



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:30 PM UTC

PDB ID : 9EY6 / pdb_00009ey6
Title : Crystal structure of human tyrosinase-related protein 1 (TYRP1)
Authors : Ng, Y.M.; Soler-Lopez, M.
Deposited on : 2024-04-09
Resolution : 2.23 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

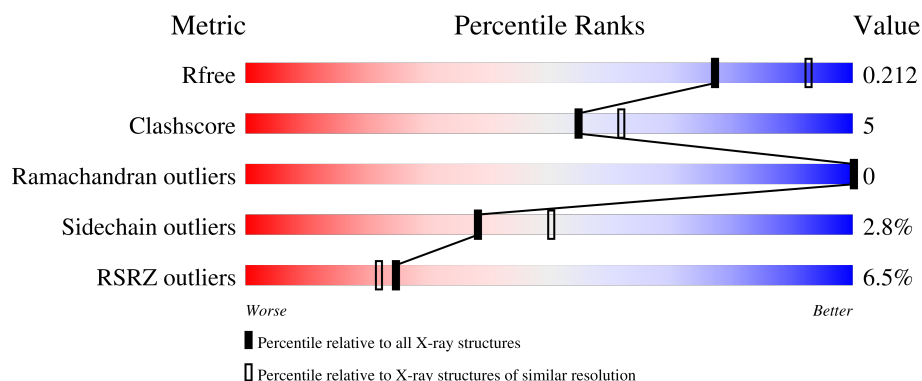
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.23 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	7682 (2.24-2.20)
Clashscore	190562	8402 (2.24-2.20)
Ramachandran outliers	187476	8303 (2.24-2.20)
Sidechain outliers	187428	8304 (2.24-2.20)
RSRZ outliers	180081	7683 (2.24-2.20)

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	447	6% 87% 12%
2	B	4	50% 50%
3	C	3	100%
3	E	3	33% 67%
4	D	3	67% 33%

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Mol	Chain	Length	Quality of chain
5	F	6	17% 33% 50%

2 Entry composition [i](#)

There are 9 unique types of molecules in this entry. The entry contains 4057 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 5,6-dihydroxyindole-2-carboxylic acid oxidase.

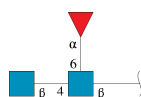
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	447	Total	C	N	O	S	0	3	0
			3596	2254	644	675	23			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



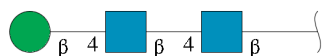
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	B	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



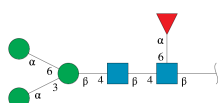
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	C	3	Total	C	N	O	0	0	0
			38	22	2	14			
3	E	3	Total	C	N	O	0	0	0
			38	22	2	14			

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



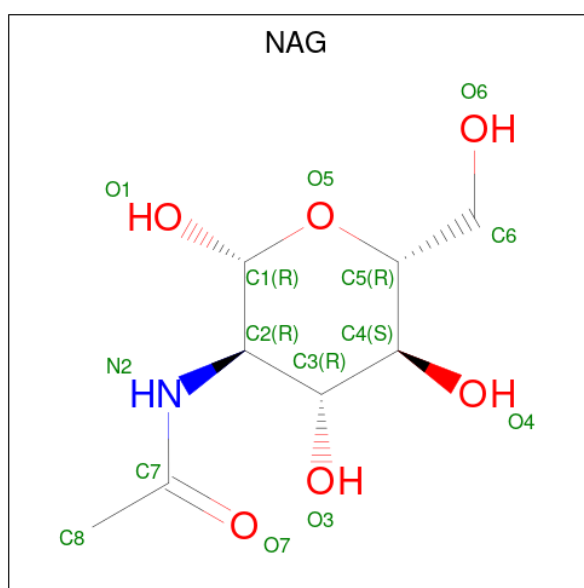
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	D	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 5 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



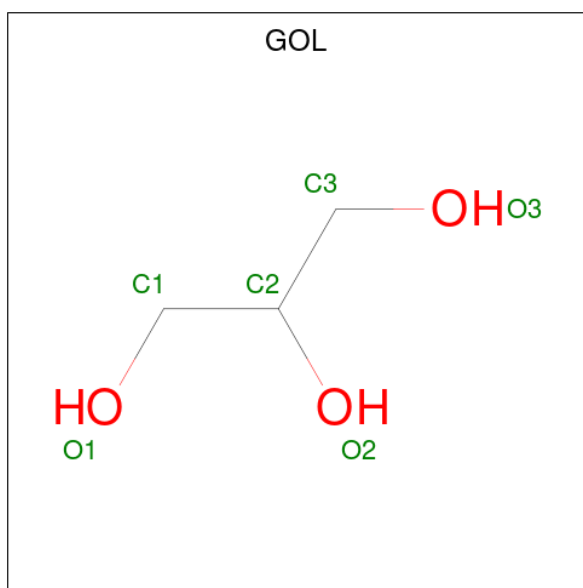
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	F	6	Total	C	N	O	0	0	0
			71	40	2	29			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 7 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 8 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	5	Total	Zn	0	0
			5	5		

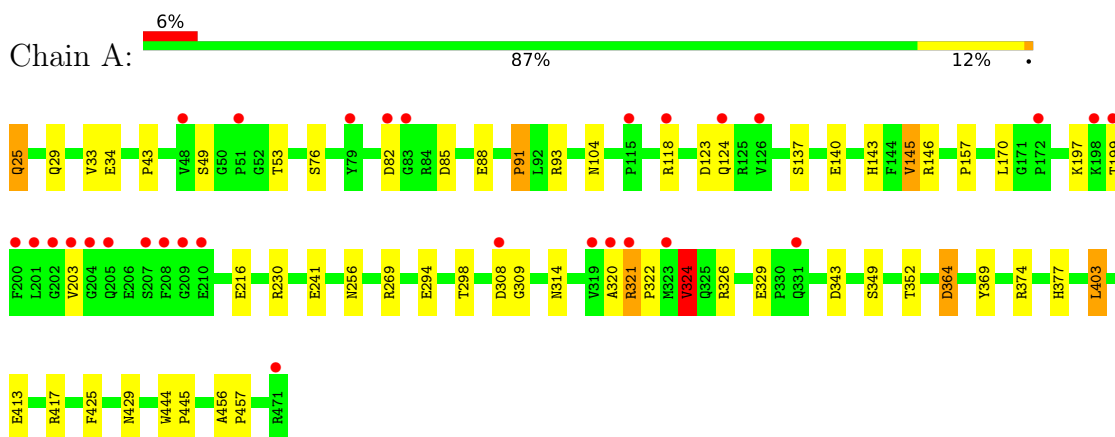
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
9	A	182	Total	O	0	0
			182	182		

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.

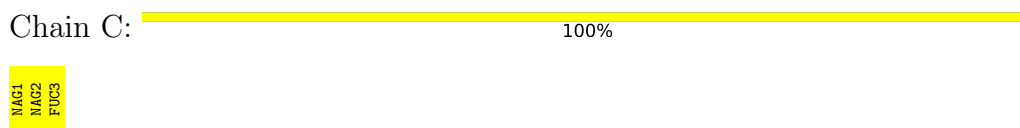
- Molecule 1: 5,6-dihydroxyindole-2-carboxylic acid oxidase



- Molecule 2: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose



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

Chain D:  67% 33%

MAG1
MAG2
BMA3

- Molecule 5: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[alpha-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  17% 33% 50%

MAG1
MAG2
BMA3
MAN4
MAN5
FUC6

4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	103.75Å 103.75Å 140.99Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	44.97 – 2.23 44.97 – 2.23	Depositor EDS
% Data completeness (in resolution range)	98.0 (44.97-2.23) 98.1 (44.97-2.23)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.10 (at 2.22Å)	Xtriage
Refinement program	REFMAC 5.8.0425	Depositor
R, R_{free}	0.170 , 0.213 (Not available) , 0.212	Depositor DCC
R_{free} test set	2108 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	44.4	Xtriage
Anisotropy	0.315	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.025 for -h,-k,l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4057	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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: MAN, NAG, GOL, BMA, FUC, ZN

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	0.66	0/3709	1.16	12/5052 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	6

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	324	VAL	N-CA-CB	-9.34	101.47	112.39
1	A	91	PRO	N-CA-CB	-7.46	94.40	102.60
1	A	53	THR	CA-CB-OG1	-6.73	99.51	109.60
1	A	145	VAL	N-CA-CB	6.50	122.53	110.77
1	A	329	GLU	CB-CG-CD	6.34	123.37	112.60
1	A	364	ASP	CA-CB-CG	6.16	118.76	112.60
1	A	241	GLU	CB-CG-CD	6.03	122.86	112.60
1	A	88	GLU	CB-CG-CD	5.86	122.56	112.60
1	A	403	LEU	N-CA-CB	-5.53	101.98	110.16
1	A	104	ASN	CB-CA-C	5.28	119.52	111.02
1	A	425	PHE	CA-CB-CG	5.25	119.05	113.80
1	A	343	ASP	CA-CB-CG	5.11	117.71	112.60

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	146[A]	ARG	Sidechain
1	A	146[B]	ARG	Sidechain
1	A	230	ARG	Sidechain
1	A	326	ARG	Sidechain
1	A	374	ARG	Sidechain
1	A	93	ARG	Sidechain

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3596	0	3371	33	1
2	B	50	0	43	0	0
3	C	38	0	34	0	0
3	E	38	0	34	2	0
4	D	39	0	34	0	0
5	F	71	0	61	2	0
6	A	14	0	13	0	0
7	A	24	0	32	4	0
8	A	5	0	0	0	0
9	A	182	0	0	10	0
All	All	4057	0	3622	35	1

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 5.

All (35) 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:25:GLN:OE1	1:A:25:GLN:N	1.90	1.03
1:A:308:ASP:O	9:A:602:HOH:O	1.97	0.82
1:A:320:ALA:O	1:A:322:PRO:HD3	1.81	0.78
1:A:309:GLY:N	9:A:601:HOH:O	1.77	0.76
1:A:140:GLU:HA	7:A:505:GOL:H31	1.74	0.69
5:F:1:NAG:H62	5:F:2:NAG:C1	2.28	0.64
1:A:140:GLU:HG2	7:A:505:GOL:H32	1.81	0.61
1:A:157:PRO:HA	3:E:3:FUC:H61	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:ASP:OD2	1:A:199:THR:HG22	2.02	0.59
1:A:308:ASP:HA	9:A:601:HOH:O	2.03	0.58
1:A:308:ASP:CA	9:A:601:HOH:O	2.51	0.58
1:A:118:ARG:NH1	1:A:124[A]:GLN:OE1	2.38	0.56
1:A:352:THR:HG21	9:A:775:HOH:O	2.06	0.54
1:A:137:SER:OG	1:A:140:GLU:HG3	2.08	0.54
1:A:143:HIS:CD2	7:A:505:GOL:H12	2.42	0.54
1:A:413:GLU:OE2	1:A:417[B]:ARG:NE	2.37	0.51
1:A:294:GLU:O	1:A:298:THR:HG23	2.11	0.51
1:A:157:PRO:HA	3:E:3:FUC:C6	2.41	0.50
1:A:29:GLN:HG3	1:A:43:PRO:HB3	1.93	0.50
1:A:216:GLU:OE1	7:A:502:GOL:H12	2.11	0.50
1:A:256:ASN:OD1	9:A:604:HOH:O	2.20	0.49
1:A:123:ASP:OD1	9:A:603:HOH:O	2.18	0.49
1:A:197:LYS:NZ	9:A:613:HOH:O	2.45	0.48
1:A:143:HIS:HD2	9:A:725:HOH:O	1.97	0.48
1:A:321:ARG:NH2	9:A:617:HOH:O	2.48	0.46
1:A:324:VAL:O	1:A:324:VAL:CG1	2.65	0.44
1:A:324:VAL:O	1:A:324:VAL:HG13	2.18	0.44
1:A:76:SER:HB3	1:A:429:ASN:O	2.17	0.43
1:A:269:ARG:NH1	1:A:314:ASN:OD1	2.44	0.42
1:A:444:TRP:CG	1:A:445:PRO:HA	2.55	0.42
1:A:33:VAL:HG22	1:A:170:LEU:HD22	2.01	0.41
1:A:352:THR:HG23	1:A:369:TYR:H	1.86	0.41
1:A:403:LEU:C	1:A:403:LEU:HD23	2.45	0.41
5:F:1:NAG:H61	5:F:6:FUC:H3	2.01	0.41
1:A:456:ALA:N	1:A:457:PRO:CD	2.84	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:ASP:O	1:A:140:GLU:OE2[6_444]	1.97	0.23

5.3 Torsion angles

5.3.1 Protein backbone

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

of similar resolution.

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	448/447 (100%)	433 (97%)	15 (3%)	0	100	100

There are no Ramachandran outliers to report.

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	399/396 (101%)	388 (97%)	11 (3%)	38	50

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	34	GLU
1	A	49	SER
1	A	91	PRO
1	A	145	VAL
1	A	203	VAL
1	A	321	ARG
1	A	324	VAL
1	A	349	SER
1	A	364	ASP
1	A	377	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	25	GLN
1	A	143	HIS
1	A	437	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

19 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	B	1	2,1	14,14,15	0.45	0	17,19,21	0.84	0
2	NAG	B	2	2	14,14,15	0.42	0	17,19,21	0.71	0
2	BMA	B	3	2	11,11,12	0.48	0	15,15,17	0.83	1 (6%)
2	MAN	B	4	2	11,11,12	0.74	0	15,15,17	1.00	1 (6%)
3	NAG	C	1	3,1	14,14,15	0.40	0	17,19,21	1.17	1 (5%)
3	NAG	C	2	3	14,14,15	0.32	0	17,19,21	1.03	1 (5%)
3	FUC	C	3	3	10,10,11	0.71	0	14,14,16	1.53	3 (21%)
4	NAG	D	1	4,1	14,14,15	0.38	0	17,19,21	0.93	1 (5%)
4	NAG	D	2	4	14,14,15	0.41	0	17,19,21	0.79	0
4	BMA	D	3	4	11,11,12	0.76	0	15,15,17	0.85	0
3	NAG	E	1	3,1	14,14,15	0.36	0	17,19,21	0.81	0
3	NAG	E	2	3	14,14,15	0.31	0	17,19,21	1.39	2 (11%)
3	FUC	E	3	3	10,10,11	0.48	0	14,14,16	0.62	0
5	NAG	F	1	5,1	14,14,15	0.71	1 (7%)	17,19,21	2.05	3 (17%)
5	NAG	F	2	5	14,14,15	0.38	0	17,19,21	1.06	2 (11%)
5	BMA	F	3	5	11,11,12	0.64	0	15,15,17	0.93	0
5	MAN	F	4	5	11,11,12	0.68	0	15,15,17	0.78	1 (6%)
5	MAN	F	5	5	11,11,12	0.76	0	15,15,17	1.55	2 (13%)
5	FUC	F	6	5	10,10,11	0.67	0	14,14,16	1.55	3 (21%)

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	B	1	2,1	-	1/6/23/26	0/1/1/1
2	NAG	B	2	2	-	0/6/23/26	0/1/1/1
2	BMA	B	3	2	-	0/2/19/22	0/1/1/1
2	MAN	B	4	2	-	0/2/19/22	1/1/1/1
3	NAG	C	1	3,1	-	1/6/23/26	0/1/1/1
3	NAG	C	2	3	-	0/6/23/26	0/1/1/1
3	FUC	C	3	3	-	-	0/1/1/1
4	NAG	D	1	4,1	-	0/6/23/26	0/1/1/1
4	NAG	D	2	4	-	2/6/23/26	0/1/1/1
4	BMA	D	3	4	-	0/2/19/22	0/1/1/1
3	NAG	E	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	E	2	3	-	2/6/23/26	0/1/1/1
3	FUC	E	3	3	-	-	0/1/1/1
5	NAG	F	1	5,1	-	0/6/23/26	0/1/1/1
5	NAG	F	2	5	-	0/6/23/26	0/1/1/1
5	BMA	F	3	5	-	2/2/19/22	0/1/1/1
5	MAN	F	4	5	-	1/2/19/22	0/1/1/1
5	MAN	F	5	5	-	1/2/19/22	0/1/1/1
5	FUC	F	6	5	-	-	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	F	1	NAG	O6-C6	-2.19	1.33	1.42

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	F	1	NAG	C1-O5-C5	6.68	121.14	112.19
5	F	5	MAN	C1-O5-C5	4.30	117.95	112.19
3	E	2	NAG	C1-C2-N2	4.11	116.91	110.43
5	F	1	NAG	C6-C5-C4	-3.71	103.92	113.02
3	C	3	FUC	C1-C2-C3	3.30	114.44	109.64
3	E	2	NAG	C1-O5-C5	3.28	116.59	112.19
5	F	6	FUC	C1-O5-C5	3.12	120.33	112.97
5	F	6	FUC	C1-C2-C3	3.12	114.18	109.64
3	C	1	NAG	C1-C2-N2	3.06	115.25	110.43
2	B	4	MAN	C1-O5-C5	2.96	116.16	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	2	NAG	C2-N2-C7	2.76	126.60	122.90
5	F	5	MAN	O2-C2-C1	2.73	115.47	109.22
3	C	3	FUC	C3-C4-C5	2.65	113.84	109.81
3	C	3	FUC	C1-O5-C5	2.57	119.03	112.97
5	F	2	NAG	O3-C3-C2	-2.56	104.08	109.40
5	F	6	FUC	O5-C1-C2	2.39	116.50	110.79
5	F	1	NAG	C3-C4-C5	2.26	114.33	110.23
4	D	1	NAG	C1-O5-C5	2.08	114.97	112.19
5	F	4	MAN	O2-C2-C3	2.02	114.34	110.15
2	B	3	BMA	C1-O5-C5	2.01	114.88	112.19
5	F	2	NAG	O5-C5-C4	-2.00	105.96	110.83

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	E	2	NAG	O5-C5-C6-O6
3	E	2	NAG	C4-C5-C6-O6
3	C	1	NAG	O5-C5-C6-O6
5	F	3	BMA	C4-C5-C6-O6
5	F	3	BMA	O5-C5-C6-O6
4	D	2	NAG	C8-C7-N2-C2
5	F	5	MAN	C4-C5-C6-O6
2	B	1	NAG	C1-C2-N2-C7
4	D	2	NAG	O7-C7-N2-C2
5	F	4	MAN	O5-C5-C6-O6

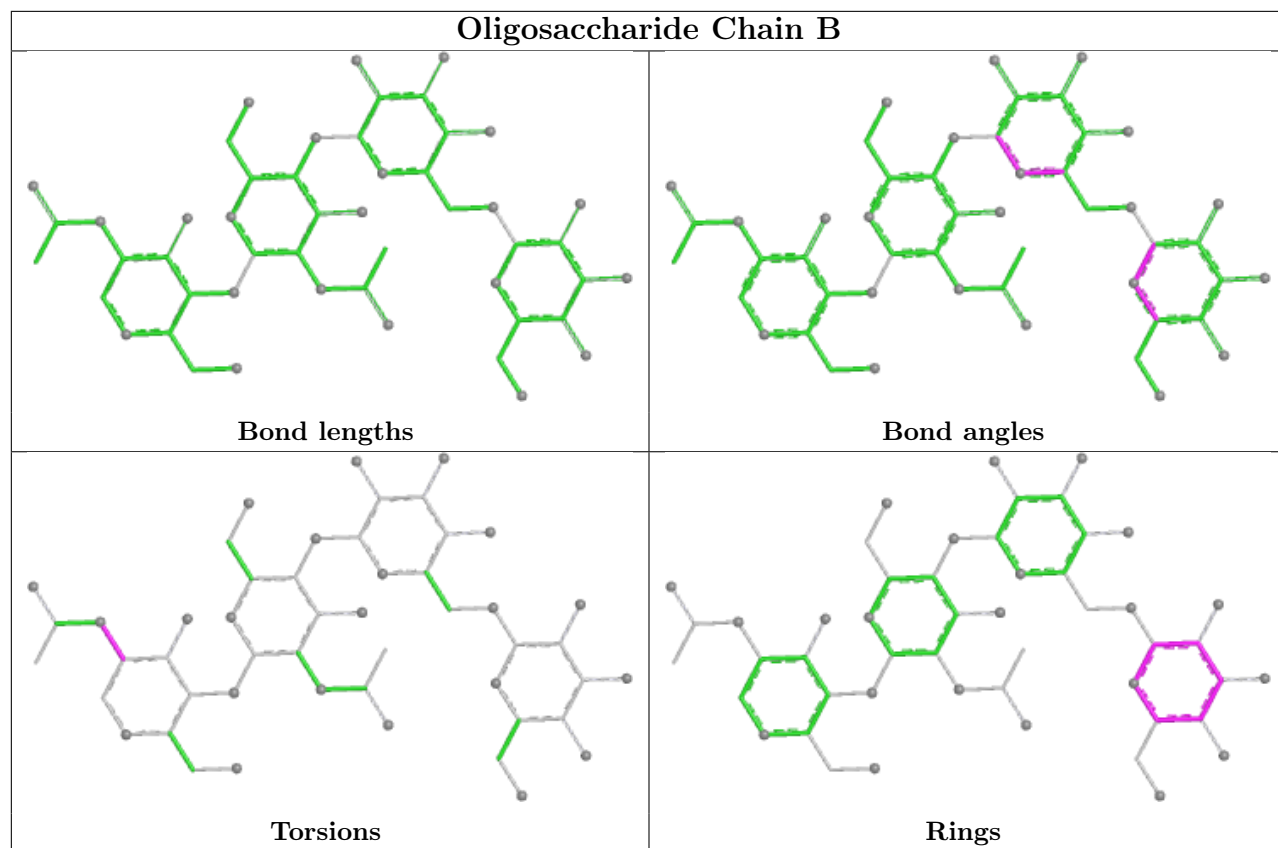
All (1) ring outliers are listed below:

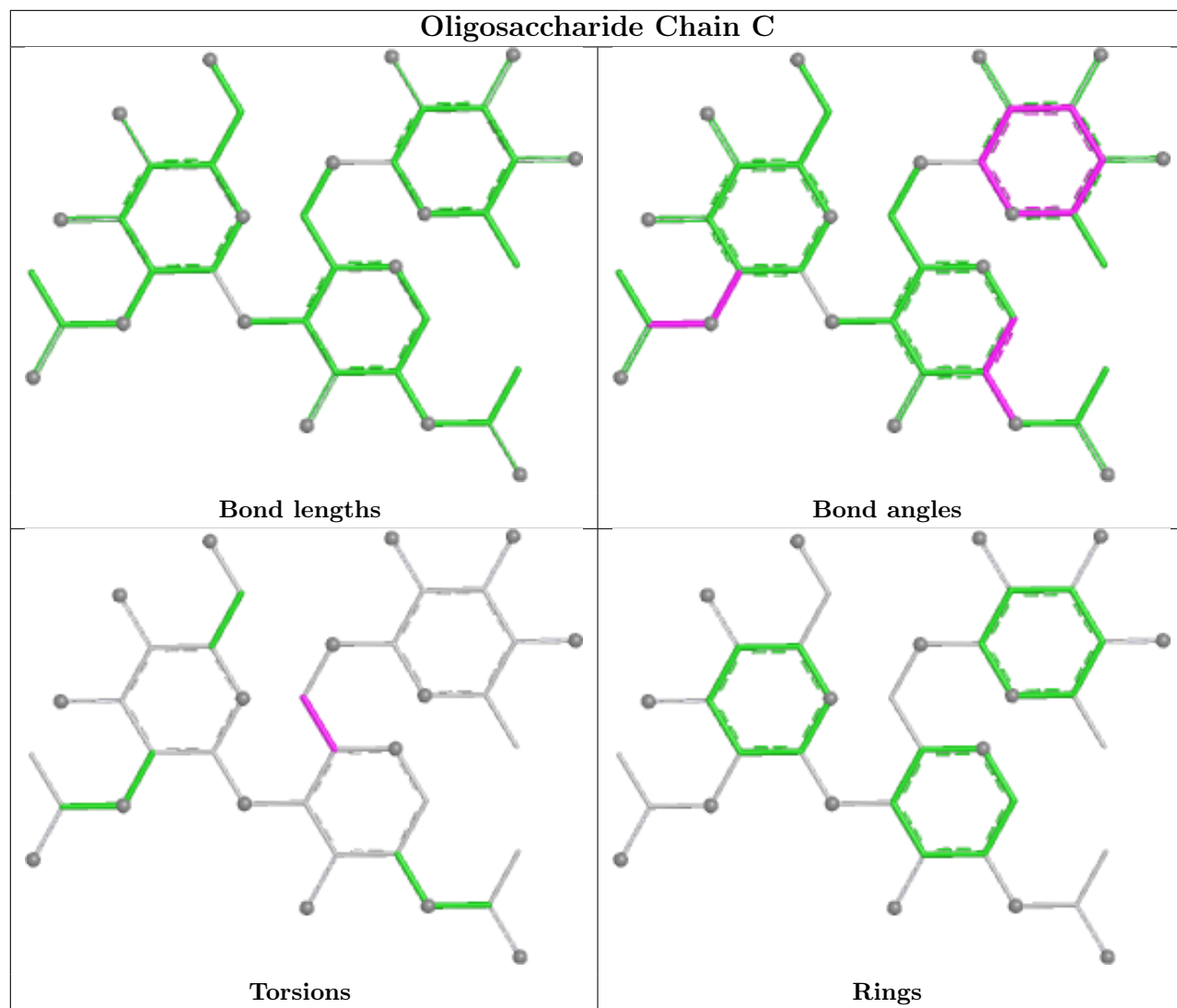
Mol	Chain	Res	Type	Atoms
2	B	4	MAN	C1-C2-C3-C4-C5-O5

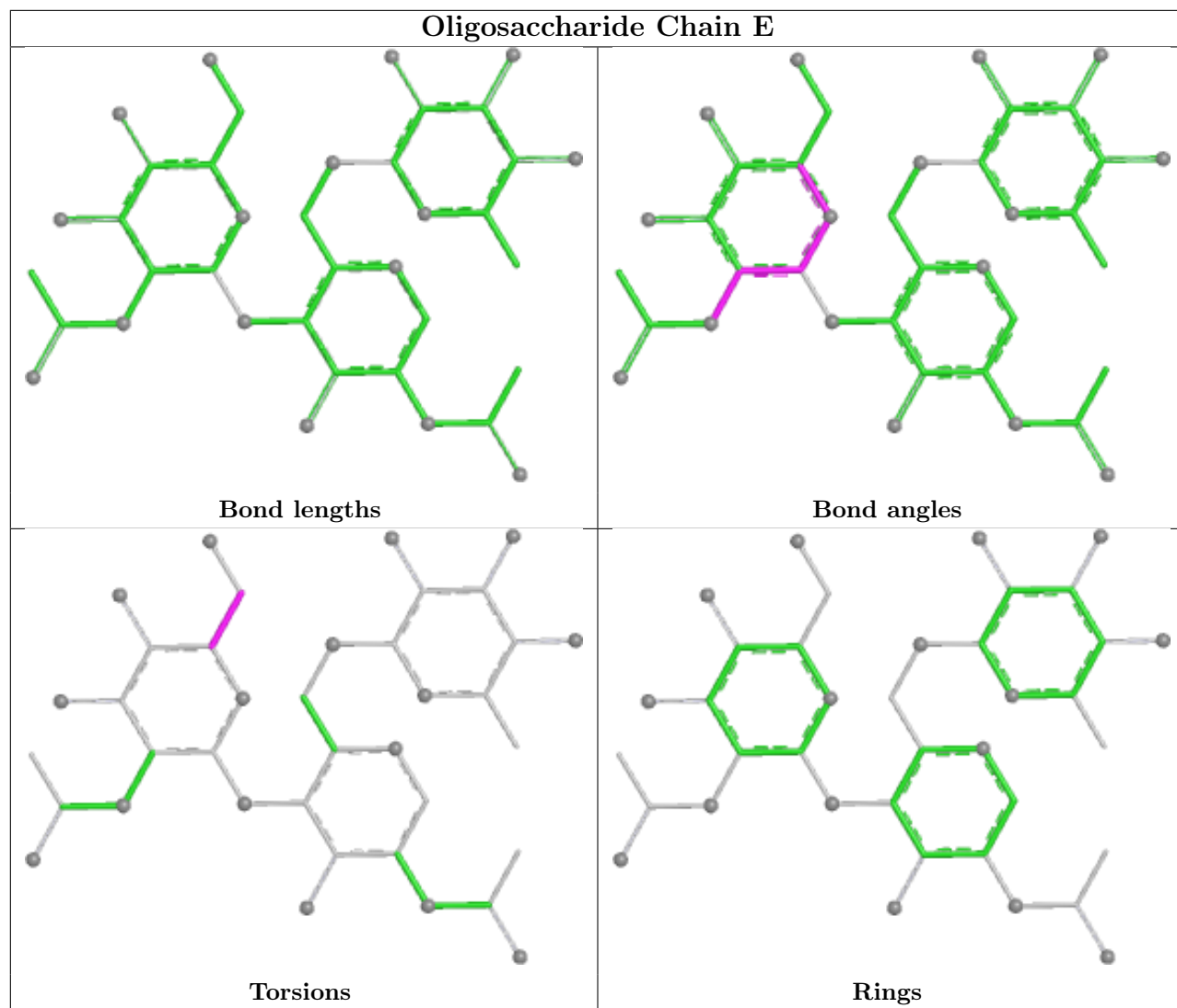
4 monomers are involved in 4 short contacts:

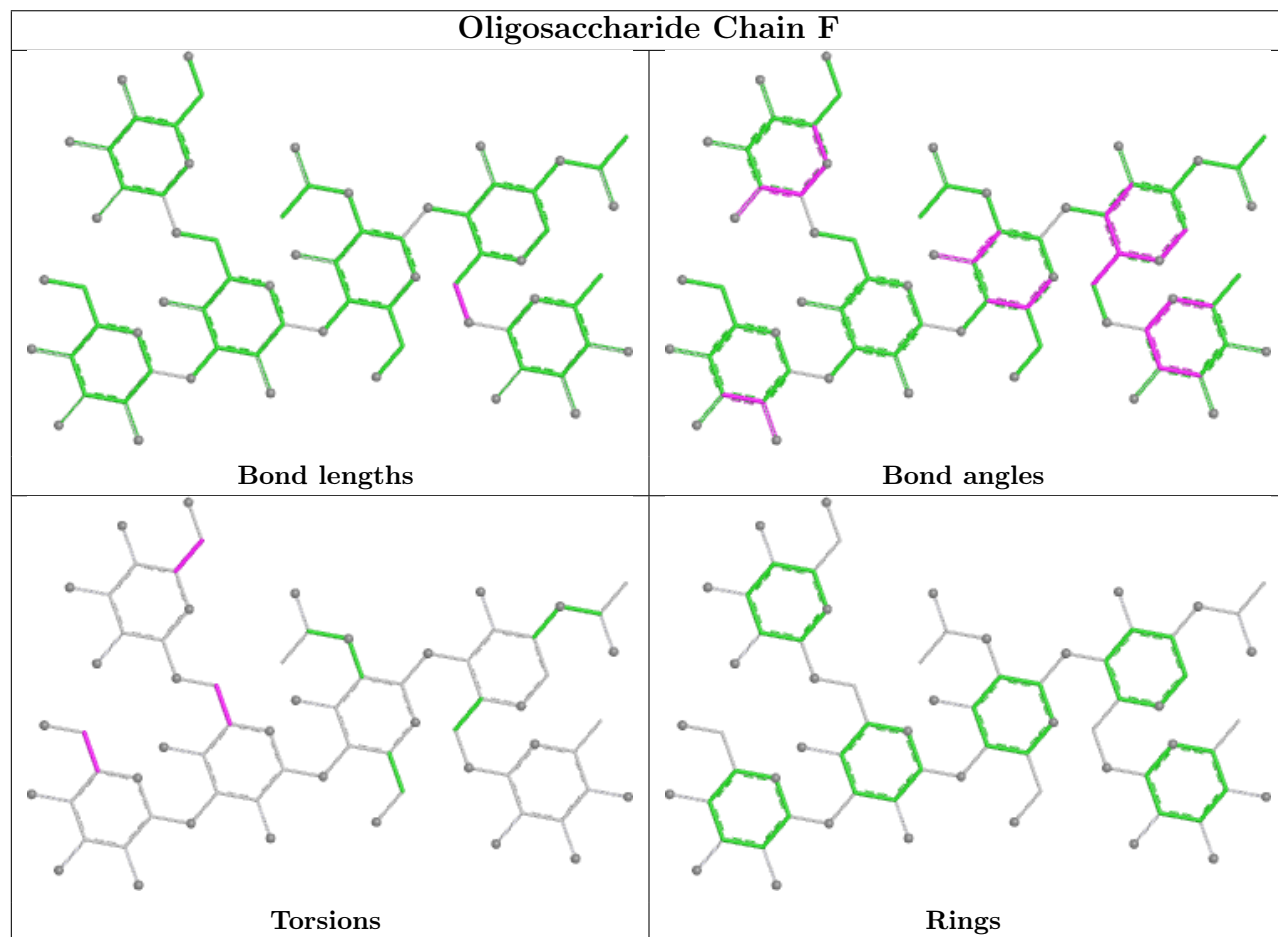
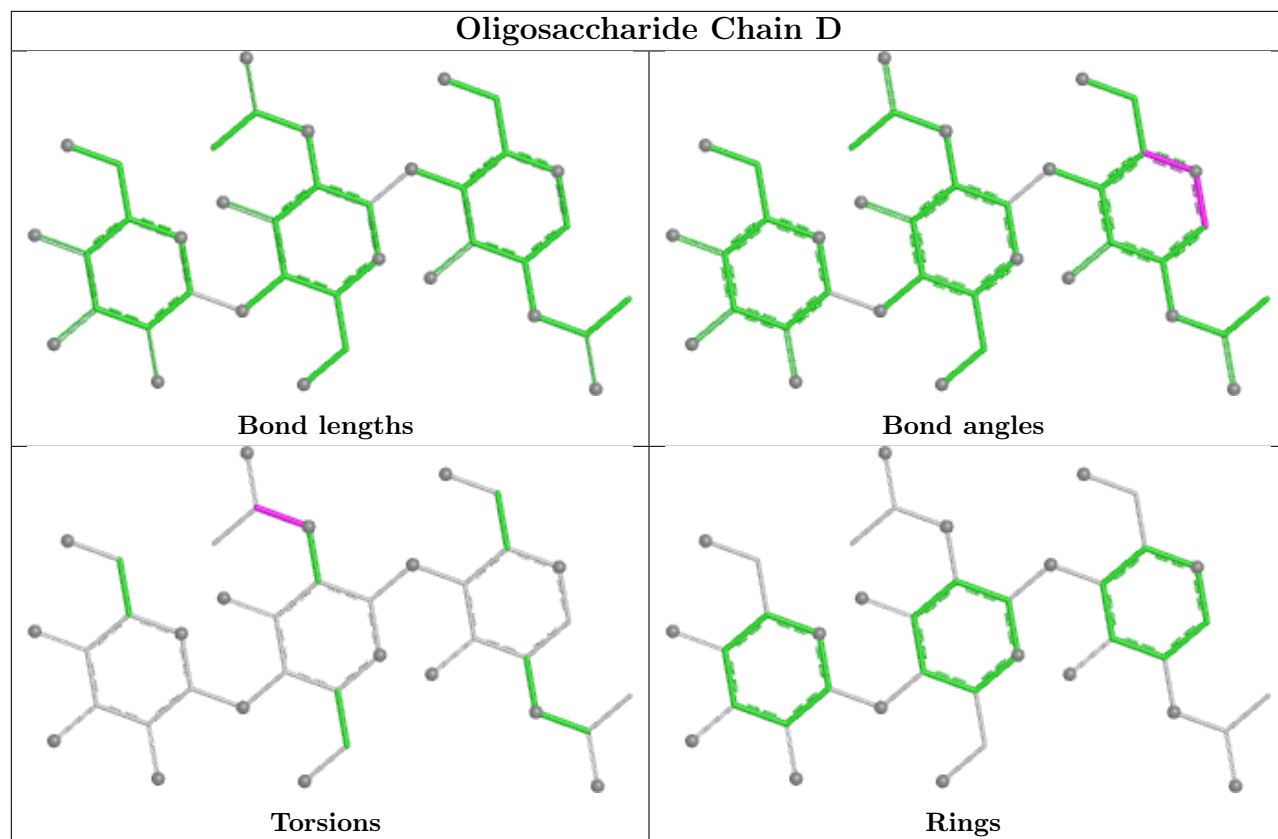
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	1	NAG	2	0
5	F	6	FUC	1	0
3	E	3	FUC	2	0
5	F	2	NAG	1	0

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









5.6 Ligand geometry

Of 10 ligands modelled in this entry, 5 are monoatomic - leaving 5 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
7	GOL	A	505	-	5,5,5	0.38	0	5,5,5	0.46	0
6	NAG	A	501	1	14,14,15	0.62	0	17,19,21	1.55	4 (23%)
7	GOL	A	503	-	5,5,5	0.14	0	5,5,5	0.32	0
7	GOL	A	504	-	5,5,5	0.12	0	5,5,5	0.39	0
7	GOL	A	502	-	5,5,5	0.22	0	5,5,5	0.52	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
7	GOL	A	505	-	-	4/4/4/4	-
6	NAG	A	501	1	-	5/6/23/26	0/1/1/1
7	GOL	A	503	-	-	2/4/4/4	-
7	GOL	A	504	-	-	4/4/4/4	-
7	GOL	A	502	-	-	2/4/4/4	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	501	NAG	O5-C1-C2	3.36	116.49	111.29
6	A	501	NAG	C1-O5-C5	3.00	116.21	112.19
6	A	501	NAG	C2-N2-C7	2.83	126.70	122.90
6	A	501	NAG	C1-C2-N2	2.46	114.30	110.43

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	502	GOL	C1-C2-C3-O3
7	A	502	GOL	O2-C2-C3-O3
7	A	503	GOL	C1-C2-C3-O3
7	A	503	GOL	O2-C2-C3-O3
7	A	504	GOL	O1-C1-C2-C3
7	A	505	GOL	O1-C1-C2-C3
7	A	505	GOL	C1-C2-C3-O3
6	A	501	NAG	C4-C5-C6-O6
6	A	501	NAG	O5-C5-C6-O6
6	A	501	NAG	O7-C7-N2-C2
6	A	501	NAG	C8-C7-N2-C2
7	A	504	GOL	C1-C2-C3-O3
7	A	504	GOL	O2-C2-C3-O3
7	A	505	GOL	O1-C1-C2-O2
7	A	505	GOL	O2-C2-C3-O3
7	A	504	GOL	O1-C1-C2-O2
6	A	501	NAG	C1-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	A	505	GOL	3	0
7	A	502	GOL	1	0

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	447/447 (100%)	0.13	29 (6%)	25 22	25, 48, 88, 145	3 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	201	LEU	6.5
1	A	203	VAL	5.3
1	A	202	GLY	4.1
1	A	82	ASP	4.1
1	A	209	GLY	4.0
1	A	79	TYR	3.5
1	A	471	ARG	3.4
1	A	208	PHE	3.4
1	A	323	MET	3.4
1	A	319	VAL	3.1
1	A	83	GLY	3.1
1	A	320	ALA	3.0
1	A	200	PHE	3.0
1	A	199	THR	2.9
1	A	308	ASP	2.9
1	A	210	GLU	2.8
1	A	118	ARG	2.7
1	A	321	ARG	2.6
1	A	126	VAL	2.6
1	A	207	SER	2.6
1	A	172	PRO	2.4
1	A	48	VAL	2.3
1	A	198	LYS	2.2
1	A	205	GLN	2.1
1	A	204	GLY	2.1
1	A	51	PRO	2.1
1	A	124[A]	GLN	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	331	GLN	2.0
1	A	115	PRO	2.0

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

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

6.3 Carbohydrates [i](#)

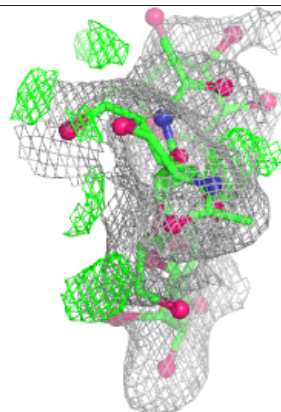
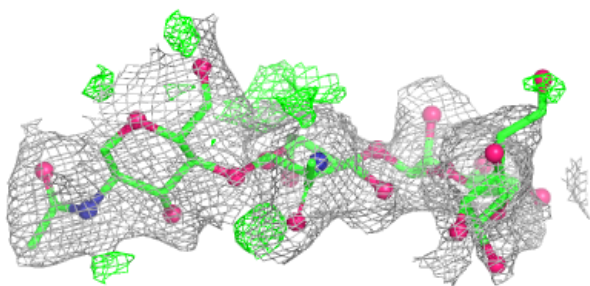
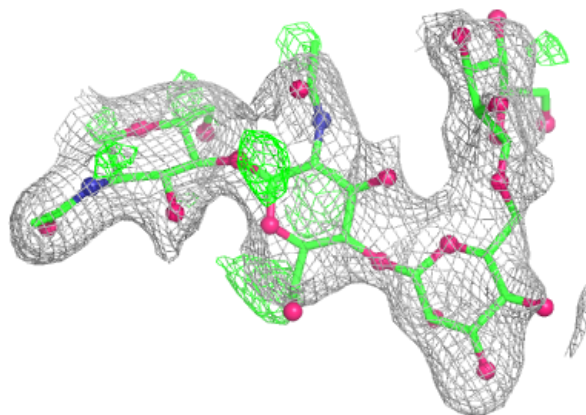
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q<0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	BMA	D	3	11/12	0.29	0.16	121,142,157,170	0
2	BMA	B	3	11/12	0.57	0.12	117,152,166,168	0
3	NAG	E	2	14/15	0.64	0.16	106,145,158,161	0
3	NAG	C	2	14/15	0.68	0.16	108,125,131,131	0
2	MAN	B	4	11/12	0.69	0.17	127,150,159,162	0
2	NAG	B	2	14/15	0.76	0.16	97,136,156,174	0
3	FUC	C	3	10/11	0.76	0.27	103,122,144,148	0
3	NAG	C	1	14/15	0.80	0.15	77,98,118,139	0
3	NAG	E	1	14/15	0.81	0.14	92,106,119,132	0
4	NAG	D	2	14/15	0.86	0.12	87,105,129,144	0
5	MAN	F	5	11/12	0.86	0.15	78,92,99,100	0
3	FUC	E	3	10/11	0.89	0.15	87,110,118,119	0
2	NAG	B	1	14/15	0.92	0.10	45,74,96,121	0
5	NAG	F	1	14/15	0.94	0.08	46,50,59,61	0
5	NAG	F	2	14/15	0.95	0.08	47,53,61,62	0
5	BMA	F	3	11/12	0.95	0.07	44,57,64,75	0
4	NAG	D	1	14/15	0.95	0.10	48,61,71,80	0
5	MAN	F	4	11/12	0.96	0.07	46,56,60,63	0
5	FUC	F	6	10/11	0.96	0.07	45,52,56,56	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.

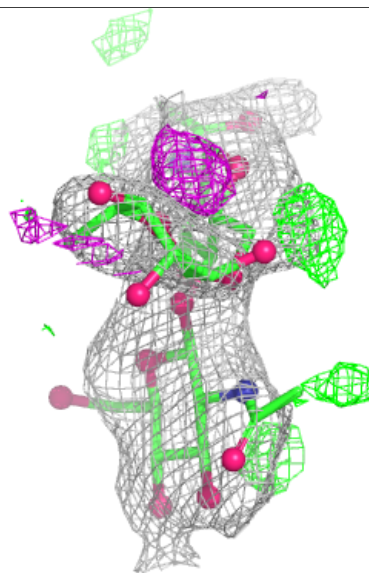
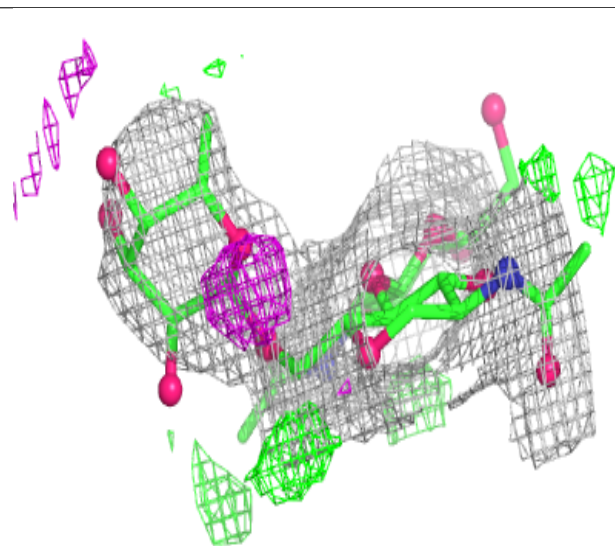
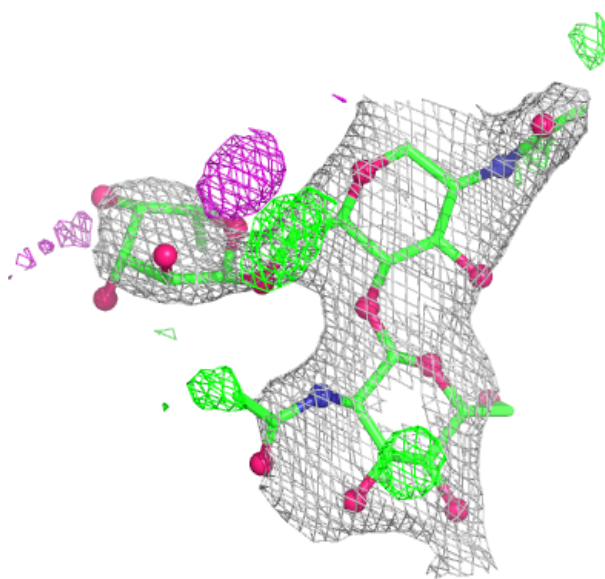
Electron density around Chain B:

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



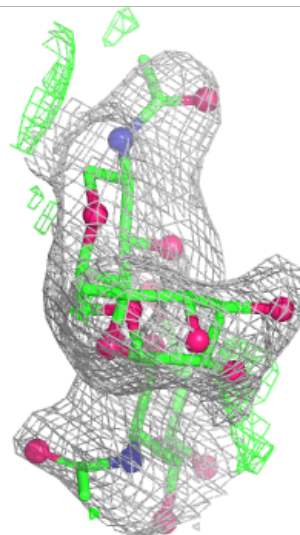
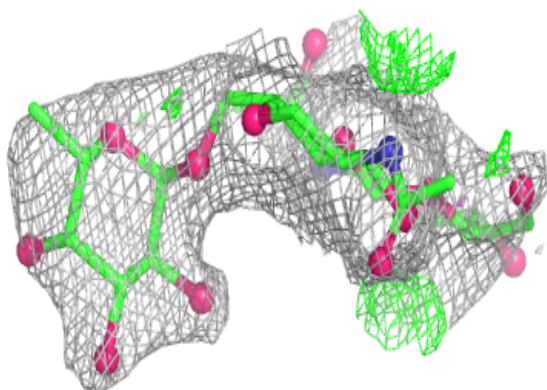
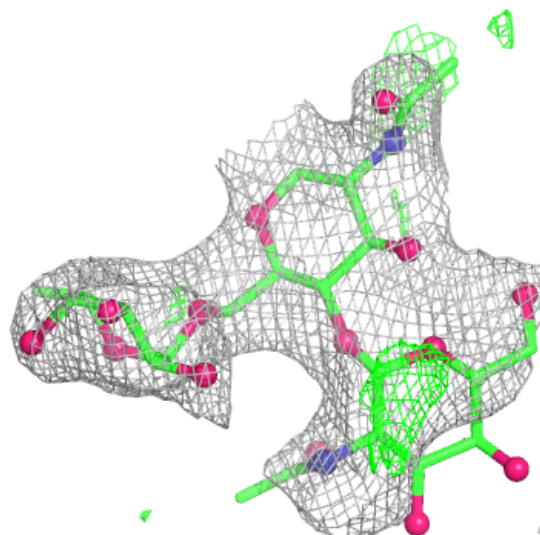
Electron density around Chain C:

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



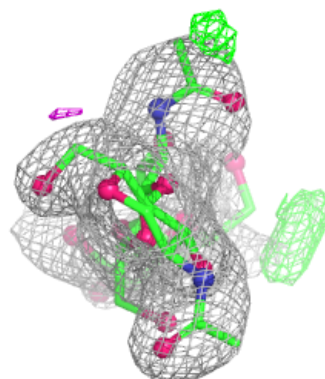
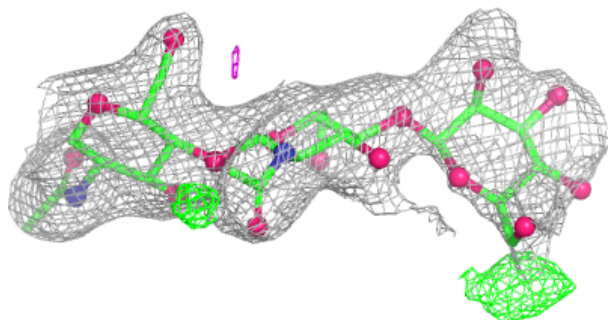
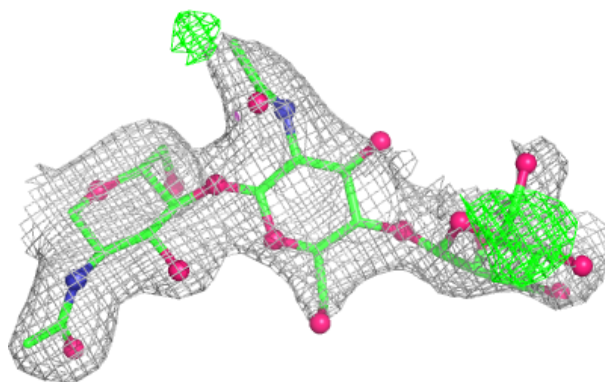
Electron density around Chain E:

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

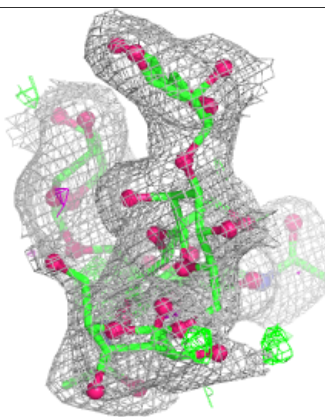
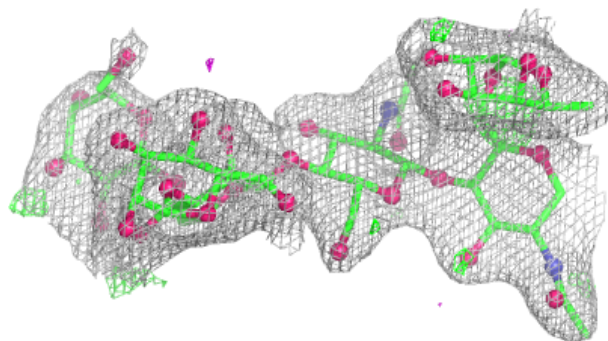
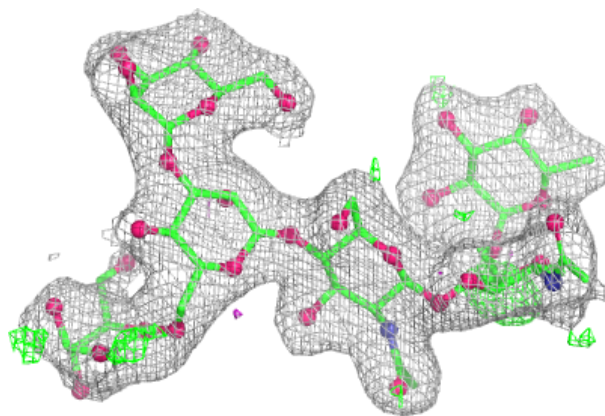


Electron density around Chain D:

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

**Electron density around Chain F:**

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



6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NAG	A	501	14/15	0.71	0.20	91,130,144,147	0
7	GOL	A	502	6/6	0.80	0.18	85,90,99,100	0
7	GOL	A	503	6/6	0.94	0.10	44,65,77,87	0
7	GOL	A	504	6/6	0.94	0.13	49,67,77,88	0
7	GOL	A	505	6/6	0.94	0.12	32,61,76,86	0
8	ZN	A	508	1/1	0.99	0.04	54,54,54,54	0
8	ZN	A	509	1/1	0.99	0.04	51,51,51,51	0
8	ZN	A	510	1/1	0.99	0.04	54,54,54,54	1
8	ZN	A	507	1/1	1.00	0.02	34,34,34,34	0
8	ZN	A	506	1/1	1.00	0.04	34,34,34,34	0

6.5 Other polymers [i](#)

There are no such residues in this entry.