



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 8, 2020 – 05:49 AM BST

PDB ID : 3UBN
Title : Influenza hemagglutinin from the 2009 pandemic in complex with ligand 6SLN
Authors : Xu, R.; Wilson, I.A.
Deposited on : 2011-10-24
Resolution : 2.51 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.13.1
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.13.1

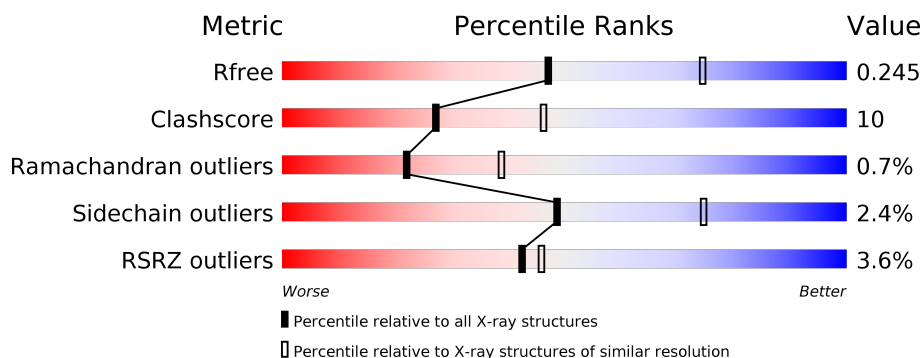
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.51 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}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 on 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	329	<div> <div>%</div> <div> <div></div> <div>78%</div> <div>19%</div> <div>..</div> </div> </div>
1	C	329	<div> <div>%</div> <div> <div></div> <div>73%</div> <div>23%</div> <div>..</div> </div> </div>
1	E	329	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>23%</div> <div>..</div> </div> </div>
1	G	329	<div> <div>%</div> <div> <div></div> <div>73%</div> <div>24%</div> <div>..</div> </div> </div>
1	I	329	<div> <div>2%</div> <div> <div></div> <div>74%</div> <div>22%</div> <div>..</div> </div> </div>
1	K	329	<div> <div>2%</div> <div> <div></div> <div>73%</div> <div>22%</div> <div>..</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	B	177	
2	D	177	
2	F	177	
2	H	177	
2	J	177	
2	L	177	
3	M	3	
3	N	3	
3	O	3	
3	R	3	
3	S	3	
4	P	2	
5	Q	2	

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
6	NAG	A	333	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 24203 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 Hemagglutinin HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	323	Total	C	N	O	S	0	0	0
			2522	1595	432	482	13			
1	C	319	Total	C	N	O	S	0	0	0
			2495	1579	428	475	13			
1	E	323	Total	C	N	O	S	0	0	0
			2522	1595	432	482	13			
1	G	321	Total	C	N	O	S	0	0	0
			2508	1586	430	479	13			
1	I	319	Total	C	N	O	S	0	0	0
			2495	1579	428	475	13			
1	K	319	Total	C	N	O	S	0	0	0
			2492	1578	428	473	13			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	9	PRO	-	expression tag	UNP C3W5S1
A	10	GLY	-	expression tag	UNP C3W5S1
A	205	CYS	GLY	engineered mutation	UNP C3W5S1
A	220	CYS	ARG	engineered mutation	UNP C3W5S1
C	9	PRO	-	expression tag	UNP C3W5S1
C	10	GLY	-	expression tag	UNP C3W5S1
C	205	CYS	GLY	engineered mutation	UNP C3W5S1
C	220	CYS	ARG	engineered mutation	UNP C3W5S1
E	9	PRO	-	expression tag	UNP C3W5S1
E	10	GLY	-	expression tag	UNP C3W5S1
E	205	CYS	GLY	engineered mutation	UNP C3W5S1
E	220	CYS	ARG	engineered mutation	UNP C3W5S1
G	9	PRO	-	expression tag	UNP C3W5S1
G	10	GLY	-	expression tag	UNP C3W5S1
G	205	CYS	GLY	engineered mutation	UNP C3W5S1
G	220	CYS	ARG	engineered mutation	UNP C3W5S1
I	9	PRO	-	expression tag	UNP C3W5S1

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Chain	Residue	Modelled	Actual	Comment	Reference
I	10	GLY	-	expression tag	UNP C3W5S1
I	205	CYS	GLY	engineered mutation	UNP C3W5S1
I	220	CYS	ARG	engineered mutation	UNP C3W5S1
K	9	PRO	-	expression tag	UNP C3W5S1
K	10	GLY	-	expression tag	UNP C3W5S1
K	205	CYS	GLY	engineered mutation	UNP C3W5S1
K	220	CYS	ARG	engineered mutation	UNP C3W5S1

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	174	Total	C	N	O	S	0	0	0
			1405	881	237	281	6			
2	D	171	Total	C	N	O	S	0	0	0
			1380	866	234	274	6			
2	F	170	Total	C	N	O	S	0	0	0
			1371	861	233	271	6			
2	H	170	Total	C	N	O	S	0	0	0
			1371	861	233	271	6			
2	J	170	Total	C	N	O	S	0	0	0
			1371	861	233	271	6			
2	L	170	Total	C	N	O	S	0	0	0
			1371	861	233	271	6			

There are 18 discrepancies between the modelled and reference sequences:

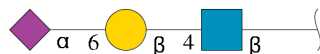
Chain	Residue	Modelled	Actual	Comment	Reference
B	175	SER	-	expression tag	UNP C3W5S1
B	176	GLY	-	expression tag	UNP C3W5S1
B	177	ARG	-	expression tag	UNP C3W5S1
D	175	SER	-	expression tag	UNP C3W5S1
D	176	GLY	-	expression tag	UNP C3W5S1
D	177	ARG	-	expression tag	UNP C3W5S1
F	175	SER	-	expression tag	UNP C3W5S1
F	176	GLY	-	expression tag	UNP C3W5S1
F	177	ARG	-	expression tag	UNP C3W5S1
H	175	SER	-	expression tag	UNP C3W5S1
H	176	GLY	-	expression tag	UNP C3W5S1
H	177	ARG	-	expression tag	UNP C3W5S1
J	175	SER	-	expression tag	UNP C3W5S1
J	176	GLY	-	expression tag	UNP C3W5S1
J	177	ARG	-	expression tag	UNP C3W5S1

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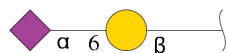
Chain	Residue	Modelled	Actual	Comment	Reference
L	175	SER	-	expression tag	UNP C3W5S1
L	176	GLY	-	expression tag	UNP C3W5S1
L	177	ARG	-	expression tag	UNP C3W5S1

- Molecule 3 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	M	3	Total	C	N	O	0	0	0
			46	25	2	19			
3	N	3	Total	C	N	O	0	0	0
			46	25	2	19			
3	O	3	Total	C	N	O	0	0	0
			46	25	2	19			
3	R	3	Total	C	N	O	0	0	0
			46	25	2	19			
3	S	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 4 is an oligosaccharide called N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose.



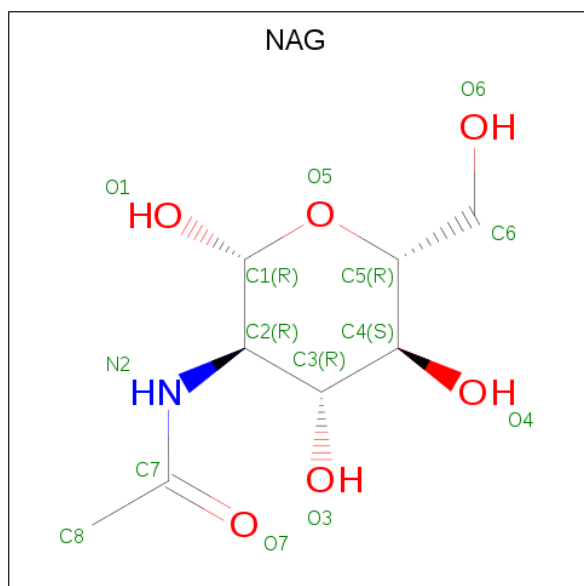
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	P	2	Total	C	N	O	0	0	0
			31	17	1	13			

- Molecule 5 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
5	Q	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



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

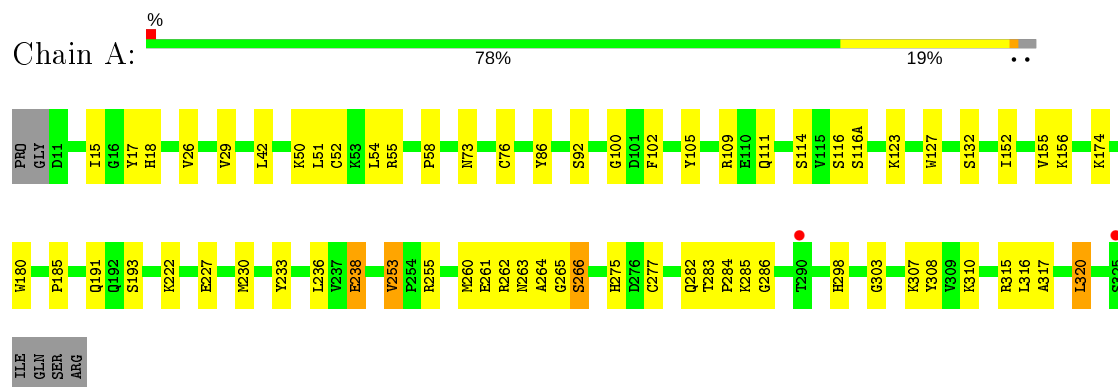
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	55	Total 55	O 55	0	0
7	B	28	Total 28	O 28	0	0
7	C	60	Total 60	O 60	0	0
7	D	31	Total 31	O 31	0	0
7	E	33	Total 33	O 33	0	0
7	F	22	Total 22	O 22	0	0
7	G	63	Total 63	O 63	0	0
7	H	27	Total 27	O 27	0	0
7	I	44	Total 44	O 44	0	0
7	J	36	Total 36	O 36	0	0
7	K	51	Total 51	O 51	0	0
7	L	35	Total 35	O 35	0	0

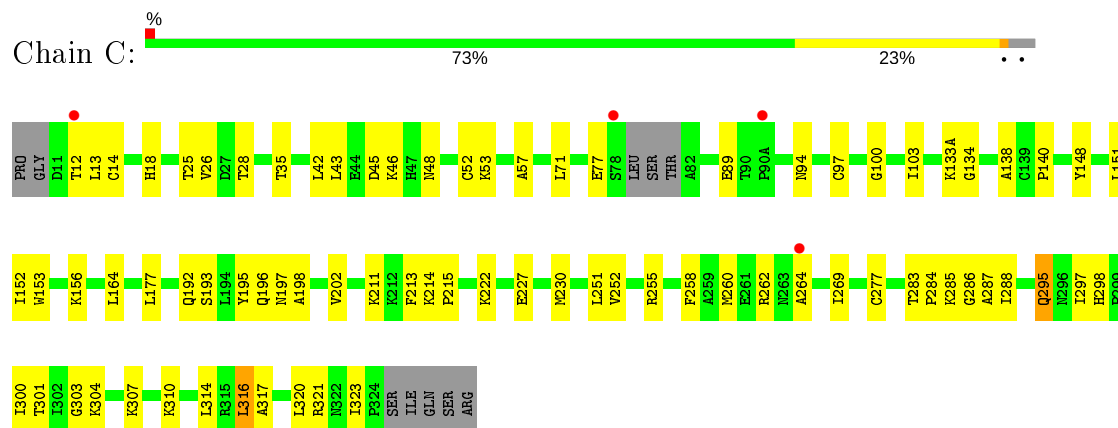
3 Residue-property plots [i](#)

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.

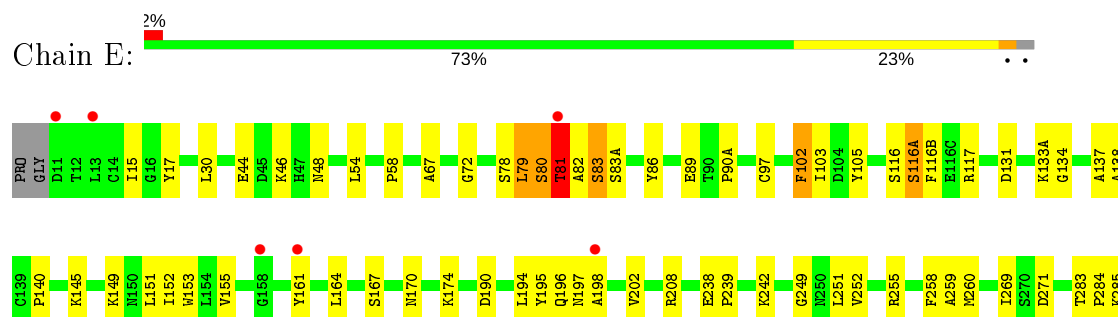
- Molecule 1: Hemagglutinin HA1

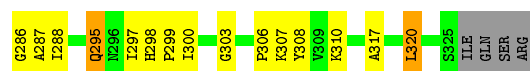


- Molecule 1: Hemagglutinin HA1

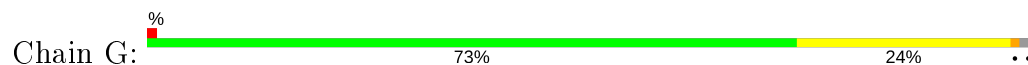


- Molecule 1: Hemagglutinin HA1

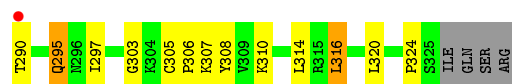
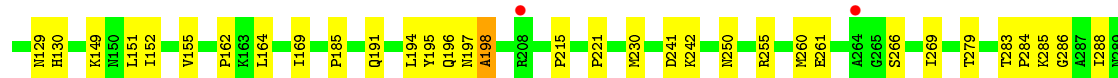
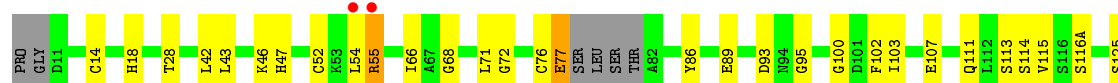




• Molecule 1: Hemagglutinin HA1



• Molecule 1: Hemagglutinin HA1

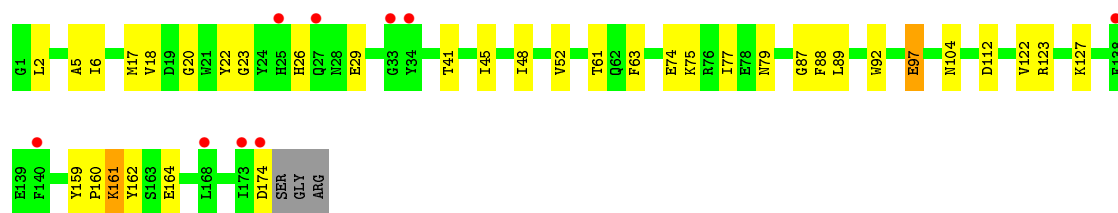


• Molecule 1: Hemagglutinin HA1

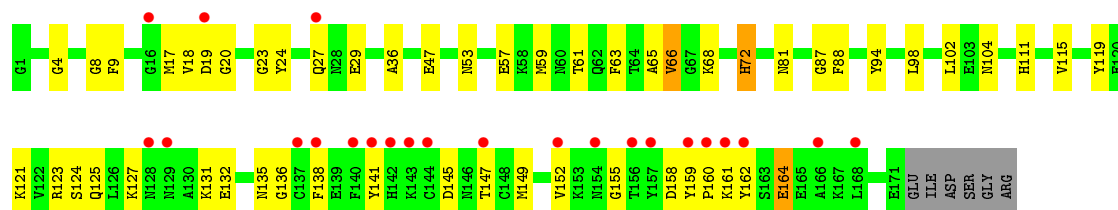


• Molecule 2: Hemagglutinin HA2

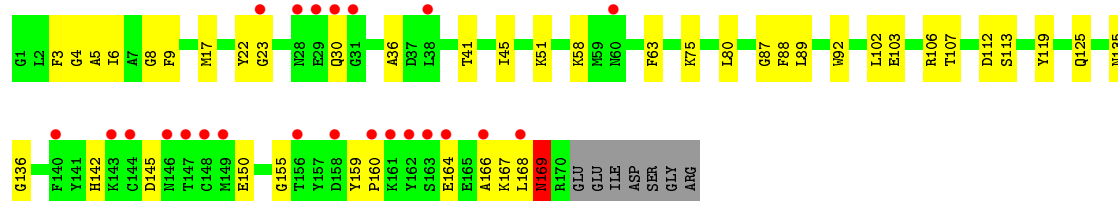




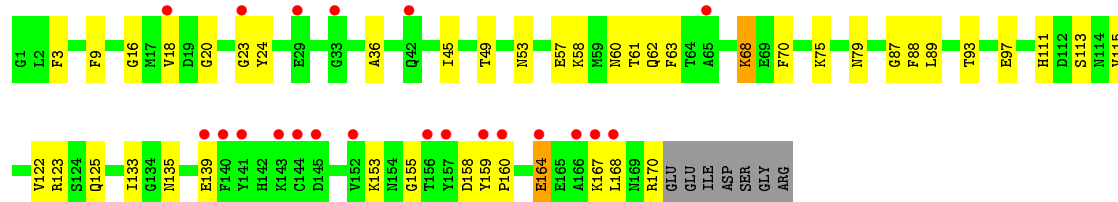
• Molecule 2: Hemagglutinin HA2



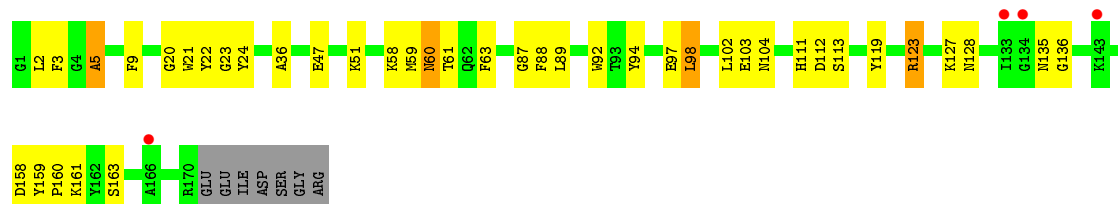
• Molecule 2: Hemagglutinin HA2



• Molecule 2: Hemagglutinin HA2

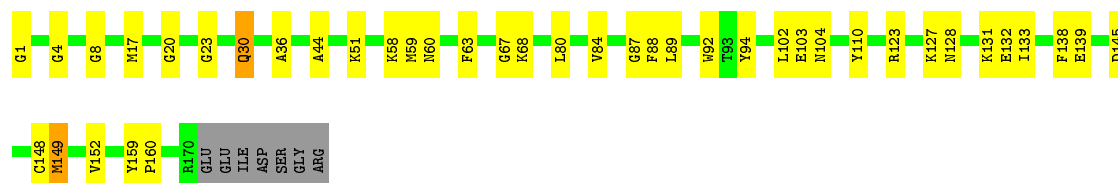


• Molecule 2: Hemagglutinin HA2



- Molecule 2: Hemagglutinin HA2

Chain L:  73% 22%



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  67% 33%



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  67% 33%



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  67% 33%



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain R:  67% 33%



- Molecule 3: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S:  67% 33%




- Molecule 4: N-acetyl-alpha-neuraminic acid-(2-6)-beta-D-galactopyranose

Chain P:  50% 50%

GAL1
SIL2

- Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain Q:  100%

NAG1
NAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	67.00Å 116.71Å 119.66Å 60.61° 77.05° 80.38°	Depositor
Resolution (Å)	49.68 – 2.51 49.68 – 2.51	Depositor EDS
% Data completeness (in resolution range)	94.2 (49.68-2.51) 94.2 (49.68-2.51)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.6.2_432	Depositor
R, R_{free}	0.205 , 0.253 0.199 , 0.245	Depositor DCC
R_{free} test set	4909 reflections (4.95%)	wwPDB-VP
Wilson B-factor (Å ²)	42.4	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	24203	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.33% 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: SIA, GAL, 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	A	0.22	0/2586	0.40	0/3516
1	C	0.22	0/2558	0.40	0/3476
1	E	0.22	0/2586	0.43	0/3516
1	G	0.22	0/2571	0.40	0/3494
1	I	0.22	0/2558	0.41	0/3476
1	K	0.22	0/2556	0.40	0/3473
2	B	0.22	0/1433	0.35	0/1931
2	D	0.22	0/1408	0.35	0/1897
2	F	0.22	0/1399	0.36	0/1885
2	H	0.22	0/1399	0.36	0/1885
2	J	0.23	0/1399	0.37	0/1885
2	L	0.23	0/1399	0.36	0/1885
All	All	0.22	0/23852	0.39	0/32319

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 [i](#)

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	2522	0	2464	47	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2495	0	2437	50	0
1	E	2522	0	2467	74	0
1	G	2508	0	2450	56	0
1	I	2495	0	2439	54	0
1	K	2492	0	2438	55	0
2	B	1405	0	1324	28	0
2	D	1380	0	1303	44	0
2	F	1371	0	1297	33	0
2	H	1371	0	1297	38	0
2	J	1371	0	1296	35	0
2	L	1371	0	1297	31	0
3	M	46	0	40	0	0
3	N	46	0	40	1	0
3	O	46	0	40	2	0
3	R	46	0	40	0	0
3	S	46	0	40	1	0
4	P	31	0	26	0	0
5	Q	28	0	25	2	0
6	A	56	0	52	0	0
6	C	28	0	26	1	0
6	E	14	0	13	1	0
6	J	14	0	13	0	0
6	K	14	0	13	0	0
7	A	55	0	0	1	0
7	B	28	0	0	0	0
7	C	60	0	0	1	0
7	D	31	0	0	0	0
7	E	33	0	0	0	0
7	F	22	0	0	1	0
7	G	63	0	0	0	0
7	H	27	0	0	0	0
7	I	44	0	0	0	0
7	J	36	0	0	1	0
7	K	51	0	0	0	0
7	L	35	0	0	1	0
All	All	24203	0	22877	470	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 10.

All (470) 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:E:79:LEU:HA	1:E:80:SER:CB	1.81	1.10
1:E:79:LEU:HA	1:E:80:SER:HB3	1.26	1.10
2:J:59:MET:HA	2:J:60:ASN:HB3	1.03	1.03
1:E:82:ALA:HB3	1:E:83:SER:HB2	1.41	1.00
2:J:59:MET:HA	2:J:60:ASN:CB	1.91	0.97
2:J:59:MET:CA	2:J:60:ASN:HB3	1.94	0.96
1:E:79:LEU:CA	1:E:80:SER:HB3	2.02	0.89
1:E:82:ALA:CB	1:E:83:SER:HB2	2.02	0.88
1:E:283:THR:HG22	1:E:285:LYS:H	1.37	0.87
2:H:97:GLU:HG2	2:L:58:LYS:HD2	1.57	0.87
1:I:283:THR:HG22	1:I:285:LYS:H	1.38	0.87
1:G:283:THR:HG22	1:G:285:LYS:H	1.41	0.85
1:C:283:THR:HG22	1:C:285:LYS:H	1.42	0.84
2:H:58:LYS:HD2	2:J:97:GLU:HG2	1.58	0.83
2:B:97:GLU:HG2	2:F:58:LYS:HD2	1.62	0.81
2:F:164:GLU:O	2:F:167:LYS:HG2	1.82	0.79
1:E:298:HIS:HD2	1:E:300:ILE:H	1.30	0.79
1:A:283:THR:HG22	1:A:285:LYS:H	1.45	0.79
1:E:72:GLY:HA3	1:E:149:LYS:H	1.48	0.77
1:A:283:THR:HB	1:A:286:GLY:O	1.87	0.75
1:I:283:THR:HB	1:I:286:GLY:O	1.87	0.74
2:H:24:TYR:CE1	2:H:153:LYS:HG3	2.22	0.74
1:C:134:GLY:HA3	1:C:153:TRP:HB3	1.73	0.71
1:E:190:ASP:O	1:E:194:LEU:HB2	1.92	0.69
2:L:23:GLY:HA3	2:L:36:ALA:HA	1.74	0.69
2:H:62:GLN:HE21	1:I:310:LYS:HZ3	1.40	0.69
1:K:114:SER:HB2	1:K:266:SER:HB2	1.75	0.69
2:H:167:LYS:HG3	2:H:170:ARG:HH12	1.59	0.68
1:K:283:THR:HG23	1:K:285:LYS:H	1.59	0.67
1:C:25:THR:HG22	1:C:35:THR:HG22	1.77	0.67
2:H:167:LYS:HG3	2:H:170:ARG:NH1	2.10	0.67
1:G:283:THR:HB	1:G:286:GLY:O	1.94	0.66
1:I:303:GLY:HA2	2:J:63:PHE:CE1	2.30	0.66
2:J:23:GLY:HA3	2:J:36:ALA:HA	1.75	0.66
2:L:131:LYS:NZ	2:L:133:ILE:HG22	2.11	0.66
1:I:283:THR:HG23	1:I:284:PRO:HD2	1.78	0.66
1:E:298:HIS:CD2	1:E:300:ILE:H	2.12	0.65
1:E:72:GLY:HA3	1:E:149:LYS:N	2.12	0.65
2:J:9:PHE:O	2:J:135:ASN:HA	1.96	0.65
1:K:307:LYS:HG3	2:L:92:TRP:CE2	2.32	0.64
1:E:174:LYS:HE3	1:E:259:ALA:HB1	1.79	0.64
1:K:13:LEU:HB2	2:L:149:MET:HE1	1.79	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:J:119:TYR:CE1	2:J:136:GLY:HA2	2.33	0.64
1:I:89:GLU:O	1:I:269:ILE:HA	1.98	0.64
2:H:53:ASN:O	2:H:57:GLU:HG2	1.97	0.64
2:D:161:LYS:HD2	2:D:162:TYR:CZ	2.34	0.63
1:G:262:ARG:HG2	1:G:262:ARG:HH11	1.64	0.63
1:E:283:THR:HB	1:E:286:GLY:O	1.99	0.62
1:K:114:SER:HB2	1:K:266:SER:CB	2.29	0.62
2:B:75:LYS:NZ	2:B:79:ASN:HD21	1.98	0.61
1:I:28:THR:HG22	2:J:104:ASN:HB3	1.80	0.61
1:K:170:ASN:HB3	1:K:239:PRO:O	2.00	0.61
2:H:23:GLY:HA3	2:H:36:ALA:HA	1.83	0.61
1:E:82:ALA:CB	1:E:83:SER:CB	2.78	0.61
1:I:169:ILE:HD13	1:I:242:LYS:HB2	1.83	0.60
1:E:80:SER:O	1:E:81:THR:HG23	2.01	0.60
1:C:303:GLY:HA2	2:D:63:PHE:CE1	2.36	0.60
1:E:310:LYS:HG2	2:F:89:LEU:HD11	1.83	0.60
1:C:283:THR:HB	1:C:286:GLY:O	2.02	0.59
1:A:283:THR:HG23	1:A:284:PRO:HD2	1.83	0.59
2:H:9:PHE:O	2:H:135:ASN:HA	2.02	0.59
1:G:97:CYS:O	1:G:224:ARG:HD3	2.02	0.59
1:G:253:VAL:HG22	1:G:254:PRO:HD2	1.85	0.59
2:H:123:ARG:HH12	2:J:123:ARG:HH21	1.50	0.59
1:A:236:LEU:HD13	1:A:262:ARG:NH1	2.18	0.58
2:B:41:THR:O	2:B:45:ILE:HG13	2.02	0.58
2:B:75:LYS:HZ3	2:B:79:ASN:HD21	1.50	0.58
2:J:59:MET:CA	2:J:60:ASN:CB	2.69	0.58
2:D:149:MET:O	2:D:152:VAL:HG22	2.04	0.58
1:E:283:THR:HG23	1:E:284:PRO:HD2	1.86	0.58
1:A:303:GLY:HA2	2:B:63:PHE:CE1	2.39	0.58
1:E:167:SER:OG	1:E:242:LYS:HE2	2.04	0.58
2:B:123:ARG:NH1	2:D:123:ARG:HH22	2.02	0.58
2:D:65:ALA:O	2:D:66:VAL:HG12	2.04	0.57
1:G:15:ILE:HD11	2:H:122:VAL:HG21	1.86	0.57
2:D:9:PHE:O	2:D:135:ASN:HA	2.04	0.57
1:G:283:THR:HG23	1:G:284:PRO:HD2	1.85	0.57
1:A:174:LYS:HD2	1:A:261:GLU:HG3	1.87	0.57
1:E:202:VAL:HG11	1:E:251:LEU:HD13	1.86	0.57
1:I:115:VAL:HG12	1:I:116(A):SER:H	1.70	0.57
2:F:30:GLN:HE21	2:F:145:ASP:HB2	1.70	0.57
2:J:119:TYR:HE1	2:J:136:GLY:HA2	1.69	0.57
1:A:116:SER:HB3	1:A:263:ASN:HD21	1.70	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:44:GLU:OE2	1:E:46:LYS:HG2	2.06	0.56
2:J:160:PRO:HA	2:J:163:SER:OG	2.06	0.56
2:B:87:GLY:HA3	2:F:88:PHE:CZ	2.41	0.56
2:D:53:ASN:O	2:D:57:GLU:HG2	2.05	0.56
1:G:307:LYS:HE2	2:H:61:THR:HG22	1.87	0.56
2:H:123:ARG:HH12	2:J:123:ARG:NH2	2.03	0.56
1:K:303:GLY:HA2	2:L:63:PHE:CE1	2.41	0.56
1:C:214:LYS:HG3	1:C:215:PRO:HD2	1.87	0.56
1:E:151:LEU:HB3	1:E:252:VAL:HG12	1.87	0.56
1:G:127:TRP:CH2	1:G:253:VAL:HG21	2.41	0.56
1:G:13:LEU:HD11	2:H:24:TYR:HB3	1.87	0.56
1:I:54:LEU:C	1:I:55:ARG:HG3	2.25	0.56
1:K:206:SER:HB2	1:K:241:ASP:OD2	2.06	0.56
1:E:79:LEU:HA	1:E:80:SER:OG	2.04	0.55
2:J:88:PHE:CZ	2:L:87:GLY:HA3	2.41	0.55
1:C:140:PRO:HD2	6:C:333:NAG:H83	1.89	0.55
1:I:52:CYS:HB2	1:I:279:THR:HG22	1.88	0.55
1:C:300:ILE:HA	2:D:66:VAL:HG11	1.88	0.55
1:G:303:GLY:HA2	2:H:63:PHE:CE1	2.41	0.55
1:K:115:VAL:HG11	1:K:116(B):PHE:HB2	1.88	0.55
1:G:200:THR:OG1	1:G:215:PRO:HG3	2.07	0.55
1:K:283:THR:HG22	1:K:286:GLY:H	1.72	0.55
1:I:307:LYS:HG3	2:J:92:TRP:CE2	2.41	0.55
1:E:161:TYR:CZ	1:E:249:GLY:HA2	2.42	0.54
1:G:170:ASN:HB3	1:G:239:PRO:O	2.07	0.54
1:C:288:ILE:HD13	1:C:295:GLN:HG3	1.90	0.54
2:D:159:TYR:HB3	2:D:160:PRO:HD3	1.89	0.54
2:L:4:GLY:O	2:L:8:GLY:HA3	2.06	0.54
1:A:111:GLN:O	1:A:262:ARG:HD2	2.07	0.54
1:I:290:THR:HG23	1:I:306:PRO:HD3	1.89	0.54
2:D:47:GLU:HB3	1:E:30:LEU:HG	1.87	0.54
1:I:197:ASN:O	1:I:198:ALA:HB3	2.07	0.54
1:G:116(B):PHE:HE1	1:G:260:MET:HE2	1.72	0.54
1:E:307:LYS:HG3	2:F:92:TRP:CE2	2.42	0.54
2:B:5:ALA:HB3	2:B:112:ASP:OD2	2.08	0.54
2:F:23:GLY:HA3	2:F:36:ALA:HA	1.90	0.54
1:K:18:HIS:HB2	2:L:20:GLY:O	2.08	0.54
1:E:208:ARG:NH2	1:E:238:GLU:H	2.06	0.53
2:L:30:GLN:OE1	2:L:145:ASP:HB2	2.08	0.53
2:D:124:SER:O	2:D:127:LYS:HE3	2.08	0.53
1:K:107:GLU:O	1:K:111:GLN:HG3	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:140:PRO:HD2	6:E:333:NAG:H83	1.90	0.53
1:A:58:PRO:HB3	1:A:86:TYR:CE1	2.44	0.53
1:C:152:ILE:HD11	1:C:255:ARG:HD2	1.90	0.53
1:G:72:GLY:O	1:G:149:LYS:HG2	2.08	0.53
2:H:88:PHE:CZ	2:J:87:GLY:HA3	2.44	0.53
2:H:159:TYR:HB3	2:H:160:PRO:HD3	1.89	0.53
1:E:82:ALA:HB1	1:E:83:SER:CA	2.39	0.53
1:G:203:PHE:HD1	1:G:204:VAL:N	2.07	0.53
1:A:310:LYS:HE3	2:B:89:LEU:HD21	1.90	0.53
1:C:18:HIS:HB2	2:D:20:GLY:O	2.08	0.53
2:D:158:ASP:OD1	2:D:161:LYS:HB2	2.08	0.53
1:C:301:THR:H	2:D:66:VAL:HG11	1.74	0.52
2:L:1:GLY:HA3	7:L:280:HOH:O	2.09	0.52
2:B:74:GLU:HB3	2:B:77:ILE:HD11	1.91	0.52
1:C:13:LEU:HD22	2:D:152:VAL:HG21	1.91	0.52
1:E:283:THR:HG22	1:E:285:LYS:N	2.16	0.52
2:B:161:LYS:HE3	2:B:162:TYR:CE1	2.43	0.52
1:E:116(A):SER:O	1:E:260:MET:HA	2.09	0.52
1:E:90(A):PRO:HD2	1:E:271:ASP:OD1	2.09	0.52
1:A:152:ILE:HG13	1:A:255:ARG:HB2	1.90	0.52
2:D:131:LYS:HE3	2:D:141:TYR:OH	2.10	0.52
1:G:283:THR:CG2	1:G:285:LYS:HG2	2.39	0.52
1:G:47:HIS:HB3	1:G:297:ILE:HD13	1.92	0.51
1:E:103:ILE:HD12	1:E:103:ILE:N	2.25	0.51
1:E:72:GLY:CA	1:E:149:LYS:H	2.21	0.51
1:E:303:GLY:HA2	2:F:63:PHE:CE1	2.45	0.51
1:I:116(A):SER:HB3	1:I:261:GLU:HG2	1.92	0.51
1:A:155:VAL:HG12	1:A:156:LYS:N	2.24	0.51
2:H:123:ARG:HH22	2:J:123:ARG:HH22	1.58	0.51
2:B:88:PHE:CZ	2:D:87:GLY:HA3	2.46	0.51
1:C:52:CYS:HB3	1:C:277:CYS:O	2.10	0.51
2:J:47:GLU:HB3	1:K:30:LEU:HG	1.93	0.51
2:L:131:LYS:HZ2	2:L:133:ILE:HG22	1.74	0.51
1:E:195:TYR:O	1:E:196:GLN:HB3	2.11	0.51
1:A:116(A):SER:O	1:A:260:MET:HA	2.11	0.51
2:L:127:LYS:HG3	2:L:128:ASN:H	1.75	0.51
1:E:170:ASN:HB3	1:E:239:PRO:O	2.11	0.51
2:D:119:TYR:CE1	2:D:136:GLY:HA2	2.46	0.51
2:H:164:GLU:H	2:H:164:GLU:CD	2.14	0.51
2:J:58:LYS:O	2:L:94:TYR:HD1	1.94	0.51
2:L:44:ALA:HA	2:L:110:TYR:OH	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:308:TYR:CD2	2:F:89:LEU:HD13	2.45	0.50
1:A:307:LYS:HE2	2:B:61:THR:HG22	1.93	0.50
2:D:59:MET:HG3	2:D:61:THR:HG23	1.93	0.50
1:A:58:PRO:HB3	1:A:86:TYR:CZ	2.46	0.50
2:B:159:TYR:HB3	2:B:160:PRO:HD3	1.92	0.50
1:A:50:LYS:HG2	1:A:275:HIS:ND1	2.27	0.50
1:C:26:VAL:HG21	1:C:317:ALA:HB2	1.94	0.50
1:E:295:GLN:O	1:E:308:TYR:HA	2.12	0.50
1:K:116(A):SER:O	1:K:260:MET:HA	2.12	0.50
2:D:132:GLU:HG2	2:D:138:PHE:CE2	2.46	0.50
1:C:28:THR:HG22	2:D:104:ASN:HB3	1.94	0.49
1:C:89:GLU:O	1:C:269:ILE:HA	2.12	0.49
1:E:48:ASN:HD21	1:E:287:ALA:HB3	1.76	0.49
1:G:89:GLU:O	1:G:269:ILE:HA	2.12	0.49
1:E:117:ARG:HD3	1:E:258:PHE:CE1	2.47	0.49
1:I:114:SER:HB2	1:I:266:SER:HB2	1.94	0.49
1:C:42:LEU:HD11	1:C:316:LEU:HG	1.93	0.49
2:D:158:ASP:OD2	2:D:160:PRO:HD2	2.13	0.49
1:E:131:ASP:HB3	1:E:155:VAL:HG23	1.95	0.49
1:C:304:LYS:HE2	2:D:61:THR:O	2.12	0.49
1:E:82:ALA:HB1	1:E:83:SER:HA	1.94	0.49
1:K:119:GLU:CD	1:K:122:PRO:HA	2.33	0.49
1:A:310:LYS:HG3	2:B:89:LEU:HD11	1.94	0.49
1:K:47:HIS:HB3	1:K:297:ILE:HD13	1.94	0.49
2:D:23:GLY:HA3	2:D:36:ALA:HA	1.95	0.49
1:E:320:LEU:HD23	1:E:320:LEU:N	2.27	0.49
1:K:86:TYR:HA	1:K:113:SER:O	2.12	0.49
1:C:283:THR:HG23	1:C:284:PRO:HD2	1.93	0.49
1:I:42:LEU:HD11	1:I:316:LEU:HG	1.95	0.49
1:K:66:ILE:HD12	1:K:109:ARG:HG2	1.95	0.49
2:H:87:GLY:HA3	2:L:88:PHE:CZ	2.48	0.48
1:E:152:ILE:HG13	1:E:255:ARG:HB2	1.96	0.48
1:K:161:TYR:CZ	1:K:249:GLY:HA2	2.48	0.48
1:A:265:GLY:O	1:A:266:SER:HB3	2.12	0.48
1:C:103:ILE:N	1:C:103:ILE:HD12	2.28	0.48
2:D:94:TYR:CZ	2:D:98:LEU:HD11	2.49	0.48
1:E:152:ILE:HD11	1:E:255:ARG:HD2	1.94	0.48
2:F:125:GLN:NE2	2:F:155:GLY:HA2	2.28	0.48
2:F:169:ASN:OD1	2:F:169:ASN:N	2.43	0.48
2:F:17:MET:SD	2:F:23:GLY:HA3	2.53	0.48
1:K:28:THR:HG22	2:L:104:ASN:HB3	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:103:ILE:H	1:E:103:ILE:HD12	1.77	0.48
1:E:82:ALA:CB	1:E:83:SER:CA	2.91	0.48
1:I:307:LYS:HE2	2:J:61:THR:HG22	1.96	0.48
1:K:156:LYS:HD2	1:K:196:GLN:HE21	1.79	0.48
5:Q:1:NAG:H62	5:Q:2:NAG:H82	1.96	0.48
2:F:6:ILE:HG13	2:F:112:ASP:HA	1.95	0.48
1:I:130:HIS:CE1	1:I:164:LEU:HB3	2.49	0.48
2:B:17:MET:SD	2:B:23:GLY:HA3	2.54	0.48
1:E:137:ALA:O	1:E:140:PRO:HD3	2.14	0.48
1:C:45:ASP:O	1:C:46:LYS:HD2	2.14	0.48
2:D:17:MET:C	2:D:19:ASP:H	2.16	0.48
1:E:283:THR:CG2	1:E:285:LYS:H	2.19	0.48
2:H:62:GLN:NE2	1:I:310:LYS:HZ3	2.10	0.48
1:G:105:TYR:CZ	1:G:109:ARG:HD2	2.49	0.47
2:H:75:LYS:HE3	2:H:79:ASN:HD21	1.79	0.47
1:I:68:GLY:HA3	1:I:95:GLY:HA2	1.95	0.47
2:F:159:TYR:N	2:F:160:PRO:HD2	2.29	0.47
1:G:103:ILE:HD12	1:G:103:ILE:N	2.28	0.47
2:H:45:ILE:O	2:H:49:THR:HG23	2.14	0.47
1:I:71:LEU:HD22	1:I:151:LEU:HD11	1.96	0.47
1:I:18:HIS:HB2	2:J:20:GLY:O	2.14	0.47
1:K:48:ASN:O	1:K:50:LYS:HG3	2.14	0.47
1:A:105:TYR:CE2	1:A:109:ARG:HD2	2.50	0.47
1:E:58:PRO:HB3	1:E:86:TYR:CZ	2.48	0.47
1:G:293:PRO:HG2	1:G:294:PHE:CD2	2.49	0.47
1:I:100:GLY:HA3	1:I:230:MET:O	2.14	0.47
1:G:202:VAL:HB	1:G:213:PHE:HB2	1.96	0.47
1:G:18:HIS:HB2	2:H:20:GLY:O	2.14	0.47
1:I:130:HIS:CE1	1:I:162:PRO:HG2	2.49	0.47
1:A:73:ASN:HB3	1:A:76:CYS:SG	2.55	0.47
1:E:48:ASN:ND2	1:E:287:ALA:HB3	2.29	0.47
1:E:82:ALA:HB1	1:E:83:SER:CB	2.44	0.47
1:G:116(B):PHE:CE1	1:G:260:MET:HE2	2.48	0.47
1:A:185:PRO:HG2	1:A:191:GLN:OE1	2.14	0.47
1:G:78:SER:HB2	1:G:149:LYS:NZ	2.29	0.47
1:G:185:PRO:HG2	1:G:191:GLN:OE1	2.14	0.47
1:I:68:GLY:CA	1:I:95:GLY:HA2	2.44	0.47
1:C:156:LYS:HE2	1:C:193:SER:O	2.15	0.47
1:C:195:TYR:O	1:C:196:GLN:HB3	2.14	0.47
2:J:94:TYR:O	2:J:98:LEU:HB2	2.15	0.47
2:H:133:ILE:HD13	2:H:139:GLU:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:316:LEU:HD22	7:C:335:HOH:O	2.14	0.47
1:I:129:ASN:HB3	1:I:162:PRO:HG3	1.97	0.47
1:I:76:CYS:O	1:I:77:GLU:C	2.54	0.47
1:C:100:GLY:HA3	1:C:230:MET:O	2.15	0.47
1:I:66:ILE:HG13	1:I:89:GLU:OE2	2.14	0.47
1:K:66:ILE:HG13	1:K:89:GLU:OE2	2.15	0.47
1:C:151:LEU:HB3	1:C:252:VAL:HG12	1.98	0.46
1:C:202:VAL:HB	1:C:213:PHE:HB2	1.95	0.46
1:E:134:GLY:HA3	1:E:153:TRP:HB3	1.95	0.46
1:E:164:LEU:C	1:E:164:LEU:HD12	2.36	0.46
1:I:107:GLU:O	1:I:111:GLN:HG3	2.16	0.46
1:A:54:LEU:O	1:A:55:ARG:HB2	2.14	0.46
1:A:307:LYS:HG3	2:B:92:TRP:CE2	2.50	0.46
1:K:54:LEU:HB3	1:K:85:SER:HB2	1.95	0.46
2:B:161:LYS:HE3	2:B:162:TYR:CZ	2.50	0.46
2:F:106:ARG:HD3	7:F:427:HOH:O	2.15	0.46
1:G:164:LEU:O	1:G:246:GLU:HA	2.15	0.46
1:I:288:ILE:HG21	1:I:297:ILE:HD12	1.97	0.46
1:K:164:LEU:HD12	1:K:164:LEU:C	2.36	0.46
2:J:5:ALA:HB3	2:J:112:ASP:OD2	2.16	0.46
2:D:19:ASP:HB3	2:D:36:ALA:HB2	1.97	0.46
1:E:89:GLU:O	1:E:269:ILE:HA	2.15	0.46
1:G:216:GLU:HB3	1:K:203:PHE:HE2	1.79	0.46
1:G:308:TYR:CD2	2:H:89:LEU:HD13	2.51	0.46
2:D:145:ASP:OD2	2:D:147:THR:HG22	2.16	0.46
2:D:88:PHE:CZ	2:F:87:GLY:HA3	2.51	0.46
1:E:116(B):PHE:CE1	1:E:260:MET:HE2	2.50	0.46
1:I:164:LEU:C	1:I:164:LEU:HD12	2.36	0.46
1:G:175:GLU:OE1	1:G:236:LEU:HD13	2.16	0.46
2:H:125:GLN:OE1	2:H:155:GLY:HA2	2.16	0.46
1:A:17:TYR:CZ	2:B:6:ILE:HG23	2.51	0.46
1:C:45:ASP:C	1:C:297:ILE:HD11	2.36	0.46
2:F:4:GLY:O	2:F:8:GLY:HA3	2.16	0.46
1:A:320:LEU:H	1:A:320:LEU:HD23	1.81	0.46
2:F:9:PHE:O	2:F:135:ASN:HA	2.16	0.46
1:C:298:HIS:CE1	1:C:300:ILE:HB	2.51	0.45
1:K:185:PRO:HG2	1:K:191:GLN:OE1	2.16	0.45
1:A:123:LYS:HE2	1:A:132:SER:O	2.16	0.45
1:C:12:THR:OG1	2:D:27:GLN:HG2	2.16	0.45
1:G:73:ASN:HB3	1:G:76:CYS:SG	2.56	0.45
1:I:47:HIS:HB3	1:I:288:ILE:HG22	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:97:GLU:HG2	2:F:58:LYS:CD	2.42	0.45
1:E:317:ALA:O	2:F:107:THR:HG21	2.15	0.45
2:H:168:LEU:HD23	2:H:168:LEU:O	2.16	0.45
1:E:197:ASN:O	1:E:198:ALA:HB3	2.17	0.45
2:F:30:GLN:NE2	2:F:145:ASP:HB2	2.30	0.45
1:E:46:LYS:N	1:E:297:ILE:HD11	2.31	0.45
2:L:148:CYS:O	2:L:152:VAL:HG23	2.17	0.45
1:A:238:GLU:HG2	1:A:238:GLU:H	1.54	0.45
1:C:71:LEU:O	1:C:148:TYR:HB3	2.16	0.45
1:E:164:LEU:O	1:E:164:LEU:HD12	2.17	0.45
1:G:134:GLY:HA3	1:G:153:TRP:HB3	1.97	0.45
1:G:320:LEU:HB3	2:H:111:HIS:CG	2.52	0.45
2:L:159:TYR:N	2:L:160:PRO:HD2	2.30	0.45
1:A:18:HIS:HB2	2:B:20:GLY:O	2.16	0.45
1:C:164:LEU:HD12	1:C:164:LEU:C	2.37	0.45
2:J:159:TYR:N	2:J:160:PRO:HD2	2.32	0.45
1:A:15:ILE:HD11	2:B:122:VAL:HG21	1.98	0.45
1:C:177:LEU:HB3	1:C:258:PHE:HB2	1.99	0.45
1:I:86:TYR:HA	1:I:113:SER:O	2.17	0.45
1:K:51:LEU:HB2	1:K:274:VAL:HA	1.99	0.45
1:A:180:TRP:CE2	1:A:233:TYR:HB2	2.51	0.45
1:C:53:LYS:HG2	1:C:57:ALA:HA	1.98	0.45
2:J:127:LYS:HG2	2:J:128:ASN:N	2.32	0.45
2:D:125:GLN:NE2	2:D:155:GLY:HA2	2.32	0.45
1:G:116:SER:HB3	1:G:263:ASN:OD1	2.16	0.45
2:L:51:LYS:HE3	2:L:103:GLU:OE2	2.17	0.45
2:F:119:TYR:CE1	2:F:136:GLY:HA2	2.51	0.44
2:J:158:ASP:CG	2:J:161:LYS:HB2	2.38	0.44
1:K:149:LYS:HE2	1:K:256:TYR:HE1	1.82	0.44
1:A:238:GLU:HG2	7:A:352:HOH:O	2.17	0.44
1:G:66:ILE:HG13	1:G:89:GLU:OE2	2.17	0.44
1:I:195:TYR:O	1:I:196:GLN:HB3	2.18	0.44
1:K:151:LEU:HB3	1:K:252:VAL:HG12	1.99	0.44
1:E:320:LEU:HD23	1:E:320:LEU:H	1.82	0.44
1:G:97:CYS:HB2	1:G:138:ALA:O	2.16	0.44
1:I:114:SER:HB2	1:I:266:SER:CB	2.47	0.44
1:K:103:ILE:HG12	1:K:233:TYR:CE2	2.52	0.44
1:K:308:TYR:CD2	2:L:89:LEU:HD13	2.53	0.44
2:L:127:LYS:HG3	2:L:128:ASN:N	2.33	0.44
1:C:43:LEU:HB2	1:C:314:LEU:HB2	2.00	0.44
1:G:161:TYR:CZ	1:G:249:GLY:HA2	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:169:ILE:HD12	1:G:169:ILE:N	2.33	0.44
2:L:67:GLY:O	2:L:68:LYS:HE3	2.18	0.44
1:A:156:LYS:HE2	1:A:193:SER:O	2.17	0.44
1:E:288:ILE:CD1	1:E:295:GLN:HG3	2.47	0.44
2:H:123:ARG:HH22	2:J:123:ARG:NH2	2.15	0.44
1:I:316:LEU:HD22	7:J:189:HOH:O	2.17	0.44
1:A:114:SER:HB2	1:A:266:SER:CB	2.48	0.44
1:A:51:LEU:HA	1:A:282:GLN:NE2	2.32	0.44
2:H:3:PHE:CZ	2:J:2:LEU:HG	2.53	0.44
1:E:288:ILE:HD11	1:E:306:PRO:HD2	1.99	0.44
1:I:116(A):SER:HB3	1:I:261:GLU:CG	2.48	0.44
2:L:133:ILE:HD13	2:L:139:GLU:HB2	2.00	0.44
2:D:18:VAL:O	2:D:18:VAL:HG22	2.18	0.43
1:K:54:LEU:O	1:K:55:ARG:HG3	2.18	0.43
1:K:97:CYS:O	1:K:224:ARG:HD3	2.19	0.43
2:B:2:LEU:HG	2:F:3:PHE:CZ	2.54	0.43
1:A:114:SER:OG	1:A:263:ASN:HB2	2.19	0.43
1:I:185:PRO:HG2	1:I:191:GLN:OE1	2.19	0.43
1:I:305:CYS:O	2:J:61:THR:HG21	2.18	0.43
1:K:175:GLU:OE1	1:K:262:ARG:NH1	2.48	0.43
1:E:145:LYS:NZ	3:O:3:SIA:H31	2.32	0.43
1:E:15:ILE:HD12	1:E:15:ILE:N	2.32	0.43
1:E:67:ALA:HB2	1:E:105:TYR:CE1	2.53	0.43
1:G:54:LEU:O	1:G:55:ARG:HB2	2.17	0.43
2:H:68:LYS:HB3	2:H:70:PHE:CE2	2.54	0.43
2:D:121:LYS:O	2:D:125:GLN:HG3	2.18	0.43
1:E:102:PHE:O	1:E:105:TYR:HB2	2.18	0.43
1:A:308:TYR:CD2	2:B:89:LEU:HD13	2.54	0.43
2:D:125:GLN:CD	2:D:155:GLY:HA2	2.39	0.43
2:D:81:ASN:OD1	2:F:80:LEU:HD13	2.19	0.43
1:E:133(A):LYS:O	3:O:3:SIA:H113	2.19	0.43
2:J:51:LYS:HE3	2:J:103:GLU:OE2	2.19	0.43
1:A:26:VAL:HB	2:B:104:ASN:ND2	2.34	0.43
2:B:127:LYS:HB2	2:B:159:TYR:CE1	2.53	0.43
2:D:4:GLY:O	2:D:8:GLY:HA3	2.19	0.43
2:F:3:PHE:CE1	2:F:113:SER:HB2	2.53	0.43
1:K:179:LEU:HD23	1:K:234:TRP:HB3	2.01	0.43
1:K:283:THR:OG1	1:K:298:HIS:HB3	2.18	0.43
1:G:120:ILE:HG13	1:G:168:TYR:CD1	2.53	0.43
1:G:206:SER:HA	1:I:221:PRO:HG2	2.01	0.43
1:K:283:THR:HG22	1:K:286:GLY:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:89:GLU:O	1:K:269:ILE:HA	2.18	0.43
2:H:93:THR:O	2:H:97:GLU:HB2	2.19	0.43
2:J:3:PHE:CE2	2:J:113:SER:HB2	2.53	0.43
1:K:54:LEU:HD21	1:K:302:ILE:HG22	2.01	0.43
2:D:131:LYS:HB3	2:D:141:TYR:CZ	2.54	0.42
2:F:75:LYS:HD3	2:F:75:LYS:HA	1.79	0.42
1:I:241:ASP:OD1	1:I:242:LYS:N	2.50	0.42
1:A:52:CYS:HB3	1:A:277:CYS:O	2.18	0.42
1:C:197:ASN:O	1:C:198:ALA:HB3	2.18	0.42
1:G:155:VAL:HG12	1:G:156:LYS:N	2.34	0.42
2:H:16:GLY:O	2:H:18:VAL:HG23	2.20	0.42
1:I:308:TYR:CD2	2:J:89:LEU:HD13	2.54	0.42
1:K:288:ILE:O	1:K:288:ILE:HG13	2.18	0.42
1:A:127:TRP:CZ2	1:A:253:VAL:HG11	2.54	0.42
1:C:307:LYS:HE2	2:D:61:THR:HG22	2.00	0.42
1:C:14:CYS:O	2:D:24:TYR:HA	2.19	0.42
2:F:41:THR:O	2:F:45:ILE:HG13	2.19	0.42
1:G:253:VAL:HA	1:G:254:PRO:HD3	1.93	0.42
1:I:72:GLY:HA3	1:I:149:LYS:H	1.85	0.42
1:I:116(A):SER:O	1:I:260:MET:HA	2.20	0.42
1:K:100:GLY:HA3	1:K:230:MET:O	2.19	0.42
1:K:288:ILE:HG21	1:K:297:ILE:HD12	2.00	0.42
1:I:14:CYS:O	2:J:24:TYR:HA	2.19	0.42
2:B:18:VAL:O	2:B:18:VAL:HG22	2.19	0.42
1:G:262:ARG:HG2	1:G:262:ARG:NH1	2.32	0.42
1:G:86:TYR:HA	1:G:113:SER:O	2.18	0.42
1:K:149:LYS:HE2	1:K:256:TYR:CE1	2.55	0.42
1:K:194:LEU:HD11	3:S:3:SIA:H91	2.00	0.42
1:C:321:ARG:HD2	1:C:323:ILE:HD11	2.01	0.42
2:H:158:ASP:OD1	2:H:160:PRO:HD2	2.19	0.42
1:K:188:SER:O	1:K:192:GLN:HG2	2.20	0.42
2:L:132:GLU:HG2	2:L:138:PHE:CE2	2.55	0.42
1:E:17:TYR:HB2	1:E:320:LEU:HD11	2.02	0.42
1:G:170:ASN:HD21	1:G:176:VAL:HG23	1.85	0.42
1:A:260:MET:HE2	1:A:262:ARG:HG2	2.02	0.42
1:I:46:LYS:N	1:I:297:ILE:HD11	2.35	0.42
2:B:48:ILE:O	2:B:52:VAL:HG23	2.20	0.42
1:G:242:LYS:HG2	1:G:243:ILE:N	2.35	0.42
1:I:155:VAL:HG13	1:I:194:LEU:O	2.20	0.42
2:L:17:MET:SD	2:L:23:GLY:HA3	2.60	0.42
1:A:42:LEU:HD11	1:A:316:LEU:HD22	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:27:GLN:HG3	2:D:27:GLN:O	2.20	0.41
1:G:164:LEU:HD23	1:G:164:LEU:C	2.39	0.41
1:K:42:LEU:O	1:K:292:LEU:HB3	2.20	0.41
2:F:51:LYS:HE3	2:F:103:GLU:OE2	2.20	0.41
1:I:152:ILE:HG13	1:I:255:ARG:HB2	2.02	0.41
1:K:103:ILE:HD12	1:K:103:ILE:N	2.35	0.41
1:K:195:TYR:O	1:K:196:GLN:HB3	2.20	0.41
2:L:131:LYS:HZ1	2:L:133:ILE:HG22	1.86	0.41
1:I:43:LEU:HB2	1:I:314:LEU:HB2	2.02	0.41
1:K:283:THR:CG2	1:K:285:LYS:H	2.31	0.41
2:D:111:HIS:O	2:D:115:VAL:HG23	2.20	0.41
1:E:83(A):SER:CB	1:E:116:SER:HA	2.50	0.41
1:G:29:VAL:CG2	2:L:51:LYS:HG3	2.51	0.41
1:A:222:LYS:HE2	1:A:227:GLU:HG3	2.03	0.41
2:F:5:ALA:HB3	2:F:112:ASP:OD2	2.20	0.41
1:E:310:LYS:HD3	2:F:89:LEU:HD21	2.03	0.41
1:G:140:PRO:HD2	5:Q:1:NAG:H83	2.02	0.41
1:C:164:LEU:O	1:C:164:LEU:HD12	2.20	0.41
1:E:116(B):PHE:HE1	1:E:260:MET:HE2	1.86	0.41
1:G:155:VAL:CG1	1:G:156:LYS:N	2.83	0.41
1:G:17:TYR:HB3	2:H:115:VAL:HG21	2.02	0.41
1:I:215:PRO:HB3	1:I:250:ASN:ND2	2.35	0.41
1:C:211:LYS:HD3	1:C:213:PHE:CZ	2.55	0.41
2:F:142:HIS:HD2	2:F:166:ALA:HB2	1.85	0.41
1:C:202:VAL:HG11	1:C:251:LEU:HD13	2.03	0.41
1:C:222:LYS:NZ	1:C:227:GLU:HG3	2.36	0.41
2:D:164:GLU:CD	2:D:164:GLU:H	2.24	0.41
1:E:97:CYS:HB2	1:E:138:ALA:O	2.21	0.41
2:F:167:LYS:HG3	2:F:168:LEU:HD12	2.02	0.41
1:G:58:PRO:HB3	1:G:86:TYR:CZ	2.56	0.41
1:K:47:HIS:HB3	1:K:288:ILE:HG22	2.03	0.41
1:A:265:GLY:O	1:A:266:SER:CB	2.69	0.41
1:C:48:ASN:ND2	1:C:287:ALA:HB3	2.36	0.41
2:D:72:HIS:CD2	2:D:72:HIS:H	2.39	0.41
1:G:26:VAL:O	1:G:32:LYS:O	2.38	0.41
2:J:21:TRP:CE3	2:J:111:HIS:HE1	2.39	0.41
1:K:310:LYS:O	1:K:310:LYS:HG2	2.21	0.41
1:C:97:CYS:HB2	1:C:138:ALA:O	2.22	0.41
2:H:62:GLN:HE21	1:I:310:LYS:NZ	2.14	0.41
1:I:295:GLN:HE21	1:I:295:GLN:HB3	1.73	0.41
1:A:284:PRO:HG2	1:A:298:HIS:CE1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:316:LEU:HD23	1:C:316:LEU:HA	1.84	0.40
2:D:125:GLN:O	2:D:127:LYS:HD2	2.21	0.40
1:G:123:LYS:HB2	1:G:152:ILE:HD11	2.03	0.40
2:L:80:LEU:O	2:L:84:VAL:HG23	2.20	0.40
1:K:134:GLY:HA3	1:K:153:TRP:HB3	2.04	0.40
1:A:100:GLY:HA3	1:A:230:MET:O	2.20	0.40
1:C:310:LYS:HB3	1:C:310:LYS:HE2	1.74	0.40
1:A:29:VAL:CG2	2:F:51:LYS:HG3	2.52	0.40
2:H:3:PHE:CE1	2:H:113:SER:HB2	2.57	0.40
2:L:123:ARG:HB2	2:L:138:PHE:HZ	1.86	0.40
1:I:103:ILE:HD12	1:I:103:ILE:N	2.37	0.40
1:K:90(A):PRO:HD2	1:K:271:ASP:OD1	2.21	0.40
1:K:90:THR:HA	1:K:90(A):PRO:HD3	1.97	0.40
2:L:149:MET:HA	2:L:149:MET:CE	2.51	0.40
1:A:26:VAL:HG21	1:A:317:ALA:HB2	2.03	0.40
1:C:133(A):LYS:O	3:N:3:SIA:H113	2.21	0.40
1:C:260:MET:CE	1:C:262:ARG:HG2	2.52	0.40
1:E:298:HIS:CG	1:E:299:PRO:HD2	2.57	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	321/329 (98%)	300 (94%)	18 (6%)	3 (1%)	17	31
1	C	315/329 (96%)	300 (95%)	13 (4%)	2 (1%)	25	43
1	E	321/329 (98%)	294 (92%)	22 (7%)	5 (2%)	9	17
1	G	317/329 (96%)	301 (95%)	16 (5%)	0	100	100
1	I	315/329 (96%)	291 (92%)	21 (7%)	3 (1%)	15	28
1	K	315/329 (96%)	297 (94%)	16 (5%)	2 (1%)	25	43

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	172/177 (97%)	167 (97%)	5 (3%)	0	100	100
2	D	169/177 (96%)	156 (92%)	13 (8%)	0	100	100
2	F	168/177 (95%)	157 (94%)	10 (6%)	1 (1%)	25	43
2	H	168/177 (95%)	159 (95%)	9 (5%)	0	100	100
2	J	168/177 (95%)	158 (94%)	8 (5%)	2 (1%)	13	24
2	L	168/177 (95%)	161 (96%)	5 (3%)	2 (1%)	13	24
All	All	2917/3036 (96%)	2741 (94%)	156 (5%)	20 (1%)	22	39

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	J	60	ASN
1	C	264	ALA
1	E	80	SER
1	I	93	ASP
1	K	265	GLY
1	A	266	SER
1	C	94	ASN
1	E	78	SER
1	E	83	SER
1	E	116(A)	SER
1	I	198	ALA
2	J	5	ALA
1	A	92	SER
1	A	264	ALA
2	F	169	ASN
1	I	324	PRO
1	K	239	PRO
2	L	60	ASN
1	E	81	THR
2	L	59	MET

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	285/290 (98%)	280 (98%)	5 (2%)	59	81
1	C	281/290 (97%)	276 (98%)	5 (2%)	59	81
1	E	285/290 (98%)	279 (98%)	6 (2%)	53	78
1	G	283/290 (98%)	277 (98%)	6 (2%)	53	78
1	I	281/290 (97%)	274 (98%)	7 (2%)	47	73
1	K	281/290 (97%)	275 (98%)	6 (2%)	53	78
2	B	150/152 (99%)	143 (95%)	7 (5%)	26	49
2	D	147/152 (97%)	141 (96%)	6 (4%)	30	55
2	F	146/152 (96%)	142 (97%)	4 (3%)	44	71
2	H	146/152 (96%)	143 (98%)	3 (2%)	53	78
2	J	146/152 (96%)	142 (97%)	4 (3%)	44	71
2	L	146/152 (96%)	143 (98%)	3 (2%)	53	78
All	All	2577/2652 (97%)	2515 (98%)	62 (2%)	49	74

All (62) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	102	PHE
1	A	238	GLU
1	A	253	VAL
1	A	315	ARG
1	A	320	LEU
2	B	22	TYR
2	B	26	HIS
2	B	29	GLU
2	B	97	GLU
2	B	161	LYS
2	B	164	GLU
2	B	174	ASP
1	C	77	GLU
1	C	192	GLN
1	C	295	GLN
1	C	316	LEU
1	C	320	LEU
2	D	29	GLU
2	D	66	VAL
2	D	68	LYS
2	D	72	HIS
2	D	102	LEU

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Mol	Chain	Res	Type
2	D	164	GLU
1	E	54	LEU
1	E	79	LEU
1	E	81	THR
1	E	102	PHE
1	E	295	GLN
1	E	320	LEU
2	F	22	TYR
2	F	102	LEU
2	F	150	GLU
2	F	169	ASN
1	G	136	THR
1	G	174	LYS
1	G	203	PHE
1	G	225	ASP
1	G	253	VAL
1	G	261	GLU
2	H	60	ASN
2	H	68	LYS
2	H	164	GLU
1	I	55	ARG
1	I	77	GLU
1	I	102	PHE
1	I	125	SER
1	I	295	GLN
1	I	316	LEU
1	I	320	LEU
2	J	22	TYR
2	J	98	LEU
2	J	102	LEU
2	J	123	ARG
1	K	55	ARG
1	K	97	CYS
1	K	102	PHE
1	K	274	VAL
1	K	283	THR
1	K	295	GLN
2	L	30	GLN
2	L	102	LEU
2	L	149	MET

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	263	ASN
2	B	26	HIS
2	B	43	ASN
2	B	79	ASN
2	B	95	ASN
2	D	72	HIS
2	D	95	ASN
1	E	298	HIS
2	F	30	GLN
2	F	50	ASN
2	H	60	ASN
2	H	62	GLN
1	I	21	ASN
1	I	289	ASN
2	J	30	GLN
2	J	95	ASN
2	J	117	ASN
1	K	196	GLN
2	L	62	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

19 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	M	1	3	15,15,15	0.45	0	21,21,21	0.58	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GAL	M	2	3	11,11,12	0.72	0	15,15,17	0.78	0
3	SIA	M	3	3	17,20,21	0.42	0	21,28,31	1.02	1 (4%)
3	NAG	N	1	3	15,15,15	0.49	0	21,21,21	0.61	0
3	GAL	N	2	3	11,11,12	0.68	0	15,15,17	0.80	0
3	SIA	N	3	3	17,20,21	0.41	0	21,28,31	0.97	2 (9%)
3	NAG	O	1	3	15,15,15	0.47	0	21,21,21	0.74	0
3	GAL	O	2	3	11,11,12	0.63	0	15,15,17	0.69	0
3	SIA	O	3	3	17,20,21	0.35	0	21,28,31	1.06	2 (9%)
4	GAL	P	1	4	11,11,12	0.64	0	15,15,17	0.57	0
4	SIA	P	2	4	17,20,21	0.35	0	21,28,31	1.06	2 (9%)
5	NAG	Q	1	1,5	14,14,15	0.58	0	17,19,21	0.75	0
5	NAG	Q	2	5	14,14,15	0.57	0	17,19,21	0.73	0
3	NAG	R	1	3	15,15,15	0.50	0	21,21,21	0.68	0
3	GAL	R	2	3	11,11,12	0.65	0	15,15,17	0.70	0
3	SIA	R	3	3	17,20,21	0.43	0	21,28,31	1.14	2 (9%)
3	NAG	S	1	3	15,15,15	0.47	0	21,21,21	0.99	1 (4%)
3	GAL	S	2	3	11,11,12	0.68	0	15,15,17	0.83	1 (6%)
3	SIA	S	3	3	17,20,21	0.43	0	21,28,31	1.07	1 (4%)

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	M	1	3	-	0/6/26/26	0/1/1/1
3	GAL	M	2	3	-	0/2/19/22	0/1/1/1
3	SIA	M	3	3	-	0/14/34/38	0/1/1/1
3	NAG	N	1	3	-	0/6/26/26	0/1/1/1
3	GAL	N	2	3	-	0/2/19/22	0/1/1/1
3	SIA	N	3	3	-	0/14/34/38	0/1/1/1
3	NAG	O	1	3	-	0/6/26/26	0/1/1/1
3	GAL	O	2	3	-	0/2/19/22	0/1/1/1
3	SIA	O	3	3	-	0/14/34/38	0/1/1/1
4	GAL	P	1	4	-	0/2/19/22	0/1/1/1
4	SIA	P	2	4	-	0/14/34/38	0/1/1/1
5	NAG	Q	1	1,5	-	2/6/23/26	0/1/1/1
5	NAG	Q	2	5	-	2/6/23/26	0/1/1/1
3	NAG	R	1	3	-	0/6/26/26	0/1/1/1
3	GAL	R	2	3	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	SIA	R	3	3	-	0/14/34/38	0/1/1/1
3	NAG	S	1	3	-	0/6/26/26	0/1/1/1
3	GAL	S	2	3	-	1/2/19/22	0/1/1/1
3	SIA	S	3	3	-	0/14/34/38	0/1/1/1

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	R	3	SIA	C6-O6-C2	3.21	118.21	111.34
3	O	3	SIA	C6-O6-C2	3.12	118.02	111.34
4	P	2	SIA	C6-O6-C2	3.10	117.97	111.34
3	S	1	NAG	O5-C1-C2	2.75	112.28	109.52
3	S	3	SIA	C6-O6-C2	2.60	116.90	111.34
3	R	3	SIA	O6-C2-C3	2.45	114.17	109.87
3	M	3	SIA	C6-O6-C2	2.44	116.55	111.34
3	N	3	SIA	C6-O6-C2	2.04	115.71	111.34
3	N	3	SIA	O6-C2-C3	2.04	113.45	109.87
3	O	3	SIA	O6-C2-C3	2.01	113.40	109.87
3	S	2	GAL	C1-O5-C5	-2.01	109.47	112.19
4	P	2	SIA	O6-C2-C3	2.01	113.40	109.87

There are no chirality outliers.

All (5) torsion outliers are listed below:

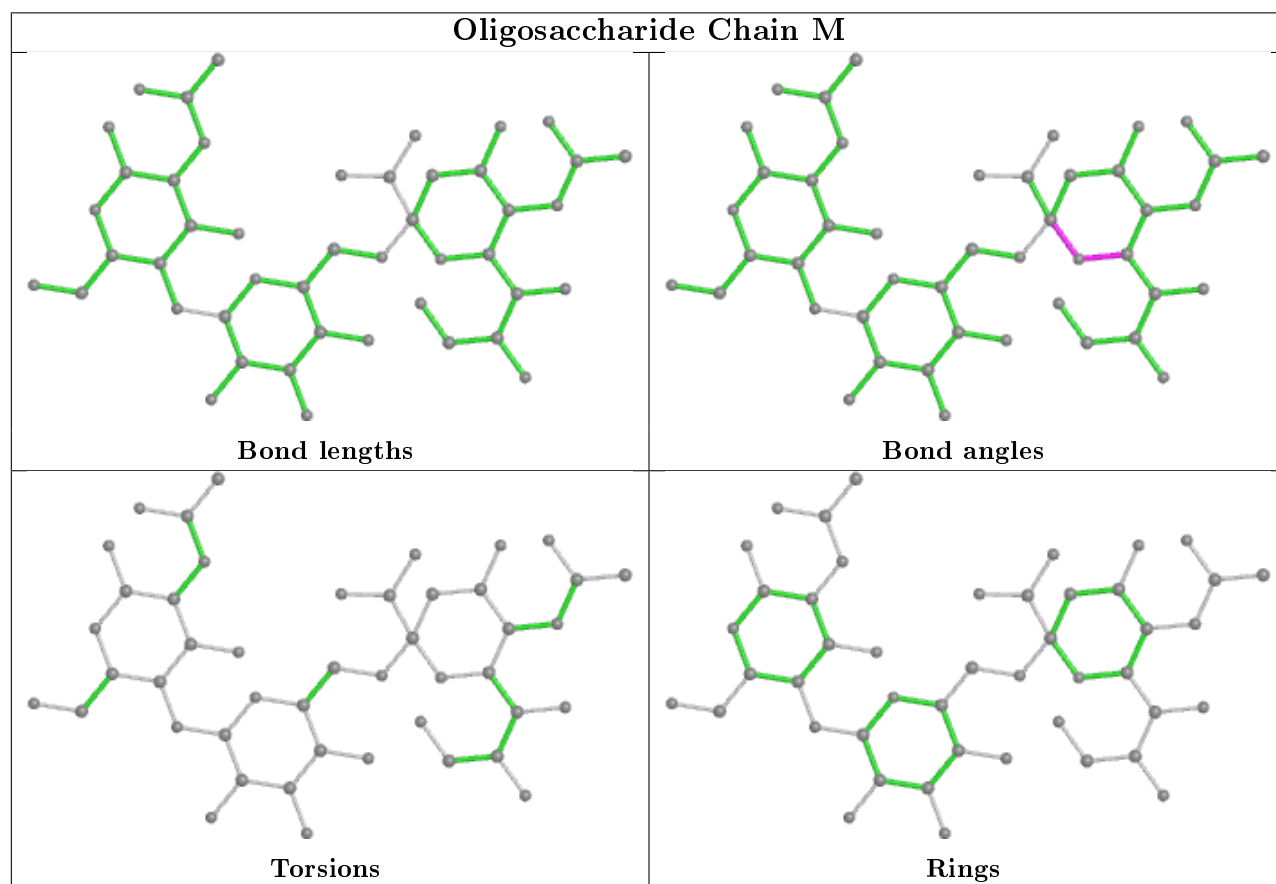
Mol	Chain	Res	Type	Atoms
5	Q	1	NAG	C8-C7-N2-C2
5	Q	1	NAG	O7-C7-N2-C2
5	Q	2	NAG	C8-C7-N2-C2
5	Q	2	NAG	O7-C7-N2-C2
3	S	2	GAL	C4-C5-C6-O6

There are no ring outliers.

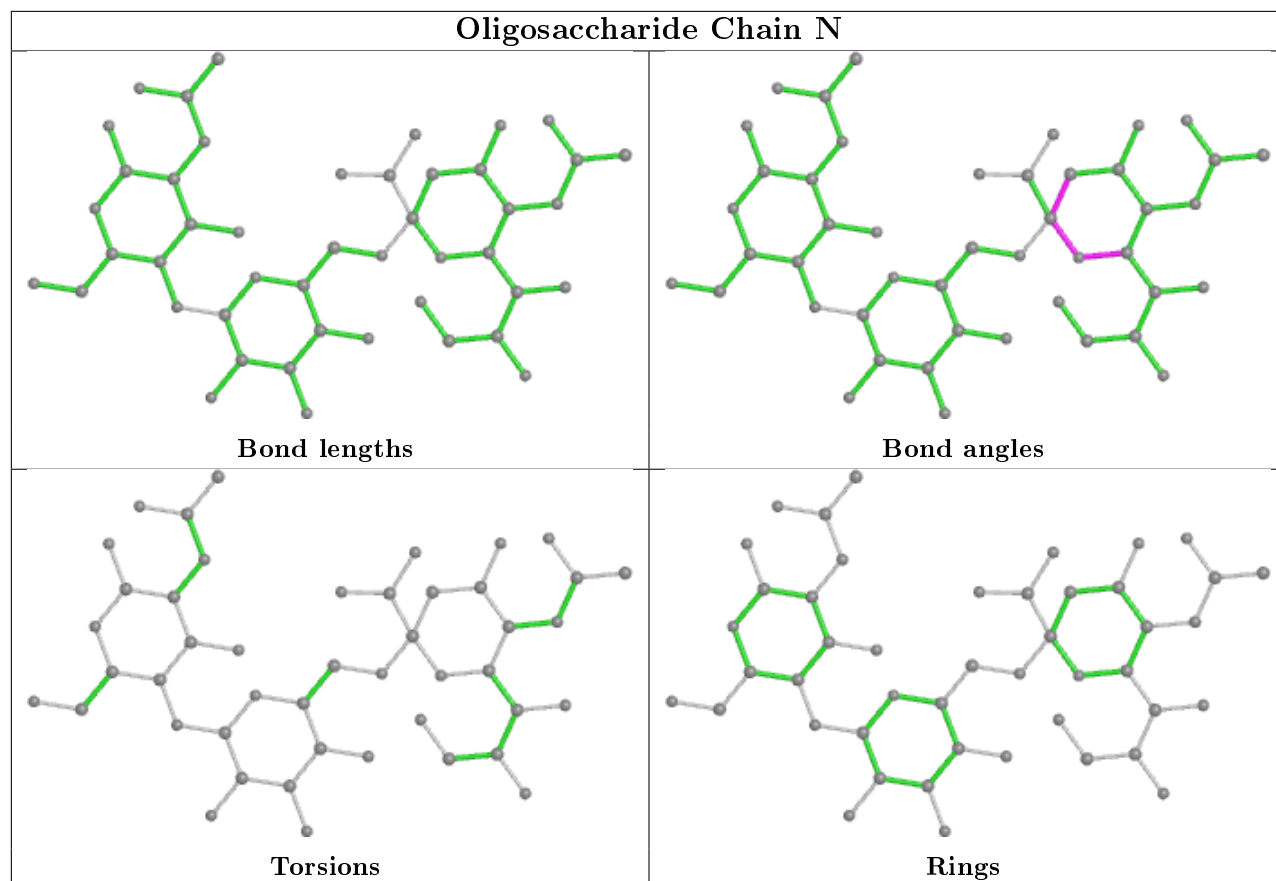
5 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	Q	1	NAG	2	0
3	S	3	SIA	1	0
3	O	3	SIA	2	0
5	Q	2	NAG	1	0
3	N	3	SIA	1	0

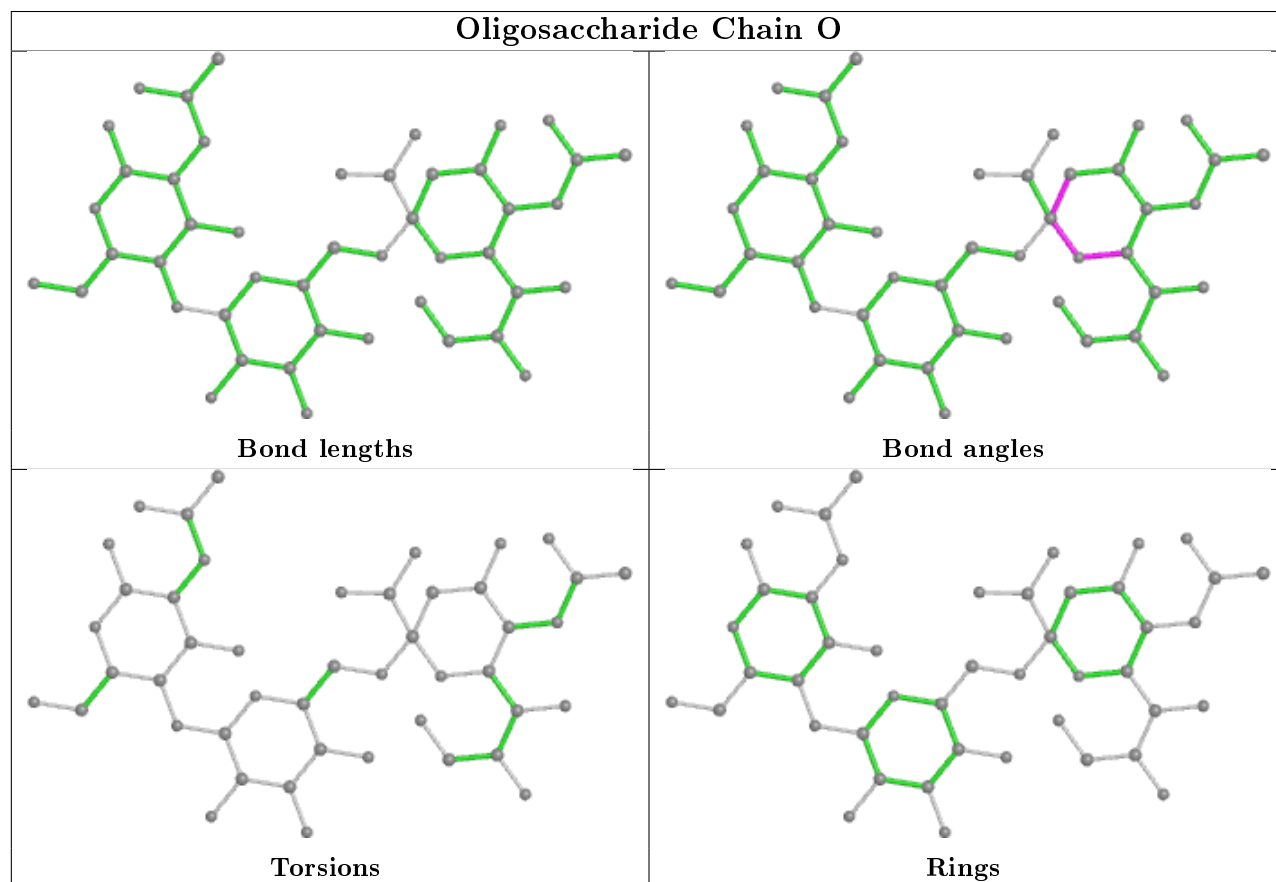
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.



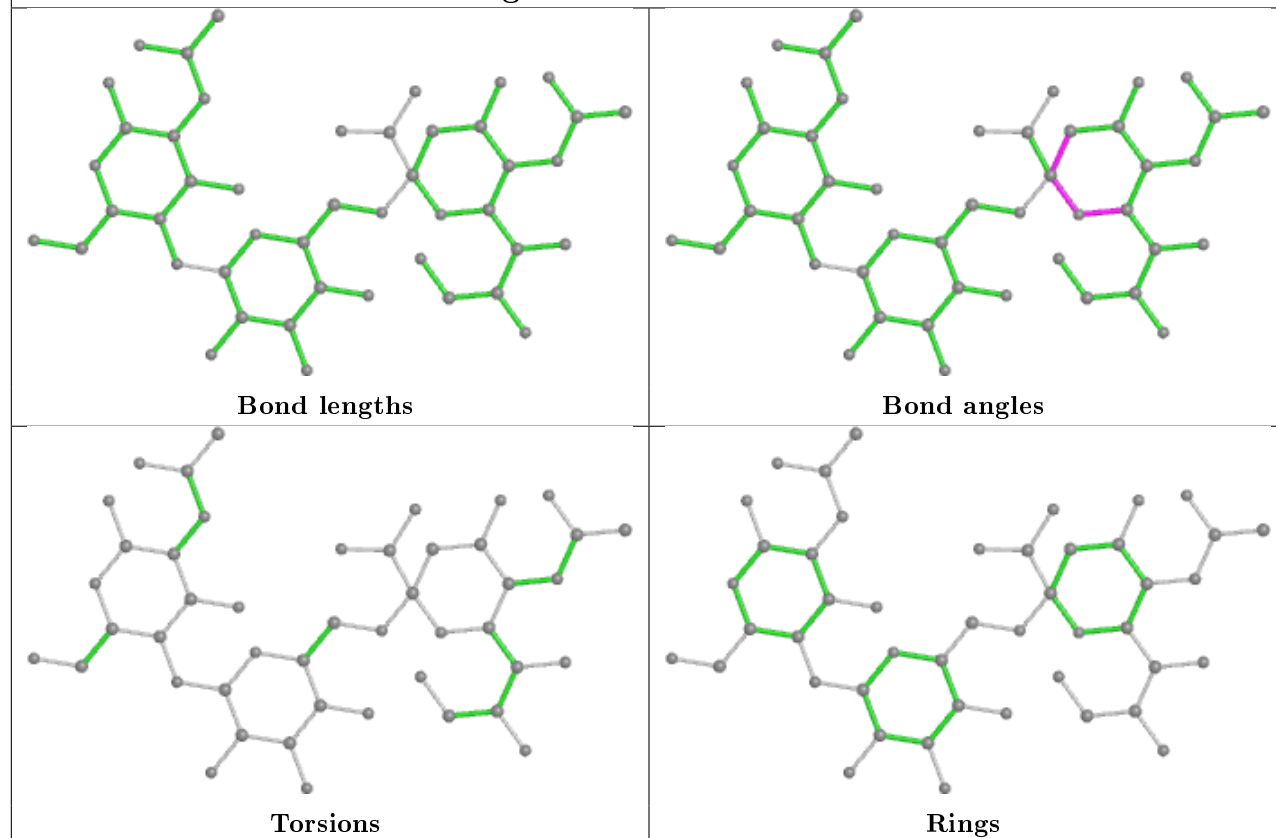
Oligosaccharide Chain N



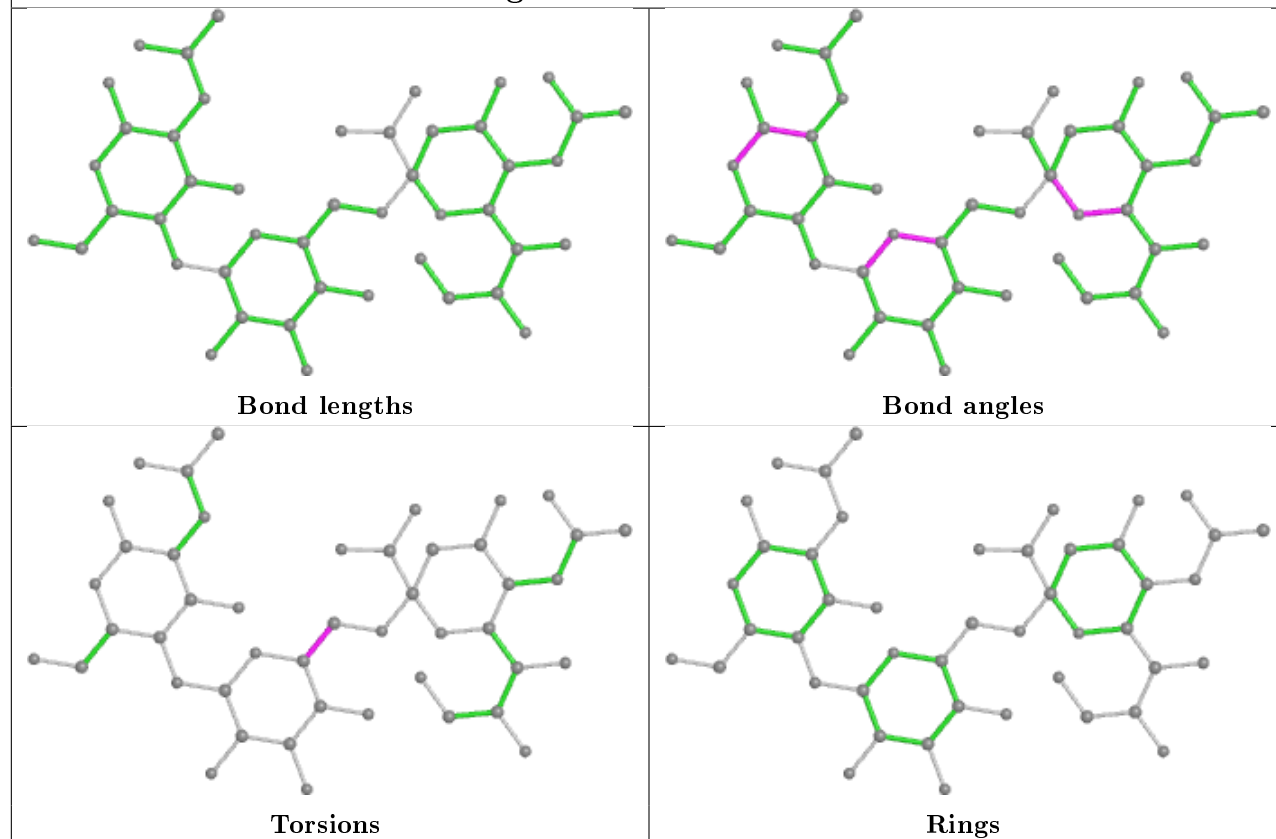
Oligosaccharide Chain O

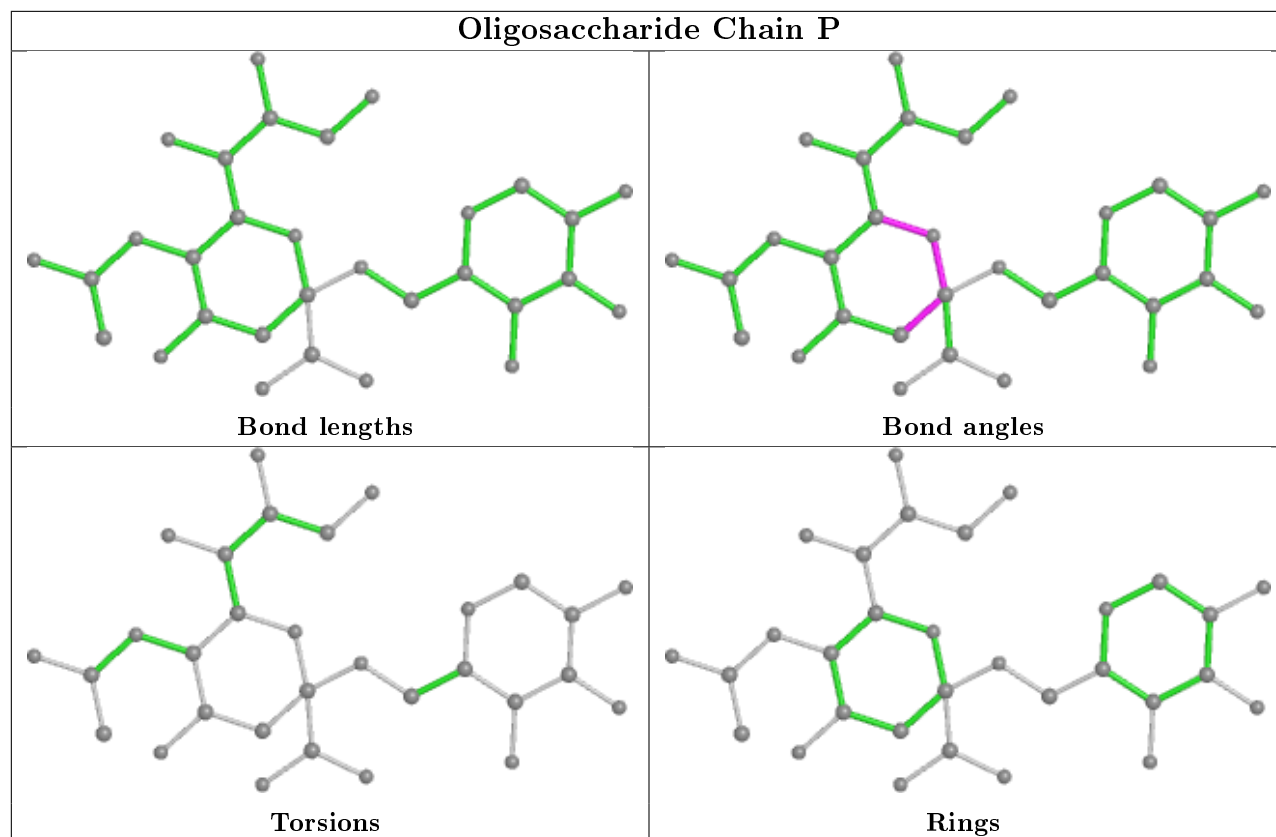


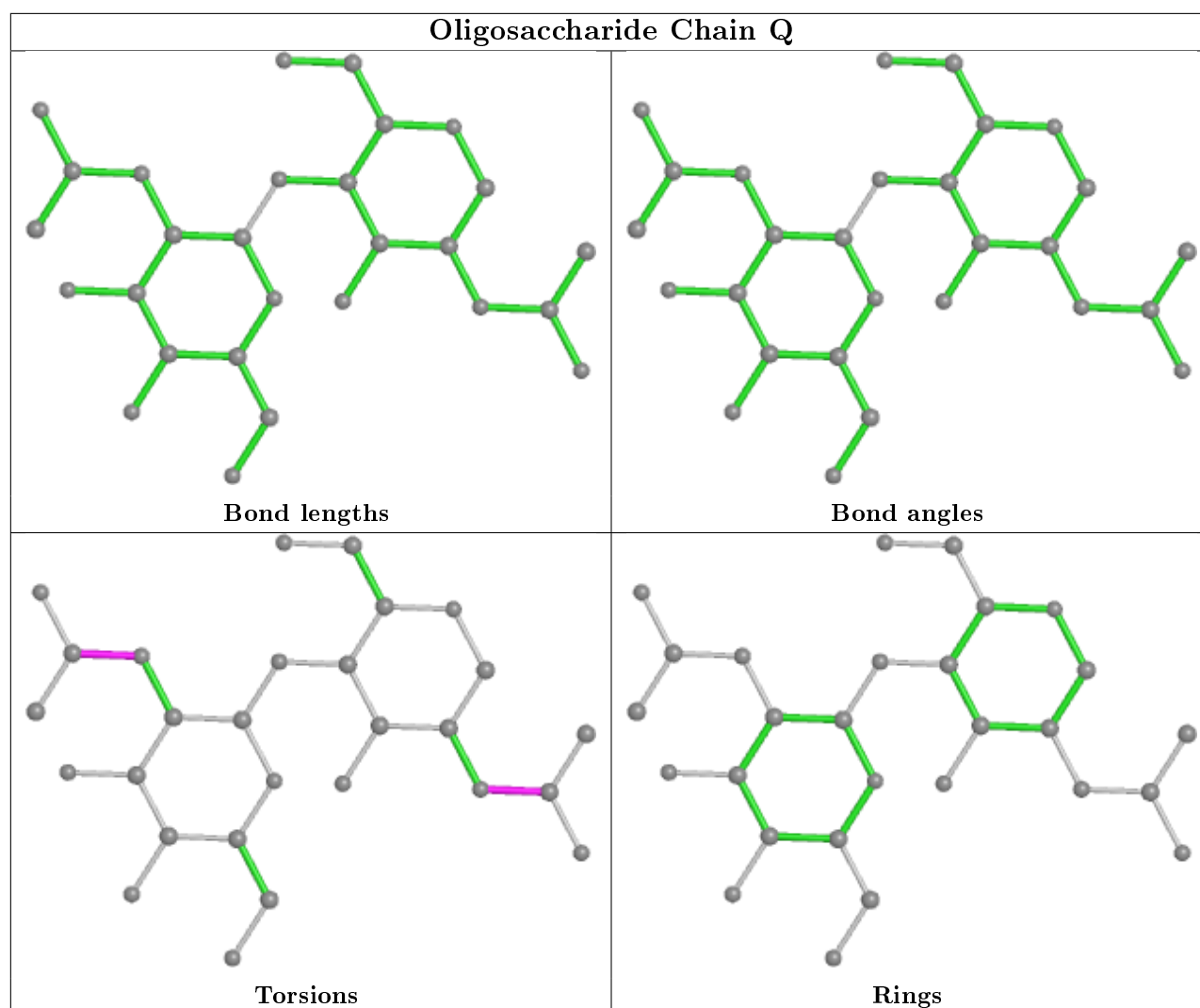
Oligosaccharide Chain R



Oligosaccharide Chain S







5.6 Ligand geometry [i](#)

9 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	$\# Z > 2$
6	NAG	C	333	1	14,14,15	0.51	0	17,19,21	0.79	0
6	NAG	K	641	1	14,14,15	0.50	0	17,19,21	0.95	0
6	NAG	A	334	1	14,14,15	0.51	0	17,19,21	0.72	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	333	1	14,14,15	0.53	0	17,19,21	0.91	1 (5%)
6	NAG	C	334	1	14,14,15	0.52	0	17,19,21	0.72	0
6	NAG	A	335	1	14,14,15	0.50	0	17,19,21	0.82	0
6	NAG	J	561	2	14,14,15	0.51	0	17,19,21	0.65	0
6	NAG	A	336	1	14,14,15	0.49	0	17,19,21	1.00	1 (5%)
6	NAG	E	333	1	14,14,15	0.54	0	17,19,21	0.67	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
6	NAG	C	333	1	-	1/6/23/26	0/1/1/1
6	NAG	K	641	1	-	0/6/23/26	0/1/1/1
6	NAG	A	334	1	-	5/6/23/26	0/1/1/1
6	NAG	A	333	1	-	4/6/23/26	0/1/1/1
6	NAG	C	334	1	-	0/6/23/26	0/1/1/1
6	NAG	A	335	1	-	0/6/23/26	0/1/1/1
6	NAG	J	561	2	-	2/6/23/26	0/1/1/1
6	NAG	A	336	1	-	6/6/23/26	0/1/1/1
6	NAG	E	333	1	-	2/6/23/26	0/1/1/1

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	333	NAG	O5-C5-C6	2.55	111.20	107.20
6	A	336	NAG	C1-O5-C5	2.06	114.99	112.19

There are no chirality outliers.

All (20) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	334	NAG	C3-C2-N2-C7
6	A	334	NAG	C8-C7-N2-C2
6	A	334	NAG	O7-C7-N2-C2
6	A	333	NAG	C8-C7-N2-C2
6	A	333	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
6	J	561	NAG	C8-C7-N2-C2
6	A	336	NAG	C8-C7-N2-C2
6	A	336	NAG	O5-C5-C6-O6
6	J	561	NAG	O7-C7-N2-C2
6	A	336	NAG	O7-C7-N2-C2
6	E	333	NAG	C8-C7-N2-C2
6	A	336	NAG	C1-C2-N2-C7
6	C	333	NAG	O5-C5-C6-O6
6	E	333	NAG	O7-C7-N2-C2
6	A	333	NAG	C4-C5-C6-O6
6	A	336	NAG	C4-C5-C6-O6
6	A	334	NAG	C4-C5-C6-O6
6	A	333	NAG	O5-C5-C6-O6
6	A	334	NAG	O5-C5-C6-O6
6	A	336	NAG	C3-C2-N2-C7

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	C	333	NAG	1	0
6	E	333	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	323/329 (98%)	-0.00	2 (0%) 89 90	30, 44, 60, 79	0
1	C	319/329 (96%)	-0.20	4 (1%) 77 79	30, 41, 60, 85	0
1	E	323/329 (98%)	-0.01	6 (1%) 66 69	35, 49, 67, 88	0
1	G	321/329 (97%)	-0.08	3 (0%) 84 86	28, 41, 61, 83	0
1	I	319/329 (96%)	-0.07	5 (1%) 72 74	32, 46, 62, 81	0
1	K	319/329 (96%)	-0.04	6 (1%) 66 69	28, 42, 60, 85	0
2	B	174/177 (98%)	0.20	9 (5%) 27 29	34, 58, 80, 95	0
2	D	171/177 (96%)	0.59	23 (13%) 3 2	30, 57, 90, 103	0
2	F	170/177 (96%)	0.49	23 (13%) 3 2	32, 55, 94, 105	0
2	H	170/177 (96%)	0.58	21 (12%) 4 3	31, 51, 88, 99	0
2	J	170/177 (96%)	0.04	4 (2%) 59 62	31, 49, 69, 81	0
2	L	170/177 (96%)	-0.07	0 100 100	30, 46, 62, 69	0
All	All	2949/3036 (97%)	0.06	106 (3%) 42 46	28, 46, 79, 105	0

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	168	LEU	5.7
2	D	147	THR	4.7
2	D	16	GLY	4.6
2	F	29	GLU	4.5
2	D	168	LEU	4.5
2	B	33	GLY	4.3
1	K	94	ASN	4.2
2	F	160	PRO	4.1
2	D	144	CYS	4.0
2	D	162	TYR	4.0
2	D	160	PRO	4.0

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Mol	Chain	Res	Type	RSRZ
2	F	161	LYS	4.0
2	H	143	LYS	3.9
2	D	157	TYR	3.9
2	J	133	ILE	3.9
1	C	12	THR	3.8
2	F	164	GLU	3.8
2	H	144	CYS	3.7
2	F	158	ASP	3.7
2	D	138	PHE	3.6
2	D	140	PHE	3.5
2	D	142	HIS	3.4
2	D	141	TYR	3.4
2	D	143	LYS	3.4
2	D	156	THR	3.3
2	D	129	ASN	3.3
1	I	264	ALA	3.2
2	D	128	ASN	3.2
2	H	29	GLU	3.2
2	H	141	TYR	3.2
2	J	143	LYS	3.2
2	F	147	THR	3.1
2	B	173	ILE	3.1
1	G	13	LEU	3.0
2	H	164	GLU	3.0
1	C	264	ALA	3.0
2	D	19	ASP	2.9
2	H	140	PHE	2.9
1	C	90(A)	PRO	2.9
2	H	167	LYS	2.9
2	F	38	LEU	2.9
2	F	166	ALA	2.9
2	H	65	ALA	2.9
1	I	54	LEU	2.9
2	D	27	GLN	2.8
2	H	23	GLY	2.8
2	F	156	THR	2.8
1	A	325	SER	2.8
2	H	159	TYR	2.7
1	G	12	THR	2.7
2	D	159	TYR	2.7
2	F	30	GLN	2.7
2	H	18	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	166	ALA	2.6
2	F	162	TYR	2.6
1	I	55	ARG	2.6
2	H	166	ALA	2.5
2	H	152	VAL	2.5
2	B	174	ASP	2.5
2	B	168	LEU	2.5
2	J	134	GLY	2.5
2	B	138	PHE	2.5
2	D	154	ASN	2.5
2	D	137	CYS	2.5
2	F	23	GLY	2.5
2	J	166	ALA	2.5
2	H	168	LEU	2.4
1	C	78	SER	2.4
1	K	93	ASP	2.4
1	I	208	ARG	2.4
1	I	290	THR	2.4
2	H	42	GLN	2.4
2	F	163	SER	2.4
2	F	140	PHE	2.4
1	E	13	LEU	2.4
1	E	161	TYR	2.4
2	H	160	PRO	2.4
1	E	198	ALA	2.3
2	D	161	LYS	2.3
2	H	33	GLY	2.3
2	H	145	ASP	2.3
1	E	81	THR	2.3
2	H	157	TYR	2.3
2	D	152	VAL	2.3
2	F	28	ASN	2.3
1	K	239	PRO	2.3
2	B	34	TYR	2.3
2	F	143	LYS	2.2
2	F	149	MET	2.2
2	F	148	CYS	2.2
2	F	60	ASN	2.2
2	F	144	CYS	2.2
1	K	275	HIS	2.2
2	B	25	HIS	2.2
2	H	156	THR	2.2

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Mol	Chain	Res	Type	RSRZ
2	F	31	GLY	2.2
1	E	11	ASP	2.2
1	G	81	THR	2.2
1	K	9	PRO	2.2
2	F	146	ASN	2.1
2	H	139	GLU	2.1
2	B	140	PHE	2.1
2	B	27	GLN	2.1
1	E	158	GLY	2.0
1	A	290	THR	2.0
1	K	208	ARG	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.

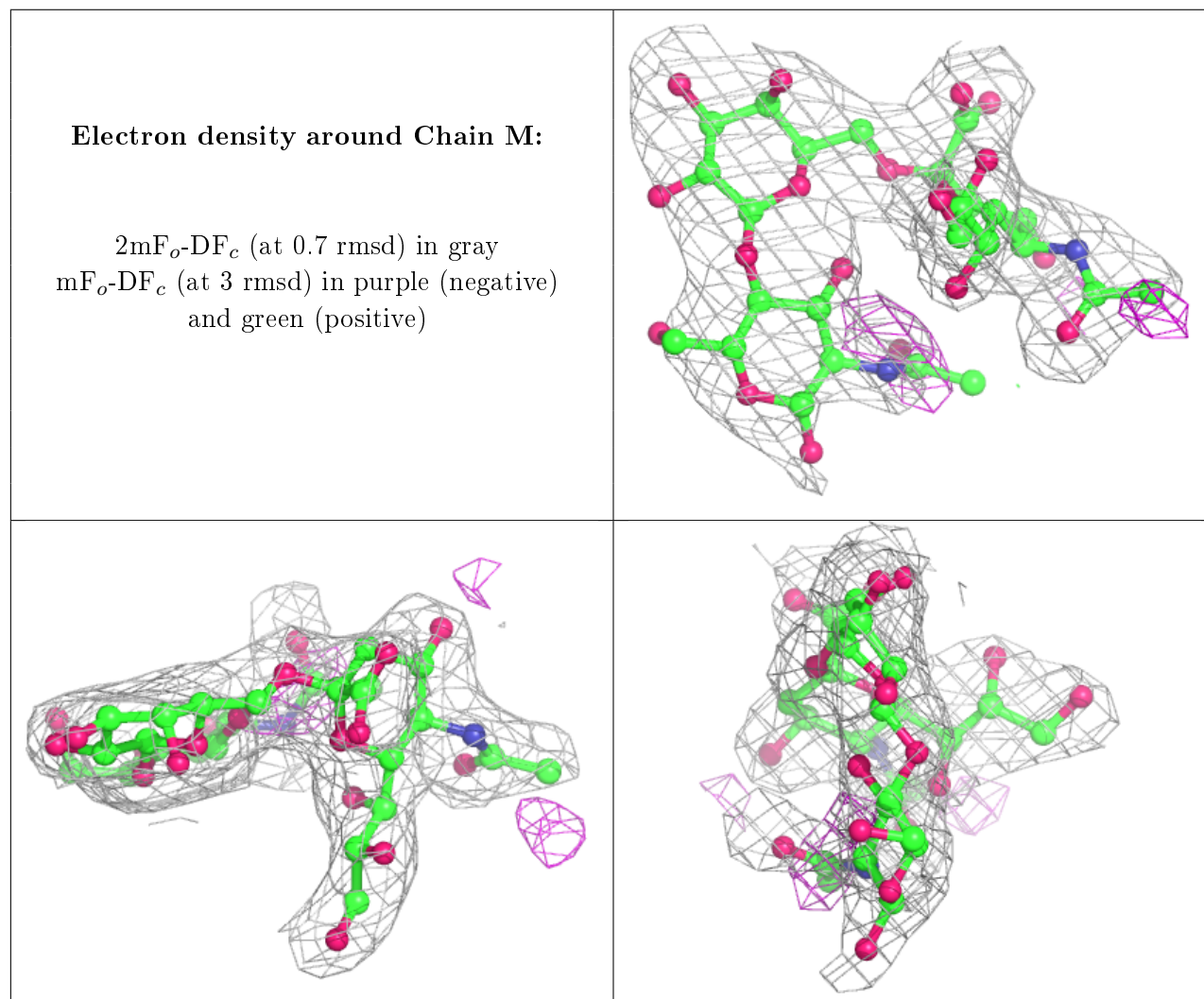
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	M	1	15/15	0.76	0.30	57,69,76,78	0
3	GAL	S	2	11/12	0.82	0.28	64,73,82,85	0
5	NAG	Q	2	14/15	0.85	0.22	48,57,64,68	0
3	NAG	S	1	15/15	0.86	0.28	58,77,81,85	0
3	NAG	R	1	15/15	0.90	0.15	45,54,65,66	0
3	NAG	O	1	15/15	0.90	0.19	55,63,66,68	0
3	GAL	M	2	11/12	0.91	0.13	49,59,66,73	0
5	NAG	Q	1	14/15	0.92	0.15	38,50,56,56	0
4	GAL	P	1	11/12	0.92	0.23	55,67,74,76	0
3	SIA	S	3	20/21	0.93	0.23	47,59,66,68	0
3	SIA	O	3	20/21	0.93	0.18	49,55,62,66	0
3	NAG	N	1	15/15	0.93	0.16	47,53,64,65	0
3	GAL	R	2	11/12	0.93	0.13	48,51,56,58	0
3	GAL	N	2	11/12	0.94	0.11	44,51,54,58	0
3	GAL	O	2	11/12	0.94	0.09	51,58,61,61	0
3	SIA	R	3	20/21	0.96	0.12	31,46,52,52	0
4	SIA	P	2	20/21	0.96	0.12	35,45,51,52	0

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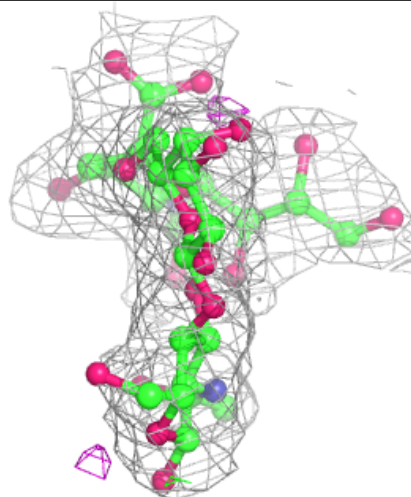
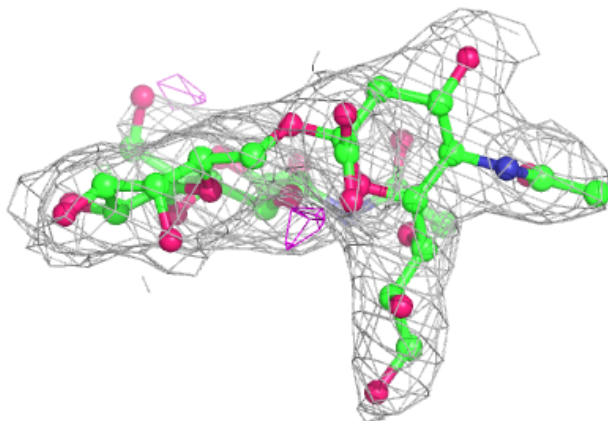
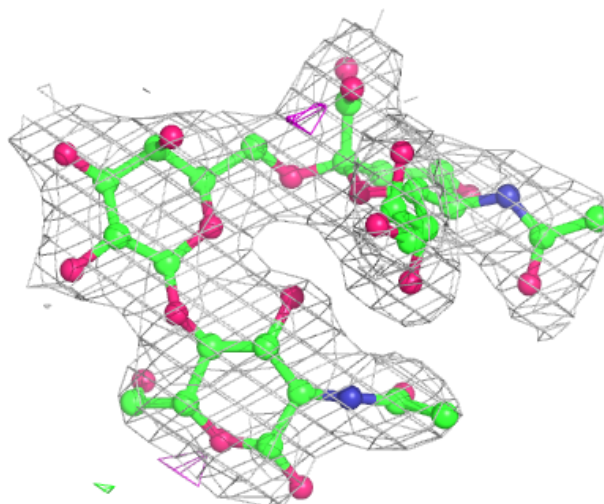
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	SIA	N	3	20/21	0.96	0.12	33,39,43,47	0
3	SIA	M	3	20/21	0.96	0.16	36,43,50,52	0

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.



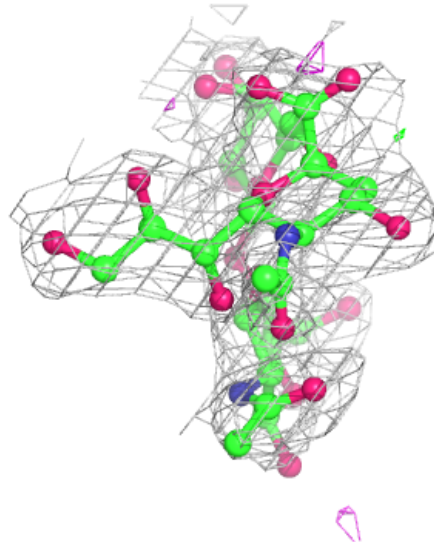
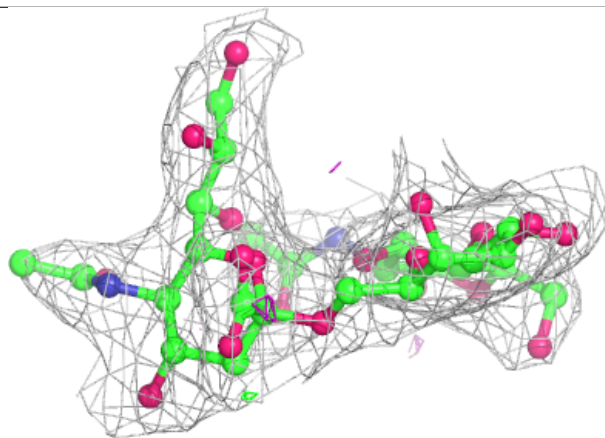
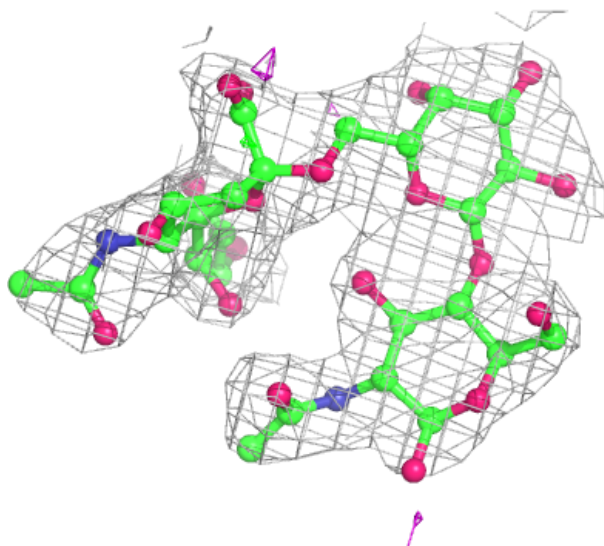
Electron density around Chain N:

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



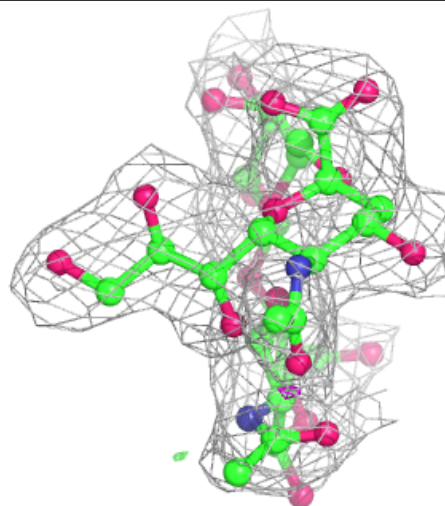
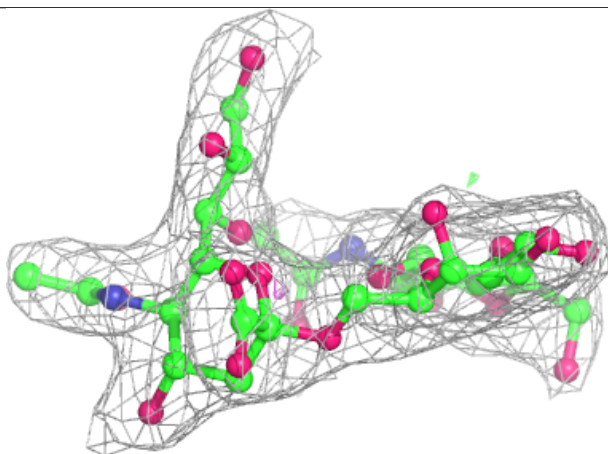
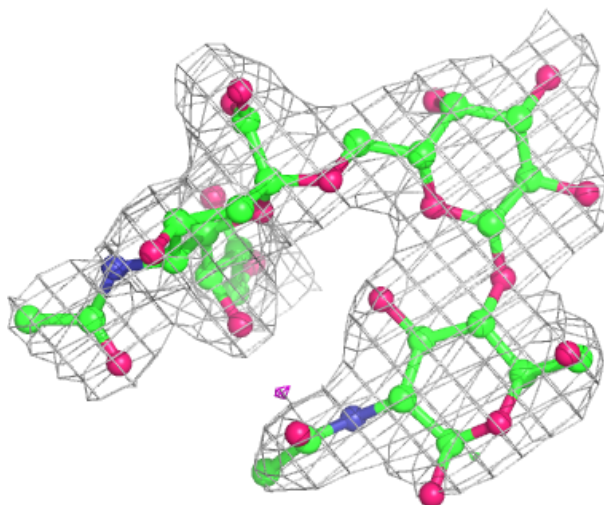
Electron density around Chain O:

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



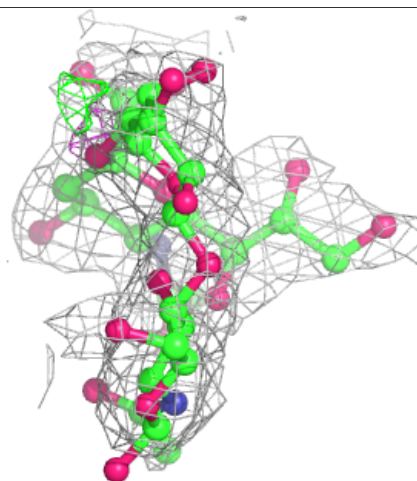
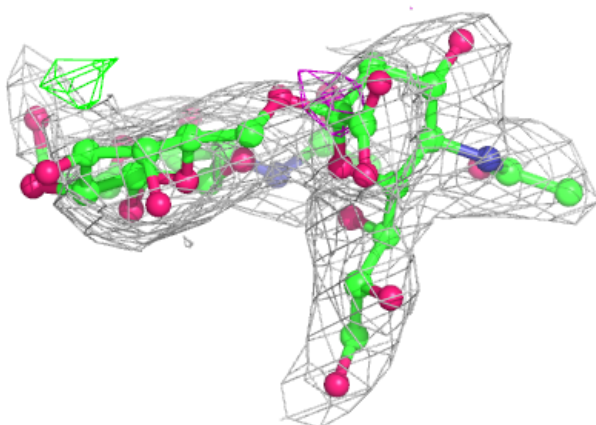
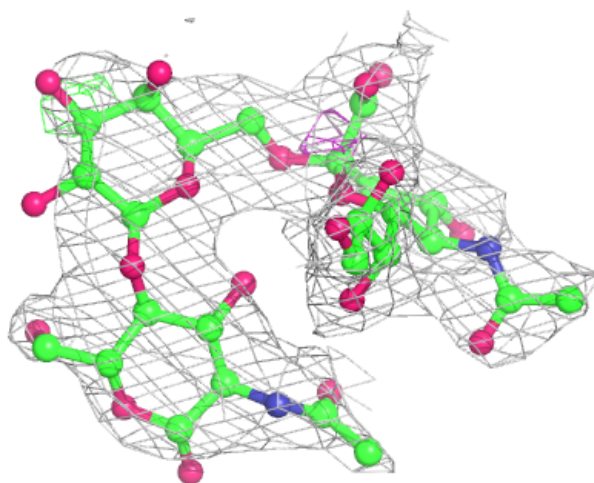
Electron density around Chain R:

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



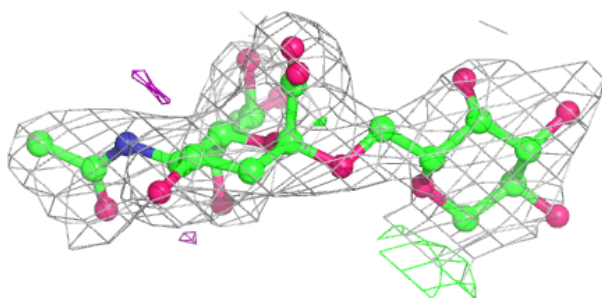
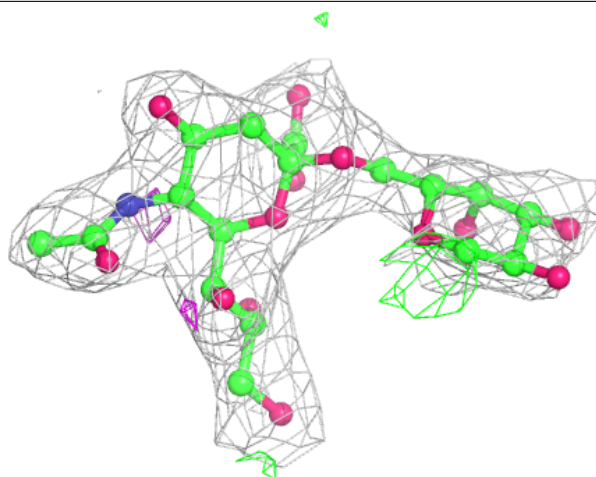
Electron density around Chain S:

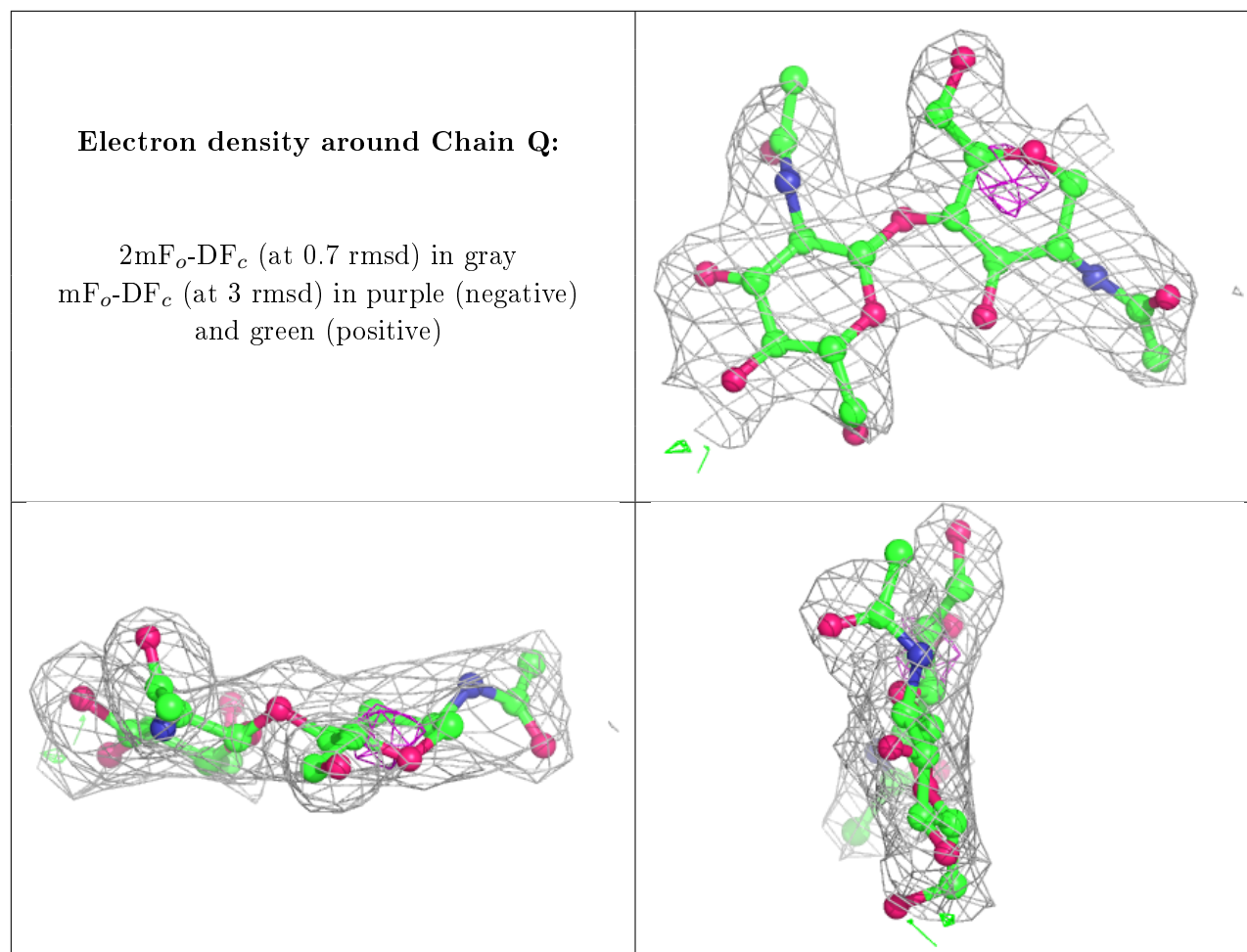
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain P:

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





6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
6	NAG	A	334	14/15	0.70	0.20	65,74,89,90	0
6	NAG	A	333	14/15	0.75	0.43	70,84,86,87	0
6	NAG	K	641	14/15	0.77	0.28	71,81,89,91	0
6	NAG	A	336	14/15	0.80	0.27	71,76,80,83	0
6	NAG	J	561	14/15	0.81	0.18	76,78,83,84	0
6	NAG	C	334	14/15	0.86	0.16	62,76,83,84	0
6	NAG	A	335	14/15	0.90	0.12	48,52,58,61	0
6	NAG	C	333	14/15	0.93	0.14	47,56,67,69	0
6	NAG	E	333	14/15	0.93	0.15	48,56,60,60	0

6.5 Other polymers

There are no such residues in this entry.