WORLDWIDE wwPDB PROTEIN DATA BANK

wwPDB X-ray Structure Validation Summary Report ⓘ

Nov 2, 2024 – 01:23 PM EDT

PDB ID	:	6NNJ
Title	:	Crystal Structure of HIV-1 BG505 SOSIP.664 Prefusion Env Trimer Bound to CH31 scFv in Complex with Crystallization Chaperones 3H109L Fab and 35O22 scFv at 3.1 Angstrom
Authors	:	Lai, Y.-T.; Kwong, P.D.
Deposited on	:	2019-01-15
Resolution	:	2.60 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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:

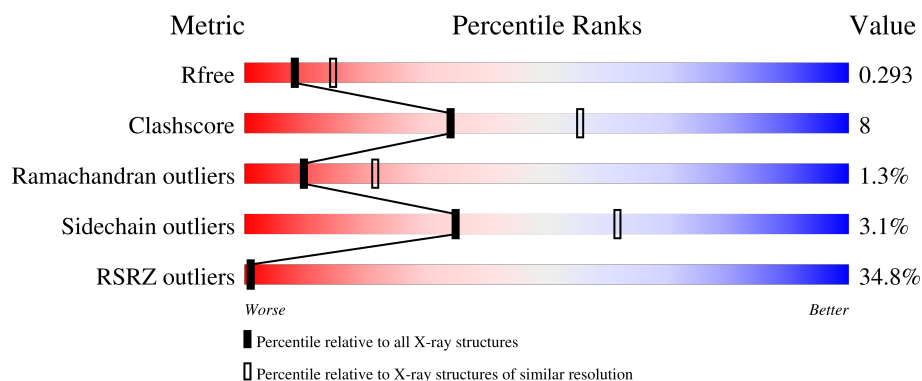|		
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.003 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.39

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 2.60 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	Percentile Ranks	Value
Rfree		0.293
Clashscore		8
Ramachandran outliers		1.3%
Sidechain outliers		3.1%
RSRZ outliers		34.8%

Worse — *Better*

▮ Percentile relative to all X-ray structures
▯ Percentile relative to X-ray structures of similar resolution

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	164625	3775 (2.60-2.60)
Clashscore	180529	4181 (2.60-2.60)
Ramachandran outliers	177936	4129 (2.60-2.60)
Sidechain outliers	177891	4129 (2.60-2.60)
RSRZ outliers	164620	3775 (2.60-2.60)

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 <=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	B	153	9% 69% 14% • 17%
2	D	153	54% 60% 22% • 16%
3	E	130	49% 69% 15% 15%
4	G	481	10% 67% 21% • 11%
5	H	244	33% 73% 18% • 7%

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Mol	Chain	Length	Quality of chain
6	L	217	13% — 74% — 21%
7	U	137	72% — 85% — 8% · 6%
8	V	117	77% — 76% — 10% — 14%
9	A	6	33% — 17% — 50%
10	C	2	100%
10	J	2	50% — 50%
10	K	2	100%
10	M	2	100%
10	O	2	50% — 50%
11	F	5	60% — 40%
12	I	4	50% — 50%
13	N	10	20% — 60% — 20%

2 Entry composition ⓘ

There are 14 unique types of molecules in this entry. The entry contains 11874 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 Envelope glycoprotein gp41.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	127	Total	C	N	O	S	0	0	0
			1009	641	173	189	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	engineered mutation	UNP Q2N0S6
B	605	CYS	THR	engineered mutation	UNP Q2N0S6

- Molecule 2 is a protein called 35O22 scFv heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	128	Total	C	N	O	S	0	0	0
			994	628	169	192	5			

- Molecule 3 is a protein called 35O22 scFv light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	E	110	Total	C	N	O	S	0	0	0
			836	525	138	167	6			

- Molecule 4 is a protein called Envelope glycoprotein gp120.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	G	430	Total	C	N	O	S	0	0	0
			3397	2140	602	628	27			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	145	ALA	ASN	engineered mutation	UNP Q2N0S6
G	332	ASN	THR	engineered mutation	UNP Q2N0S6

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Chain	Residue	Modelled	Actual	Comment	Reference
G	501	CYS	ALA	engineered mutation	UNP Q2N0S6
G	509	ARG	-	expression tag	UNP Q2N0S6
G	510	ARG	-	expression tag	UNP Q2N0S6
G	511	ARG	-	expression tag	UNP Q2N0S6
G	512	ARG	-	expression tag	UNP Q2N0S6
G	513	ARG	-	expression tag	UNP Q2N0S6

- Molecule 5 is a protein called 3H109L Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
5	H	226	1715	1093	278	338	6	0	0	0

- Molecule 6 is a protein called 3H109L Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
6	L	210	1598	1006	275	310	7	0	0	0

- Molecule 7 is a protein called CH31 scFv heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
7	U	129	1027	652	184	187	4	0	0	0

- Molecule 8 is a protein called CH31 scFv light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
8	V	101	761	472	130	156	3	0	0	0

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.
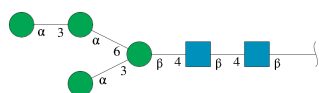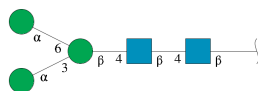
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	A	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 10 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
10	C	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	J	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	K	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	M	2	Total	C	N	O	0	0	0
			28	16	2	10			
10	O	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 11 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.
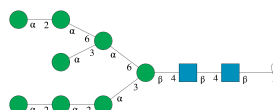

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
11	F	5	Total	C	N	O	0	0	0
			61	34	2	25			

- Molecule 12 is an oligosaccharide called alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.
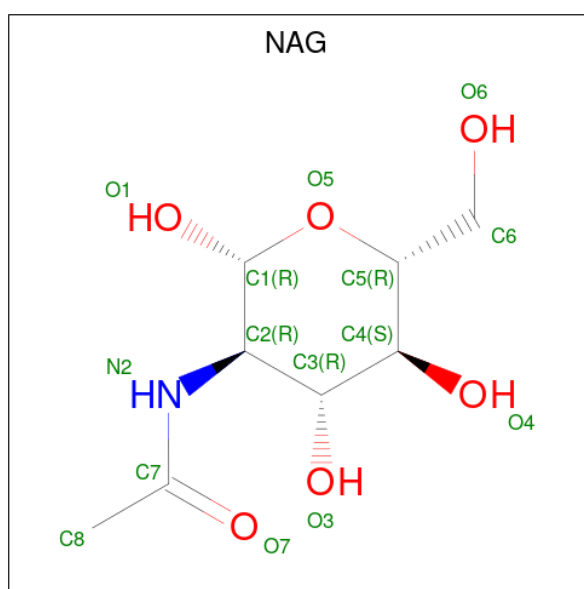
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
12	I	4	Total 50	C 28	N 2	O 20	0	0	0

- Molecule 13 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
13	N	10	Total 116	C 64	N 2	O 50	0	0	0

- Molecule 14 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).

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

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| Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
14	G	1	Total 14	C 8	N 1	O 5	0	0
14	G	1	Total 14	C 8	N 1	O 5	0	0
14	G	1	Total 14	C 8	N 1	O 5	0	0
14	G	1	Total 14	C 8	N 1	O 5	0	0

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.
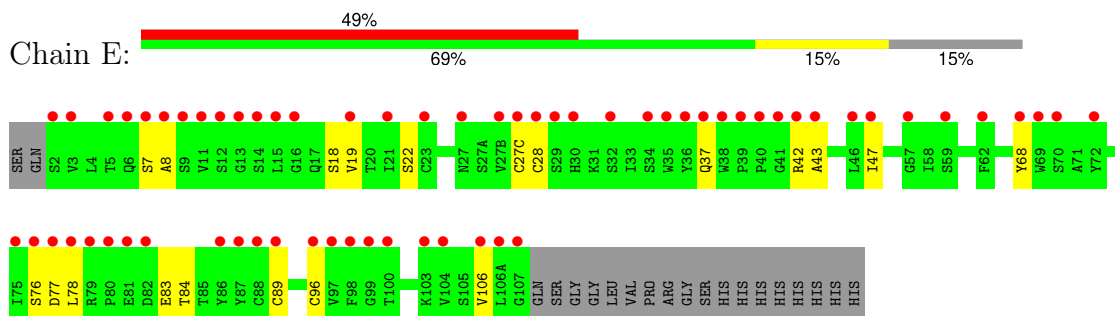
- Molecule 1: Envelope glycoprotein gp41
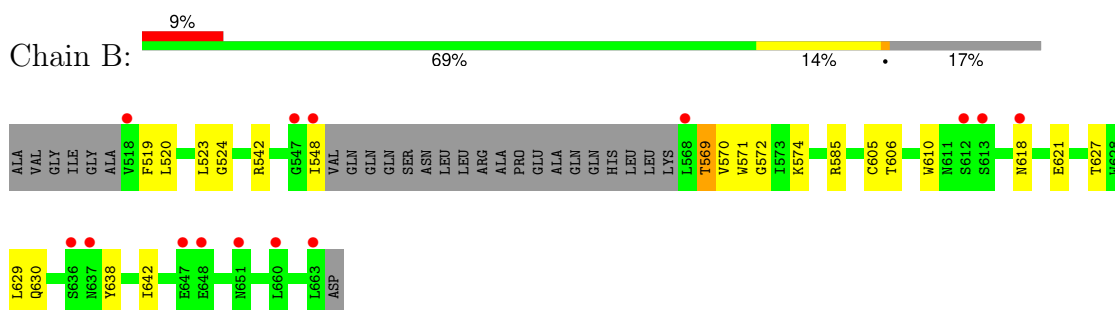
Chain B:

- Molecule 2: 35O22 scFv heavy chain
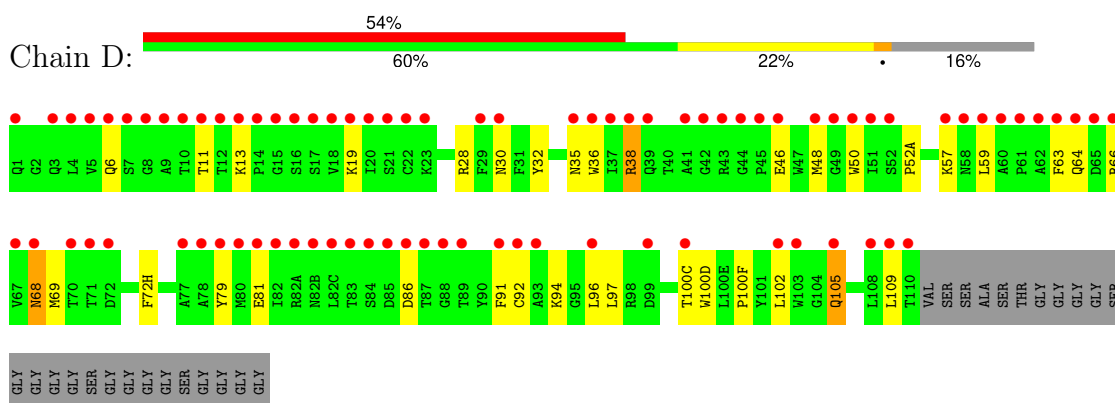
Chain D:

- Molecule 3: 35O22 scFv light chain

Chain E:

- Molecule 4: Envelope glycoprotein gp120

Chain G:

10%

67% 21% 11%
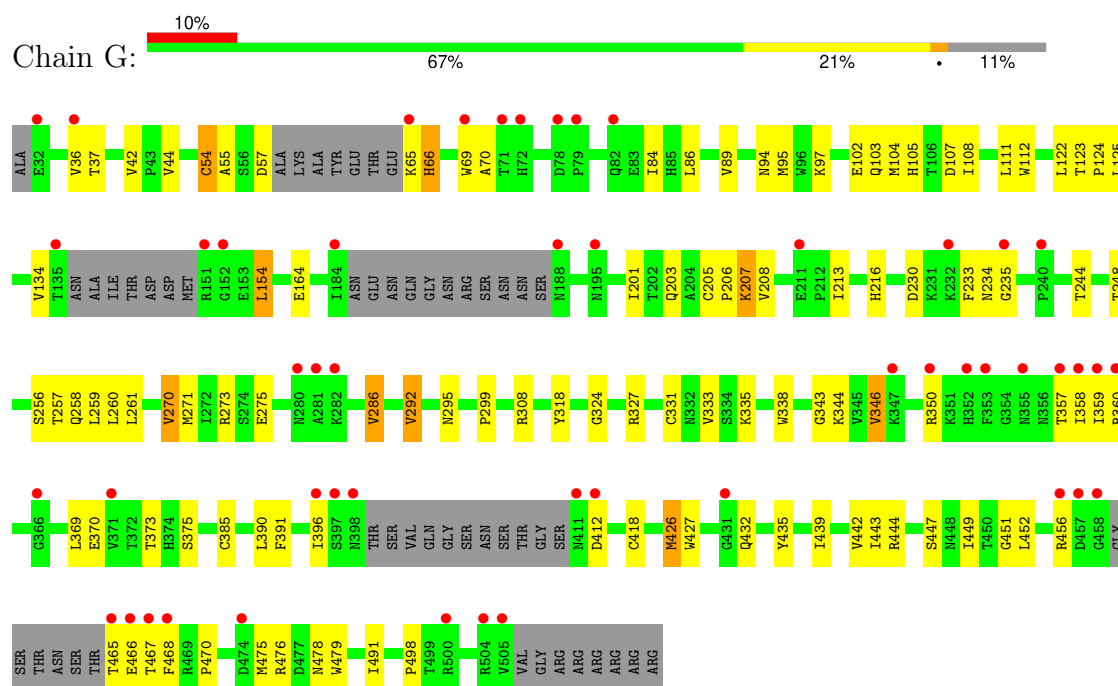
• Molecule 5: 3H109L Fab heavy chain

Chain H:

33%

73% 18% 7%
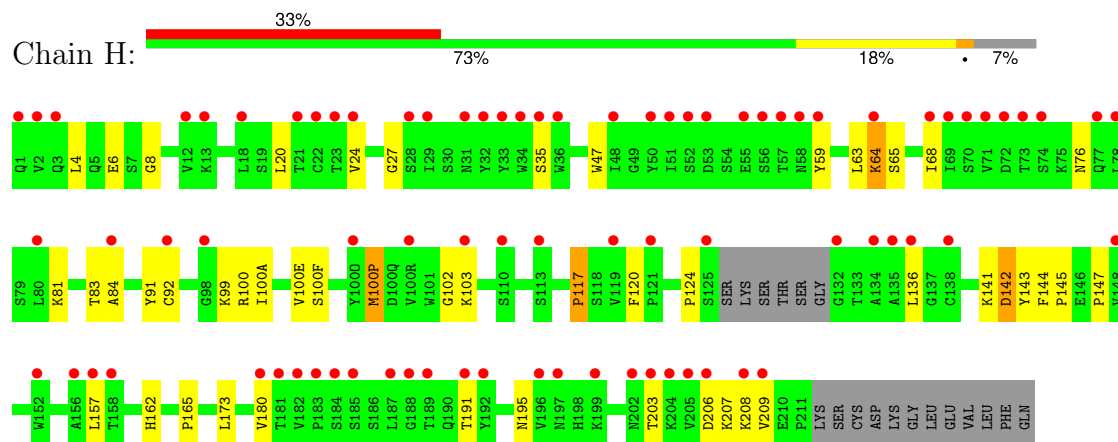
• Molecule 6: 3H109L Fab light chain

Chain L:

13%

74% 21%
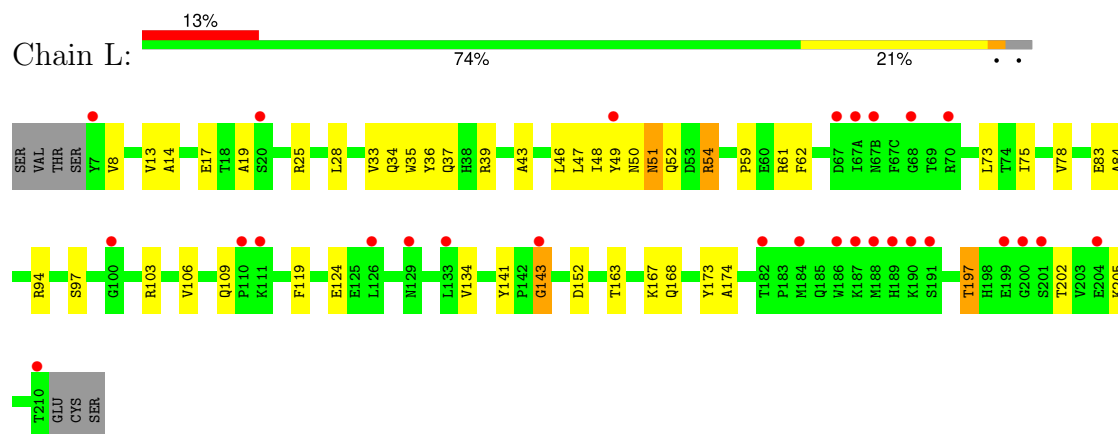
• Molecule 7: CH31 scFv heavy chain

Chain U:

72%

85%

8% · 6%

GLN V2 Q3 L4 V5 Q6 S7 G8 A9 A10 V11 R12 K13 P14 G15 A16 S17 V18 T19 V20 K23 F24 A25 E26 D27 D28 D29 Y30 S31 P31A Y31B W31C V31D N31E P31F A31G F36 L37 R38 Q39 A40 P41 G42 Q43 Q44 L45 E46 W47 L48 A49 W50 M51 N58 Y59 A60 W61 Y62

L63 N64 G65 R66 V67 T68 A69 T70 R71 D72 R73 T76 T77 A78 F79 L80 E81 V82 K82A S82B L82C R83 S84 D85 D86 T87 A88 V89 Y90 Y91 C92 A93 R94 A95 Q96 K97 R98 G99 R100 S100A E100B W100C A100D Y100E W103 G104 Q105 G106 T107 P108 V109 V110 V111 S112 SER GLY LEU

VAL PRO ARG

• Molecule 8: CH31 scFv light chain

Chain V:

77%

76%

10%   14%

D1 I2 Q3 M4 T5 Q6 S7 P8 S9 S10 L11 S12 A13 S14 L15 G16 D17 R18 V19 T20 I21 T22 C23 Q24 A25 S26 R27 G28 I29 G30 K31 D32 L33 N34 W35 Y36 Q37 Q38 K39 A40 G41 K42 A43 P44 K45 L46 L47 V48 S49 D50 A51 L54 E55 G56 G57 V58 P59 S60 R61

F62 S63 G64 S65 G66 F67 H68 Q69 N70 F71 S72 L73 T74 I75 S76 S77 L78 Q79 A80 E81 D82 V83 A84 T85 Y86 F87 C88 Q89 T97 F98 G99 Q100 G101 T102 K103 V104 D105 ILE GLY GLY GLY GLY SER GLY GLY GLY SER GLY GLY GLY SER

• Molecule 9: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyran ose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetam ido-2-deoxy-beta-D-glucopyranose

Chain A:

33%   17%   50%

NAG1 NAG2 BMA3 MAN4 MAN5 MAN6

• Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glu copyranose

Chain C:

100%

NAG1 NAG2

• Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glu copyranose

Chain J:

50%   50%

NAG1 NAG2
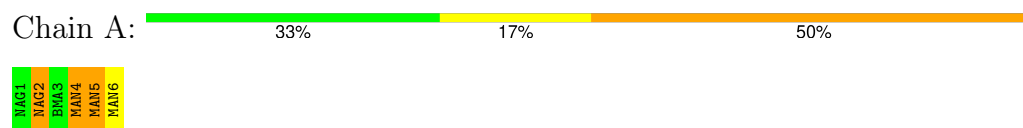
• Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glu copyranose

Chain K:

100%

NAG1
NAG2

• Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glu copyranose

Chain M:

100%

NAG1
NAG2

• Molecule 10: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glu copyranose

Chain O:

50%50%

NAG1
NAG2

• Molecule 11: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyran ose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyra nose

Chain F:

60%40%

NAG1
NAG2
BMA3
MAN4
MAN5

• Molecule 12: alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy- beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:

50%50%

NAG1
NAG2
BMA3
MAN4

• Molecule 13: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyra nose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranos e-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-g lucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose
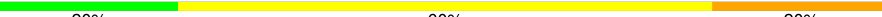
Chain N:

20%60%20%

NAG1
NAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α, β, γ	132.15Å 132.15Å 315.95Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	43.26 – 2.60 43.26 – 2.60	Depositor EDS
% Data completeness (in resolution range)	41.7 (43.26-2.60) 41.7 (43.26-2.60)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.54 (at 2.58Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.248 , 0.293 0.248 , 0.293	Depositor DCC
R_{free} test set	93476 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å^2)	31.5	Xtriage
Anisotropy	0.107	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 34.6	EDS
L-test for twinning[2]	$< \lvert L \rvert > = 0.50, < L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.048 for h,-h-k,-l	Xtriage
F_o,F_c correlation	0.82	EDS
Total number of atoms	11874	wwPDB-VP
Average B, all atoms (Å^2)	55.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.01% of the height of the origin peak. No significant pseudotranslation is detected.*

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[1]Intensities estimated from amplitudes.
[2]Theoretical values of $< \lvert L \rvert >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: MAN, BMA, NAG

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	B	0.25	0/1027	0.39	0/1393
2	D	0.25	0/1021	0.48	0/1390
3	E	0.25	0/860	0.42	0/1175
4	G	0.27	0/3467	0.47	0/4702
5	H	0.26	0/1758	0.46	0/2397
6	L	0.25	0/1641	0.45	0/2239
7	U	0.24	0/1059	0.43	0/1447
8	V	0.25	0/774	0.46	0/1045
All	All	0.25	0/11607	0.45	0/15788

There are no bond length outliers.

There are no bond angle outliers.

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	B	1009	0	994	16	0
2	D	994	0	953	20	0
3	E	836	0	785	8	0
4	G	3397	0	3349	73	0
5	H	1715	0	1685	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
6	L	1598	0	1548	27	0
7	U	1027	0	973	4	0
8	V	761	0	735	7	0
9	A	72	0	61	4	0
10	C	28	0	25	0	0
10	J	28	0	25	2	0
10	K	28	0	25	0	0
10	M	28	0	25	0	0
10	O	28	0	25	0	0
11	F	61	0	52	0	0
12	I	50	0	43	1	0
13	N	116	0	97	2	0
14	B	28	0	26	0	0
14	G	70	0	65	0	0
All	All	11874	0	11491	176	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 8.

The worst 5 of 176 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)
5:H:59:TYR:HB2	5:H:64:LYS:HD2	1.56	0.87
4:G:230:ASP:HB3	4:G:233:PHE:HB2	1.62	0.81
4:G:475:MET:SD	4:G:478:ASN:ND2	2.55	0.78
4:G:358:ILE:HG22	4:G:396:ILE:HA	1.68	0.73
4:G:259:LEU:HD23	4:G:449:ILE:HG21	1.70	0.72

There are no symmetry-related clashes.

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	B	123/153 (80%)	116 (94%)	7 (6%)	0	100	100
2	D	126/153 (82%)	115 (91%)	10 (8%)	1 (1%)	16	34
3	E	108/130 (83%)	89 (82%)	17 (16%)	2 (2%)	6	13
4	G	418/481 (87%)	383 (92%)	31 (7%)	4 (1%)	13	29
5	H	222/244 (91%)	200 (90%)	18 (8%)	4 (2%)	7	14
6	L	208/217 (96%)	194 (93%)	11 (5%)	3 (1%)	9	19
7	U	127/137 (93%)	120 (94%)	4 (3%)	3 (2%)	5	9
8	V	99/117 (85%)	91 (92%)	7 (7%)	1 (1%)	13	29
All	All	1431/1632 (88%)	1308 (91%)	105 (7%)	18 (1%)	10	21

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
6	L	51	ASN
7	U	31	SER
4	G	427	TRP
6	L	143	GLY
4	G	70	ALA

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	B	109/129 (84%)	107 (98%)	2 (2%)	54	77
2	D	107/115 (93%)	102 (95%)	5 (5%)	22	45
3	E	96/113 (85%)	92 (96%)	4 (4%)	25	50
4	G	385/427 (90%)	372 (97%)	13 (3%)	32	58
5	H	196/212 (92%)	190 (97%)	6 (3%)	35	62
6	L	174/181 (96%)	169 (97%)	5 (3%)	37	64
7	U	105/111 (95%)	101 (96%)	4 (4%)	28	54
8	V	84/88 (96%)	84 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
All	All	1256/1376 (91%)	1217 (97%)	39 (3%)	35 62

5 of 39 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
5	H	203	THR
7	U	28	ASP
5	H	207	LYS
6	L	97	SER
7	U	30	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
8	V	24	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

35 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	$\#|Z| > 2$
9	NAG	A	1	9,4	14,14,15	0.39	0	17,19,21	0.49	0
9	NAG	A	2	9	14,14,15	0.42	0	17,19,21	1.40	2 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	#|Z| > 2	Counts	RMSZ	#|Z| > 2
9	BMA	A	3	9	11,11,12	0.71	0	15,15,17	0.85	0
9	MAN	A	4	9	11,11,12	0.69	0	15,15,17	1.05	2 (13%)
9	MAN	A	5	9	11,11,12	1.66	1 (9%)	15,15,17	1.54	2 (13%)
9	MAN	A	6	9	11,11,12	0.69	0	15,15,17	0.97	2 (13%)
10	NAG	C	1	10,4	14,14,15	0.23	0	17,19,21	0.46	0
10	NAG	C	2	10	14,14,15	0.29	0	17,19,21	0.49	0
11	NAG	F	1	11,4	14,14,15	0.31	0	17,19,21	0.48	0
11	NAG	F	2	11	14,14,15	0.16	0	17,19,21	0.43	0
11	BMA	F	3	11	11,11,12	0.70	0	15,15,17	0.72	0
11	MAN	F	4	11	11,11,12	0.72	0	15,15,17	0.99	2 (13%)
11	MAN	F	5	11	11,11,12	0.75	0	15,15,17	0.98	2 (13%)
12	NAG	I	1	12,4	14,14,15	0.35	0	17,19,21	0.49	0
12	NAG	I	2	12	14,14,15	0.24	0	17,19,21	0.60	0
12	BMA	I	3	12	11,11,12	0.71	0	15,15,17	0.68	0
12	MAN	I	4	12	11,11,12	0.75	1 (9%)	15,15,17	1.07	2 (13%)
10	NAG	J	1	10,4	14,14,15	0.34	0	17,19,21	0.39	0
10	NAG	J	2	10	14,14,15	0.37	0	17,19,21	0.46	0
10	NAG	K	1	10,4	14,14,15	0.36	0	17,19,21	0.40	0
10	NAG	K	2	10	14,14,15	0.20	0	17,19,21	0.47	0
10	NAG	M	1	10,4	14,14,15	0.31	0	17,19,21	0.48	0
10	NAG	M	2	10	14,14,15	0.28	0	17,19,21	0.53	0
13	NAG	N	1	13,4	14,14,15	0.29	0	17,19,21	1.59	2 (11%)
13	MAN	N	10	13	11,11,12	1.07	2 (18%)	15,15,17	1.43	3 (20%)
13	NAG	N	2	13	14,14,15	0.52	0	17,19,21	0.48	0
13	BMA	N	3	13	11,11,12	0.99	0	15,15,17	1.00	0
13	MAN	N	4	13	11,11,12	0.80	1 (9%)	15,15,17	1.30	1 (6%)
13	MAN	N	5	13	11,11,12	0.75	0	15,15,17	1.11	1 (6%)
13	MAN	N	6	13	11,11,12	0.88	0	15,15,17	0.88	1 (6%)
13	MAN	N	7	13	11,11,12	0.72	0	15,15,17	0.96	2 (13%)
13	MAN	N	8	13	11,11,12	0.69	0	15,15,17	0.98	1 (6%)
13	MAN	N	9	13	11,11,12	0.87	1 (9%)	15,15,17	1.23	2 (13%)
10	NAG	O	1	10,4	14,14,15	0.88	1 (7%)	17,19,21	1.50	1 (5%)
10	NAG	O	2	10	14,14,15	0.27	0	17,19,21	0.39	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
9	NAG	A	1	9,4	-	0/6/23/26	0/1/1/1
9	NAG	A	2	9	-	6/6/23/26	0/1/1/1
9	BMA	A	3	9	-	2/2/19/22	0/1/1/1
9	MAN	A	4	9	-	0/2/19/22	0/1/1/1
9	MAN	A	5	9	-	2/2/19/22	0/1/1/1
9	MAN	A	6	9	-	0/2/19/22	0/1/1/1
10	NAG	C	1	10,4	-	0/6/23/26	0/1/1/1
10	NAG	C	2	10	-	2/6/23/26	0/1/1/1
11	NAG	F	1	11,4	-	0/6/23/26	0/1/1/1
11	NAG	F	2	11	-	0/6/23/26	0/1/1/1
11	BMA	F	3	11	-	2/2/19/22	0/1/1/1
11	MAN	F	4	11	-	2/2/19/22	0/1/1/1
11	MAN	F	5	11	-	2/2/19/22	0/1/1/1
12	NAG	I	1	12,4	-	2/6/23/26	0/1/1/1
12	NAG	I	2	12	-	2/6/23/26	0/1/1/1
12	BMA	I	3	12	-	2/2/19/22	0/1/1/1
12	MAN	I	4	12	-	2/2/19/22	0/1/1/1
10	NAG	J	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	J	2	10	-	2/6/23/26	0/1/1/1
10	NAG	K	1	10,4	-	2/6/23/26	0/1/1/1
10	NAG	K	2	10	-	0/6/23/26	0/1/1/1
10	NAG	M	1	10,4	-	1/6/23/26	0/1/1/1
10	NAG	M	2	10	-	0/6/23/26	0/1/1/1
13	NAG	N	1	13,4	-	6/6/23/26	0/1/1/1
13	MAN	N	10	13	-	0/2/19/22	0/1/1/1
13	NAG	N	2	13	-	2/6/23/26	0/1/1/1
13	BMA	N	3	13	-	0/2/19/22	0/1/1/1
13	MAN	N	4	13	-	2/2/19/22	0/1/1/1
13	MAN	N	5	13	-	0/2/19/22	0/1/1/1
13	MAN	N	6	13	-	0/2/19/22	0/1/1/1
13	MAN	N	7	13	-	0/2/19/22	0/1/1/1
13	MAN	N	8	13	-	0/2/19/22	0/1/1/1
13	MAN	N	9	13	-	0/2/19/22	0/1/1/1
10	NAG	O	1	10,4	-	4/6/23/26	0/1/1/1
10	NAG	O	2	10	-	2/6/23/26	0/1/1/1

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	A	5	MAN	C1-C2	4.83	1.63	1.52
10	O	1	NAG	O5-C1	3.17	1.49	1.43
13	N	9	MAN	C1-C2	2.38	1.57	1.52
13	N	10	MAN	C1-C2	2.31	1.57	1.52
12	I	4	MAN	C1-C2	2.17	1.57	1.52

The worst 5 of 28 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
10	O	1	NAG	C1-O5-C5	5.80	119.96	112.19
13	N	1	NAG	C2-N2-C7	5.01	129.61	122.90
9	A	2	NAG	C2-N2-C7	4.66	129.14	122.90
13	N	4	MAN	C1-O5-C5	3.88	117.39	112.19
9	A	5	MAN	C1-C2-C3	3.86	115.27	109.64

There are no chirality outliers.
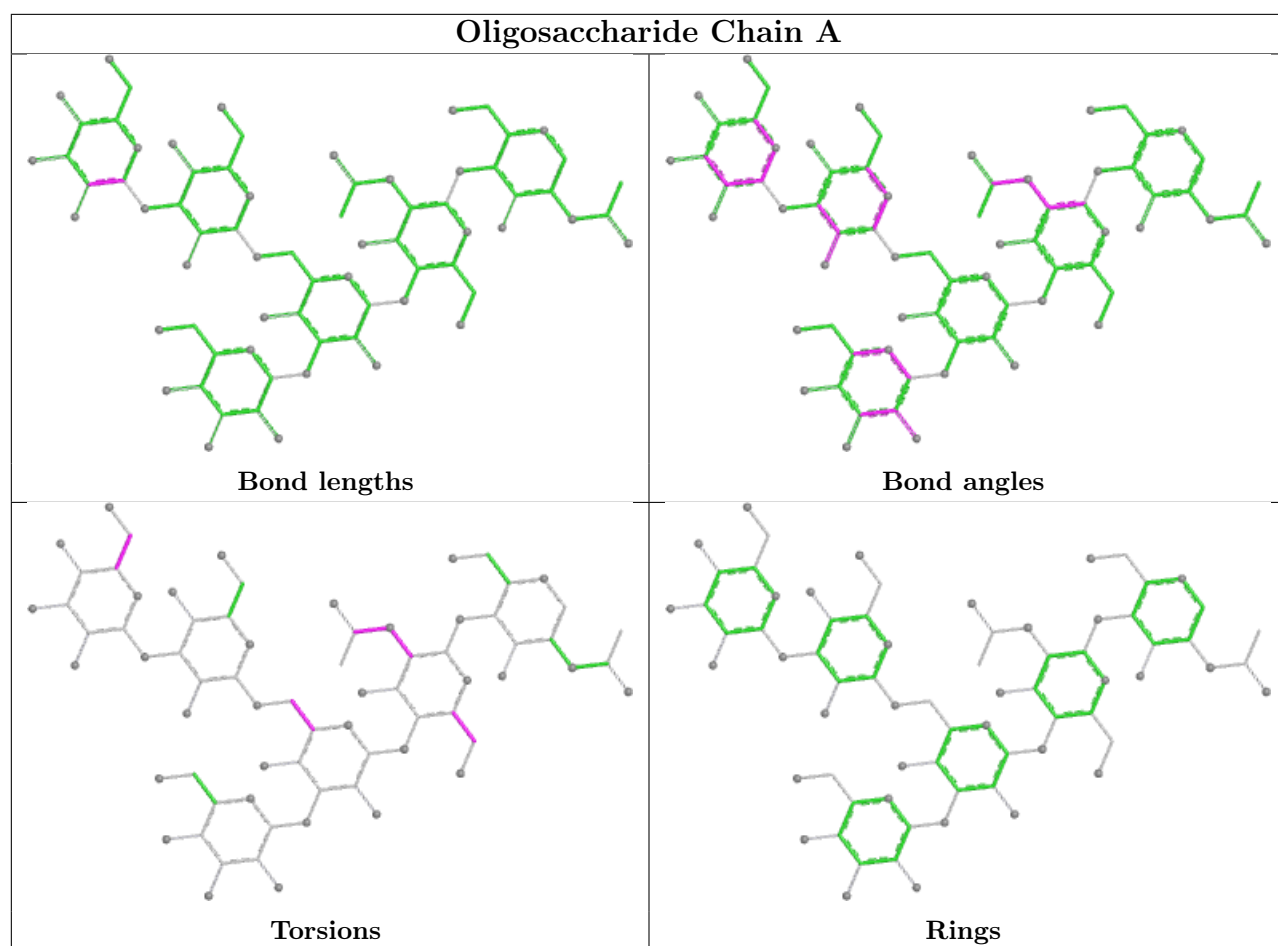
5 of 49 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
10	O	2	NAG	O5-C5-C6-O6
12	I	2	NAG	O5-C5-C6-O6
9	A	3	BMA	O5-C5-C6-O6
10	K	1	NAG	O5-C5-C6-O6
12	I	1	NAG	O5-C5-C6-O6

There are no ring outliers.

7 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	A	2	NAG	1	0
9	A	4	MAN	3	0
10	J	2	NAG	2	0
9	A	5	MAN	1	0
13	N	4	MAN	1	0
12	I	1	NAG	1	0
13	N	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.
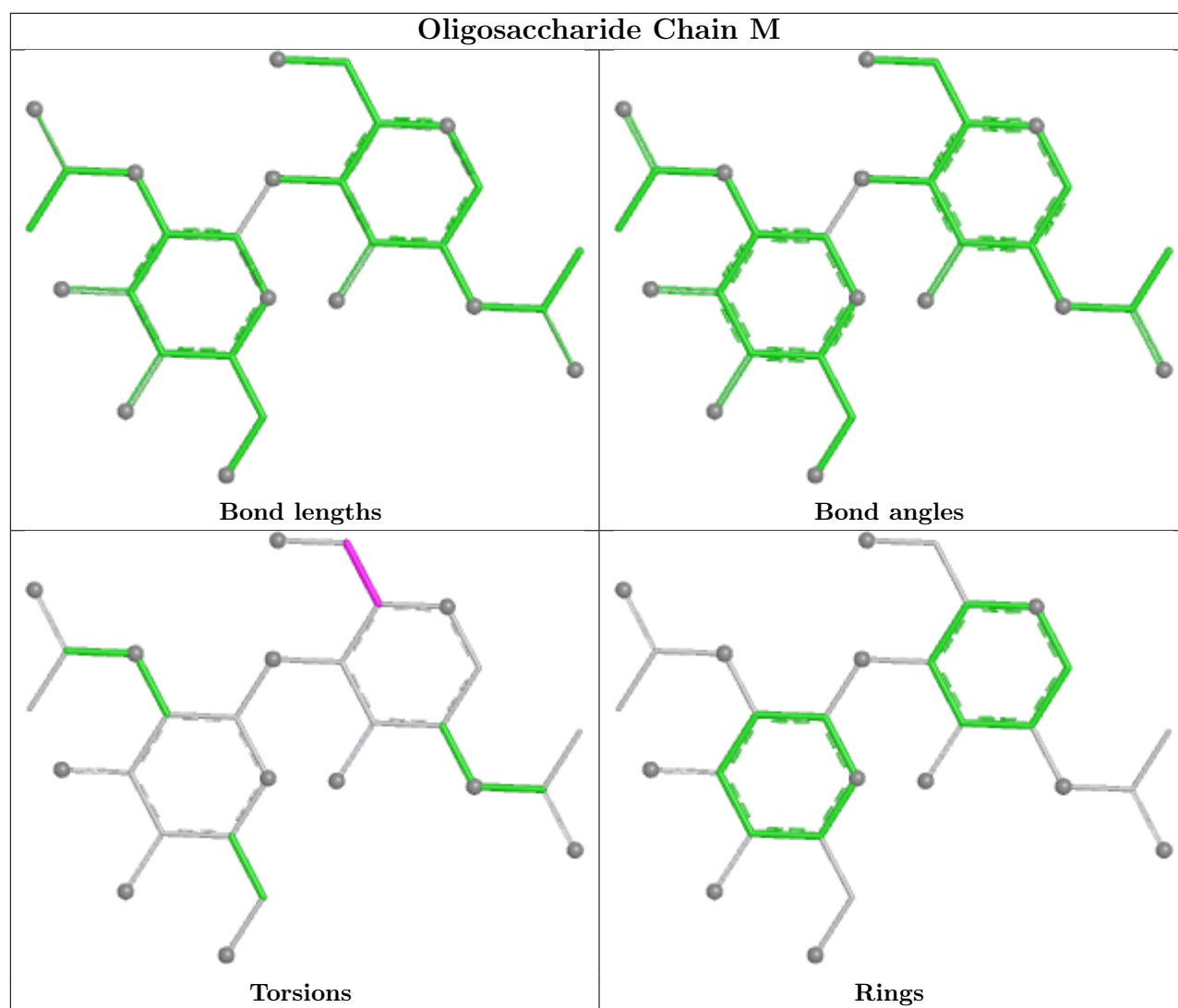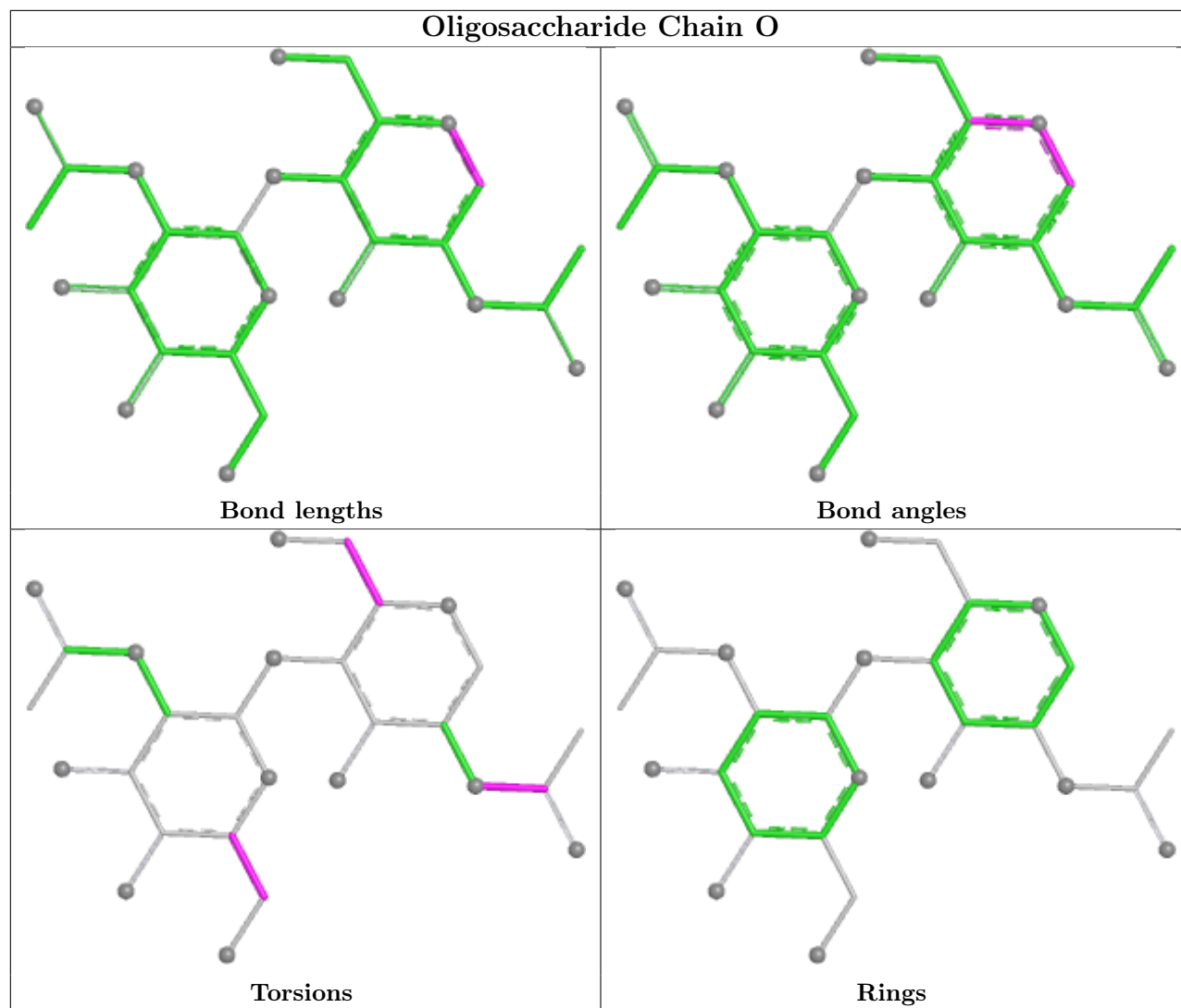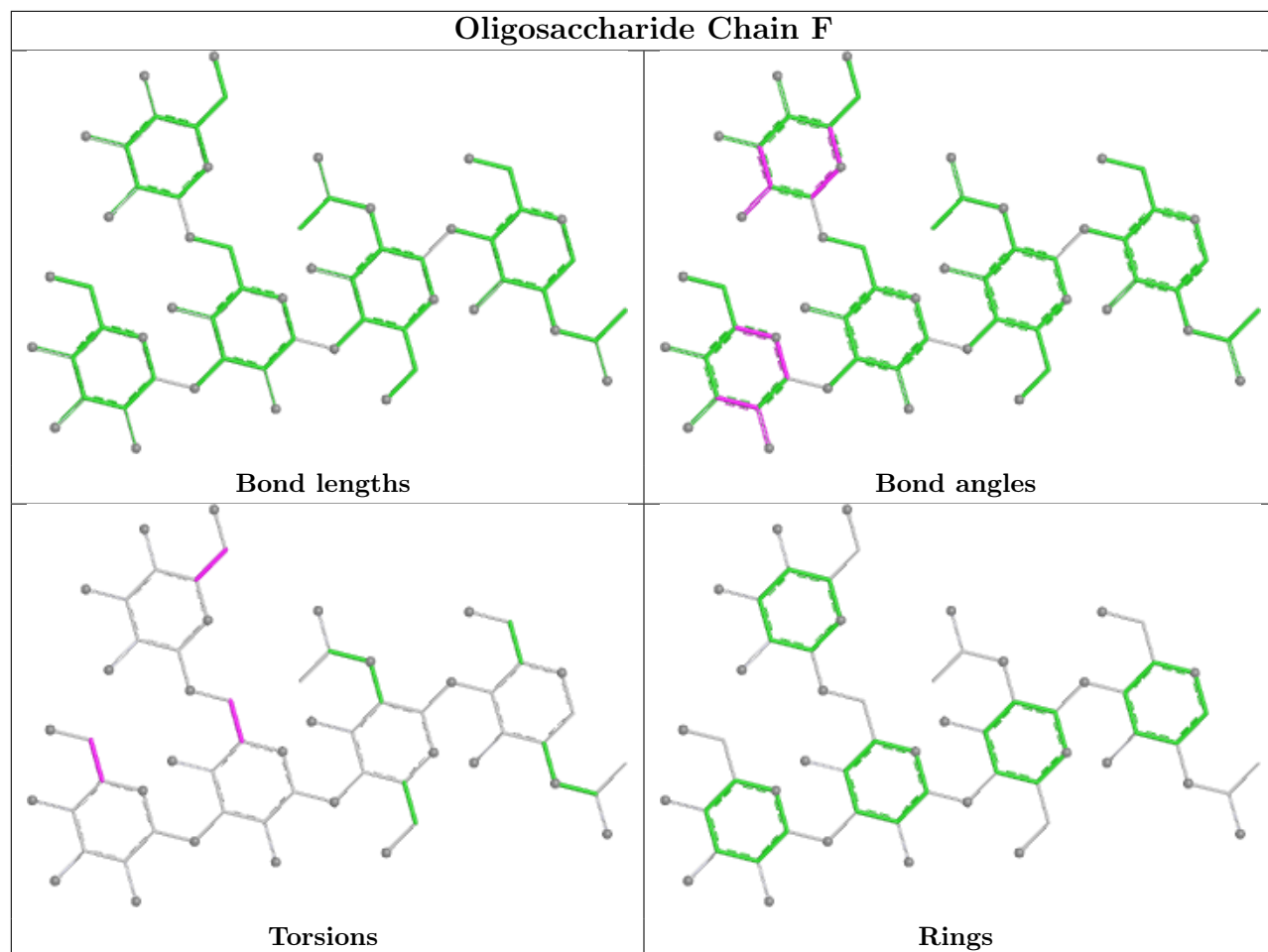
Oligosaccharide Chain A

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain C

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain J

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain K

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain M

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain O

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain F

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain I

Bond lengths

Bond angles

Torsions

Rings

Oligosaccharide Chain N

Bond lengths

Bond angles

Torsions

Rings
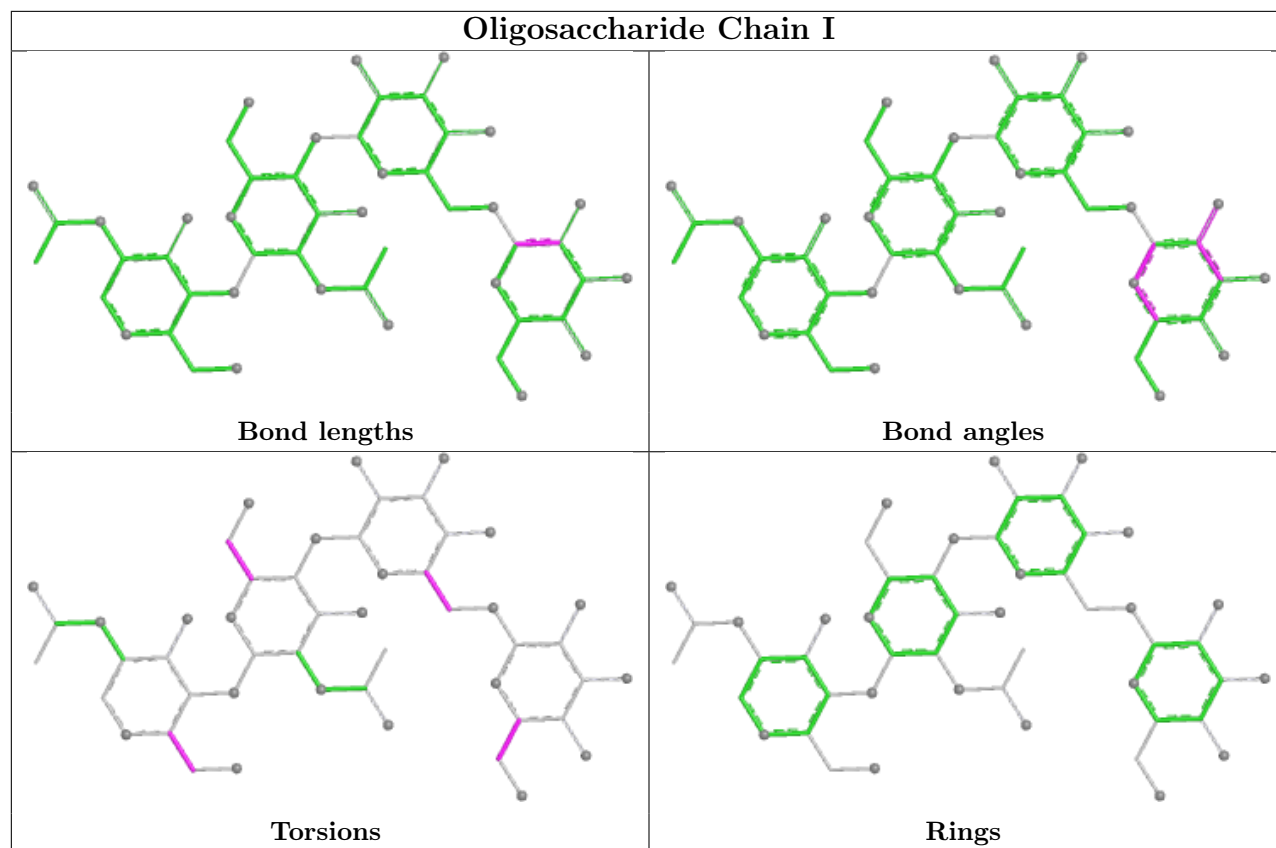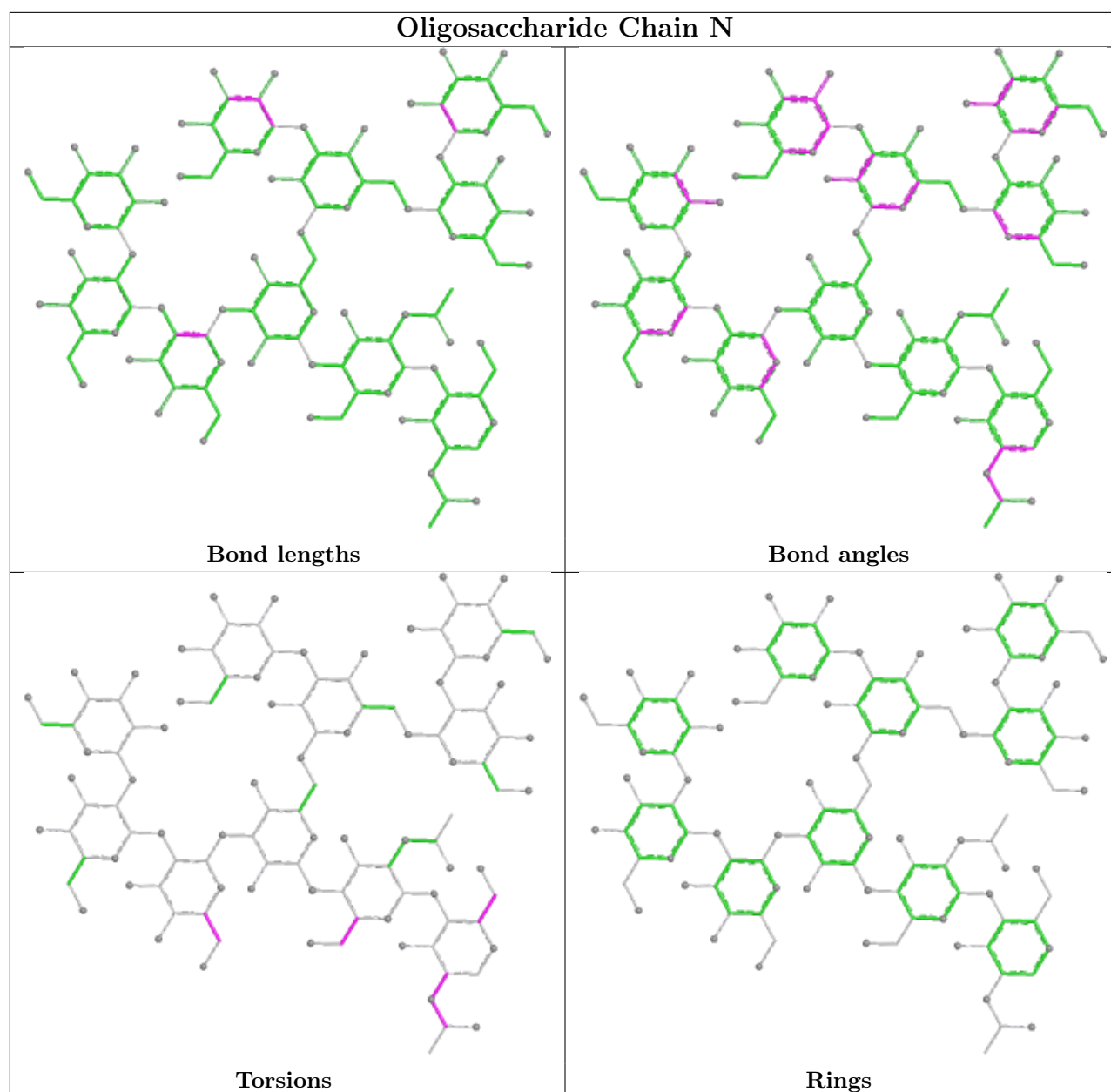
5.6 Ligand geometry (i)

7 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
14	NAG	G	614	4	14,14,15	0.29	0	17,19,21	0.51	0
14	NAG	B	702	1	14,14,15	0.33	0	17,19,21	0.44	0
14	NAG	B	701	1	14,14,15	0.27	0	17,19,21	0.58	0
14	NAG	G	637	4	14,14,15	0.27	0	17,19,21	0.52	0
14	NAG	G	615	4	14,14,15	0.78	1 (7%)	17,19,21	0.59	0
14	NAG	G	616	4	14,14,15	0.29	0	17,19,21	0.51	0
14	NAG	G	638	4	14,14,15	0.26	0	17,19,21	0.53	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
14	NAG	G	614	4	-	2/6/23/26	0/1/1/1
14	NAG	B	702	1	-	2/6/23/26	0/1/1/1
14	NAG	B	701	1	-	2/6/23/26	0/1/1/1
14	NAG	G	637	4	-	2/6/23/26	0/1/1/1
14	NAG	G	615	4	-	2/6/23/26	0/1/1/1
14	NAG	G	616	4	-	2/6/23/26	0/1/1/1
14	NAG	G	638	4	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	G	615	NAG	C1-C2	2.71	1.56	1.52

There are no bond angle outliers.

There are no chirality outliers.

5 of 14 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
14	G	638	NAG	O5-C5-C6-O6
14	G	614	NAG	O5-C5-C6-O6
14	G	616	NAG	O5-C5-C6-O6
14	G	637	NAG	O5-C5-C6-O6
14	G	615	NAG	O5-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	B	127/153 (83%)	0.69	14 (11%)	12	9	8, 31, 61, 80	0
2	D	128/153 (83%)	2.62	83 (64%)	0	0	38, 83, 116, 127	0
3	E	110/130 (84%)	2.48	64 (58%)	0	0	42, 72, 104, 116	0
4	G	430/481 (89%)	0.67	50 (11%)	11	8	8, 26, 71, 109	0
5	H	226/244 (92%)	1.62	81 (35%)	1	1	23, 61, 93, 105	0
6	L	210/217 (96%)	0.90	28 (13%)	8	6	22, 46, 68, 95	0
7	U	129/137 (94%)	2.91	98 (75%)	0	0	34, 85, 116, 132	0
8	V	101/117 (86%)	3.62	90 (89%)	0	0	64, 101, 118, 123	0
All	All	1461/1632 (89%)	1.56	508 (34%)	1	1	8, 52, 107, 132	0

The worst 5 of 508 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	E	68	TYR	9.4
3	E	2	SER	8.1
8	V	2	ILE	8.0
2	D	10	THR	7.8
4	G	465	THR	7.1

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

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

6.3 Carbohydrates ⓘ

SUGAR-RSR INFOmissingINFO

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
14	NAG	B	701	14/15	0.60	0.23	72,99,112,124	0
14	NAG	G	616	14/15	0.68	0.20	66,88,101,107	0
14	NAG	G	614	14/15	0.74	0.17	21,55,84,89	0
14	NAG	G	615	14/15	0.79	0.20	37,60,79,82	0
14	NAG	B	702	14/15	0.83	0.19	78,92,101,102	0
14	NAG	G	637	14/15	0.88	0.14	36,47,59,63	0
14	NAG	G	638	14/15	0.92	0.11	36,42,51,66	0

6.5 Other polymers ⓘ

There are no such residues in this entry.