



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 04:15 PM UTC

PDB ID : 1N9E / pdb_00001n9e
Title : Crystal structure of Pichia pastoris Lysyl Oxidase PPLO
Authors : Guss, J.M.; Duff, A.P.
Deposited on : 2002-11-24
Resolution : 1.65 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

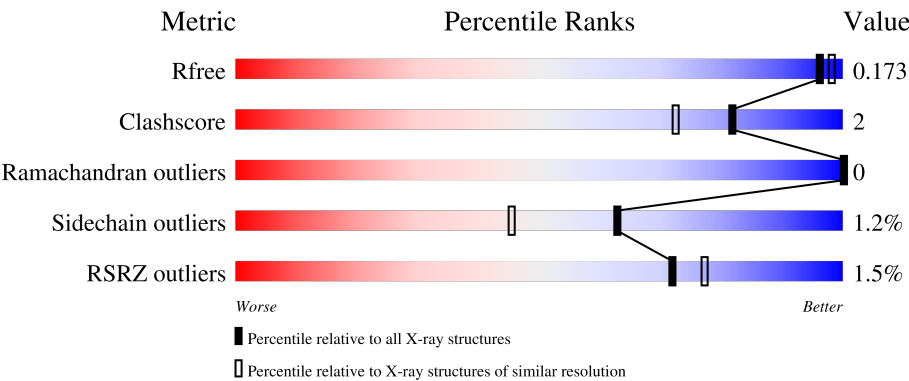
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.65 Å.

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



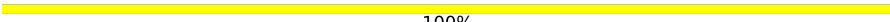
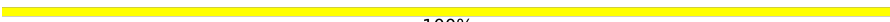
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	2563 (1.66-1.66)
Clashscore	190562	2662 (1.66-1.66)
Ramachandran outliers	187476	2621 (1.66-1.66)
Sidechain outliers	187428	2621 (1.66-1.66)
RSRZ outliers	180081	2564 (1.66-1.66)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	787	% 87%6% • 7%
1	B	787	2% 88%5% • 7%
1	C	787	2% 87%6% • 7%
1	D	787	% 87%6% • 7%
2	E	2	100%

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	G	2	 100%
2	H	2	 100%

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 28271 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 LYSYL OXIDASE.

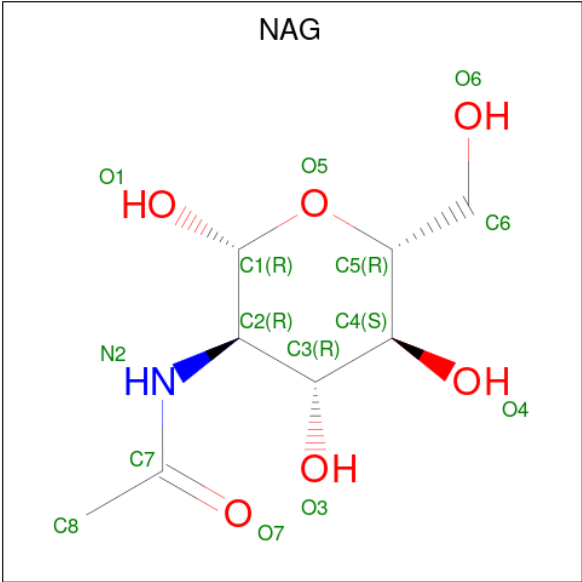
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	735	Total	C	N	O	S	5	8	0
			5951	3772	962	1201	16			
1	B	735	Total	C	N	O	S	4	7	0
			5950	3772	962	1200	16			
1	C	735	Total	C	N	O	S	5	7	0
			5950	3772	962	1200	16			
1	D	735	Total	C	N	O	S	5	8	0
			5951	3772	962	1201	16			

- 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	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
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 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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

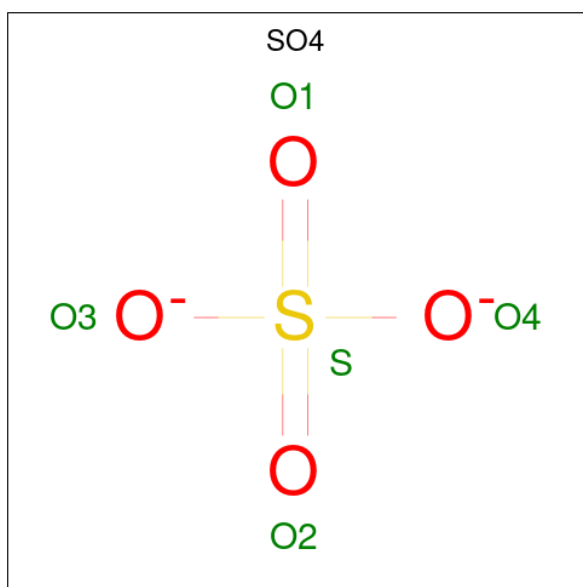
- Molecule 4 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

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

- Molecule 5 is CALCIUM ION (CCD ID: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	2	Total	Ca	0	0
			2	2		
5	B	2	Total	Ca	0	0
			2	2		
5	C	2	Total	Ca	0	0
			2	2		
5	D	2	Total	Ca	0	0
			2	2		

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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

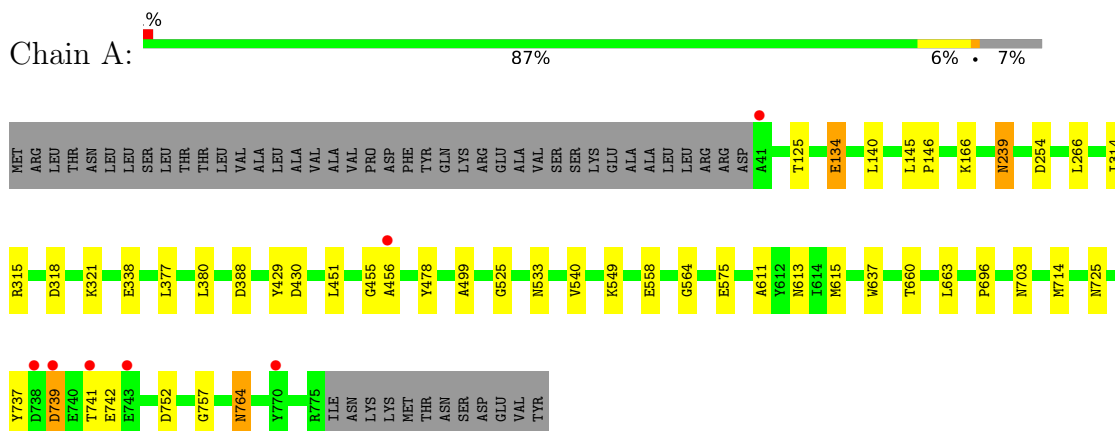
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	961	Total	O	0	0
			961	961		
7	B	1048	Total	O	0	0
			1048	1048		
7	C	1036	Total	O	0	0
			1036	1036		
7	D	936	Total	O	0	0
			936	936		

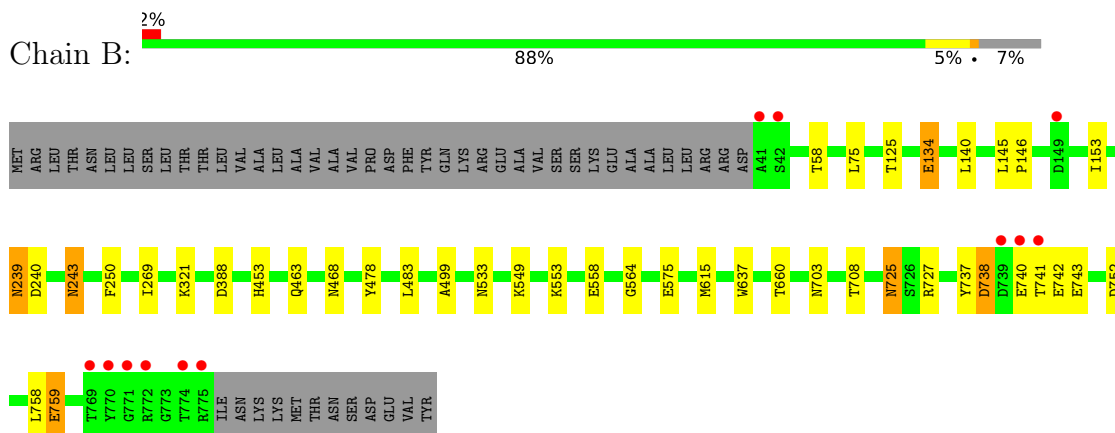
3 Residue-property plots [i](#)

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

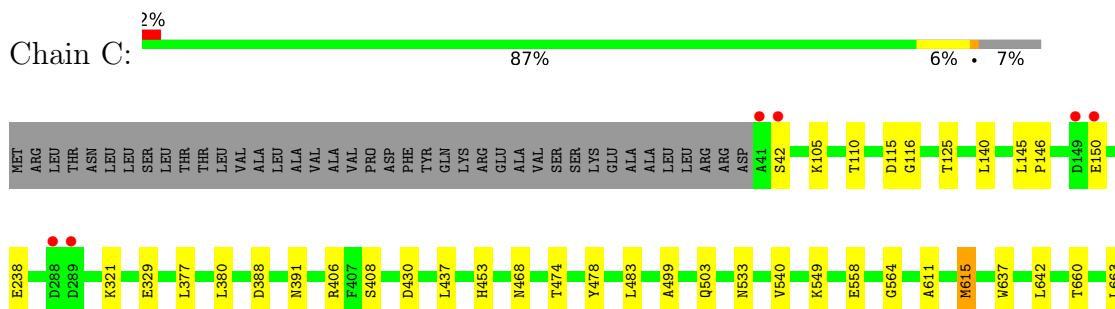
• Molecule 1: LYSYL OXIDASE

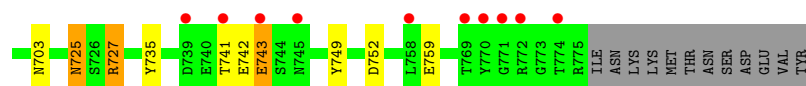


• Molecule 1: LYSYL OXIDASE

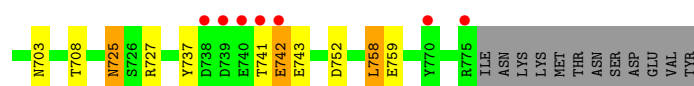
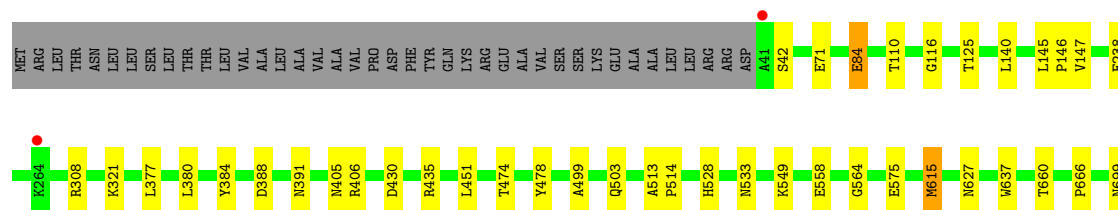
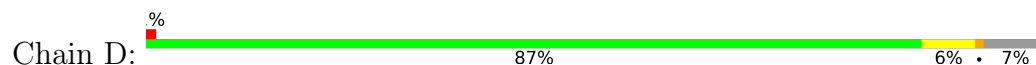


• Molecule 1: LYSYL OXIDASE





• Molecule 1: LYSYL OXIDASE



• 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



• 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	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	248.44Å 121.12Å 151.84Å 90.00° 124.64° 90.00°	Depositor
Resolution (Å)	24.60 – 1.65 24.60 – 1.65	Depositor EDS
% Data completeness (in resolution range)	94.7 (24.60-1.65) 90.5 (24.60-1.65)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.94 (at 1.65Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.161 , 0.187 0.166 , 0.173	Depositor DCC
R_{free} test set	39120 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	13.4	Xtriage
Anisotropy	0.180	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.010 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	28271	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.37% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: TPQ, CA, CU, NAG, SO4

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	0.57	0/6132	0.95	2/8355 (0.0%)
1	B	0.59	2/6126 (0.0%)	0.95	5/8347 (0.1%)
1	C	0.59	2/6126 (0.0%)	0.95	5/8347 (0.1%)
1	D	0.56	2/6132 (0.0%)	0.95	6/8355 (0.1%)
All	All	0.58	6/24516 (0.0%)	0.95	18/33404 (0.1%)

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	615[A]	MET	SD-CE	5.82	1.94	1.79
1	C	615[B]	MET	SD-CE	5.82	1.94	1.79
1	D	615[A]	MET	SD-CE	5.24	1.92	1.79
1	D	615[B]	MET	SD-CE	5.24	1.92	1.79
1	B	615[A]	MET	SD-CE	5.02	1.92	1.79
1	B	615[B]	MET	SD-CE	5.02	1.92	1.79

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	663	LEU	N-CA-C	7.37	119.31	111.28
1	C	725	ASN	N-CA-C	7.14	120.54	112.97
1	C	150	GLU	N-CA-C	-6.18	105.69	113.72
1	D	725	ASN	N-CA-C	6.12	119.49	112.57
1	C	660	THR	N-CA-C	6.02	117.84	111.28
1	B	725	ASN	N-CA-C	6.00	119.56	112.72
1	D	660	THR	N-CA-C	5.82	117.62	111.28
1	A	660	THR	N-CA-C	5.73	117.53	111.28
1	A	663	LEU	N-CA-C	5.72	117.51	111.28
1	D	742	GLU	N-CA-C	-5.70	105.63	112.58
1	B	553	LYS	CB-CG-CD	5.65	124.30	111.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	666	PRO	O-C-N	5.57	123.77	121.15
1	B	660	THR	N-CA-C	5.53	117.30	111.28
1	D	615[A]	MET	CB-CA-C	-5.35	103.99	110.15
1	D	615[B]	MET	CB-CA-C	-5.35	103.99	110.15
1	C	105	LYS	N-CA-C	5.30	117.13	111.36
1	B	737	TYR	CA-C-N	-5.26	115.78	122.77
1	B	737	TYR	C-N-CA	-5.26	115.78	122.77

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	5951	0	5497	37	0
1	B	5950	0	5496	29	0
1	C	5950	0	5496	27	0
1	D	5951	0	5497	36	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0
2	G	28	0	25	0	0
2	H	28	0	25	0	0
3	A	56	0	52	0	0
3	B	56	0	52	0	0
3	C	56	0	52	0	0
3	D	56	0	52	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	2	0	0	0	0
5	B	2	0	0	0	0
5	C	2	0	0	0	0
5	D	2	0	0	0	0
6	A	35	0	0	0	0
6	B	35	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	C	35	0	0	0	0
6	D	35	0	0	0	0
7	A	961	0	0	4	0
7	B	1048	0	0	3	0
7	C	1036	0	0	3	0
7	D	936	0	0	3	0
All	All	28271	0	22294	116	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 2.

All (116) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:764:ASN:HD22	1:A:764:ASN:H	1.16	0.90
7:C:8739:HOH:O	1:D:615[A]:MET:HE1	1.70	0.90
1:C:743:GLU:OE2	1:C:743:GLU:HA	1.71	0.90
1:A:549:LYS:NZ	1:A:575:GLU:OE1	2.06	0.88
1:A:499:ALA:HB1	7:A:2670:HOH:O	1.74	0.85
1:B:239:ASN:H	1:B:239:ASN:HD22	1.24	0.84
1:C:499:ALA:HB1	7:C:6670:HOH:O	1.86	0.74
1:B:759:GLU:HA	1:B:759:GLU:OE2	1.87	0.73
1:B:499:ALA:HB1	7:B:4670:HOH:O	1.88	0.73
1:B:738:ASP:OD1	1:B:738:ASP:C	2.33	0.72
1:A:239:ASN:HD22	1:A:239:ASN:H	1.37	0.72
1:A:764:ASN:H	1:A:764:ASN:ND2	1.89	0.70
1:B:738:ASP:OD1	1:B:740:GLU:N	2.27	0.66
1:B:239:ASN:HD22	1:B:239:ASN:N	1.94	0.66
1:D:391:ASN:ND2	1:D:503:GLN:HE21	1.94	0.64
1:D:549:LYS:NZ	1:D:575:GLU:OE1	2.31	0.64
1:D:741:THR:O	1:D:742:GLU:HB2	1.97	0.63
1:D:499:ALA:HB1	7:D:8670:HOH:O	1.97	0.63
1:C:391:ASN:ND2	1:C:503:GLN:HE21	1.99	0.61
1:D:627:ASN:ND2	7:D:9062:HOH:O	2.35	0.59
1:A:166:LYS:NZ	7:A:2939:HOH:O	2.35	0.59
1:A:558[B]:GLU:CD	1:A:564:GLY:H	2.11	0.58
1:C:42:SER:OG	1:D:759:GLU:OE2	2.20	0.57
1:B:558[B]:GLU:CD	1:B:564:GLY:H	2.12	0.57
1:D:84:GLU:OE2	1:D:84:GLU:CA	2.52	0.57
1:C:558[B]:GLU:CD	1:C:564:GLY:H	2.12	0.56
1:A:377:LEU:HD21	1:A:380:LEU:HD11	1.87	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:741:THR:O	1:B:742:GLU:HB2	2.06	0.56
1:D:110:THR:HG22	1:D:116:GLY:HA3	1.88	0.55
1:D:558[B]:GLU:CD	1:D:564:GLY:H	2.14	0.54
1:A:613:ASN:ND2	1:A:615[B]:MET:HG3	2.22	0.54
1:B:463:GLN:NE2	7:B:2441:HOH:O	2.41	0.54
1:B:758:LEU:C	1:B:758:LEU:HD12	2.33	0.54
1:C:741:THR:O	1:C:742:GLU:HB2	2.08	0.53
1:C:615[B]:MET:HE1	1:D:708:THR:O	2.09	0.53
1:D:84:GLU:OE2	1:D:84:GLU:N	2.43	0.52
1:D:758:LEU:C	1:D:758:LEU:HD12	2.34	0.52
1:A:451:LEU:HD12	1:A:451:LEU:C	2.35	0.52
1:C:145:LEU:HA	1:C:146:PRO:C	2.35	0.52
1:D:84:GLU:OE2	1:D:84:GLU:HA	2.11	0.51
1:D:451:LEU:HD12	1:D:451:LEU:C	2.36	0.51
1:A:741:THR:O	1:A:742:GLU:HB2	2.11	0.51
1:A:239:ASN:HD22	1:A:239:ASN:N	2.03	0.51
1:A:737:TYR:CZ	1:A:742:GLU:HA	2.45	0.50
1:A:558[B]:GLU:OE1	7:A:2647:HOH:O	2.20	0.50
1:B:727:ARG:CZ	1:B:727:ARG:HA	2.42	0.50
1:A:145:LEU:HA	1:A:146:PRO:C	2.37	0.50
1:B:145:LEU:HA	1:B:146:PRO:C	2.37	0.49
1:D:405:ASN:HA	1:D:435:ARG:CZ	2.42	0.49
1:B:58:THR:HG21	7:B:5161:HOH:O	2.12	0.49
1:D:145:LEU:HA	1:D:146:PRO:C	2.38	0.49
1:A:739:ASP:N	1:A:739:ASP:OD1	2.45	0.49
1:C:321:LYS:HE2	1:D:752:ASP:O	2.12	0.48
1:D:377:LEU:HD21	1:D:380:LEU:HD11	1.95	0.48
1:C:377:LEU:HD21	1:C:380:LEU:HD11	1.94	0.48
1:A:456:ALA:HB2	1:B:240:ASP:HB2	1.95	0.48
1:C:238:GLU:HA	1:C:238:GLU:OE1	2.13	0.48
1:D:125[B]:THR:CG2	1:D:140:LEU:HB2	2.44	0.48
1:C:743:GLU:OE2	1:C:743:GLU:CA	2.54	0.47
1:A:125[B]:THR:HG22	1:A:140:LEU:O	2.14	0.47
1:A:703:ASN:HD21	1:B:453:HIS:HB2	1.79	0.47
1:C:533:ASN:HB2	1:C:637:TRP:CE3	2.50	0.47
1:B:533:ASN:HB2	1:B:637:TRP:CE3	2.50	0.47
1:A:533:ASN:HB2	1:A:637:TRP:CE3	2.51	0.46
1:A:757:GLY:HA2	1:B:759:GLU:OE2	2.14	0.46
1:C:453:HIS:HB2	1:D:703:ASN:HD21	1.80	0.46
1:B:125[B]:THR:CG2	1:B:140:LEU:HB2	2.46	0.46
1:D:533:ASN:HB2	1:D:637:TRP:CE3	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:615[A]:MET:HE1	1:B:708:THR:O	2.15	0.46
1:D:741:THR:OG1	1:D:743:GLU:HB2	2.16	0.45
1:C:125[B]:THR:HG22	1:C:140:LEU:HB2	1.99	0.45
1:C:749:TYR:CE2	1:D:308:ARG:HB3	2.52	0.45
1:D:238:GLU:H	1:D:238:GLU:HG2	1.66	0.45
1:D:727:ARG:CZ	1:D:727:ARG:HA	2.47	0.45
1:A:321:LYS:HE2	1:B:752:ASP:O	2.17	0.45
1:C:727:ARG:CZ	1:C:727:ARG:HA	2.48	0.44
1:A:314:ILE:HG23	1:A:318:ASP:HB3	1.99	0.44
1:A:737:TYR:OH	1:A:742:GLU:HA	2.18	0.44
1:A:540:VAL:HG11	1:A:611:ALA:HA	1.98	0.44
1:C:430:ASP:N	7:C:6516:HOH:O	2.36	0.43
1:A:254:ASP:HB2	1:A:266:LEU:HD11	2.00	0.43
1:C:125[B]:THR:CG2	1:C:140:LEU:HB2	2.48	0.43
1:D:384:TYR:CD2	1:D:528:HIS:HB3	2.54	0.43
1:C:540:VAL:HG11	1:C:611:ALA:HA	2.00	0.43
1:D:513:ALA:N	1:D:514:PRO:CD	2.82	0.43
1:D:758:LEU:C	1:D:758:LEU:CD1	2.92	0.43
1:C:406:ARG:HD2	1:C:474:THR:O	2.18	0.43
1:C:329:GLU:CD	1:D:321:LYS:HG3	2.44	0.42
1:C:468:ASN:HB3	1:C:483:LEU:HD11	2.01	0.42
1:B:134:GLU:OE2	1:B:134:GLU:HA	2.17	0.42
1:B:243:ASN:HD22	1:B:243:ASN:HA	1.61	0.42
1:B:250:PHE:HB2	1:B:269:ILE:HB	2.02	0.42
1:C:408:SER:HB3	1:C:437:LEU:HB2	2.02	0.42
1:D:737:TYR:CZ	1:D:742:GLU:HA	2.54	0.42
1:C:752:ASP:O	1:D:321:LYS:HE2	2.19	0.42
1:B:75:LEU:HG	1:B:153:ILE:HD11	2.02	0.41
1:C:735:TYR:CD1	1:C:735:TYR:C	2.98	0.41
1:D:125[B]:THR:HG22	1:D:140:LEU:O	2.20	0.41
1:D:406:ARG:HD2	1:D:474:THR:O	2.20	0.41
1:A:615[B]:MET:HE2	1:A:714:MET:HB2	2.02	0.41
1:A:637:TRP:CD1	1:A:637:TRP:H	2.38	0.41
1:B:758:LEU:HD12	1:B:758:LEU:O	2.21	0.41
1:B:549:LYS:NZ	1:B:575:GLU:OE1	2.54	0.41
1:A:338:GLU:HG2	7:A:3127:HOH:O	2.20	0.41
1:A:455:GLY:O	1:A:456:ALA:C	2.63	0.41
1:A:525[A]:GLY:O	1:A:696:PRO:HD2	2.20	0.41
1:A:613:ASN:ND2	1:A:615[B]:MET:CG	2.84	0.41
1:B:468:ASN:HB3	1:B:483:LEU:HD11	2.02	0.41
1:D:71:GLU:HB3	1:D:147:VAL:CG2	2.51	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:134:GLU:OE2	1:A:134:GLU:HA	2.19	0.41
1:A:752:ASP:O	1:B:321:LYS:HE2	2.21	0.41
1:C:110:THR:HG22	1:C:116:GLY:HA3	2.03	0.41
1:B:758:LEU:C	1:B:758:LEU:CD1	2.93	0.41
1:A:429:TYR:O	1:A:430:ASP:HB2	2.22	0.40
1:D:430:ASP:N	7:D:8516:HOH:O	2.44	0.40
1:A:314:ILE:HG22	1:A:315:ARG:O	2.22	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	740/787 (94%)	716 (97%)	24 (3%)	0	100	100
1	B	739/787 (94%)	717 (97%)	22 (3%)	0	100	100
1	C	739/787 (94%)	715 (97%)	24 (3%)	0	100	100
1	D	740/787 (94%)	715 (97%)	25 (3%)	0	100	100
All	All	2958/3148 (94%)	2863 (97%)	95 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	652/692 (94%)	646 (99%)	6 (1%)	70	56
1	B	651/692 (94%)	642 (99%)	9 (1%)	59	39
1	C	651/692 (94%)	642 (99%)	9 (1%)	59	39
1	D	652/692 (94%)	646 (99%)	6 (1%)	70	56
All	All	2606/2768 (94%)	2576 (99%)	30 (1%)	63	45

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	134	GLU
1	A	239	ASN
1	A	388	ASP
1	A	725	ASN
1	A	739	ASP
1	A	764	ASN
1	B	134	GLU
1	B	239	ASN
1	B	243	ASN
1	B	388	ASP
1	B	703	ASN
1	B	725	ASN
1	B	738	ASP
1	B	743	GLU
1	B	759	GLU
1	C	115	ASP
1	C	388	ASP
1	C	549	LYS
1	C	642	LEU
1	C	703	ASN
1	C	725	ASN
1	C	727	ARG
1	C	743	GLU
1	C	759	GLU
1	D	42	SER
1	D	84	GLU
1	D	388	ASP
1	D	699	ASN
1	D	725	ASN
1	D	758	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	239	ASN
1	A	243	ASN
1	A	613	ASN
1	A	627	ASN
1	A	703	ASN
1	A	710	HIS
1	A	732	GLN
1	A	745	ASN
1	A	764	ASN
1	B	70	GLN
1	B	239	ASN
1	B	243	ASN
1	B	272	ASN
1	B	608	ASN
1	B	710	HIS
1	B	745	ASN
1	C	286	GLN
1	C	391	ASN
1	C	571	GLN
1	C	583	ASN
1	C	613	ASN
1	C	745	ASN
1	D	243	ASN
1	D	391	ASN
1	D	583	ASN
1	D	627	ASN
1	D	703	ASN
1	D	710	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
1	TPQ	C	478	4,1	13,14,15	1.47	3 (23%)	13,19,21	0.99	0
1	TPQ	A	478	4,1	13,14,15	1.58	1 (7%)	13,19,21	0.78	0
1	TPQ	B	478	4,1	13,14,15	1.45	2 (15%)	13,19,21	0.99	1 (7%)
1	TPQ	D	478	4,1	13,14,15	1.47	2 (15%)	13,19,21	0.93	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
1	TPQ	C	478	4,1	-	2/5/22/24	0/1/1/1
1	TPQ	A	478	4,1	-	2/5/22/24	0/1/1/1
1	TPQ	B	478	4,1	-	2/5/22/24	0/1/1/1
1	TPQ	D	478	4,1	-	2/5/22/24	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	478	TPQ	O4-C4	-3.63	1.24	1.33
1	D	478	TPQ	O4-C4	-3.58	1.24	1.33
1	C	478	TPQ	O4-C4	-3.33	1.25	1.33
1	B	478	TPQ	O4-C4	-3.29	1.25	1.33
1	D	478	TPQ	C6-C1	2.18	1.40	1.34
1	C	478	TPQ	C6-C1	2.10	1.39	1.34
1	C	478	TPQ	C3-C4	2.08	1.39	1.36
1	B	478	TPQ	C6-C1	2.04	1.39	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	478	TPQ	C6-C5-C4	2.01	120.31	116.80

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	478	TPQ	C-CA-CB-C1
1	B	478	TPQ	C-CA-CB-C1
1	C	478	TPQ	C-CA-CB-C1
1	D	478	TPQ	C-CA-CB-C1
1	A	478	TPQ	N-CA-CB-C1
1	B	478	TPQ	N-CA-CB-C1
1	C	478	TPQ	N-CA-CB-C1
1	D	478	TPQ	N-CA-CB-C1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates ⓘ

8 monosaccharides are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	NAG	E	1	2,1	14,14,15	0.63	0	17,19,21	1.10	0
2	NAG	E	2	2	14,14,15	0.58	0	17,19,21	0.79	0
2	NAG	F	1	2,1	14,14,15	0.75	1 (7%)	17,19,21	1.24	3 (17%)
2	NAG	F	2	2	14,14,15	0.67	0	17,19,21	1.02	1 (5%)
2	NAG	G	1	2,1	14,14,15	0.71	1 (7%)	17,19,21	1.37	3 (17%)
2	NAG	G	2	2	14,14,15	0.49	0	17,19,21	3.19	5 (29%)
2	NAG	H	1	2,1	14,14,15	0.79	1 (7%)	17,19,21	1.00	2 (11%)
2	NAG	H	2	2	14,14,15	0.48	0	17,19,21	1.08	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	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1
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	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1	NAG	O5-C1	-2.19	1.40	1.43
2	G	1	NAG	O5-C1	-2.12	1.40	1.43
2	H	1	NAG	O5-C1	-2.02	1.40	1.43

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	NAG	C2-N2-C7	-7.64	112.67	122.90
2	G	2	NAG	C1-C2-N2	5.96	119.83	110.43
2	G	2	NAG	C4-C3-C2	-5.68	102.69	111.02
2	G	2	NAG	C1-O5-C5	4.47	118.17	112.19
2	G	2	NAG	O5-C1-C2	-3.82	105.37	111.29
2	G	1	NAG	C2-N2-C7	-3.17	118.65	122.90
2	F	2	NAG	C1-O5-C5	2.79	115.92	112.19
2	H	2	NAG	C1-O5-C5	2.71	115.82	112.19
2	F	1	NAG	C3-C4-C5	-2.66	105.41	110.23
2	F	1	NAG	O5-C1-C2	-2.65	107.19	111.29
2	G	1	NAG	C3-C4-C5	-2.59	105.54	110.23
2	G	1	NAG	O5-C1-C2	-2.52	107.39	111.29
2	H	1	NAG	C2-N2-C7	-2.50	119.55	122.90
2	F	1	NAG	C2-N2-C7	-2.45	119.61	122.90
2	H	1	NAG	O5-C1-C2	-2.37	107.62	111.29

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	2	NAG	O5-C5-C6-O6
2	F	2	NAG	O5-C5-C6-O6

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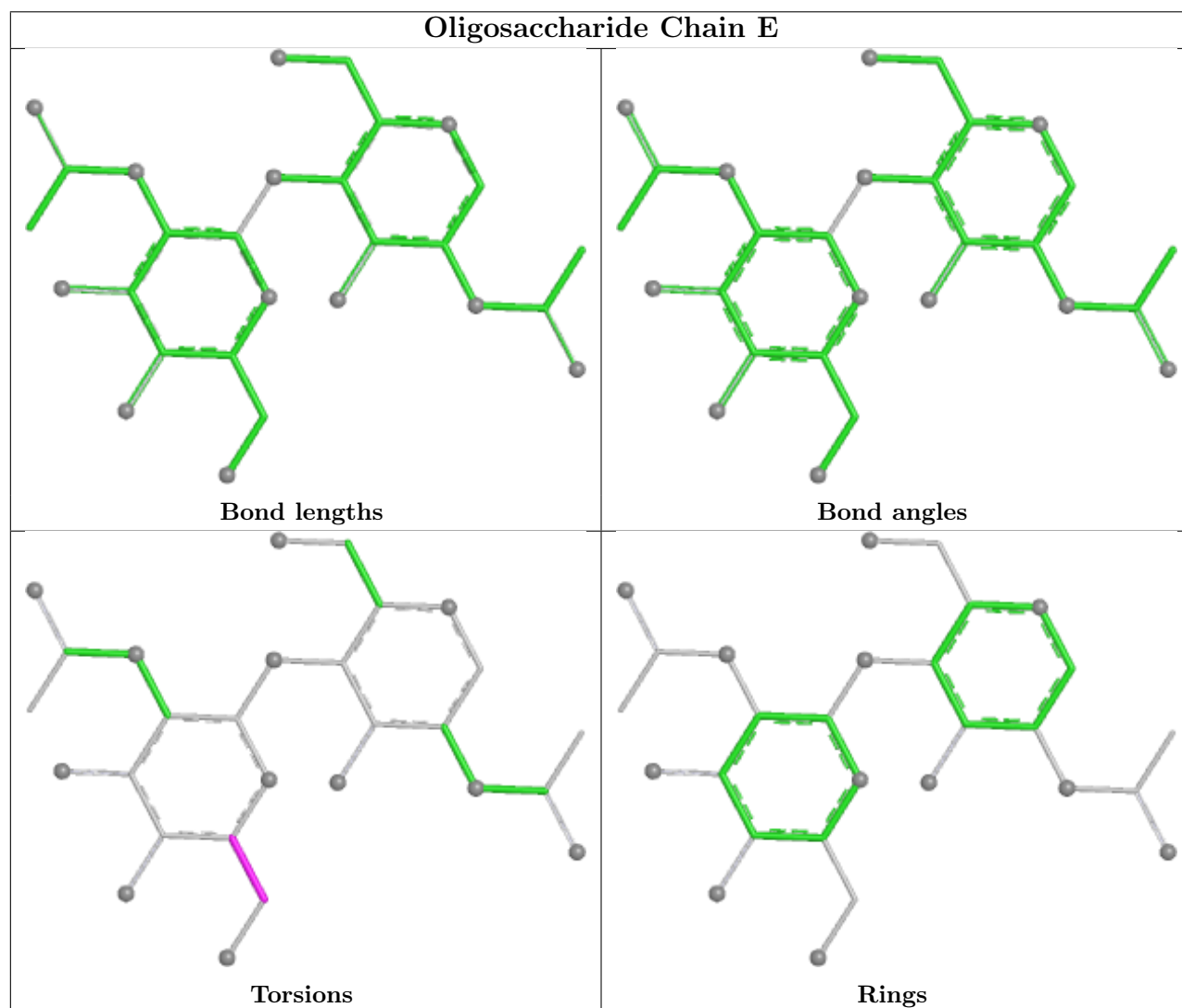
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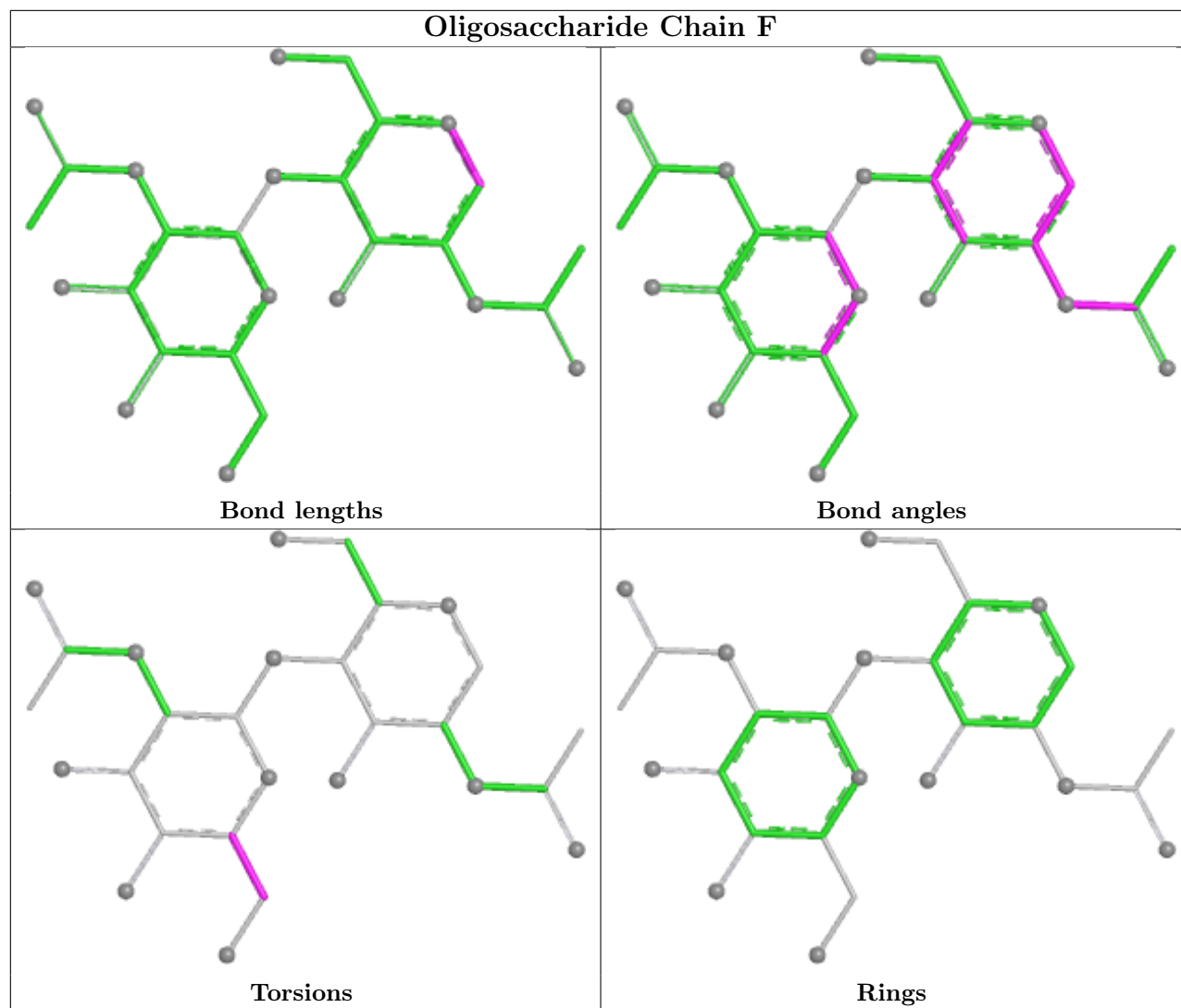
Mol	Chain	Res	Type	Atoms
2	F	2	NAG	C4-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6
2	H	2	NAG	O5-C5-C6-O6

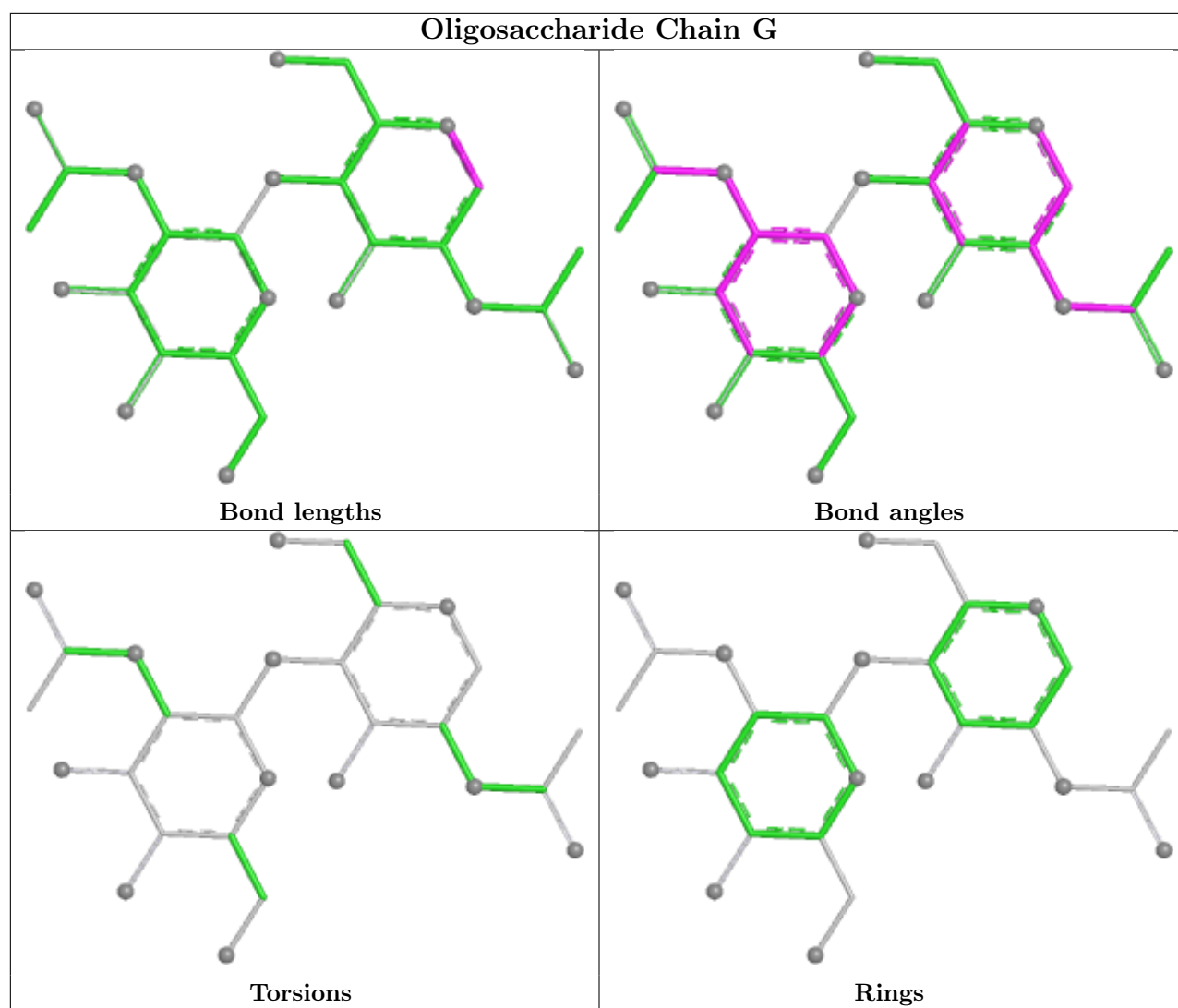
There are no ring outliers.

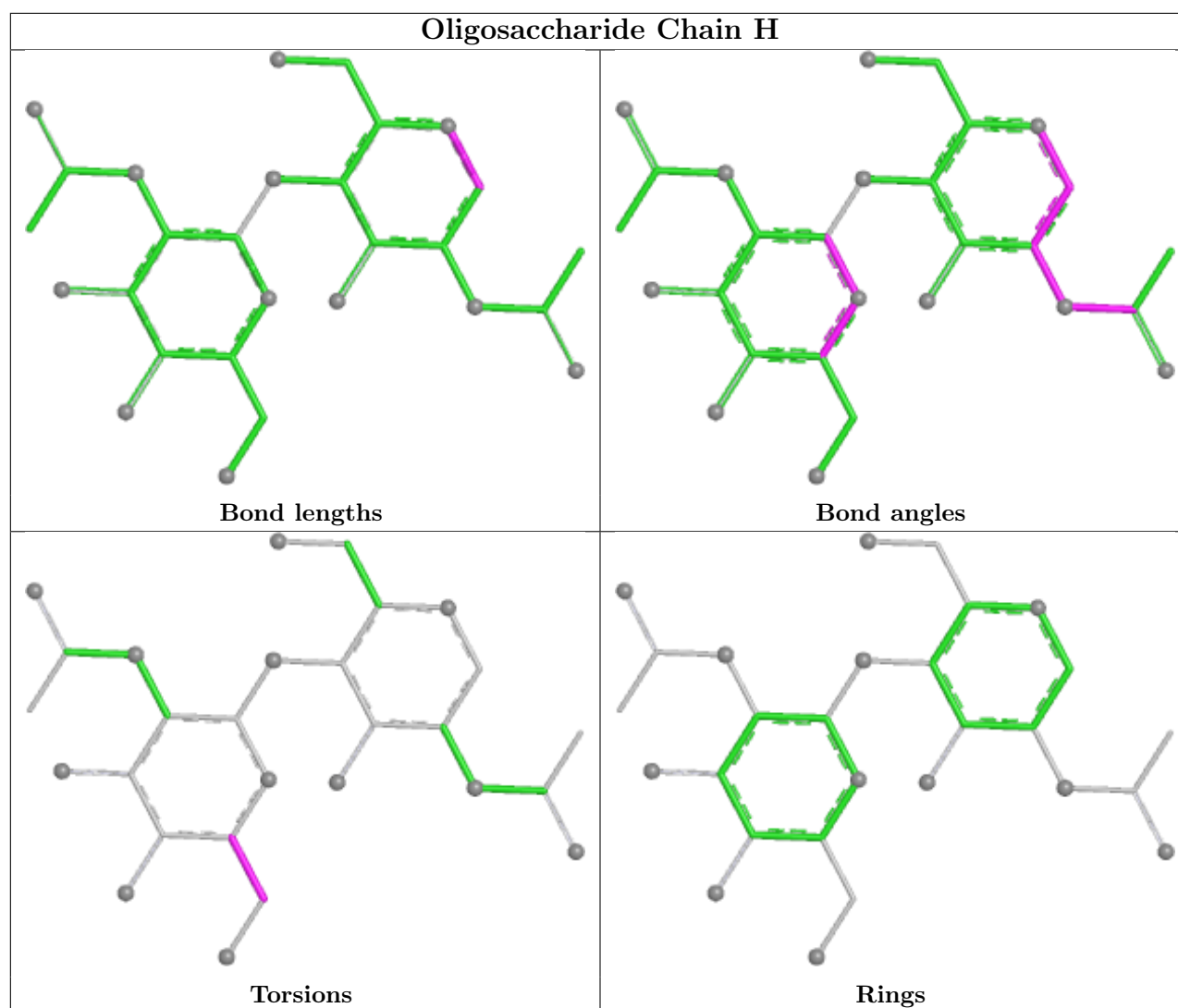
No monomer is involved in short contacts.

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









5.6 Ligand geometry [i](#)

Of 56 ligands modelled in this entry, 12 are monoatomic - leaving 44 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	$\# Z > 2$
3	NAG	C	1081	1	14,14,15	0.51	0	17,19,21	1.33	2 (11%)
6	SO4	A	811	-	4,4,4	0.47	0	6,6,6	0.76	0
6	SO4	B	820	-	4,4,4	0.32	0	6,6,6	0.21	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	D	842	-	4,4,4	0.25	0	6,6,6	0.23	0
3	NAG	A	1434	1	14,14,15	0.53	0	17,19,21	1.13	2 (11%)
6	SO4	B	824	-	4,4,4	0.30	0	6,6,6	0.19	0
3	NAG	A	1081	1	14,14,15	0.65	0	17,19,21	0.98	0
3	NAG	D	1191	1	14,14,15	0.49	0	17,19,21	0.76	0
6	SO4	C	835	-	4,4,4	0.34	0	6,6,6	0.08	0
6	SO4	D	841	-	4,4,4	0.36	0	6,6,6	0.60	0
6	SO4	A	816	-	4,4,4	0.27	0	6,6,6	0.13	0
6	SO4	D	846	-	4,4,4	0.25	0	6,6,6	0.11	0
3	NAG	C	1309	1	14,14,15	0.62	0	17,19,21	1.07	2 (11%)
3	NAG	C	1191	1	14,14,15	0.57	0	17,19,21	1.19	2 (11%)
6	SO4	A	814	-	4,4,4	0.32	0	6,6,6	0.33	0
6	SO4	A	812	-	4,4,4	0.32	0	6,6,6	0.14	0
3	NAG	B	1434	1	14,14,15	0.57	0	17,19,21	0.87	2 (11%)
6	SO4	B	825	-	4,4,4	0.27	0	6,6,6	0.12	0
6	SO4	C	830	-	4,4,4	0.38	0	6,6,6	0.20	0
6	SO4	C	831	-	4,4,4	0.37	0	6,6,6	0.40	0
6	SO4	D	840	-	4,4,4	0.30	0	6,6,6	0.23	0
3	NAG	A	1191	1	14,14,15	0.70	0	17,19,21	1.05	0
3	NAG	B	1081	1	14,14,15	0.48	0	17,19,21	1.44	2 (11%)
3	NAG	D	1081	1	14,14,15	0.72	0	17,19,21	1.12	2 (11%)
6	SO4	B	821	-	4,4,4	0.49	0	6,6,6	0.83	0
6	SO4	C	832	-	4,4,4	0.30	0	6,6,6	0.21	0
6	SO4	D	843	-	4,4,4	0.26	0	6,6,6	0.36	0
3	NAG	C	1434	1	14,14,15	0.54	0	17,19,21	0.75	0
3	NAG	D	1309	1	14,14,15	0.62	0	17,19,21	1.02	0
6	SO4	B	822	-	4,4,4	0.32	0	6,6,6	0.13	0
6	SO4	B	826	-	4,4,4	0.26	0	6,6,6	0.06	0
6	SO4	D	845	-	4,4,4	0.32	0	6,6,6	0.14	0
3	NAG	D	1434	1	14,14,15	0.55	0	17,19,21	1.05	1 (5%)
6	SO4	B	823	-	4,4,4	0.44	0	6,6,6	0.53	0
6	SO4	A	815	-	4,4,4	0.37	0	6,6,6	0.28	0
6	SO4	A	813	-	4,4,4	0.30	0	6,6,6	0.21	0
3	NAG	B	1309	1	14,14,15	0.56	0	17,19,21	0.70	0
6	SO4	D	844	-	4,4,4	0.22	0	6,6,6	0.23	0
6	SO4	C	833	-	4,4,4	0.23	0	6,6,6	0.23	0
6	SO4	C	836	-	4,4,4	0.30	0	6,6,6	0.17	0
6	SO4	A	810	-	4,4,4	0.44	0	6,6,6	0.18	0
6	SO4	C	834	-	4,4,4	0.22	0	6,6,6	0.22	0
3	NAG	B	1191	1	14,14,15	0.56	0	17,19,21	0.98	1 (5%)
3	NAG	A	1309	1	14,14,15	0.56	0	17,19,21	0.91	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	1081	1	-	0/6/23/26	0/1/1/1
3	NAG	D	1081	1	-	0/6/23/26	0/1/1/1
3	NAG	D	1191	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1309	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1434	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1081	1	-	2/6/23/26	0/1/1/1
3	NAG	D	1309	1	-	0/6/23/26	0/1/1/1
3	NAG	A	1081	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1191	1	-	0/6/23/26	0/1/1/1
3	NAG	C	1309	1	-	2/6/23/26	0/1/1/1
3	NAG	D	1434	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1434	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1191	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1434	1	-	2/6/23/26	0/1/1/1
3	NAG	A	1191	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1309	1	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1081	NAG	O5-C1-C2	-3.93	105.22	111.29
3	B	1081	NAG	C1-C2-N2	2.88	114.97	110.43
3	D	1081	NAG	C4-C3-C2	2.87	115.23	111.02
3	C	1081	NAG	C3-C4-C5	2.82	115.34	110.23
3	C	1081	NAG	O5-C1-C2	-2.64	107.21	111.29
3	C	1191	NAG	O5-C1-C2	-2.57	107.31	111.29
3	D	1434	NAG	O5-C1-C2	-2.46	107.48	111.29
3	C	1191	NAG	C6-C5-C4	-2.44	107.03	113.02
3	C	1309	NAG	O5-C1-C2	-2.43	107.53	111.29
3	A	1434	NAG	C2-N2-C7	-2.40	119.69	122.90
3	C	1309	NAG	C1-O5-C5	2.33	115.31	112.19
3	A	1309	NAG	O5-C1-C2	-2.28	107.76	111.29
3	B	1191	NAG	O5-C5-C6	2.25	112.04	107.66
3	A	1434	NAG	O5-C1-C2	-2.19	107.91	111.29
3	B	1434	NAG	O5-C1-C2	-2.09	108.05	111.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1434	NAG	C2-N2-C7	-2.05	120.16	122.90
3	D	1081	NAG	O5-C1-C2	2.02	114.42	111.29

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1191	NAG	O5-C5-C6-O6
3	A	1434	NAG	O5-C5-C6-O6
3	C	1434	NAG	O5-C5-C6-O6
3	A	1434	NAG	C4-C5-C6-O6
3	B	1191	NAG	C4-C5-C6-O6
3	D	1434	NAG	O5-C5-C6-O6
3	C	1434	NAG	C4-C5-C6-O6
3	D	1434	NAG	C4-C5-C6-O6
3	A	1081	NAG	C8-C7-N2-C2
3	A	1191	NAG	O5-C5-C6-O6
3	A	1191	NAG	C4-C5-C6-O6
3	A	1081	NAG	O7-C7-N2-C2
3	C	1309	NAG	O5-C5-C6-O6
3	C	1309	NAG	C4-C5-C6-O6
3	B	1434	NAG	C4-C5-C6-O6
3	B	1434	NAG	O5-C5-C6-O6
3	C	1081	NAG	C3-C2-N2-C7
3	C	1081	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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	734/787 (93%)	-0.42	7 (0%) 79 84	6, 14, 30, 68	10 (1%)
1	B	734/787 (93%)	-0.50	12 (1%) 70 75	5, 13, 30, 63	9 (1%)
1	C	734/787 (93%)	-0.45	16 (2%) 62 68	5, 13, 30, 65	10 (1%)
1	D	734/787 (93%)	-0.40	9 (1%) 76 81	6, 15, 30, 71	10 (1%)
All	All	2936/3148 (93%)	-0.44	44 (1%) 72 77	5, 14, 30, 71	39 (1%)

All (44) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	41	ALA	6.0
1	B	41	ALA	5.8
1	C	41	ALA	5.8
1	A	739	ASP	5.3
1	A	741	THR	5.0
1	D	741	THR	4.5
1	D	739	ASP	4.3
1	B	769	THR	3.7
1	D	41	ALA	3.7
1	B	739	ASP	3.4
1	C	770	TYR	3.3
1	C	739	ASP	3.2
1	B	771	GLY	3.0
1	A	770	TYR	3.0
1	A	456	ALA	3.0
1	B	741	THR	2.8
1	D	770	TYR	2.8
1	C	774	THR	2.7
1	C	288	ASP	2.6
1	B	770	TYR	2.6
1	C	42	SER	2.5

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Mol	Chain	Res	Type	RSRZ
1	B	42	SER	2.5
1	D	740	GLU	2.5
1	C	149	ASP	2.4
1	C	769	THR	2.4
1	A	738	ASP	2.4
1	D	775	ARG	2.4
1	B	772	ARG	2.3
1	C	771	GLY	2.3
1	C	772	ARG	2.3
1	D	738	ASP	2.3
1	C	741	THR	2.3
1	C	745	ASN	2.2
1	D	742	GLU	2.2
1	A	743	GLU	2.1
1	C	758	LEU	2.1
1	C	743	GLU	2.1
1	B	774	THR	2.1
1	D	264	LYS	2.1
1	B	740	GLU	2.1
1	C	289	ASP	2.1
1	B	149	ASP	2.0
1	C	150	GLU	2.0
1	B	775	ARG	2.0

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	TPQ	C	478	14/15	0.93	0.14	13,23,46,53	0
1	TPQ	B	478	14/15	0.94	0.11	12,23,45,52	0
1	TPQ	A	478	14/15	0.94	0.12	12,26,46,52	0
1	TPQ	D	478	14/15	0.94	0.13	14,28,48,51	0

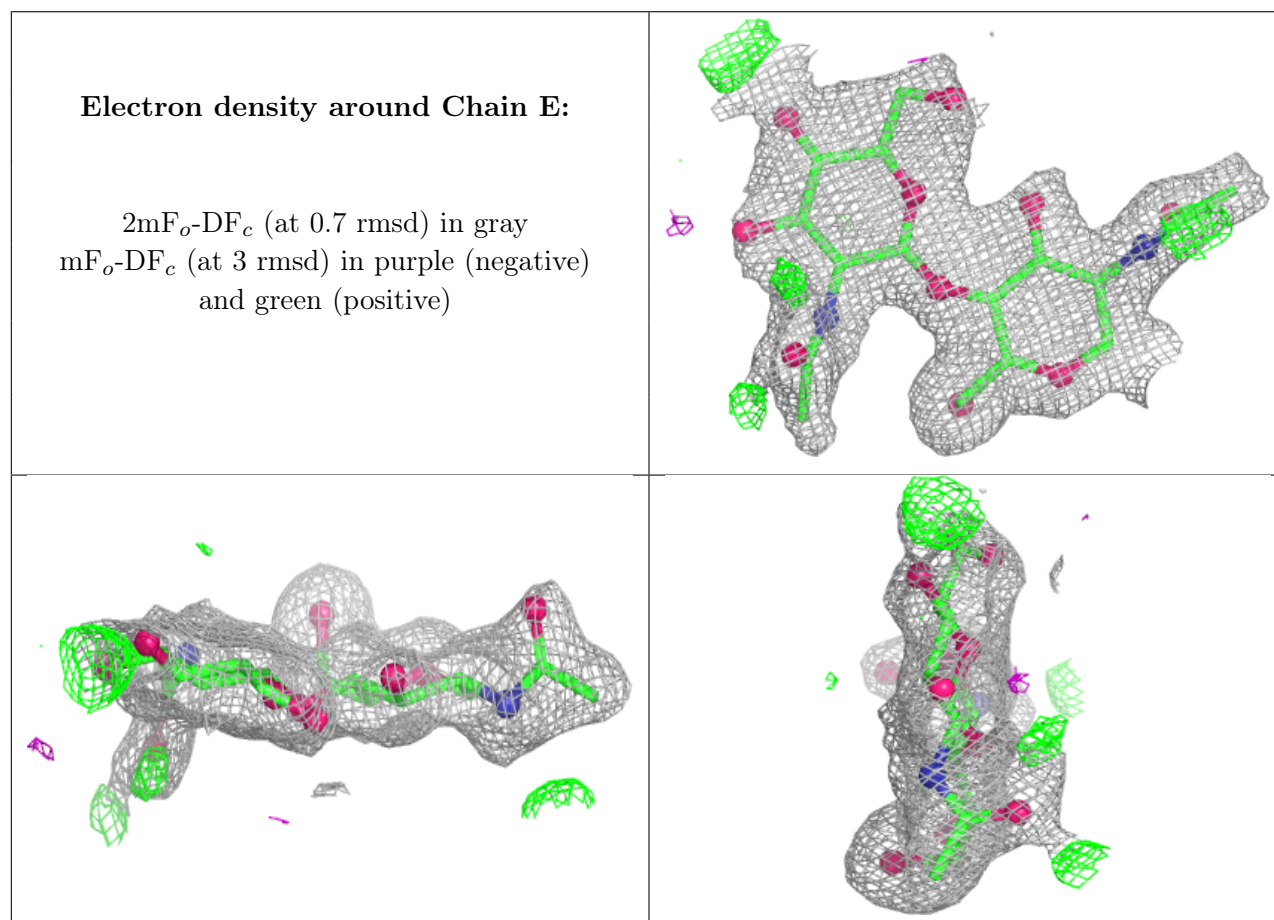
6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

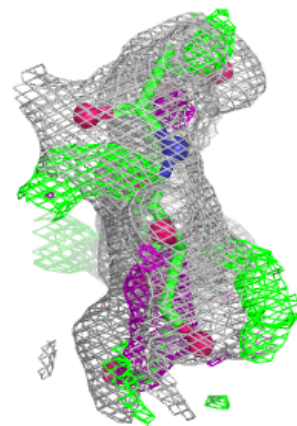
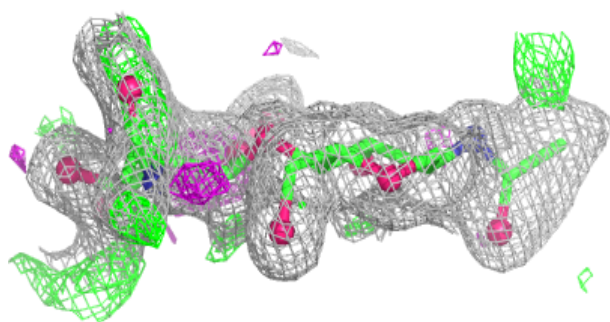
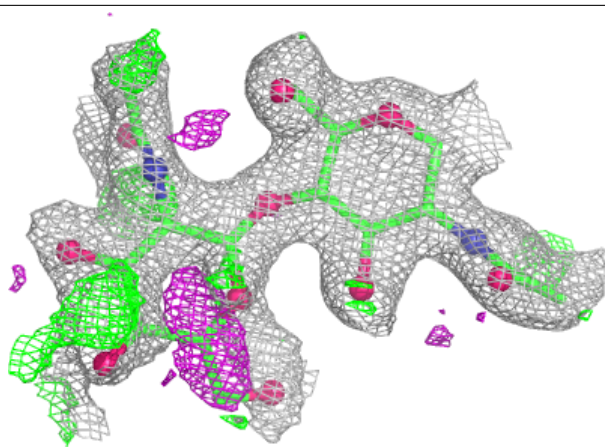
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	F	2	14/15	0.71	0.17	30,44,53,56	0
2	NAG	G	2	14/15	0.77	0.14	28,38,56,58	0
2	NAG	H	2	14/15	0.79	0.13	38,46,53,54	0
2	NAG	E	2	14/15	0.83	0.12	36,44,52,54	0
2	NAG	H	1	14/15	0.90	0.10	16,30,38,39	0
2	NAG	E	1	14/15	0.92	0.09	22,30,39,45	0
2	NAG	F	1	14/15	0.93	0.08	15,23,30,37	0
2	NAG	G	1	14/15	0.94	0.08	16,24,36,37	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.



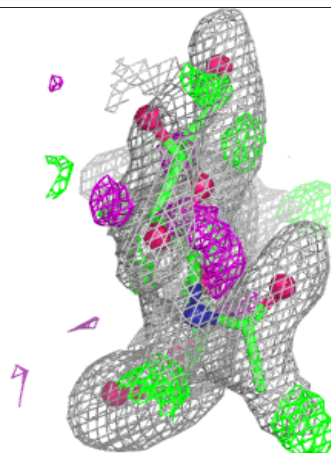
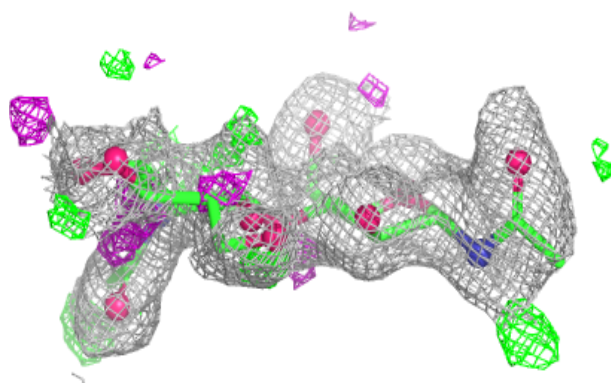
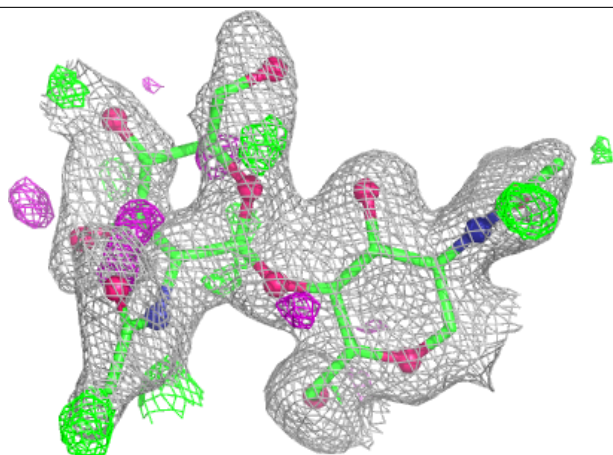
Electron density around Chain F:

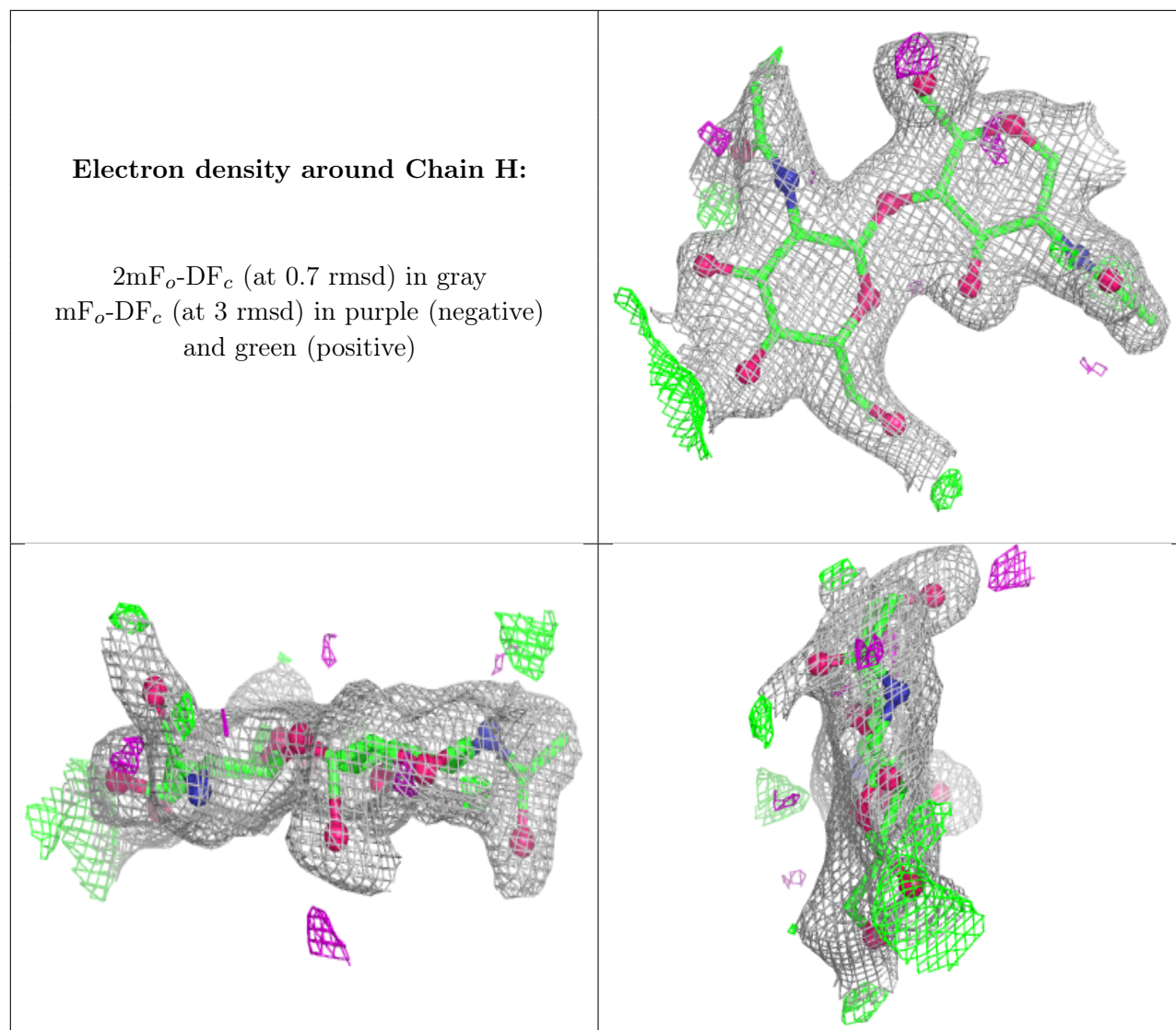
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain 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 ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	B	1309	14/15	0.49	0.19	67,75,77,79	0
3	NAG	C	1309	14/15	0.51	0.18	65,69,73,76	0
3	NAG	C	1081	14/15	0.57	0.22	55,66,74,77	0
3	NAG	D	1434	14/15	0.57	0.21	50,63,75,77	0
3	NAG	C	1434	14/15	0.58	0.18	51,64,68,71	0
3	NAG	A	1434	14/15	0.63	0.19	51,62,70,72	0
3	NAG	A	1081	14/15	0.64	0.16	45,51,68,71	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	B	1434	14/15	0.64	0.18	48,61,66,67	0
3	NAG	D	1081	14/15	0.67	0.16	61,70,76,76	0
6	SO4	C	832	5/5	0.68	0.18	38,49,51,59	0
6	SO4	A	816	5/5	0.69	0.15	52,58,63,65	0
6	SO4	C	836	5/5	0.70	0.16	44,51,59,61	0
3	NAG	B	1081	14/15	0.73	0.14	36,44,48,50	0
6	SO4	D	846	5/5	0.73	0.16	61,62,69,70	0
6	SO4	B	822	5/5	0.74	0.17	33,44,49,55	0
6	SO4	B	826	5/5	0.78	0.13	35,52,56,58	0
6	SO4	A	815	5/5	0.78	0.14	23,32,44,48	0
6	SO4	D	840	5/5	0.82	0.18	22,27,32,36	0
6	SO4	A	814	5/5	0.82	0.13	26,31,45,47	0
3	NAG	D	1309	14/15	0.83	0.12	30,37,45,46	0
6	SO4	C	835	5/5	0.84	0.13	18,30,45,46	0
6	SO4	D	842	5/5	0.84	0.12	39,46,51,54	0
6	SO4	D	845	5/5	0.84	0.12	20,36,45,47	0
6	SO4	A	812	5/5	0.84	0.11	30,41,45,48	0
6	SO4	A	810	5/5	0.85	0.16	21,26,29,34	0
6	SO4	D	843	5/5	0.87	0.11	29,39,45,47	0
6	SO4	D	844	5/5	0.87	0.10	34,34,43,46	0
6	SO4	B	820	5/5	0.88	0.18	20,26,32,35	0
3	NAG	A	1309	14/15	0.88	0.11	23,33,42,43	0
6	SO4	A	811	5/5	-	-	10,36,47,52	5
6	SO4	C	833	5/5	0.89	0.11	23,38,42,42	0
6	SO4	C	830	5/5	0.89	0.16	23,25,26,33	0
6	SO4	B	825	5/5	0.89	0.24	15,30,41,43	0
3	NAG	A	1191	14/15	0.90	0.09	23,27,37,37	0
3	NAG	B	1191	14/15	0.91	0.09	15,25,30,41	0
3	NAG	D	1191	14/15	0.91	0.09	23,28,34,39	0
6	SO4	B	821	5/5	-	-	7,34,44,51	5
6	SO4	B	824	5/5	0.92	0.20	26,29,36,38	0
6	SO4	C	834	5/5	0.92	0.19	24,27,35,37	0
6	SO4	A	813	5/5	0.93	0.18	22,31,42,42	0
6	SO4	B	823	5/5	0.94	0.15	19,30,33,35	0
3	NAG	C	1191	14/15	0.95	0.06	12,19,23,25	0
4	CU	B	801	1/1	0.98	0.03	16,16,16,16	0
6	SO4	C	831	5/5	-	-	7,39,48,50	5
4	CU	A	801	1/1	0.99	0.03	16,16,16,16	0
4	CU	C	801	1/1	0.99	0.03	16,16,16,16	0
4	CU	D	801	1/1	0.99	0.03	17,17,17,17	0
5	CA	A	802	1/1	0.99	0.02	11,11,11,11	0
5	CA	D	5802	1/1	0.99	0.02	12,12,12,12	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	CA	D	5803	1/1	0.99	0.03	14,14,14,14	0
6	SO4	D	841	5/5	-	-	14,37,52,58	5
5	CA	B	1802	1/1	1.00	0.02	8,8,8,8	0
5	CA	B	1803	1/1	1.00	0.02	10,10,10,10	0
5	CA	C	3802	1/1	1.00	0.01	10,10,10,10	0
5	CA	C	3803	1/1	1.00	0.01	10,10,10,10	0
5	CA	A	803	1/1	1.00	0.01	13,13,13,13	0

6.5 Other polymers [i](#)

There are no such residues in this entry.