



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 06:23 PM UTC

PDB ID : 5M8Q / pdb_00005m8q
Title : Crystal structure of human tyrosinase related protein 1 mutant (T391V-R374S-Y362F) in complex with kojic acid
Authors : Lai, X.; Soler-Lopez, M.; Wichers, H.J.; Dijkstra, B.W.
Deposited on : 2016-10-29
Resolution : 2.85 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

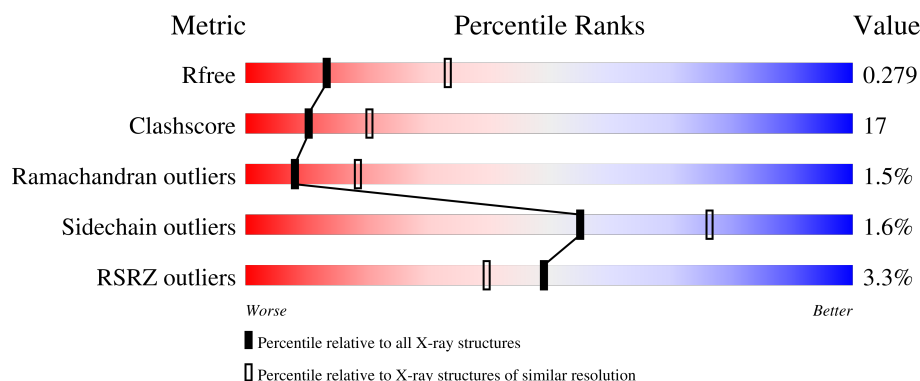
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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.



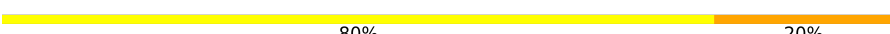
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	446	% 67% 32% .
1	B	446	11% 54% 40% 6% .
1	C	446	% 72% 26% .
1	D	446	74% 25% .
2	E	3	67% 33%

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Mol	Chain	Length	Quality of chain
3	F	2	 50% 50%
3	H	2	 100%
3	K	2	 50% 50%
3	O	2	 100%
4	G	5	 80% 20%
5	I	2	 100%
5	J	2	 50% 50%
5	M	2	 100%
5	N	2	 100%
6	L	5	 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
9	KOJ	B	508	-	X	-	-
9	KOJ	C	514	-	X	-	-
9	KOJ	D	510	-	X	-	-

2 Entry composition [i](#)

There are 10 unique types of molecules in this entry. The entry contains 14735 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.

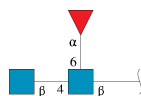
- Molecule 1 is a protein called 5,6-dihydroxyindole-2-carboxylic acid oxidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	446	Total	C	N	O	S	0	0	0
			3554	2231	629	671	23			
1	B	446	Total	C	N	O	S	0	0	0
			3554	2231	629	671	23			
1	C	446	Total	C	N	O	S	0	0	0
			3554	2231	629	671	23			
1	D	446	Total	C	N	O	S	0	0	0
			3554	2231	629	671	23			

There are 12 discrepancies between the modelled and reference sequences:

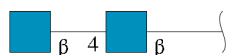
Chain	Residue	Modelled	Actual	Comment	Reference
A	362	PHE	TYR	engineered mutation	UNP P17643
A	374	SER	ARG	engineered mutation	UNP P17643
A	391	VAL	THR	engineered mutation	UNP P17643
B	362	PHE	TYR	engineered mutation	UNP P17643
B	374	SER	ARG	engineered mutation	UNP P17643
B	391	VAL	THR	engineered mutation	UNP P17643
C	362	PHE	TYR	engineered mutation	UNP P17643
C	374	SER	ARG	engineered mutation	UNP P17643
C	391	VAL	THR	engineered mutation	UNP P17643
D	362	PHE	TYR	engineered mutation	UNP P17643
D	374	SER	ARG	engineered mutation	UNP P17643
D	391	VAL	THR	engineered mutation	UNP P17643

- Molecule 2 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.



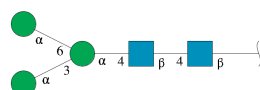
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			38	22	2	14			

- Molecule 3 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
3	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	H	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	O	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-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	G	5	Total	C	N	O	0	0	0
			61	34	2	25			

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



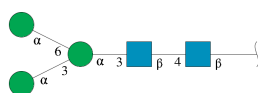
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	2	Total	C	N	O	0	0	0
			24	14	1	9			

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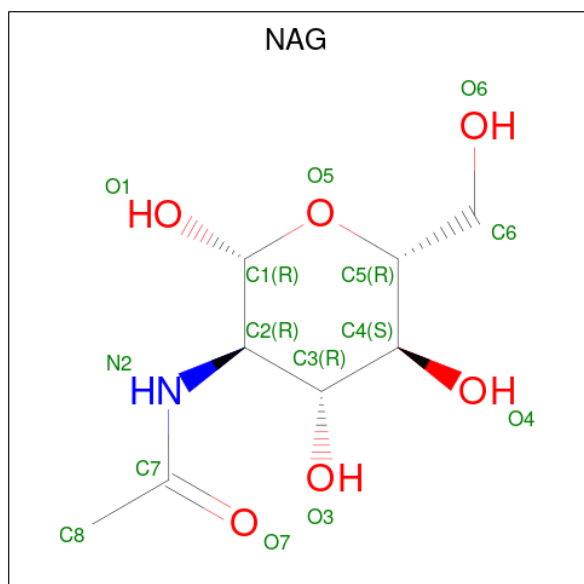
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	J	2	Total	C	N	O	0	0	0
			24	14	1	9			
5	M	2	Total	C	N	O	0	0	0
			24	14	1	9			
5	N	2	Total	C	N	O	0	0	0
			24	14	1	9			

- Molecule 6 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(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
6	L	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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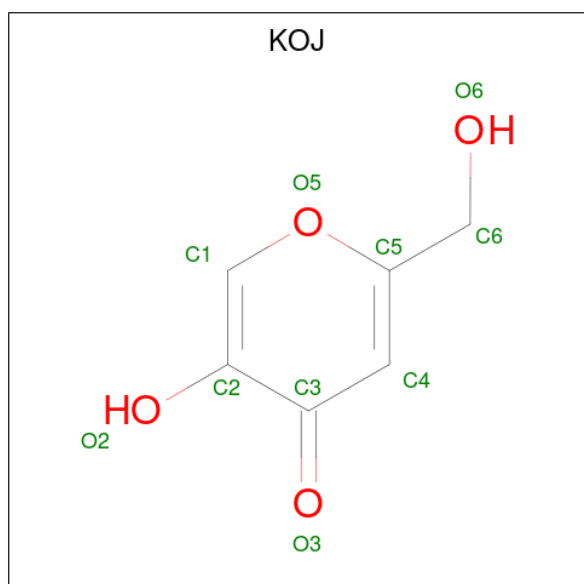
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	B	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		
7	C	1	Total	C	N	O	0	0
			14	8	1	5		
7	D	1	Total	C	N	O	0	0
	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

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	2	Total	Zn	0	0
			2	2		
8	B	2	Total	Zn	0	0
			2	2		
8	C	3	Total	Zn	0	0
			3	3		
8	D	2	Total	Zn	0	0
			2	2		

- Molecule 9 is 5-HYDROXY-2-(HYDROXYMETHYL)-4H-PYRAN-4-ONE (CCD ID: KOJ) (formula: C₆H₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
9	A	1	Total	C	O	0	0
			10	6	4		
9	B	1	Total	C	O	0	0
			10	6	4		
9	C	1	Total	C	O	0	0
			10	6	4		
9	D	1	Total	C	O	0	0
			10	6	4		

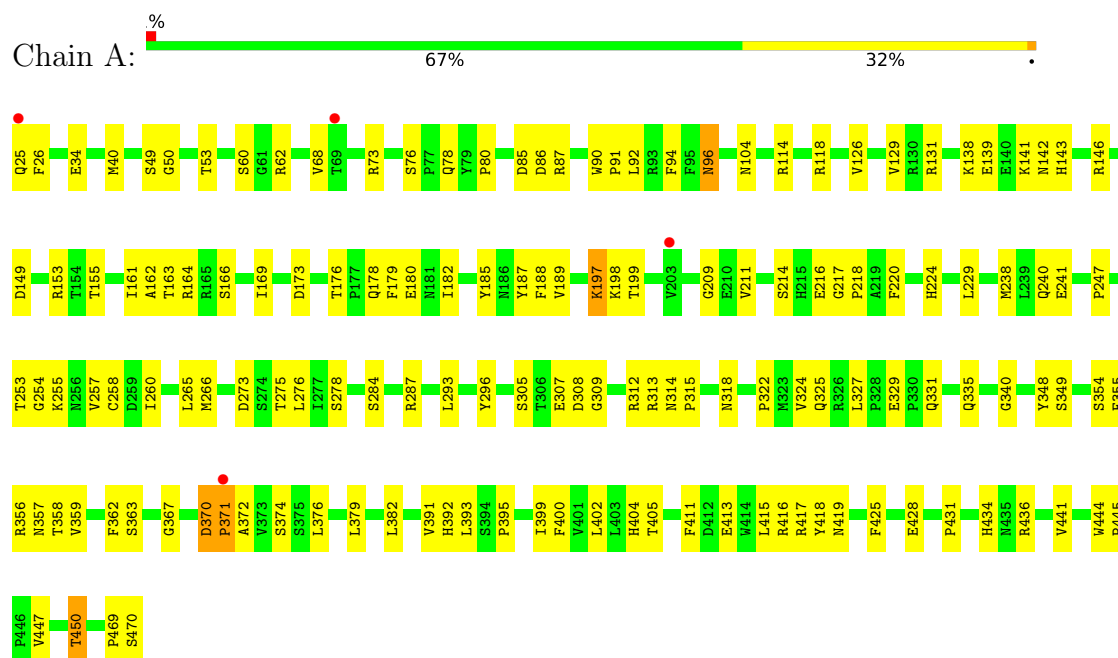
- Molecule 10 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
10	A	1	Total	O	0	0
			1	1		
10	B	1	Total	O	0	0
			1	1		
10	C	1	Total	O	0	0
			1	1		
10	D	1	Total	O	0	0
			1	1		

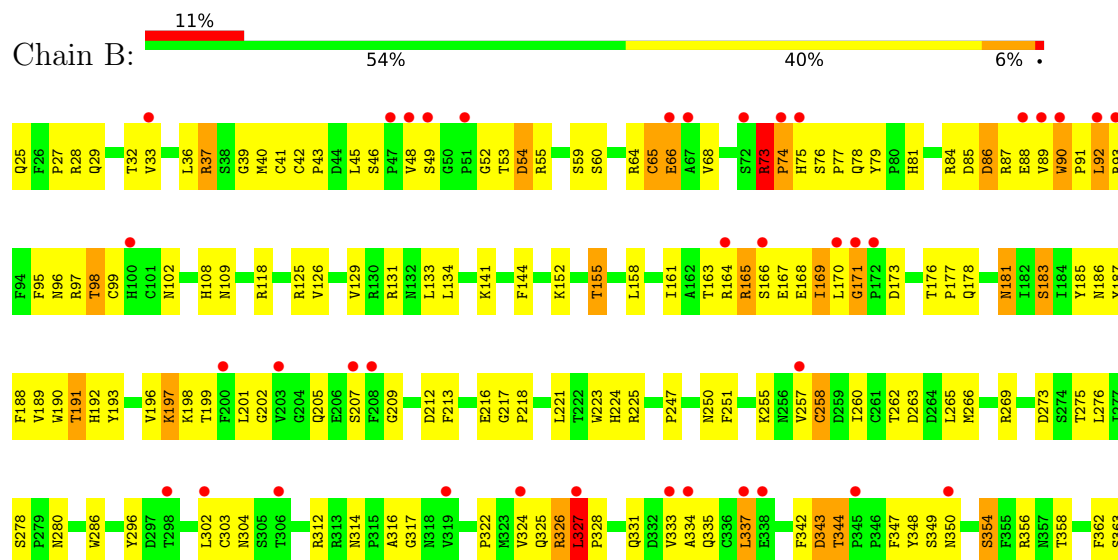
3 Residue-property plots

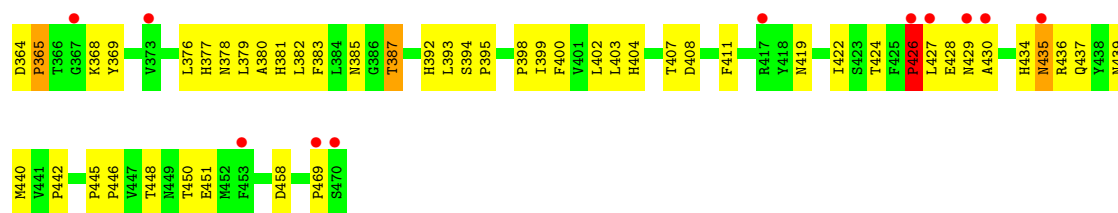
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: 5,6-dihydroxyindole-2-carboxylic acid oxidase

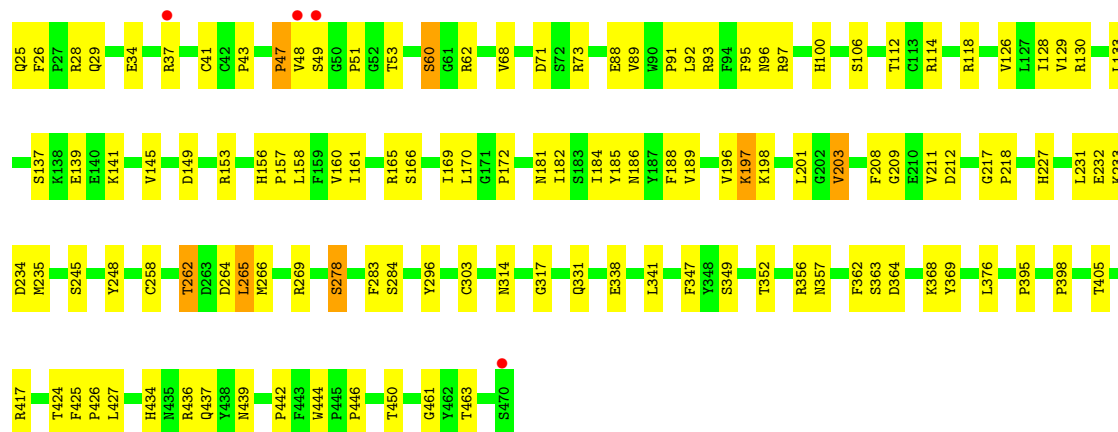
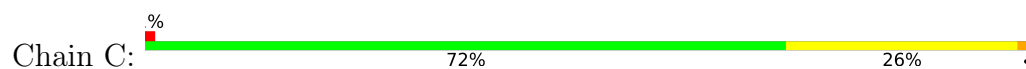


- Molecule 1: 5,6-dihydroxyindole-2-carboxylic acid oxidase

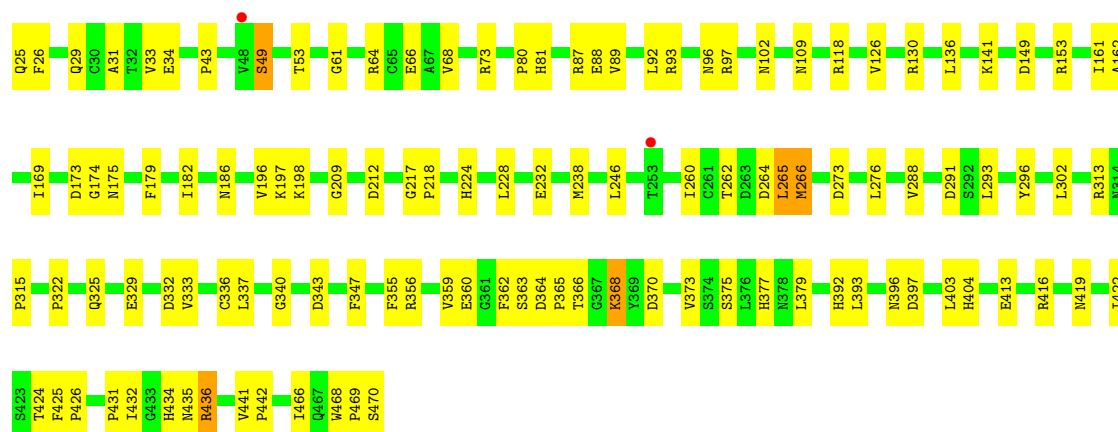




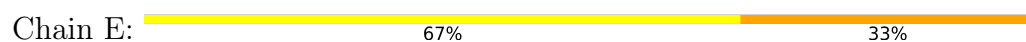
• Molecule 1: 5,6-dihydroxyindole-2-carboxylic acid oxidase



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



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

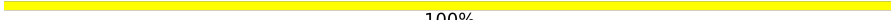


- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  50% 50%

 NAG1
NAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  100%

 NAG1
NAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  50% 50%

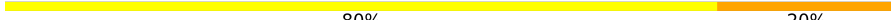
 NAG1
NAG2

- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  100%

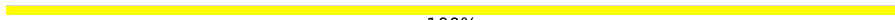
 NAG1
NAG2

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

Chain G:  80% 20%

 NAG1
NAG2
MAN3
MAN4
MAN5

- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  100%

 NAG1
FUC2

- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  50% 50%



MAG1
FUC2

- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  100%



MAG1
FUC2

- Molecule 5: alpha-L-fucopyranose-(1-6)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  100%



MAG1
FUC2

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

Chain L:  100%



MAG1
MAG2
MAN3
MAN4
MAN5

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	89.64Å 140.43Å 191.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.88 – 2.85 47.88 – 2.85	Depositor EDS
% Data completeness (in resolution range)	100.0 (47.88-2.85) 92.9 (47.88-2.85)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.19 (at 2.86Å)	Xtriage
Refinement program	REFMAC, PHENIX	Depositor
R, R_{free}	0.208 , 0.275 0.217 , 0.279	Depositor DCC
R_{free} test set	2855 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	53.5	Xtriage
Anisotropy	0.474	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 40.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14735	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.04% 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: KOJ, MAN, FUC, NAG, 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	0.55	3/3661 (0.1%)	0.75	4/4990 (0.1%)
1	B	0.60	6/3661 (0.2%)	0.94	14/4990 (0.3%)
1	C	0.42	0/3661	0.68	3/4990 (0.1%)
1	D	0.45	0/3661	0.65	0/4990
All	All	0.51	9/14644 (0.1%)	0.76	21/19960 (0.1%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	B	0	4
1	C	0	1
All	All	0	5

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	73	ARG	NE-CZ	-10.17	1.21	1.33
1	A	371	PRO	CA-CB	9.75	1.67	1.53
1	B	86	ASP	CB-CG	-6.79	1.35	1.52
1	B	73	ARG	CZ-NH2	-6.66	1.24	1.33
1	B	169	ILE	CG1-CD1	-6.08	1.28	1.51
1	B	73	ARG	CD-NE	-5.83	1.38	1.46
1	A	370	ASP	N-CA	5.43	1.53	1.46
1	B	73	ARG	CZ-NH1	-5.18	1.25	1.32
1	A	371	PRO	N-CD	5.09	1.54	1.47

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	327	LEU	CA-CB-CG	9.61	149.94	116.30
1	B	37	ARG	NE-CZ-NH2	9.28	127.55	119.20
1	B	337	LEU	CB-CG-CD1	-8.84	84.19	110.70
1	B	73	ARG	NE-CZ-NH2	8.61	126.95	119.20
1	B	169	ILE	CG1-CB-CG2	-7.21	89.06	110.70
1	B	98	THR	CA-C-N	-7.10	109.81	122.74
1	B	98	THR	C-N-CA	-7.10	109.81	122.74
1	A	96	ASN	CB-CA-C	7.08	121.76	109.15
1	A	96	ASN	CA-CB-CG	6.59	119.19	112.60
1	B	92	LEU	CB-CG-CD1	-6.18	92.16	110.70
1	B	344	THR	CA-C-N	-5.74	114.46	120.38
1	B	344	THR	C-N-CA	-5.74	114.46	120.38
1	A	96	ASN	N-CA-C	-5.74	106.92	114.04
1	A	370	ASP	CB-CG-OD1	5.58	131.24	118.40
1	B	90	TRP	C-N-CD	-5.56	108.36	120.60
1	B	37	ARG	CG-CD-NE	5.32	123.71	112.00
1	C	265	LEU	CA-CB-CG	5.26	134.71	116.30
1	C	262	THR	CA-C-N	-5.25	113.53	122.56
1	C	262	THR	C-N-CA	-5.25	113.53	122.56
1	B	165	ARG	CB-CA-C	5.12	118.15	109.50
1	B	165	ARG	CB-CG-CD	-5.07	99.65	111.30

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	426	PRO	Peptide
1	B	435	ASN	Peptide
1	B	65	CYS	Peptide
1	B	99	CYS	Peptide
1	C	48	VAL	Peptide

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	3554	0	3319	112	0
1	B	3554	0	3322	204	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	3554	0	3321	87	0
1	D	3554	0	3321	86	0
2	E	38	0	34	1	0
3	F	28	0	25	0	0
3	H	28	0	25	1	0
3	K	28	0	25	0	0
3	O	28	0	25	0	0
4	G	61	0	52	1	0
5	I	24	0	22	0	0
5	J	24	0	22	0	0
5	M	24	0	22	1	0
5	N	24	0	22	0	0
6	L	61	0	52	0	0
7	A	14	0	13	0	0
7	B	42	0	39	3	0
7	C	28	0	26	1	0
7	D	14	0	13	0	0
8	A	2	0	0	0	0
8	B	2	0	0	0	0
8	C	3	0	0	0	0
8	D	2	0	0	0	0
9	A	10	0	6	3	0
9	B	10	0	6	3	0
9	C	10	0	6	1	0
9	D	10	0	6	0	0
10	A	1	0	0	2	0
10	B	1	0	0	1	0
10	C	1	0	0	1	0
10	D	1	0	0	0	0
All	All	14735	0	13724	483	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 17.

All (483) 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:165:ARG:NH2	1:B:167:GLU:OE2	1.62	1.31
1:B:92:LEU:HD11	1:B:96:ASN:HA	1.24	1.20
1:B:327:LEU:HD12	1:B:380:ALA:HA	1.19	1.15
1:B:255:LYS:O	1:B:326:ARG:NH1	1.85	1.09

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:350:ASN:HA	1:B:369:TYR:CE2	2.02	0.94
1:B:165:ARG:NH2	1:B:167:GLU:CD	2.27	0.92
9:A:516:KOJ:O3	10:A:601:HOH:O	1.89	0.91
1:B:439:ASN:HD22	1:B:446:PRO:HB3	1.35	0.90
1:B:331:GLN:HA	1:B:334:ALA:HB3	1.54	0.87
1:B:85:ASP:OD1	1:B:86:ASP:N	2.06	0.87
1:B:337:LEU:N	1:B:337:LEU:HD12	1.90	0.85
1:B:327:LEU:HD13	1:B:383:PHE:HB3	1.58	0.85
1:B:66:GLU:OE1	1:B:98:THR:N	2.12	0.83
1:B:166:SER:HA	1:B:169:ILE:HD13	1.61	0.83
1:C:141:LYS:HD2	1:C:265:LEU:HD11	1.61	0.83
1:B:334:ALA:HA	1:B:337:LEU:HD11	1.62	0.82
1:C:338:GLU:OE1	1:C:417:ARG:NH2	2.12	0.81
1:A:254:GLY:HA3	1:A:329:GLU:OE2	1.81	0.81
1:B:342:PHE:H	1:B:424:THR:HG21	1.43	0.81
1:B:73:ARG:HB3	1:B:74:PRO:HD2	1.63	0.80
1:B:327:LEU:HD12	1:B:380:ALA:CA	2.07	0.80
1:B:337:LEU:HD12	1:B:337:LEU:H	1.42	0.79
1:B:378:ASN:ND2	9:B:508:KOJ:H4	1.96	0.79
1:B:439:ASN:ND2	1:B:446:PRO:HB3	1.97	0.79
1:B:75:HIS:NE2	1:B:79:TYR:HB3	1.97	0.79
1:D:425:PHE:O	1:D:436:ARG:NH1	2.16	0.79
1:B:66:GLU:N	1:B:66:GLU:OE2	2.16	0.79
1:C:89:VAL:HG12	1:C:92:LEU:HB2	1.64	0.78
1:B:218:PRO:HG3	1:B:426:PRO:HG2	1.65	0.78
9:B:508:KOJ:O3	10:B:601:HOH:O	2.00	0.77
1:A:49:SER:HB2	1:A:53:THR:HG21	1.66	0.77
1:A:436:ARG:HG2	1:A:450:THR:HG22	1.65	0.77
1:D:364:ASP:OD1	1:D:368:LYS:HB3	1.84	0.77
1:B:326:ARG:NH2	1:B:327:LEU:HB2	1.99	0.76
1:B:92:LEU:HD11	1:B:96:ASN:CA	2.11	0.76
1:C:73:ARG:HH22	1:D:97:ARG:HH22	1.33	0.75
1:C:28:ARG:NH1	1:C:157:PRO:O	2.20	0.74
1:B:28:ARG:NH2	1:B:181:ASN:OD1	2.20	0.74
1:D:43:PRO:HG2	1:D:109:ASN:HB3	1.68	0.73
1:B:216:GLU:HG2	1:B:362:PHE:HE2	1.52	0.73
1:C:25:GLN:HG2	1:C:161:ILE:HA	1.70	0.73
1:D:425:PHE:HB3	1:D:436:ARG:HH11	1.52	0.73
1:C:363:SER:HB3	1:C:369:TYR:HA	1.71	0.73
1:B:29:GLN:HG3	1:B:43:PRO:HB3	1.72	0.72
1:B:25:GLN:HG2	1:B:161:ILE:HD12	1.72	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:327:LEU:CD1	1:B:380:ALA:HA	2.12	0.71
1:A:293:LEU:HD11	1:A:392:HIS:NE2	2.05	0.71
1:A:118:ARG:HG3	1:A:126:VAL:HG11	1.71	0.71
1:B:378:ASN:HD22	9:B:508:KOJ:H4	1.55	0.71
1:B:85:ASP:HB3	1:B:88:GLU:OE2	1.91	0.71
1:D:89:VAL:HG13	1:D:92:LEU:HB2	1.72	0.71
1:B:430:ALA:HB3	1:B:435:ASN:OD1	1.91	0.71
1:B:86:ASP:OD2	1:B:165:ARG:NH1	2.24	0.71
1:D:49:SER:HB2	1:D:53:THR:HG21	1.73	0.70
1:B:66:GLU:OE1	1:B:97:ARG:HA	1.91	0.70
1:B:76:SER:OG	1:B:78:GLN:OE1	2.08	0.70
1:B:76:SER:CB	1:B:78:GLN:OE1	2.40	0.70
1:A:25:GLN:HG2	1:A:161:ILE:HA	1.72	0.70
1:A:413:GLU:OE2	1:A:416:ARG:NH2	2.24	0.70
1:B:96:ASN:HD22	7:B:501:NAG:H83	1.57	0.70
1:B:358:THR:HA	1:B:363:SER:HB2	1.74	0.69
1:C:34:GLU:HG3	1:C:37:ARG:HH21	1.56	0.69
1:A:60:SER:HB2	1:A:62:ARG:HG3	1.75	0.69
1:D:413:GLU:OE2	1:D:416:ARG:NH1	2.26	0.69
1:B:327:LEU:HB3	1:B:328:PRO:HD2	1.75	0.68
1:A:163:THR:HG23	1:A:164:ARG:HG2	1.76	0.68
1:A:73:ARG:NH1	1:A:428:GLU:OE2	2.27	0.68
1:A:138:LYS:H	1:A:138:LYS:HD3	1.58	0.68
1:A:354:SER:O	1:A:358:THR:HG23	1.92	0.68
1:C:106:SER:HB2	1:C:114:ARG:HG2	1.76	0.68
1:B:201:LEU:HD12	1:B:205:GLN:OE1	1.94	0.67
1:B:165:ARG:NH1	1:B:166:SER:OG	2.28	0.67
1:B:250:ASN:O	1:B:250:ASN:ND2	2.29	0.66
1:B:428:GLU:CD	1:B:435:ASN:HD22	2.03	0.66
1:A:104:ASN:OD1	1:A:114:ARG:NH1	2.24	0.66
1:B:52:GLY:HA2	1:B:55:ARG:NH1	2.11	0.65
1:A:265:LEU:HD12	1:A:265:LEU:O	1.97	0.65
1:A:392:HIS:CD2	1:A:393:LEU:HG	2.31	0.65
1:C:112:THR:OG1	1:C:233:LYS:NZ	2.28	0.65
1:A:218:PRO:HD2	1:A:434:HIS:HB3	1.79	0.65
1:B:152:LYS:HD3	1:B:286:TRP:CD2	2.30	0.65
1:B:349:SER:O	1:B:369:TYR:HE2	1.80	0.64
1:D:25:GLN:HG3	1:D:26:PHE:H	1.62	0.64
1:B:161:ILE:HD11	1:B:190:TRP:CD2	2.31	0.64
1:B:257:VAL:HG12	1:B:258:CYS:H	1.62	0.64
1:B:33:VAL:HG13	1:B:177:PRO:HG3	1.78	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:64:ARG:NH2	1:D:66:GLU:OE2	2.31	0.64
1:C:29:GLN:HG3	1:C:43:PRO:HB3	1.79	0.64
1:C:218:PRO:HD2	1:C:434:HIS:HB3	1.79	0.64
1:D:260:ILE:HD12	1:D:266:MET:HG3	1.80	0.64
1:B:165:ARG:HD2	1:B:166:SER:H	1.63	0.64
1:D:61:GLY:O	1:D:102:ASN:ND2	2.20	0.64
1:B:65:CYS:HA	1:B:66:GLU:OE2	1.97	0.63
1:B:170:LEU:H	1:B:170:LEU:HD23	1.63	0.63
1:B:328:PRO:CG	1:B:376:LEU:HD23	2.28	0.63
1:B:75:HIS:CE1	1:B:76:SER:O	2.51	0.63
1:A:358:THR:HG22	1:A:363:SER:HB2	1.81	0.63
1:B:32:THR:O	1:B:36:LEU:HD23	1.98	0.63
1:B:87:ARG:HB2	1:B:213:PHE:CZ	2.33	0.63
1:A:85:ASP:HB2	1:A:199:THR:HG23	1.81	0.62
1:C:269:ARG:NH1	1:C:314:ASN:OD1	2.33	0.62
1:A:49:SER:HB3	1:D:93:ARG:HE	1.63	0.62
1:C:436:ARG:HG2	1:C:450:THR:HG23	1.81	0.62
1:B:327:LEU:HA	1:B:379:LEU:HG	1.80	0.62
1:B:108:HIS:HA	1:B:445:PRO:HG3	1.80	0.62
1:C:118:ARG:HD3	1:C:126:VAL:HG11	1.80	0.62
1:D:25:GLN:HG2	1:D:161:ILE:HA	1.82	0.61
1:B:324:VAL:HG11	1:B:382:LEU:HB3	1.81	0.61
1:B:322:PRO:HA	1:B:325:GLN:CD	2.26	0.61
1:A:49:SER:HB2	1:A:53:THR:CG2	2.30	0.61
1:D:149:ASP:OD1	1:D:153:ARG:NH2	2.34	0.61
1:B:164:ARG:HG3	1:B:178:GLN:HG3	1.83	0.60
1:D:198:LYS:HG2	1:D:209:GLY:HA2	1.82	0.60
1:A:293:LEU:HD11	1:A:392:HIS:CD2	2.35	0.60
1:D:422:ILE:O	1:D:436:ARG:NH1	2.35	0.60
1:D:34:GLU:OE2	1:D:34:GLU:N	2.35	0.60
1:A:273:ASP:HB3	1:A:276:LEU:HD12	1.83	0.60
1:B:266:MET:HE1	1:B:398:PRO:HB2	1.82	0.60
1:C:25:GLN:HG3	1:C:26:PHE:H	1.67	0.60
1:B:66:GLU:CD	1:B:98:THR:O	2.45	0.60
1:C:352:THR:HG23	1:C:369:TYR:H	1.65	0.60
1:B:64:ARG:NH2	1:D:366:THR:O	2.32	0.59
1:C:341:LEU:HD22	1:C:424:THR:HG21	1.84	0.59
1:D:96:ASN:HD22	5:M:1:NAG:H83	1.67	0.59
1:D:118:ARG:HD2	1:D:126:VAL:HG21	1.83	0.59
1:A:370:ASP:OD1	1:A:371:PRO:HD2	2.02	0.59
1:B:347:PHE:CG	1:B:426:PRO:HG3	2.37	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:141:LYS:HD3	1:B:265:LEU:HD23	1.84	0.59
1:A:166:SER:HA	1:A:169:ILE:HB	1.85	0.59
1:B:161:ILE:HD11	1:B:190:TRP:CE3	2.38	0.59
1:C:197:LYS:NZ	1:C:198:LYS:O	2.35	0.59
1:B:197:LYS:NZ	1:B:302:LEU:HD13	2.17	0.59
1:B:75:HIS:NE2	1:B:76:SER:O	2.36	0.58
1:B:40:MET:HA	1:B:97:ARG:HB2	1.84	0.58
1:B:49:SER:HB2	1:B:53:THR:HG21	1.84	0.58
1:B:322:PRO:HA	1:B:325:GLN:NE2	2.17	0.58
1:B:327:LEU:HD13	1:B:383:PHE:CB	2.31	0.58
1:C:262:THR:HG22	1:C:264:ASP:H	1.66	0.58
1:B:43:PRO:HG2	1:B:109:ASN:HB3	1.86	0.58
1:D:29:GLN:HG3	1:D:43:PRO:HB3	1.85	0.58
1:B:166:SER:HA	1:B:169:ILE:CD1	2.32	0.58
1:B:91:PRO:HG2	1:B:446:PRO:HG3	1.86	0.57
1:C:73:ARG:NH2	1:D:97:ARG:HH22	2.02	0.57
1:B:218:PRO:HG2	1:B:434:HIS:HB3	1.86	0.57
1:A:358:THR:HG21	1:A:367:GLY:HA2	1.87	0.57
1:B:263:ASP:OD2	1:B:280:ASN:HB3	2.05	0.57
1:A:218:PRO:HB2	1:A:425:PHE:CE1	2.39	0.57
1:B:76:SER:C	1:B:78:GLN:OE1	2.48	0.57
1:B:269:ARG:NH1	1:B:314:ASN:OD1	2.37	0.57
1:B:328:PRO:HG3	1:B:376:LEU:HD23	1.87	0.57
1:B:190:TRP:O	1:B:193:TYR:N	2.38	0.57
1:B:304:ASN:HD21	7:B:504:NAG:C1	2.18	0.57
1:B:326:ARG:HH21	1:B:327:LEU:HB2	1.70	0.57
1:D:364:ASP:CG	1:D:368:LYS:HB3	2.29	0.56
1:B:276:LEU:HD21	1:B:312:ARG:CZ	2.36	0.56
1:C:425:PHE:O	1:C:436:ARG:NH1	2.39	0.56
1:D:343:ASP:O	1:D:424:THR:HB	2.06	0.56
1:B:163:THR:OG1	1:B:178:GLN:O	2.24	0.56
1:B:364:ASP:CG	1:B:368:LYS:HG2	2.31	0.56
1:D:392:HIS:CD2	1:D:393:LEU:HG	2.41	0.56
1:B:260:ILE:HD12	1:B:266:MET:HG3	1.88	0.56
1:A:25:GLN:N	1:A:187:TYR:HH	2.04	0.55
1:B:66:GLU:OE1	1:B:97:ARG:C	2.49	0.55
1:A:25:GLN:OE1	1:A:25:GLN:HA	2.06	0.55
1:B:66:GLU:OE1	1:B:97:ARG:CA	2.54	0.55
1:D:291:ASP:HB2	1:D:393:LEU:HD21	1.87	0.55
1:C:182:ILE:HD11	1:C:186:ASN:HB3	1.89	0.55
1:C:196:VAL:HG12	1:C:212:ASP:CG	2.31	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:161:ILE:HD13	1:A:182:ILE:HD13	1.87	0.55
1:A:241:GLU:HG3	1:C:203:VAL:HG12	1.89	0.55
1:B:328:PRO:HG3	1:B:376:LEU:HA	1.89	0.55
1:A:78:GLN:O	1:A:80:PRO:HD3	2.07	0.54
1:D:173:ASP:O	1:D:175:ASN:N	2.39	0.54
1:D:218:PRO:HD2	1:D:434:HIS:HB3	1.88	0.54
1:D:232:GLU:HG3	1:D:246:LEU:HG	1.88	0.54
1:B:201:LEU:HB2	1:B:205:GLN:HG2	1.89	0.54
1:B:73:ARG:HB3	1:B:74:PRO:CD	2.37	0.54
1:C:129:VAL:HG22	1:C:463:THR:HG22	1.89	0.54
1:A:331:GLN:O	1:A:335:GLN:HG3	2.07	0.54
1:A:138:LYS:H	1:A:138:LYS:CD	2.21	0.53
1:A:217:GLY:O	1:A:356:ARG:HD3	2.08	0.53
1:B:81:HIS:ND1	1:B:84:ARG:HG3	2.24	0.53
1:D:322:PRO:HA	1:D:325:GLN:HG3	1.90	0.53
1:A:209:GLY:O	1:A:211:VAL:N	2.40	0.53
1:B:88:GLU:O	1:B:90:TRP:N	2.41	0.53
1:C:248:TYR:HA	1:C:405:THR:OG1	2.09	0.53
1:A:370:ASP:C	1:A:372:ALA:H	2.15	0.53
1:D:80:PRO:HG2	1:D:81:HIS:CE1	2.44	0.53
1:A:169:ILE:HD11	1:A:179:PHE:CZ	2.44	0.53
1:D:293:LEU:HD23	1:D:296:TYR:HD2	1.74	0.53
1:A:141:LYS:HD3	1:A:265:LEU:HD23	1.90	0.53
1:B:273:ASP:HB3	1:B:276:LEU:HG	1.91	0.53
1:B:190:TRP:O	1:B:192:HIS:N	2.42	0.52
1:C:34:GLU:OE2	1:C:37:ARG:NE	2.42	0.52
1:B:52:GLY:HA2	1:B:55:ARG:CZ	2.40	0.52
1:B:85:ASP:C	1:B:86:ASP:OD1	2.52	0.52
1:A:104:ASN:CG	1:A:114:ARG:HH11	2.16	0.52
1:A:413:GLU:OE2	1:A:417:ARG:NE	2.42	0.52
1:D:198:LYS:HG2	1:D:209:GLY:CA	2.39	0.52
1:B:327:LEU:HB3	1:B:328:PRO:CD	2.37	0.52
1:B:185:TYR:O	1:B:188:PHE:HB3	2.08	0.52
1:B:29:GLN:CG	1:B:43:PRO:HB3	2.39	0.52
1:B:216:GLU:HG2	1:B:362:PHE:CE2	2.38	0.52
1:C:141:LYS:O	1:C:145:VAL:HG23	2.10	0.52
1:A:275:THR:HG22	1:A:314:ASN:HD22	1.74	0.52
1:B:197:LYS:HZ1	1:B:302:LEU:HD13	1.74	0.52
1:D:31:ALA:HB1	1:D:179:PHE:CD2	2.45	0.52
1:B:163:THR:OG1	1:B:178:GLN:HB3	2.10	0.52
1:B:73:ARG:HD3	1:B:73:ARG:N	2.24	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:149:ASP:CG	1:D:153:ARG:HH21	2.18	0.51
1:D:228:LEU:HD11	1:D:404:HIS:HB3	1.91	0.51
1:B:428:GLU:OE1	1:B:429:ASN:N	2.25	0.51
1:A:355:PHE:O	1:A:359:VAL:HG12	2.11	0.51
1:B:39:GLY:O	1:B:96:ASN:N	2.43	0.51
1:A:189:VAL:HG22	1:A:395:PRO:HB2	1.92	0.51
1:C:60:SER:HB2	1:C:62:ARG:HG3	1.91	0.51
1:C:425:PHE:HB3	1:C:436:ARG:HD3	1.92	0.51
1:A:185:TYR:O	1:A:188:PHE:HB3	2.10	0.51
1:C:26:PHE:CD1	1:C:95:PHE:HE2	2.28	0.51
1:A:68:VAL:HG12	1:A:96:ASN:O	2.11	0.51
1:A:293:LEU:HD12	1:A:293:LEU:H	1.75	0.51
1:B:344:THR:HG23	1:B:354:SER:HB2	1.92	0.51
1:B:155:THR:HA	1:B:183:SER:HA	1.91	0.51
1:B:364:ASP:OD1	1:B:368:LYS:HG2	2.11	0.51
1:B:328:PRO:CG	1:B:376:LEU:HA	2.41	0.51
1:B:342:PHE:HB3	1:B:424:THR:HG23	1.93	0.50
1:C:118:ARG:HD3	1:C:126:VAL:CG1	2.41	0.50
1:C:436:ARG:HG2	1:C:450:THR:CG2	2.42	0.50
1:B:28:ARG:HH22	1:B:181:ASN:CG	2.19	0.50
1:B:152:LYS:NZ	1:B:186:ASN:OD1	2.44	0.50
1:A:247:PRO:O	1:A:402:LEU:HD23	2.11	0.50
1:A:340:GLY:HA2	1:A:418:TYR:CE2	2.47	0.50
1:B:41:CYS:HB2	1:B:95:PHE:CD1	2.47	0.50
1:D:333:VAL:O	1:D:337:LEU:HD12	2.12	0.50
1:C:88:GLU:OE1	1:C:93:ARG:NH2	2.45	0.50
1:C:185:TYR:O	1:C:188:PHE:HB3	2.12	0.50
1:C:427:LEU:HD11	1:C:436:ARG:NH2	2.26	0.50
1:B:88:GLU:O	1:B:89:VAL:C	2.54	0.50
1:C:137:SER:O	1:C:141:LYS:HG3	2.12	0.50
1:A:173:ASP:OD2	1:A:176:THR:N	2.44	0.49
1:B:187:TYR:O	1:B:191:THR:HG23	2.11	0.49
1:C:89:VAL:HG12	1:C:89:VAL:O	2.12	0.49
1:A:322:PRO:HA	1:A:325:GLN:HG3	1.94	0.49
1:B:337:LEU:H	1:B:337:LEU:CD1	2.17	0.49
1:D:329:GLU:N	1:D:332:ASP:OD2	2.43	0.49
1:A:293:LEU:HD21	1:A:392:HIS:CG	2.48	0.49
1:B:403:LEU:O	1:B:407:THR:HG23	2.13	0.49
1:B:437:GLN:OE1	1:B:450:THR:HG21	2.12	0.49
1:B:86:ASP:CG	1:B:165:ARG:NH1	2.70	0.49
1:D:25:GLN:HE21	1:D:162:ALA:H	1.60	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:278:SER:O	1:C:284:SER:HB3	2.13	0.49
1:A:197:LYS:HE2	1:A:198:LYS:O	2.12	0.49
1:B:347:PHE:CD1	1:B:426:PRO:HG3	2.46	0.49
4:G:2:NAG:O7	4:G:2:NAG:O3	2.29	0.49
1:A:293:LEU:HA	1:A:296:TYR:CD2	2.48	0.49
1:D:118:ARG:CD	1:D:126:VAL:HG21	2.42	0.49
1:D:466:ILE:HD13	1:D:468:TRP:CH2	2.47	0.49
1:B:448:THR:OG1	1:B:451:GLU:HG3	2.13	0.49
1:D:336:CYS:SG	1:D:359:VAL:HG22	2.52	0.49
1:B:337:LEU:N	1:B:337:LEU:CD1	2.72	0.48
1:B:176:THR:OG1	1:B:178:GLN:NE2	2.46	0.48
1:B:217:GLY:O	1:B:356:ARG:HD3	2.13	0.48
1:A:214:SER:O	1:A:220:PHE:HA	2.13	0.48
1:B:73:ARG:O	1:B:74:PRO:O	2.30	0.48
1:B:224:HIS:CD2	1:B:404:HIS:CE1	3.01	0.48
1:C:28:ARG:HH21	1:C:181:ASN:ND2	2.10	0.48
1:C:133:LEU:HD21	1:C:265:LEU:HG	1.96	0.48
1:C:181:ASN:OD1	7:C:503:NAG:H2	2.13	0.48
1:A:275:THR:OG1	1:A:312:ARG:NH1	2.46	0.48
1:B:86:ASP:CG	1:B:165:ARG:HH11	2.22	0.48
1:A:25:GLN:HG3	1:A:26:PHE:H	1.78	0.48
1:B:350:ASN:HA	1:B:369:TYR:HE2	1.74	0.48
1:B:422:ILE:O	1:B:436:ARG:NH2	2.47	0.48
1:D:260:ILE:O	1:D:266:MET:HB2	2.14	0.48
1:A:260:ILE:HB	1:A:265:LEU:CD1	2.43	0.48
1:C:34:GLU:CG	1:C:37:ARG:HH21	2.25	0.48
1:B:66:GLU:N	1:B:66:GLU:CD	2.71	0.48
1:D:33:VAL:HG13	1:D:34:GLU:OE2	2.13	0.48
1:A:469:PRO:O	1:A:470:SER:HB2	2.14	0.47
1:B:93:ARG:HE	1:B:93:ARG:HA	1.78	0.47
1:A:216:GLU:HB3	1:A:349:SER:HB3	1.95	0.47
1:B:125:ARG:HD2	1:C:49:SER:HA	1.97	0.47
1:C:201:LEU:HD11	1:C:208:PHE:HB2	1.96	0.47
1:A:182:ILE:HD12	1:A:305:SER:HB3	1.96	0.47
1:A:85:ASP:OD1	1:A:86:ASP:N	2.47	0.47
1:B:85:ASP:C	1:B:88:GLU:OE1	2.58	0.47
1:C:71:ASP:HB2	1:C:437:GLN:O	2.15	0.47
1:D:182:ILE:HD11	1:D:186:ASN:HB3	1.97	0.47
1:D:291:ASP:OD1	1:D:393:LEU:HD11	2.14	0.47
1:C:196:VAL:HG12	1:C:212:ASP:OD2	2.15	0.47
1:C:149:ASP:CG	1:C:153:ARG:HH12	2.23	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:SER:HB2	1:B:78:GLN:OE1	2.14	0.47
1:B:134:LEU:HD21	1:B:260:ILE:HG22	1.96	0.47
1:D:375:SER:HA	1:D:379:LEU:HB2	1.95	0.47
1:B:41:CYS:HB2	1:B:95:PHE:CG	2.50	0.46
1:B:42:CYS:O	1:B:55:ARG:HG3	2.16	0.46
1:C:28:ARG:HH12	1:C:157:PRO:C	2.23	0.46
1:C:130:ARG:HH21	1:C:232:GLU:CD	2.17	0.46
1:B:255:LYS:N	1:B:326:ARG:NH1	2.64	0.46
1:B:27:PRO:HG3	1:B:158:LEU:HD11	1.97	0.46
1:B:273:ASP:OD1	1:B:275:THR:OG1	2.33	0.46
1:B:133:LEU:HD22	1:B:402:LEU:HD13	1.98	0.46
1:B:430:ALA:H	1:B:435:ASN:HD21	1.63	0.46
1:B:212:ASP:O	1:B:434:HIS:NE2	2.45	0.46
1:A:50:GLY:O	1:A:53:THR:OG1	2.31	0.46
1:B:37:ARG:C	1:B:39:GLY:H	2.24	0.46
1:D:347:PHE:CG	1:D:426:PRO:HG3	2.51	0.46
1:C:166:SER:HA	1:C:169:ILE:HB	1.98	0.46
1:D:313:ARG:HD3	1:D:397:ASP:OD2	2.16	0.46
1:A:416:ARG:NH2	1:A:417:ARG:HE	2.14	0.46
1:C:47:PRO:HB2	1:C:53:THR:OG1	2.15	0.46
1:C:145:VAL:HG13	1:C:283:PHE:CZ	2.51	0.46
1:B:45:LEU:HB3	1:B:54:ASP:OD1	2.16	0.45
1:B:66:GLU:OE2	1:B:98:THR:O	2.33	0.45
1:B:314:ASN:ND2	1:B:317:GLY:HA3	2.31	0.45
1:C:156:HIS:CE1	1:C:158:LEU:HB3	2.51	0.45
1:B:49:SER:CB	1:B:53:THR:HG21	2.45	0.45
1:A:138:LYS:HD3	1:A:138:LYS:N	2.28	0.45
1:A:318:ASN:HA	3:H:1:NAG:O6	2.17	0.45
1:B:46:SER:O	1:B:48:VAL:HG23	2.17	0.45
1:B:68:VAL:HG23	1:B:96:ASN:C	2.40	0.45
1:B:196:VAL:HG12	1:B:212:ASP:OD1	2.17	0.45
1:C:25:GLN:N	1:C:442:PRO:O	2.49	0.45
1:A:163:THR:HG22	1:A:178:GLN:O	2.16	0.45
1:B:161:ILE:HD12	1:B:161:ILE:HA	1.58	0.45
1:A:90:TRP:HA	1:A:91:PRO:HA	1.78	0.45
1:B:37:ARG:HD2	1:B:37:ARG:HA	1.27	0.45
1:D:355:PHE:CE1	1:D:359:VAL:HG21	2.52	0.45
1:D:288:VAL:HA	1:D:396:ASN:OD1	2.15	0.45
1:D:403:LEU:O	1:D:403:LEU:HD23	2.17	0.45
1:A:266:MET:HE3	1:A:266:MET:HA	1.99	0.45
1:B:316:ALA:O	1:B:325:GLN:HB3	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:91:PRO:CG	1:B:446:PRO:HG3	2.47	0.45
1:C:89:VAL:HG13	1:C:92:LEU:HD12	1.99	0.45
1:D:238:MET:HE2	1:D:238:MET:HB3	1.89	0.45
1:B:334:ALA:HB2	1:B:469:PRO:HB2	1.99	0.44
1:D:68:VAL:CG2	1:D:92:LEU:HD21	2.47	0.44
1:B:76:SER:O	1:B:78:GLN:N	2.50	0.44
1:B:85:ASP:HB2	1:B:199:THR:HG23	1.98	0.44
1:B:66:GLU:CD	1:B:66:GLU:C	2.83	0.44
1:D:169:ILE:HD11	1:D:179:PHE:CZ	2.53	0.44
1:A:118:ARG:HG3	1:A:126:VAL:CG1	2.42	0.44
1:A:224:HIS:CD2	1:A:404:HIS:CE1	3.05	0.44
1:B:54:ASP:OD2	1:B:59:SER:HB3	2.17	0.44
1:B:129:VAL:HG12	1:B:131:ARG:HG2	1.99	0.44
1:D:92:LEU:HA	1:D:92:LEU:HD23	1.64	0.44
1:C:73:ARG:HH22	1:D:97:ARG:NH2	2.09	0.44
1:D:370:ASP:HB3	1:D:373:VAL:CG2	2.47	0.44
1:A:173:ASP:OD2	1:A:176:THR:HG23	2.18	0.44
1:B:202:GLY:O	1:B:205:GLN:HG3	2.18	0.44
1:C:264:ASP:OD2	1:C:265:LEU:HD13	2.17	0.44
1:B:144:PHE:HE2	1:B:402:LEU:HD11	1.83	0.44
1:B:173:ASP:HB3	1:B:176:THR:HB	1.99	0.44
1:B:221:LEU:HD22	1:B:411:PHE:HB2	1.99	0.44
1:B:324:VAL:HG22	1:B:379:LEU:HD12	2.00	0.44
1:B:265:LEU:HD12	1:B:265:LEU:O	2.17	0.44
1:B:296:TYR:CZ	1:B:303:CYS:HA	2.52	0.44
1:A:260:ILE:HD12	1:A:266:MET:HG3	1.98	0.44
1:C:169:ILE:HG23	1:C:170:LEU:HG	2.00	0.44
1:C:296:TYR:CZ	1:C:303:CYS:HA	2.53	0.44
1:A:391:VAL:HG13	1:A:392:HIS:N	2.32	0.43
1:B:152:LYS:HD3	1:B:286:TRP:CE3	2.53	0.43
1:B:165:ARG:HD2	1:B:166:SER:N	2.30	0.43
1:B:168:GLU:C	1:B:171:GLY:H	2.26	0.43
1:B:387:THR:O	1:B:394:SER:HA	2.18	0.43
1:B:86:ASP:N	1:B:86:ASP:OD1	2.50	0.43
1:B:225:ARG:HB2	1:B:408:ASP:OD1	2.18	0.43
1:A:357:ASN:OD1	1:A:362:PHE:HB2	2.17	0.43
1:C:189:VAL:HG22	1:C:395:PRO:HB2	1.99	0.43
1:D:419:ASN:OD1	1:D:419:ASN:O	2.36	0.43
1:A:253:THR:OG1	1:A:255:LYS:HG3	2.19	0.43
1:A:324:VAL:HA	1:A:379:LEU:CD1	2.48	0.43
1:A:180:GLU:OE1	1:A:305:SER:OG	2.36	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:257:VAL:HG12	1:B:258:CYS:N	2.29	0.43
1:C:357:ASN:OD1	1:C:362:PHE:HB2	2.18	0.43
1:B:189:VAL:HA	1:B:395:PRO:HB2	1.99	0.43
1:A:92:LEU:HD23	1:A:92:LEU:HA	1.73	0.43
1:C:68:VAL:HG11	1:C:91:PRO:HD2	2.01	0.43
1:C:331:GLN:CD	1:C:331:GLN:H	2.25	0.43
1:D:196:VAL:HG12	1:D:212:ASP:OD1	2.19	0.43
1:A:118:ARG:CZ	1:A:126:VAL:HG11	2.49	0.43
1:C:139:GLU:OE1	1:C:139:GLU:N	2.47	0.42
1:B:343:ASP:O	1:B:424:THR:OG1	2.37	0.42
1:B:392:HIS:CD2	1:B:393:LEU:HD22	2.54	0.42
1:C:128:ILE:HG12	1:C:461:GLY:O	2.18	0.42
1:D:332:ASP:OD1	1:D:365:PRO:HG2	2.19	0.42
1:C:129:VAL:HG22	1:C:463:THR:CG2	2.50	0.42
1:D:273:ASP:HB3	1:D:276:LEU:HD12	2.02	0.42
1:A:87:ARG:HG2	1:A:441:VAL:HG11	2.01	0.42
1:A:142:ASN:OD1	1:A:146:ARG:HD2	2.20	0.42
1:A:162:ALA:HB2	1:A:179:PHE:CE2	2.54	0.42
1:C:232:GLU:OE2	1:C:245:SER:HA	2.20	0.42
1:D:87:ARG:NH1	1:D:197:LYS:HG2	2.35	0.42
1:A:129:VAL:HG12	1:A:131:ARG:HG2	2.01	0.42
1:A:358:THR:HG22	1:A:363:SER:CB	2.48	0.42
1:B:324:VAL:O	1:B:324:VAL:CG1	2.68	0.42
1:D:265:LEU:HD12	1:D:265:LEU:O	2.19	0.42
1:A:238:MET:HB3	1:A:238:MET:HE2	1.75	0.42
1:A:308:ASP:HB3	1:A:309:GLY:H	1.76	0.42
1:A:376:LEU:HA	1:A:376:LEU:HD23	1.80	0.42
1:C:266:MET:CE	1:C:398:PRO:HB2	2.49	0.42
1:B:65:CYS:CA	1:B:66:GLU:OE2	2.67	0.42
9:C:514:KOJ:O2	10:C:601:HOH:O	2.21	0.42
1:A:149:ASP:O	1:A:153:ARG:HG3	2.19	0.42
1:A:444:TRP:HA	1:A:445:PRO:C	2.44	0.42
1:C:347:PHE:CG	1:C:426:PRO:HG3	2.55	0.42
1:A:76:SER:OG	1:A:431:PRO:HA	2.20	0.42
9:A:516:KOJ:O2	10:A:601:HOH:O	2.21	0.42
1:C:439:ASN:CG	1:C:446:PRO:HB3	2.45	0.42
1:A:143:HIS:CE1	1:A:241:GLU:OE2	2.73	0.42
1:A:229:LEU:HD12	1:A:229:LEU:HA	1.87	0.42
1:A:348:TYR:CG	1:A:349:SER:N	2.88	0.42
1:A:392:HIS:NE2	1:A:393:LEU:HG	2.34	0.42
1:B:102:ASN:OD1	1:D:340:GLY:HA3	2.20	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:41:CYS:HB3	1:C:444:TRP:CZ2	2.55	0.42
1:D:224:HIS:CD2	1:D:404:HIS:CE1	3.08	0.42
1:A:278:SER:O	1:A:284:SER:HB3	2.20	0.41
1:B:190:TRP:C	1:B:192:HIS:N	2.77	0.41
1:B:436:ARG:HG2	1:B:450:THR:CG2	2.50	0.41
1:C:96:ASN:C	1:C:97:ARG:HG3	2.45	0.41
1:A:34:GLU:H	1:A:34:GLU:HG3	1.71	0.41
1:A:257:VAL:HG22	1:A:258:CYS:N	2.35	0.41
1:B:458:ASP:OD2	1:C:51:PRO:HG2	2.20	0.41
1:A:25:GLN:HG3	1:A:26:PHE:N	2.34	0.41
1:D:136:LEU:O	1:D:141:LYS:HE3	2.20	0.41
1:D:313:ARG:HG2	1:D:315:PRO:HD3	2.02	0.41
1:A:313:ARG:HG2	1:A:315:PRO:HD3	2.03	0.41
1:B:247:PRO:O	1:B:402:LEU:HD23	2.20	0.41
1:C:100:HIS:ND1	1:D:64:ARG:NH2	2.69	0.41
1:C:217:GLY:O	1:C:356:ARG:HD3	2.20	0.41
1:B:348:TYR:CG	1:B:349:SER:N	2.89	0.41
1:D:435:ASN:O	1:D:436:ARG:C	2.64	0.41
1:A:68:VAL:HG13	1:A:92:LEU:HD21	2.03	0.41
1:C:160:VAL:HG12	1:C:181:ASN:HD22	1.86	0.41
1:C:188:PHE:CE1	1:C:227:HIS:CE1	3.08	0.41
1:A:143:HIS:HE1	1:A:241:GLU:OE2	2.04	0.41
1:A:260:ILE:O	1:A:266:MET:HB2	2.21	0.41
1:B:37:ARG:C	1:B:39:GLY:N	2.75	0.41
1:D:130:ARG:HH21	1:D:232:GLU:CD	2.28	0.41
1:D:262:THR:OG1	1:D:264:ASP:OD1	2.33	0.41
1:D:359:VAL:O	1:D:377:HIS:HB3	2.20	0.41
1:D:441:VAL:HA	1:D:442:PRO:HA	1.90	0.41
1:A:138:LYS:HE2	1:A:139:GLU:OE2	2.21	0.41
1:A:382:LEU:CD2	9:A:516:KOJ:H6C2	2.51	0.41
1:A:391:VAL:CG1	1:A:392:HIS:N	2.84	0.41
1:B:198:LYS:HD2	1:B:209:GLY:HA3	2.03	0.41
1:B:223:TRP:CD1	1:B:440:MET:HB3	2.56	0.41
1:B:304:ASN:HD21	7:B:504:NAG:C2	2.33	0.41
1:B:364:ASP:HB2	1:B:365:PRO:HD2	2.02	0.41
1:B:399:ILE:O	1:B:400:PHE:C	2.63	0.41
1:C:231:LEU:O	1:C:235:MET:HG3	2.21	0.41
1:C:368:LYS:HE2	1:C:368:LYS:HB3	1.94	0.41
1:D:80:PRO:HG2	1:D:81:HIS:ND1	2.36	0.41
1:D:197:LYS:HE3	1:D:302:LEU:HD23	2.02	0.41
1:A:287:ARG:HB2	1:A:307:GLU:OE1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:418:TYR:CD1	1:A:418:TYR:N	2.89	0.41
1:B:170:LEU:HA	1:B:176:THR:O	2.21	0.41
1:B:440:MET:H	1:B:440:MET:HG3	1.69	0.41
1:C:89:VAL:O	1:C:89:VAL:CG1	2.69	0.41
1:C:314:ASN:ND2	1:C:317:GLY:HA3	2.36	0.41
1:B:118:ARG:HG3	1:B:126:VAL:CG1	2.51	0.40
1:C:209:GLY:C	1:C:211:VAL:H	2.28	0.40
1:D:73:ARG:H	1:D:73:ARG:HG2	1.64	0.40
1:B:251:PHE:HB2	1:B:327:LEU:HG	2.03	0.40
1:B:333:VAL:C	1:B:335:GLN:H	2.29	0.40
1:D:173:ASP:C	1:D:175:ASN:N	2.79	0.40
1:D:431:PRO:O	1:D:432:ILE:C	2.65	0.40
1:D:469:PRO:O	1:D:470:SER:HB2	2.21	0.40
1:A:240:GLN:OE1	1:C:165:ARG:NH2	2.53	0.40
1:B:328:PRO:HG3	1:B:376:LEU:H	1.86	0.40
1:D:25:GLN:HG3	1:D:26:PHE:N	2.31	0.40
1:D:217:GLY:O	1:D:356:ARG:HD3	2.22	0.40
1:A:94:PHE:O	2:E:3:FUC:H61	2.21	0.40
1:A:399:ILE:O	1:A:400:PHE:C	2.64	0.40
1:A:411:PHE:CZ	1:A:415:LEU:HD11	2.56	0.40
1:B:192:HIS:O	1:B:196:VAL:HG22	2.21	0.40
1:C:92:LEU:HD23	1:C:92:LEU:HA	1.69	0.40
1:C:184:ILE:CG2	1:C:234:ASP:HB3	2.51	0.40
1:D:88:GLU:O	1:D:89:VAL:C	2.63	0.40
1:D:360:GLU:OE2	1:D:362:PHE:HE2	2.05	0.40
1:D:370:ASP:HB3	1:D:373:VAL:HG23	2.03	0.40
1:A:254:GLY:HA2	1:A:327:LEU:O	2.21	0.40
1:B:377:HIS:O	1:B:381:HIS:HD2	2.04	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	444/446 (100%)	402 (90%)	40 (9%)	2 (0%)	24	42
1	B	444/446 (100%)	377 (85%)	51 (12%)	16 (4%)	2	5
1	C	444/446 (100%)	409 (92%)	31 (7%)	4 (1%)	14	28
1	D	444/446 (100%)	403 (91%)	37 (8%)	4 (1%)	14	28
All	All	1776/1784 (100%)	1591 (90%)	159 (9%)	26 (2%)	8	18

All (26) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	74	PRO
1	B	191	THR
1	B	197	LYS
1	B	327	LEU
1	B	427	LEU
1	D	174	GLY
1	A	197	LYS
1	B	171	GLY
1	B	385	ASN
1	B	419	ASN
1	A	419	ASN
1	B	54	ASP
1	B	181	ASN
1	B	326	ARG
1	B	387	THR
1	C	47	PRO
1	C	172	PRO
1	D	368	LYS
1	D	436	ARG
1	C	376	LEU
1	B	426	PRO
1	D	266	MET
1	B	77	PRO
1	C	364	ASP
1	B	365	PRO
1	B	442	PRO

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	395/395 (100%)	389 (98%)	6 (2%)	57	77
1	B	395/395 (100%)	384 (97%)	11 (3%)	38	62
1	C	395/395 (100%)	389 (98%)	6 (2%)	57	77
1	D	395/395 (100%)	392 (99%)	3 (1%)	73	86
All	All	1580/1580 (100%)	1554 (98%)	26 (2%)	55	76

All (26) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	MET
1	A	155	THR
1	A	374	SER
1	A	405	THR
1	A	447	VAL
1	A	450	THR
1	B	60	SER
1	B	66	GLU
1	B	73	ARG
1	B	155	THR
1	B	183	SER
1	B	207	SER
1	B	258	CYS
1	B	262	THR
1	B	278	SER
1	B	343	ASP
1	B	354	SER
1	C	60	SER
1	C	197	LYS
1	C	203	VAL
1	C	258	CYS
1	C	278	SER
1	C	349	SER
1	D	49	SER
1	D	265	LEU
1	D	363	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	GLN
1	A	143	HIS
1	A	205	GLN
1	A	224	HIS
1	B	178	GLN
1	B	304	ASN
1	B	378	ASN
1	C	78	GLN
1	C	181	ASN
1	D	78	GLN
1	D	250	ASN
1	D	419	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

29 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	1.15	2 (14%)	17,19,21	1.30	2 (11%)
2	NAG	E	2	2	14,14,15	1.01	1 (7%)	17,19,21	0.59	0
2	FUC	E	3	2	10,10,11	1.42	0	14,14,16	1.25	2 (14%)
3	NAG	F	1	1,3	14,14,15	0.94	1 (7%)	17,19,21	0.91	1 (5%)
3	NAG	F	2	3	14,14,15	0.46	0	17,19,21	0.72	0
4	NAG	G	1	1,4	14,14,15	0.83	1 (7%)	17,19,21	0.90	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	G	2	4	14,14,15	0.67	1 (7%)	17,19,21	1.32	2 (11%)
4	MAN	G	3	4	11,11,12	1.16	1 (9%)	15,15,17	1.73	2 (13%)
4	MAN	G	4	4	11,11,12	0.98	1 (9%)	15,15,17	1.80	4 (26%)
4	MAN	G	5	4	11,11,12	1.52	3 (27%)	15,15,17	1.62	3 (20%)
3	NAG	H	1	1,3	14,14,15	0.53	0	17,19,21	0.61	0
3	NAG	H	2	3	14,14,15	0.83	1 (7%)	17,19,21	0.55	0
5	NAG	I	1	1,5	14,14,15	1.23	3 (21%)	17,19,21	1.21	1 (5%)
5	FUC	I	2	5	10,10,11	1.84	4 (40%)	14,14,16	1.44	3 (21%)
5	NAG	J	1	1,5	14,14,15	0.65	0	17,19,21	0.61	0
5	FUC	J	2	5	10,10,11	1.52	3 (30%)	14,14,16	1.88	2 (14%)
3	NAG	K	1	1,3	14,14,15	0.79	1 (7%)	17,19,21	0.74	0
3	NAG	K	2	3	14,14,15	0.54	0	17,19,21	0.45	0
6	NAG	L	1	1,6	14,14,15	1.81	2 (14%)	17,19,21	2.42	4 (23%)
6	NAG	L	2	6	14,14,15	0.97	1 (7%)	17,19,21	1.24	2 (11%)
6	MAN	L	3	6	11,11,12	1.30	1 (9%)	15,15,17	1.64	2 (13%)
6	MAN	L	4	6	11,11,12	0.71	0	15,15,17	1.59	3 (20%)
6	MAN	L	5	6	11,11,12	1.25	1 (9%)	15,15,17	1.60	1 (6%)
5	NAG	M	1	1,5	14,14,15	0.55	0	17,19,21	0.77	0
5	FUC	M	2	5	10,10,11	2.35	4 (40%)	14,14,16	1.95	4 (28%)
5	NAG	N	1	1,5	14,14,15	1.96	3 (21%)	17,19,21	1.24	2 (11%)
5	FUC	N	2	5	10,10,11	2.03	3 (30%)	14,14,16	1.52	3 (21%)
3	NAG	O	1	1,3	14,14,15	0.33	0	17,19,21	0.67	0
3	NAG	O	2	3	14,14,15	0.56	0	17,19,21	0.55	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	FUC	E	3	2	-	-	0/1/1/1
3	NAG	F	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
4	NAG	G	1	1,4	-	4/6/23/26	0/1/1/1
4	NAG	G	2	4	-	3/6/23/26	0/1/1/1
4	MAN	G	3	4	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	MAN	G	4	4	-	2/2/19/22	0/1/1/1
4	MAN	G	5	4	-	2/2/19/22	0/1/1/1
3	NAG	H	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
5	NAG	I	1	1,5	-	2/6/23/26	0/1/1/1
5	FUC	I	2	5	-	-	0/1/1/1
5	NAG	J	1	1,5	-	2/6/23/26	0/1/1/1
5	FUC	J	2	5	-	-	0/1/1/1
3	NAG	K	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	K	2	3	-	1/6/23/26	0/1/1/1
6	NAG	L	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	L	2	6	-	2/6/23/26	0/1/1/1
6	MAN	L	3	6	-	0/2/19/22	1/1/1/1
6	MAN	L	4	6	-	0/2/19/22	0/1/1/1
6	MAN	L	5	6	-	2/2/19/22	0/1/1/1
5	NAG	M	1	1,5	-	2/6/23/26	0/1/1/1
5	FUC	M	2	5	-	-	0/1/1/1
5	NAG	N	1	1,5	-	2/6/23/26	0/1/1/1
5	FUC	N	2	5	-	-	0/1/1/1
3	NAG	O	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	O	2	3	-	1/6/23/26	0/1/1/1

All (38) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	L	1	NAG	O5-C1	5.46	1.52	1.43
5	N	1	NAG	C1-C2	5.32	1.59	1.52
5	M	2	FUC	C2-C3	4.50	1.59	1.52
5	N	1	NAG	O5-C1	-4.31	1.36	1.43
5	N	2	FUC	C1-C2	3.76	1.61	1.52
6	L	1	NAG	C1-C2	3.68	1.57	1.52
2	E	2	NAG	C1-C2	3.62	1.57	1.52
5	N	2	FUC	C4-C5	3.53	1.60	1.52
6	L	2	NAG	O5-C1	-3.52	1.37	1.43
6	L	3	MAN	O5-C5	3.44	1.50	1.43
5	J	2	FUC	O5-C5	3.42	1.50	1.43
5	I	2	FUC	C1-C2	3.41	1.60	1.52
3	F	1	NAG	O5-C1	-3.38	1.38	1.43
5	M	2	FUC	C1-C2	3.23	1.59	1.52
5	I	2	FUC	O5-C5	3.12	1.49	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	G	1	NAG	O5-C1	3.00	1.48	1.43
5	I	1	NAG	O5-C1	-2.92	1.38	1.43
2	E	1	NAG	O5-C1	2.87	1.48	1.43
3	H	2	NAG	C1-C2	2.86	1.56	1.52
6	L	5	MAN	O5-C5	2.83	1.49	1.43
5	M	2	FUC	O4-C4	2.79	1.49	1.43
3	K	1	NAG	O5-C1	-2.78	1.39	1.43
5	I	1	NAG	C1-C2	2.71	1.56	1.52
4	G	5	MAN	C2-C3	2.62	1.56	1.52
4	G	5	MAN	C4-C3	2.58	1.59	1.52
4	G	4	MAN	C2-C3	2.39	1.56	1.52
4	G	2	NAG	C1-C2	-2.35	1.49	1.52
5	N	2	FUC	C2-C3	2.32	1.56	1.52
5	N	1	NAG	C3-C2	2.29	1.57	1.52
4	G	5	MAN	C1-C2	2.28	1.57	1.52
5	M	2	FUC	C4-C3	2.23	1.58	1.52
2	E	1	NAG	C8-C7	2.22	1.55	1.50
5	J	2	FUC	C4-C5	2.21	1.57	1.52
4	G	3	MAN	O5-C5	2.14	1.47	1.43
5	I	1	NAG	C3-C2	2.14	1.57	1.52
5	I	2	FUC	C4-C5	2.08	1.57	1.52
5	J	2	FUC	C2-C3	2.07	1.55	1.52
5	I	2	FUC	O5-C1	2.05	1.47	1.43

All (44) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	L	1	NAG	C1-O5-C5	5.83	120.00	112.19
6	L	1	NAG	O4-C4-C5	5.69	123.33	109.32
6	L	5	MAN	C1-O5-C5	5.48	119.54	112.19
5	J	2	FUC	O5-C5-C4	5.17	118.86	109.55
6	L	3	MAN	C1-O5-C5	4.84	118.67	112.19
5	M	2	FUC	C1-C2-C3	4.65	116.42	109.64
4	G	4	MAN	C1-O5-C5	4.30	117.95	112.19
4	G	3	MAN	C1-O5-C5	4.30	117.95	112.19
5	I	1	NAG	C4-C3-C2	3.87	116.68	111.02
5	N	1	NAG	C4-C3-C2	3.80	116.58	111.02
5	J	2	FUC	C1-O5-C5	3.62	121.50	112.97
4	G	5	MAN	C2-C3-C4	3.49	117.00	110.86
6	L	4	MAN	O2-C2-C3	-3.49	102.92	110.15
4	G	3	MAN	O2-C2-C3	-3.49	102.93	110.15
6	L	1	NAG	C1-C2-N2	3.42	115.81	110.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	G	2	NAG	C2-N2-C7	3.28	127.30	122.90
5	N	2	FUC	O2-C2-C1	3.26	116.69	109.22
6	L	1	NAG	C4-C3-C2	-3.20	106.32	111.02
4	G	4	MAN	C1-C2-C3	3.15	114.23	109.64
6	L	4	MAN	C1-O5-C5	3.09	116.32	112.19
6	L	3	MAN	O2-C2-C3	-3.05	103.83	110.15
4	G	5	MAN	C1-C2-C3	3.00	114.02	109.64
2	E	3	FUC	O5-C5-C4	2.99	114.93	109.55
5	N	1	NAG	O5-C5-C4	-2.86	103.86	110.83
5	I	2	FUC	O5-C5-C4	2.85	114.67	109.55
6	L	4	MAN	C1-C2-C3	2.84	113.78	109.64
4	G	5	MAN	C3-C4-C5	2.83	115.36	110.23
5	I	2	FUC	C1-O5-C5	2.78	119.51	112.97
5	M	2	FUC	O5-C5-C4	2.75	114.50	109.55
4	G	2	NAG	O4-C4-C5	-2.68	102.72	109.32
5	M	2	FUC	C2-C3-C4	2.52	115.30	110.86
5	N	2	FUC	C1-O5-C5	2.48	118.82	112.97
2	E	1	NAG	O4-C4-C5	2.46	115.37	109.32
2	E	3	FUC	C1-O5-C5	2.42	118.67	112.97
6	L	2	NAG	C3-C4-C5	-2.39	105.89	110.23
4	G	1	NAG	O4-C4-C3	-2.39	104.75	110.38
5	M	2	FUC	C1-O5-C5	2.35	118.50	112.97
3	F	1	NAG	C1-O5-C5	2.32	115.29	112.19
5	N	2	FUC	O5-C5-C4	2.30	113.69	109.55
4	G	4	MAN	O5-C1-C2	2.30	116.27	110.79
4	G	4	MAN	C3-C4-C5	-2.25	106.16	110.23
5	I	2	FUC	O2-C2-C1	2.12	114.07	109.22
6	L	2	NAG	C4-C3-C2	-2.08	107.98	111.02
2	E	1	NAG	C8-C7-N2	2.02	119.47	116.12

There are no chirality outliers.

All (35) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	G	5	MAN	O5-C5-C6-O6
3	F	2	NAG	O5-C5-C6-O6
4	G	1	NAG	O5-C5-C6-O6
4	G	2	NAG	O5-C5-C6-O6
6	L	5	MAN	O5-C5-C6-O6
3	F	2	NAG	C4-C5-C6-O6
5	I	1	NAG	O5-C5-C6-O6
5	N	1	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
4	G	5	MAN	C4-C5-C6-O6
4	G	1	NAG	C4-C5-C6-O6
5	I	1	NAG	C4-C5-C6-O6
5	N	1	NAG	C4-C5-C6-O6
6	L	5	MAN	C4-C5-C6-O6
4	G	2	NAG	C4-C5-C6-O6
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
4	G	1	NAG	C8-C7-N2-C2
4	G	1	NAG	O7-C7-N2-C2
5	J	1	NAG	C8-C7-N2-C2
5	J	1	NAG	O7-C7-N2-C2
5	M	1	NAG	C8-C7-N2-C2
5	M	1	NAG	O7-C7-N2-C2
6	L	1	NAG	C8-C7-N2-C2
6	L	1	NAG	O7-C7-N2-C2
2	E	1	NAG	O5-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
4	G	4	MAN	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
4	G	4	MAN	C4-C5-C6-O6
3	K	2	NAG	O5-C5-C6-O6
4	G	2	NAG	C3-C2-N2-C7
6	L	2	NAG	C3-C2-N2-C7
3	O	2	NAG	O5-C5-C6-O6
6	L	2	NAG	C1-C2-N2-C7

All (1) ring outliers are listed below:

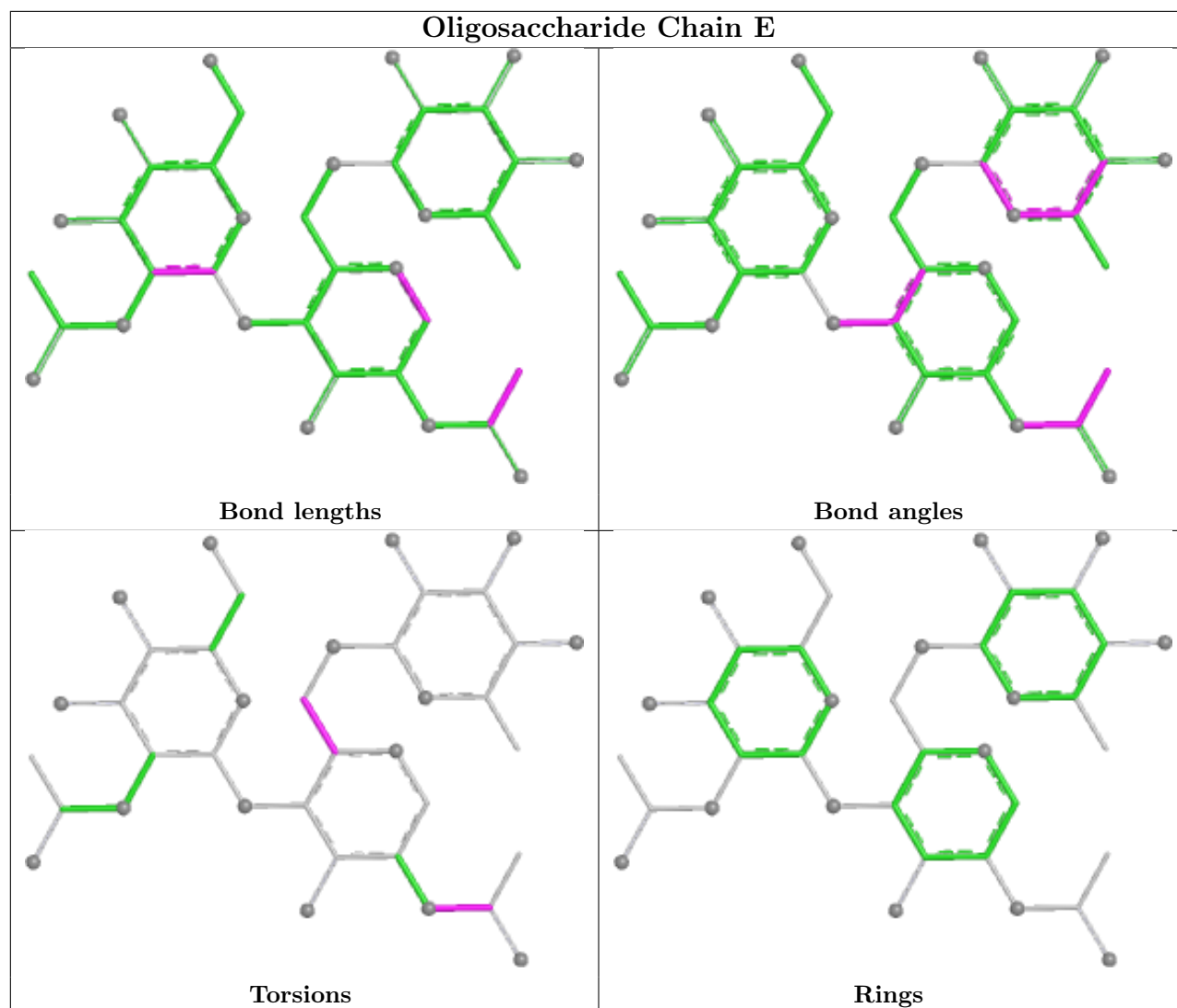
Mol	Chain	Res	Type	Atoms
6	L	3	MAN	C1-C2-C3-C4-C5-O5

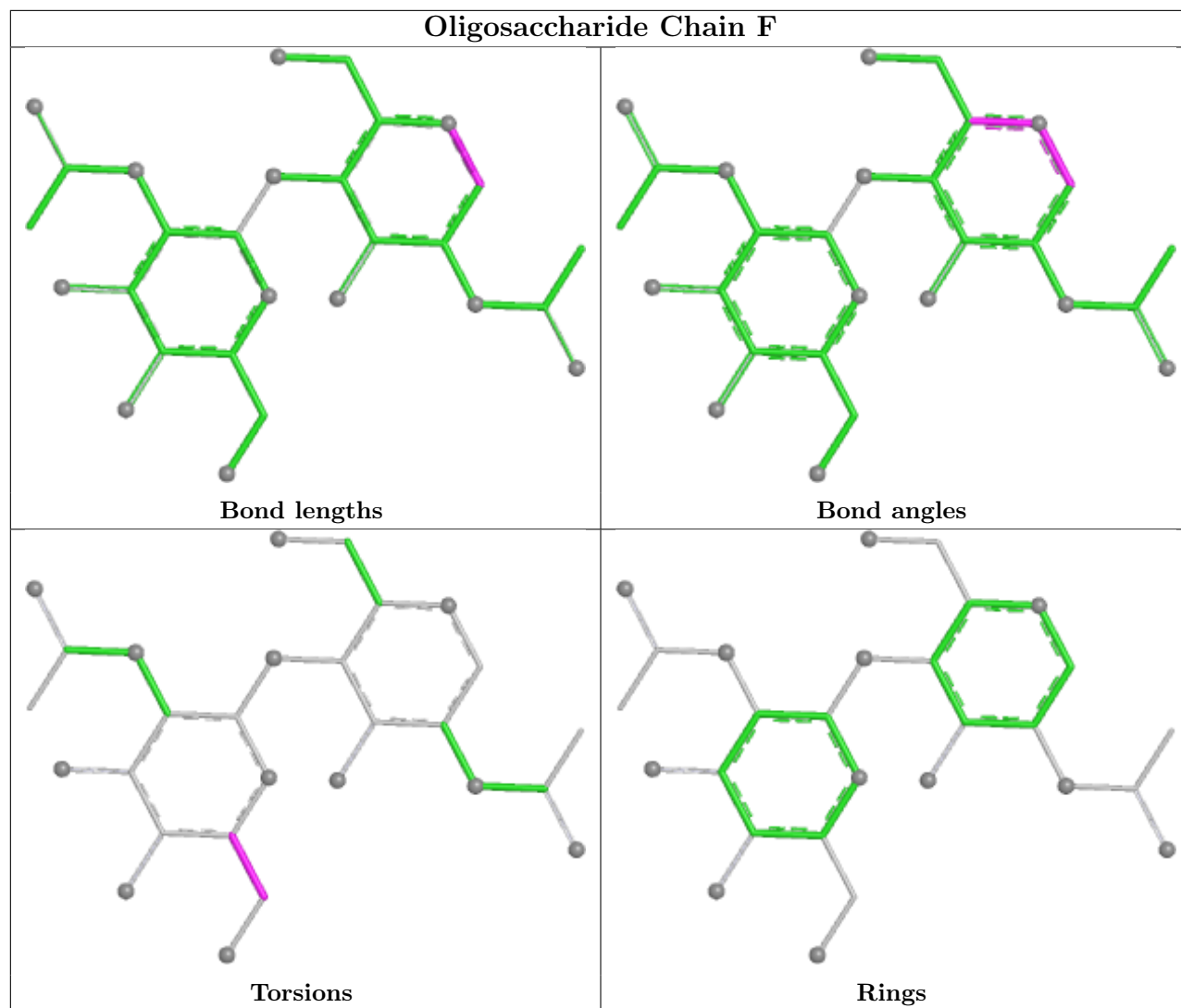
4 monomers are involved in 4 short contacts:

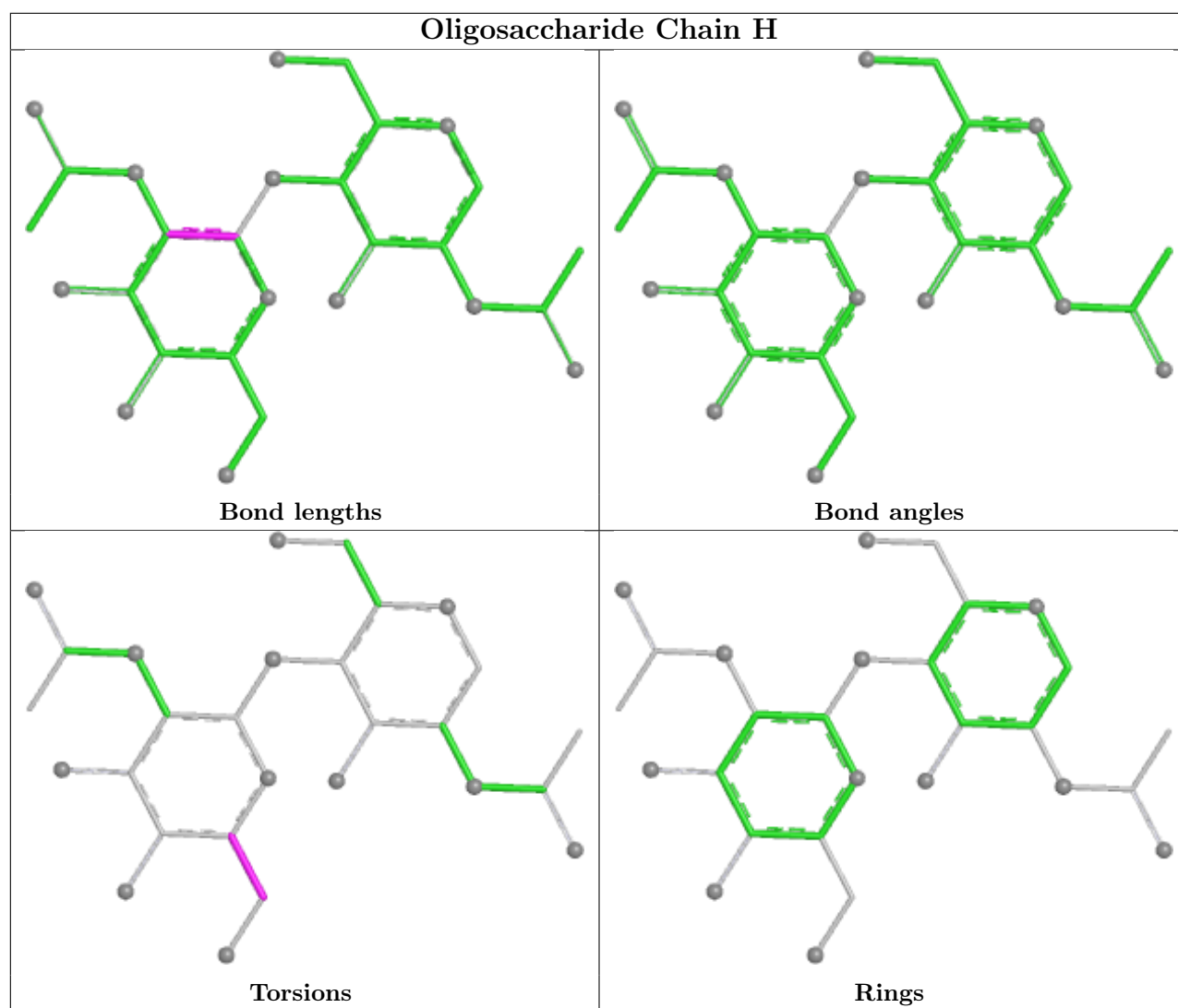
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	1	NAG	1	0
2	E	3	FUC	1	0
5	M	1	NAG	1	0
4	G	2	NAG	1	0

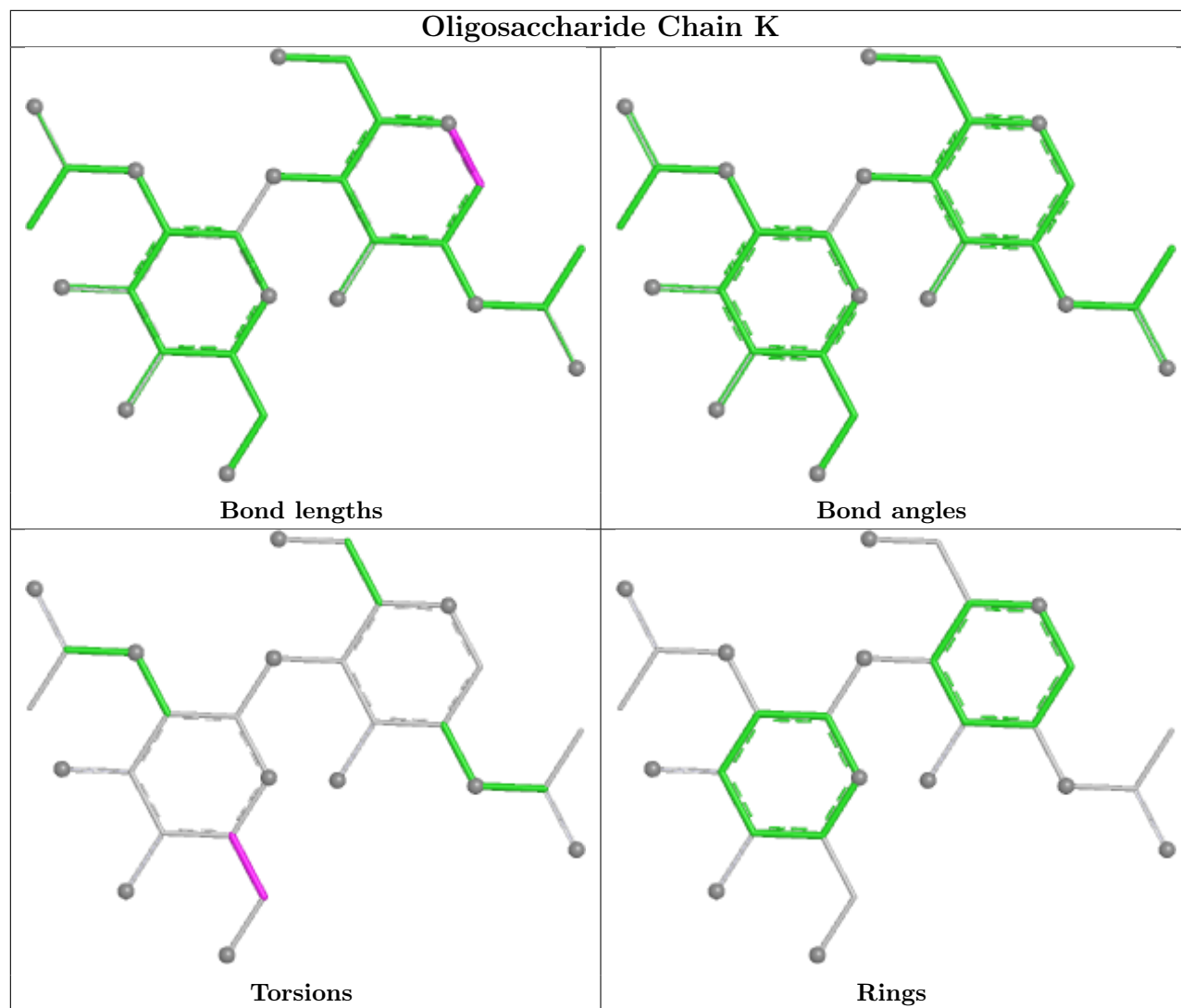
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

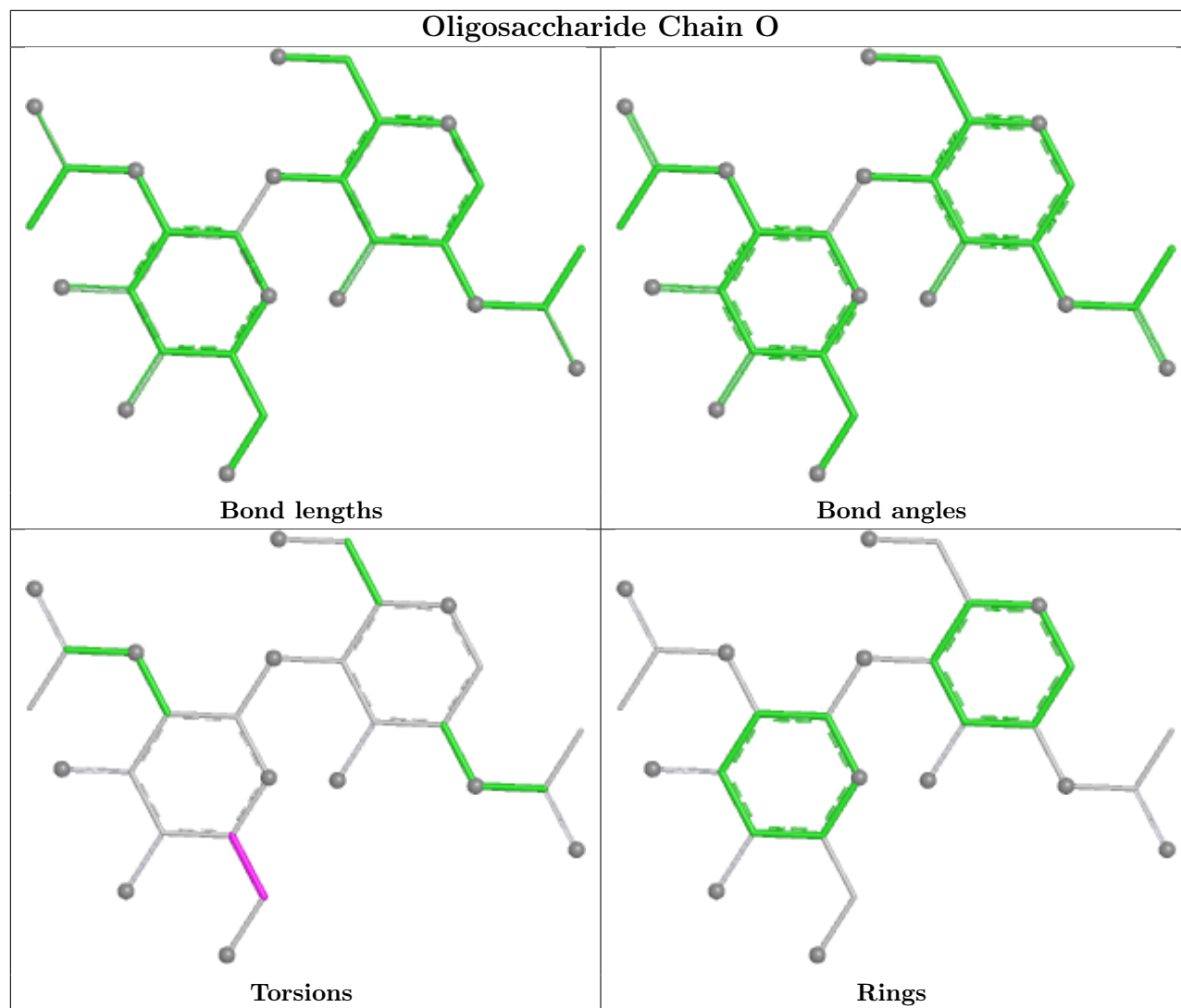
bond angles, torsion angles, and ring geometry for oligosaccharide.



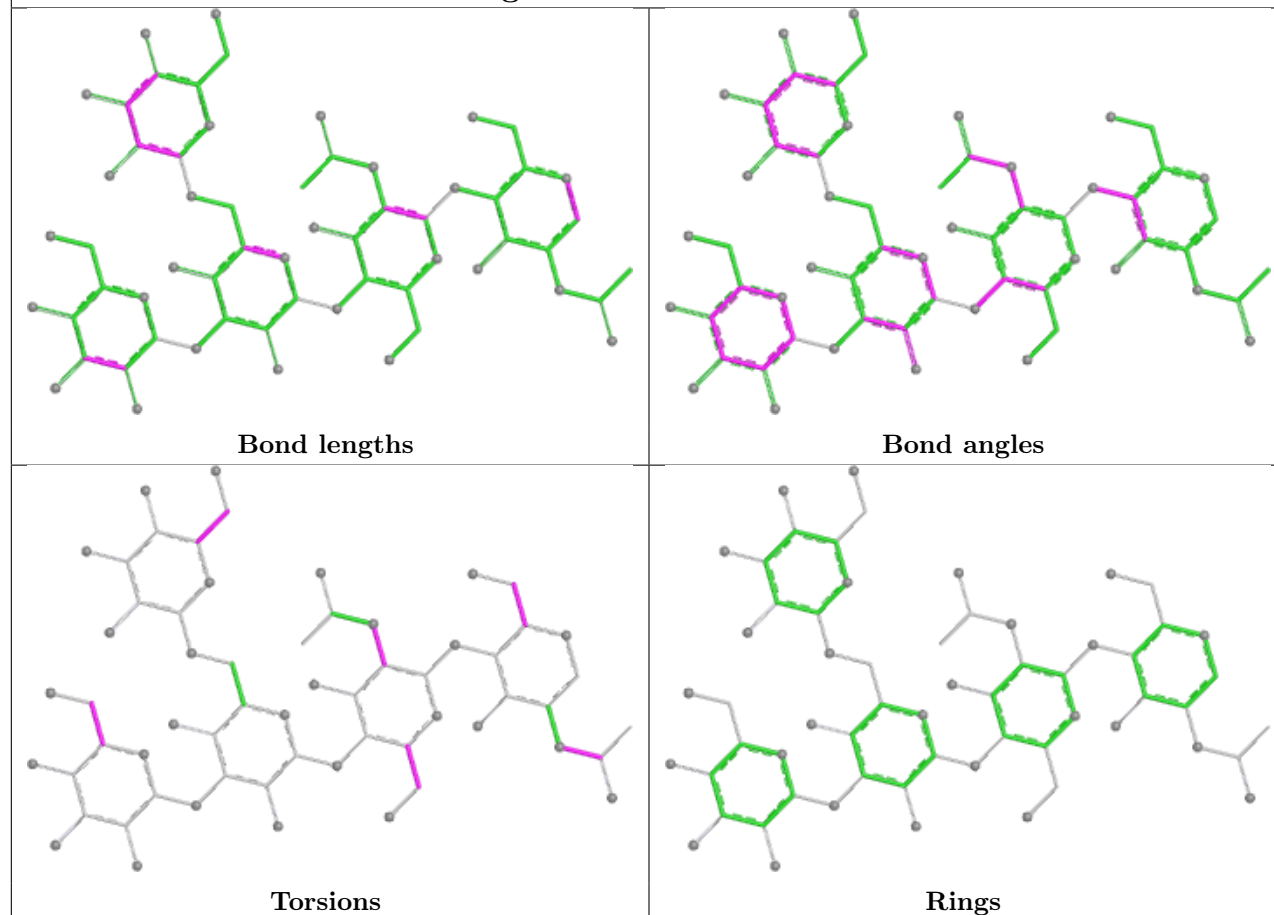




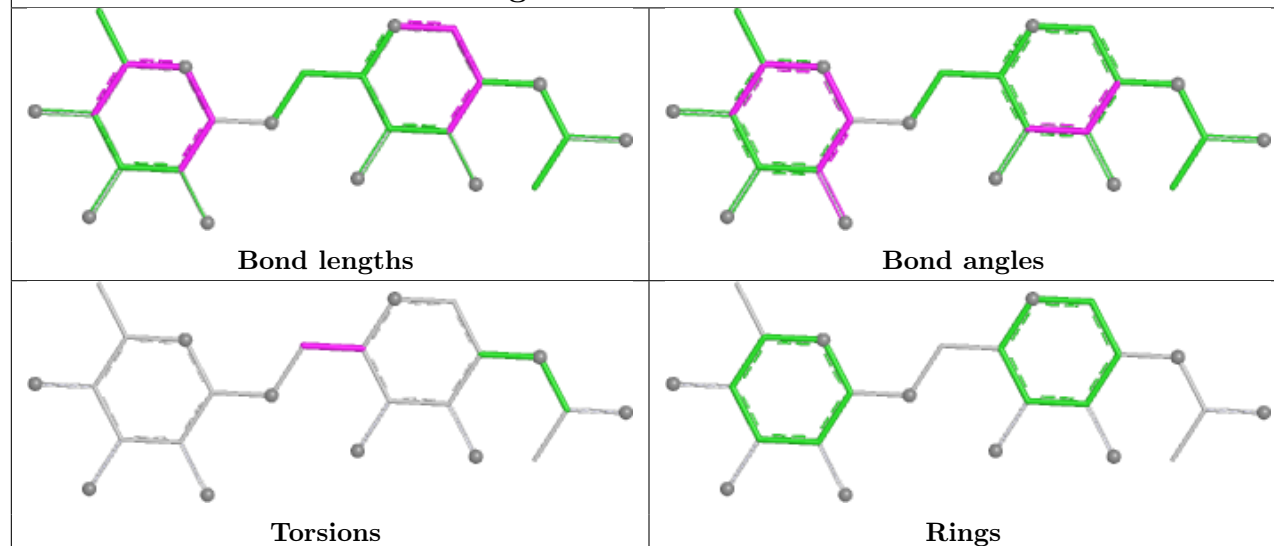


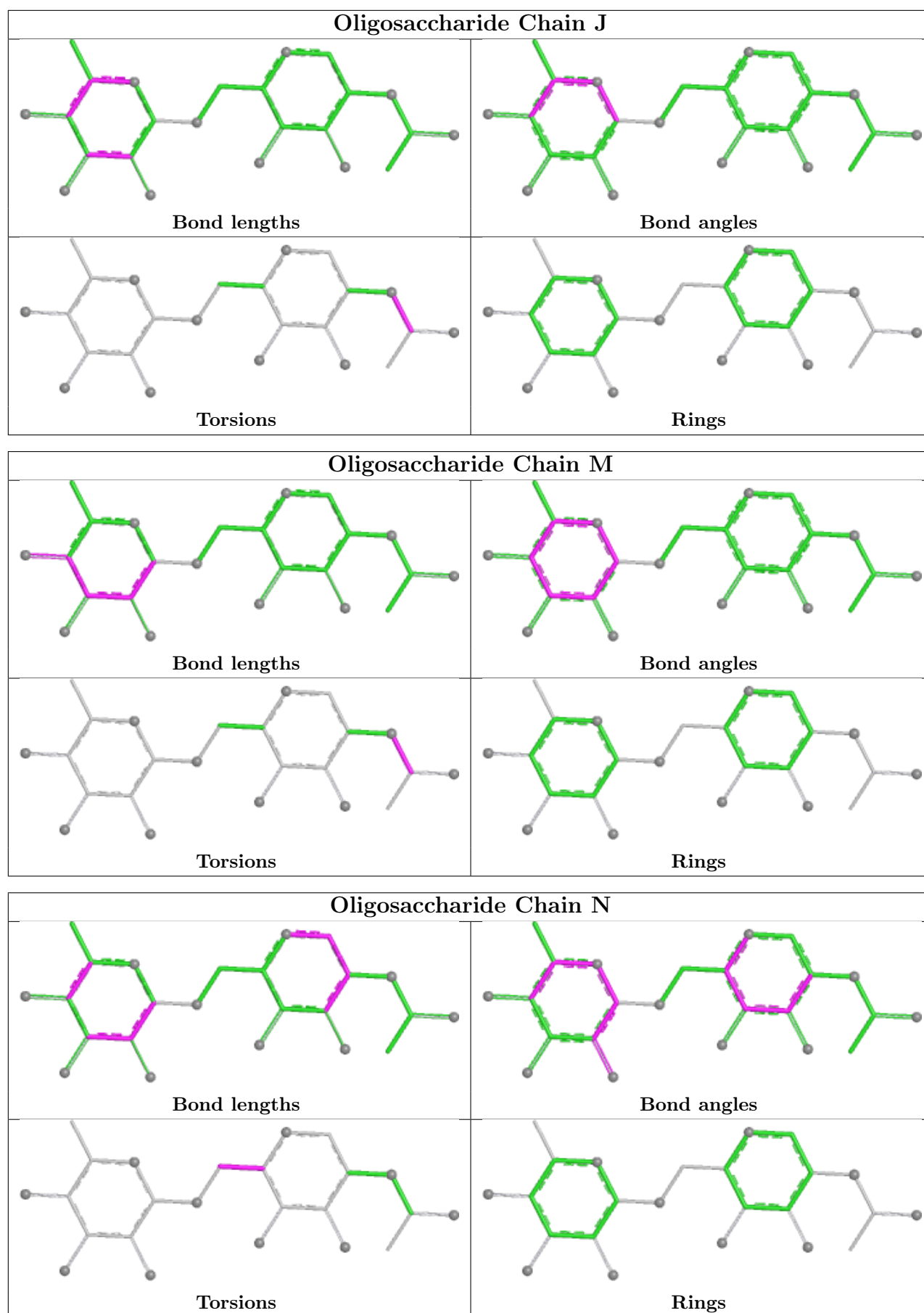


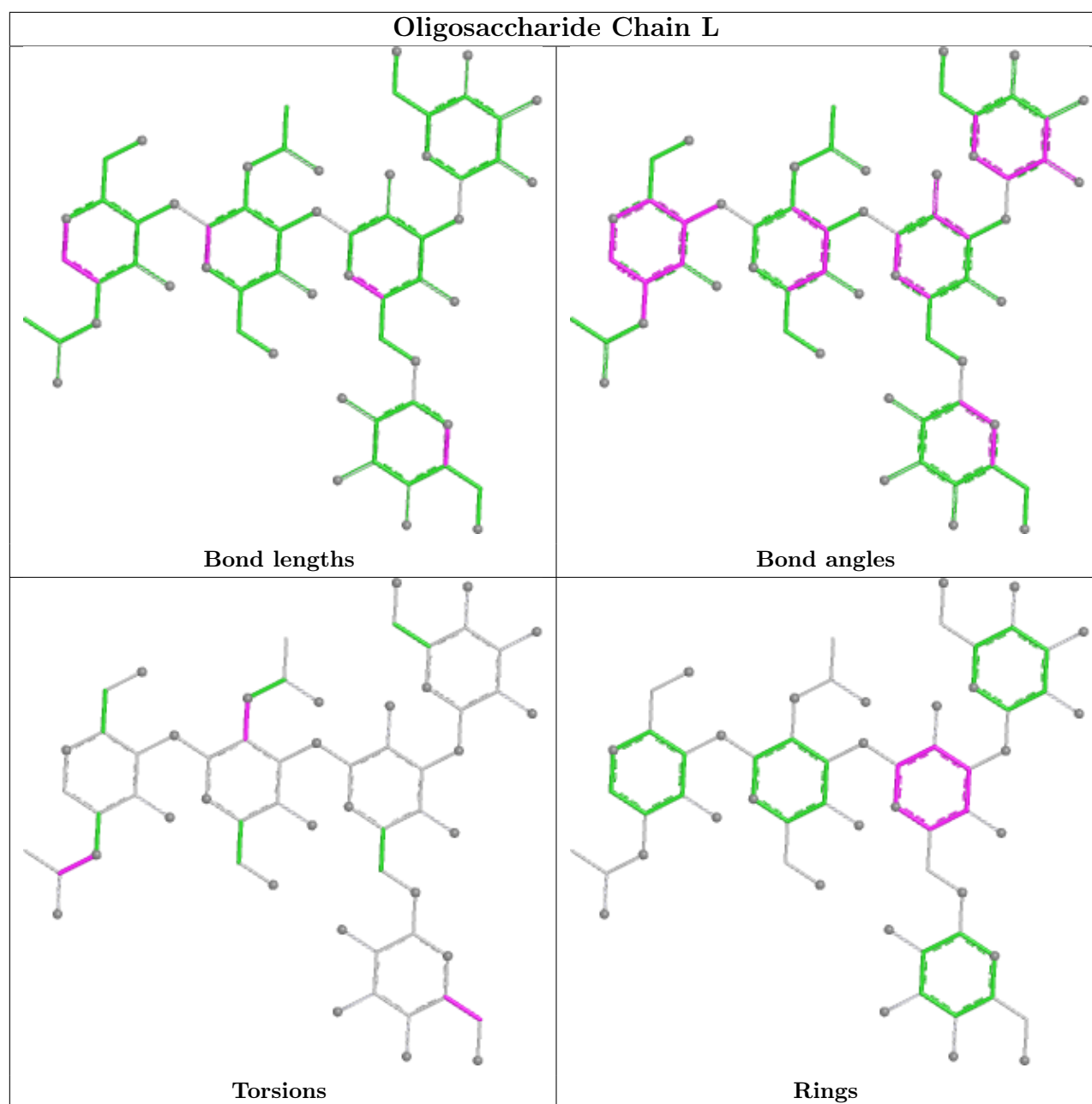
Oligosaccharide Chain G



Oligosaccharide Chain I







5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 9 are monoatomic - leaving 11 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
9	KOJ	C	514	-	10,10,10	2.92	4 (40%)	12,13,13	2.44	5 (41%)
9	KOJ	D	510	-	10,10,10	2.98	3 (30%)	12,13,13	3.23	6 (50%)
9	KOJ	A	516	-	10,10,10	3.00	3 (30%)	12,13,13	3.50	6 (50%)
7	NAG	B	501	1	14,14,15	1.80	1 (7%)	17,19,21	1.82	4 (23%)
7	NAG	D	507	1	14,14,15	1.03	1 (7%)	17,19,21	0.93	1 (5%)
7	NAG	C	503	-	14,14,15	1.04	1 (7%)	17,19,21	1.59	3 (17%)
7	NAG	B	504	-	14,14,15	1.14	1 (7%)	17,19,21	0.63	0
7	NAG	B	505	1	14,14,15	0.81	1 (7%)	17,19,21	0.53	0
7	NAG	C	511	1	14,14,15	0.66	1 (7%)	17,19,21	0.73	1 (5%)
9	KOJ	B	508	-	10,10,10	2.90	4 (40%)	12,13,13	2.45	5 (41%)
7	NAG	A	504	1	14,14,15	1.71	2 (14%)	17,19,21	1.15	3 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	KOJ	C	514	-	-	2/2/2/2	0/1/1/1
9	KOJ	D	510	-	-	2/2/2/2	0/1/1/1
9	KOJ	A	516	-	-	0/2/2/2	0/1/1/1
7	NAG	B	501	1	-	4/6/23/26	0/1/1/1
7	NAG	D	507	1	-	2/6/23/26	0/1/1/1
7	NAG	C	503	-	-	0/6/23/26	0/1/1/1
7	NAG	B	504	-	-	2/6/23/26	0/1/1/1
7	NAG	B	505	1	-	2/6/23/26	0/1/1/1
7	NAG	C	511	1	-	2/6/23/26	0/1/1/1
9	KOJ	B	508	-	-	2/2/2/2	0/1/1/1
7	NAG	A	504	1	-	2/6/23/26	0/1/1/1

All (22) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	B	501	NAG	O5-C1	-6.20	1.33	1.43
9	A	516	KOJ	O5-C5	6.01	1.48	1.35
9	A	516	KOJ	O5-C1	5.95	1.48	1.37
9	D	510	KOJ	O5-C5	5.90	1.47	1.35
7	A	504	NAG	C1-C2	5.87	1.60	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	D	510	KOJ	O5-C1	5.75	1.48	1.37
9	B	508	KOJ	O5-C5	5.65	1.47	1.35
9	C	514	KOJ	O5-C1	5.63	1.47	1.37
9	B	508	KOJ	O5-C1	5.55	1.47	1.37
9	C	514	KOJ	O5-C5	5.12	1.46	1.35
7	B	504	NAG	O5-C1	-4.09	1.36	1.43
9	C	514	KOJ	O2-C2	3.42	1.42	1.33
7	C	503	NAG	O5-C1	-3.33	1.38	1.43
7	D	507	NAG	O5-C1	3.28	1.49	1.43
9	A	516	KOJ	O2-C2	3.15	1.42	1.33
9	D	510	KOJ	O2-C2	2.94	1.41	1.33
9	B	508	KOJ	O2-C2	2.79	1.41	1.33
9	C	514	KOJ	C1-C2	-2.66	1.32	1.34
9	B	508	KOJ	C1-C2	-2.45	1.32	1.34
7	B	505	NAG	O5-C1	2.42	1.47	1.43
7	A	504	NAG	C3-C2	2.09	1.56	1.52
7	C	511	NAG	O5-C1	2.01	1.47	1.43

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	A	516	KOJ	O5-C5-C6	6.75	122.05	111.06
9	A	516	KOJ	O5-C1-C2	-6.54	118.76	122.75
9	D	510	KOJ	O5-C5-C6	6.34	121.38	111.06
9	D	510	KOJ	O5-C1-C2	-5.72	119.26	122.75
7	B	501	NAG	O5-C5-C6	-5.44	97.07	107.66
9	C	514	KOJ	O5-C5-C6	5.34	119.75	111.06
9	A	516	KOJ	C6-C5-C4	-5.31	118.67	126.78
9	B	508	KOJ	O5-C5-C6	5.27	119.64	111.06
9	D	510	KOJ	C6-C5-C4	-4.74	119.54	126.78
9	B	508	KOJ	O5-C1-C2	-4.06	120.27	122.75
9	C	514	KOJ	O5-C5-C4	-4.05	117.45	121.92
9	B	508	KOJ	C6-C5-C4	-3.69	121.14	126.78
7	C	503	NAG	C3-C4-C5	3.58	116.73	110.23
7	C	503	NAG	C1-O5-C5	-3.43	107.58	112.19
9	A	516	KOJ	O2-C2-C1	3.38	124.90	120.15
7	C	503	NAG	C4-C3-C2	3.20	115.70	111.02
7	A	504	NAG	C4-C3-C2	2.81	115.14	111.02
7	B	501	NAG	O4-C4-C5	2.79	116.20	109.32
7	D	507	NAG	C1-O5-C5	2.77	115.90	112.19
7	B	501	NAG	O6-C6-C5	-2.73	102.03	111.33
9	D	510	KOJ	O2-C2-C1	2.65	123.88	120.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	C	514	KOJ	C6-C5-C4	-2.63	122.76	126.78
9	C	514	KOJ	O5-C1-C2	-2.63	121.15	122.75
9	D	510	KOJ	C4-C3-C2	2.59	121.34	116.80
9	D	510	KOJ	O5-C5-C4	-2.58	119.08	121.92
7	A	504	NAG	C2-N2-C7	2.50	126.26	122.90
9	B	508	KOJ	O5-C5-C4	-2.45	119.22	121.92
9	A	516	KOJ	C4-C3-C2	2.44	121.07	116.80
7	A	504	NAG	O5-C5-C4	-2.40	104.99	110.83
9	A	516	KOJ	O5-C5-C4	-2.40	119.28	121.92
9	C	514	KOJ	C1-O5-C5	2.34	121.75	119.33
7	C	511	NAG	C1-O5-C5	2.14	115.06	112.19
7	B	501	NAG	C1-O5-C5	-2.08	109.40	112.19
9	B	508	KOJ	C4-C3-C2	2.05	120.38	116.80

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	C	511	NAG	O5-C5-C6-O6
7	B	505	NAG	O5-C5-C6-O6
7	B	504	NAG	O5-C5-C6-O6
7	C	511	NAG	C4-C5-C6-O6
7	D	507	NAG	O5-C5-C6-O6
7	B	504	NAG	C4-C5-C6-O6
7	B	505	NAG	C4-C5-C6-O6
7	D	507	NAG	C4-C5-C6-O6
7	B	501	NAG	C8-C7-N2-C2
7	B	501	NAG	O7-C7-N2-C2
7	B	501	NAG	O5-C5-C6-O6
9	B	508	KOJ	O5-C5-C6-O6
9	C	514	KOJ	O5-C5-C6-O6
9	D	510	KOJ	C4-C5-C6-O6
9	B	508	KOJ	C4-C5-C6-O6
9	C	514	KOJ	C4-C5-C6-O6
7	B	501	NAG	C4-C5-C6-O6
7	A	504	NAG	O5-C5-C6-O6
9	D	510	KOJ	O5-C5-C6-O6
7	A	504	NAG	C4-C5-C6-O6

There are no ring outliers.

6 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	C	514	KOJ	1	0
9	A	516	KOJ	3	0
7	B	501	NAG	1	0
7	C	503	NAG	1	0
7	B	504	NAG	2	0
9	B	508	KOJ	3	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	446/446 (100%)	0.00	4 (0%) 81 76	33, 51, 66, 79	0
1	B	446/446 (100%)	0.74	49 (10%) 10 8	42, 71, 95, 108	0
1	C	446/446 (100%)	0.02	4 (0%) 81 76	38, 54, 69, 90	0
1	D	446/446 (100%)	-0.10	2 (0%) 88 87	34, 49, 66, 81	0
All	All	1784/1784 (100%)	0.16	59 (3%) 49 40	33, 54, 87, 108	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	48	VAL	5.3
1	A	371	PRO	4.5
1	B	327	LEU	4.4
1	B	427	LEU	3.8
1	B	170	LEU	3.6
1	B	430	ALA	3.5
1	C	48	VAL	3.3
1	B	75	HIS	3.1
1	B	373	VAL	3.1
1	B	334	ALA	3.1
1	B	49	SER	3.0
1	B	100	HIS	2.9
1	B	453	PHE	2.9
1	C	49	SER	2.8
1	B	164	ARG	2.8
1	B	435	ASN	2.8
1	B	350	ASN	2.8
1	B	302	LEU	2.8
1	B	166	SER	2.8
1	D	253	THR	2.7
1	B	33	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	203	VAL	2.6
1	A	69	THR	2.6
1	C	470	SER	2.5
1	B	51	PRO	2.5
1	B	171	GLY	2.5
1	B	257	VAL	2.5
1	B	426	PRO	2.4
1	B	470	SER	2.4
1	B	429	ASN	2.4
1	B	47	PRO	2.4
1	B	324	VAL	2.4
1	B	72	SER	2.4
1	B	172	PRO	2.4
1	B	367	GLY	2.4
1	B	92	LEU	2.3
1	B	207	SER	2.3
1	B	417	ARG	2.3
1	B	319	VAL	2.3
1	B	333	VAL	2.3
1	B	208	PHE	2.3
1	B	298	THR	2.3
1	B	469	PRO	2.2
1	A	25	GLN	2.2
1	C	37	ARG	2.2
1	D	48	VAL	2.2
1	B	90	TRP	2.2
1	B	338	GLU	2.2
1	B	345	PRO	2.1
1	B	306	THR	2.1
1	B	66	GLU	2.1
1	B	67	ALA	2.1
1	A	203	VAL	2.1
1	B	88	GLU	2.1
1	B	89	VAL	2.0
1	B	337	LEU	2.0
1	B	200	PHE	2.0
1	B	74	PRO	2.0
1	B	93	ARG	2.0

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

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

6.3 Carbohydrates

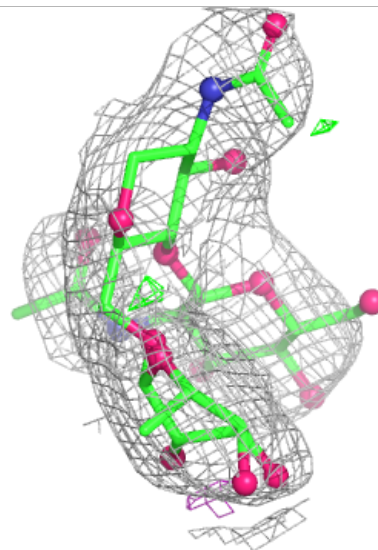
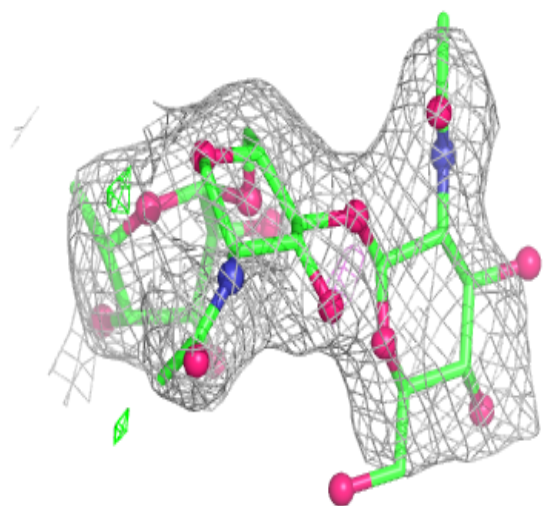
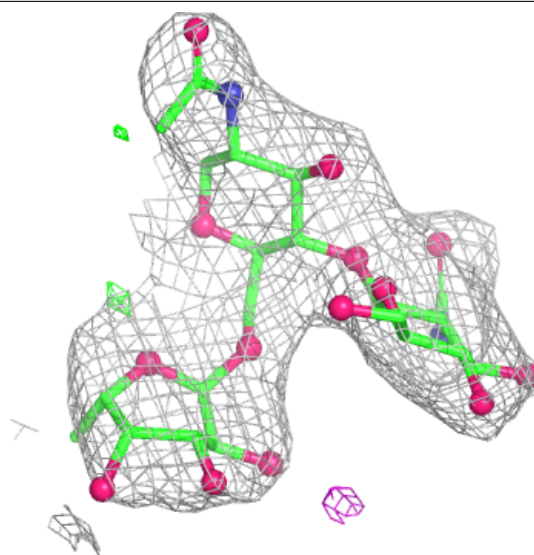
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	H	2	14/15	0.67	0.17	77,84,93,93	0
5	FUC	I	2	10/11	0.68	0.17	69,83,89,90	0
5	NAG	I	1	14/15	0.71	0.14	71,78,86,88	0
5	NAG	N	1	14/15	0.74	0.14	53,68,79,79	0
3	NAG	O	2	14/15	0.75	0.14	64,72,75,76	0
2	NAG	E	2	14/15	0.79	0.12	66,78,87,88	0
5	FUC	J	2	10/11	0.80	0.16	65,68,72,84	0
2	FUC	E	3	10/11	0.80	0.14	59,61,67,72	0
4	MAN	G	4	11/12	0.81	0.15	65,71,85,88	0
5	NAG	J	1	14/15	0.82	0.13	54,64,71,72	0
3	NAG	F	2	14/15	0.82	0.13	57,66,73,78	0
2	NAG	E	1	14/15	0.82	0.12	60,70,72,73	0
6	NAG	L	1	14/15	0.84	0.15	43,49,57,58	0
3	NAG	K	2	14/15	0.85	0.12	56,63,68,69	0
5	FUC	N	2	10/11	0.87	0.11	58,76,83,87	0
5	FUC	M	2	10/11	0.88	0.12	57,63,67,67	0
5	NAG	M	1	14/15	0.88	0.10	44,60,69,71	0
6	NAG	L	2	14/15	0.88	0.13	33,53,62,68	0
6	MAN	L	4	11/12	0.88	0.12	52,59,63,67	0
4	MAN	G	5	11/12	0.89	0.11	63,67,75,75	0
6	MAN	L	5	11/12	0.89	0.10	50,58,65,67	0
3	NAG	K	1	14/15	0.90	0.10	53,58,65,66	0
4	NAG	G	1	14/15	0.90	0.10	44,49,52,52	0
6	MAN	L	3	11/12	0.92	0.10	45,52,58,60	0
3	NAG	H	1	14/15	0.92	0.08	55,61,71,75	0
3	NAG	F	1	14/15	0.92	0.10	56,61,64,65	0
4	MAN	G	3	11/12	0.93	0.10	53,58,67,68	0
4	NAG	G	2	14/15	0.93	0.10	45,50,58,70	0
3	NAG	O	1	14/15	0.94	0.09	49,56,62,62	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.

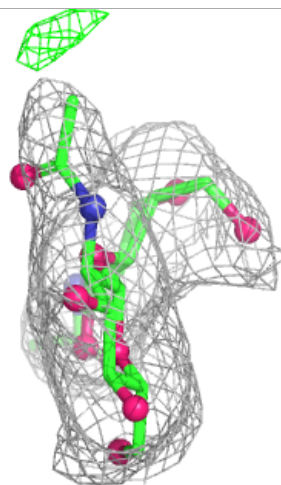
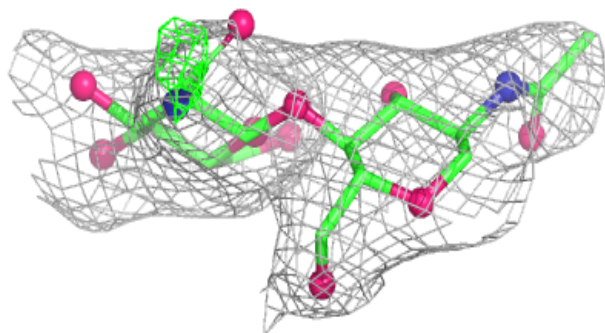
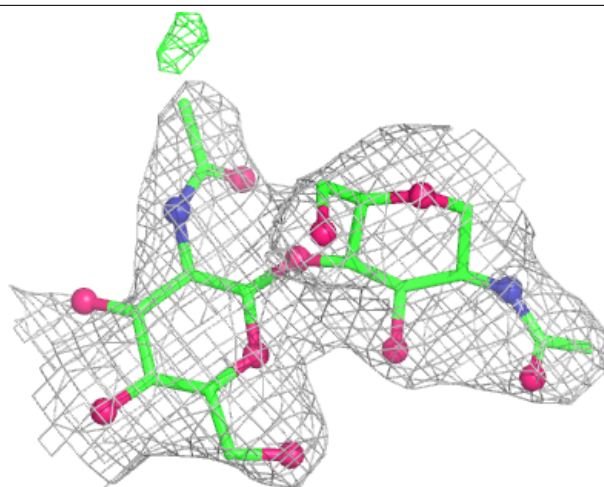
Electron density around Chain E:

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



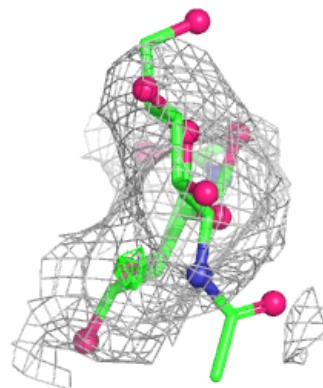
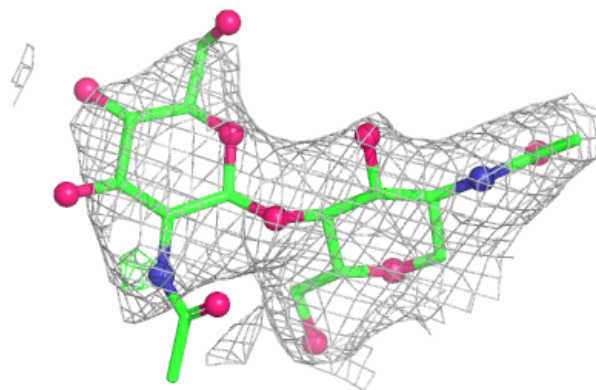
Electron density around Chain F:

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

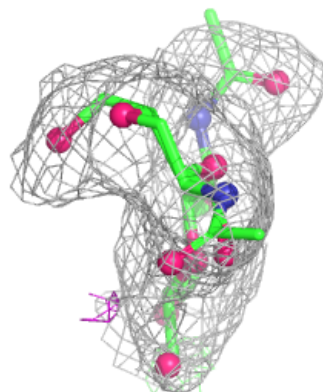
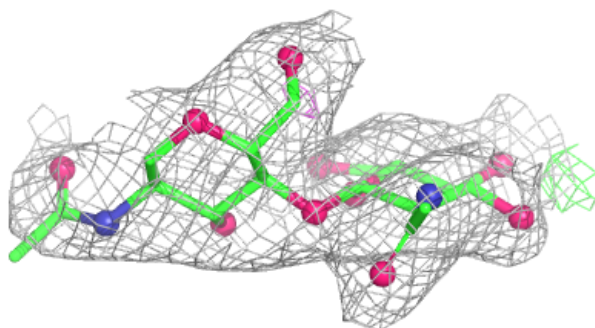
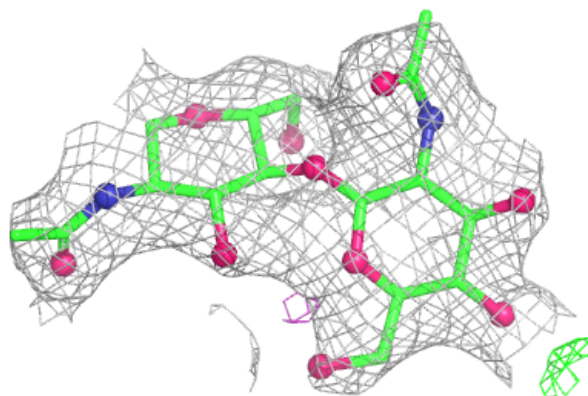


Electron density around Chain H:

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

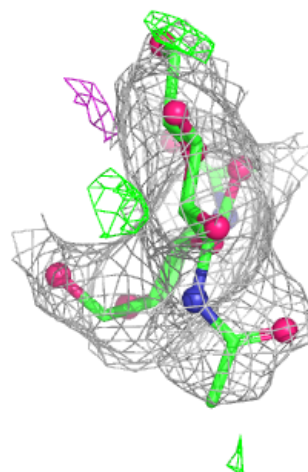
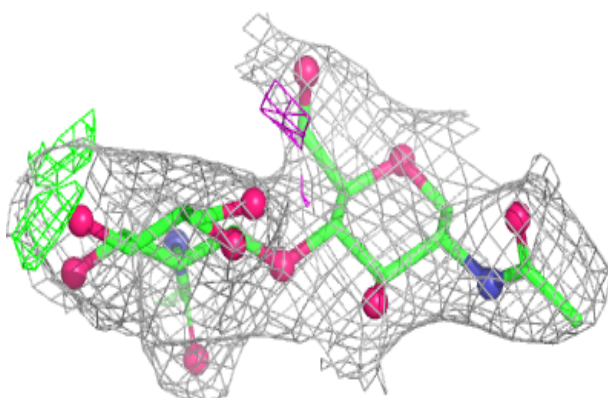
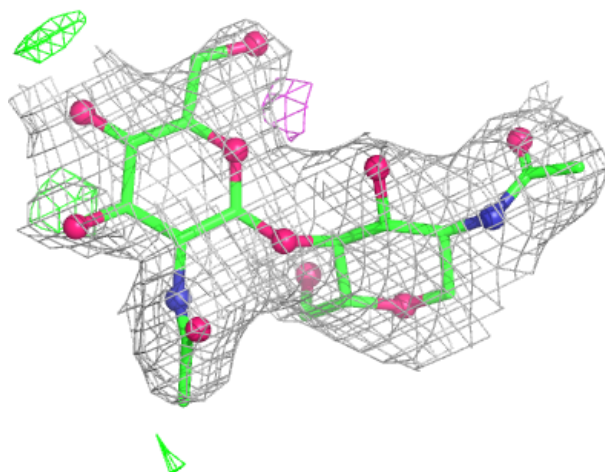
**Electron density around Chain K:**

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



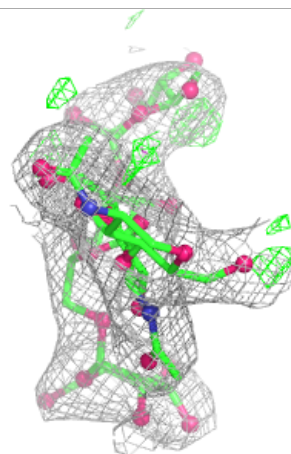
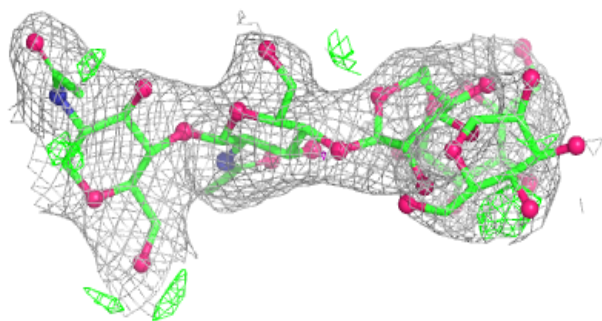
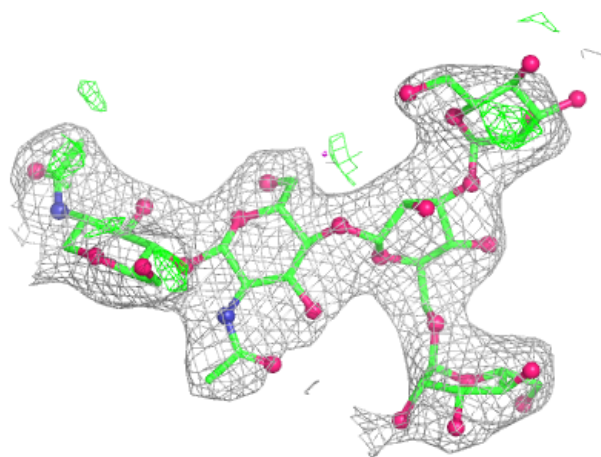
Electron density around Chain O:

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



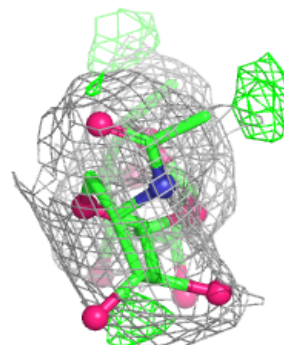
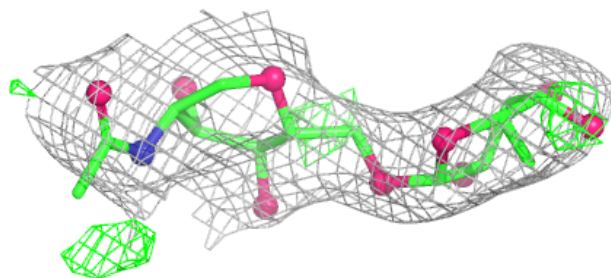
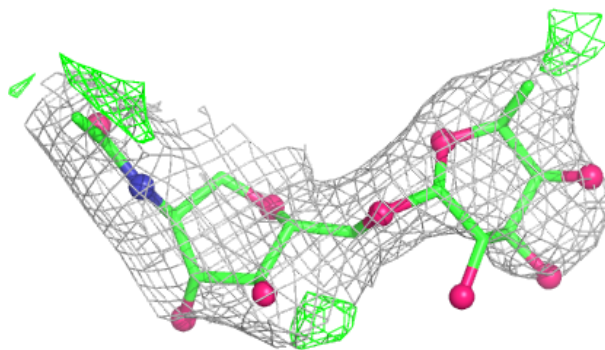
Electron density around Chain G:

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



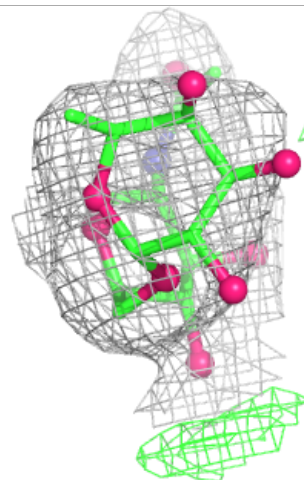
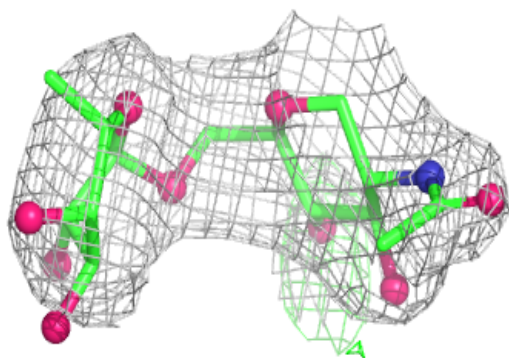
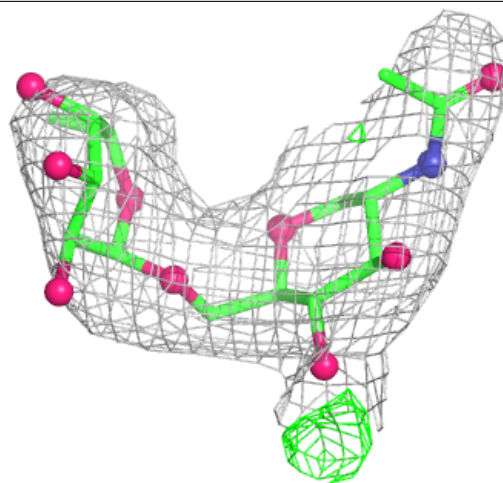
Electron density around Chain I:

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



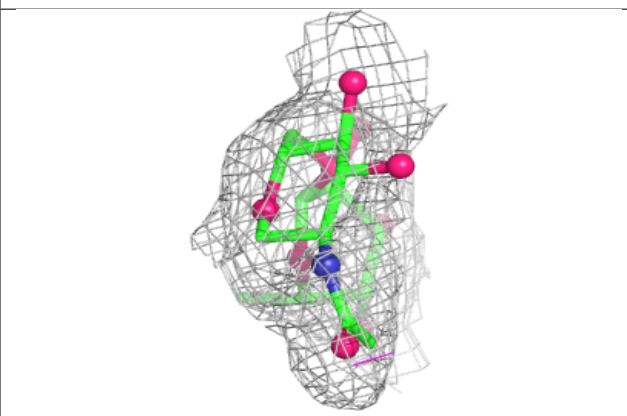
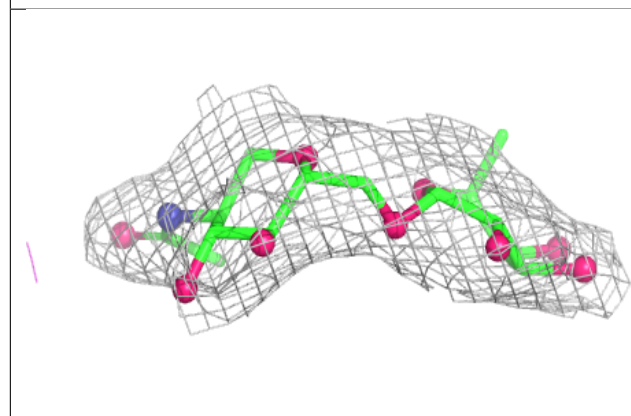
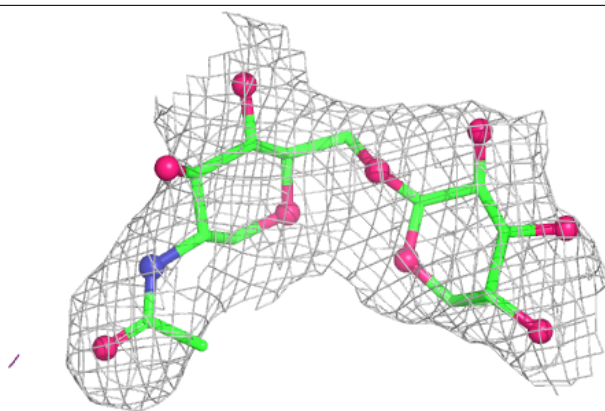
Electron density around Chain J:

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

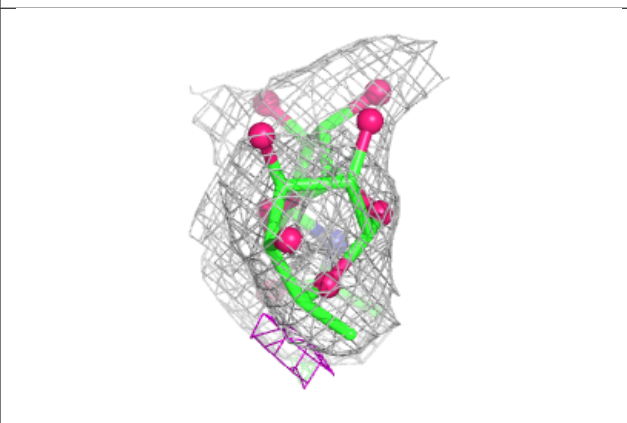
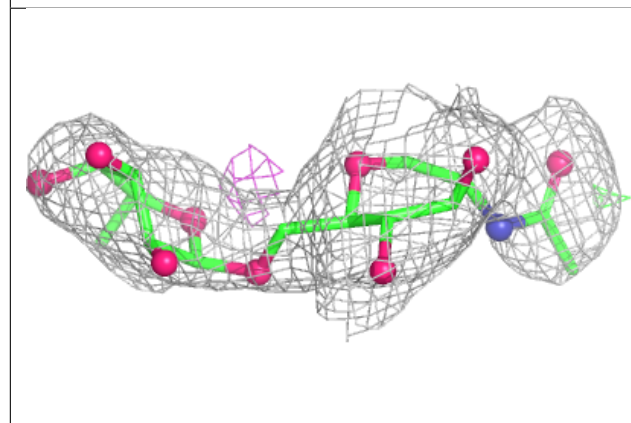
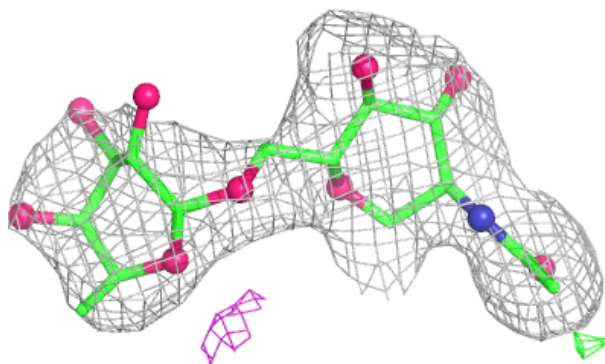


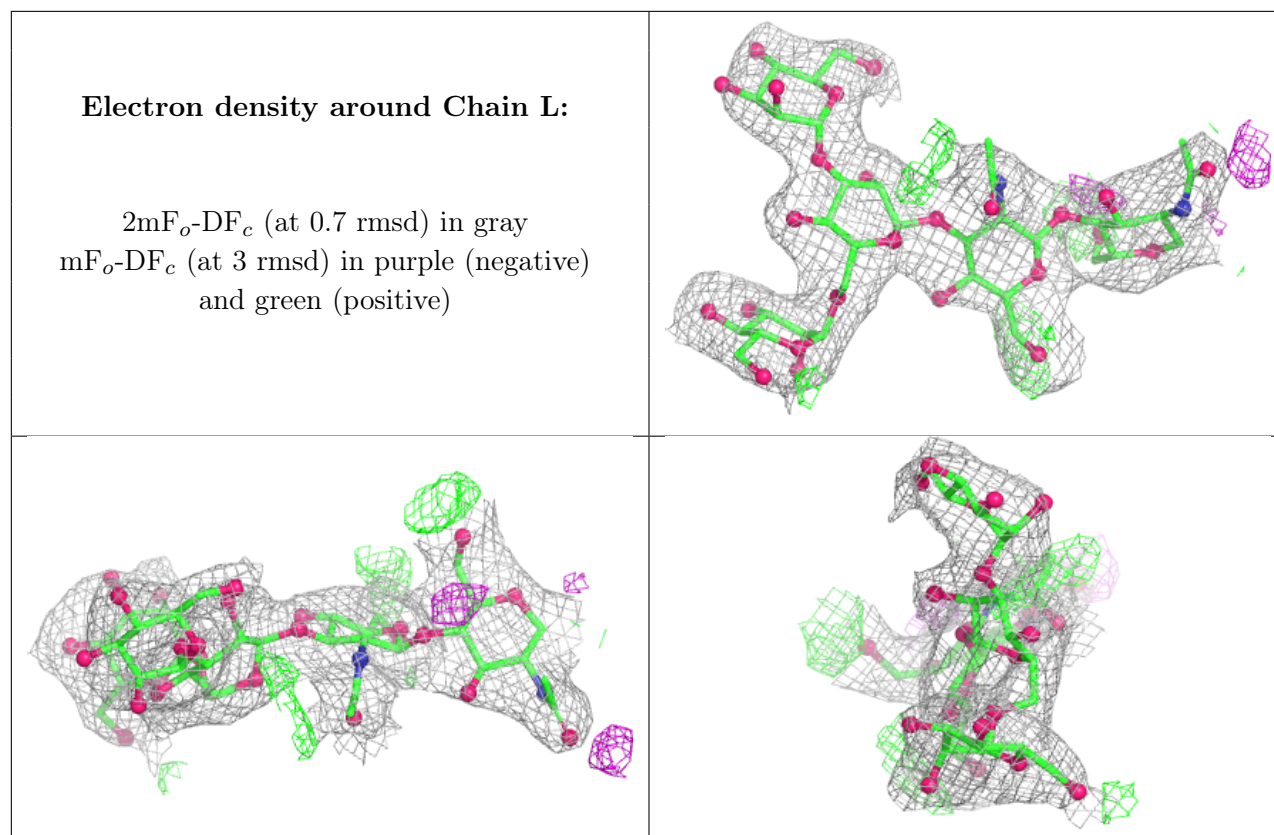
Electron density around Chain M:

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

**Electron density around Chain N:**

$2mF_o-DF_c$ (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($\text{\AA}^2$)	Q<0.9
7	NAG	B	501	14/15	0.38	0.17	68,87,95,95	0
7	NAG	C	503	14/15	0.68	0.20	54,71,85,87	0
7	NAG	B	505	14/15	0.72	0.15	76,87,93,94	0
7	NAG	A	504	14/15	0.72	0.16	41,67,75,76	0
9	KOJ	A	516	10/10	0.82	0.17	50,54,56,62	0
7	NAG	C	511	14/15	0.85	0.12	50,57,67,68	0
7	NAG	D	507	14/15	0.86	0.12	43,59,64,64	0
9	KOJ	B	508	10/10	0.86	0.12	74,75,82,83	0
9	KOJ	C	514	10/10	0.86	0.15	50,54,57,59	0
7	NAG	B	504	14/15	0.88	0.12	59,70,81,82	0
9	KOJ	D	510	10/10	0.88	0.17	45,50,56,56	0
8	ZN	B	506	1/1	0.97	0.04	78,78,78,78	0
8	ZN	A	514	1/1	0.98	0.04	40,40,40,40	0
8	ZN	B	507	1/1	0.98	0.04	68,68,68,68	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	ZN	A	515	1/1	0.98	0.04	51,51,51,51	0
8	ZN	C	513	1/1	0.99	0.03	60,60,60,60	0
8	ZN	D	509	1/1	0.99	0.02	53,53,53,53	0
8	ZN	D	508	1/1	1.00	0.02	39,39,39,39	0
8	ZN	C	512	1/1	1.00	0.01	42,42,42,42	0
8	ZN	C	515	1/1	1.00	0.02	58,58,58,58	0

6.5 Other polymers [i](#)

There are no such residues in this entry.