



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 09:54 PM UTC

PDB ID : 4DT2 / pdb_00004dt2
Title : Crystal structure of red kidney bean purple acid phosphatase in complex with Maybridge fragment CC27209
Authors : Feder, D.; Hussein, W.M.; Clayton, D.J.; Kan, M.; Schenk, G.; McGeary, R.P.; Guddat, L.W.
Deposited on : 2012-02-20
Resolution : 2.70 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

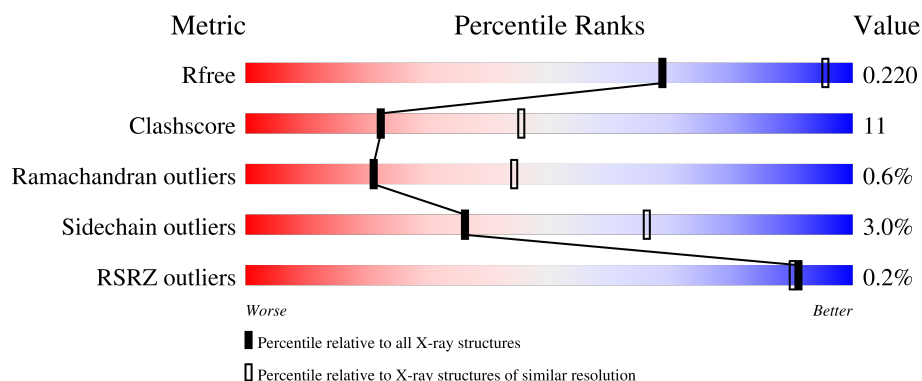
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

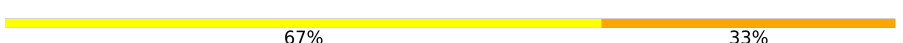
The reported resolution of this entry is 2.70 Å.

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.



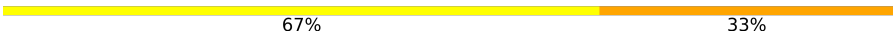
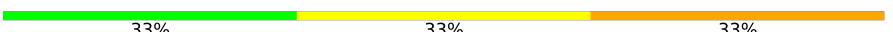
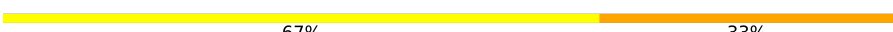
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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	426	 77% 23% .
1	B	426	 77% 22% .
1	C	426	 71% 27% .
1	D	426	 79% 20% .
2	E	3	 67% 33%

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Mol	Chain	Length	Quality of chain
2	F	3	 67% 33%
2	G	3	 33% 67%
2	H	3	 33% 33% 33%
2	J	3	 67% 33%
2	K	3	 67% 33%
2	L	3	 33% 67%
2	M	3	 33% 67%
3	I	2	 100%

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
10	EDO	A	512	-	-	X	-
7	OLV	A	504	-	-	X	-
7	OLV	B	504	-	-	X	-
7	OLV	C	512	-	-	X	-
7	OLV	D	511	-	-	X	-

2 Entry composition [i](#)

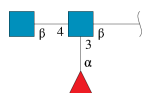
There are 12 unique types of molecules in this entry. The entry contains 15025 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Purple acid phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	426	Total	C	N	O	S	0	0	0
			3510	2253	610	637	10			
1	B	426	Total	C	N	O	S	0	0	0
			3510	2253	610	637	10			
1	C	426	Total	C	N	O	S	1	0	0
			3510	2253	610	637	10			
1	D	426	Total	C	N	O	S	0	0	0
			3510	2253	610	637	10			

- Molecule 2 is an oligosaccharide called alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	F	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	G	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	H	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	J	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	K	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	L	3	Total	C	N	O	0	0	0
			38	22	2	14			
2	M	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 3 is an oligosaccharide called alpha-L-fucopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	I	2	Total	C	N	O	0	0	0
			24	14	1	9			

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		
4	B	1	Total	Zn	0	0
			1	1		
4	C	1	Total	Zn	0	0
			1	1		
4	D	1	Total	Zn	0	0
			1	1		

- Molecule 5 is FE (III) ION (CCD ID: FE) (formula: Fe).

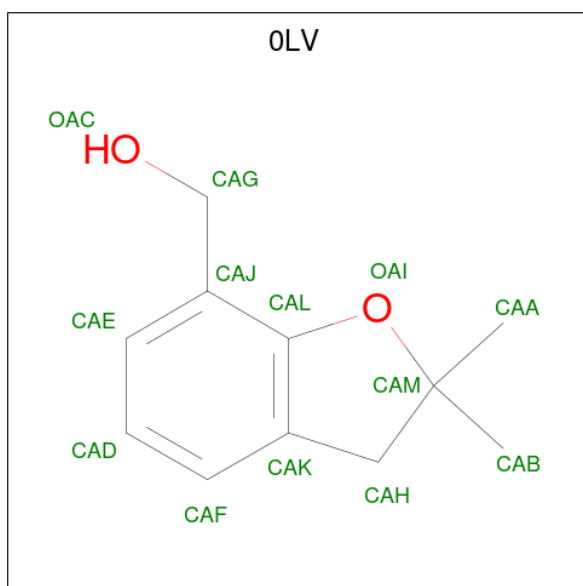
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total	Fe	0	0
			1	1		
5	B	1	Total	Fe	0	0
			1	1		
5	C	1	Total	Fe	0	0
			1	1		
5	D	1	Total	Fe	0	0
			1	1		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



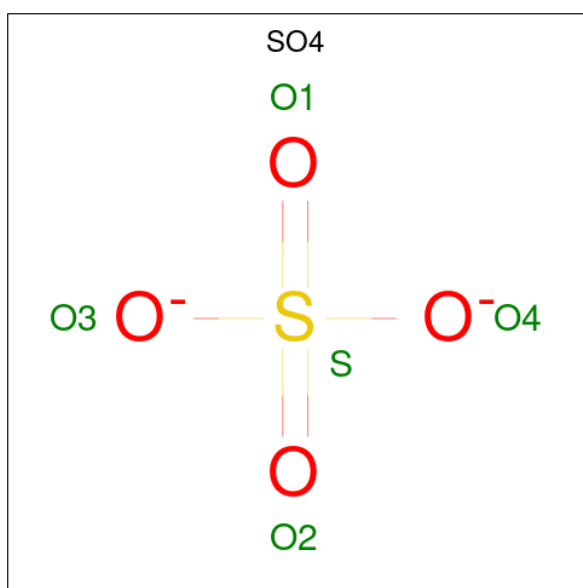
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		
6	C	1	Total	C	O	0	0
			6	3	3		
6	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is (2,2-dimethyl-2,3-dihydro-1-benzofuran-7-yl)methanol (CCD ID: 0LV) (formula: C₁₁H₁₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			13	11	2		
7	B	1	Total	C	O	0	0
			13	11	2		
7	C	1	Total	C	O	0	0
			13	11	2		
7	D	1	Total	C	O	0	0
			13	11	2		

- Molecule 8 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



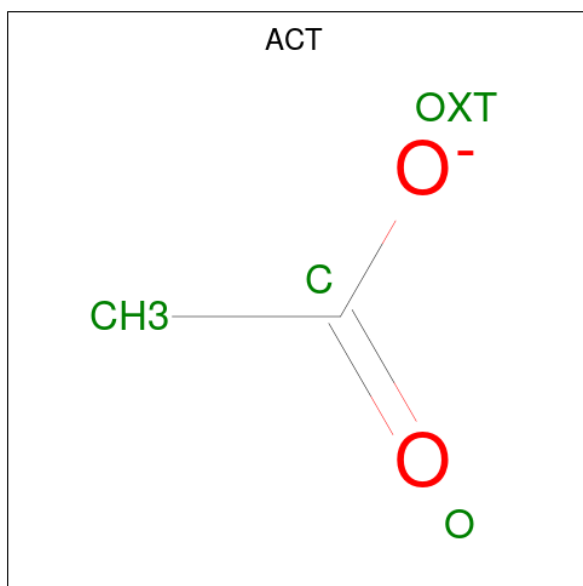
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	A	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		
8	B	1	Total	O	S	0	0
			5	4	1		

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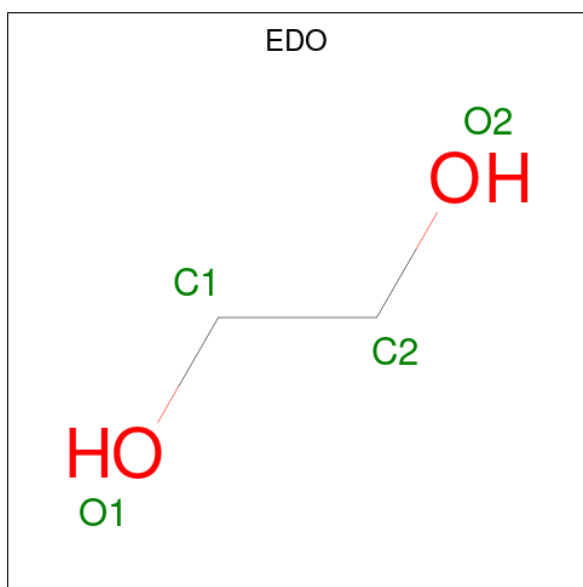
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	B	1	Total	O	S	0	0
			5	4	1		
8	C	1	Total	O	S	0	0
			5	4	1		
8	C	1	Total	O	S	0	0
			5	4	1		
8	C	1	Total	O	S	0	0
			5	4	1		
8	D	1	Total	O	S	0	0
			5	4	1		
8	D	1	Total	O	S	0	0
			5	4	1		
8	D	1	Total	O	S	0	0
			5	4	1		
8	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 9 is ACETATE ION (CCD ID: ACT) (formula: $C_2H_3O_2$).



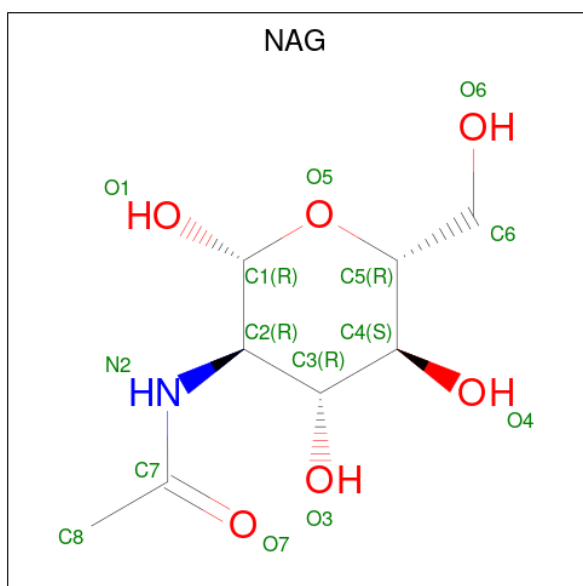
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 10 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
10	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 11 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	A	1	Total	C	N	O	0	0
			14	8	1	5		
11	B	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
11	B	1	Total	C	N	O	0	0
			14	8	1	5		
11	C	1	Total	C	N	O	0	0
			14	8	1	5		
11	C	1	Total	C	N	O	0	0
			14	8	1	5		
11	D	1	Total	C	N	O	0	0
			14	8	1	5		
11	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 12 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
12	A	127	Total	O	0	0
			127	127		
12	B	85	Total	O	0	0
			85	85		
12	C	85	Total	O	0	0
			85	85		
12	D	96	Total	O	0	0
			96	96		

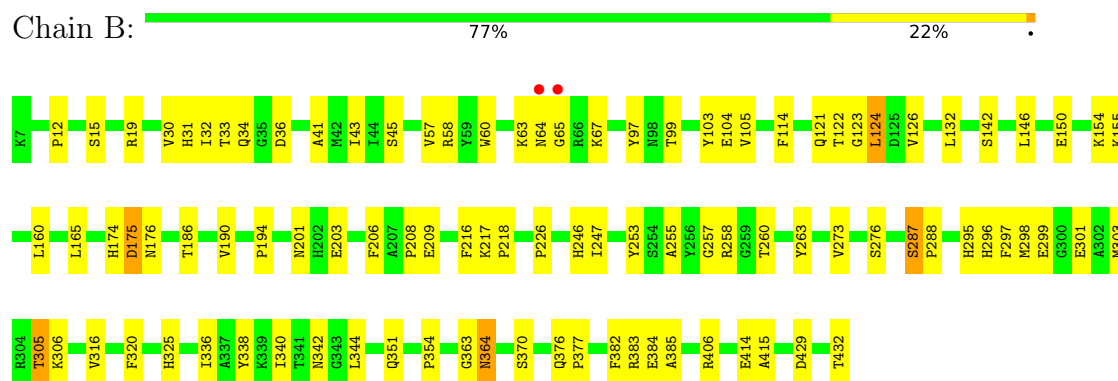
3 Residue-property plots

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.

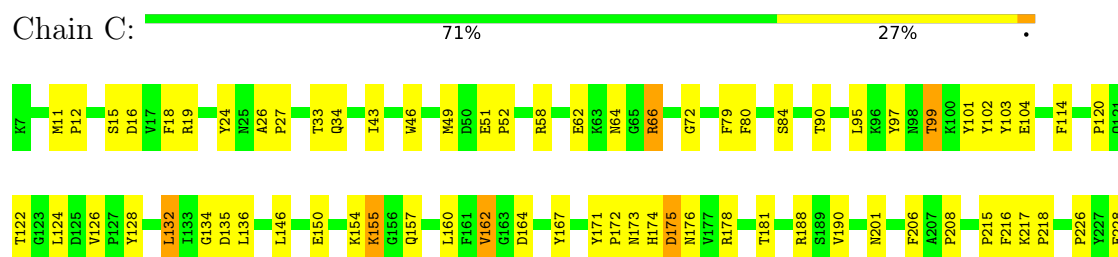
• Molecule 1: Purple acid phosphatase

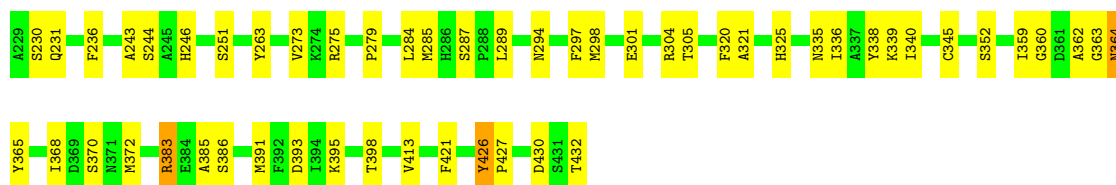


• Molecule 1: Purple acid phosphatase



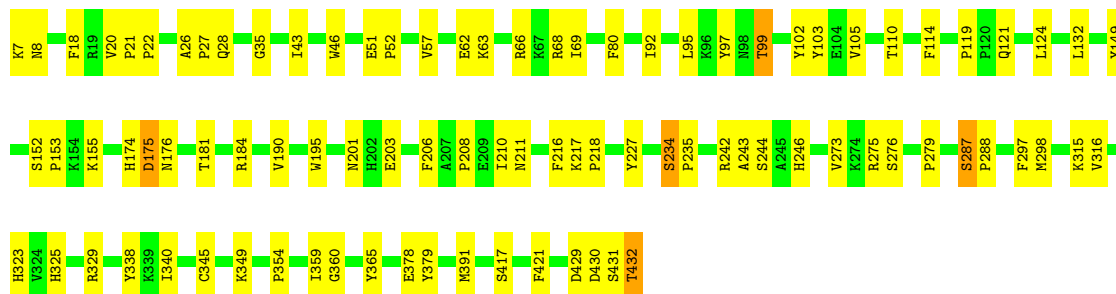
• Molecule 1: Purple acid phosphatase





- Molecule 1: Purple acid phosphatase

Chain D: 79% 20% .



- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E: 67% 33%



- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F: 67% 33%



- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain G: 33% 67%



- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H: 33% 33% 33%



- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  67% 33%



- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  67% 33%



- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  33% 67%

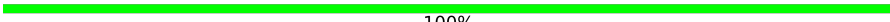


- Molecule 2: alpha-L-fucopyranose-(1-3)-[2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  33% 67%



- Molecule 3: alpha-L-fucopyranose-(1-3)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	126.76Å 126.76Å 298.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	19.72 – 2.70 19.72 – 2.70	Depositor EDS
% Data completeness (in resolution range)	99.8 (19.72-2.70) 99.5 (19.72-2.70)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.67 (at 2.71Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.7_650)	Depositor
R, R_{free}	0.175 , 0.228 0.170 , 0.220	Depositor DCC
R_{free} test set	3861 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	45.4	Xtriage
Anisotropy	0.049	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 26.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.034 for -h,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	15025	wwPDB-VP
Average B, all atoms (Å ²)	42.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.41% 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: FE, SO4, ACT, GOL, OLV, NAG, EDO, ZN, FUC

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.54	0/3629	0.84	2/4934 (0.0%)
1	B	0.51	0/3629	0.81	3/4934 (0.1%)
1	C	0.51	0/3629	0.82	5/4934 (0.1%)
1	D	0.53	0/3629	0.84	7/4934 (0.1%)
All	All	0.52	0/14516	0.83	17/19736 (0.1%)

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	426	TYR	CA-C-N	7.02	128.61	119.84
1	C	426	TYR	C-N-CA	7.02	128.61	119.84
1	B	287	SER	CA-C-N	6.27	126.48	119.90
1	B	287	SER	C-N-CA	6.27	126.48	119.90
1	D	227	TYR	N-CA-C	6.22	118.14	111.36
1	A	278	THR	CA-C-N	6.16	125.88	119.05
1	A	278	THR	C-N-CA	6.16	125.88	119.05
1	C	287	SER	CA-C-N	6.02	125.96	119.76
1	C	287	SER	C-N-CA	6.02	125.96	119.76
1	D	234	SER	CA-C-N	5.73	125.83	119.87
1	D	234	SER	C-N-CA	5.73	125.83	119.87
1	D	287	SER	CA-C-N	5.71	125.62	119.85
1	D	287	SER	C-N-CA	5.71	125.62	119.85
1	D	20	VAL	N-CA-C	5.25	113.06	107.76
1	D	243	ALA	CB-CA-C	-5.15	110.65	116.63
1	B	124	LEU	N-CA-C	5.12	117.27	111.02
1	C	413	VAL	N-CA-C	5.06	115.14	107.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3510	0	3326	77	0
1	B	3510	0	3327	71	0
1	C	3510	0	3327	84	0
1	D	3510	0	3327	63	0
2	E	38	0	34	2	0
2	F	38	0	34	1	0
2	G	38	0	34	1	0
2	H	38	0	34	1	0
2	J	38	0	34	5	0
2	K	38	0	34	0	0
2	L	38	0	34	2	0
2	M	38	0	34	0	0
3	I	24	0	22	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	1	0	0	0	0
5	B	1	0	0	0	0
5	C	1	0	0	0	0
5	D	1	0	0	0	0
6	A	6	0	8	2	0
6	C	6	0	8	1	0
6	D	6	0	8	0	0
7	A	13	0	13	9	0
7	B	13	0	13	10	0
7	C	13	0	14	13	0
7	D	13	0	14	13	0
8	A	30	0	0	2	0
8	B	15	0	0	1	0
8	C	15	0	0	0	0
8	D	20	0	0	0	0
9	A	4	0	3	0	0
10	A	4	0	6	6	0
11	A	14	0	13	0	0
11	B	28	0	26	0	0
11	C	28	0	26	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
11	D	28	0	26	0	0
12	A	127	0	0	3	0
12	B	85	0	0	0	0
12	C	85	0	0	2	0
12	D	96	0	0	0	0
All	All	15025	0	13779	301	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 11.

All (301) 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:258:ARG:HH22	7:C:512:0LV:H4	1.13	1.08
1:D:323:HIS:CE1	7:D:511:0LV:H9	1.93	1.03
1:C:124:LEU:HD12	1:C:279:PRO:HG3	1.46	0.95
1:A:296:HIS:HE1	7:A:504:0LV:H2	1.30	0.95
1:C:325:HIS:CE1	7:C:512:0LV:H11	2.02	0.94
1:D:7:LYS:HG2	1:D:8:ASN:H	1.30	0.94
1:C:325:HIS:HE1	7:C:512:0LV:H11	1.34	0.92
1:B:258:ARG:HH12	7:C:512:0LV:H2	1.38	0.87
1:C:66:ARG:HG2	1:C:66:ARG:HH11	1.39	0.87
1:A:258:ARG:HH22	7:D:511:0LV:H5	1.41	0.85
1:C:201:ASN:HD21	7:C:512:0LV:H9	1.44	0.83
1:B:258:ARG:NH2	7:C:512:0LV:H4	1.94	0.81
1:C:201:ASN:HD21	7:C:512:0LV:CAF	1.94	0.80
1:A:296:HIS:CE1	7:A:504:0LV:H2	2.17	0.78
1:D:124:LEU:HD12	1:D:279:PRO:HG3	1.66	0.77
1:C:66:ARG:HH11	1:C:66:ARG:CG	1.99	0.74
1:D:323:HIS:HE1	7:D:511:0LV:H9	1.53	0.73
1:C:372:MET:SD	1:C:383:ARG:HD3	2.28	0.73
1:D:365:TYR:HE1	7:D:511:0LV:H12	1.54	0.73
1:A:217:LYS:HB3	1:A:218:PRO:HD3	1.70	0.73
1:B:325:HIS:CE1	7:B:504:0LV:H11	2.24	0.72
1:C:325:HIS:HE1	7:C:512:0LV:CAE	2.01	0.72
1:B:217:LYS:HB3	1:B:218:PRO:HD3	1.72	0.72
1:A:124:LEU:HD12	1:A:279:PRO:HG3	1.72	0.72
1:A:323:HIS:CE1	7:A:504:0LV:H9	2.26	0.70
1:C:325:HIS:HD2	1:C:364:ASN:ND2	1.89	0.70
1:A:258:ARG:NH2	7:D:511:0LV:H5	2.06	0.70
1:C:154:LYS:HB2	1:C:391:MET:HE2	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:57:VAL:HG11	1:D:92:ILE:HD11	1.75	0.69
1:D:217:LYS:HB3	1:D:218:PRO:HD3	1.76	0.68
1:C:421:PHE:CD2	1:C:430:ASP:HB3	2.29	0.68
1:C:201:ASN:ND2	7:C:512:OLV:H9	2.08	0.67
1:B:34:GLN:O	1:B:226:PRO:HG3	1.93	0.67
1:D:421:PHE:CD2	1:D:430:ASP:HB3	2.30	0.67
1:D:323:HIS:ND1	7:D:511:OLV:H9	2.08	0.67
1:B:316:VAL:HB	1:B:354:PRO:HG3	1.75	0.67
1:B:258:ARG:HH22	7:C:512:OLV:CAA	2.01	0.66
2:J:3:NAG:H82	2:J:3:NAG:H3	1.76	0.66
1:C:132:LEU:HD22	1:C:320:PHE:CD1	2.31	0.65
1:B:301:GLU:O	1:B:305:THR:HG23	1.96	0.65
1:C:285:MET:O	1:C:321:ALA:HA	1.97	0.65
1:B:124:LEU:HD11	1:B:276:SER:O	1.97	0.65
1:B:342:ASN:ND2	1:B:344:LEU:HD12	2.11	0.65
1:D:174:HIS:O	1:D:175:ASP:C	2.40	0.65
1:D:7:LYS:HG2	1:D:8:ASN:N	2.08	0.65
1:B:406:ARG:HD3	1:B:415:ALA:CB	2.27	0.64
1:A:336:ILE:HA	6:A:503:GOL:H32	1.80	0.64
1:C:188:ARG:HD3	12:C:635:HOH:O	1.98	0.64
1:A:371:ASN:H	10:A:512:EDO:H21	1.62	0.64
1:C:12:PRO:HD2	1:C:15:SER:OG	1.97	0.64
1:A:201:ASN:HD21	7:A:504:OLV:CAF	2.11	0.63
1:B:33:THR:HA	1:B:194:PRO:HB3	1.80	0.62
1:D:244:SER:O	1:D:279:PRO:HD2	2.00	0.62
1:A:124:LEU:HD13	1:A:276:SER:O	1.98	0.62
1:A:174:HIS:O	1:A:175:ASP:C	2.43	0.62
1:C:217:LYS:HB3	1:C:218:PRO:HD3	1.82	0.62
1:C:16:ASP:HA	1:C:19:ARG:HD3	1.81	0.61
1:A:103:TYR:CE1	1:A:114:PHE:HB2	2.36	0.61
1:B:203:GLU:O	1:B:216:PHE:HA	2.00	0.61
1:D:325:HIS:CE1	7:D:511:OLV:H11	2.36	0.61
1:D:103:TYR:CZ	1:D:114:PHE:HB2	2.36	0.60
1:D:275:ARG:NH2	1:D:315:LYS:O	2.33	0.60
1:A:371:ASN:H	10:A:512:EDO:C2	2.13	0.60
1:C:134:GLY:HA3	1:C:162:VAL:HG23	1.82	0.60
1:C:72:GLY:HA3	1:C:90:THR:OG1	2.01	0.60
1:A:34:GLN:O	1:A:226:PRO:HG3	2.01	0.60
1:D:119:PRO:HG3	1:D:242:ARG:CZ	2.31	0.60
1:B:295:HIS:HE1	7:B:504:OLV:H13	1.67	0.59
1:A:258:ARG:HH12	7:D:511:OLV:H2	1.66	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:201:ASN:HD21	7:B:504:0LV:CAF	2.15	0.59
1:B:325:HIS:NE2	7:B:504:0LV:H11	2.17	0.59
1:C:24:TYR:CD2	2:J:1:NAG:H82	2.38	0.59
1:C:244:SER:HB2	1:C:279:PRO:HD2	1.83	0.58
1:A:370:SER:N	10:A:512:EDO:H22	2.17	0.58
1:C:206:PHE:CZ	1:C:208:PRO:HG3	2.38	0.58
1:D:338:TYR:CZ	1:D:340:ILE:HA	2.38	0.58
1:C:160:LEU:HB3	1:C:284:LEU:HD11	1.86	0.58
1:B:32:ILE:HA	1:B:43:ILE:O	2.05	0.57
1:B:146:LEU:O	1:B:150:GLU:HG3	2.04	0.57
1:D:325:HIS:HA	1:D:360:GLY:O	2.04	0.57
1:C:174:HIS:O	1:C:175:ASP:C	2.47	0.57
1:B:160:LEU:HD21	1:B:247:ILE:HD13	1.87	0.57
1:C:155:LYS:HD3	1:C:155:LYS:C	2.29	0.57
1:D:246:HIS:CE1	1:D:273:VAL:HG22	2.39	0.57
2:E:2:FUC:H61	2:E:3:NAG:H83	1.87	0.57
1:C:27:PRO:HB3	1:C:46:TRP:CD1	2.40	0.56
2:L:2:FUC:H5	2:L:3:NAG:H2	1.86	0.56
1:D:201:ASN:ND2	7:D:511:0LV:H1	2.19	0.56
1:D:153:PRO:O	1:D:155:LYS:HG3	2.05	0.56
1:B:165:LEU:HD13	1:B:203:GLU:CD	2.31	0.56
1:D:429:ASP:OD2	1:D:431:SER:HB3	2.05	0.56
1:D:27:PRO:HB3	1:D:46:TRP:CD1	2.40	0.56
1:D:206:PHE:CZ	1:D:208:PRO:HG3	2.40	0.56
1:A:132:LEU:HD22	1:A:320:PHE:CD1	2.41	0.56
1:A:258:ARG:HH12	7:D:511:0LV:CAH	2.19	0.56
1:D:201:ASN:HD21	7:D:511:0LV:H1	1.70	0.56
1:D:275:ARG:NH1	1:D:279:PRO:O	2.38	0.56
1:B:296:HIS:O	1:B:299:GLU:HB2	2.06	0.55
1:A:97:TYR:O	1:A:99:THR:HG23	2.06	0.55
1:A:325:HIS:CE1	7:A:504:0LV:H11	2.42	0.55
1:A:336:ILE:HA	6:A:503:GOL:C3	2.36	0.55
1:C:336:ILE:HA	6:C:508:GOL:H12	1.88	0.55
1:D:97:TYR:O	1:D:99:THR:HG22	2.07	0.55
1:B:12:PRO:HD2	1:B:15:SER:OG	2.07	0.54
1:C:426:TYR:N	1:C:427:PRO:HD3	2.21	0.54
1:D:18:PHE:CD1	1:D:181:THR:HB	2.42	0.54
1:B:201:ASN:HD21	7:B:504:0LV:H9	1.73	0.54
1:C:171:TYR:CE1	1:C:178:ARG:HG3	2.42	0.54
1:C:359:ILE:O	1:C:359:ILE:HG13	2.07	0.54
1:B:338:TYR:CZ	1:B:340:ILE:HA	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:107:LEU:HD23	1:A:107:LEU:N	2.23	0.54
1:D:51:GLU:HB2	1:D:52:PRO:HD2	1.90	0.54
2:E:2:FUC:C6	2:E:3:NAG:H83	2.38	0.54
1:D:124:LEU:HB2	1:D:276:SER:O	2.08	0.53
1:D:316:VAL:HB	1:D:354:PRO:HG3	1.91	0.53
1:C:128:TYR:CE2	1:C:157:GLN:HB2	2.43	0.53
1:A:421:PHE:CD2	1:A:430:ASP:HB3	2.45	0.52
1:B:406:ARG:HD3	1:B:415:ALA:HB2	1.90	0.52
1:C:325:HIS:CD2	1:C:364:ASN:CG	2.87	0.52
1:B:103:TYR:CE2	1:B:114:PHE:HB2	2.44	0.52
1:B:297:PHE:CD2	1:B:298:MET:HG3	2.44	0.52
1:B:123:GLY:O	1:B:126:VAL:HG23	2.09	0.52
1:D:62:GLU:HB3	1:D:102:TYR:HE2	1.74	0.52
1:D:365:TYR:CE1	7:D:511:OLV:H12	2.40	0.52
1:D:18:PHE:CE1	1:D:181:THR:HB	2.45	0.52
1:A:166:SER:O	1:A:167:TYR:HB2	2.11	0.51
1:C:201:ASN:ND2	7:C:512:OLV:H1	2.25	0.51
1:A:25:ASN:OD1	1:A:48:THR:HB	2.11	0.51
1:C:325:HIS:HD2	1:C:364:ASN:CG	2.19	0.51
1:B:351:GLN:OE1	1:B:429:ASP:HA	2.11	0.51
1:A:338:TYR:CZ	1:A:340:ILE:HA	2.46	0.51
1:B:257:GLY:HA2	1:B:306:LYS:HE3	1.93	0.51
1:C:301:GLU:O	1:C:305:THR:HG23	2.10	0.51
1:C:325:HIS:CD2	1:C:364:ASN:ND2	2.76	0.51
1:B:60:TRP:HB3	1:B:67:LYS:HA	1.93	0.50
1:B:103:TYR:CZ	1:B:114:PHE:HB2	2.47	0.50
1:A:258:ARG:HH22	7:D:511:OLV:CAA	2.17	0.50
1:A:155:LYS:HE3	12:A:724:HOH:O	2.12	0.50
1:A:413:VAL:HG21	12:A:686:HOH:O	2.12	0.50
1:B:63:LYS:O	1:B:65:GLY:N	2.45	0.50
1:C:18:PHE:CD1	1:C:181:THR:HB	2.47	0.50
1:C:325:HIS:HA	1:C:360:GLY:O	2.11	0.50
1:B:201:ASN:ND2	7:B:504:OLV:H9	2.26	0.50
1:A:27:PRO:HB3	1:A:46:TRP:CD1	2.47	0.50
1:B:57:VAL:HG22	1:B:105:VAL:HG12	1.94	0.50
1:C:172:PRO:O	1:C:173:ASN:HB2	2.12	0.50
1:B:97:TYR:CE2	1:B:121:GLN:HG3	2.47	0.50
1:B:174:HIS:O	1:B:175:ASP:C	2.55	0.50
1:A:36:ASP:HB3	1:A:41:ALA:HB3	1.94	0.49
1:A:323:HIS:HE1	7:A:504:OLV:H9	1.77	0.49
1:D:429:ASP:CG	1:D:431:SER:HB3	2.38	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:295:HIS:HD2	1:A:369:ASP:OD2	1.94	0.49
1:B:190:VAL:HG22	1:B:190:VAL:O	2.13	0.49
1:C:190:VAL:O	1:C:190:VAL:HG22	2.13	0.49
1:A:97:TYR:CD2	1:A:121:GLN:HB2	2.47	0.49
1:A:301:GLU:O	1:A:305:THR:HG23	2.12	0.49
1:C:62:GLU:HG3	1:C:102:TYR:HE1	1.77	0.49
1:C:338:TYR:CZ	1:C:340:ILE:HA	2.48	0.49
1:D:297:PHE:CE2	1:D:298:MET:HG3	2.47	0.49
1:A:370:SER:H	10:A:512:EDO:H11	1.77	0.48
1:D:27:PRO:HG2	1:D:105:VAL:HG23	1.95	0.48
1:B:201:ASN:OD1	7:B:504:OLV:H9	2.12	0.48
1:D:329:ARG:HH22	1:D:432:THR:HG21	1.77	0.48
1:A:51:GLU:HB2	1:A:52:PRO:HD2	1.95	0.48
1:C:51:GLU:HB2	1:C:52:PRO:HD2	1.95	0.48
1:D:246:HIS:ND1	1:D:273:VAL:HG22	2.29	0.48
1:A:215:PRO:O	1:A:216:PHE:HB2	2.13	0.48
1:B:363:GLY:O	1:B:364:ASN:C	2.55	0.48
1:C:18:PHE:CE1	1:C:181:THR:HB	2.48	0.48
1:C:66:ARG:CG	1:C:66:ARG:NH1	2.67	0.48
1:C:103:TYR:CZ	1:C:114:PHE:HB2	2.49	0.48
1:C:246:HIS:CE1	1:C:273:VAL:HG22	2.49	0.47
1:C:79:PHE:O	1:C:80:PHE:HB3	2.14	0.47
1:D:35:GLY:HA2	1:D:43:ILE:HG13	1.96	0.47
1:C:393:ASP:OD2	1:C:395:LYS:HE2	2.15	0.47
1:D:203:GLU:O	1:D:216:PHE:HA	2.14	0.47
1:A:339:LYS:HG2	1:A:342:ASN:ND2	2.29	0.47
1:B:31:HIS:CD2	1:B:45:SER:HB2	2.50	0.47
1:A:74:MET:HA	1:A:87:ILE:O	2.15	0.47
1:C:95:LEU:HD22	1:C:101:TYR:CZ	2.50	0.47
1:B:58:ARG:O	1:B:103:TYR:HA	2.15	0.47
1:B:208:PRO:HD2	1:B:209:GLU:OE1	2.15	0.47
1:D:244:SER:HB2	1:D:279:PRO:HD2	1.97	0.47
1:B:201:ASN:HB3	1:B:253:TYR:CE1	2.51	0.46
1:B:31:HIS:NE2	1:B:45:SER:HB2	2.30	0.46
1:C:304:ARG:NH1	1:C:335:ASN:HB3	2.31	0.46
1:D:97:TYR:CE1	1:D:121:GLN:HA	2.50	0.46
1:B:295:HIS:CE1	7:B:504:OLV:H13	2.48	0.46
1:D:329:ARG:HH22	1:D:432:THR:CG2	2.28	0.46
1:D:234:SER:HA	1:D:235:PRO:HD3	1.80	0.45
1:B:201:ASN:ND2	7:B:504:OLV:H1	2.31	0.45
1:C:275:ARG:HA	1:C:275:ARG:HD2	1.63	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:209:GLU:HG2	1:A:210:ILE:HG23	1.98	0.45
1:A:78:ARG:NH2	1:C:84:SER:O	2.49	0.45
1:D:80:PHE:CD1	1:D:210:ILE:HB	2.52	0.45
1:A:289:LEU:HD13	1:A:304:ARG:HA	1.99	0.45
1:C:58:ARG:HB3	1:C:104:GLU:HB2	1.99	0.45
1:A:78:ARG:NH1	12:A:691:HOH:O	2.49	0.45
1:A:165:LEU:HD13	1:A:203:GLU:CD	2.42	0.45
1:B:342:ASN:HD22	1:B:344:LEU:HD12	1.82	0.45
1:C:236:PHE:O	1:C:251:SER:HB2	2.16	0.45
1:D:190:VAL:HG23	1:D:195:TRP:CD1	2.52	0.44
1:A:120:PRO:HB2	1:A:126:VAL:HG11	1.99	0.44
2:J:1:NAG:H61	2:J:3:NAG:C7	2.48	0.44
1:A:21:PRO:HA	1:A:22:PRO:HD3	1.84	0.44
1:B:255:ALA:HB1	1:B:260:THR:HG21	2.00	0.44
1:D:152:SER:HA	1:D:153:PRO:HD3	1.88	0.44
1:B:298:MET:SD	1:C:339:LYS:HA	2.57	0.44
2:H:1:NAG:O4	2:H:2:FUC:H5	2.17	0.44
1:B:297:PHE:CE2	1:B:298:MET:HG3	2.53	0.44
1:A:372:MET:SD	1:A:383:ARG:HD3	2.58	0.43
1:B:186:THR:O	1:B:190:VAL:HG12	2.18	0.43
1:A:325:HIS:HE1	7:A:504:OLV:H13	1.83	0.43
1:B:287:SER:HA	1:B:288:PRO:HD3	1.76	0.43
2:J:3:NAG:H3	2:J:3:NAG:C8	2.46	0.43
1:A:421:PHE:HB3	1:A:426:TYR:O	2.19	0.43
1:C:26:ALA:HA	1:C:27:PRO:HD3	1.89	0.43
1:C:58:ARG:O	1:C:103:TYR:HA	2.18	0.43
1:C:134:GLY:CA	1:C:162:VAL:HG23	2.48	0.43
1:C:174:HIS:N	1:C:174:HIS:CD2	2.86	0.43
1:C:363:GLY:O	1:C:364:ASN:C	2.62	0.43
1:A:217:LYS:HB3	1:A:218:PRO:CD	2.46	0.43
1:C:146:LEU:O	1:C:150:GLU:HG3	2.19	0.43
2:J:2:FUC:H5	2:J:3:NAG:C1	2.49	0.43
1:B:154:LYS:HE3	1:B:414:GLU:OE2	2.18	0.43
1:A:234:SER:HA	1:A:235:PRO:HD3	1.83	0.43
1:B:246:HIS:ND1	1:B:273:VAL:HG22	2.34	0.43
1:B:298:MET:HE2	1:B:298:MET:HB3	1.85	0.43
1:C:263:TYR:CD1	1:C:263:TYR:C	2.97	0.43
1:D:338:TYR:OH	1:D:340:ILE:HA	2.19	0.43
1:C:201:ASN:HD22	7:C:512:OLV:H1	1.83	0.42
1:A:201:ASN:HA	1:A:204:ILE:HG13	2.01	0.42
1:C:33:THR:HG22	1:C:43:ILE:HB	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:34:GLN:O	1:C:226:PRO:HG3	2.19	0.42
1:D:21:PRO:HA	1:D:22:PRO:HD3	1.95	0.42
1:D:379:TYR:CD1	1:D:379:TYR:C	2.97	0.42
1:A:164:ASP:HB3	1:A:167:TYR:CE1	2.54	0.42
1:C:66:ARG:HD3	1:C:66:ARG:C	2.45	0.42
1:D:66:ARG:HD2	1:D:68:ARG:CZ	2.49	0.42
1:A:45:SER:HA	1:A:88:HIS:O	2.19	0.42
1:B:201:ASN:CG	7:B:504:OLV:H9	2.45	0.42
1:C:11:MET:HA	1:C:12:PRO:HD3	1.83	0.42
1:C:215:PRO:O	1:C:216:PHE:HB2	2.20	0.42
1:C:97:TYR:O	1:C:99:THR:HG23	2.20	0.42
1:C:289:LEU:HG	12:C:614:HOH:O	2.20	0.42
1:D:287:SER:HA	1:D:288:PRO:HD3	1.77	0.42
1:C:135:ASP:HB3	1:C:167:TYR:OH	2.20	0.42
1:C:136:LEU:HD13	1:C:362:ALA:HB1	2.02	0.42
1:C:370:SER:HA	1:C:385:ALA:HB2	2.02	0.42
1:A:122:THR:HA	1:A:243:ALA:O	2.20	0.42
1:B:376:GLN:HA	1:B:377:PRO:HD3	1.81	0.42
1:D:97:TYR:CD1	1:D:121:GLN:HA	2.55	0.42
2:G:1:NAG:C6	2:G:3:NAG:C1	2.98	0.42
1:A:79:PHE:O	1:A:80:PHE:HB3	2.20	0.41
1:A:289:LEU:N	1:A:289:LEU:HD23	2.35	0.41
1:A:338:TYR:CE1	1:D:345:CYS:HB2	2.55	0.41
1:B:382:PHE:HA	8:B:509:SO4:O2	2.20	0.41
1:C:421:PHE:HB3	1:C:426:TYR:O	2.19	0.41
1:A:66:ARG:HB2	8:A:505:SO4:O1	2.21	0.41
1:D:28:GLN:OE1	1:D:184:ARG:HD3	2.20	0.41
1:B:206:PHE:CZ	1:B:208:PRO:HG3	2.54	0.41
1:D:103:TYR:CE2	1:D:114:PHE:HB2	2.56	0.41
1:D:329:ARG:NH2	1:D:432:THR:HB	2.35	0.41
1:A:32:ILE:HA	1:A:43:ILE:O	2.19	0.41
1:B:132:LEU:HD13	1:B:320:PHE:CE1	2.56	0.41
1:B:165:LEU:N	1:B:165:LEU:HD12	2.35	0.41
1:C:120:PRO:HB2	1:C:126:VAL:HG11	2.02	0.41
1:D:149:TYR:CD1	1:D:391:MET:HE2	2.56	0.41
1:A:325:HIS:CE1	7:A:504:OLV:CAE	3.03	0.41
1:A:371:ASN:H	10:A:512:EDO:H22	1.85	0.41
1:B:30:VAL:HG21	1:B:114:PHE:CZ	2.55	0.41
1:C:122:THR:HA	1:C:243:ALA:O	2.21	0.41
1:A:132:LEU:HD22	1:A:320:PHE:CG	2.55	0.41
1:C:230:SER:O	1:C:231:GLN:HB2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:SER:HA	1:A:288:PRO:HD3	1.79	0.41
1:A:365:TYR:OH	7:A:504:OLV:OAC	2.38	0.41
1:A:376:GLN:HA	1:A:377:PRO:HD3	1.88	0.41
1:B:160:LEU:HD21	1:B:247:ILE:CD1	2.50	0.41
1:C:398:THR:HG22	1:C:426:TYR:HB2	2.01	0.41
1:D:57:VAL:CG1	1:D:92:ILE:HD11	2.47	0.41
1:B:258:ARG:HA	1:B:263:TYR:CE1	2.56	0.41
1:B:297:PHE:C	1:B:299:GLU:H	2.29	0.41
1:A:316:VAL:HB	1:A:354:PRO:HG3	2.03	0.41
1:A:370:SER:H	10:A:512:EDO:H22	1.82	0.41
1:B:36:ASP:HB3	1:B:41:ALA:HB3	2.03	0.41
1:B:338:TYR:CE1	1:C:345:CYS:HB2	2.55	0.41
1:C:297:PHE:CE2	1:C:298:MET:HG3	2.56	0.41
1:D:26:ALA:HA	1:D:110:THR:HB	2.02	0.41
1:A:33:THR:HA	1:A:194:PRO:HB3	2.03	0.41
1:A:103:TYR:CZ	1:A:114:PHE:HB2	2.54	0.41
2:L:2:FUC:H5	2:L:3:NAG:C2	2.50	0.40
1:A:399:HIS:CG	2:F:1:NAG:H5	2.56	0.40
1:B:296:HIS:HA	1:B:299:GLU:OE2	2.21	0.40
1:D:92:ILE:HG22	1:D:95:LEU:HD21	2.02	0.40
1:A:26:ALA:HA	1:A:110:THR:HB	2.03	0.40
1:A:68:ARG:NH1	8:A:508:SO4:O2	2.49	0.40
1:A:196:ILE:N	1:A:196:ILE:HD13	2.35	0.40
1:A:206:PHE:CZ	1:A:208:PRO:HG3	2.55	0.40
1:B:384:GLU:HG2	1:B:385:ALA:N	2.35	0.40
1:C:365:TYR:OH	7:C:512:OLV:H12	2.21	0.40
1:D:51:GLU:O	1:D:52:PRO:C	2.64	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	A	424/426 (100%)	401 (95%)	22 (5%)	1 (0%)	43	68
1	B	424/426 (100%)	387 (91%)	33 (8%)	4 (1%)	14	35
1	C	424/426 (100%)	399 (94%)	22 (5%)	3 (1%)	18	41
1	D	424/426 (100%)	396 (93%)	26 (6%)	2 (0%)	24	48
All	All	1696/1704 (100%)	1583 (93%)	103 (6%)	10 (1%)	21	44

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	64	ASN
1	A	175	ASP
1	B	175	ASP
1	C	175	ASP
1	D	63	LYS
1	C	164	ASP
1	D	175	ASP
1	B	155	LYS
1	B	364	ASN
1	C	364	ASN

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	375/375 (100%)	367 (98%)	8 (2%)	47	75
1	B	375/375 (100%)	363 (97%)	12 (3%)	34	64
1	C	375/375 (100%)	360 (96%)	15 (4%)	28	56
1	D	375/375 (100%)	365 (97%)	10 (3%)	39	69
All	All	1500/1500 (100%)	1455 (97%)	45 (3%)	36	66

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

Mol	Chain	Res	Type
1	A	19	ARG

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Mol	Chain	Res	Type
1	A	32	ILE
1	A	132	LEU
1	A	147	SER
1	A	154	LYS
1	A	176	ASN
1	A	383	ARG
1	A	417	SER
1	B	19	ARG
1	B	99	THR
1	B	104	GLU
1	B	122	THR
1	B	142	SER
1	B	176	ASN
1	B	303	MET
1	B	305	THR
1	B	336	ILE
1	B	370	SER
1	B	383	ARG
1	B	432	THR
1	C	49	MET
1	C	64	ASN
1	C	66	ARG
1	C	99	THR
1	C	132	LEU
1	C	155	LYS
1	C	162	VAL
1	C	176	ASN
1	C	228	GLU
1	C	294	ASN
1	C	352	SER
1	C	368	ILE
1	C	383	ARG
1	C	386	SER
1	C	432	THR
1	D	69	ILE
1	D	99	THR
1	D	132	LEU
1	D	176	ASN
1	D	211	ASN
1	D	349	LYS
1	D	359	ILE
1	D	378	GLU

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Mol	Chain	Res	Type
1	D	417	SER
1	D	432	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	HIS
1	A	193	GLN
1	A	231	GLN
1	A	295	HIS
1	A	405	ASN
1	A	424	HIS
1	B	295	HIS
1	B	376	GLN
1	B	401	HIS
1	C	193	GLN
1	C	202	HIS
1	C	211	ASN
1	C	246	HIS
1	C	294	ASN
1	C	325	HIS
1	D	335	ASN
1	D	401	HIS

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

26 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	E	1	2,1	14,14,15	0.63	0	17,19,21	1.04	1 (5%)
2	FUC	E	2	2	10,10,11	0.64	0	14,14,16	0.98	0
2	NAG	E	3	2	14,14,15	0.75	0	17,19,21	2.70	4 (23%)
2	NAG	F	1	2,1	14,14,15	0.67	0	17,19,21	1.25	1 (5%)
2	FUC	F	2	2	10,10,11	0.86	0	14,14,16	1.61	4 (28%)
2	NAG	F	3	2	14,14,15	0.56	0	17,19,21	1.29	2 (11%)
2	NAG	G	1	2,1	14,14,15	0.60	0	17,19,21	1.04	1 (5%)
2	FUC	G	2	2	10,10,11	0.57	0	14,14,16	1.06	0
2	NAG	G	3	2	14,14,15	0.74	0	17,19,21	1.55	3 (17%)
2	NAG	H	1	2,1	14,14,15	0.52	0	17,19,21	1.42	3 (17%)
2	FUC	H	2	2	10,10,11	0.52	0	14,14,16	1.05	0
2	NAG	H	3	2	14,14,15	0.51	0	17,19,21	0.96	0
3	NAG	I	1	3,1	14,14,15	0.54	0	17,19,21	0.94	0
3	FUC	I	2	3	10,10,11	0.65	0	14,14,16	0.69	0
2	NAG	J	1	2,1	14,14,15	0.52	0	17,19,21	0.90	0
2	FUC	J	2	2	10,10,11	0.58	0	14,14,16	0.79	0
2	NAG	J	3	2	14,14,15	0.80	0	17,19,21	1.87	3 (17%)
2	NAG	K	1	2,1	14,14,15	0.54	0	17,19,21	0.78	0
2	FUC	K	2	2	10,10,11	0.60	0	14,14,16	1.13	0
2	NAG	K	3	2	14,14,15	0.60	0	17,19,21	1.09	1 (5%)
2	NAG	L	1	2,1	14,14,15	0.60	0	17,19,21	0.85	0
2	FUC	L	2	2	10,10,11	0.61	0	14,14,16	1.19	1 (7%)
2	NAG	L	3	2	14,14,15	0.61	0	17,19,21	1.55	2 (11%)
2	NAG	M	1	2,1	14,14,15	0.61	0	17,19,21	0.93	2 (11%)
2	FUC	M	2	2	10,10,11	0.59	0	14,14,16	0.46	0
2	NAG	M	3	2	14,14,15	0.67	0	17,19,21	1.19	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
2	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	E	2	2	-	-	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	3	2	-	5/6/23/26	0/1/1/1
2	NAG	F	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	F	2	2	-	-	0/1/1/1
2	NAG	F	3	2	-	0/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	1/6/23/26	0/1/1/1
2	FUC	G	2	2	-	-	0/1/1/1
2	NAG	G	3	2	-	2/6/23/26	0/1/1/1
2	NAG	H	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	H	2	2	-	-	0/1/1/1
2	NAG	H	3	2	-	0/6/23/26	0/1/1/1
3	NAG	I	1	3,1	-	0/6/23/26	0/1/1/1
3	FUC	I	2	3	-	-	0/1/1/1
2	NAG	J	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	J	2	2	-	-	0/1/1/1
2	NAG	J	3	2	-	4/6/23/26	0/1/1/1
2	NAG	K	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	K	2	2	-	-	0/1/1/1
2	NAG	K	3	2	-	0/6/23/26	0/1/1/1
2	NAG	L	1	2,1	-	0/6/23/26	0/1/1/1
2	FUC	L	2	2	-	-	0/1/1/1
2	NAG	L	3	2	-	5/6/23/26	0/1/1/1
2	NAG	M	1	2,1	-	1/6/23/26	0/1/1/1
2	FUC	M	2	2	-	-	0/1/1/1
2	NAG	M	3	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	3	NAG	C1-O5-C5	6.97	121.53	112.19
2	E	3	NAG	C2-N2-C7	-6.61	114.04	122.90
2	J	3	NAG	C2-N2-C7	-4.63	116.70	122.90
2	J	3	NAG	C1-O5-C5	4.49	118.20	112.19
2	L	3	NAG	C2-N2-C7	3.99	128.25	122.90
2	F	2	FUC	C2-C3-C4	3.67	117.32	110.86
2	M	3	NAG	C2-N2-C7	-3.21	118.60	122.90
2	E	3	NAG	O3-C3-C2	-3.19	102.78	109.40
2	L	3	NAG	O5-C1-C2	-3.10	106.50	111.29
2	G	3	NAG	O4-C4-C5	3.09	116.92	109.32
2	G	3	NAG	C2-N2-C7	-3.08	118.78	122.90
2	H	1	NAG	C1-O5-C5	2.99	116.19	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	2	FUC	C3-C4-C5	2.91	114.23	109.81
2	L	2	FUC	C3-C4-C5	2.81	114.09	109.81
2	K	3	NAG	C4-C3-C2	2.74	115.03	111.02
2	E	3	NAG	O3-C3-C4	-2.73	103.94	110.38
2	F	3	NAG	O5-C5-C6	2.67	112.87	107.66
2	E	1	NAG	C1-O5-C5	2.61	115.68	112.19
2	H	1	NAG	C2-N2-C7	-2.52	119.53	122.90
2	G	3	NAG	C3-C4-C5	-2.49	105.72	110.23
2	M	1	NAG	C1-O5-C5	2.37	115.36	112.19
2	G	1	NAG	C4-C3-C2	-2.34	107.59	111.02
2	J	3	NAG	C3-C4-C5	-2.33	106.02	110.23
2	H	1	NAG	C4-C3-C2	-2.27	107.69	111.02
2	M	1	NAG	C2-N2-C7	-2.20	119.95	122.90
2	F	3	NAG	O7-C7-C8	-2.19	118.16	122.05
2	F	2	FUC	C1-C2-C3	2.16	112.80	109.64
2	F	2	FUC	C1-O5-C5	-2.11	107.98	112.97
2	F	1	NAG	O3-C3-C4	-2.07	105.50	110.38

There are no chirality outliers.

All (18) torsion outliers are listed below:

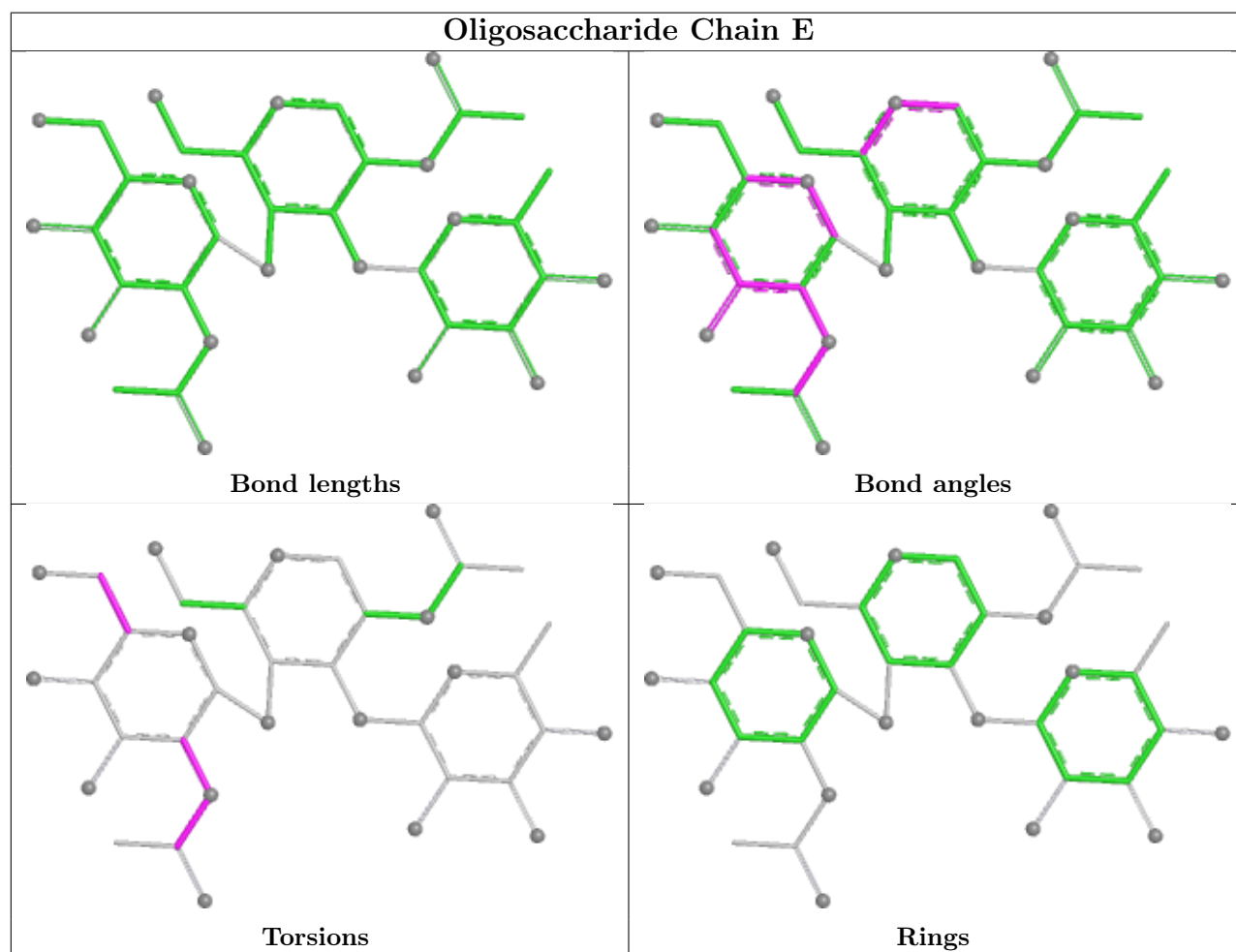
Mol	Chain	Res	Type	Atoms
2	E	3	NAG	C8-C7-N2-C2
2	E	3	NAG	O7-C7-N2-C2
2	J	3	NAG	C1-C2-N2-C7
2	J	3	NAG	C8-C7-N2-C2
2	J	3	NAG	O7-C7-N2-C2
2	G	3	NAG	O5-C5-C6-O6
2	G	3	NAG	C4-C5-C6-O6
2	E	3	NAG	O5-C5-C6-O6
2	L	3	NAG	O5-C5-C6-O6
2	E	3	NAG	C4-C5-C6-O6
2	L	3	NAG	C4-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	J	3	NAG	O5-C5-C6-O6
2	L	3	NAG	C3-C2-N2-C7
2	L	3	NAG	C8-C7-N2-C2
2	L	3	NAG	O7-C7-N2-C2
2	E	3	NAG	C1-C2-N2-C7
2	M	1	NAG	C1-C2-N2-C7

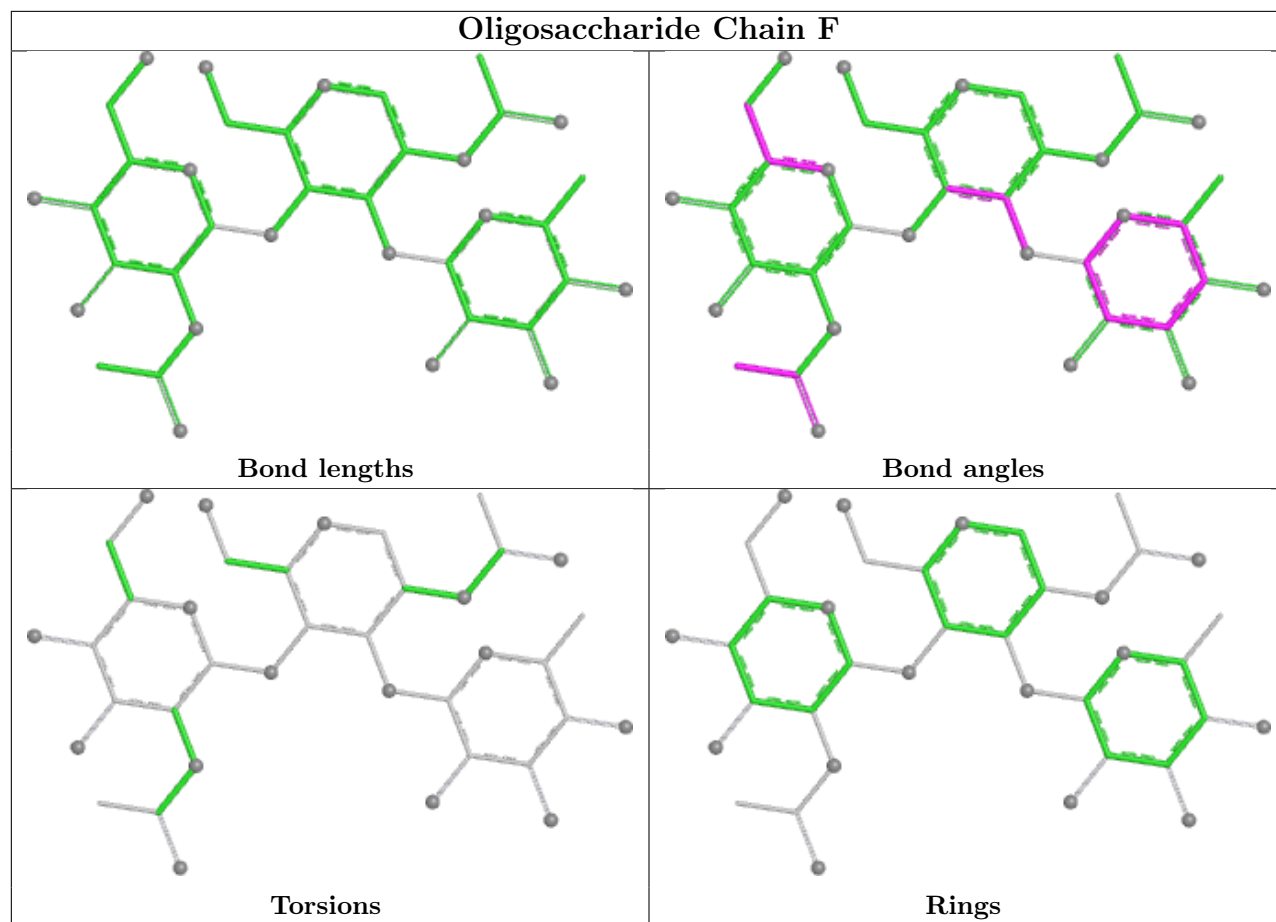
There are no ring outliers.

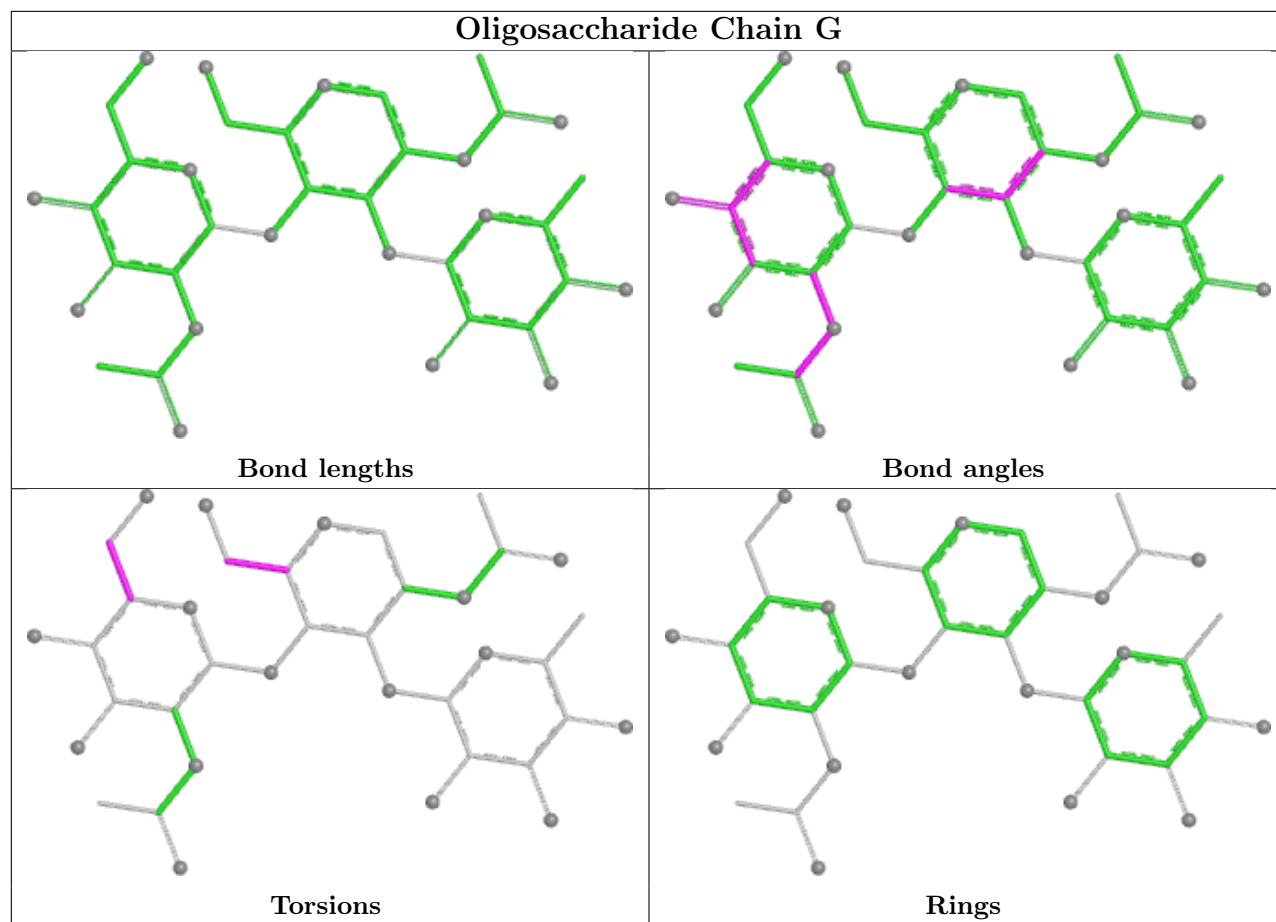
12 monomers are involved in 12 short contacts:

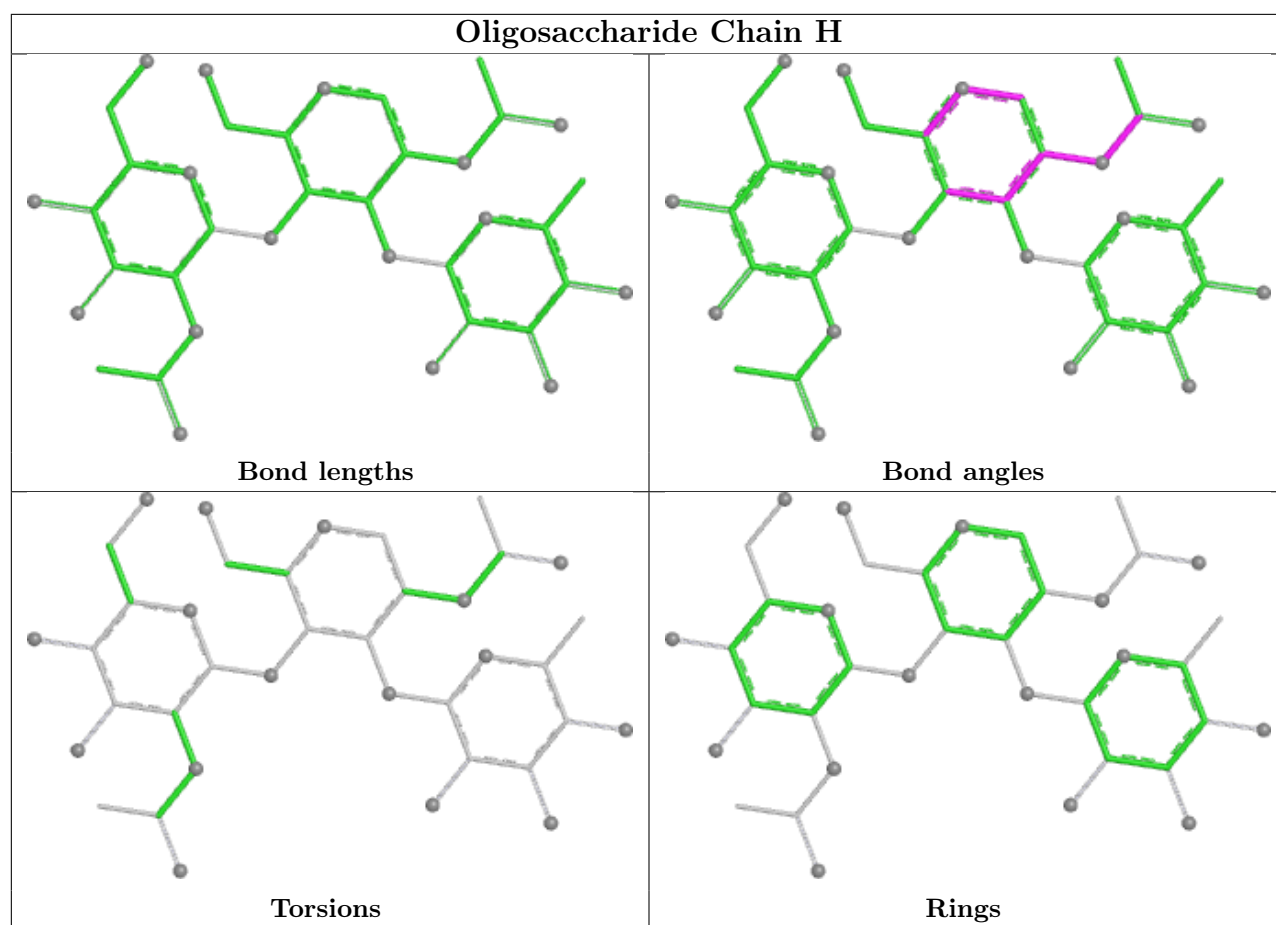
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	F	1	NAG	1	0
2	H	1	NAG	1	0
2	E	3	NAG	2	0
2	J	2	FUC	1	0
2	J	1	NAG	2	0
2	J	3	NAG	4	0
2	H	2	FUC	1	0
2	G	1	NAG	1	0
2	E	2	FUC	2	0
2	L	3	NAG	2	0
2	G	3	NAG	1	0
2	L	2	FUC	2	0

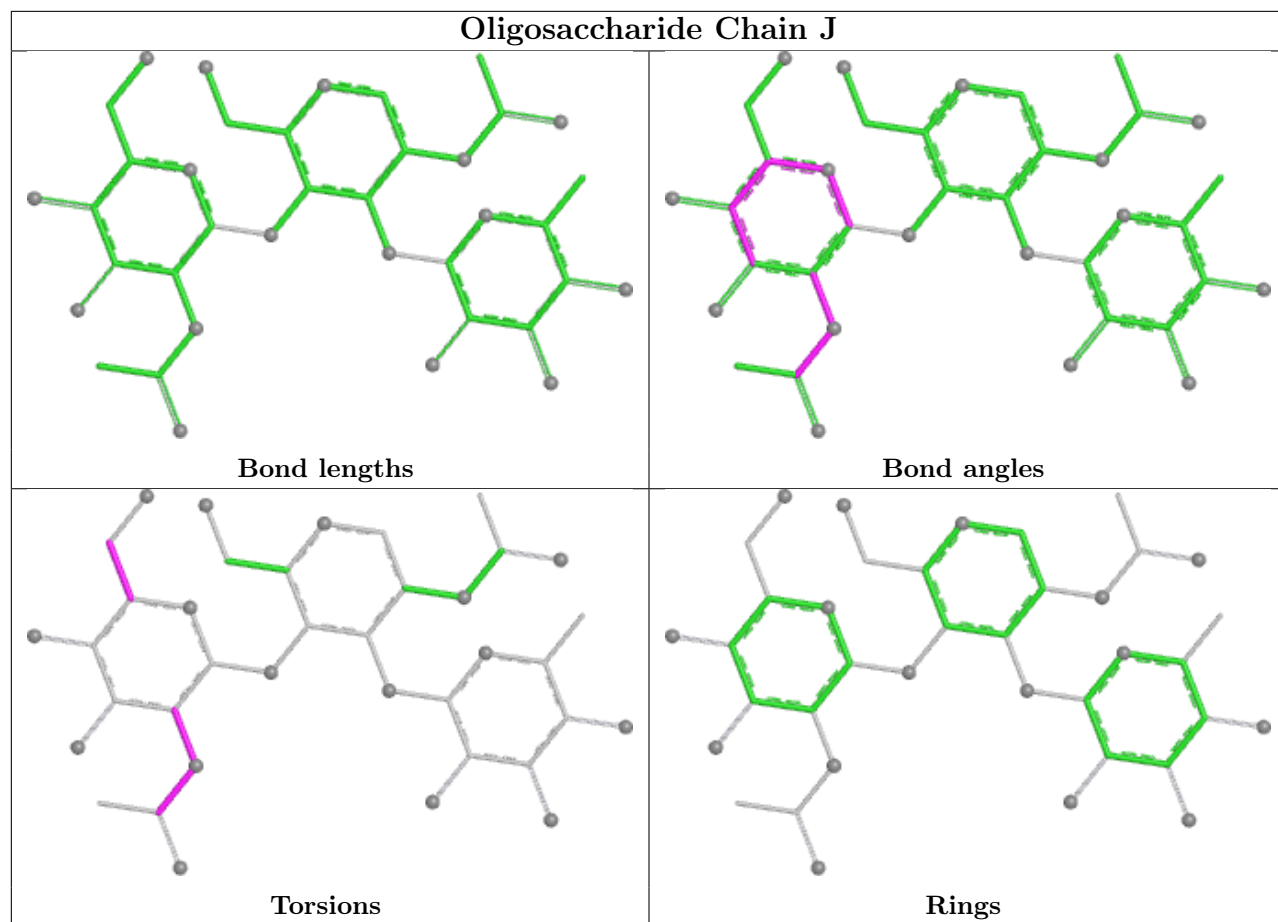
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

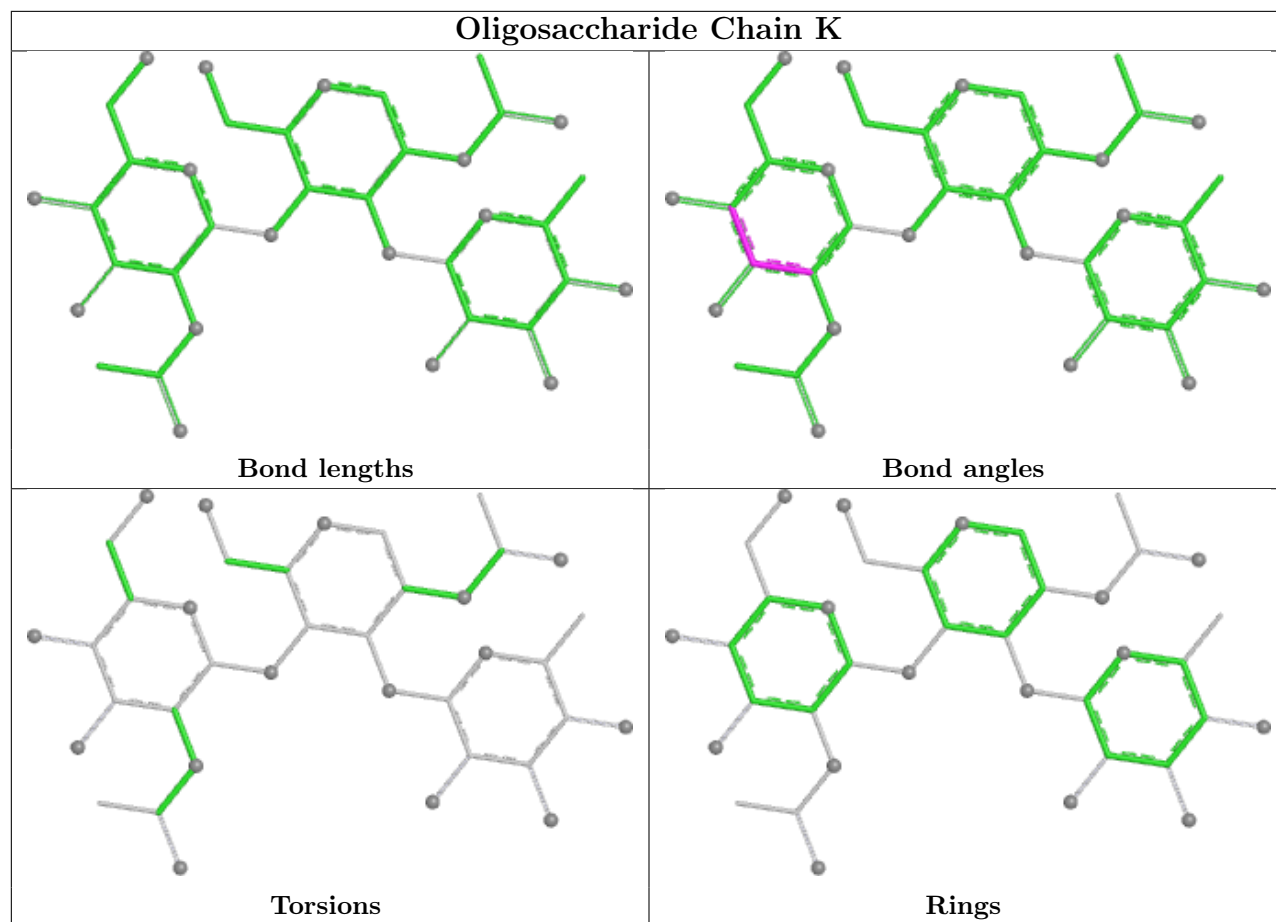


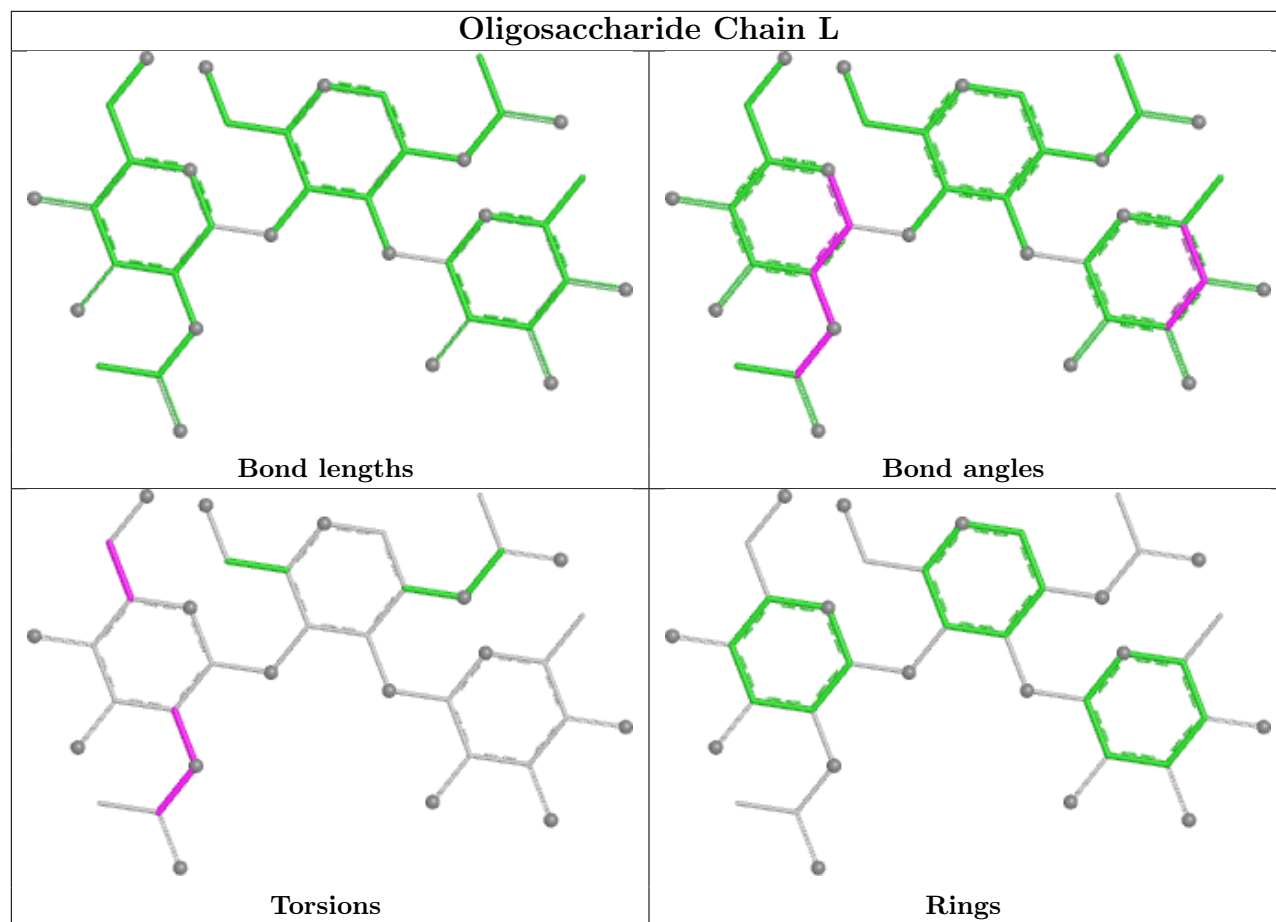


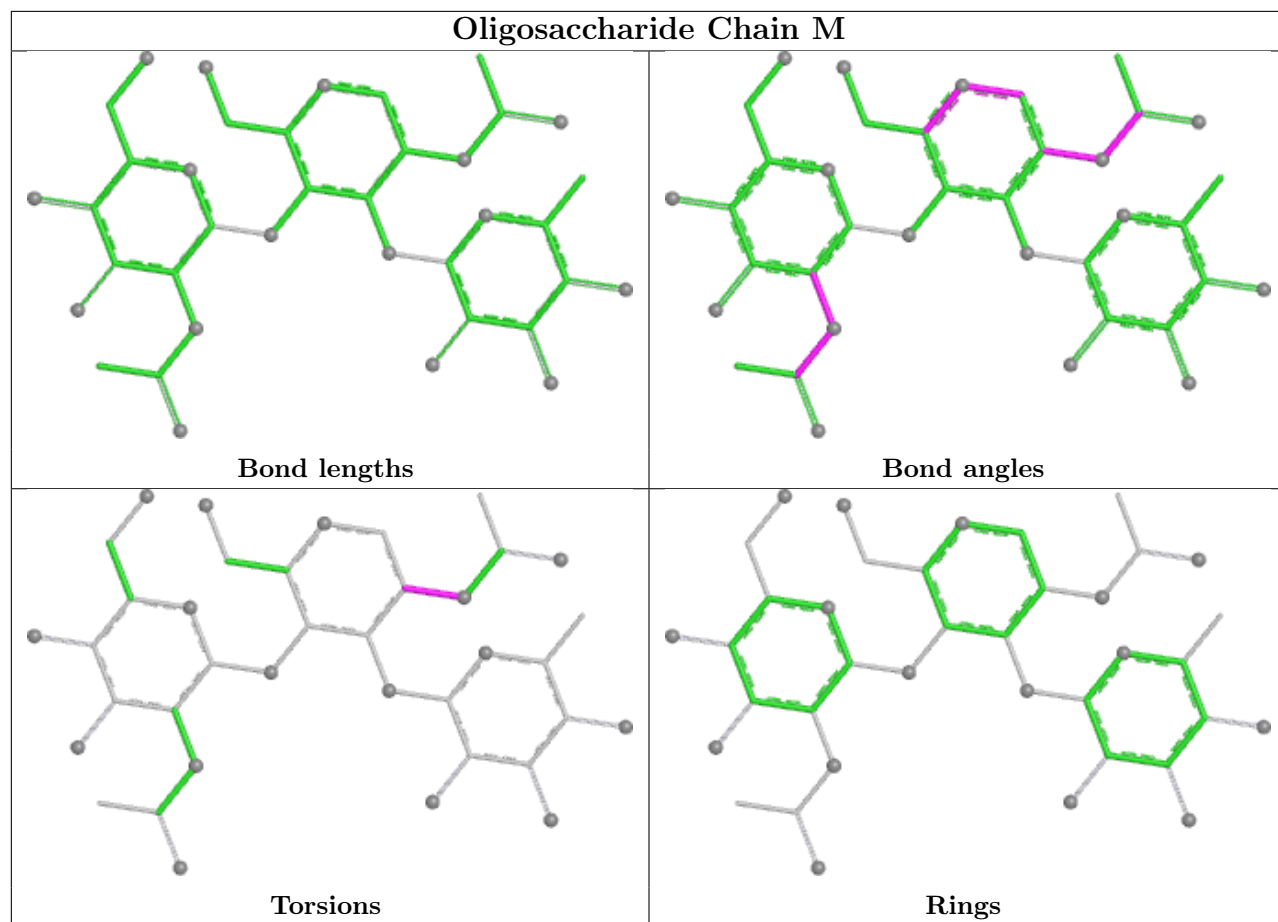


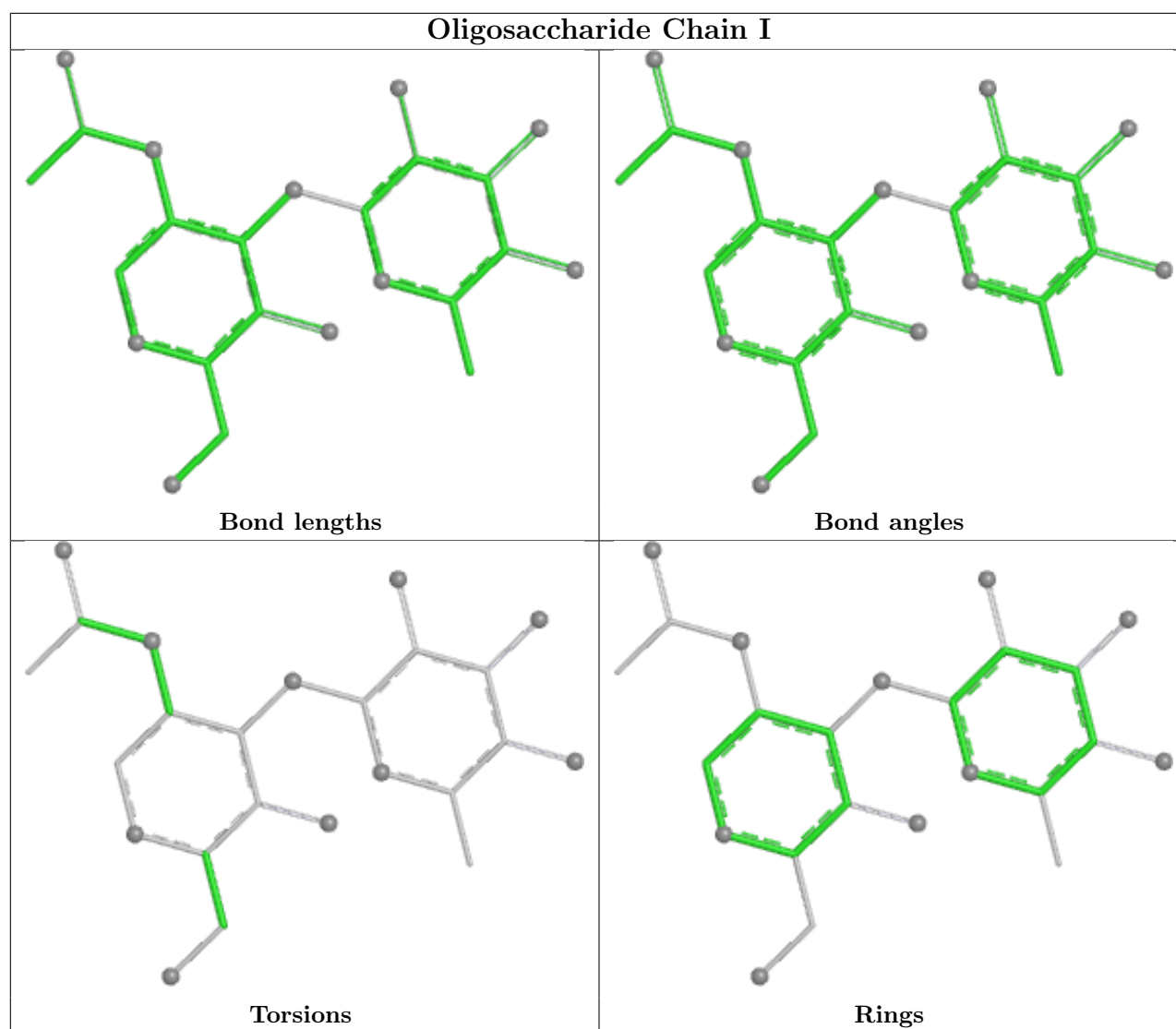












5.6 Ligand geometry [i](#)

Of 40 ligands modelled in this entry, 8 are monoatomic - leaving 32 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	$\# Z > 2$
11	NAG	C	507	1	14,14,15	0.52	0	17,19,21	1.68	2 (11%)
7	OLV	D	511	-	14,14,14	3.20	6 (42%)	21,21,21	2.34	5 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
8	SO4	B	508	-	4,4,4	0.30	0	6,6,6	0.25	0
11	NAG	B	503	1	14,14,15	0.54	0	17,19,21	1.15	1 (5%)
11	NAG	B	511	1	14,14,15	0.76	0	17,19,21	1.42	4 (23%)
8	SO4	A	508	-	4,4,4	0.25	0	6,6,6	0.13	0
10	EDO	A	512	-	3,3,3	0.50	0	2,2,2	0.30	0
8	SO4	A	506	-	4,4,4	0.43	0	6,6,6	0.52	0
8	SO4	A	510	-	4,4,4	0.43	0	6,6,6	0.12	0
8	SO4	D	513	-	4,4,4	0.38	0	6,6,6	0.18	0
7	OLV	B	504	-	14,14,14	3.06	5 (35%)	21,21,21	2.73	6 (28%)
8	SO4	C	515	-	4,4,4	0.30	0	6,6,6	0.29	0
8	SO4	A	507	-	4,4,4	0.30	0	6,6,6	0.15	0
8	SO4	C	513	-	4,4,4	0.36	0	6,6,6	0.26	0
8	SO4	C	514	-	4,4,4	0.31	0	6,6,6	0.23	0
8	SO4	D	515	-	4,4,4	0.25	0	6,6,6	0.28	0
7	OLV	C	512	-	14,14,14	3.31	6 (42%)	21,21,21	2.91	8 (38%)
8	SO4	A	509	-	4,4,4	0.34	0	6,6,6	0.17	0
8	SO4	A	505	-	4,4,4	0.34	0	6,6,6	0.24	0
6	GOL	D	510	-	5,5,5	0.34	0	5,5,5	0.44	0
6	GOL	A	503	-	5,5,5	0.38	0	5,5,5	0.26	0
8	SO4	D	512	-	4,4,4	0.33	0	6,6,6	0.16	0
9	ACT	A	511	-	3,3,3	0.95	0	3,3,3	1.33	0
8	SO4	B	510	-	4,4,4	0.34	0	6,6,6	0.19	0
11	NAG	A	516	1	14,14,15	0.62	0	17,19,21	1.27	2 (11%)
8	SO4	B	509	-	4,4,4	0.34	0	6,6,6	0.26	0
8	SO4	D	514	-	4,4,4	0.35	0	6,6,6	0.42	0
6	GOL	C	508	-	5,5,5	0.32	0	5,5,5	0.46	0
11	NAG	D	504	1	14,14,15	0.57	0	17,19,21	1.43	4 (23%)
11	NAG	D	516	1	14,14,15	0.55	0	17,19,21	1.08	1 (5%)
11	NAG	C	501	1	14,14,15	0.51	0	17,19,21	0.81	0
7	OLV	A	504	-	14,14,14	2.98	5 (35%)	21,21,21	2.77	9 (42%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	GOL	C	508	-	-	2/4/4/4	-
6	GOL	D	510	-	-	3/4/4/4	-
6	GOL	A	503	-	-	0/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	NAG	C	501	1	-	0/6/23/26	0/1/1/1
11	NAG	C	507	1	-	2/6/23/26	0/1/1/1
11	NAG	A	516	1	-	2/6/23/26	0/1/1/1
7	OLV	B	504	-	-	0/2/12/12	0/2/2/2
11	NAG	D	504	1	-	1/6/23/26	0/1/1/1
7	OLV	D	511	-	-	0/2/12/12	0/2/2/2
11	NAG	B	503	1	-	2/6/23/26	0/1/1/1
11	NAG	D	516	1	-	1/6/23/26	0/1/1/1
7	OLV	C	512	-	-	2/2/12/12	0/2/2/2
11	NAG	B	511	1	-	3/6/23/26	0/1/1/1
10	EDO	A	512	-	-	0/1/1/1	-
7	OLV	A	504	-	-	0/2/12/12	0/2/2/2

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	511	OLV	OAI-CAM	-7.33	1.41	1.48
7	C	512	OLV	OAI-CAM	-7.05	1.41	1.48
7	A	504	OLV	OAI-CAM	-6.50	1.42	1.48
7	B	504	OLV	OAI-CAM	-6.44	1.42	1.48
7	C	512	OLV	CAH-CAM	-5.78	1.47	1.54
7	B	504	OLV	CAH-CAM	-5.52	1.47	1.54
7	D	511	OLV	CAH-CAK	-5.52	1.44	1.50
7	C	512	OLV	CAH-CAK	-5.31	1.44	1.50
7	D	511	OLV	CAH-CAM	-5.25	1.48	1.54
7	A	504	OLV	CAH-CAM	-5.09	1.48	1.54
7	B	504	OLV	CAH-CAK	-4.90	1.44	1.50
7	A	504	OLV	CAH-CAK	-4.83	1.45	1.50
7	C	512	OLV	CAF-CAK	3.82	1.45	1.39
7	C	512	OLV	OAI-CAL	3.59	1.44	1.37
7	B	504	OLV	OAI-CAL	3.46	1.43	1.37
7	A	504	OLV	CAF-CAK	3.45	1.45	1.39
7	B	504	OLV	CAF-CAK	3.36	1.45	1.39
7	D	511	OLV	CAF-CAK	3.16	1.44	1.39
7	D	511	OLV	OAI-CAL	3.11	1.43	1.37
7	A	504	OLV	OAI-CAL	3.00	1.43	1.37
7	C	512	OLV	CAE-CAJ	2.31	1.43	1.39
7	D	511	OLV	CAE-CAJ	2.26	1.43	1.39

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	A	504	0LV	CAM-CAH-CAK	6.81	107.03	103.07
7	C	512	0LV	OAI-CAL-CAJ	6.56	131.67	123.32
7	C	512	0LV	OAI-CAL-CAK	-6.45	108.88	113.82
7	B	504	0LV	OAI-CAL-CAJ	6.05	131.02	123.32
7	B	504	0LV	OAI-CAL-CAK	-5.92	109.29	113.82
7	D	511	0LV	CAM-CAH-CAK	5.63	106.34	103.07
7	C	512	0LV	OAI-CAM-CAA	5.46	111.38	106.53
7	B	504	0LV	CAM-CAH-CAK	5.41	106.21	103.07
7	A	504	0LV	OAI-CAL-CAJ	5.28	130.05	123.32
7	D	511	0LV	OAI-CAL-CAJ	5.05	129.75	123.32
7	B	504	0LV	OAI-CAM-CAA	5.01	110.98	106.53
7	D	511	0LV	OAI-CAL-CAK	-4.35	110.49	113.82
7	C	512	0LV	CAM-OAI-CAL	-4.25	104.64	107.16
11	C	507	NAG	C1-O5-C5	4.24	117.87	112.19
7	A	504	0LV	OAI-CAL-CAK	-3.78	110.93	113.82
7	A	504	0LV	OAI-CAM-CAB	3.76	109.88	106.53
7	A	504	0LV	OAI-CAM-CAA	3.67	109.79	106.53
7	A	504	0LV	CAK-CAL-CAJ	-3.62	119.02	122.92
7	C	512	0LV	OAI-CAM-CAB	3.52	109.66	106.53
11	D	504	NAG	C1-O5-C5	3.30	116.60	112.19
7	C	512	0LV	CAK-CAL-CAJ	-3.22	119.44	122.92
7	B	504	0LV	CAK-CAL-CAJ	-2.99	119.69	122.92
11	C	507	NAG	C4-C3-C2	-2.99	106.63	111.02
11	A	516	NAG	C2-N2-C7	-2.98	118.91	122.90
7	D	511	0LV	OAI-CAM-CAB	2.95	109.15	106.53
7	D	511	0LV	CAK-CAL-CAJ	-2.93	119.76	122.92
7	A	504	0LV	CAB-CAM-CAH	-2.67	108.59	112.67
7	B	504	0LV	OAI-CAM-CAB	2.60	108.84	106.53
11	B	511	NAG	C6-C5-C4	-2.60	106.64	113.02
11	D	516	NAG	C2-N2-C7	-2.58	119.44	122.90
7	A	504	0LV	CAE-CAJ-CAL	2.57	121.65	117.88
7	C	512	0LV	OAC-CAG-CAJ	2.31	118.38	111.94
11	B	511	NAG	O3-C3-C4	-2.25	105.07	110.38
11	D	504	NAG	C4-C3-C2	-2.22	107.76	111.02
11	B	511	NAG	O7-C7-C8	-2.21	118.11	122.05
11	D	504	NAG	O3-C3-C2	2.15	113.86	109.40
7	C	512	0LV	OAI-CAM-CAH	-2.09	103.48	105.15
11	B	511	NAG	O5-C5-C4	-2.06	105.81	110.83
7	A	504	0LV	CAG-CAJ-CAL	-2.06	115.45	119.62
11	B	503	NAG	C1-O5-C5	2.06	114.94	112.19
11	D	504	NAG	C1-C2-N2	-2.05	107.21	110.43
11	A	516	NAG	C1-C2-N2	-2.00	107.28	110.43

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	D	510	GOL	O1-C1-C2-C3
11	C	507	NAG	C8-C7-N2-C2
11	B	503	NAG	C8-C7-N2-C2
11	B	503	NAG	O7-C7-N2-C2
11	C	507	NAG	O7-C7-N2-C2
11	B	511	NAG	C8-C7-N2-C2
11	B	511	NAG	O7-C7-N2-C2
6	D	510	GOL	O1-C1-C2-O2
6	C	508	GOL	O1-C1-C2-C3
11	A	516	NAG	O5-C5-C6-O6
11	A	516	NAG	C4-C5-C6-O6
11	D	504	NAG	O5-C5-C6-O6
6	C	508	GOL	O1-C1-C2-O2
6	D	510	GOL	O2-C2-C3-O3
7	C	512	0LV	OAC-CAG-CAJ-CAE
11	B	511	NAG	O5-C5-C6-O6
11	D	516	NAG	C4-C5-C6-O6
7	C	512	0LV	OAC-CAG-CAJ-CAL

There are no ring outliers.

10 monomers are involved in 57 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	D	511	0LV	13	0
8	A	508	SO4	1	0
10	A	512	EDO	6	0
7	B	504	0LV	10	0
7	C	512	0LV	13	0
8	A	505	SO4	1	0
6	A	503	GOL	2	0
8	B	509	SO4	1	0
6	C	508	GOL	1	0
7	A	504	0LV	9	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	426/426 (100%)	-0.83	1 (0%) 91 90	29, 37, 51, 78	0
1	B	426/426 (100%)	-0.67	2 (0%) 87 86	30, 44, 59, 88	0
1	C	426/426 (100%)	-0.74	0 100 100	31, 41, 58, 89	1 (0%)
1	D	426/426 (100%)	-0.78	0 100 100	28, 40, 58, 85	0
All	All	1704/1704 (100%)	-0.76	3 (0%) 91 90	28, 40, 58, 89	1 (0%)

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	65	GLY	2.4
1	B	64	ASN	2.3
1	A	432	THR	2.1

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
2	NAG	J	3	14/15	0.73	0.12	58,64,67,69	0
2	NAG	L	3	14/15	0.73	0.14	80,90,96,98	0
2	FUC	K	2	10/11	0.76	0.12	61,70,74,78	0
2	FUC	E	2	10/11	0.76	0.12	56,67,73,78	0

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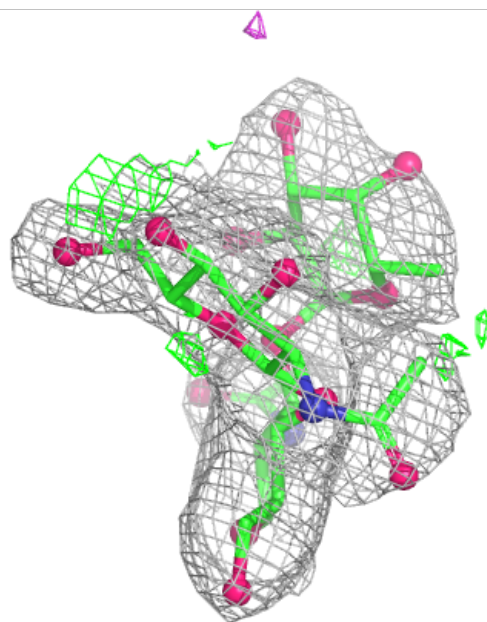
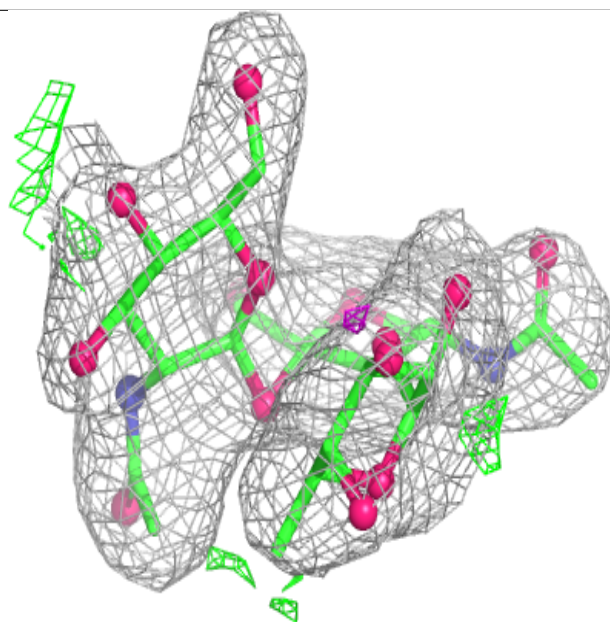
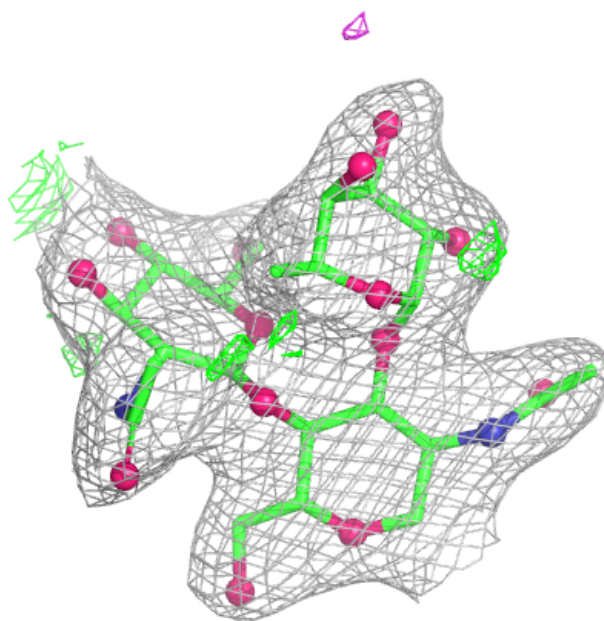
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	FUC	G	2	10/11	0.78	0.11	58,62,69,69	0
3	FUC	I	2	10/11	0.79	0.11	69,75,76,78	0
2	NAG	E	3	14/15	0.80	0.10	53,61,68,68	0
2	NAG	G	3	14/15	0.81	0.11	51,59,64,65	0
2	FUC	J	2	10/11	0.84	0.11	68,70,79,81	0
2	FUC	L	2	10/11	0.84	0.10	70,79,86,88	0
2	FUC	M	2	10/11	0.85	0.10	57,66,71,71	0
3	NAG	I	1	14/15	0.85	0.09	52,64,69,69	0
2	NAG	K	3	14/15	0.85	0.09	50,68,78,81	0
2	NAG	M	3	14/15	0.86	0.09	54,65,74,75	0
2	FUC	F	2	10/11	0.89	0.09	35,46,52,57	0
2	NAG	L	1	14/15	0.90	0.09	50,70,79,79	0
2	NAG	J	1	14/15	0.91	0.08	48,59,64,68	0
2	NAG	H	3	14/15	0.91	0.07	46,53,55,56	0
2	NAG	F	3	14/15	0.92	0.07	32,40,43,47	0
2	FUC	H	2	10/11	0.92	0.08	49,56,58,59	0
2	NAG	E	1	14/15	0.92	0.08	47,53,62,62	0
2	NAG	G	1	14/15	0.94	0.07	44,53,58,58	0
2	NAG	K	1	14/15	0.95	0.07	47,53,62,66	0
2	NAG	H	1	14/15	0.96	0.07	38,46,52,58	0
2	NAG	F	1	14/15	0.97	0.05	34,37,44,47	0
2	NAG	M	1	14/15	0.97	0.06	49,53,63,68	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.

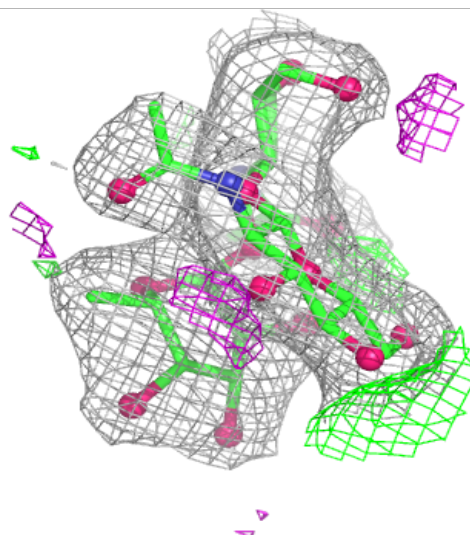
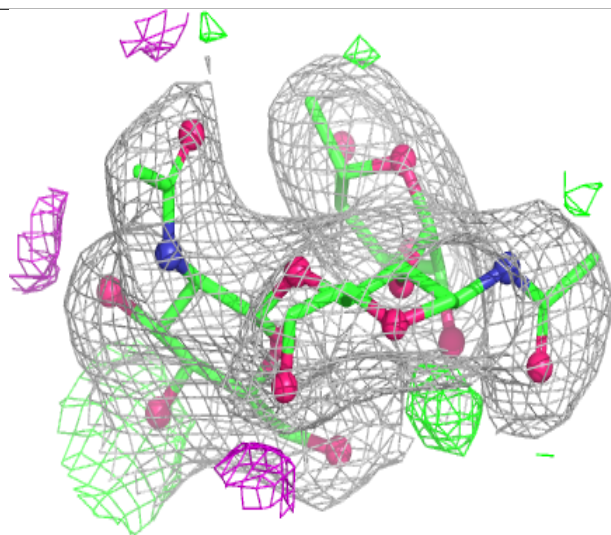
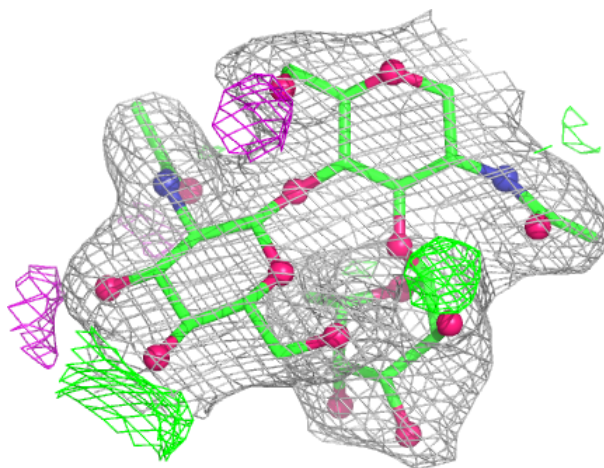
Electron density around Chain E:

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



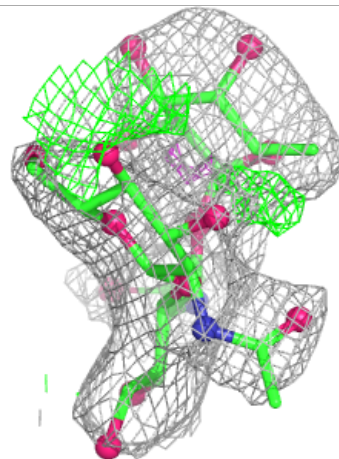
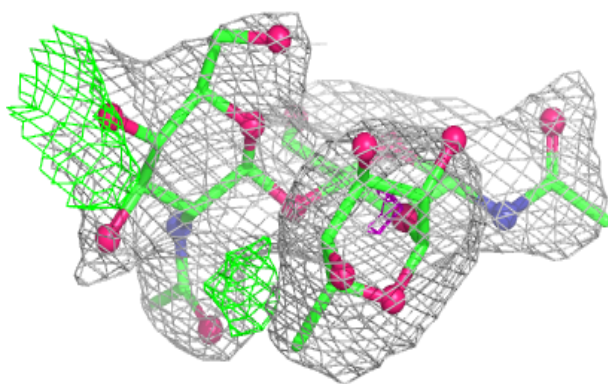
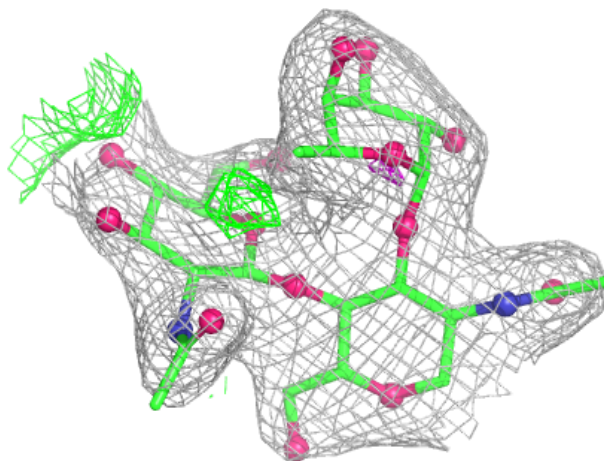
Electron density around Chain F:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



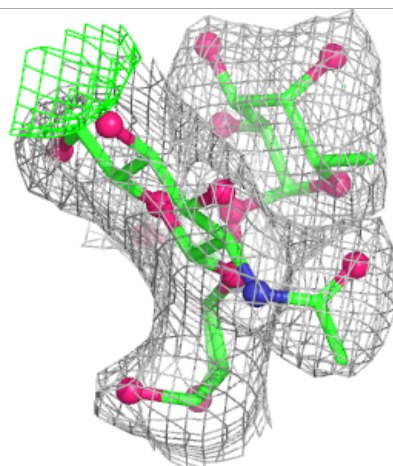
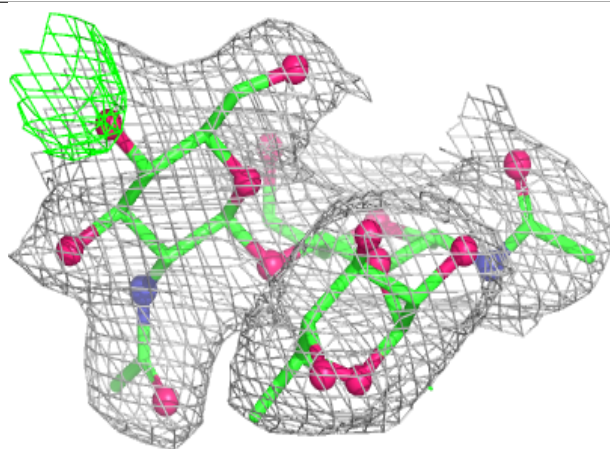
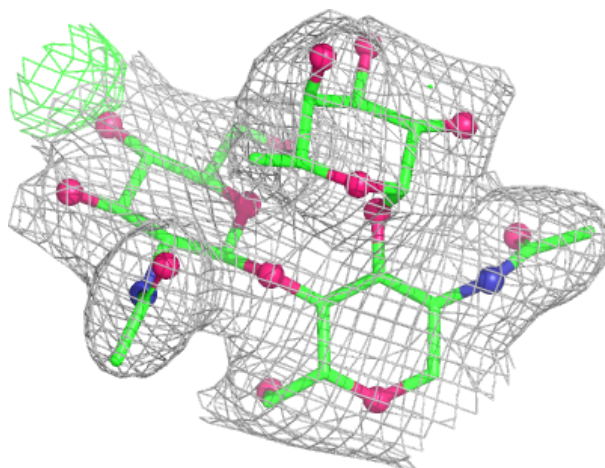
Electron density around Chain G:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



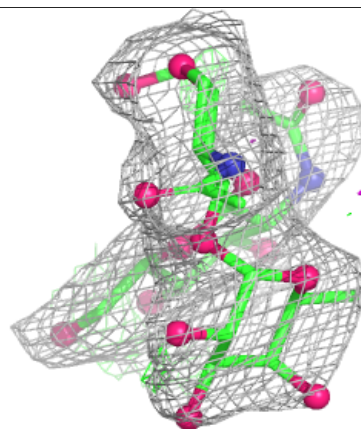
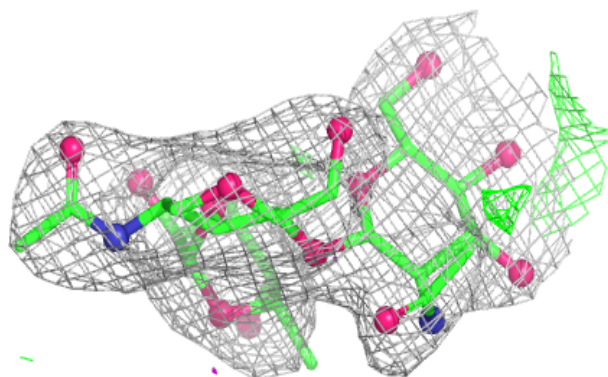
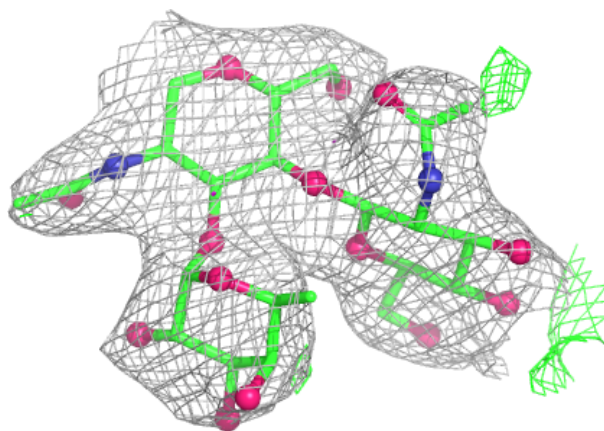
Electron density around Chain H:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



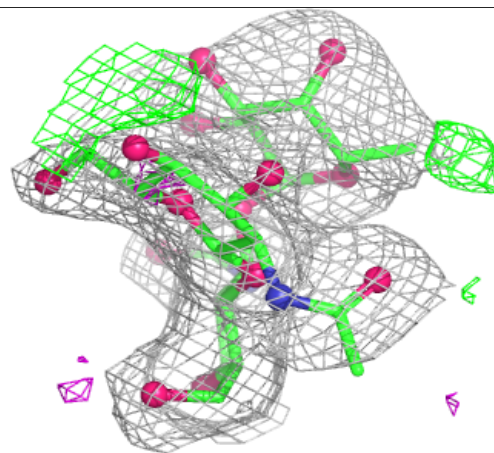
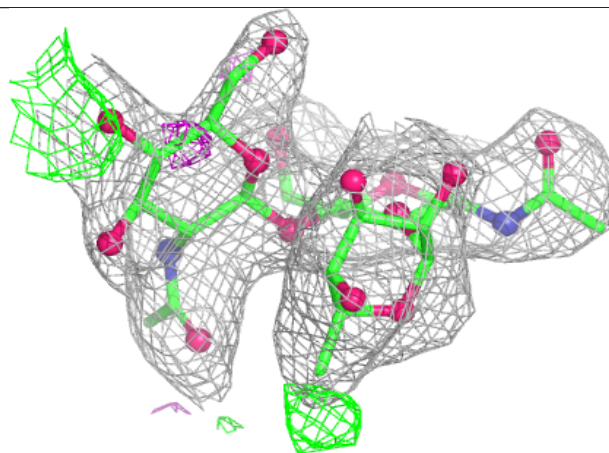
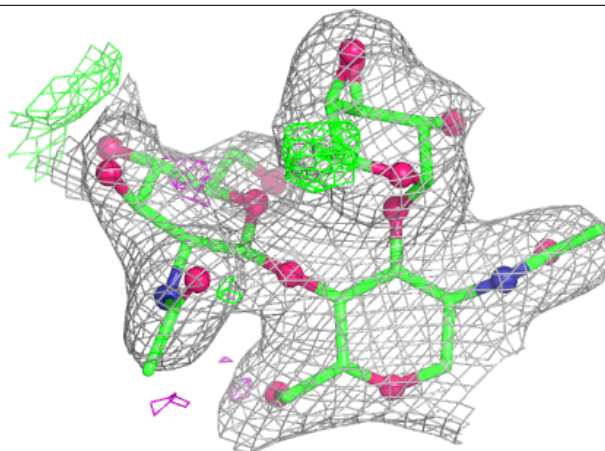
Electron density around Chain J:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



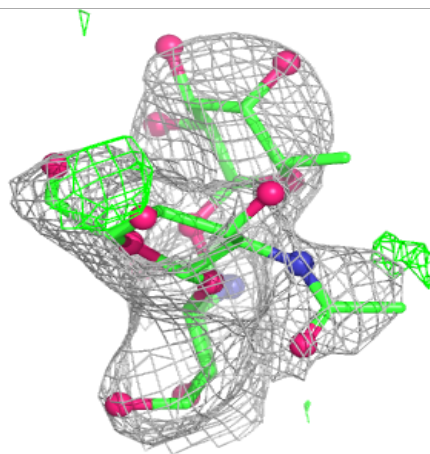
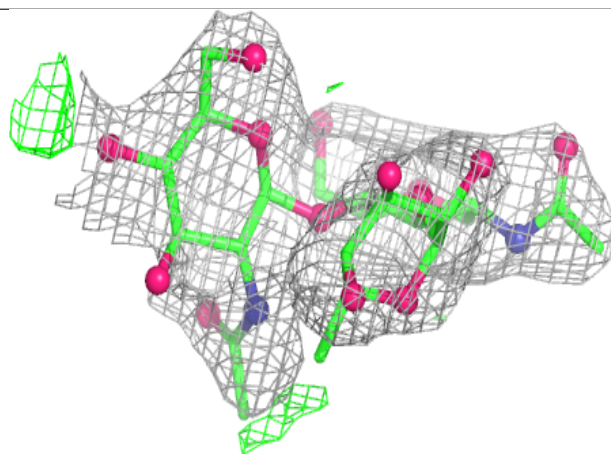
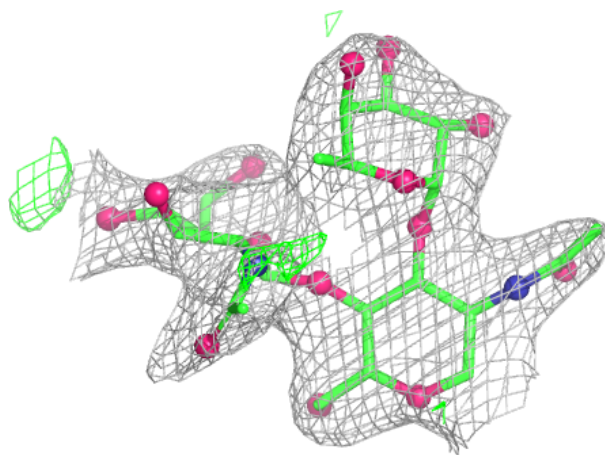
Electron density around Chain K:

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and green (positive)



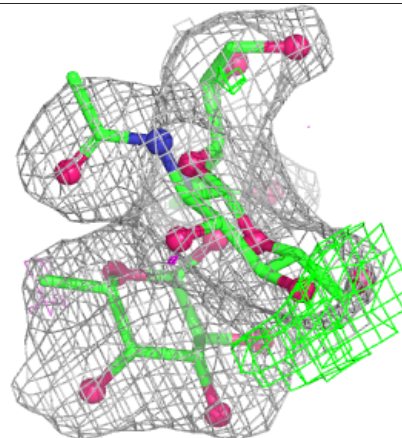
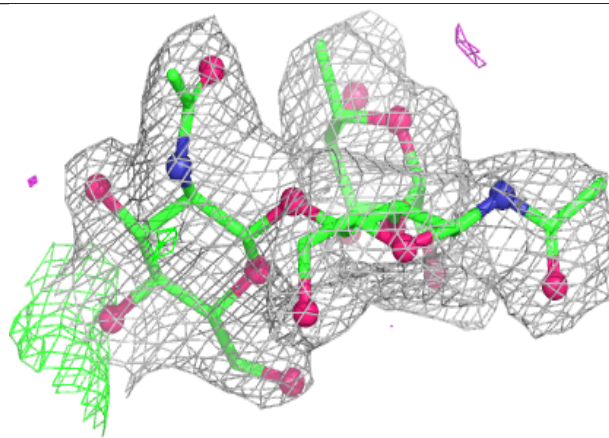
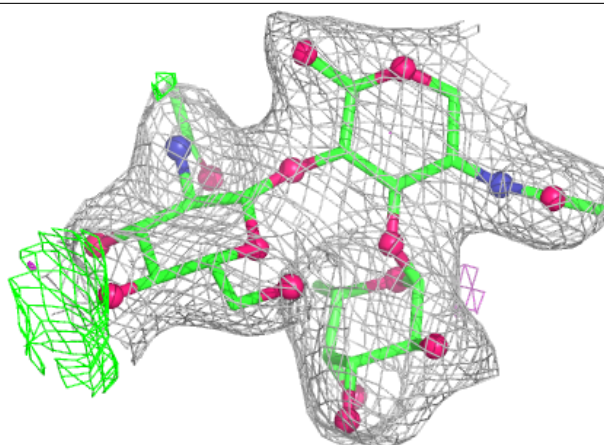
Electron density around Chain L:

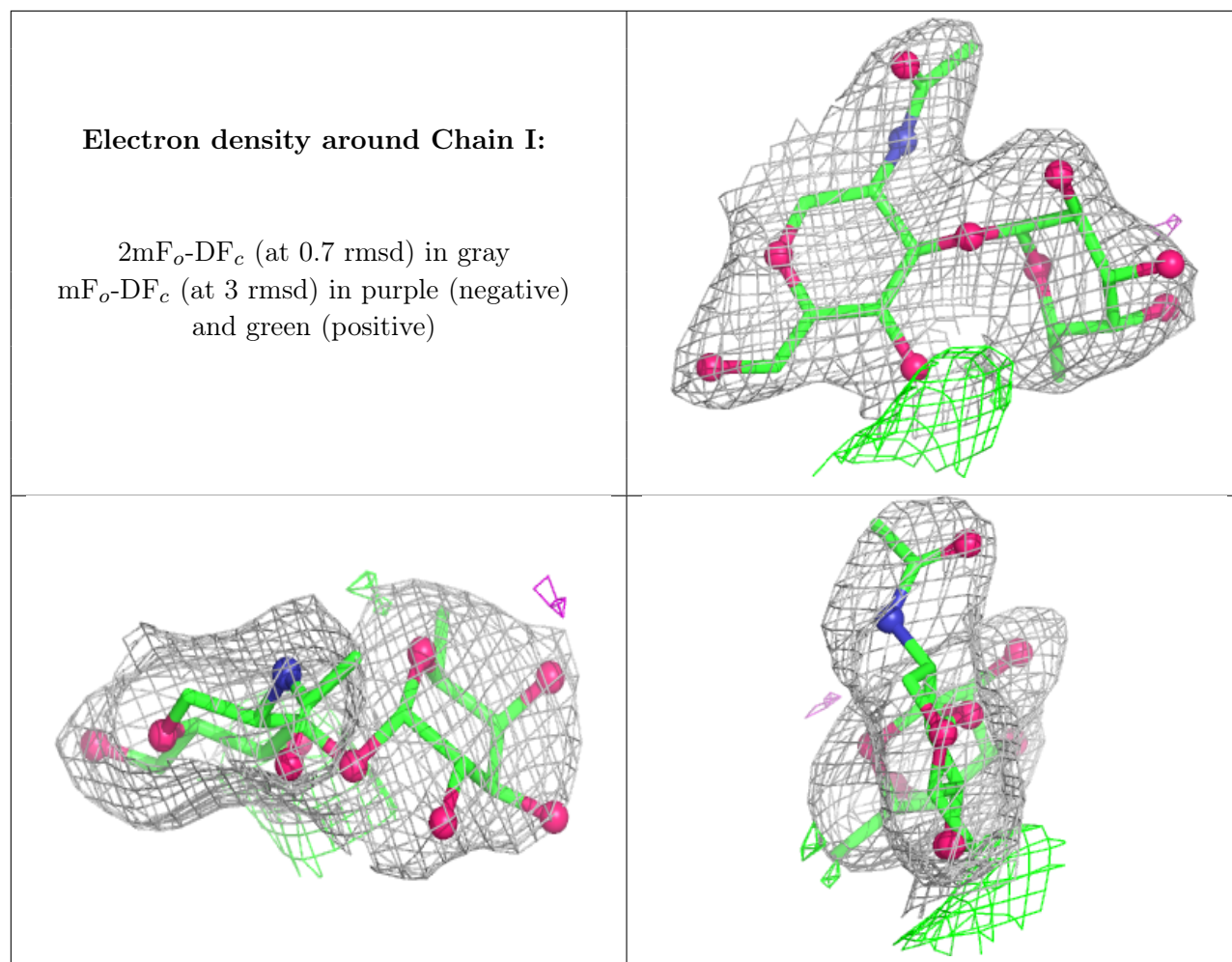
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain M:

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





6.4 Ligands ⓘ

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
7	OLV	A	504	13/13	0.70	0.24	35,56,71,72	0
7	OLV	B	504	13/13	0.71	0.21	46,62,75,76	0
7	OLV	C	512	13/13	0.73	0.19	48,64,72,75	0
7	OLV	D	511	13/13	0.75	0.20	38,60,72,74	0
11	NAG	C	507	14/15	0.76	0.11	57,64,68,69	0
9	ACT	A	511	4/4	0.81	0.17	41,46,47,52	0
8	SO4	D	515	5/5	0.81	0.16	67,72,88,96	0
8	SO4	A	505	5/5	0.83	0.31	59,66,86,90	0
8	SO4	D	514	5/5	0.85	0.33	57,62,62,77	0
6	GOL	D	510	6/6	0.85	0.12	53,55,61,66	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	SO4	A	509	5/5	0.86	0.31	57,63,84,90	0
8	SO4	B	509	5/5	0.86	0.20	63,67,83,95	0
11	NAG	B	503	14/15	0.86	0.09	54,60,64,65	0
8	SO4	A	506	5/5	0.86	0.21	51,53,69,90	0
8	SO4	B	510	5/5	0.88	0.14	66,70,93,98	0
8	SO4	A	507	5/5	0.88	0.25	60,69,78,86	0
11	NAG	D	504	14/15	0.88	0.09	57,66,69,74	0
8	SO4	B	508	5/5	0.89	0.24	67,74,86,93	0
8	SO4	C	514	5/5	0.89	0.22	61,72,78,90	0
8	SO4	C	515	5/5	0.89	0.27	71,76,79,97	0
8	SO4	D	512	5/5	0.90	0.25	59,62,75,83	0
10	EDO	A	512	4/4	0.91	0.08	44,45,47,47	0
8	SO4	A	510	5/5	0.91	0.12	54,58,76,91	0
11	NAG	D	516	14/15	0.91	0.07	40,48,54,58	0
11	NAG	B	511	14/15	0.92	0.07	47,51,55,58	0
8	SO4	D	513	5/5	0.92	0.23	52,52,58,77	0
8	SO4	C	513	5/5	0.92	0.21	52,63,81,84	0
6	GOL	A	503	6/6	0.92	0.11	45,48,49,50	0
6	GOL	C	508	6/6	0.94	0.10	49,49,57,58	0
11	NAG	A	516	14/15	0.94	0.06	38,44,51,54	0
11	NAG	C	501	14/15	0.94	0.06	46,51,55,55	0
8	SO4	A	508	5/5	0.98	0.06	43,46,51,57	0
4	ZN	C	502	1/1	0.99	0.01	42,42,42,42	0
5	FE	A	502	1/1	0.99	0.02	45,45,45,45	1
5	FE	C	506	1/1	0.99	0.02	56,56,56,56	0
5	FE	D	509	1/1	0.99	0.02	48,48,48,48	1
4	ZN	B	501	1/1	1.00	0.01	51,51,51,51	0
5	FE	B	502	1/1	1.00	0.01	48,48,48,48	1
4	ZN	A	501	1/1	1.00	0.02	44,44,44,44	0
4	ZN	D	508	1/1	1.00	0.01	41,41,41,41	0

6.5 Other polymers ⓘ

There are no such residues in this entry.