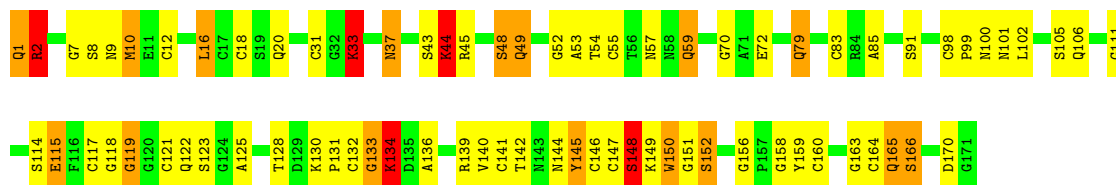
- Molecule 1: AGGLUTININ ISOLECTIN 1 (WGA1)

Chain A: 



- Molecule 1: AGGLUTININ ISOLECTIN 1 (WGA1)

Chain B: 



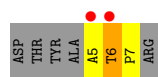
- Molecule 2: T5 SIALOGLYCOPEPTIDE OF GLYCOPHORIN A

Chain D: 



- Molecule 2: T5 SIALOGLYCOPEPTIDE OF GLYCOPHORIN A

Chain E: 



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)]2-acetamido-2-deoxy-alpha-D-glucopyranose

Chain C: 



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-3)-beta-D-galactopyranose-(1-3)-[N-acetyl-alpha-neuraminic acid-(2-6)]2-acetamido-2-deoxy-alpha-D-glucopyranose

Chain F:  50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	111.00Å 50.40Å 63.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00 10.00 – 2.00	Depositor EDS
% Data completeness (in resolution range)	(Not available) (10.00-2.00) 91.8 (10.00-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.74 (at 1.99Å)	Xtriage
Refinement program	PROLSQ, X-PLOR	Depositor
R, R_{free}	0.171 , (Not available) 0.168 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	26.4	Xtriage
Anisotropy	0.254	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , 87.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2935	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.24% 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: SIA, GAL, NDG, PCA

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	1.26	1/1181 (0.1%)	2.21	54/1580 (3.4%)
1	B	1.27	0/1181	2.38	76/1580 (4.8%)
2	D	1.20	0/64	2.33	3/86 (3.5%)
2	E	1.34	0/19	2.04	0/26
All	All	1.26	1/2445 (0.0%)	2.29	133/3272 (4.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	70	GLY	N-CA	5.69	1.50	1.45

All (133) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	59	GLN	OE1-CD-NE2	-9.85	112.75	122.60
1	B	145	TYR	N-CA-C	-9.50	98.07	110.33
1	B	122	GLN	OE1-CD-NE2	-9.35	113.25	122.60
1	B	148	SER	CA-C-N	9.26	139.63	121.58
1	B	148	SER	C-N-CA	9.26	139.63	121.58
1	A	2	ARG	NE-CZ-NH2	8.88	127.19	119.20
1	B	106	GLN	CA-CB-CG	8.87	131.85	114.10
1	B	57	ASN	CA-CB-CG	8.75	121.35	112.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	70	GLY	CA-C-O	-8.64	113.71	122.26
2	D	9	ALA	CA-C-O	8.57	129.75	120.58
1	A	147	CYS	CA-C-O	8.51	129.39	120.54
1	B	45	ARG	NE-CZ-NH1	-7.95	113.55	121.50
1	B	150	TRP	CA-CB-CG	7.93	128.67	113.60
1	B	72	GLU	CB-CG-CD	7.91	126.04	112.60
1	B	45	ARG	NE-CZ-NH2	7.83	126.25	119.20
1	B	117	CYS	CA-C-N	7.79	132.07	121.01
1	B	117	CYS	C-N-CA	7.79	132.07	121.01
1	A	92	GLN	CB-CG-CD	7.67	125.63	112.60
1	B	121	CYS	CA-C-O	-7.58	112.69	120.96
1	B	145	TYR	CA-C-O	-7.51	113.55	122.03
1	A	123	SER	CA-C-O	-7.39	113.55	121.38
1	B	114	SER	CA-CB-OG	-7.39	96.33	111.10
1	A	113	GLY	CA-C-O	-7.32	117.02	122.37
1	A	2	ARG	NE-CZ-NH1	-7.09	114.41	121.50
1	A	86	ASP	CA-CB-CG	-6.98	105.62	112.60
1	A	170	ASP	CA-CB-CG	6.89	119.50	112.60
1	B	130	LYS	CA-C-N	6.85	126.88	119.89
1	B	130	LYS	C-N-CA	6.85	126.88	119.89
1	B	83	CYS	CA-C-O	-6.77	114.43	121.87
1	A	167	GLY	O-C-N	6.71	128.45	122.88
1	A	152	SER	CA-C-O	-6.68	113.56	121.11
2	D	5	ASP	CA-CB-CG	6.67	119.27	112.60
1	B	166	SER	N-CA-C	6.65	119.05	109.14
1	B	2	ARG	NE-CZ-NH1	6.64	128.14	121.50
1	B	149	LYS	CA-CB-CG	6.62	127.35	114.10
1	B	98	CYS	CA-C-N	6.56	126.74	120.31
1	B	98	CYS	C-N-CA	6.56	126.74	120.31
1	B	2	ARG	NH1-CZ-NH2	-6.55	110.78	119.30
1	B	122	GLN	N-CA-C	6.50	121.33	113.41
1	B	156	GLY	O-C-N	6.46	128.23	121.77
1	B	101	ASN	OD1-CG-ND2	-6.46	116.14	122.60
1	B	91	SER	CA-CB-OG	6.41	123.92	111.10
1	B	133	GLY	N-CA-C	6.41	122.66	110.86
1	B	142	THR	N-CA-C	-6.39	101.93	110.55
1	A	46	CYS	CA-C-O	-6.38	114.49	121.31
1	A	45	ARG	CD-NE-CZ	-6.35	115.51	124.40
1	A	121	CYS	CB-CA-C	6.27	119.25	109.84
1	B	100	ASN	CA-C-O	-6.23	114.31	121.54
1	B	163	GLY	N-CA-C	-6.21	106.51	115.27
1	B	145	TYR	O-C-N	6.21	130.00	122.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	2	ARG	CB-CA-C	6.19	121.85	110.10
1	A	157	PRO	CB-CA-C	6.16	121.83	112.21
1	A	64	TYR	CA-C-O	-6.15	111.84	119.28
1	A	63	GLN	CG-CD-NE2	-6.14	107.19	116.40
1	B	20	GLN	CB-CA-C	-6.14	100.60	110.79
1	A	33	LYS	O-C-N	6.13	130.45	122.97
1	B	166	SER	CA-C-N	6.11	127.33	120.42
1	B	166	SER	C-N-CA	6.11	127.33	120.42
1	A	49	GLN	CA-C-O	-6.05	113.26	120.10
1	A	96	LYS	CA-CB-CG	-6.04	102.01	114.10
1	A	36	GLN	CA-C-O	-6.01	112.50	119.78
1	B	31	CYS	O-C-N	5.99	130.26	122.42
1	A	14	ASN	CA-C-O	-5.97	114.61	121.54
1	A	52	GLY	N-CA-C	5.96	123.29	115.36
1	B	165	GLN	CA-C-O	-5.95	114.57	120.82
1	B	43	SER	CA-CB-OG	5.95	123.00	111.10
1	B	9	ASN	CA-C-N	5.93	129.71	120.75
1	B	9	ASN	C-N-CA	5.93	129.71	120.75
1	B	33	LYS	CB-CA-C	-5.89	101.87	110.06
1	B	33	LYS	CA-C-O	-5.88	114.45	120.80
1	A	140	VAL	N-CA-C	-5.87	100.13	109.17
1	B	100	ASN	CA-CB-CG	5.86	118.46	112.60
1	A	139	ARG	CD-NE-CZ	5.86	132.60	124.40
1	A	143	ASN	CA-C-O	-5.85	113.06	120.55
1	A	43	SER	N-CA-CB	5.83	118.96	109.69
1	A	52	GLY	CA-C-N	5.81	129.01	120.82
1	A	52	GLY	C-N-CA	5.81	129.01	120.82
1	A	106	GLN	OE1-CD-NE2	5.79	128.39	122.60
1	B	20	GLN	CA-CB-CG	5.75	125.60	114.10
1	B	158	GLY	N-CA-C	5.75	120.78	114.40
1	B	111	GLY	CA-C-N	5.74	132.24	122.37
1	B	111	GLY	C-N-CA	5.74	132.24	122.37
1	A	139	ARG	NE-CZ-NH1	5.72	127.22	121.50
1	A	122	GLN	OE1-CD-NE2	-5.70	116.90	122.60
1	A	20	GLN	OE1-CD-NE2	5.69	128.29	122.60
1	B	44	LYS	CA-C-O	-5.67	115.30	121.87
1	A	108	GLY	O-C-N	5.67	130.06	122.70
1	A	139	ARG	N-CA-CB	5.64	118.33	110.04
1	B	52	GLY	CA-C-O	-5.64	112.58	119.06
2	D	5	ASP	O-C-N	5.62	131.99	123.00
1	A	113	GLY	N-CA-C	-5.60	106.43	112.08
1	A	75	GLY	CA-C-N	5.58	129.03	121.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	75	GLY	C-N-CA	5.58	129.03	121.05
1	B	85	ALA	CA-C-N	5.58	128.76	120.95
1	B	85	ALA	C-N-CA	5.58	128.76	120.95
1	B	72	GLU	CG-CD-OE1	5.58	131.22	118.40
1	B	54	THR	CA-CB-CG2	5.55	119.94	110.50
1	A	37	ASN	CA-C-O	-5.55	115.50	121.38
1	B	141	CYS	N-CA-C	-5.54	100.87	109.24
1	A	57	ASN	CA-C-O	-5.53	114.17	121.08
1	B	134	LYS	N-CA-C	5.50	117.74	111.02
1	A	98	CYS	CA-C-O	5.46	126.26	120.64
1	B	114	SER	CA-C-O	-5.45	112.90	119.49
1	B	44	LYS	N-CA-C	-5.43	103.36	110.53
1	A	162	ALA	CA-C-O	-5.41	112.77	120.51
1	A	140	VAL	CB-CA-C	5.41	118.95	111.38
1	B	48	SER	CA-C-N	5.41	128.27	120.38
1	B	48	SER	C-N-CA	5.41	128.27	120.38
1	A	81	GLY	N-CA-C	-5.40	101.33	112.34
1	B	102	LEU	N-CA-CB	5.38	117.92	110.17
1	A	123	SER	N-CA-CB	-5.33	103.25	111.46
1	B	128	THR	CA-CB-OG1	-5.33	101.61	109.60
1	A	5	GLU	CB-CG-CD	5.32	121.65	112.60
1	B	16	LEU	N-CA-C	-5.31	103.13	110.35
1	A	147	CYS	CB-CA-C	5.28	119.23	110.78
1	A	61	CYS	CA-C-O	5.28	126.11	120.46
1	B	105	SER	O-C-N	5.26	128.85	122.85
1	B	166	SER	CA-C-O	5.20	127.33	121.51
1	B	53	ALA	CA-C-N	5.18	130.40	120.97
1	B	53	ALA	C-N-CA	5.18	130.40	120.97
1	B	150	TRP	N-CA-C	-5.18	106.36	113.30
1	A	43	SER	CA-CB-OG	5.17	121.44	111.10
1	A	104	CYS	CA-C-O	-5.16	115.18	120.54
1	A	72	GLU	CG-CD-OE1	5.15	130.25	118.40
1	B	165	GLN	N-CA-C	5.14	116.57	111.07
1	A	5	GLU	CA-CB-CG	5.13	124.36	114.10
1	A	56	THR	N-CA-C	-5.11	103.19	110.59
1	B	123	SER	CA-C-O	-5.10	115.83	121.28
1	B	134	LYS	CA-C-O	-5.09	115.47	121.07
1	A	41	TRP	CB-CA-C	5.06	120.39	110.67
1	A	42	THR	CA-CB-OG1	-5.06	102.02	109.60
1	B	115	GLU	CB-CG-CD	5.05	121.18	112.60
1	B	8	SER	CA-C-O	-5.03	115.70	121.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	2	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	A	1167	0	999	24	0
1	B	1167	0	999	33	0
2	D	63	0	55	9	0
2	E	19	0	18	8	0
3	C	65	0	53	2	0
3	F	65	0	53	6	0
4	A	187	0	0	6	0
4	B	176	0	0	9	0
4	D	24	0	0	0	0
4	E	2	0	0	0	0
All	All	2935	0	2177	69	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 15.

All (69) 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:A:2:ARG:O	1:B:10:MET:HE3	1.63	0.99
2:E:6:THR:OG1	3:F:1:NDG:C1	2.12	0.98
2:D:6:THR:O	2:D:7:TYR:HB2	1.65	0.94
1:B:1:PCA:HB3	2:E:7:PRO:O	1.72	0.88
2:E:6:THR:N	2:E:7:PRO:HD3	1.92	0.83
2:D:5:ASP:O	2:D:6:THR:HG23	1.82	0.79
1:A:2:ARG:C	1:B:10:MET:HE3	2.09	0.78
4:B:866:HOH:O	3:F:1:NDG:H8C1	1.90	0.71
1:B:48:SER:H	1:B:79:GLN:HE22	1.39	0.71
1:B:1:PCA:CB	2:E:7:PRO:O	2.39	0.69
1:A:2:ARG:NH2	1:A:21:TYR:O	2.26	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:866:HOH:O	3:F:1:NDG:C8	2.40	0.67
1:A:170:ASP:O	4:A:865:HOH:O	2.12	0.67
1:A:162:ALA:HB3	4:A:877:HOH:O	1.95	0.66
1:B:44:LYS:HG3	4:B:839:HOH:O	1.96	0.65
2:D:7:TYR:O	2:D:8:ALA:HB3	1.99	0.63
4:B:732:HOH:O	3:F:1:NDG:H6C1	2.01	0.61
2:E:6:THR:N	2:E:7:PRO:CD	2.65	0.60
1:B:7:GLY:O	1:B:10:MET:HB2	2.02	0.60
1:A:96:LYS:HG2	1:A:97:LEU:N	2.16	0.59
1:A:1:PCA:HB2	1:A:2:ARG:HG2	1.85	0.58
1:A:132:CYS:HA	1:A:139:ARG:HG3	1.84	0.58
1:A:2:ARG:C	1:B:10:MET:CE	2.78	0.56
2:D:5:ASP:HB3	1:B:115:GLU:HG3	1.88	0.56
1:A:7:GLY:O	1:A:10:MET:HB2	2.05	0.56
2:D:12:ARG:HB3	3:C:1:NDG:H8C3	1.88	0.55
1:B:132:CYS:SG	1:B:151:GLY:C	2.90	0.55
1:B:49:GLN:HE21	1:B:49:GLN:H	1.53	0.55
1:B:144:ASN:C	1:B:145:TYR:O	2.47	0.53
2:E:5:ALA:N	3:F:2:GAL:HO2	2.06	0.52
1:B:133:GLY:N	1:B:139:ARG:O	2.35	0.52
1:A:91:SER:C	1:A:93:ALA:H	2.18	0.51
1:B:132:CYS:SG	1:B:151:GLY:HA2	2.50	0.51
1:B:18:CYS:HB2	1:B:37:ASN:ND2	2.27	0.50
1:B:49:GLN:H	1:B:49:GLN:NE2	2.10	0.50
1:B:132:CYS:SG	1:B:152:SER:N	2.85	0.49
2:D:9:ALA:HB1	3:C:3:SIA:O9	2.13	0.49
2:E:5:ALA:C	2:E:7:PRO:HD3	2.37	0.49
1:A:131:PRO:HA	1:A:151:GLY:O	2.12	0.48
1:B:33:LYS:HD3	1:B:33:LYS:HA	1.68	0.47
1:B:118:GLY:O	1:B:119:GLY:C	2.58	0.47
1:B:147:CYS:SG	1:B:151:GLY:HA2	2.55	0.46
2:D:6:THR:O	2:D:7:TYR:CB	2.50	0.46
1:B:160:CYS:O	1:B:170:ASP:HB2	2.16	0.46
1:B:134:LYS:N	4:B:669:HOH:O	2.49	0.46
1:A:55:CYS:HB3	1:A:59:GLN:HB2	1.98	0.45
1:A:82:PRO:HA	4:A:879:HOH:O	2.16	0.45
1:B:99:PRO:HG3	4:B:867:HOH:O	2.17	0.45
2:D:7:TYR:O	2:D:8:ALA:CB	2.64	0.44
2:E:5:ALA:N	3:F:2:GAL:O2	2.50	0.44
1:B:44:LYS:HD2	4:B:839:HOH:O	2.17	0.44
1:A:10:MET:CG	1:B:10:MET:HG2	2.48	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:55:CYS:HB3	1:B:59:GLN:HB2	1.99	0.43
1:A:10:MET:HG3	1:B:10:MET:HG2	2.00	0.43
1:B:12:CYS:HB2	4:B:599:HOH:O	2.18	0.43
1:A:106:GLN:HG2	4:A:540:HOH:O	2.18	0.43
1:A:131:PRO:O	1:A:139:ARG:HD3	2.18	0.43
1:A:45:ARG:NH2	1:A:49:GLN:HG2	2.34	0.42
1:B:147:CYS:O	1:B:148:SER:O	2.37	0.42
2:D:10:THR:HA	2:D:11:PRO:HD3	1.89	0.41
1:B:146:CYS:HB3	1:B:159:TYR:O	2.19	0.41
1:A:106:GLN:HG3	1:A:107:TRP:CE3	2.56	0.41
1:A:132:CYS:HA	1:A:139:ARG:CG	2.49	0.41
1:A:106:GLN:HB2	4:A:670:HOH:O	2.20	0.41
1:A:2:ARG:HE	1:A:2:ARG:HB3	1.71	0.41
1:B:134:LYS:C	1:B:136:ALA:H	2.29	0.40
1:A:9:ASN:ND2	4:A:546:HOH:O	2.51	0.40
1:B:37:ASN:ND2	4:B:599:HOH:O	2.49	0.40
1:B:147:CYS:C	1:B:148:SER:O	2.63	0.40

There are no symmetry-related clashes.

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	169/171 (99%)	159 (94%)	7 (4%)	3 (2%)	6	3
1	B	169/171 (99%)	154 (91%)	10 (6%)	5 (3%)	3	1
2	D	6/8 (75%)	3 (50%)	1 (17%)	2 (33%)	0	0
2	E	1/8 (12%)	0	0	1 (100%)	0	0
All	All	345/358 (96%)	316 (92%)	18 (5%)	11 (3%)	3	1

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	7	TYR
1	B	140	VAL
1	B	148	SER
1	A	92	GLN
1	A	119	GLY
2	D	6	THR
1	B	119	GLY
1	B	125	ALA
1	A	93	ALA
1	B	164	CYS
2	E	6	THR

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	119/119 (100%)	113 (95%)	6 (5%)	22	20
1	B	119/119 (100%)	105 (88%)	14 (12%)	5	3
2	D	6/6 (100%)	5 (83%)	1 (17%)	2	1
2	E	2/6 (33%)	2 (100%)	0	100	100
All	All	246/250 (98%)	225 (92%)	21 (8%)	10	7

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	44	LYS
1	A	48	SER
1	A	106	GLN
1	A	149	LYS
1	A	152	SER
2	D	7	TYR
1	B	2	ARG
1	B	10	MET
1	B	16	LEU
1	B	33	LYS

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Mol	Chain	Res	Type
1	B	37	ASN
1	B	44	LYS
1	B	49	GLN
1	B	79	GLN
1	B	131	PRO
1	B	134	LYS
1	B	150	TRP
1	B	152	SER
1	B	165	GLN
1	B	166	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	14	ASN
1	A	37	ASN
1	A	59	GLN
1	A	63	GLN
1	A	106	GLN
1	B	37	ASN
1	B	49	GLN
1	B	79	GLN
1	B	100	ASN
1	B	165	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

2 non-standard protein/DNA/RNA residues 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
1	PCA	B	1	1	7,8,9	0.85	0	9,10,12	2.00	2 (22%)
1	PCA	A	1	1	7,8,9	1.14	1 (14%)	9,10,12	2.37	4 (44%)

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
1	PCA	B	1	1	-	0/0/11/13	0/1/1/1
1	PCA	A	1	1	-	0/0/11/13	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	PCA	O-C	2.05	1.27	1.20

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	1	PCA	CB-CA-C	-4.78	106.10	112.66
1	B	1	PCA	CB-CA-C	4.65	119.04	112.66
1	A	1	PCA	O-C-CA	-3.41	116.00	124.77
1	B	1	PCA	O-C-CA	-3.01	117.02	124.77
1	A	1	PCA	OE-CD-CG	-2.83	121.67	126.72
1	A	1	PCA	OE-CD-N	2.39	130.12	124.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	B	1	PCA	2	0
1	A	1	PCA	1	0

5.5 Carbohydrates

8 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	NDG	C	1	3	14,14,15	1.12	2 (14%)	17,19,21	2.11	6 (35%)
3	GAL	C	2	3	11,11,12	1.01	0	15,15,17	1.18	3 (20%)
3	SIA	C	3	3	20,20,21	1.26	3 (15%)	21,28,31	1.63	5 (23%)
3	SIA	C	4	3	20,20,21	1.32	2 (10%)	21,28,31	2.04	8 (38%)
3	NDG	F	1	3	14,14,15	0.90	1 (7%)	17,19,21	2.28	8 (47%)
3	GAL	F	2	3	11,11,12	1.36	1 (9%)	15,15,17	2.38	4 (26%)
3	SIA	F	3	3	20,20,21	1.32	3 (15%)	21,28,31	1.78	5 (23%)
3	SIA	F	4	3	20,20,21	1.49	4 (20%)	21,28,31	1.59	6 (28%)

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	NDG	C	1	3	1/1/5/7	0/6/23/26	0/1/1/1
3	GAL	C	2	3	-	0/2/19/22	0/1/1/1
3	SIA	C	3	3	-	0/18/34/38	0/1/1/1
3	SIA	C	4	3	-	3/18/34/38	0/1/1/1
3	NDG	F	1	3	1/1/5/7	0/6/23/26	0/1/1/1
3	GAL	F	2	3	-	2/2/19/22	0/1/1/1
3	SIA	F	3	3	-	7/18/34/38	0/1/1/1
3	SIA	F	4	3	-	2/18/34/38	0/1/1/1

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	3	SIA	C3-C2	3.54	1.58	1.52
3	C	4	SIA	C4-C5	-3.28	1.50	1.53
3	F	2	GAL	O3-C3	-3.25	1.34	1.43
3	F	4	SIA	O1A-C1	3.17	1.31	1.22
3	F	4	SIA	O1B-C1	-3.12	1.20	1.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	F	4	SIA	C6-C5	2.81	1.57	1.53
3	F	1	NDG	C2-N2	2.76	1.50	1.46
3	F	3	SIA	O10-C10	2.43	1.28	1.23
3	C	1	NDG	O5-C1	-2.41	1.39	1.43
3	C	3	SIA	O6-C6	-2.39	1.40	1.44
3	C	3	SIA	O10-C10	2.38	1.28	1.23
3	F	3	SIA	O6-C6	-2.36	1.40	1.44
3	C	3	SIA	C3-C4	-2.14	1.48	1.52
3	C	4	SIA	C3-C2	2.09	1.55	1.52
3	C	1	NDG	C1-C2	2.01	1.55	1.52
3	F	4	SIA	O6-C2	-2.01	1.39	1.43

All (45) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	2	GAL	O3-C3-C4	5.85	124.17	110.38
3	F	1	NDG	O4-C4-C5	-5.12	96.72	109.32
3	C	1	NDG	C1-O5-C5	4.68	118.45	112.19
3	F	2	GAL	O3-C3-C2	4.61	119.46	110.05
3	F	1	NDG	C3-C4-C5	-4.44	102.19	110.23
3	C	4	SIA	O6-C2-C1	4.25	115.73	107.72
3	C	3	SIA	O9-C9-C8	-3.85	103.06	111.16
3	C	4	SIA	O1B-C1-C2	3.73	122.41	112.71
3	F	2	GAL	C6-C5-C4	3.44	121.45	113.02
3	F	3	SIA	O1B-C1-O1A	-3.39	116.39	124.08
3	C	4	SIA	C8-C7-C6	-3.37	106.72	113.05
3	F	3	SIA	C5-N5-C10	3.36	130.97	123.11
3	F	1	NDG	O5-C5-C4	-3.26	102.91	110.83
3	F	3	SIA	O6-C2-C3	-2.99	106.53	110.56
3	C	1	NDG	O7-C7-C8	-2.94	116.83	122.05
3	C	1	NDG	O4-C4-C3	2.93	117.27	110.38
3	C	4	SIA	O9-C9-C8	-2.92	105.02	111.16
3	C	1	NDG	C2-N2-C7	2.92	126.81	122.90
3	F	4	SIA	C11-C10-N5	-2.89	111.33	116.12
3	F	4	SIA	O6-C2-C1	2.79	112.98	107.72
3	C	1	NDG	O4-C4-C5	-2.78	102.47	109.32
3	C	4	SIA	C6-C5-N5	-2.66	106.66	110.91
3	C	3	SIA	C9-C8-C7	-2.56	106.96	112.17
3	C	1	NDG	C1-C2-N2	2.55	114.46	110.43
3	F	3	SIA	O6-C2-C1	2.53	112.50	107.72
3	C	3	SIA	O4-C4-C3	2.53	116.14	109.86
3	C	4	SIA	C5-N5-C10	2.49	128.93	123.11

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	GAL	O2-C2-C1	-2.46	103.58	109.22
3	F	1	NDG	C6-C5-C4	-2.43	107.04	113.02
3	F	4	SIA	O8-C8-C7	-2.43	103.57	109.25
3	C	3	SIA	O8-C8-C9	-2.40	103.58	109.03
3	C	4	SIA	O1B-C1-O1A	-2.39	118.66	124.08
3	F	3	SIA	O7-C7-C8	2.38	114.33	108.93
3	F	1	NDG	C2-N2-C7	-2.37	119.73	122.90
3	F	4	SIA	C4-C5-N5	2.34	115.04	110.44
3	F	1	NDG	C1-C2-N2	-2.33	106.76	110.43
3	C	3	SIA	O7-C7-C6	2.28	114.37	109.44
3	C	2	GAL	C1-C2-C3	-2.22	106.42	109.64
3	F	1	NDG	O7-C7-N2	-2.17	118.15	121.98
3	F	2	GAL	O5-C1-C2	2.17	115.96	110.79
3	F	4	SIA	C5-N5-C10	2.16	128.15	123.11
3	C	2	GAL	O5-C5-C6	-2.10	103.57	107.66
3	C	4	SIA	O1A-C1-C2	-2.08	118.35	122.85
3	F	4	SIA	C6-C5-N5	2.08	114.22	110.91
3	F	1	NDG	C8-C7-N2	2.04	119.51	116.12

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
3	C	1	NDG	C4
3	F	1	NDG	C4

All (14) torsion outliers are listed below:

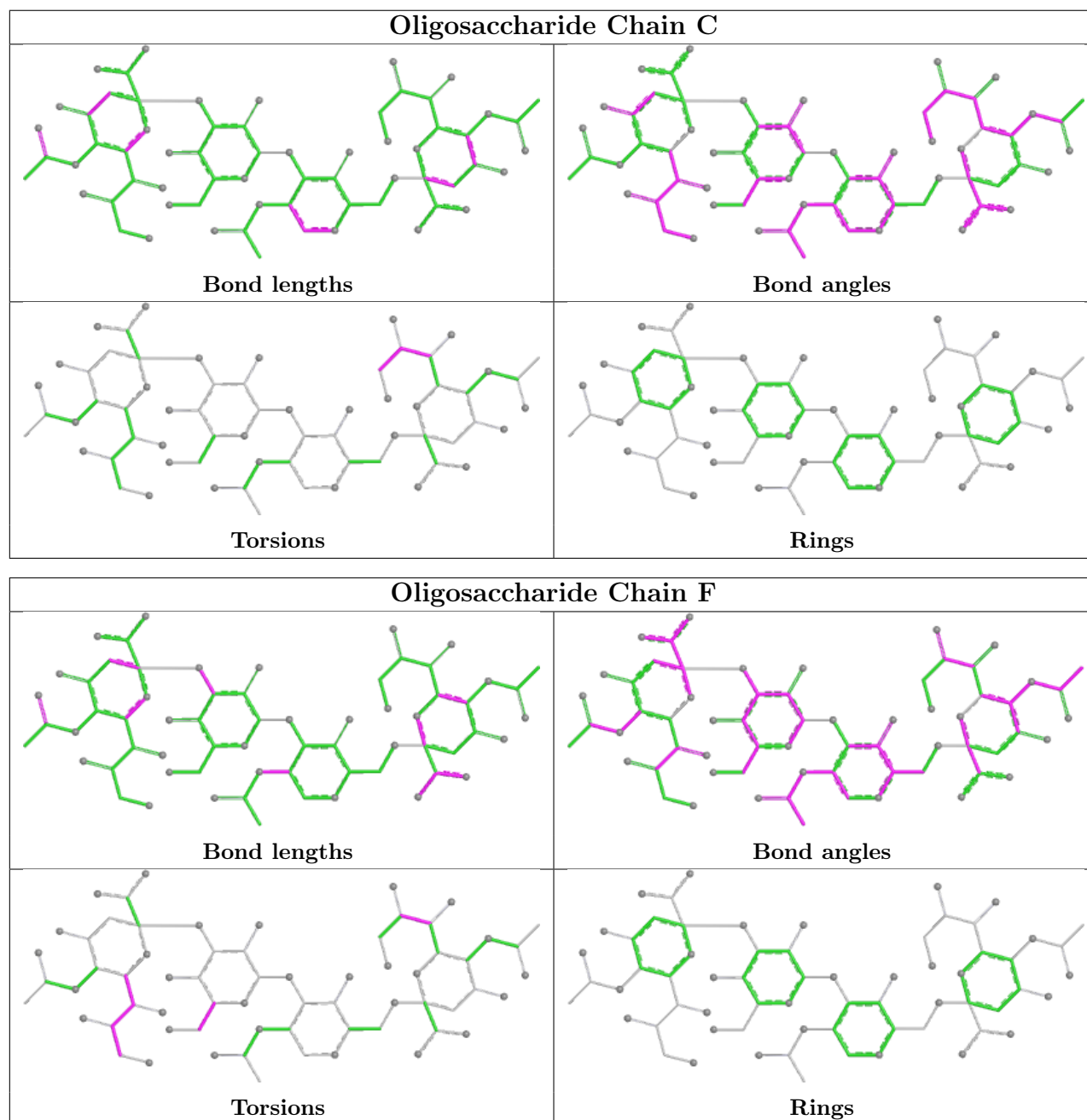
Mol	Chain	Res	Type	Atoms
3	F	3	SIA	C5-C6-C7-C8
3	F	3	SIA	C5-C6-C7-O7
3	F	3	SIA	O6-C6-C7-C8
3	F	3	SIA	O6-C6-C7-O7
3	F	3	SIA	C7-C8-C9-O9
3	F	3	SIA	O8-C8-C9-O9
3	C	4	SIA	O8-C8-C9-O9
3	C	4	SIA	C7-C8-C9-O9
3	F	2	GAL	C4-C5-C6-O6
3	F	2	GAL	O5-C5-C6-O6
3	F	3	SIA	C6-C7-C8-O8
3	F	4	SIA	O7-C7-C8-O8
3	F	4	SIA	C6-C7-C8-O8
3	C	4	SIA	O7-C7-C8-C9

There are no ring outliers.

4 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	3	SIA	1	0
3	F	2	GAL	2	0
3	C	1	NDG	1	0
3	F	1	NDG	4	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 [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	170/171 (99%)	-0.90	0	100 100	11, 22, 40, 55	0
1	B	170/171 (99%)	-0.67	0	100 100	13, 25, 56, 67	0
2	D	8/8 (100%)	0.50	0	100 100	51, 74, 87, 92	0
2	E	3/8 (37%)	2.46	2 (66%)	0 0	99, 99, 99, 100	0
All	All	351/358 (98%)	-0.73	2 (0%)	85 85	11, 24, 56, 100	0

All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	E	5	ALA	3.3
2	E	6	THR	2.9

6.2 Non-standard residues in protein, DNA, RNA chains [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.

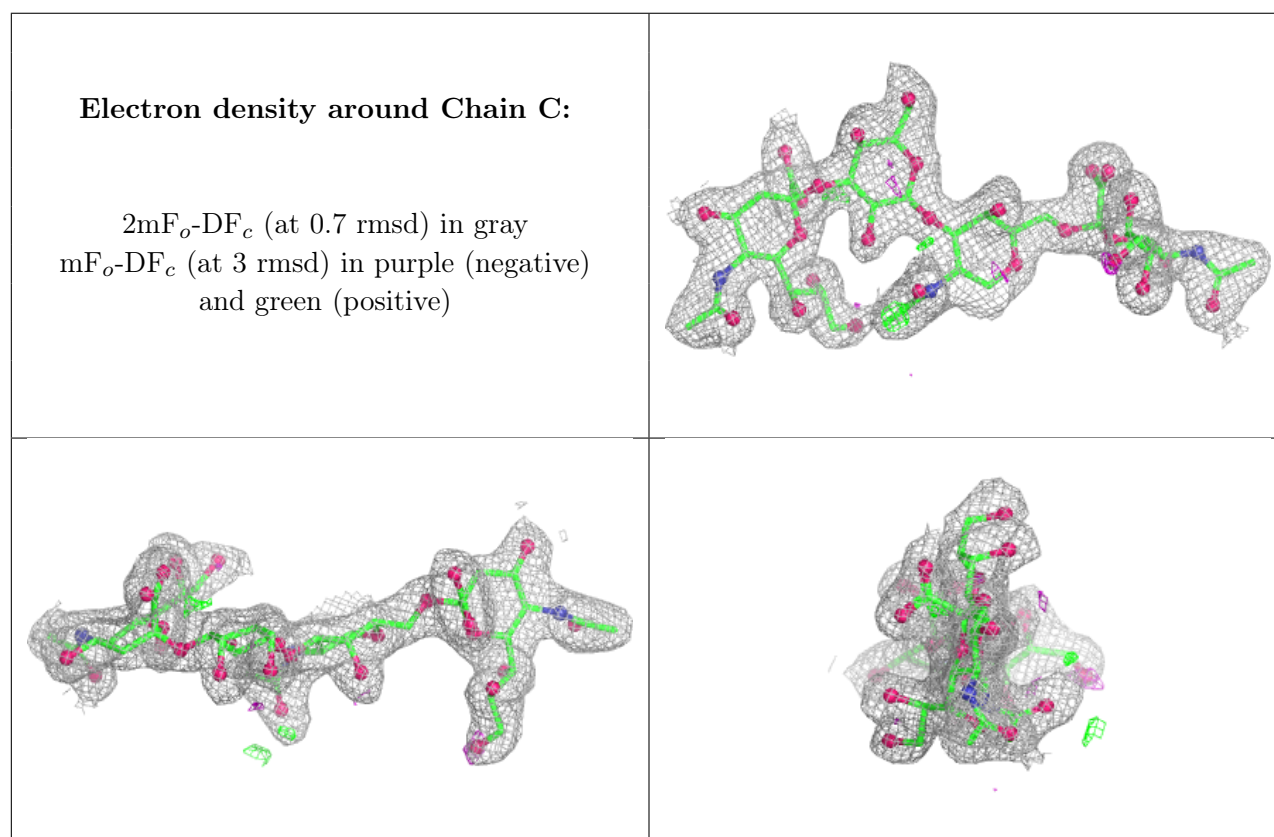
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	PCA	A	1	8/9	0.86	0.07	35,36,37,38	0
1	PCA	B	1	8/9	0.92	0.05	29,32,34,34	0

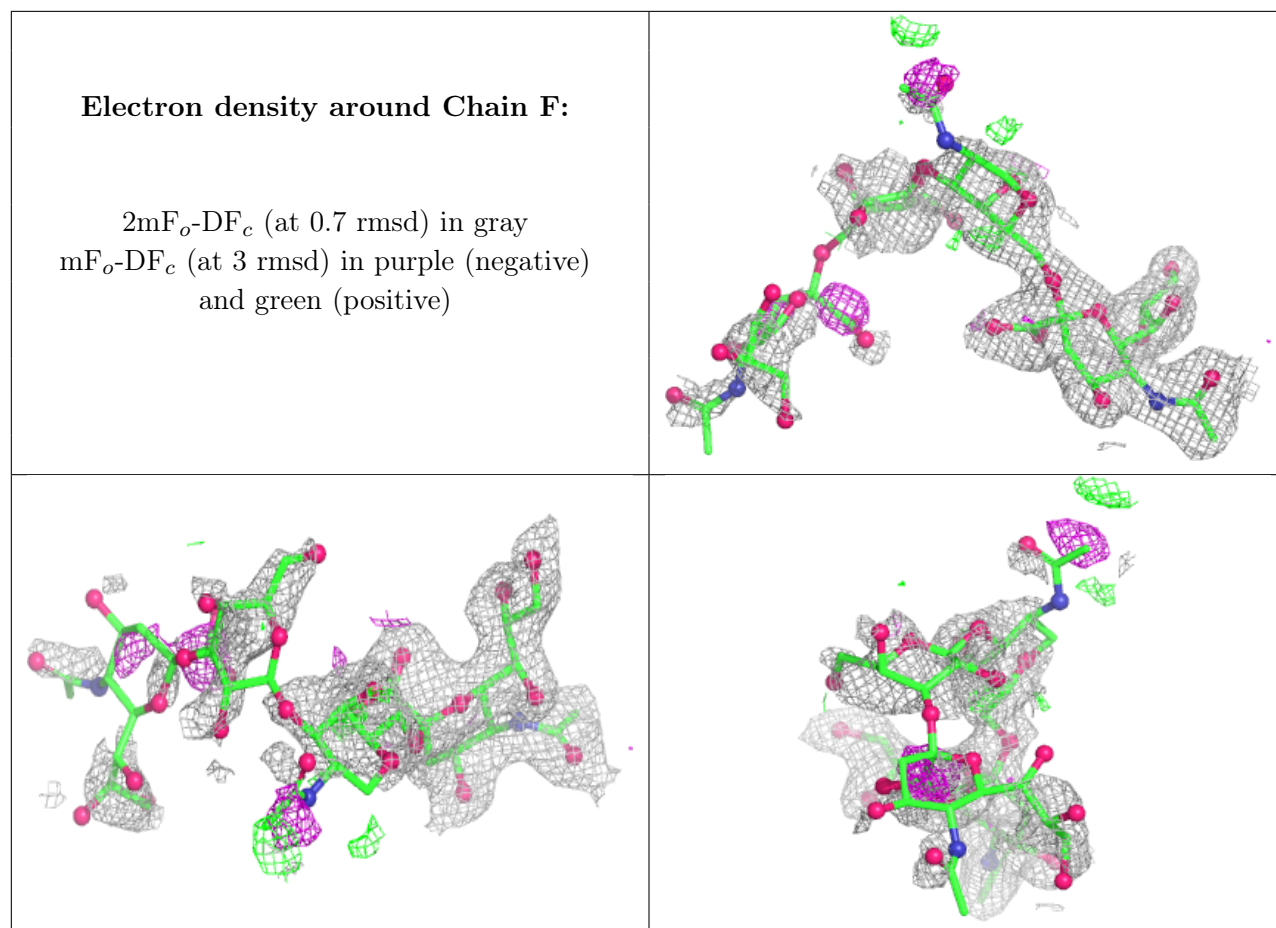
6.3 Carbohydrates [i](#)

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
3	NDG	C	1	14/15	-	-	18,24,27,29	0
3	GAL	C	2	11/12	-	-	20,29,36,40	0
3	SIA	C	3	20/21	-	-	16,21,28,30	0
3	SIA	C	4	20/21	-	-	13,20,30,36	0
3	GAL	F	2	11/12	0.35	0.11	93,96,99,99	0
3	SIA	F	3	20/21	0.51	0.10	99,99,99,99	0
3	NDG	F	1	14/15	0.63	0.13	52,77,86,86	0
3	SIA	F	4	20/21	0.89	0.07	26,42,46,48	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.