



Full wwPDB X-ray Structure Validation Report ⓘ

May 7, 2026 – 09:25 AM EDT

PDB ID : 2J3U / pdb_00002j3u
Title : L-ficolin complexed to galactose
Authors : Garlatti, V.; Gaboriaud, C.
Deposited on : 2006-08-23
Resolution : 2.15 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
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

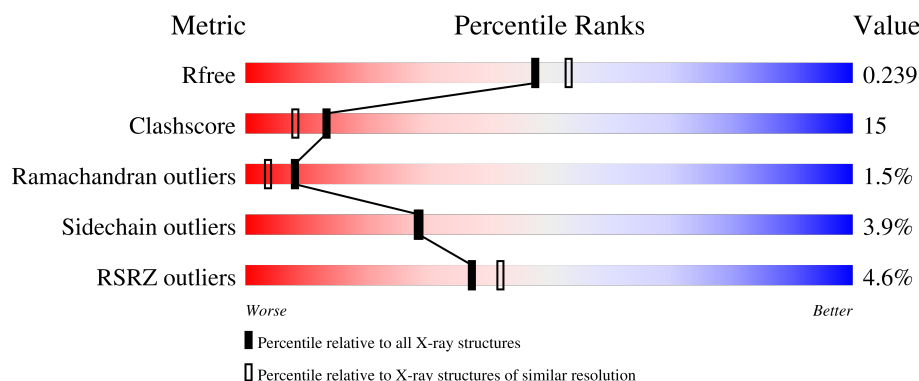
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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	2057 (2.16-2.16)
Clashscore	190562	2159 (2.16-2.16)
Ramachandran outliers	187476	2134 (2.16-2.16)
Sidechain outliers	187428	2133 (2.16-2.16)
RSRZ outliers	180081	2059 (2.16-2.16)

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	218	13% 65% 27% 6% .
1	B	218	81% 19%
1	C	218	83% 14% ..
1	D	218	11% 68% 25% 5% ..
1	E	218	79% 21%

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Mol	Chain	Length	Quality of chain
1	F	218	
2	G	2	
2	H	2	

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
4	ACT	B	1290	-	-	X	-
4	ACT	E	1290	-	-	X	-
6	GLA	E	1289	-	-	X	-

2 Entry composition

There are 10 unique types of molecules in this entry. The entry contains 11129 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 FICOLIN-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	217	Total	C	N	O	S	1	3	0
			1767	1112	312	334	9			
1	B	218	Total	C	N	O	S	3	1	0
			1752	1101	310	332	9			
1	C	214	Total	C	N	O	S	5	1	0
			1721	1082	302	329	8			
1	D	216	Total	C	N	O	S	2	1	0
			1738	1092	306	331	9			
1	E	218	Total	C	N	O	S	0	1	0
			1752	1100	309	334	9			
1	F	217	Total	C	N	O	S	0	0	0
			1736	1092	305	330	9			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	168	THR	VAL	conflict	UNP Q15485
A	247	THR	VAL	conflict	UNP Q15485
B	168	THR	VAL	conflict	UNP Q15485
B	247	THR	VAL	conflict	UNP Q15485
C	168	THR	VAL	conflict	UNP Q15485
C	247	THR	VAL	conflict	UNP Q15485
D	168	THR	VAL	conflict	UNP Q15485
D	247	THR	VAL	conflict	UNP Q15485
E	168	THR	VAL	conflict	UNP Q15485
E	247	THR	VAL	conflict	UNP Q15485
F	168	THR	VAL	conflict	UNP Q15485
F	247	THR	VAL	conflict	UNP Q15485

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

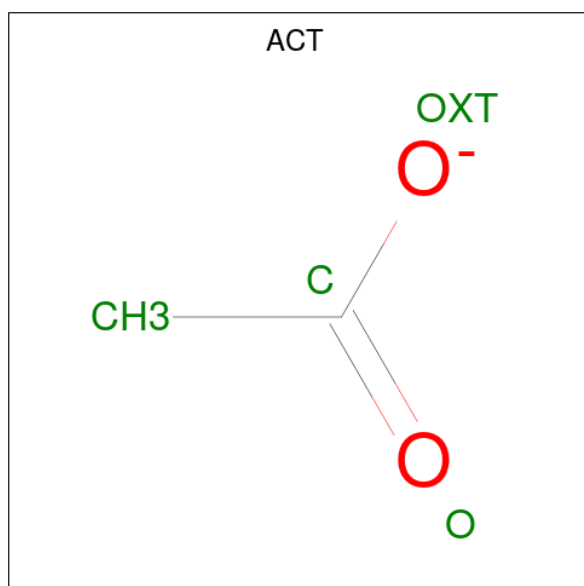


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is CALCIUM ION (CCD ID: CA) (formula: Ca).

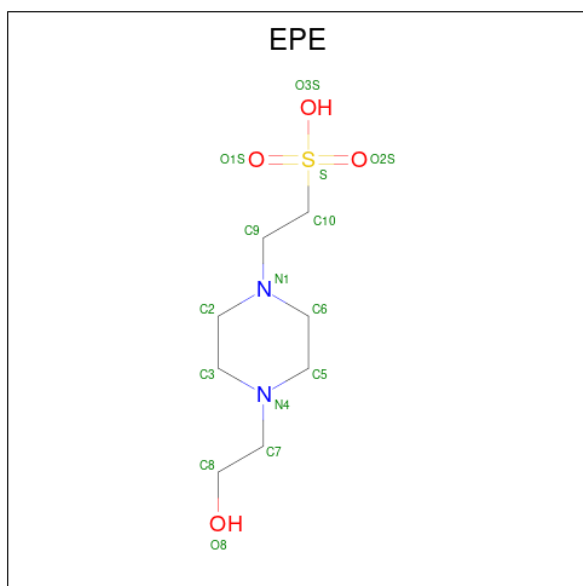
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Ca	0	0
			1	1		
3	B	1	Total	Ca	0	0
			1	1		
3	C	1	Total	Ca	0	0
			1	1		
3	D	1	Total	Ca	0	0
			1	1		
3	E	1	Total	Ca	0	0
			1	1		
3	F	1	Total	Ca	0	0
			1	1		

- Molecule 4 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



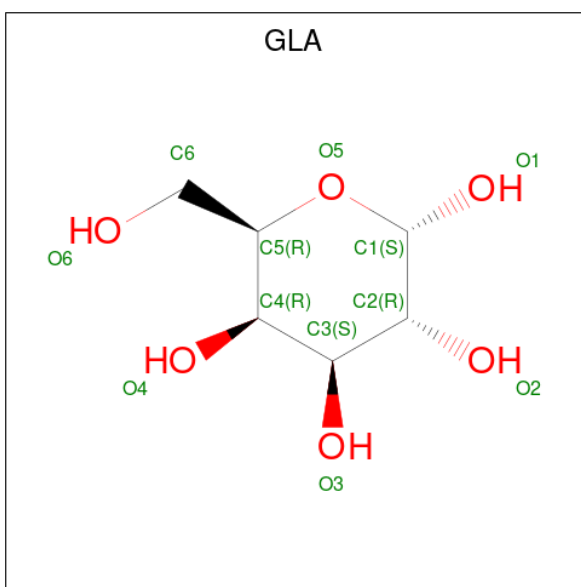
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total C O 4 2 2	0	0
4	B	1	Total C O 4 2 2	0	0
4	C	1	Total C O 4 2 2	0	0
4	E	1	Total C O 4 2 2	0	0
4	F	1	Total C O 4 2 2	0	0

- Molecule 5 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (CCD ID: EPE) (formula: $C_8H_{18}N_2O_4S$).



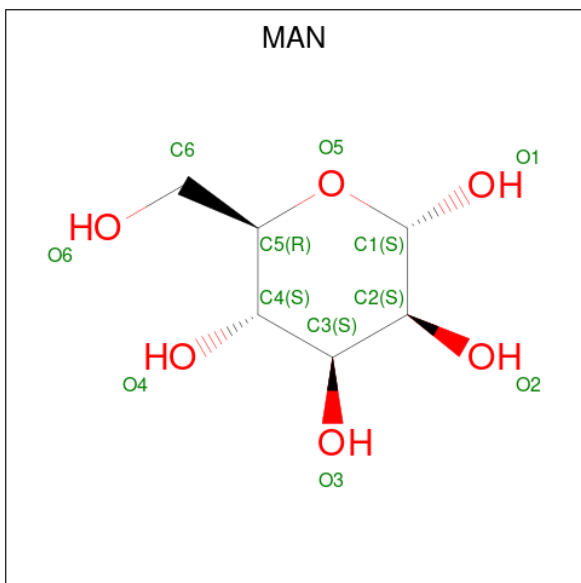
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	B	1	Total C N O S 12 6 2 3 1	0	0
5	C	1	Total C N O S 12 6 2 3 1	0	0
5	F	1	Total C N O S 12 6 2 3 1	0	0

- Molecule 6 is alpha-D-galactopyranose (CCD ID: GLA) (formula: $C_6H_{12}O_6$).



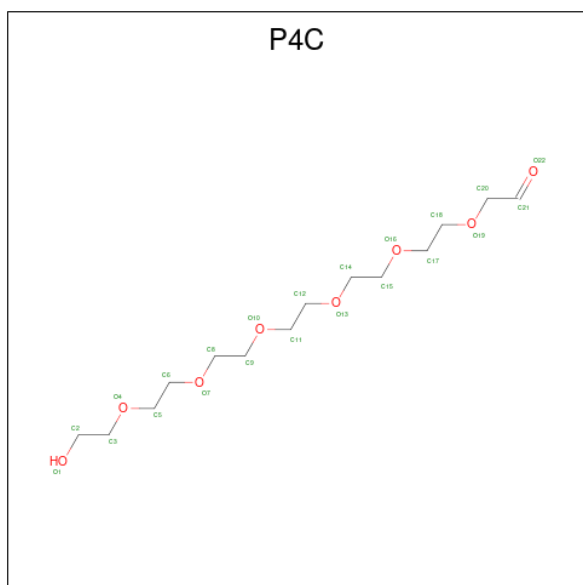
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			12	6	6		
6	C	1	Total	C	O	0	0
			12	6	6		
6	E	1	Total	C	O	0	0
			12	6	6		
6	F	1	Total	C	O	0	0
			12	6	6		

- Molecule 7 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	B	1	Total	C	O	0	0
			11	6	5		
7	E	1	Total	C	O	0	0
			11	6	5		

- Molecule 8 is O-ACETALDEHYDYL-HEXAETHYLENE GLYCOL (CCD ID: P4C) (formula: C₁₄H₂₈O₈).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	C	O	0	0
			22	14	8		
8	C	1	Total	C	O	0	0
			22	14	8		
8	E	1	Total	C	O	0	0
			22	14	8		
8	E	1	Total	C	O	0	0
			22	14	8		
8	F	1	Total	C	O	0	0
			22	14	8		

- Molecule 9 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
9	C	1	Total	C	N	O	0	0
			15	8	1	6		
9	F	1	Total	C	N	O	0	0
			15	8	1	6		

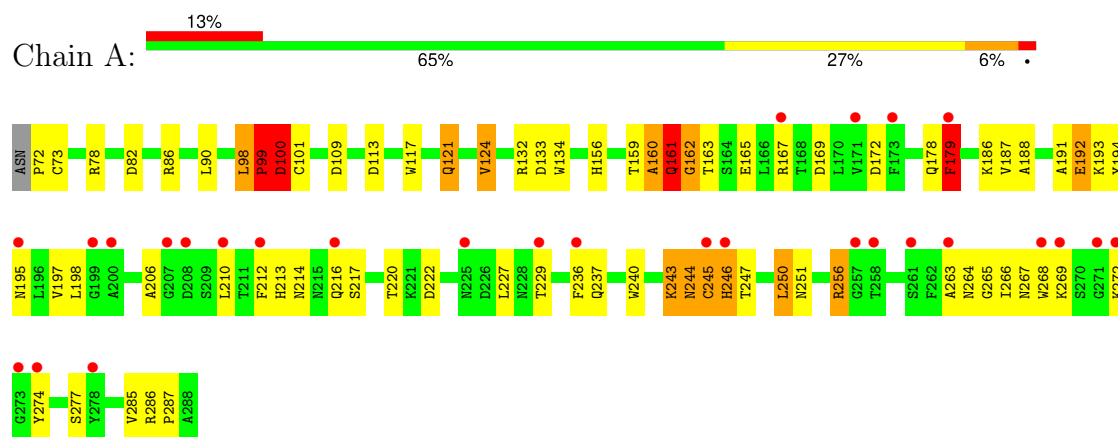
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	29	Total	O	0	0
			29	29		
10	B	89	Total	O	0	0
			89	89		
10	C	59	Total	O	0	0
			59	59		
10	D	32	Total	O	0	0
			32	32		
10	E	64	Total	O	0	0
			64	64		
10	F	62	Total	O	0	0
			62	62		

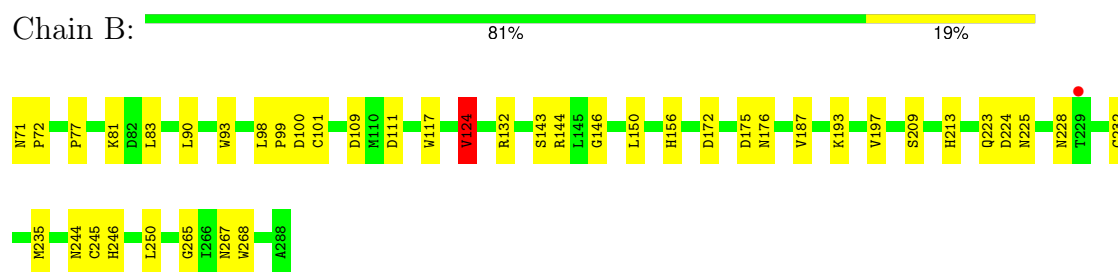
3 Residue-property plots

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.

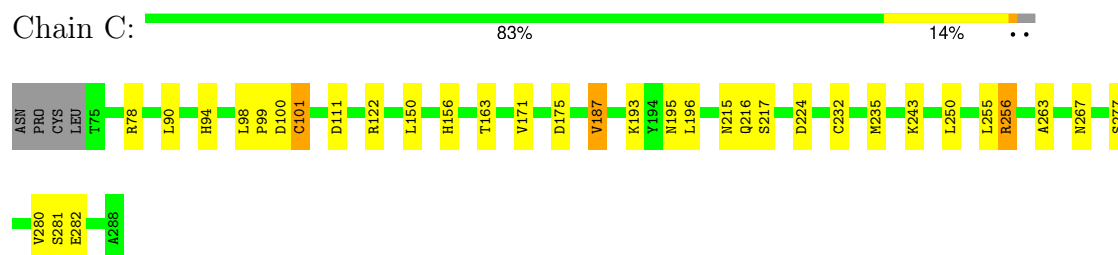
• Molecule 1: FICOLIN-2



• Molecule 1: FICOLIN-2

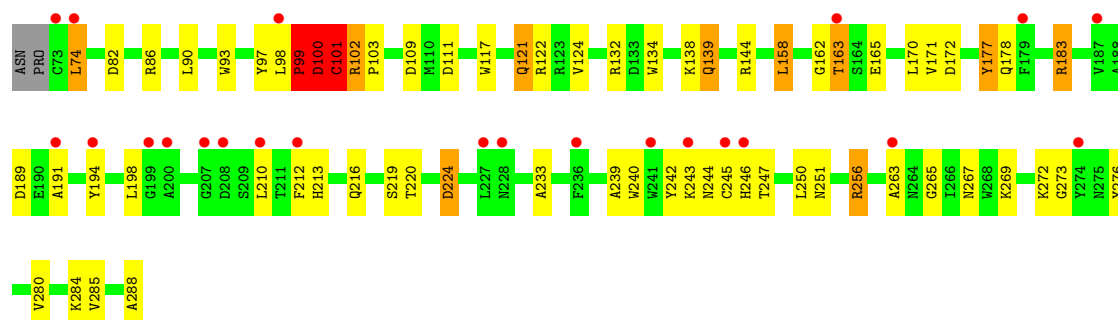


• Molecule 1: FICOLIN-2

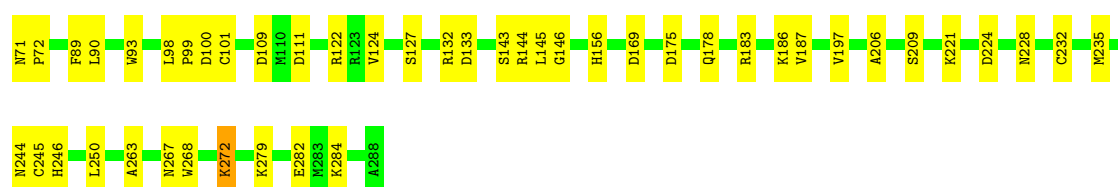
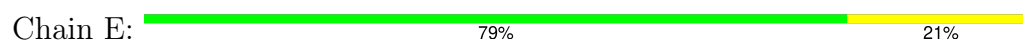


• Molecule 1: FICOLIN-2

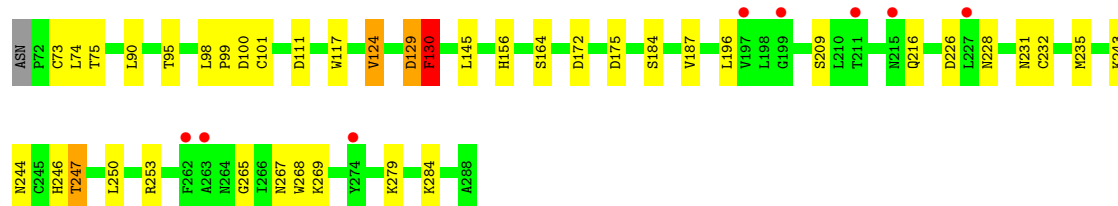
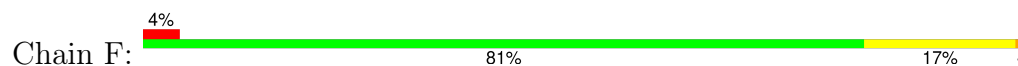




• Molecule 1: FICOLIN-2



• Molecule 1: FICOLIN-2



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



• Molecule 2: 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 32	Depositor
Cell constants a, b, c, α , β , γ	96.08Å 96.08Å 140.98Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.95 – 2.15 19.95 – 2.15	Depositor EDS
% Data completeness (in resolution range)	100.0 (19.95-2.15) 99.1 (19.95-2.15)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.51 (at 2.15Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.188 , 0.237 0.191 , 0.239	Depositor DCC
R_{free} test set	3923 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	30.9	Xtriage
Anisotropy	0.286	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 36.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.029 for -h,-k,l 0.470 for h,-h-k,-l 0.029 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.89	EDS
Total number of atoms	11129	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.70% 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: GLA, P4C, EPE, ACT, MAN, CA, NAG

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	1.21	9/1816 (0.5%)	1.26	16/2455 (0.7%)
1	B	0.53	1/1803 (0.1%)	0.77	1/2439 (0.0%)
1	C	0.49	0/1771	0.77	1/2395 (0.0%)
1	D	1.88	2/1785 (0.1%)	0.90	3/2414 (0.1%)
1	E	0.51	0/1800	0.75	0/2436
1	F	0.50	0/1784	0.82	2/2413 (0.1%)
All	All	1.00	12/10759 (0.1%)	0.90	23/14552 (0.2%)

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	A	2	5
1	D	2	4
1	F	0	1
All	All	4	10

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	74	LEU	C-O	73.63	2.33	1.24
1	A	100	ASP	CA-CB	-25.49	1.10	1.53
1	A	256	ARG	C-O	24.39	1.61	1.23
1	D	256	ARG	C-O	23.68	1.53	1.24
1	A	99	PRO	N-CA	-18.10	1.24	1.47
1	A	99	PRO	CA-CB	-15.31	1.32	1.53
1	A	162	GLY	N-CA	13.05	1.66	1.45
1	A	100	ASP	CA-C	11.50	1.68	1.52
1	A	161	GLN	C-N	7.53	1.43	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	161	GLN	C-O	-7.02	1.15	1.24
1	A	100	ASP	C-O	6.54	1.32	1.24
1	B	81	LYS	CG-CD	6.07	1.70	1.52

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	99	PRO	N-CA-CB	25.34	129.86	103.25
1	A	256	ARG	CA-C-O	-19.44	96.77	122.44
1	D	256	ARG	CA-C-O	-18.88	93.51	120.51
1	D	74	LEU	CA-C-O	-17.33	99.68	119.78
1	A	100	ASP	N-CA-CB	15.59	136.84	110.49
1	A	256	ARG	O-C-N	-14.77	105.31	123.33
1	A	98	LEU	CA-C-N	13.15	136.27	119.84
1	A	98	LEU	C-N-CA	13.15	136.27	119.84
1	A	100	ASP	CB-CA-C	12.14	134.58	110.42
1	A	99	PRO	CA-N-CD	-11.96	95.26	112.00
1	A	100	ASP	N-CA-C	-10.52	88.40	110.80
1	A	161	GLN	CA-C-O	-9.90	106.35	120.51
1	A	99	PRO	CA-CB-CG	-8.26	88.80	104.50
1	F	130	PHE	N-CA-C	-7.43	94.98	110.80
1	A	179	PHE	N-CA-C	7.03	119.07	108.46
1	A	100	ASP	CA-CB-CG	6.99	119.58	112.60
1	A	161	GLN	O-C-N	-6.72	113.66	122.59
1	D	177	TYR	N-CA-C	6.59	121.45	113.41
1	B	81	LYS	CB-CG-CD	-5.97	97.58	111.30
1	A	100	ASP	CA-C-O	5.71	128.68	120.51
1	A	246	HIS	N-CA-C	5.42	117.42	109.24
1	C	78	ARG	CB-CG-CD	-5.31	99.10	111.30
1	F	246	HIS	N-CA-C	5.16	116.91	109.07

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	99	PRO	CA
1	A	100	ASP	CA
1	D	100	ASP	CA
1	D	101	CYS	CA

All (10) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	100	ASP	Peptide
1	A	161	GLN	Mainchain,Peptide
1	A	162	GLY	Peptide
1	A	256	ARG	Mainchain
1	D	100	ASP	Peptide
1	D	256	ARG	Mainchain
1	D	74	LEU	Mainchain
1	D	99	PRO	Peptide
1	F	129	ASP	Peptide

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	1767	0	1636	79	0
1	B	1752	0	1624	40	0
1	C	1721	0	1591	25	0
1	D	1738	0	1608	69	0
1	E	1752	0	1616	47	0
1	F	1736	0	1609	43	0
2	G	28	0	25	0	0
2	H	28	0	25	2	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
4	B	8	0	6	10	0
4	C	4	0	3	0	0
4	E	4	0	3	14	0
4	F	4	0	3	0	0
5	B	12	0	12	1	0
5	C	12	0	12	0	0
5	F	12	0	12	1	0
6	B	12	0	12	0	0
6	C	12	0	12	1	0
6	E	12	0	12	6	0
6	F	12	0	12	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
7	B	11	0	10	0	0
7	E	11	0	10	2	0
8	B	22	0	27	3	0
8	C	22	0	27	1	0
8	E	44	0	54	3	0
8	F	22	0	27	4	0
9	C	15	0	15	1	0
9	F	15	0	15	0	0
10	A	29	0	0	10	0
10	B	89	0	0	3	0
10	C	59	0	0	3	0
10	D	32	0	0	13	0
10	E	64	0	0	0	0
10	F	62	0	0	2	0
All	All	11129	0	10018	306	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 (306) 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:D:101:CYS:CB	1:D:102:ARG:HA	1.24	1.44
1:D:101:CYS:HB3	1:D:102:ARG:CA	1.52	1.36
1:D:101:CYS:CB	1:D:102:ARG:CA	2.15	1.16
1:D:101:CYS:HB2	1:D:102:ARG:HA	1.21	1.13
1:C:255:LEU:O	1:C:277:SER:HB3	1.49	1.13
1:F:73:CYS:CB	1:F:74:LEU:HB2	1.82	1.10
1:F:74:LEU:N	1:F:75:THR:HA	1.58	1.10
1:D:269:LYS:HB3	10:D:2030:HOH:O	1.51	1.09
1:C:243:LYS:HE2	10:C:2036:HOH:O	1.52	1.08
1:F:73:CYS:HB2	1:F:74:LEU:HB2	1.06	1.05
1:A:98:LEU:O	1:A:101:CYS:HA	1.57	1.04
1:A:243:LYS:C	1:A:245:CYS:H	1.63	1.02
1:D:102:ARG:HB3	1:D:103:PRO:HD2	1.38	1.02
1:A:265:GLY:HA3	10:A:2026:HOH:O	1.62	1.00
1:A:269:LYS:HE3	10:A:2020:HOH:O	1.61	1.00
1:F:216:GLN:HG2	1:F:243:LYS:HE3	1.45	0.97
1:D:101:CYS:HB3	1:D:102:ARG:C	1.93	0.93
1:D:97:TYR:HD1	1:D:101:CYS:HG	1.15	0.92
1:D:101:CYS:HB3	1:D:102:ARG:HA	0.92	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:ASP:OD1	1:A:179:PHE:CD2	2.25	0.89
1:A:198:LEU:H	1:A:214:ASN:HD21	1.19	0.88
1:B:245:CYS:H	4:B:1290:ACT:CH3	1.87	0.87
1:D:99:PRO:HA	1:D:100:ASP:OD1	1.75	0.87
1:A:198:LEU:H	1:A:214:ASN:ND2	1.73	0.86
1:B:223:GLN:HE21	1:B:225:ASN:HD21	1.21	0.86
1:A:191:ALA:O	1:A:192:GLU:HB2	1.75	0.84
1:F:73:CYS:HB2	1:F:74:LEU:CB	2.01	0.84
1:A:179:PHE:HA	10:A:2016:HOH:O	1.77	0.84
1:C:215:ASN:ND2	9:C:1293:NAG:O5	2.10	0.83
1:A:243:LYS:C	1:A:245:CYS:N	2.36	0.82
1:C:100:ASP:O	1:C:101:CYS:HB2	1.76	0.82
1:D:138:LYS:HB2	10:D:2022:HOH:O	1.81	0.81
1:A:244:ASN:O	1:A:245:CYS:HB2	1.80	0.81
1:A:192:GLU:O	1:A:193:LYS:HG2	1.81	0.80
1:E:71:ASN:N	1:E:72:PRO:HD2	1.96	0.79
1:A:265:GLY:CA	10:A:2026:HOH:O	2.22	0.77
1:E:246:HIS:H	4:E:1290:ACT:CH3	1.97	0.77
1:E:71:ASN:N	1:E:72:PRO:CD	2.49	0.76
1:B:245:CYS:H	4:B:1290:ACT:H3	1.50	0.76
1:A:169:ASP:HA	1:A:179:PHE:HB3	1.66	0.76
1:A:188:ALA:CB	10:A:2017:HOH:O	2.33	0.76
1:C:156:HIS:HE1	10:C:2029:HOH:O	1.69	0.75
1:B:71:ASN:N	1:B:72:PRO:CD	2.49	0.74
1:F:73:CYS:HG	1:F:101:CYS:HG	1.33	0.74
1:A:277:SER:HB2	10:A:2026:HOH:O	1.86	0.74
1:D:244:ASN:N	1:D:245:CYS:HA	2.02	0.74
1:A:160:ALA:O	1:A:186:LYS:HD3	1.88	0.73
1:F:100:ASP:O	1:F:101:CYS:HB2	1.88	0.71
1:A:169:ASP:OD1	1:A:179:PHE:HD2	1.72	0.70
1:B:246:HIS:H	4:B:1290:ACT:CH3	2.05	0.69
1:D:224:ASP:HB3	10:D:2028:HOH:O	1.90	0.69
1:E:127:SER:HA	8:F:1294:P4C:H121	1.74	0.69
1:D:102:ARG:CB	1:D:103:PRO:HD2	2.20	0.69
1:F:73:CYS:C	1:F:75:THR:HA	2.18	0.69
1:A:266:ILE:N	10:A:2026:HOH:O	2.26	0.68
1:C:232:CYS:HA	1:C:235:MET:HE2	1.74	0.68
1:E:246:HIS:H	4:E:1290:ACT:H1	1.58	0.68
1:A:179:PHE:N	1:A:179:PHE:HD1	1.93	0.67
1:A:237:GLN:NE2	1:A:264:ASN:HD22	1.93	0.67
1:F:216:GLN:CG	1:F:243:LYS:HG2	2.24	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:98:LEU:HB3	1:C:99:PRO:HD2	1.77	0.66
1:F:232:CYS:HA	1:F:235:MET:HE3	1.75	0.66
1:B:197:VAL:HG22	10:B:2040:HOH:O	1.97	0.65
1:D:233:ALA:HB2	10:D:2026:HOH:O	1.96	0.65
1:E:127:SER:HA	8:F:1294:P4C:C12	2.27	0.65
1:E:156:HIS:HD2	1:E:187:VAL:O	1.79	0.65
1:A:172:ASP:HB3	1:A:272:LYS:HE2	1.77	0.65
1:E:221:LYS:HZ1	6:E:1289:GLA:H62	1.62	0.64
1:B:156:HIS:HD2	1:B:187:VAL:O	1.80	0.64
1:E:122:ARG:HG3	1:E:282:GLU:HG2	1.80	0.64
1:A:210:LEU:HG	1:A:268:TRP:CH2	2.33	0.64
1:D:102:ARG:HB3	1:D:103:PRO:CD	2.22	0.64
1:B:124:VAL:HG22	8:B:1296:P4C:H172	1.80	0.63
1:A:216:GLN:HB3	1:A:243:LYS:HD3	1.79	0.63
1:E:228:ASN:ND2	1:E:244[A]:ASN:OD1	2.26	0.63
1:D:243:LYS:HG2	1:D:244:ASN:N	2.14	0.63
1:D:212:PHE:HD1	10:D:2025:HOH:O	1.82	0.62
1:B:246:HIS:H	4:B:1290:ACT:H1	1.63	0.62
1:C:122:ARG:HD2	1:C:282:GLU:OE2	1.99	0.62
1:F:73:CYS:HG	1:F:101:CYS:CB	2.12	0.62
1:C:255:LEU:O	1:C:277:SER:CB	2.37	0.61
1:A:179:PHE:N	1:A:179:PHE:CD1	2.67	0.61
1:B:245:CYS:HB2	4:B:1290:ACT:CH3	2.31	0.61
1:D:134:TRP:HA	1:D:220:THR:HG21	1.82	0.61
1:F:216:GLN:HG2	1:F:243:LYS:CE	2.25	0.61
1:E:246:HIS:H	4:E:1290:ACT:H3	1.65	0.61
1:A:188:ALA:HB3	10:A:2017:HOH:O	1.94	0.60
1:A:167:ARG:NH2	1:A:179:PHE:CD2	2.69	0.60
1:E:183:ARG:HD3	1:E:183:ARG:O	2.01	0.60
1:D:183:ARG:HG3	1:D:183:ARG:HH11	1.67	0.59
1:F:98:LEU:HB3	1:F:99:PRO:HD2	1.83	0.59
1:F:156:HIS:HE1	10:F:2036:HOH:O	1.84	0.59
1:D:246:HIS:ND1	10:D:2029:HOH:O	2.32	0.59
1:D:82:ASP:O	1:D:86:ARG:HB2	2.02	0.59
1:F:231:ASN:C	1:F:235:MET:HE2	2.27	0.59
1:B:143:SER:HB2	6:C:1290:GLA:O5	2.03	0.59
1:F:253:ARG:NH2	10:F:2056:HOH:O	2.35	0.59
1:D:194:TYR:CE2	10:D:2022:HOH:O	2.52	0.58
1:B:245:CYS:HB2	4:B:1290:ACT:H1	1.84	0.58
1:D:219:SER:HB2	10:D:2026:HOH:O	2.03	0.58
1:E:132:ARG:HB3	6:E:1289:GLA:HO1	1.69	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:188:ALA:HB1	1:A:192:GLU:OE1	2.04	0.58
1:A:191:ALA:O	1:A:192:GLU:CB	2.47	0.58
1:A:132:ARG:NH2	1:B:111:ASP:OD1	2.37	0.57
1:F:74:LEU:N	1:F:75:THR:CA	2.48	0.57
1:F:100:ASP:O	1:F:101:CYS:CB	2.53	0.56
1:F:156:HIS:HD2	1:F:187:VAL:O	1.88	0.56
1:A:101:CYS:O	1:A:101:CYS:SG	2.63	0.56
1:D:194:TYR:HE2	10:D:2022:HOH:O	1.86	0.56
1:E:221:LYS:HZ3	6:E:1289:GLA:H5	1.71	0.56
1:E:245:CYS:HB2	4:E:1290:ACT:H1	1.86	0.56
1:D:272:LYS:HD3	1:D:276:TYR:CD2	2.40	0.56
1:B:223:GLN:NE2	1:B:225:ASN:HD21	1.97	0.56
1:B:245:CYS:H	4:B:1290:ACT:H2	1.67	0.56
1:D:172:ASP:CB	1:D:272:LYS:HE3	2.37	0.55
1:D:121[A]:GLN:OE1	1:D:251:ASN:OD1	2.24	0.55
1:D:97:TYR:HD1	1:D:101:CYS:SG	2.26	0.55
1:E:143:SER:HB2	6:F:1290:GLA:O5	2.05	0.55
1:B:71:ASN:N	1:B:72:PRO:HD3	2.20	0.55
1:B:209:SER:HB2	1:B:268:TRP:CE2	2.42	0.54
1:E:245:CYS:N	4:E:1290:ACT:H2	2.16	0.54
1:F:74:LEU:H	1:F:75:THR:HA	1.61	0.54
1:B:132:ARG:HH22	1:C:111:ASP:CG	2.15	0.54
1:C:156:HIS:HD2	1:C:187:VAL:O	1.89	0.54
1:F:279:LYS:HE3	8:F:1294:P4C:H52	1.90	0.54
1:D:98:LEU:O	1:D:100:ASP:N	2.36	0.53
1:A:167:ARG:NH2	1:A:179:PHE:CE2	2.57	0.53
1:A:210:LEU:HG	1:A:268:TRP:HH2	1.71	0.53
1:E:272:LYS:HE3	8:E:1296:P4C:H92	1.91	0.53
1:A:73:CYS:O	1:A:78:ARG:HD3	2.08	0.53
1:B:245:CYS:N	4:B:1290:ACT:CH3	2.67	0.53
1:D:170:LEU:O	1:D:177:TYR:O	2.25	0.53
1:D:219:SER:CB	10:D:2026:HOH:O	2.57	0.52
1:B:144:ARG:NH2	1:C:94:HIS:NE2	2.57	0.52
1:B:246:HIS:H	4:B:1290:ACT:H3	1.75	0.52
1:C:193:LYS:CB	1:C:217:SER:HB3	2.39	0.52
1:A:247:THR:O	1:A:269:LYS:HG2	2.10	0.52
1:D:121[A]:GLN:HG3	1:D:122:ARG:N	2.17	0.52
1:D:172:ASP:HB3	1:D:272:LYS:HE3	1.92	0.52
1:A:172:ASP:CB	1:A:272:LYS:HE2	2.39	0.52
1:F:209:SER:HB3	1:F:247:THR:HG22	1.92	0.52
1:E:132:ARG:HD2	6:E:1289:GLA:H3	1.91	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:132:ARG:NH2	1:F:111:ASP:OD2	2.41	0.51
1:A:72:PRO:HA	1:A:99:PRO:HB3	1.91	0.51
1:A:192:GLU:C	1:A:193:LYS:HG2	2.35	0.51
1:E:186:LYS:HB3	1:E:197:VAL:HB	1.92	0.50
1:E:169:ASP:OD2	1:E:284:LYS:NZ	2.42	0.50
1:B:244:ASN:HB3	10:B:2063:HOH:O	2.10	0.50
1:E:209:SER:HB2	1:E:268:TRP:CE2	2.46	0.50
1:E:244[A]:ASN:N	1:E:245:CYS:HA	2.27	0.50
1:D:224:ASP:OD1	1:D:224:ASP:C	2.54	0.50
1:A:216:GLN:HG2	1:A:243:LYS:NZ	2.27	0.50
1:F:73:CYS:HG	1:F:101:CYS:HB3	1.77	0.49
1:D:246:HIS:HE1	1:D:263:ALA:O	1.95	0.49
1:E:245:CYS:N	4:E:1290:ACT:CH3	2.62	0.49
1:A:117:TRP:CZ3	1:A:286:ARG:HG2	2.48	0.49
1:A:243:LYS:O	1:A:245:CYS:N	2.40	0.49
1:E:244[B]:ASN:N	1:E:245:CYS:HA	2.28	0.49
1:A:160:ALA:O	1:A:161:GLN:HG2	2.12	0.49
1:C:98:LEU:HB3	1:C:99:PRO:CD	2.43	0.49
1:F:265:GLY:H	1:F:267:ASN:HD21	1.60	0.49
1:F:117:TRP:HB3	1:F:284:LYS:HB2	1.95	0.48
1:E:245:CYS:HB2	4:E:1290:ACT:CH3	2.42	0.48
1:F:232:CYS:HA	1:F:235:MET:CE	2.43	0.48
1:A:82:ASP:O	1:A:86:ARG:HG3	2.13	0.48
1:D:265:GLY:H	1:D:267:ASN:HD21	1.62	0.48
1:E:132:ARG:HH22	1:F:111:ASP:CG	2.21	0.48
1:C:193:LYS:HB2	1:C:217:SER:HB3	1.94	0.48
1:E:178:GLN:HB3	1:E:206:ALA:HB2	1.95	0.48
1:F:175:ASP:OD2	8:F:1294:P4C:H151	2.13	0.48
1:F:98:LEU:HB3	1:F:99:PRO:CD	2.43	0.48
1:A:188:ALA:O	1:A:194:TYR:HA	2.13	0.47
1:C:263:ALA:HA	1:C:267:ASN:ND2	2.28	0.47
1:E:100:ASP:O	1:E:101:CYS:HB2	2.14	0.47
1:F:209:SER:HB2	1:F:268:TRP:CE2	2.49	0.47
1:A:212[B]:PHE:N	1:A:212[B]:PHE:CD2	2.82	0.47
1:A:198:LEU:N	1:A:214:ASN:ND2	2.53	0.47
1:A:227:LEU:HG	1:A:243:LYS:HG3	1.96	0.47
1:A:187:VAL:HA	1:A:195:ASN:O	2.15	0.47
1:A:113:ASP:OD2	1:A:167:ARG:NH1	2.47	0.47
1:A:121[A]:GLN:OE1	1:A:251:ASN:OD1	2.33	0.47
1:B:100:ASP:O	1:B:101:CYS:HB2	2.15	0.47
1:D:244:ASN:N	1:D:245:CYS:CA	2.76	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:99:PRO:CA	1:D:100:ASP:OD1	2.56	0.47
1:E:175:ASP:OD1	8:E:1295:P4C:H112	2.14	0.47
1:A:159:THR:O	1:A:160:ALA:C	2.58	0.46
1:A:240:TRP:CH2	1:A:250:LEU:HB2	2.50	0.46
1:A:263:ALA:HB2	1:A:274:TYR:HB3	1.96	0.46
1:A:156:HIS:HD2	1:A:187:VAL:O	1.98	0.46
1:D:98:LEU:HB3	1:D:99:PRO:HD2	1.96	0.46
1:D:243:LYS:CG	1:D:244:ASN:N	2.77	0.46
7:E:1294:MAN:H2	2:H:2:NAG:O4	2.15	0.46
1:A:178:GLN:HB3	1:A:206:ALA:HB2	1.97	0.46
1:A:195:ASN:HB3	10:A:2017:HOH:O	2.15	0.46
1:B:193:LYS:HD3	1:B:225:ASN:CG	2.41	0.46
1:A:165:GLU:O	1:A:285:VAL:HA	2.15	0.46
1:A:263:ALA:HA	1:A:267:ASN:ND2	2.31	0.46
1:D:242:TYR:HD2	10:D:2026:HOH:O	1.98	0.46
1:F:172:ASP:OD1	1:F:172:ASP:C	2.58	0.46
1:B:144:ARG:NH1	10:B:2028:HOH:O	2.48	0.46
1:D:101:CYS:HB2	1:D:102:ARG:CA	2.14	0.46
1:D:102:ARG:CB	1:D:103:PRO:CD	2.88	0.46
1:A:212[B]:PHE:N	1:A:212[B]:PHE:HD2	2.14	0.46
1:A:286:ARG:O	1:A:287:PRO:C	2.58	0.46
1:D:243:LYS:C	1:D:245:CYS:HA	2.40	0.46
1:A:214:ASN:O	1:A:216:GLN:HG3	2.16	0.45
1:F:216:GLN:HG3	1:F:243:LYS:HG2	1.96	0.45
1:E:246:HIS:CD2	1:E:246:HIS:C	2.93	0.45
1:A:244:ASN:O	1:A:245:CYS:CB	2.60	0.45
1:C:171:VAL:HB	1:C:280:VAL:HB	1.98	0.45
1:A:134:TRP:HA	1:A:220:THR:HG21	1.99	0.45
1:F:73:CYS:H	1:F:74:LEU:C	2.25	0.45
1:E:221:LYS:NZ	6:E:1289:GLA:H62	2.30	0.45
1:B:232:CYS:HA	1:B:235:MET:HE3	1.99	0.45
1:F:164:SER:O	1:F:184:SER:HA	2.17	0.45
1:A:213:HIS:CE1	1:A:246:HIS:HA	2.52	0.44
1:B:265:GLY:H	1:B:267:ASN:HD21	1.65	0.44
1:D:139:GLN:HG2	1:E:89:PHE:CE2	2.52	0.44
1:E:133:ASP:H	6:E:1289:GLA:HO1	1.65	0.44
1:A:188:ALA:N	10:A:2017:HOH:O	2.21	0.44
1:B:109:ASP:HB3	1:B:117:TRP:HB2	1.99	0.44
1:D:100:ASP:OD1	1:D:100:ASP:N	2.50	0.44
1:E:109:ASP:OD1	1:E:109:ASP:C	2.60	0.44
1:A:193:LYS:HG3	1:A:217:SER:CB	2.47	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:111:ASP:CG	10:D:2007:HOH:O	2.60	0.44
1:B:71:ASN:N	1:B:72:PRO:HD2	2.30	0.44
8:B:1296:P4C:H21A	8:B:1296:P4C:H52	1.37	0.44
1:D:172:ASP:OD1	1:D:172:ASP:C	2.60	0.44
1:D:246:HIS:CD2	1:D:246:HIS:C	2.95	0.44
1:D:98:LEU:CD2	1:D:158:LEU:HD13	2.47	0.44
1:D:240:TRP:CH2	1:D:250:LEU:HB2	2.52	0.44
1:D:247:THR:HB	10:D:2030:HOH:O	2.17	0.44
7:E:1294:MAN:C2	2:H:2:NAG:O4	2.65	0.44
1:C:175:ASP:HB3	8:C:1294:P4C:H171	1.99	0.44
1:F:216:GLN:HE21	1:F:243:LYS:HE2	1.83	0.44
1:D:189:ASP:OD1	1:D:191:ALA:HB3	2.18	0.44
1:F:129:ASP:OD1	1:F:130:PHE:N	2.51	0.44
1:D:109:ASP:OD1	1:D:109:ASP:C	2.61	0.43
1:D:171:VAL:HB	1:D:280:VAL:HB	1.99	0.43
1:F:216:GLN:CG	1:F:243:LYS:CG	2.96	0.43
1:A:109:ASP:OD1	1:A:109:ASP:C	2.61	0.43
1:F:216:GLN:HG2	1:F:243:LYS:HG2	2.01	0.43
1:B:93:TRP:CZ2	1:B:144:ARG:HA	2.53	0.43
1:D:165:GLU:O	1:D:285:VAL:HA	2.18	0.43
1:D:162:GLY:O	1:D:163:THR:HG22	2.19	0.43
1:A:247:THR:HB	1:A:269:LYS:HE2	1.99	0.43
1:D:269:LYS:HA	1:D:273:GLY:HA2	2.01	0.43
5:B:1289:EPE:H102	5:B:1289:EPE:H62	1.56	0.43
1:A:98:LEU:O	1:A:101:CYS:CA	2.48	0.42
1:D:213:HIS:CE1	1:D:246:HIS:HA	2.54	0.42
1:F:226:ASP:C	1:F:226:ASP:OD1	2.61	0.42
5:F:1289:EPE:H101	5:F:1289:EPE:H21	1.68	0.42
1:B:132:ARG:NH2	1:C:111:ASP:OD2	2.48	0.42
1:D:132:ARG:NH2	1:E:111:ASP:OD1	2.52	0.42
1:C:224:ASP:OD1	1:C:224:ASP:C	2.62	0.42
1:A:227:LEU:CD1	1:A:243:LYS:HG3	2.49	0.42
1:E:279:LYS:HZ3	8:E:1295:P4C:H142	1.84	0.42
1:B:77:PRO:HG3	1:B:83:LEU:HD21	2.00	0.42
1:C:195:ASN:HD22	1:C:216:GLN:C	2.28	0.42
1:E:246:HIS:N	4:E:1290:ACT:H1	2.29	0.42
1:A:178:GLN:C	1:A:179:PHE:CD1	2.98	0.42
1:B:175:ASP:OD1	8:B:1296:P4C:H112	2.20	0.42
1:C:187:VAL:HA	1:C:195:ASN:O	2.20	0.42
1:A:73:CYS:HB3	1:A:78:ARG:HG2	2.01	0.42
1:D:243:LYS:N	1:D:245:CYS:O	2.26	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:243:LYS:HA	1:A:243:LYS:HD2	1.82	0.42
1:C:216:GLN:HB2	1:C:243:LYS:HE3	2.02	0.42
1:D:117:TRP:HB3	1:D:284:LYS:HB2	2.02	0.42
1:E:122:ARG:HD2	1:E:282:GLU:OE2	2.20	0.42
1:E:263:ALA:HA	1:E:267:ASN:ND2	2.34	0.42
1:A:133:ASP:HB2	1:A:222:ASP:OD2	2.19	0.41
1:A:167:ARG:HE	1:A:179:PHE:HD2	1.68	0.41
1:D:93:TRP:CZ2	1:D:144:ARG:HA	2.55	0.41
1:F:145:LEU:N	1:F:145:LEU:HD12	2.34	0.41
1:E:93:TRP:CZ2	1:E:144:ARG:HA	2.56	0.41
1:E:156:HIS:CD2	1:E:187:VAL:O	2.67	0.41
1:B:172:ASP:OD2	1:B:176:ASN:HB2	2.21	0.41
1:C:256:ARG:H	1:C:256:ARG:HG3	1.52	0.41
1:E:232:CYS:HA	1:E:235:MET:HE3	2.03	0.41
1:B:224:ASP:OD1	1:B:224:ASP:C	2.63	0.41
1:A:109:ASP:HB3	1:A:117:TRP:HB2	2.01	0.41
1:A:236:PHE:HE2	1:A:245:CYS:SG	2.43	0.41
1:B:245:CYS:HB2	4:B:1290:ACT:H2	2.03	0.41
1:C:122:ARG:NH2	10:C:2015:HOH:O	2.38	0.41
1:B:98:LEU:HB3	1:B:99:PRO:HD2	2.03	0.41
1:D:239:ALA:O	1:D:251:ASN:HB2	2.19	0.41
1:A:172:ASP:OD1	1:A:172:ASP:C	2.64	0.41
1:B:213:HIS:CE1	1:B:246:HIS:HA	2.56	0.41
1:D:198:LEU:HD22	1:D:210:LEU:HG	2.03	0.41
1:F:228:ASN:ND2	1:F:244:ASN:OD1	2.53	0.41
1:D:165:GLU:HG3	1:D:288:ALA:CB	2.51	0.40
1:E:272:LYS:HE2	1:E:272:LYS:HB3	1.83	0.40
1:B:228:ASN:ND2	1:B:244:ASN:OD1	2.50	0.40
1:B:246:HIS:CD2	1:B:246:HIS:C	2.98	0.40
1:D:216:GLN:HE22	1:D:243:LYS:HG3	1.86	0.40
1:F:247:THR:HG23	1:F:269:LYS:HB3	2.02	0.40
1:D:109:ASP:HB3	1:D:117:TRP:HB2	2.03	0.40
1:A:156:HIS:CD2	1:A:187:VAL:O	2.74	0.40
1:E:98:LEU:HB3	1:E:99:PRO:HD2	2.03	0.40
1:E:224:ASP:C	1:E:224:ASP:OD1	2.64	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	218/218 (100%)	188 (86%)	22 (10%)	8 (4%)	2	0
1	B	217/218 (100%)	207 (95%)	8 (4%)	2 (1%)	14	9
1	C	213/218 (98%)	201 (94%)	11 (5%)	1 (0%)	24	20
1	D	215/218 (99%)	192 (89%)	18 (8%)	5 (2%)	5	1
1	E	217/218 (100%)	204 (94%)	11 (5%)	2 (1%)	14	9
1	F	215/218 (99%)	202 (94%)	12 (6%)	1 (0%)	24	20
All	All	1295/1308 (99%)	1194 (92%)	82 (6%)	19 (2%)	8	4

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	192	GLU
1	A	244	ASN
1	A	245	CYS
1	D	100	ASP
1	D	178	GLN
1	C	256	ARG
1	A	99	PRO
1	A	160	ALA
1	A	100	ASP
1	D	99	PRO
1	D	101	CYS
1	A	163	THR
1	A	124	VAL
1	B	124	VAL
1	E	124	VAL
1	B	146	GLY
1	F	124	VAL
1	D	124	VAL
1	E	146	GLY

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	185/183 (101%)	174 (94%)	11 (6%)	18	13
1	B	184/183 (100%)	180 (98%)	4 (2%)	45	51
1	C	180/183 (98%)	172 (96%)	8 (4%)	25	23
1	D	182/183 (100%)	171 (94%)	11 (6%)	17	13
1	E	184/183 (100%)	180 (98%)	4 (2%)	45	51
1	F	182/183 (100%)	175 (96%)	7 (4%)	29	29
All	All	1097/1098 (100%)	1052 (96%)	45 (4%)	28	26

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	90	LEU
1	A	99	PRO
1	A	121[A]	GLN
1	A	121[B]	GLN
1	A	124	VAL
1	A	161	GLN
1	A	179	PHE
1	A	197	VAL
1	A	229	THR
1	A	243	LYS
1	A	250	LEU
1	B	90	LEU
1	B	124	VAL
1	B	150	LEU
1	B	250	LEU
1	C	90	LEU
1	C	101	CYS
1	C	150	LEU
1	C	163	THR
1	C	187	VAL
1	C	196	LEU
1	C	250	LEU

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Mol	Chain	Res	Type
1	C	281	SER
1	D	90	LEU
1	D	100	ASP
1	D	101	CYS
1	D	102	ARG
1	D	121[A]	GLN
1	D	121[B]	GLN
1	D	139	GLN
1	D	158	LEU
1	D	163	THR
1	D	183	ARG
1	D	224	ASP
1	E	90	LEU
1	E	145	LEU
1	E	250	LEU
1	E	272	LYS
1	F	90	LEU
1	F	95	THR
1	F	124	VAL
1	F	130	PHE
1	F	196	LEU
1	F	247	THR
1	F	250	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	156	HIS
1	A	161	GLN
1	A	214	ASN
1	A	228	ASN
1	A	237	GLN
1	A	267	ASN
1	B	71	ASN
1	B	94	HIS
1	B	156	HIS
1	B	225	ASN
1	B	246	HIS
1	B	267	ASN
1	C	156	HIS
1	C	195	ASN
1	C	223	GLN

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Mol	Chain	Res	Type
1	C	267	ASN
1	D	94	HIS
1	D	156	HIS
1	D	216	GLN
1	D	246	HIS
1	D	267	ASN
1	D	275	ASN
1	E	71	ASN
1	E	94	HIS
1	E	156	HIS
1	E	246	HIS
1	E	267	ASN
1	F	156	HIS
1	F	176	ASN
1	F	195	ASN
1	F	216	GLN
1	F	267	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	$\# Z > 2$
2	NAG	G	1	2,1	14,14,15	0.59	0	17,19,21	1.14	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	G	2	2	14,14,15	0.38	0	17,19,21	0.88	0
2	NAG	H	1	2,1	14,14,15	0.68	0	17,19,21	1.12	2 (11%)
2	NAG	H	2	2	14,14,15	0.52	0	17,19,21	0.81	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	NAG	H	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1	NAG	C3-C4-C5	-2.86	105.05	110.23
2	G	1	NAG	C3-C4-C5	-2.65	105.42	110.23
2	G	1	NAG	O5-C1-C2	-2.63	107.22	111.29
2	H	1	NAG	O5-C1-C2	-2.12	108.01	111.29

There are no chirality outliers.

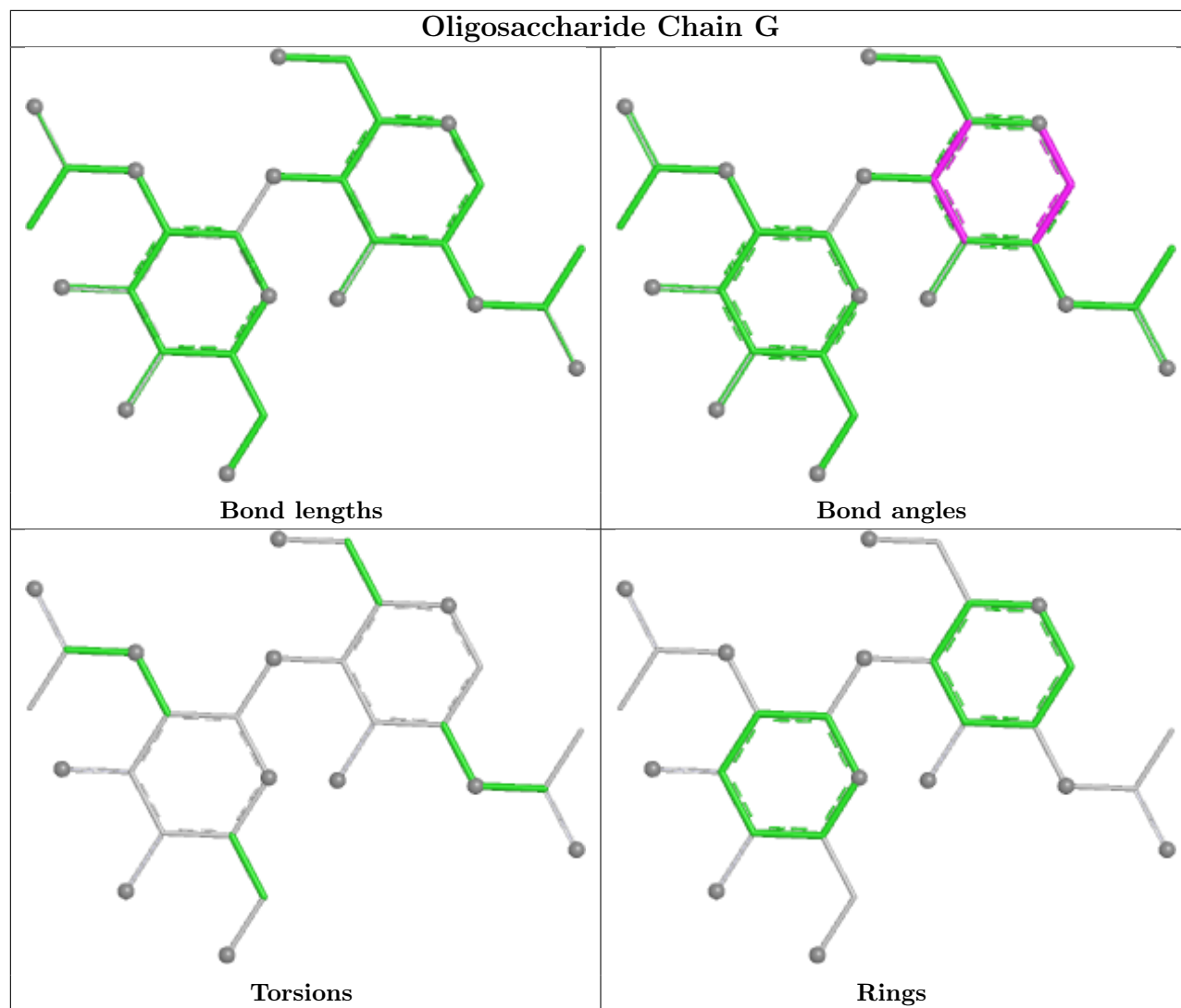
There are no torsion outliers.

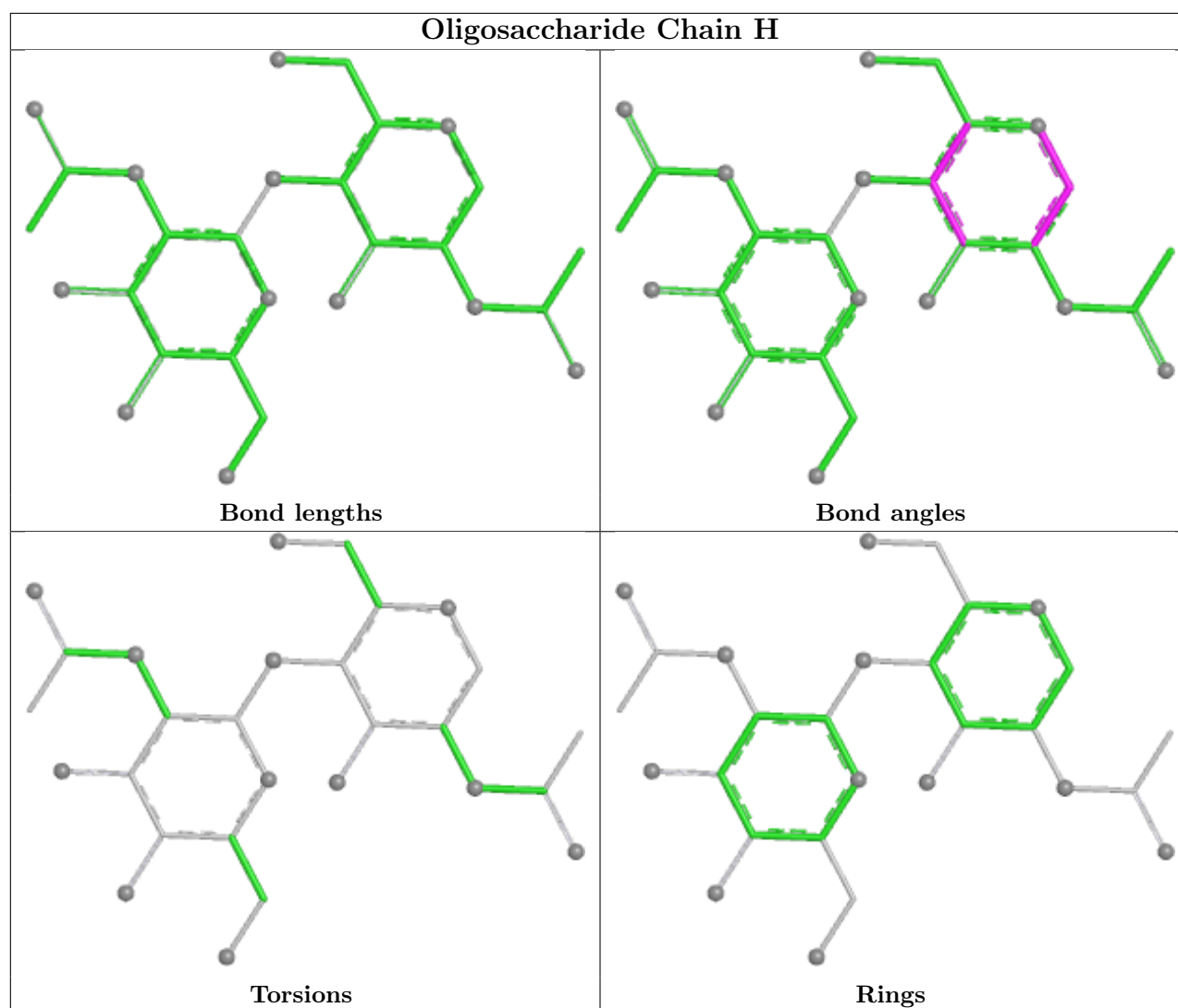
There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	2	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](#)

Of 27 ligands modelled in this entry, 6 are monoatomic - leaving 21 for Mogul analysis.

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	GLA	C	1290	-	12,12,12	0.55	0	17,17,17	0.53	0
4	ACT	C	1291	-	3,3,3	0.83	0	3,3,3	1.18	0
8	P4C	B	1296	-	21,21,21	1.35	2 (9%)	20,20,20	1.32	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	ACT	B	407	-	3,3,3	0.84	0	3,3,3	1.22	0
7	MAN	E	1294	-	11,11,12	0.59	0	15,15,17	1.17	3 (20%)
6	GLA	F	1290	-	12,12,12	0.52	0	17,17,17	0.72	0
4	ACT	B	1290	-	3,3,3	0.85	0	3,3,3	0.93	0
6	GLA	E	1289	-	12,12,12	0.76	0	17,17,17	1.50	3 (17%)
8	P4C	F	1294	-	21,21,21	1.37	2 (9%)	20,20,20	1.29	1 (5%)
8	P4C	E	1296	-	21,21,21	1.38	2 (9%)	20,20,20	1.24	1 (5%)
6	GLA	B	1291	-	12,12,12	0.68	0	17,17,17	0.83	0
8	P4C	C	1294	-	21,21,21	1.37	2 (9%)	20,20,20	1.27	1 (5%)
4	ACT	E	1290	-	3,3,3	0.85	0	3,3,3	0.90	0
5	EPE	B	1289	-	12,12,15	1.10	1 (8%)	15,16,20	1.45	2 (13%)
8	P4C	E	1295	-	21,21,21	1.35	2 (9%)	20,20,20	1.28	1 (5%)
4	ACT	F	1291	-	3,3,3	0.78	0	3,3,3	1.43	0
7	MAN	B	1295	-	11,11,12	0.56	0	15,15,17	0.55	0
9	NAG	C	1293	-	15,15,15	0.46	0	21,21,21	0.72	0
5	EPE	C	1289	-	12,12,15	1.15	1 (8%)	15,16,20	1.22	1 (6%)
9	NAG	F	1293	-	15,15,15	0.52	0	21,21,21	0.98	1 (4%)
5	EPE	F	1289	-	12,12,15	0.99	1 (8%)	15,16,20	1.34	1 (6%)

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
9	NAG	C	1293	-	-	3/6/26/26	0/1/1/1
7	MAN	E	1294	-	-	2/2/19/22	0/1/1/1
6	GLA	C	1290	-	-	1/2/22/22	0/1/1/1
6	GLA	B	1291	-	-	1/2/22/22	0/1/1/1
6	GLA	F	1290	-	-	1/2/22/22	0/1/1/1
8	P4C	C	1294	-	-	8/18/19/19	-
6	GLA	E	1289	-	-	2/2/22/22	0/1/1/1
8	P4C	F	1294	-	-	14/18/19/19	-
5	EPE	C	1289	-	-	1/6/14/19	0/1/1/1
9	NAG	F	1293	-	-	6/6/26/26	0/1/1/1
8	P4C	B	1296	-	-	13/18/19/19	-
5	EPE	B	1289	-	-	5/6/14/19	0/1/1/1
8	P4C	E	1296	-	-	12/18/19/19	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
8	P4C	E	1295	-	-	8/18/19/19	-
7	MAN	B	1295	-	-	2/2/19/22	0/1/1/1
5	EPE	F	1289	-	-	0/6/14/19	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	B	1296	P4C	O22-C21	4.29	1.44	1.20
8	E	1295	P4C	O22-C21	4.27	1.44	1.20
8	E	1296	P4C	O22-C21	4.26	1.44	1.20
8	C	1294	P4C	O22-C21	4.26	1.44	1.20
8	F	1294	P4C	O22-C21	4.25	1.44	1.20
8	F	1294	P4C	O1-C2	-3.85	1.22	1.42
8	E	1296	P4C	O1-C2	-3.82	1.22	1.42
8	C	1294	P4C	O1-C2	-3.74	1.23	1.42
8	E	1295	P4C	O1-C2	-3.72	1.23	1.42
8	B	1296	P4C	O1-C2	-3.70	1.23	1.42
5	C	1289	EPE	C10-S	3.61	1.82	1.77
5	B	1289	EPE	C10-S	3.43	1.82	1.77
5	F	1289	EPE	C10-S	3.14	1.82	1.77

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	B	1296	P4C	O22-C21-C20	-5.57	109.63	126.10
8	F	1294	P4C	O22-C21-C20	-5.51	109.82	126.10
8	C	1294	P4C	O22-C21-C20	-5.40	110.13	126.10
8	E	1295	P4C	O22-C21-C20	-5.29	110.45	126.10
8	E	1296	P4C	O22-C21-C20	-5.23	110.64	126.10
5	B	1289	EPE	C5-N4-C3	3.79	121.12	110.40
6	E	1289	GLA	C3-C4-C5	3.64	116.83	110.23
5	F	1289	EPE	C5-N4-C3	3.41	120.02	110.40
6	E	1289	GLA	C4-C3-C2	3.40	116.80	110.83
5	C	1289	EPE	C5-N4-C3	3.01	118.89	110.40
5	B	1289	EPE	O2S-S-C10	2.95	111.18	106.73
7	E	1294	MAN	C3-C4-C5	2.50	114.77	110.23
7	E	1294	MAN	C2-C3-C4	2.25	114.82	110.86
9	F	1293	NAG	O5-C1-C2	2.23	111.75	109.52
6	E	1289	GLA	O5-C1-C2	-2.06	106.67	110.30
7	E	1294	MAN	C1-C2-C3	2.01	112.57	109.64

There are no chirality outliers.

All (79) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	B	1289	EPE	C10-C9-N1-C6
5	C	1289	EPE	C10-C9-N1-C6
9	C	1293	NAG	C3-C2-N2-C7
9	C	1293	NAG	C8-C7-N2-C2
9	C	1293	NAG	O7-C7-N2-C2
9	F	1293	NAG	C1-C2-N2-C7
9	F	1293	NAG	C8-C7-N2-C2
9	F	1293	NAG	O7-C7-N2-C2
8	E	1296	P4C	C5-C6-O7-C8
8	B	1296	P4C	C2-C3-O4-C5
7	E	1294	MAN	C4-C5-C6-O6
9	F	1293	NAG	O5-C5-C6-O6
9	F	1293	NAG	C4-C5-C6-O6
8	C	1294	P4C	O13-C14-C15-O16
8	E	1296	P4C	O13-C14-C15-O16
7	E	1294	MAN	O5-C5-C6-O6
8	B	1296	P4C	O4-C5-C6-O7
8	E	1296	P4C	O7-C8-C9-O10
8	E	1296	P4C	O16-C17-C18-O19
8	F	1294	P4C	O10-C11-C12-O13
7	B	1295	MAN	C4-C5-C6-O6
6	E	1289	GLA	O5-C5-C6-O6
8	C	1294	P4C	O7-C8-C9-O10
8	C	1294	P4C	O10-C11-C12-O13
8	E	1295	P4C	O13-C14-C15-O16
7	B	1295	MAN	O5-C5-C6-O6
8	B	1296	P4C	O1-C2-C3-O4
8	F	1294	P4C	O1-C2-C3-O4
8	E	1295	P4C	O16-C17-C18-O19
8	E	1295	P4C	C21-C20-O19-C18
8	E	1296	P4C	O4-C5-C6-O7
6	C	1290	GLA	O5-C5-C6-O6
6	F	1290	GLA	O5-C5-C6-O6
5	B	1289	EPE	C9-C10-S-O3S
8	F	1294	P4C	C21-C20-O19-C18
8	C	1294	P4C	C11-C12-O13-C14
6	E	1289	GLA	C4-C5-C6-O6
6	B	1291	GLA	O5-C5-C6-O6
5	B	1289	EPE	C9-C10-S-O1S
5	B	1289	EPE	C9-C10-S-O2S
8	F	1294	P4C	C11-C12-O13-C14
8	B	1296	P4C	C17-C18-O19-C20

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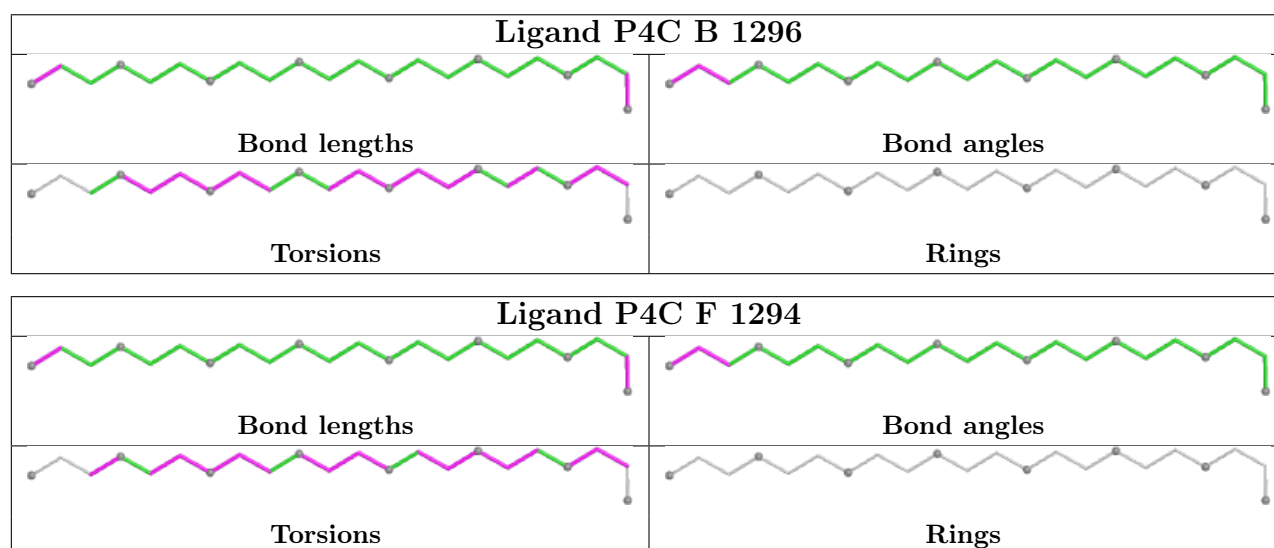
Mol	Chain	Res	Type	Atoms
8	E	1295	P4C	C11-C12-O13-C14
8	E	1296	P4C	C11-C12-O13-C14
8	B	1296	P4C	C18-C17-O16-C15
8	B	1296	P4C	C9-C8-O7-C6
8	B	1296	P4C	C12-C11-O10-C9
8	E	1295	P4C	C2-C3-O4-C5
8	E	1296	P4C	C9-C8-O7-C6
8	F	1294	P4C	C2-C3-O4-C5
8	F	1294	P4C	C12-C11-O10-C9
8	F	1294	P4C	C5-C6-O7-C8
8	B	1296	P4C	O13-C14-C15-O16
8	B	1296	P4C	C14-C15-O16-C17
8	E	1296	P4C	C2-C3-O4-C5
8	F	1294	P4C	O13-C14-C15-O16
8	B	1296	P4C	O7-C8-C9-O10
8	C	1294	P4C	C15-C14-O13-C12
8	F	1294	P4C	C9-C8-O7-C6
8	E	1296	P4C	C21-C20-O19-C18
8	E	1296	P4C	C6-C5-O4-C3
8	E	1295	P4C	C5-C6-O7-C8
8	E	1296	P4C	C17-C18-O19-C20
8	B	1296	P4C	O10-C11-C12-O13
8	F	1294	P4C	O7-C8-C9-O10
8	C	1294	P4C	C5-C6-O7-C8
8	F	1294	P4C	O4-C5-C6-O7
8	F	1294	P4C	C14-C15-O16-C17
9	F	1293	NAG	C3-C2-N2-C7
8	F	1294	P4C	C18-C17-O16-C15
8	C	1294	P4C	C21-C20-O19-C18
8	E	1295	P4C	O1-C2-C3-O4
8	B	1296	P4C	C8-C9-O10-C11
5	B	1289	EPE	C10-C9-N1-C2
8	E	1296	P4C	C18-C17-O16-C15
8	E	1295	P4C	C14-C15-O16-C17
8	C	1294	P4C	C17-C18-O19-C20
8	F	1294	P4C	O16-C17-C18-O19
8	B	1296	P4C	O16-C17-C18-O19

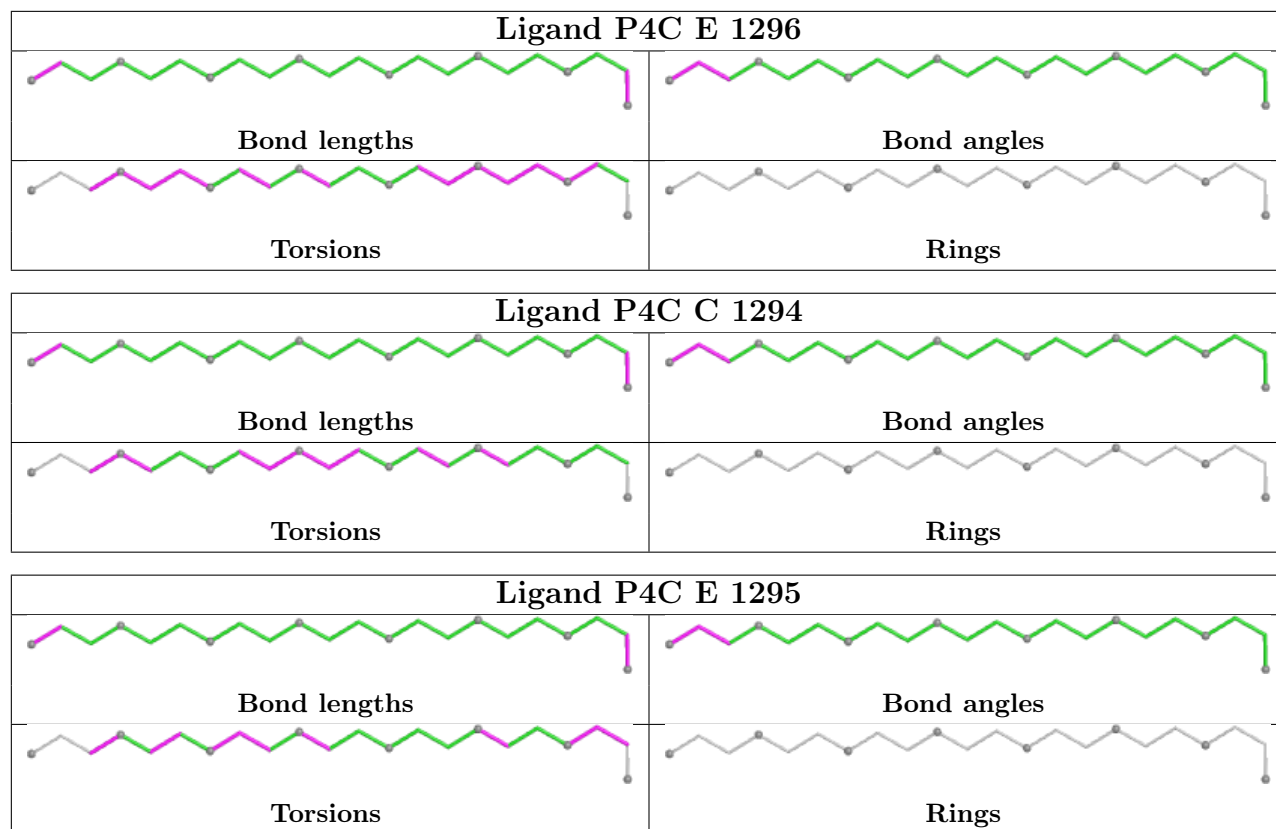
There are no ring outliers.

14 monomers are involved in 48 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	1290	GLA	1	0
8	B	1296	P4C	3	0
7	E	1294	MAN	2	0
6	F	1290	GLA	1	0
4	B	1290	ACT	10	0
6	E	1289	GLA	6	0
8	F	1294	P4C	4	0
8	E	1296	P4C	1	0
8	C	1294	P4C	1	0
4	E	1290	ACT	14	0
5	B	1289	EPE	1	0
8	E	1295	P4C	2	0
9	C	1293	NAG	1	0
5	F	1289	EPE	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 all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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 ⓘ

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	217/218 (99%)	0.55	28 (12%) 7 8	18, 38, 45, 52	3 (1%)
1	B	218/218 (100%)	-0.65	1 (0%) 87 89	23, 38, 44, 53	3 (1%)
1	C	214/218 (98%)	-0.46	0 100 100	20, 38, 45, 51	2 (0%)
1	D	216/218 (99%)	0.45	23 (10%) 11 12	18, 38, 45, 58	2 (0%)
1	E	218/218 (100%)	-0.92	0 100 100	19, 38, 45, 53	1 (0%)
1	F	217/218 (99%)	-0.04	8 (3%) 45 49	32, 38, 46, 63	0
All	All	1300/1308 (99%)	-0.18	60 (4%) 37 42	18, 38, 45, 63	11 (0%)

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	258	THR	3.9
1	A	199	GLY	3.8
1	A	257	GLY	3.6
1	A	263	ALA	3.5
1	D	212	PHE	3.3
1	D	227	LEU	3.2
1	A	246	HIS	3.0
1	A	173	PHE	3.0
1	D	74	LEU	2.9
1	A	245	CYS	2.8
1	F	274	TYR	2.8
1	A	278	TYR	2.8
1	D	246	HIS	2.7
1	D	98	LEU	2.7
1	D	187	VAL	2.7
1	D	191	ALA	2.7
1	D	194	TYR	2.7
1	A	272	LYS	2.7
1	A	195	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	A	208	ASP	2.6
1	A	269	LYS	2.6
1	A	236	PHE	2.6
1	A	207	GLY	2.6
1	A	229	THR	2.6
1	F	199	GLY	2.5
1	A	268	TRP	2.5
1	D	243	LYS	2.4
1	A	273	GLY	2.4
1	D	200	ALA	2.4
1	A	271	GLY	2.3
1	A	216	GLN	2.3
1	A	212[A]	PHE	2.3
1	F	211	THR	2.3
1	A	210	LEU	2.3
1	D	73	CYS	2.3
1	D	207	GLY	2.2
1	D	263	ALA	2.2
1	F	262	PHE	2.2
1	B	229	THR	2.2
1	F	263	ALA	2.2
1	D	163	THR	2.2
1	A	200	ALA	2.2
1	A	261	SER	2.2
1	A	225	ASN	2.2
1	A	179	PHE	2.1
1	D	236	PHE	2.1
1	D	199	GLY	2.1
1	A	167	ARG	2.1
1	D	210	LEU	2.1
1	F	227	LEU	2.1
1	D	274	TYR	2.1
1	D	241	TRP	2.1
1	A	171	VAL	2.1
1	F	197	VAL	2.1
1	D	245	CYS	2.1
1	A	274	TYR	2.1
1	D	228	ASN	2.0
1	D	208	ASP	2.0
1	D	179	PHE	2.0
1	F	215	ASN	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 ⓘ

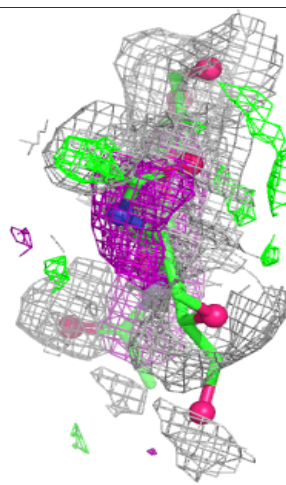
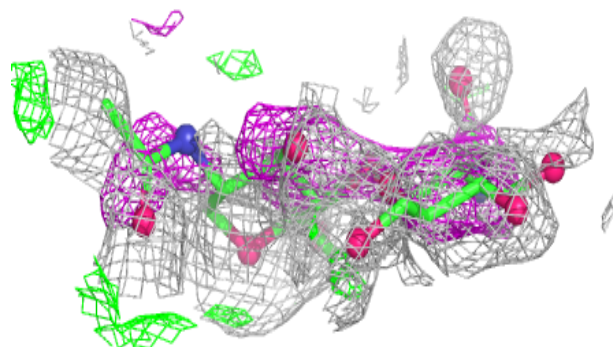
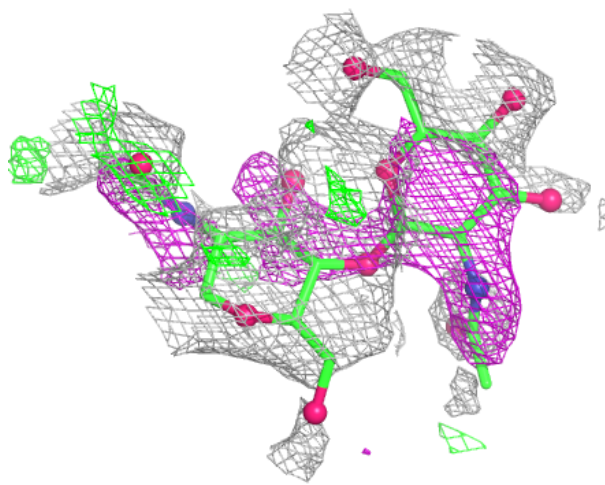
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

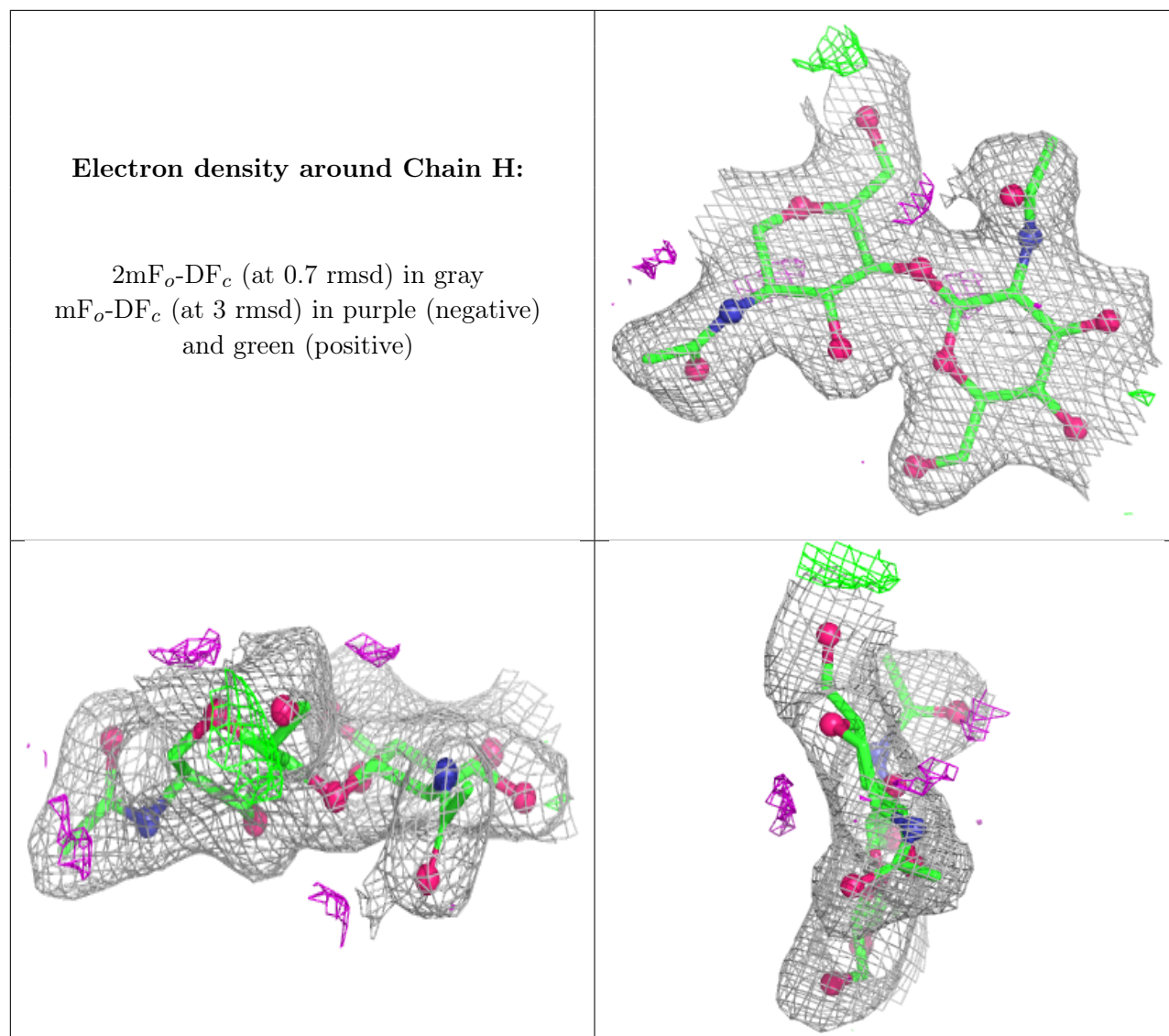
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	G	2	14/15	0.90	0.12	44,45,47,47	0
2	NAG	G	1	14/15	0.93	0.10	41,43,44,45	0
2	NAG	H	2	14/15	0.98	0.05	38,40,42,42	0
2	NAG	H	1	14/15	0.99	0.04	35,36,39,41	0

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.

Electron density around Chain G:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	ACT	B	407	4/4	0.95	0.08	67,67,67,67	0
6	GLA	C	1290	12/12	0.95	0.07	52,54,55,55	0
5	EPE	B	1289	12/15	0.96	0.07	59,64,65,65	0
5	EPE	C	1289	12/15	0.96	0.08	51,53,53,54	0
5	EPE	F	1289	12/15	0.96	0.08	47,49,51,52	0
4	ACT	B	1290	4/4	0.96	0.09	36,37,37,38	0
6	GLA	F	1290	12/12	0.96	0.07	51,53,54,55	0

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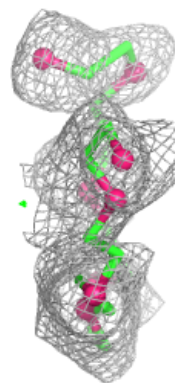
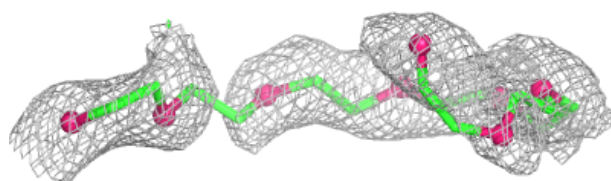
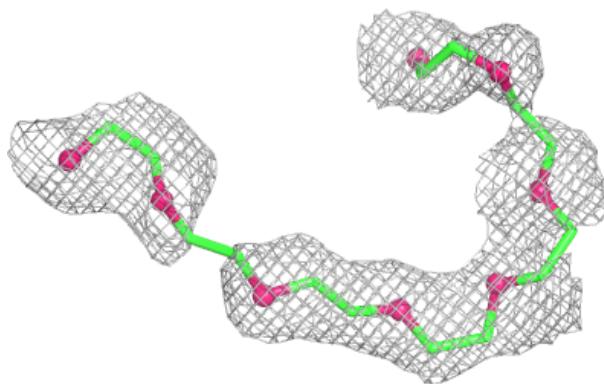
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
7	MAN	B	1295	11/12	0.96	0.08	82,83,83,83	0
8	P4C	B	1296	22/22	0.96	0.07	47,51,56,57	0
8	P4C	E	1295	22/22	0.96	0.06	50,53,62,62	0
9	NAG	C	1293	15/15	0.96	0.05	64,65,66,66	0
9	NAG	F	1293	15/15	0.96	0.07	67,69,70,70	0
7	MAN	E	1294	11/12	0.97	0.07	59,59,60,60	0
4	ACT	F	1291	4/4	0.97	0.12	36,36,37,38	0
8	P4C	C	1294	22/22	0.97	0.06	62,66,69,70	0
6	GLA	E	1289	12/12	0.97	0.07	53,57,59,59	0
8	P4C	E	1296	22/22	0.97	0.06	50,54,56,57	0
8	P4C	F	1294	22/22	0.97	0.06	59,59,62,62	0
3	CA	A	1289	1/1	0.97	0.06	83,83,83,83	0
6	GLA	B	1291	12/12	0.97	0.05	45,47,47,48	0
4	ACT	E	1290	4/4	0.98	0.07	37,38,38,39	0
3	CA	C	1292	1/1	0.99	0.05	42,42,42,42	0
4	ACT	C	1291	4/4	0.99	0.08	36,36,36,37	0
3	CA	D	1289	1/1	0.99	0.03	62,62,62,62	0
3	CA	F	1292	1/1	0.99	0.07	40,40,40,40	0
3	CA	B	1292	1/1	0.99	0.09	26,26,26,26	0
3	CA	E	1291	1/1	1.00	0.07	23,23,23,23	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

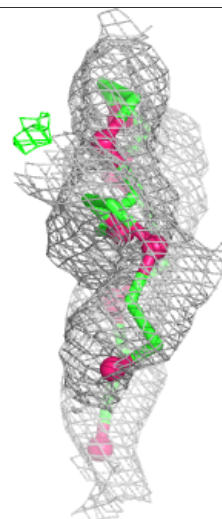
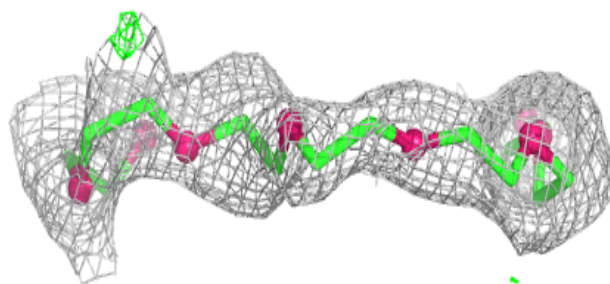
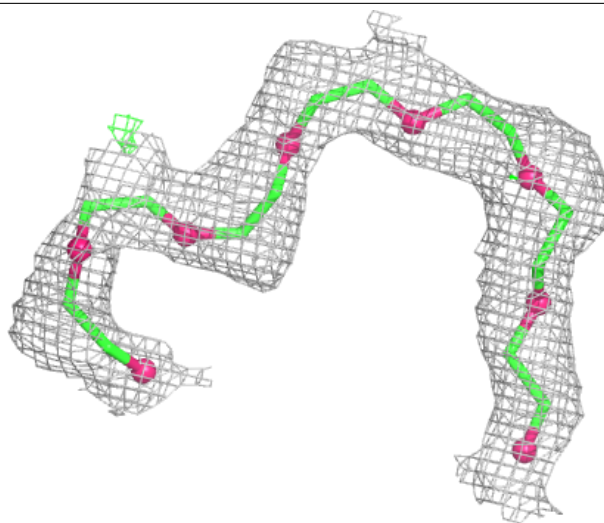
Electron density around P4C B 1296:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



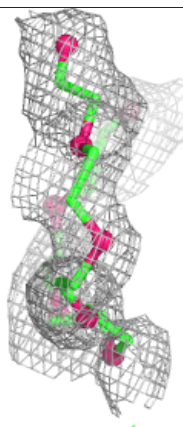
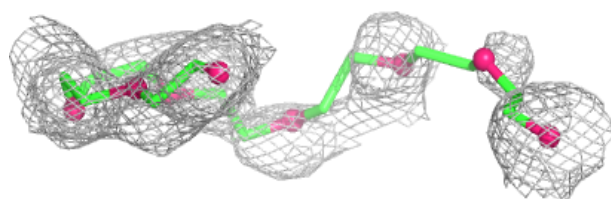
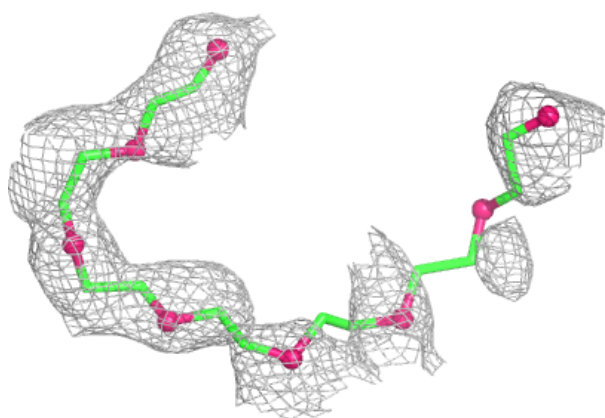
Electron density around P4C E 1295:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

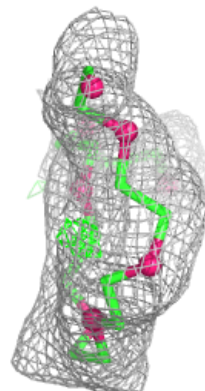
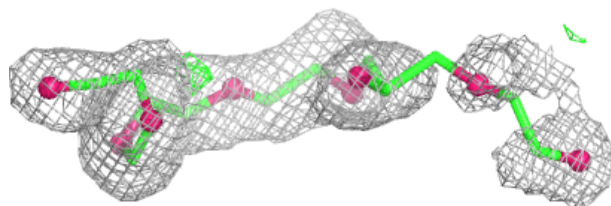
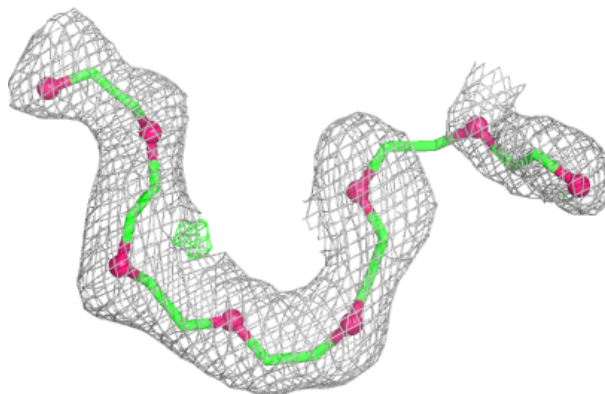


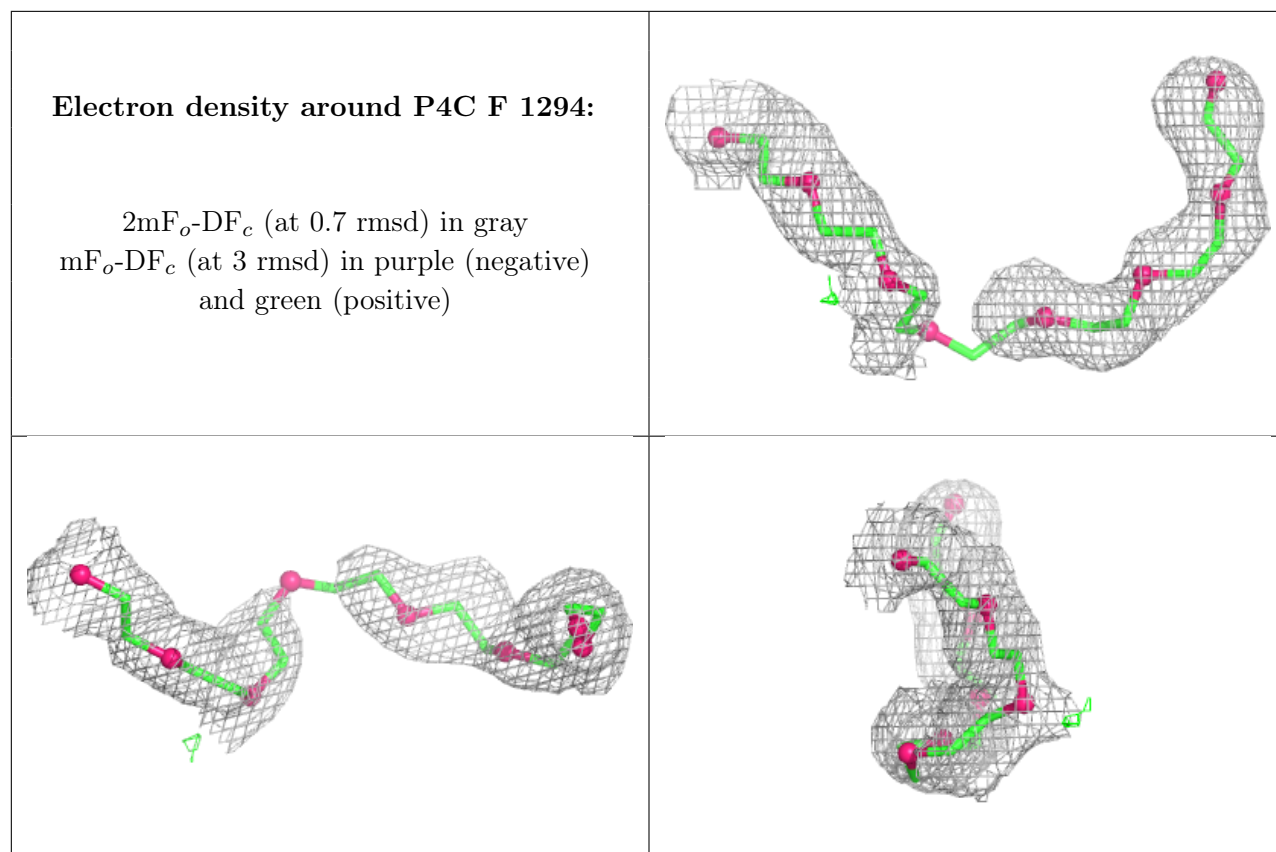
Electron density around P4C C 1294:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around P4C E 1296:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





6.5 Other polymers [i](#)

There are no such residues in this entry.