



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 5, 2026 – 11:17 AM EDT

PDB ID : 4OFY / pdb_00004ofy
Title : Crystal Structure of the Complex of SYG-1 D1-D2 and SYG-2 D1-D4
Authors : Ozkan, E.; Garcia, K.C.
Deposited on : 2014-01-15
Resolution : 3.30 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 : 9.0.015 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

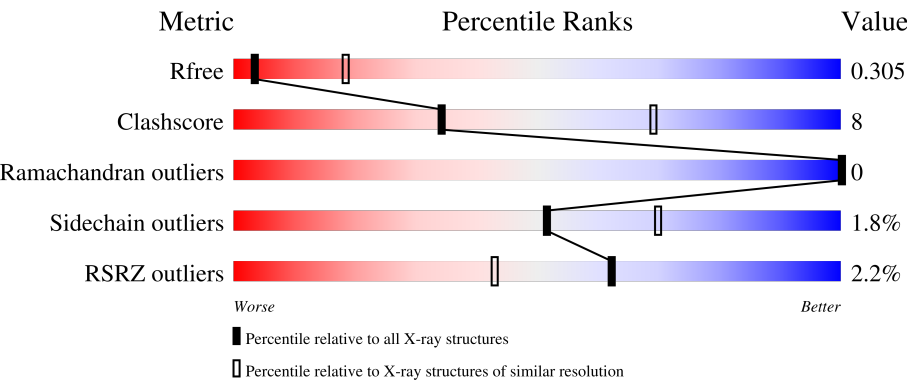
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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	1169 (3.32-3.28)
Clashscore	190562	1209 (3.32-3.28)
Ramachandran outliers	187476	1188 (3.32-3.28)
Sidechain outliers	187428	1187 (3.32-3.28)
RSRZ outliers	180081	1169 (3.32-3.28)

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	262	2% 75% 13% 12%
1	B	262	2% 75% 12% 13%
1	C	262	2% 75% 11% 13%
2	D	420	2% 75% 21% ••
2	E	420	3% 74% 21% ••

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Mol	Chain	Length	Quality of chain
2	F	420	2% 75% 20%
3	G	2	50% 50%
3	H	2	50% 50%
3	I	2	50% 50%
3	J	2	100%
3	K	2	100%
3	L	2	100%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14816 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 Protein SYG-1, isoform b.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	230	Total	C	N	O	S	0	0	0
			1767	1109	310	340	8			
1	B	227	Total	C	N	O	S	0	0	0
			1752	1100	307	337	8			
1	C	227	Total	C	N	O	S	0	0	0
			1752	1100	307	337	8			

There are 27 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	16	ALA	-	expression tag	UNP B1Q236
A	17	ASP	-	expression tag	UNP B1Q236
A	18	PRO	-	expression tag	UNP B1Q236
A	272	HIS	-	expression tag	UNP B1Q236
A	273	HIS	-	expression tag	UNP B1Q236
A	274	HIS	-	expression tag	UNP B1Q236
A	275	HIS	-	expression tag	UNP B1Q236
A	276	HIS	-	expression tag	UNP B1Q236
A	277	HIS	-	expression tag	UNP B1Q236
B	16	ALA	-	expression tag	UNP B1Q236
B	17	ASP	-	expression tag	UNP B1Q236
B	18	PRO	-	expression tag	UNP B1Q236
B	272	HIS	-	expression tag	UNP B1Q236
B	273	HIS	-	expression tag	UNP B1Q236
B	274	HIS	-	expression tag	UNP B1Q236
B	275	HIS	-	expression tag	UNP B1Q236
B	276	HIS	-	expression tag	UNP B1Q236
B	277	HIS	-	expression tag	UNP B1Q236
C	16	ALA	-	expression tag	UNP B1Q236
C	17	ASP	-	expression tag	UNP B1Q236
C	18	PRO	-	expression tag	UNP B1Q236
C	272	HIS	-	expression tag	UNP B1Q236
C	273	HIS	-	expression tag	UNP B1Q236

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Chain	Residue	Modelled	Actual	Comment	Reference
C	274	HIS	-	expression tag	UNP B1Q236
C	275	HIS	-	expression tag	UNP B1Q236
C	276	HIS	-	expression tag	UNP B1Q236
C	277	HIS	-	expression tag	UNP B1Q236

- Molecule 2 is a protein called Protein SYG-2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	407	Total	C	N	O	S	0	0	0
			3045	1912	520	598	15			
2	E	407	Total	C	N	O	S	0	0	0
			3042	1910	519	598	15			
2	F	407	Total	C	N	O	S	0	0	0
			3057	1921	522	599	15			

There are 30 discrepancies between the modelled and reference sequences:

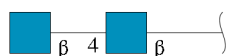
Chain	Residue	Modelled	Actual	Comment	Reference
D	17	ALA	-	expression tag	UNP Q9U3P2
D	18	ASP	-	expression tag	UNP Q9U3P2
D	19	PRO	-	expression tag	UNP Q9U3P2
D	391	CYS	ASN	engineered mutation	UNP Q9U3P2
D	431	HIS	-	expression tag	UNP Q9U3P2
D	432	HIS	-	expression tag	UNP Q9U3P2
D	433	HIS	-	expression tag	UNP Q9U3P2
D	434	HIS	-	expression tag	UNP Q9U3P2
D	435	HIS	-	expression tag	UNP Q9U3P2
D	436	HIS	-	expression tag	UNP Q9U3P2
E	17	ALA	-	expression tag	UNP Q9U3P2
E	18	ASP	-	expression tag	UNP Q9U3P2
E	19	PRO	-	expression tag	UNP Q9U3P2
E	391	CYS	ASN	engineered mutation	UNP Q9U3P2
E	431	HIS	-	expression tag	UNP Q9U3P2
E	432	HIS	-	expression tag	UNP Q9U3P2
E	433	HIS	-	expression tag	UNP Q9U3P2
E	434	HIS	-	expression tag	UNP Q9U3P2
E	435	HIS	-	expression tag	UNP Q9U3P2
E	436	HIS	-	expression tag	UNP Q9U3P2
F	17	ALA	-	expression tag	UNP Q9U3P2
F	18	ASP	-	expression tag	UNP Q9U3P2
F	19	PRO	-	expression tag	UNP Q9U3P2
F	391	CYS	ASN	engineered mutation	UNP Q9U3P2

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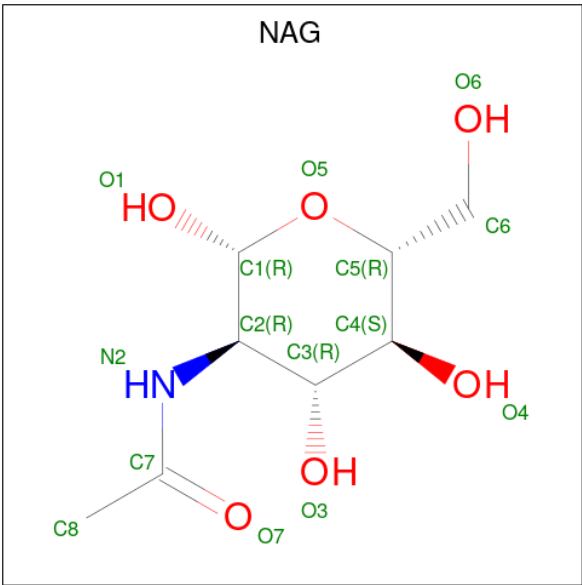
Chain	Residue	Modelled	Actual	Comment	Reference
F	431	HIS	-	expression tag	UNP Q9U3P2
F	432	HIS	-	expression tag	UNP Q9U3P2
F	433	HIS	-	expression tag	UNP Q9U3P2
F	434	HIS	-	expression tag	UNP Q9U3P2
F	435	HIS	-	expression tag	UNP Q9U3P2
F	436	HIS	-	expression tag	UNP Q9U3P2

- 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	G	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	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	J	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	L	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



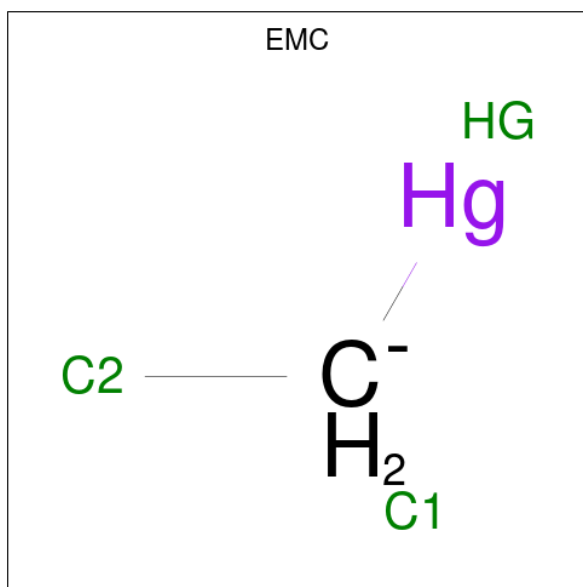
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		
4	E	1	Total	C	N	O	0	0
			14	8	1	5		
4	E	1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

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	F	1	Total	C	N	O	0	0
			14	8	1	5		
4	F	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is ETHYL MERCURY ION (CCD ID: EMC) (formula: C_2H_5Hg).

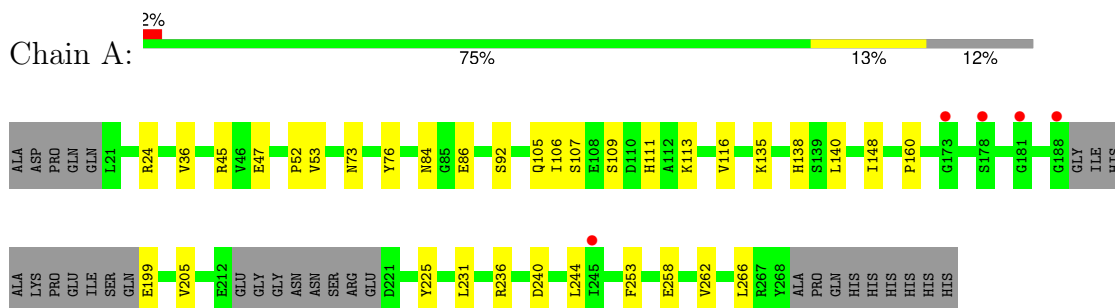


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	D	1	Total	C	Hg	0	0
			3	2	1		
5	E	1	Total	C	Hg	0	0
			3	2	1		
5	F	1	Total	C	Hg	0	0
			3	2	1		

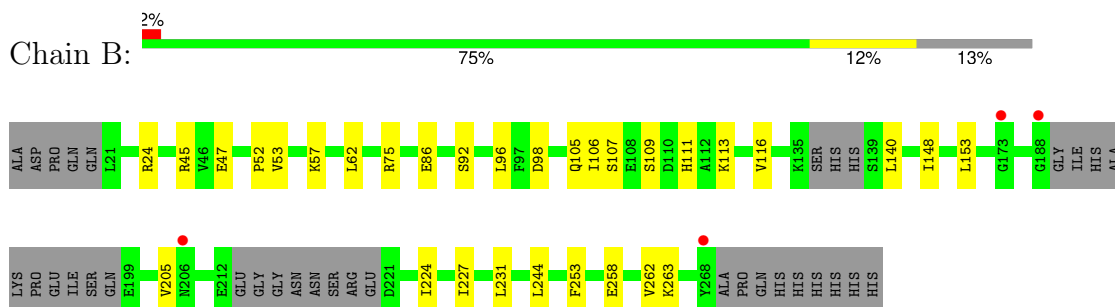
3 Residue-property plots [i](#)

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.

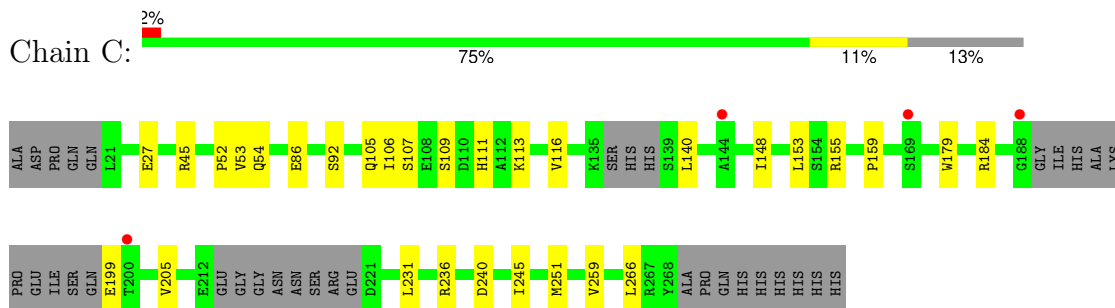
- Molecule 1: Protein SYG-1, isoform b



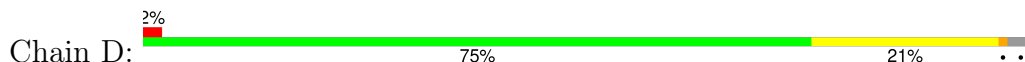
- Molecule 1: Protein SYG-1, isoform b

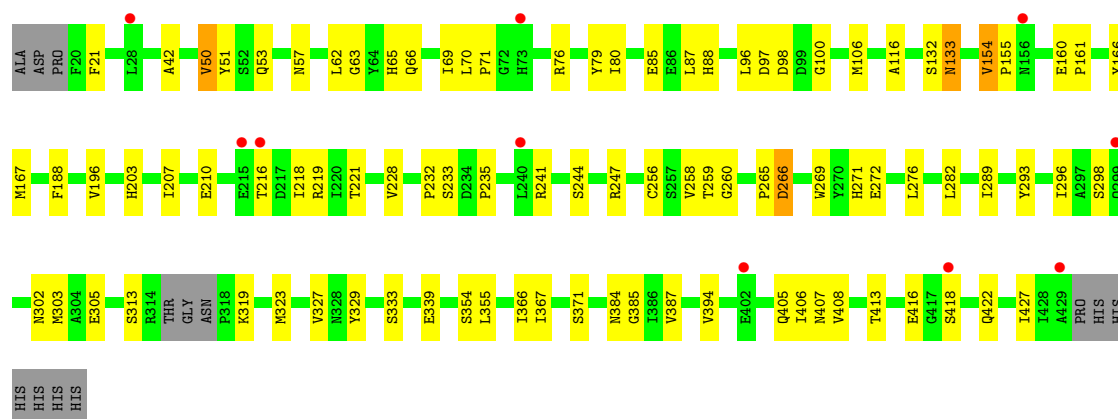


- Molecule 1: Protein SYG-1, isoform b

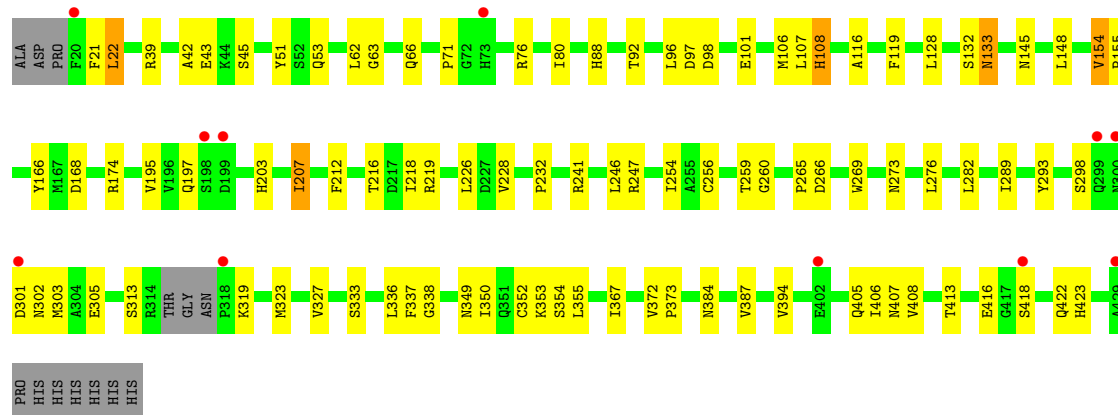
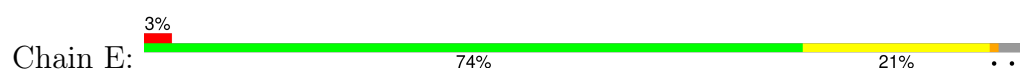


- Molecule 2: Protein SYG-2

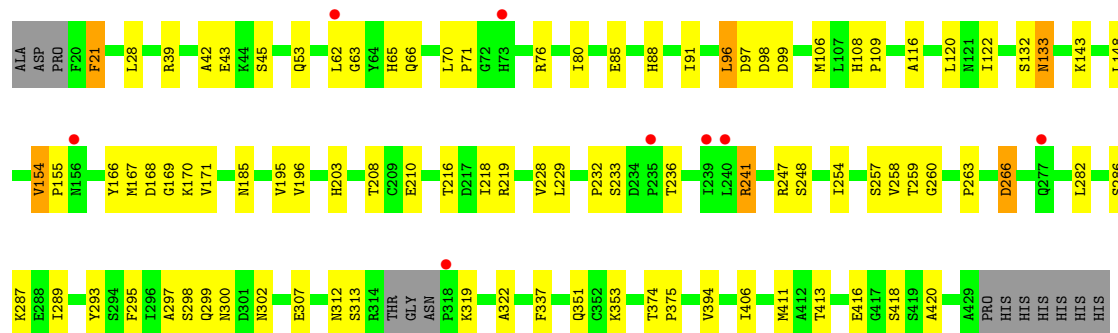
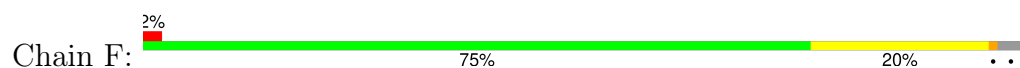




- Molecule 2: Protein SYG-2



- Molecule 2: Protein SYG-2



- Molecule 3: 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)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  50% 50%



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

Chain I:  50% 50%



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

Chain J:  100%



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

Chain K:  100%



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

Chain L:  100%



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	286.79Å 116.14Å 98.03Å 90.00° 107.62° 90.00°	Depositor
Resolution (Å)	41.06 – 3.30 41.06 – 3.30	Depositor EDS
% Data completeness (in resolution range)	95.3 (41.06-3.30) 96.0 (41.06-3.30)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.55 (at 3.32Å)	Xtriage
Refinement program	PHENIX (phenix.refine: dev_1593)	Depositor
R, R_{free}	0.259 , 0.306 0.263 , 0.305	Depositor DCC
R_{free} test set	2266 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	90.9	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 54.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.000 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	14816	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 52.25 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 4.9989e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: NAG, EMC

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	0.26	0/1801	0.71	0/2439
1	B	0.27	0/1785	0.70	0/2415
1	C	0.27	0/1785	0.73	0/2415
2	D	0.32	0/3116	0.85	6/4266 (0.1%)
2	E	0.34	0/3113	0.86	8/4263 (0.2%)
2	F	0.34	0/3128	0.85	2/4280 (0.0%)
All	All	0.31	0/14728	0.81	16/20078 (0.1%)

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	22	LEU	N-CA-C	6.80	119.32	111.02
2	E	303	MET	N-CA-C	-6.80	104.30	112.59
2	D	313	SER	CB-CA-C	-5.95	109.72	116.63
2	F	313	SER	CB-CA-C	-5.75	109.96	116.63
2	E	168	ASP	CB-CA-C	-5.72	109.45	117.23
2	F	66	GLN	CB-CA-C	-5.54	109.69	117.23
2	D	66	GLN	CB-CA-C	-5.45	109.81	117.23
2	E	66	GLN	CB-CA-C	-5.42	109.86	117.23
2	E	108	HIS	CA-C-N	5.40	125.48	119.32
2	E	108	HIS	C-N-CA	5.40	125.48	119.32
2	D	260	GLY	N-CA-C	5.29	118.15	112.33
2	E	174	ARG	N-CA-C	5.25	116.70	110.97
2	D	244	SER	N-CA-C	-5.23	106.23	113.18
2	E	313	SER	CB-CA-C	-5.21	110.59	116.63
2	D	385	GLY	N-CA-C	5.14	117.68	110.38
2	D	427	ILE	N-CA-C	5.10	115.19	107.80

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1767	0	1751	20	0
1	B	1752	0	1744	17	0
1	C	1752	0	1743	18	0
2	D	3045	0	2835	52	2
2	E	3042	0	2820	51	3
2	F	3057	0	2857	62	0
3	G	28	0	25	1	0
3	H	28	0	25	1	0
3	I	28	0	25	1	0
3	J	28	0	25	0	0
3	K	28	0	25	0	0
3	L	28	0	25	0	0
4	C	14	0	13	0	0
4	D	70	0	65	1	0
4	E	70	0	65	0	0
4	F	70	0	65	0	0
5	D	3	0	0	0	0
5	E	3	0	0	0	0
5	F	3	0	0	0	0
All	All	14816	0	14108	219	3

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 8.

All (219) 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)
2:E:282:LEU:HD13	2:E:289:ILE:HG13	1.61	0.82
2:F:39:ARG:HA	2:F:88:HIS:HD2	1.49	0.78
2:E:276:LEU:HD13	2:E:293:TYR:CE2	2.19	0.77
2:F:233:SER:HG	2:F:259:THR:HG1	1.25	0.76
2:F:53:GLN:HG3	2:F:63:GLY:HA3	1.68	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:140:LEU:HD21	1:A:148:ILE:HG21	1.67	0.75
2:F:21:PHE:HA	2:F:43:GLU:HG3	1.67	0.74
1:A:53:VAL:HG22	1:A:106:ILE:HG12	1.71	0.72
2:F:96:LEU:HA	2:F:122:ILE:HD13	1.72	0.72
2:F:21:PHE:HZ	2:F:116:ALA:HB2	1.54	0.71
2:D:53:GLN:HG3	2:D:63:GLY:HA3	1.74	0.70
2:D:216:THR:HB	2:D:218:ILE:HG22	1.73	0.70
2:D:233:SER:HG	2:D:259:THR:HG1	1.35	0.70
2:F:167:MET:HE1	2:F:196:VAL:HG21	1.72	0.69
1:B:53:VAL:HG22	1:B:106:ILE:HG12	1.73	0.69
2:D:282:LEU:HD13	2:D:289:ILE:HG13	1.74	0.68
2:F:282:LEU:HD13	2:F:289:ILE:HG13	1.74	0.68
2:D:62:LEU:HG	2:D:70:LEU:HD23	1.76	0.68
2:F:216:THR:HB	2:F:218:ILE:HG22	1.75	0.67
2:E:216:THR:HB	2:E:218:ILE:HG22	1.75	0.67
1:B:140:LEU:HD21	1:B:148:ILE:HG21	1.76	0.66
2:D:21:PHE:HB2	2:D:42:ALA:HA	1.75	0.66
2:F:21:PHE:HB2	2:F:42:ALA:HA	1.77	0.66
2:E:21:PHE:HZ	2:E:116:ALA:HB2	1.61	0.66
2:E:21:PHE:HB2	2:E:42:ALA:HA	1.77	0.66
2:D:62:LEU:HD21	2:D:71:PRO:HD2	1.78	0.65
2:E:53:GLN:HG3	2:E:63:GLY:HA3	1.78	0.65
1:C:52:PRO:HD2	1:C:107:SER:HB2	1.80	0.64
2:D:276:LEU:HD13	2:D:293:TYR:CE2	2.34	0.62
2:D:247:ARG:NH1	2:D:416:GLU:OE2	2.32	0.62
2:E:51:TYR:HB2	2:E:107:LEU:HD13	1.81	0.62
1:A:52:PRO:HD2	1:A:107:SER:HB2	1.81	0.62
2:D:413:THR:HG23	2:D:418:SER:HB3	1.82	0.62
2:F:108:HIS:HD2	2:F:109:PRO:HD2	1.62	0.62
2:F:96:LEU:HD23	2:F:97:ASP:H	1.65	0.62
2:D:298:SER:N	2:D:302:ASN:OD1	2.25	0.61
1:A:24:ARG:NH1	1:A:47:GLU:OE2	2.34	0.61
2:E:39:ARG:HA	2:E:88:HIS:HD2	1.66	0.60
2:E:413:THR:HG23	2:E:418:SER:HB3	1.84	0.60
2:D:21:PHE:HZ	2:D:116:ALA:HB2	1.67	0.59
2:F:413:THR:HG23	2:F:418:SER:HB3	1.85	0.59
1:B:205:VAL:HG22	1:B:231:LEU:HA	1.85	0.59
2:D:210:GLU:CD	2:D:219:ARG:HD2	2.28	0.59
2:F:80:ILE:HB	2:F:88:HIS:HB2	1.84	0.59
1:C:53:VAL:HG22	1:C:106:ILE:HG12	1.84	0.58
1:B:45:ARG:HG2	1:B:86:GLU:HG2	1.84	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:407:ASN:HB3	2:E:422:GLN:NE2	2.18	0.58
2:F:21:PHE:CZ	2:F:116:ALA:HB2	2.37	0.58
1:B:109:SER:O	1:B:111:HIS:N	2.36	0.57
2:E:247:ARG:NH1	2:E:416:GLU:OE2	2.37	0.57
1:C:140:LEU:HD21	1:C:148:ILE:HG21	1.87	0.57
2:D:42:ALA:HB2	2:D:106:MET:HE3	1.87	0.57
2:E:96:LEU:HD23	2:E:97:ASP:H	1.69	0.57
2:D:266:ASP:OD1	2:D:266:ASP:N	2.30	0.56
2:F:62:LEU:HD21	2:F:71:PRO:HD2	1.87	0.56
1:C:199:GLU:O	1:C:236:ARG:NH1	2.39	0.55
2:D:210:GLU:HG3	2:D:221:THR:HG22	1.88	0.55
1:B:140:LEU:N	1:B:263:LYS:O	2.36	0.55
2:D:256:CYS:HB2	2:D:269:TRP:CZ2	2.42	0.55
2:D:394:VAL:HG11	2:D:406:ILE:HG21	1.89	0.54
1:C:109:SER:O	1:C:111:HIS:N	2.41	0.54
2:E:336:LEU:HD13	2:E:423:HIS:HB2	1.90	0.54
1:A:199:GLU:O	1:A:236:ARG:NH1	2.40	0.53
1:A:45:ARG:HG2	1:A:86:GLU:HG2	1.89	0.53
2:D:271:HIS:CD2	2:D:272:GLU:HG3	2.44	0.53
2:F:210:GLU:CD	2:F:219:ARG:HD2	2.34	0.53
2:F:247:ARG:NH1	2:F:416:GLU:OE2	2.41	0.53
2:D:63:GLY:O	2:D:65:HIS:ND1	2.38	0.53
1:C:45:ARG:HG2	1:C:86:GLU:HG2	1.91	0.52
2:D:232:PRO:O	2:D:319:LYS:HD2	2.09	0.52
2:D:367:ILE:HD12	2:D:408:VAL:HG22	1.90	0.52
2:E:232:PRO:O	2:E:319:LYS:HD2	2.10	0.52
2:E:232:PRO:HG3	2:E:265:PRO:HG3	1.91	0.52
2:D:296:ILE:HD11	4:D:1506:NAG:H82	1.92	0.51
2:F:337:PHE:HE2	2:F:353:LYS:HG2	1.76	0.51
2:E:80:ILE:HB	2:E:88:HIS:HB2	1.93	0.51
1:A:109:SER:O	1:A:111:HIS:N	2.40	0.51
2:F:166:TYR:CE2	2:F:171:VAL:HG12	2.46	0.51
1:C:105:GLN:HG3	1:C:116:VAL:HG22	1.93	0.51
2:D:235:PRO:HB3	2:D:258:VAL:HG22	1.92	0.51
2:E:212:PHE:CD1	2:E:219:ARG:HD3	2.46	0.51
2:F:39:ARG:HA	2:F:88:HIS:CD2	2.39	0.51
2:F:229:LEU:HD22	2:F:263:PRO:HD3	1.92	0.51
2:E:62:LEU:HD21	2:E:71:PRO:HD2	1.93	0.51
2:F:286:SER:O	2:F:287:LYS:HG2	2.11	0.51
2:D:167:MET:HE1	2:D:196:VAL:HG21	1.93	0.50
2:F:411:MET:HG3	2:F:420:ALA:HB2	1.93	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:92:SER:HB2	3:I:1:NAG:H82	1.94	0.50
2:E:407:ASN:HB3	2:E:422:GLN:HE21	1.75	0.50
2:F:42:ALA:HB2	2:F:106:MET:HE3	1.93	0.50
2:F:21:PHE:CD1	2:F:21:PHE:N	2.79	0.50
2:E:203:HIS:HB2	2:E:228:VAL:HG22	1.93	0.50
1:A:205:VAL:HG22	1:A:231:LEU:HA	1.93	0.49
2:E:305:GLU:HA	2:E:323:MET:O	2.11	0.49
2:D:203:HIS:HB2	2:D:228:VAL:HG22	1.94	0.49
2:F:298:SER:O	2:F:302:ASN:HB2	2.13	0.49
1:B:92:SER:HB2	3:H:1:NAG:H82	1.94	0.48
2:F:233:SER:OG	2:F:259:THR:OG1	2.08	0.48
2:F:232:PRO:O	2:F:319:LYS:HD2	2.14	0.48
1:B:52:PRO:HD2	1:B:107:SER:HB2	1.96	0.48
2:E:39:ARG:HA	2:E:88:HIS:CD2	2.48	0.48
2:F:210:GLU:OE2	2:F:219:ARG:HD2	2.13	0.48
2:F:236:THR:OG1	2:F:257:SER:HB3	2.13	0.48
2:F:394:VAL:HG11	2:F:406:ILE:HG21	1.96	0.48
1:C:27:GLU:OE1	1:C:45:ARG:NE	2.36	0.47
2:E:22:LEU:H	2:E:43:GLU:N	2.11	0.47
2:D:50:VAL:HG23	2:D:51:TYR:CD2	2.49	0.47
2:D:333:SER:H	2:D:355:LEU:HD12	1.80	0.47
2:F:45:SER:HB3	2:F:108:HIS:NE2	2.30	0.47
1:C:205:VAL:HG22	1:C:231:LEU:HA	1.97	0.47
1:C:54:GLN:NE2	2:F:53:GLN:OE1	2.44	0.47
1:B:24:ARG:NH1	1:B:47:GLU:OE2	2.46	0.47
2:D:305:GLU:HA	2:D:323:MET:O	2.15	0.47
2:E:42:ALA:HB2	2:E:106:MET:HE3	1.97	0.47
2:F:76:ARG:NH1	2:F:98:ASP:OD2	2.48	0.47
2:F:299:GLN:HG3	2:F:300:ASN:H	1.80	0.47
1:A:73:ASN:HB3	1:A:76:TYR:CG	2.50	0.46
1:C:240:ASP:HB2	1:C:266:LEU:HD22	1.97	0.46
2:D:302:ASN:HB3	2:D:327:VAL:HG21	1.97	0.46
2:E:337:PHE:HE2	2:E:353:LYS:HG2	1.80	0.46
2:F:203:HIS:HB2	2:F:228:VAL:HG22	1.97	0.46
2:F:307:GLU:HA	2:F:322:ALA:HA	1.97	0.46
2:E:354:SER:O	2:E:387:VAL:HG13	2.16	0.46
2:E:298:SER:O	2:E:302:ASN:HB2	2.15	0.46
2:E:132:SER:HA	2:E:133:ASN:HA	1.49	0.46
2:F:99:ASP:HB2	2:F:122:ILE:HD12	1.97	0.46
2:D:354:SER:O	2:D:387:VAL:HG13	2.16	0.46
2:E:101:GLU:HB3	2:E:119:PHE:CE2	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:259:THR:OG1	2:E:260:GLY:N	2.48	0.46
2:F:298:SER:N	2:F:302:ASN:OD1	2.39	0.45
1:B:75:ARG:HH22	1:B:98:ASP:CG	2.24	0.45
2:F:63:GLY:O	2:F:65:HIS:ND1	2.45	0.45
1:C:153:LEU:HD21	1:C:155:ARG:HD2	1.99	0.45
2:E:96:LEU:C	2:E:98:ASP:H	2.23	0.45
2:E:154:VAL:HA	2:E:155:PRO:HD3	1.75	0.45
2:E:246:LEU:HB2	2:E:327:VAL:HG22	1.98	0.45
2:F:185:ASN:OD1	2:F:185:ASN:N	2.49	0.45
1:A:240:ASP:HB2	1:A:266:LEU:HD22	1.99	0.45
2:D:232:PRO:HG3	2:D:265:PRO:HG3	1.99	0.45
2:E:254:ILE:HG13	2:E:293:TYR:HB3	1.99	0.45
1:A:253:PHE:CE2	1:A:258:GLU:HG3	2.52	0.45
2:D:166:TYR:O	2:D:207:ILE:HA	2.17	0.45
1:A:240:ASP:HA	1:A:266:LEU:HD13	1.99	0.44
1:B:153:LEU:HD12	1:B:227:ILE:O	2.17	0.44
1:B:253:PHE:CE2	1:B:258:GLU:HG3	2.52	0.44
2:F:374:THR:HA	2:F:375:PRO:HD3	1.81	0.44
1:B:96:LEU:HD11	1:B:224:ILE:HG23	1.99	0.44
2:F:154:VAL:HA	2:F:155:PRO:HD3	1.83	0.44
2:F:299:GLN:HG3	2:F:300:ASN:N	2.32	0.44
2:D:303:MET:HG3	2:D:384:ASN:HA	1.99	0.44
2:D:160:GLU:HB3	2:D:188:PHE:CG	2.52	0.44
2:D:80:ILE:HB	2:D:88:HIS:HB2	1.99	0.44
2:D:329:TYR:OH	2:D:384:ASN:O	2.30	0.44
2:F:248:SER:O	2:F:297:ALA:HB3	2.17	0.44
1:A:84:ASN:HB3	1:A:86:GLU:HG3	1.99	0.44
1:C:245:ILE:HG23	1:C:259:VAL:HG13	1.99	0.43
2:D:96:LEU:HD23	2:D:97:ASP:H	1.83	0.43
2:D:366:ILE:HG12	2:D:371:SER:HA	2.00	0.43
2:E:336:LEU:HD23	2:E:352:CYS:HB2	1.99	0.43
2:D:96:LEU:C	2:D:98:ASP:H	2.26	0.43
2:E:333:SER:H	2:E:355:LEU:HD12	1.82	0.43
2:F:28:LEU:O	2:F:120:LEU:HA	2.18	0.43
2:F:254:ILE:HG13	2:F:293:TYR:HB3	2.00	0.43
2:F:132:SER:HA	2:F:133:ASN:HA	1.53	0.43
2:F:241:ARG:HE	2:F:241:ARG:HB2	1.38	0.43
2:E:21:PHE:CZ	2:E:116:ALA:HB2	2.49	0.43
2:E:76:ARG:HB3	2:E:92:THR:O	2.19	0.43
2:D:160:GLU:HA	2:D:161:PRO:HD3	1.86	0.43
2:E:256:CYS:HB2	2:E:269:TRP:CZ2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:258:VAL:HG12	2:F:260:GLY:H	1.83	0.43
2:D:333:SER:N	2:D:355:LEU:HD12	2.34	0.43
2:E:301:ASP:H	2:E:384:ASN:ND2	2.17	0.43
2:D:339:GLU:H	2:D:339:GLU:HG3	1.64	0.42
1:A:105:GLN:HG3	1:A:116:VAL:HG22	2.02	0.42
2:D:154:VAL:HA	2:D:155:PRO:HD3	1.78	0.42
2:F:293:TYR:HE1	2:F:295:PHE:HB3	1.84	0.42
2:F:143:LYS:HA	2:F:229:LEU:HB2	2.01	0.42
2:E:45:SER:HB3	2:E:108:HIS:NE2	2.35	0.42
1:A:105:GLN:HE21	1:A:106:ILE:C	2.27	0.42
2:E:372:VAL:HB	2:E:373:PRO:HD2	2.01	0.42
2:F:62:LEU:HG	2:F:70:LEU:HD23	2.01	0.42
2:F:166:TYR:HB2	2:F:208:THR:HB	2.02	0.42
2:F:232:PRO:HD3	2:F:312:ASN:ND2	2.34	0.42
1:C:159:PRO:HG2	1:C:251:MET:HE2	2.02	0.41
2:F:148:LEU:O	2:F:195:VAL:HA	2.19	0.41
2:F:266:ASP:OD1	2:F:266:ASP:N	2.43	0.41
2:D:271:HIS:CD2	2:D:271:HIS:C	2.98	0.41
2:E:21:PHE:CD1	2:E:21:PHE:N	2.89	0.41
2:E:96:LEU:HD23	2:E:97:ASP:N	2.34	0.41
2:E:333:SER:N	2:E:355:LEU:HD12	2.35	0.41
1:A:92:SER:HB2	3:G:1:NAG:H82	2.02	0.41
2:E:338:GLY:HA3	2:E:350:ILE:HD13	2.03	0.41
1:A:105:GLN:HG2	1:A:106:ILE:N	2.34	0.41
2:F:96:LEU:C	2:F:98:ASP:H	2.27	0.41
1:B:57:LYS:HB3	1:B:62:LEU:HD21	2.03	0.41
1:B:244:LEU:HB3	1:B:262:VAL:HG12	2.02	0.41
2:F:96:LEU:HD23	2:F:97:ASP:N	2.32	0.41
2:E:394:VAL:HG11	2:E:406:ILE:HG21	2.03	0.41
1:B:105:GLN:HG2	1:B:106:ILE:N	2.35	0.41
1:C:179:TRP:CD1	1:C:184:ARG:HB2	2.55	0.41
2:D:407:ASN:HB3	2:D:422:GLN:NE2	2.35	0.41
1:A:244:LEU:HB3	1:A:262:VAL:HG12	2.03	0.41
1:C:105:GLN:HE21	1:C:106:ILE:C	2.29	0.41
1:C:105:GLN:HG2	1:C:106:ILE:N	2.36	0.41
2:D:76:ARG:NH1	2:D:98:ASP:OD2	2.52	0.41
1:A:160:PRO:HG3	1:A:225:TYR:CD1	2.56	0.40
1:B:105:GLN:HG3	1:B:116:VAL:HG22	2.03	0.40
2:D:21:PHE:CZ	2:D:116:ALA:HB2	2.51	0.40
2:D:210:GLU:OE2	2:D:219:ARG:HD2	2.20	0.40
2:F:337:PHE:HB2	2:F:351:GLN:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:57:ASN:HB2	2:D:100:GLY:HA3	2.03	0.40
2:E:148:LEU:O	2:E:195:VAL:HA	2.21	0.40
2:F:168:ASP:HB3	2:F:169:GLY:H	1.62	0.40
2:D:79:TYR:CD2	2:D:87:LEU:HD11	2.56	0.40
2:D:132:SER:HA	2:D:133:ASN:HA	1.47	0.40
2:E:166:TYR:O	2:E:207:ILE:HA	2.20	0.40
2:E:226:LEU:HD12	2:E:226:LEU:HA	1.85	0.40
2:E:367:ILE:HD12	2:E:408:VAL:HG22	2.03	0.40
1:A:135:LYS:HD2	1:A:138:HIS:HA	2.03	0.40

All (3) 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)
2:D:405:GLN:NE2	2:E:422:GLN:OE1[1_452]	2.09	0.11
2:D:422:GLN:OE1	2:E:405:GLN:NE2[1_452]	2.09	0.11
2:E:145:ASN:O	2:E:273:ASN:ND2[2_658]	2.10	0.10

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	224/262 (86%)	212 (95%)	12 (5%)	0	100	100
1	B	219/262 (84%)	208 (95%)	11 (5%)	0	100	100
1	C	219/262 (84%)	208 (95%)	11 (5%)	0	100	100
2	D	403/420 (96%)	382 (95%)	21 (5%)	0	100	100
2	E	403/420 (96%)	382 (95%)	21 (5%)	0	100	100
2	F	403/420 (96%)	383 (95%)	20 (5%)	0	100	100
All	All	1871/2046 (91%)	1775 (95%)	96 (5%)	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	197/228 (86%)	195 (99%)	2 (1%)	68	76
1	B	197/228 (86%)	196 (100%)	1 (0%)	81	83
1	C	197/228 (86%)	196 (100%)	1 (0%)	81	83
2	D	331/380 (87%)	324 (98%)	7 (2%)	47	67
2	E	329/380 (87%)	321 (98%)	8 (2%)	43	65
2	F	333/380 (88%)	324 (97%)	9 (3%)	39	63
All	All	1584/1824 (87%)	1556 (98%)	28 (2%)	51	70

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	VAL
1	A	113	LYS
1	B	113	LYS
1	C	113	LYS
2	D	50	VAL
2	D	69	ILE
2	D	85	GLU
2	D	133	ASN
2	D	154	VAL
2	D	241	ARG
2	D	266	ASP
2	E	128	LEU
2	E	133	ASN
2	E	154	VAL
2	E	197	GLN
2	E	207	ILE
2	E	241	ARG
2	E	266	ASP
2	E	349	ASN

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Mol	Chain	Res	Type
2	F	21	PHE
2	F	85	GLU
2	F	91	ILE
2	F	96	LEU
2	F	133	ASN
2	F	154	VAL
2	F	170	LYS
2	F	241	ARG
2	F	266	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
2	D	242	ASN
2	D	262	ASN
2	D	328	ASN
2	D	362	GLN
2	D	405	GLN
2	D	422	GLN
2	E	88	HIS
2	E	242	ASN
2	E	262	ASN
2	E	271	HIS
2	E	328	ASN
2	E	362	GLN
2	E	405	GLN
2	E	422	GLN
2	F	88	HIS

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 ⓘ

12 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	$\# Z > 2$
3	NAG	G	1	1,3	14,14,15	0.41	0	17,19,21	0.58	0
3	NAG	G	2	3	14,14,15	0.23	0	17,19,21	0.38	0
3	NAG	H	1	1,3	14,14,15	0.43	0	17,19,21	0.52	0
3	NAG	H	2	3	14,14,15	0.23	0	17,19,21	0.37	0
3	NAG	I	1	1,3	14,14,15	0.39	0	17,19,21	0.54	0
3	NAG	I	2	3	14,14,15	0.29	0	17,19,21	0.40	0
3	NAG	J	1	2,3	14,14,15	0.29	0	17,19,21	0.49	0
3	NAG	J	2	3	14,14,15	0.30	0	17,19,21	0.50	0
3	NAG	K	1	2,3	14,14,15	0.43	0	17,19,21	0.47	0
3	NAG	K	2	3	14,14,15	0.33	0	17,19,21	0.51	0
3	NAG	L	1	2,3	14,14,15	0.20	0	17,19,21	0.53	0
3	NAG	L	2	3	14,14,15	0.25	0	17,19,21	0.43	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
3	NAG	G	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
3	NAG	H	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	H	2	3	-	1/6/23/26	0/1/1/1
3	NAG	I	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
3	NAG	J	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	J	2	3	-	2/6/23/26	0/1/1/1
3	NAG	K	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	K	2	3	-	2/6/23/26	0/1/1/1
3	NAG	L	1	2,3	-	2/6/23/26	0/1/1/1
3	NAG	L	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

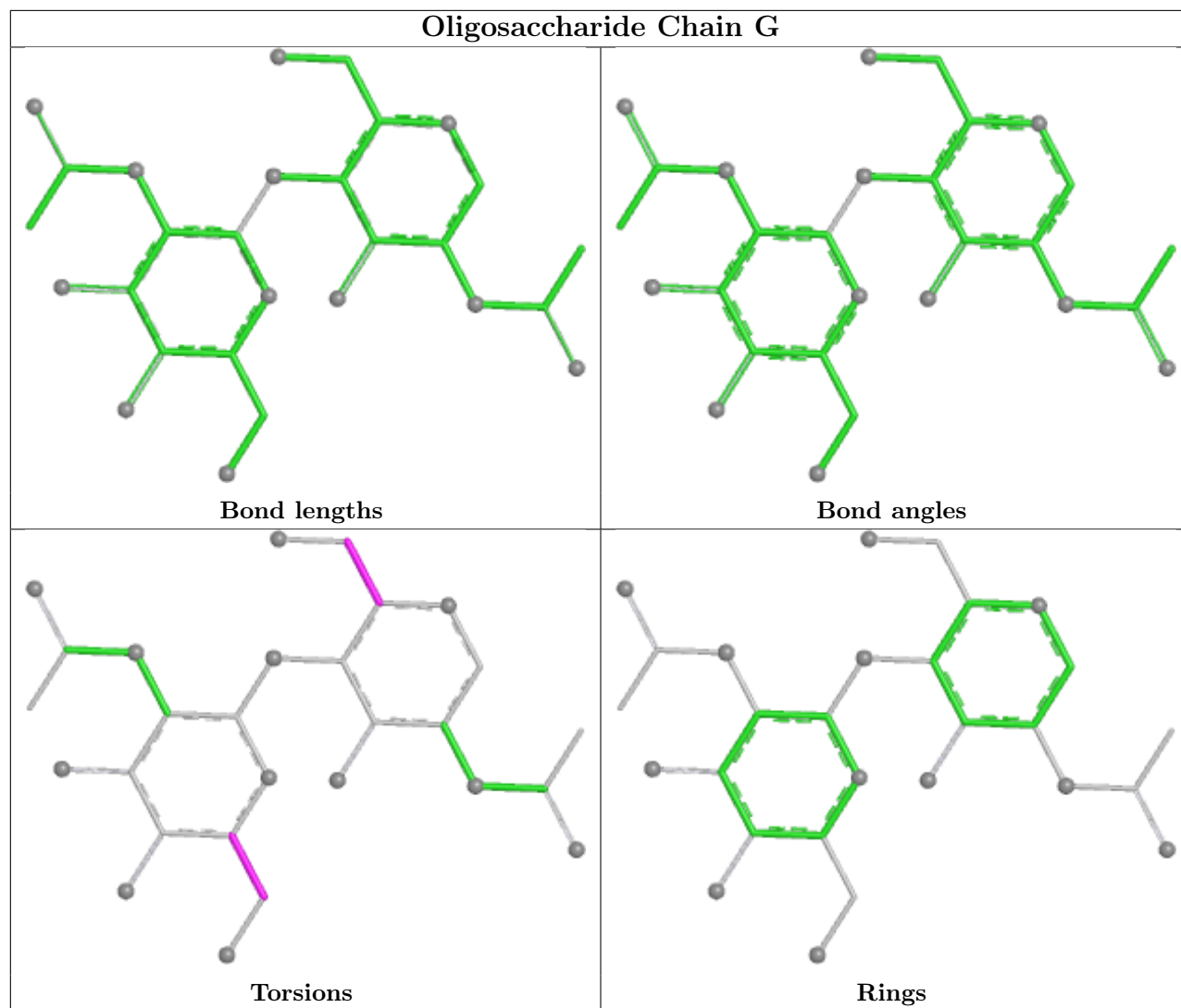
Mol	Chain	Res	Type	Atoms
3	J	2	NAG	O5-C5-C6-O6
3	K	2	NAG	O5-C5-C6-O6
3	H	1	NAG	O5-C5-C6-O6
3	J	1	NAG	O5-C5-C6-O6
3	K	1	NAG	O5-C5-C6-O6
3	I	1	NAG	O5-C5-C6-O6
3	K	2	NAG	C4-C5-C6-O6
3	J	1	NAG	C4-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
3	K	1	NAG	C4-C5-C6-O6
3	J	2	NAG	C4-C5-C6-O6
3	G	1	NAG	O5-C5-C6-O6
3	H	1	NAG	C4-C5-C6-O6
3	I	1	NAG	C4-C5-C6-O6
3	L	1	NAG	C4-C5-C6-O6
3	L	1	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
3	H	2	NAG	C4-C5-C6-O6

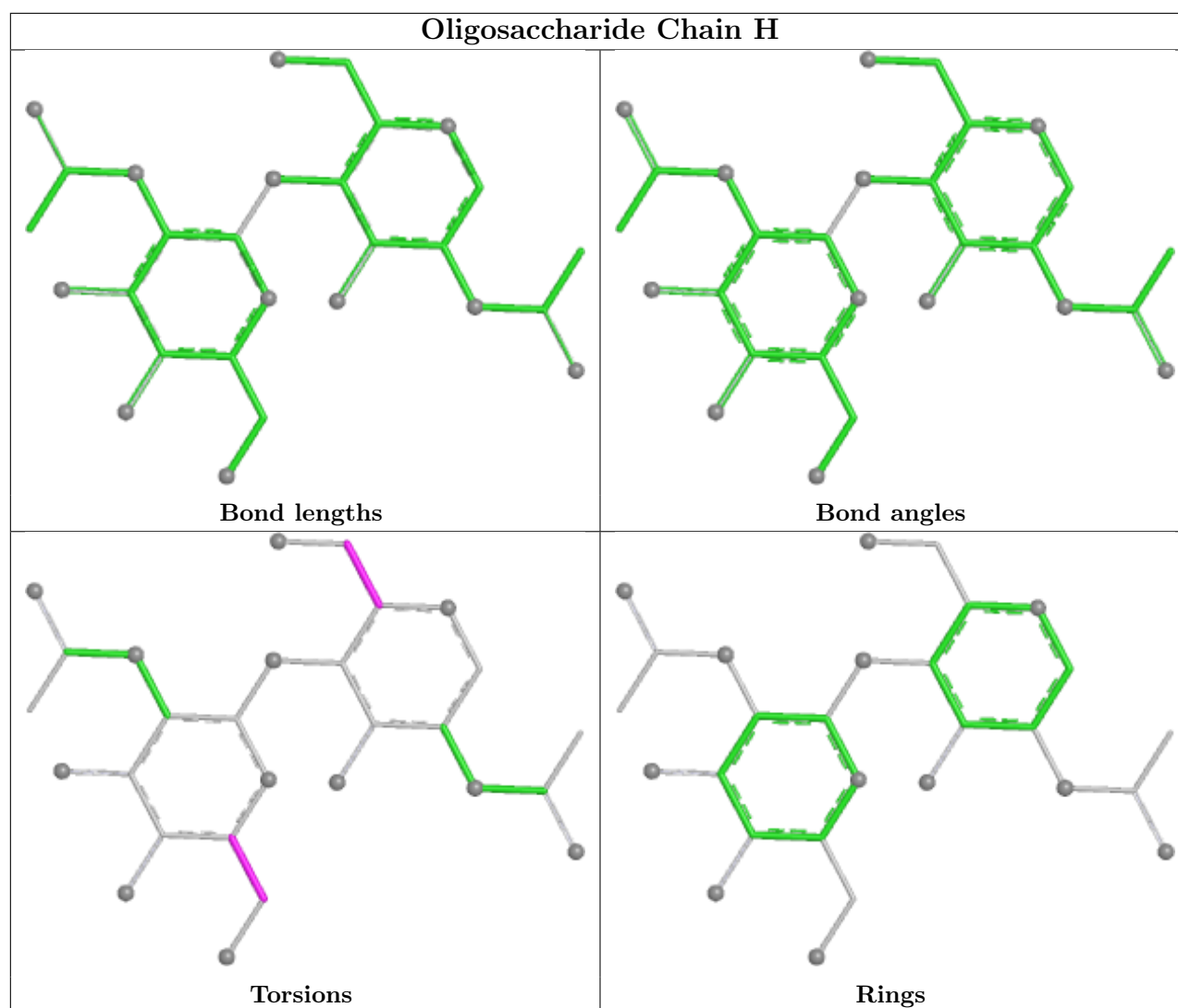
There are no ring outliers.

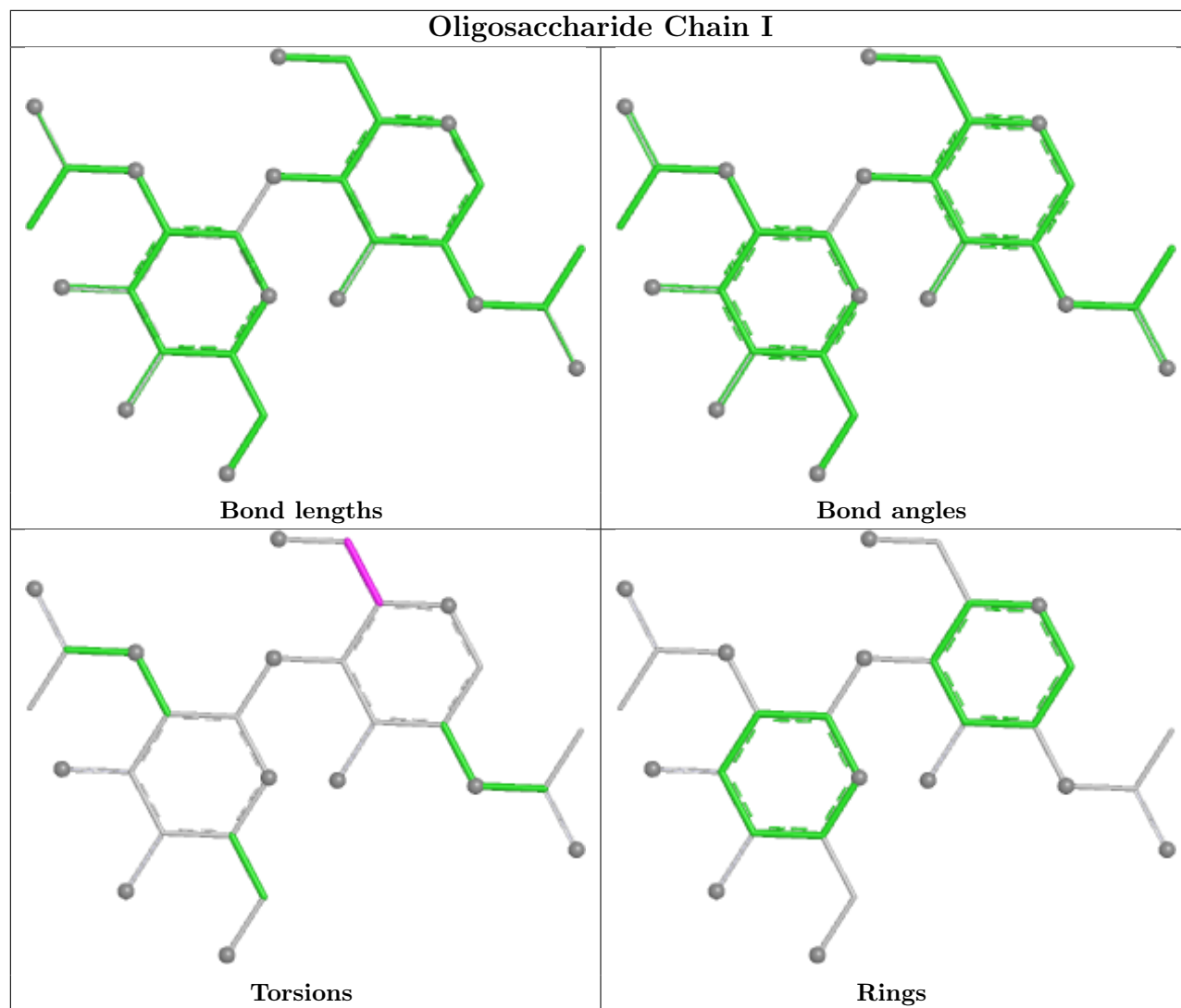
3 monomers are involved in 3 short contacts:

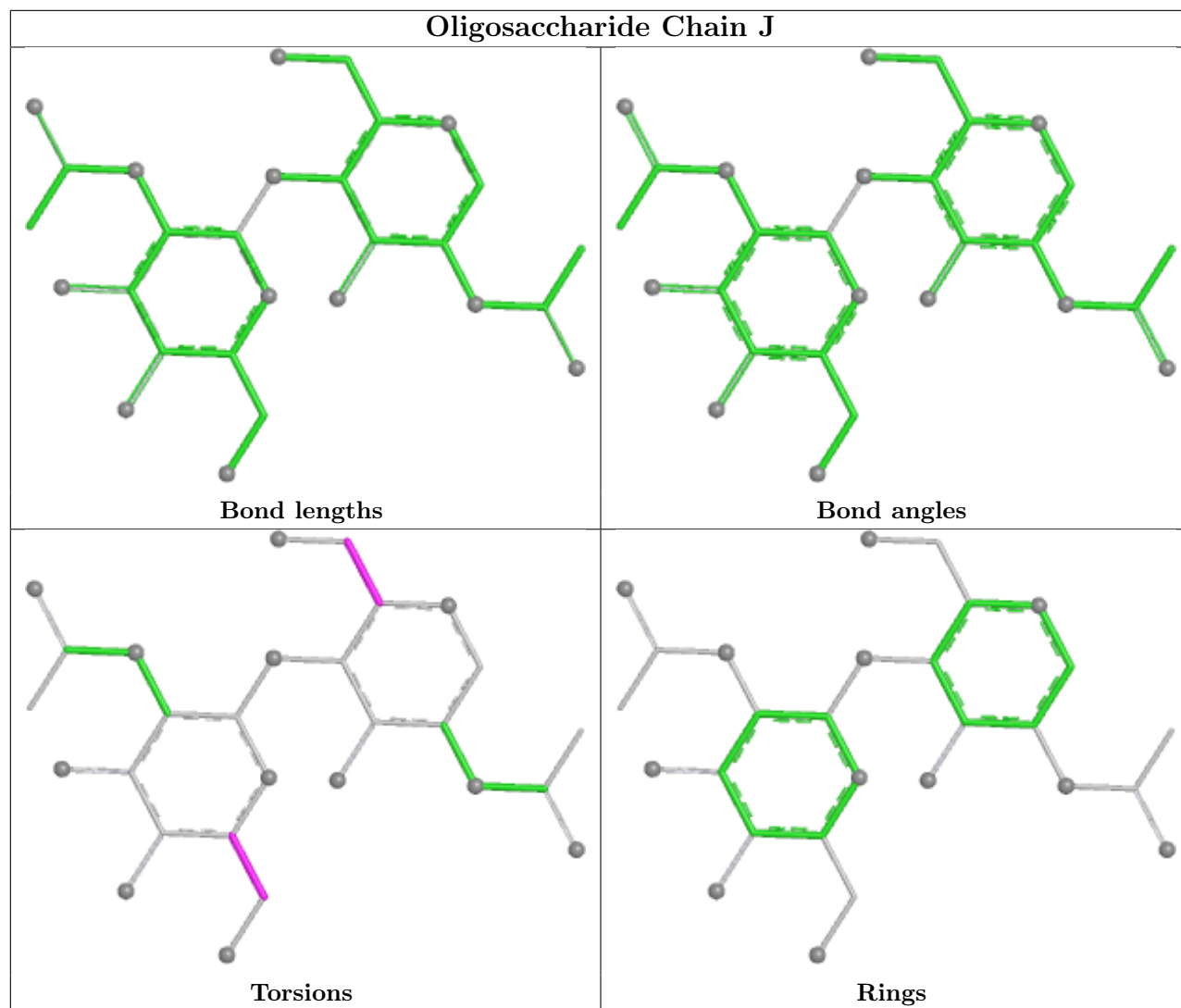
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	1	NAG	1	0
3	H	1	NAG	1	0
3	I	1	NAG	1	0

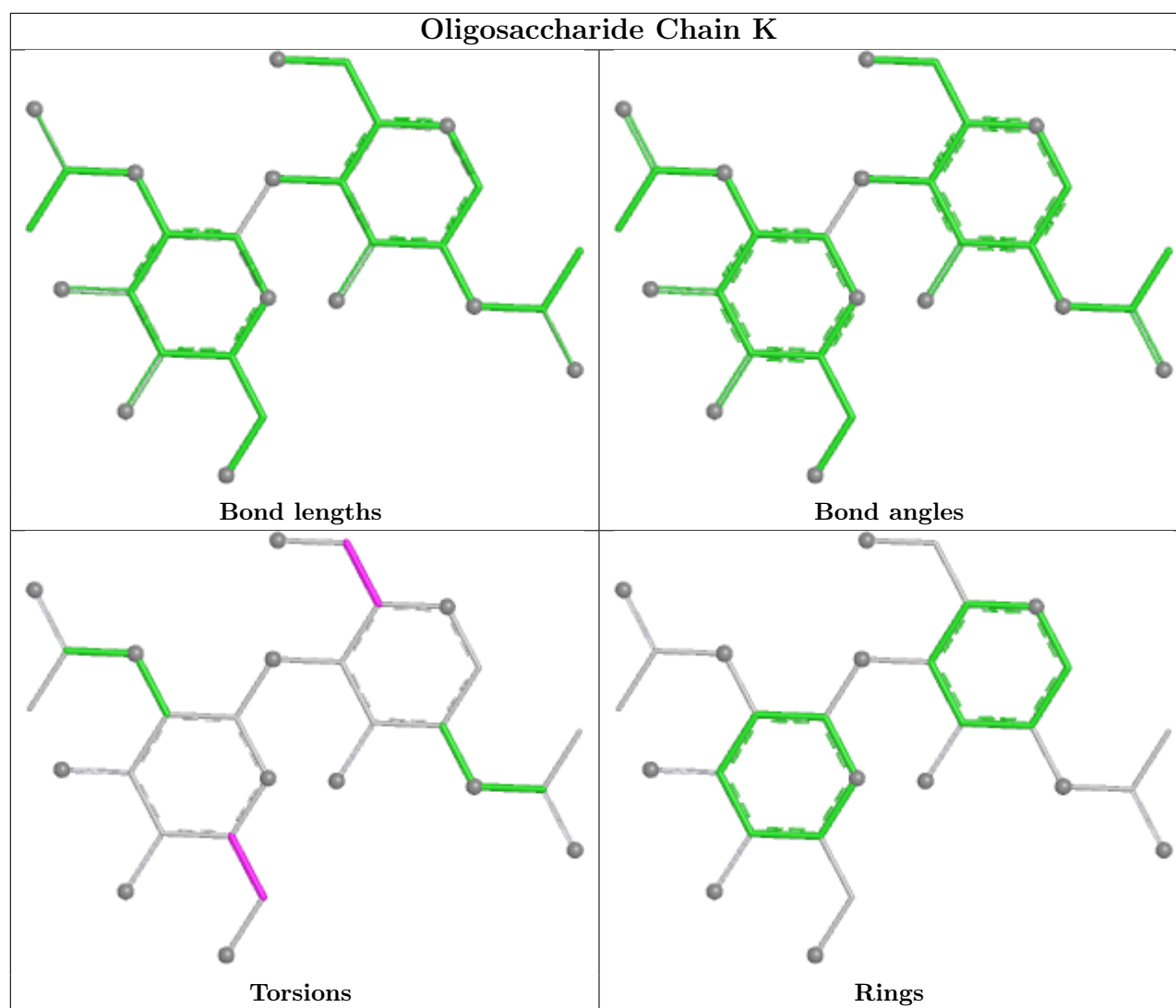
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

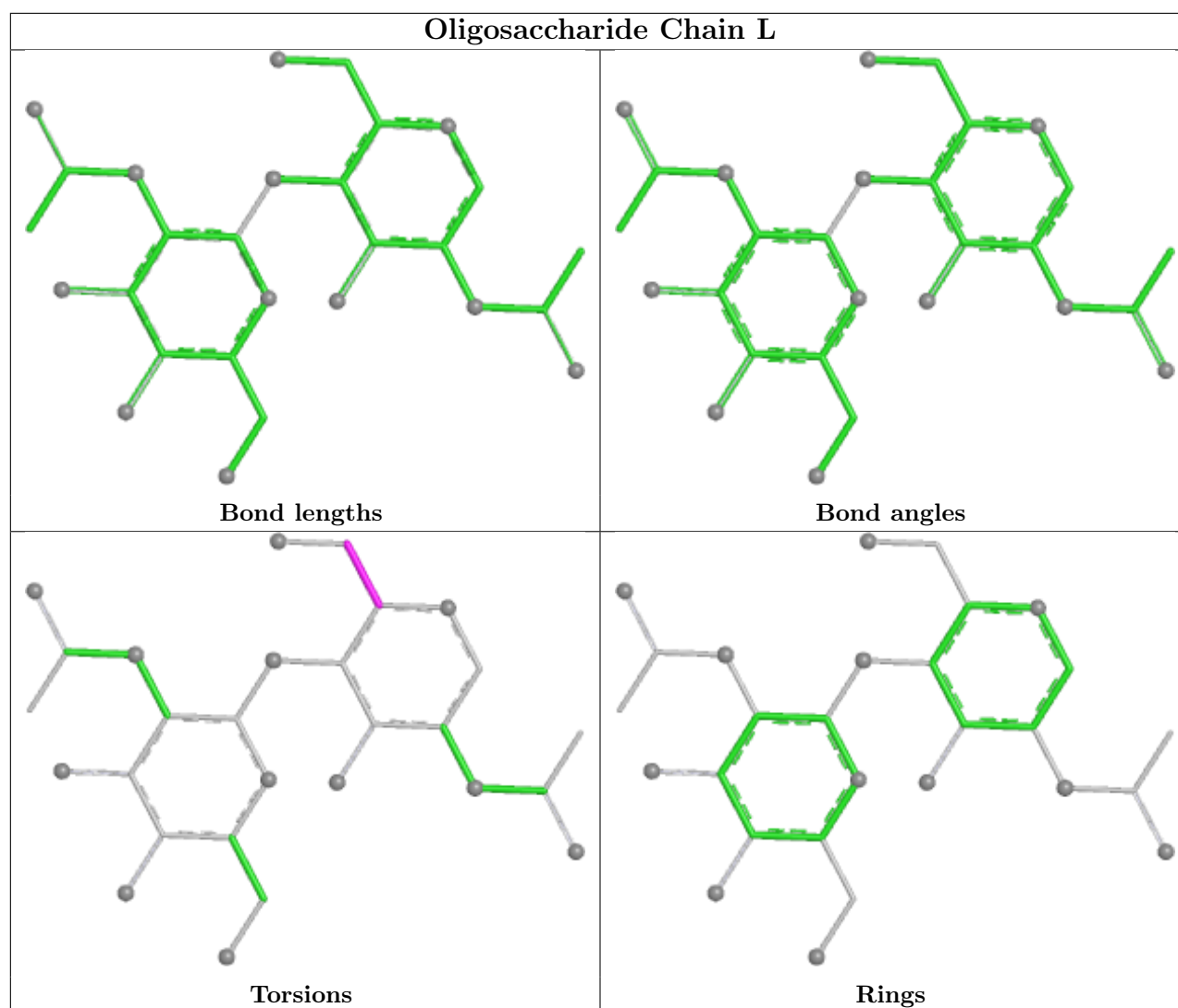












5.6 Ligand geometry [i](#)

19 ligands 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	$\# Z > 2$
4	NAG	E	1505	2	14,14,15	0.28	0	17,19,21	0.49	0
4	NAG	E	1506	2	14,14,15	0.35	0	17,19,21	0.59	0
5	EMC	F	1508	2	1,2,2	0.04	0	-		

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	F	1501	2	14,14,15	0.30	0	17,19,21	0.39	0
4	NAG	D	1507	2	14,14,15	0.30	0	17,19,21	0.54	0
4	NAG	F	1506	2	14,14,15	0.24	0	17,19,21	0.57	0
4	NAG	D	1505	2	14,14,15	0.27	0	17,19,21	0.47	0
4	NAG	F	1507	2	14,14,15	0.35	0	17,19,21	0.50	0
4	NAG	D	1501	2	14,14,15	0.29	0	17,19,21	0.50	0
4	NAG	E	1507	2	14,14,15	0.22	0	17,19,21	0.53	0
4	NAG	F	1502	2	14,14,15	0.32	0	17,19,21	0.55	0
4	NAG	F	1505	2	14,14,15	0.26	0	17,19,21	0.50	0
4	NAG	C	803	1	14,14,15	0.36	0	17,19,21	0.41	0
4	NAG	D	1506	2	14,14,15	0.33	0	17,19,21	0.51	0
4	NAG	E	1501	2	14,14,15	0.29	0	17,19,21	0.50	0
4	NAG	E	1502	2	14,14,15	0.34	0	17,19,21	0.54	0
4	NAG	D	1502	2	14,14,15	0.32	0	17,19,21	0.48	0
5	EMC	D	1508	2	1,2,2	0.04	0	-		
5	EMC	E	1508	2	1,2,2	0.06	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
4	NAG	E	1505	2	-	0/6/23/26	0/1/1/1
4	NAG	C	803	1	-	0/6/23/26	0/1/1/1
4	NAG	D	1501	2	-	0/6/23/26	0/1/1/1
4	NAG	D	1506	2	-	1/6/23/26	0/1/1/1
4	NAG	E	1501	2	-	0/6/23/26	0/1/1/1
4	NAG	E	1502	2	-	0/6/23/26	0/1/1/1
4	NAG	D	1502	2	-	2/6/23/26	0/1/1/1
4	NAG	E	1507	2	-	0/6/23/26	0/1/1/1
4	NAG	E	1506	2	-	0/6/23/26	0/1/1/1
4	NAG	F	1506	2	-	2/6/23/26	0/1/1/1
4	NAG	F	1501	2	-	0/6/23/26	0/1/1/1
4	NAG	F	1502	2	-	0/6/23/26	0/1/1/1
4	NAG	D	1505	2	-	0/6/23/26	0/1/1/1
4	NAG	D	1507	2	-	0/6/23/26	0/1/1/1
4	NAG	F	1505	2	-	0/6/23/26	0/1/1/1
4	NAG	F	1507	2	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (5) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	F	1506	NAG	C4-C5-C6-O6
4	F	1506	NAG	O5-C5-C6-O6
4	D	1502	NAG	C4-C5-C6-O6
4	D	1502	NAG	O5-C5-C6-O6
4	D	1506	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	D	1506	NAG	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	230/262 (87%)	0.22	5 (2%) 62 43	58, 91, 153, 173	0
1	B	227/262 (86%)	0.11	4 (1%) 67 49	56, 82, 134, 153	0
1	C	227/262 (86%)	0.13	4 (1%) 67 49	57, 93, 136, 160	0
2	D	407/420 (96%)	0.19	10 (2%) 58 39	59, 89, 123, 163	0
2	E	407/420 (96%)	0.14	11 (2%) 56 37	58, 89, 131, 168	0
2	F	407/420 (96%)	0.25	8 (1%) 65 46	56, 92, 131, 181	0
All	All	1905/2046 (93%)	0.18	42 (2%) 62 43	56, 90, 137, 181	0

All (42) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	268	TYR	4.2
1	C	188	GLY	4.1
2	D	73	HIS	3.7
2	F	73	HIS	3.7
2	E	300	ASN	3.6
2	D	299	GLN	3.6
1	A	188	GLY	3.3
2	E	20	PHE	3.1
2	D	418	SER	3.0
2	E	402	GLU	2.8
1	C	169	SER	2.8
2	D	402	GLU	2.8
2	E	301	ASP	2.7
1	B	188	GLY	2.7
2	F	239	ILE	2.7
2	E	73	HIS	2.7
2	E	318	PRO	2.6
2	F	318	PRO	2.5
1	C	144	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	245	ILE	2.4
2	E	429	ALA	2.4
2	F	156	ASN	2.3
2	F	62	LEU	2.3
1	A	173	GLY	2.3
2	F	240	LEU	2.3
2	E	198	SER	2.2
2	D	216	THR	2.2
2	E	418	SER	2.2
1	A	181	GLY	2.2
1	B	206	ASN	2.2
2	D	28	LEU	2.2
2	D	240	LEU	2.1
1	C	200	THR	2.1
2	E	299	GLN	2.1
1	A	178	SER	2.1
2	D	215	GLU	2.1
2	F	235	PRO	2.1
1	B	173	GLY	2.1
2	D	429	ALA	2.1
2	F	277	GLN	2.1
2	D	156	ASN	2.0
2	E	199	ASP	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.

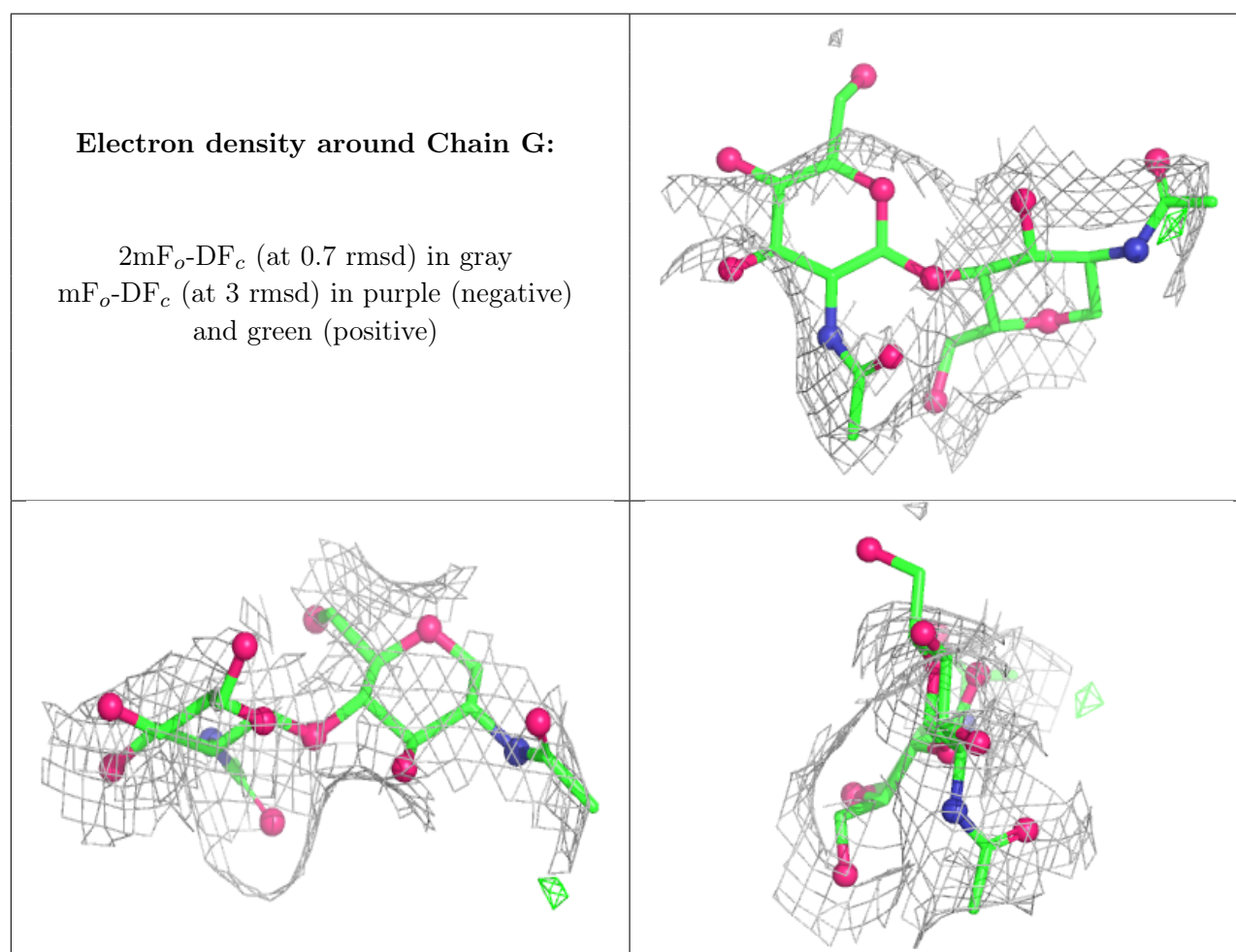
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	G	2	14/15	0.41	0.10	124,148,160,162	0
3	NAG	I	2	14/15	0.41	0.12	138,160,172,174	0
3	NAG	I	1	14/15	0.45	0.13	85,142,157,158	0
3	NAG	G	1	14/15	0.49	0.12	102,136,142,151	0
3	NAG	K	2	14/15	0.53	0.11	107,118,133,146	0

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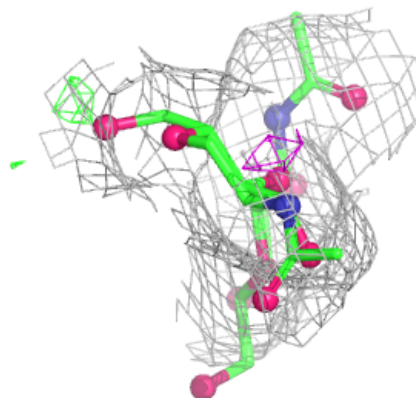
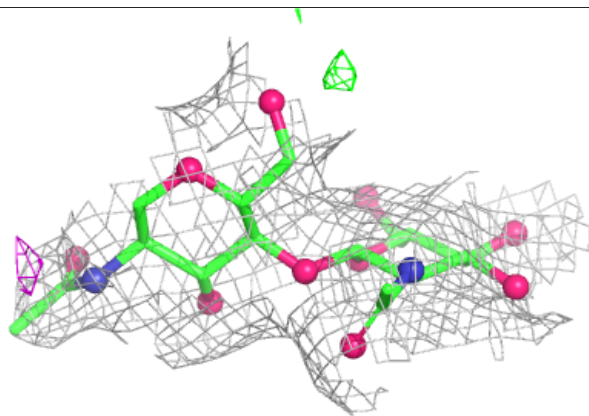
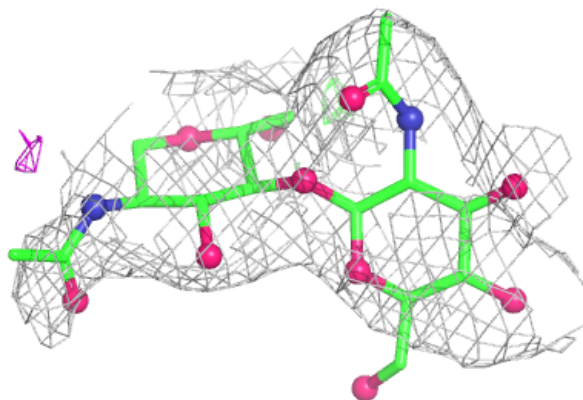
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	H	2	14/15	0.54	0.10	114,131,140,142	0
3	NAG	H	1	14/15	0.62	0.11	87,120,134,137	0
3	NAG	J	2	14/15	0.75	0.08	116,122,124,125	0
3	NAG	L	2	14/15	0.78	0.10	102,111,131,135	0
3	NAG	K	1	14/15	0.85	0.10	70,109,120,129	0
3	NAG	L	1	14/15	0.88	0.10	62,93,115,118	0
3	NAG	J	1	14/15	0.91	0.08	76,104,113,115	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.



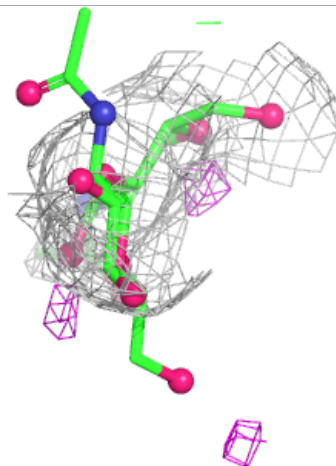
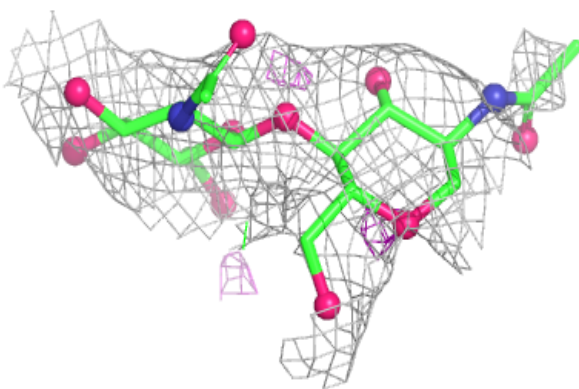
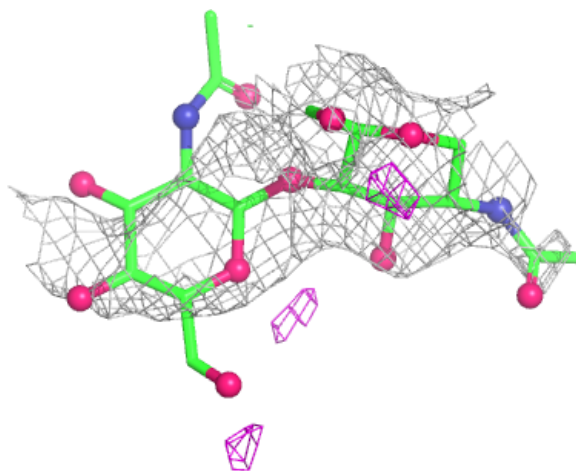
Electron density around Chain H:

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



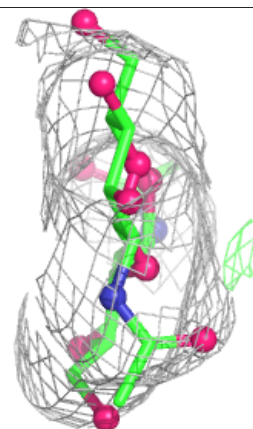
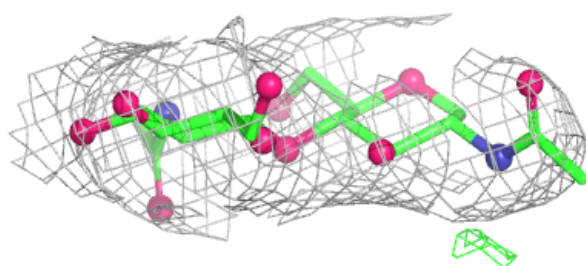
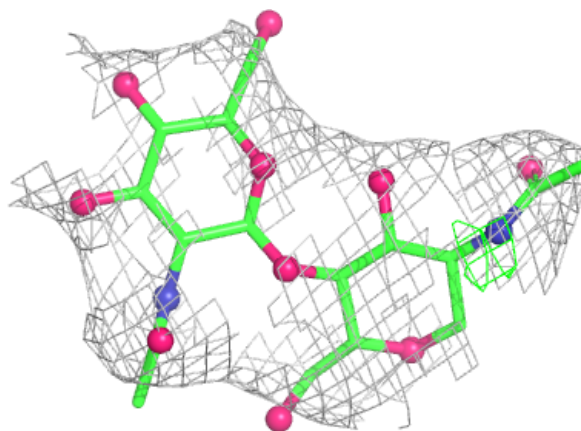
Electron density around Chain I:

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



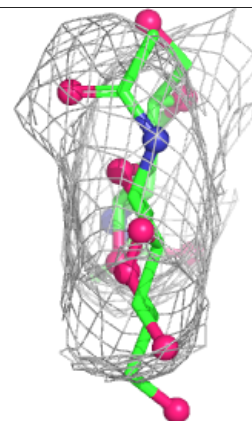
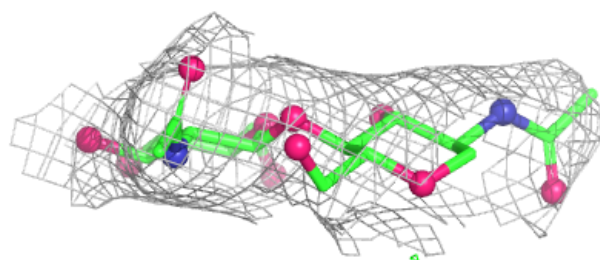
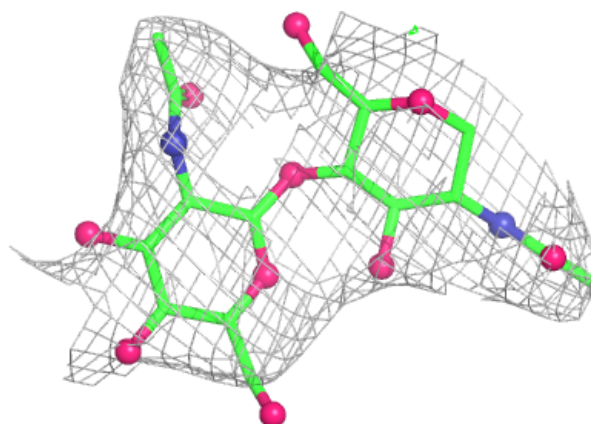
Electron density around Chain J:

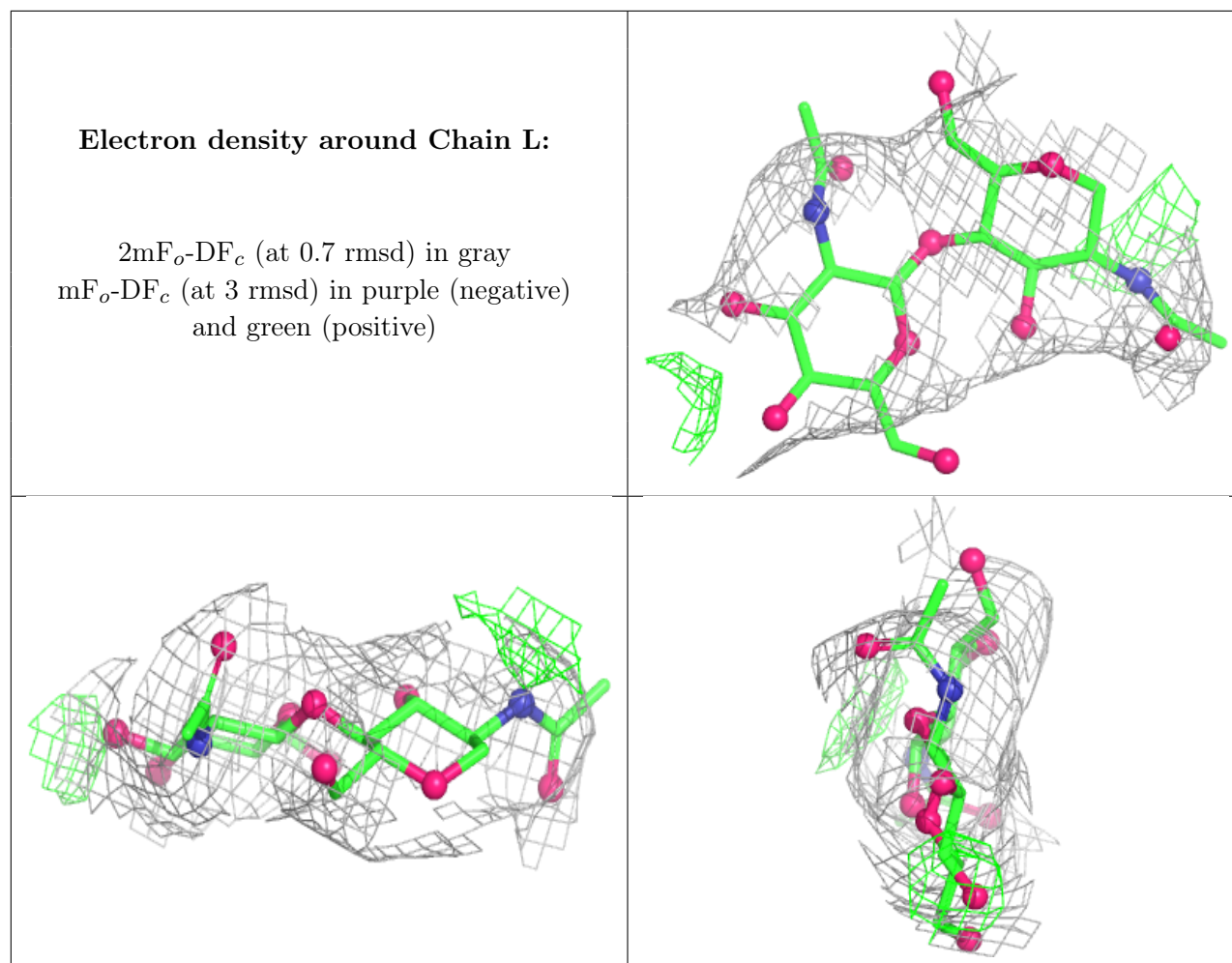
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain K:

$2mF_o - DF_c$ (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
4	NAG	F	1506	14/15	0.37	0.14	98,137,147,148	0
4	NAG	F	1507	14/15	0.48	0.15	99,130,145,152	0
4	NAG	E	1506	14/15	0.52	0.13	100,127,140,140	0
5	EMC	D	1508	3/3	0.55	0.25	115,115,133,307	0
4	NAG	C	803	14/15	0.56	0.14	94,155,166,171	0
5	EMC	E	1508	3/3	0.61	0.24	108,108,117,281	3
4	NAG	D	1507	14/15	0.62	0.15	122,139,155,168	0
4	NAG	D	1506	14/15	0.64	0.13	113,135,139,141	0
4	NAG	F	1501	14/15	0.68	0.13	80,103,114,127	0
4	NAG	D	1501	14/15	0.69	0.12	99,109,130,138	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	D	1505	14/15	0.73	0.11	71,96,112,118	0
4	NAG	E	1502	14/15	0.74	0.13	101,117,135,135	0
4	NAG	E	1507	14/15	0.75	0.12	118,135,147,157	0
4	NAG	E	1505	14/15	0.75	0.10	95,104,119,127	0
5	EMC	F	1508	3/3	0.77	0.19	166,166,187,321	0
4	NAG	F	1502	14/15	0.78	0.11	83,104,111,115	0
4	NAG	F	1505	14/15	0.79	0.11	74,83,95,107	0
4	NAG	D	1502	14/15	0.79	0.12	76,117,128,128	0
4	NAG	E	1501	14/15	0.81	0.11	87,106,123,131	0

6.5 Other polymers [i](#)

There are no such residues in this entry.