



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 05:44 AM UTC

PDB ID : 5THC / pdb_00005thc
Title : Crystal structure of H10 hemagglutinin mutant (T193D-Q226L-G228S) from Jiangxi-Donghu (2013) H10N8 influenza virus in complex with 6'-SLNLN
Authors : Tzarum, N.; Wilson, I.A.
Deposited on : 2016-09-29
Resolution : 2.79 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

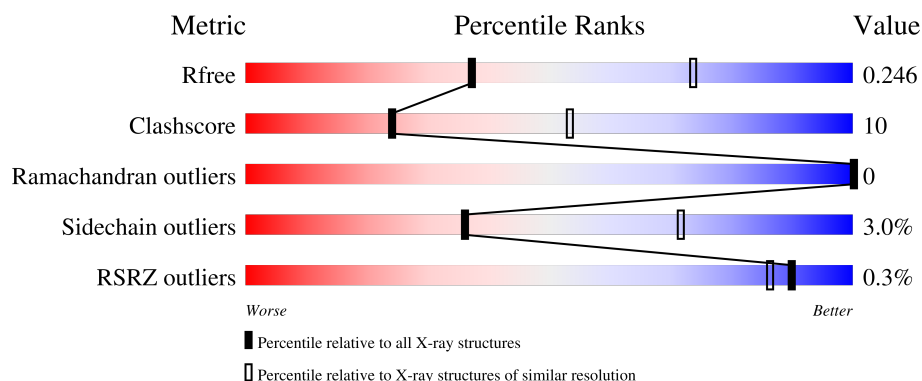
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	323	 76% 22% ..
1	C	323	 78% 19% ..
1	E	323	 72% 26% ..
2	B	180	 % 68% 26% • 5%
2	D	180	 73% 22% 5%

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Mol	Chain	Length	Quality of chain
2	F	180	% 71%21%6%
3	G	2	100%
4	H	3	33%33%33%
5	I	4	25%50%25%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 11801 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 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	318	Total	C	N	O	S	0	0	0
			2437	1508	448	464	17			
1	C	318	Total	C	N	O	S	0	0	0
			2437	1508	448	464	17			
1	E	318	Total	C	N	O	S	0	0	0
			2437	1508	448	464	17			

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	193	THR	ASP	engineered mutation	UNP A0A0J9X252
A	226	LEU	GLN	engineered mutation	UNP A0A0J9X252
A	228	SER	GLY	engineered mutation	UNP A0A0J9X252
C	193	THR	ASP	engineered mutation	UNP A0A0J9X252
C	226	LEU	GLN	engineered mutation	UNP A0A0J9X252
C	228	SER	GLY	engineered mutation	UNP A0A0J9X252
E	193	THR	ASP	engineered mutation	UNP A0A0J9X252
E	226	LEU	GLN	engineered mutation	UNP A0A0J9X252
E	228	SER	GLY	engineered mutation	UNP A0A0J9X252

- Molecule 2 is a protein called Hemagglutinin HA2 chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	171	Total	C	N	O	S	0	0	0
			1382	854	239	281	8			
2	D	171	Total	C	N	O	S	0	0	0
			1382	854	239	281	8			
2	F	169	Total	C	N	O	S	0	0	0
			1371	848	237	278	8			

There are 21 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	175	SER	-	expression tag	UNP A0A0J9X253
B	176	GLY	-	expression tag	UNP A0A0J9X253
B	177	ARG	-	expression tag	UNP A0A0J9X253
B	178	LEU	-	expression tag	UNP A0A0J9X253
B	179	VAL	-	expression tag	UNP A0A0J9X253
B	180	PRO	-	expression tag	UNP A0A0J9X253
B	181	ARG	-	expression tag	UNP A0A0J9X253
D	175	SER	-	expression tag	UNP A0A0J9X253
D	176	GLY	-	expression tag	UNP A0A0J9X253
D	177	ARG	-	expression tag	UNP A0A0J9X253
D	178	LEU	-	expression tag	UNP A0A0J9X253
D	179	VAL	-	expression tag	UNP A0A0J9X253
D	180	PRO	-	expression tag	UNP A0A0J9X253
D	181	ARG	-	expression tag	UNP A0A0J9X253
F	175	SER	-	expression tag	UNP A0A0J9X253
F	176	GLY	-	expression tag	UNP A0A0J9X253
F	177	ARG	-	expression tag	UNP A0A0J9X253
F	178	LEU	-	expression tag	UNP A0A0J9X253
F	179	VAL	-	expression tag	UNP A0A0J9X253
F	180	PRO	-	expression tag	UNP A0A0J9X253
F	181	ARG	-	expression tag	UNP A0A0J9X253

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



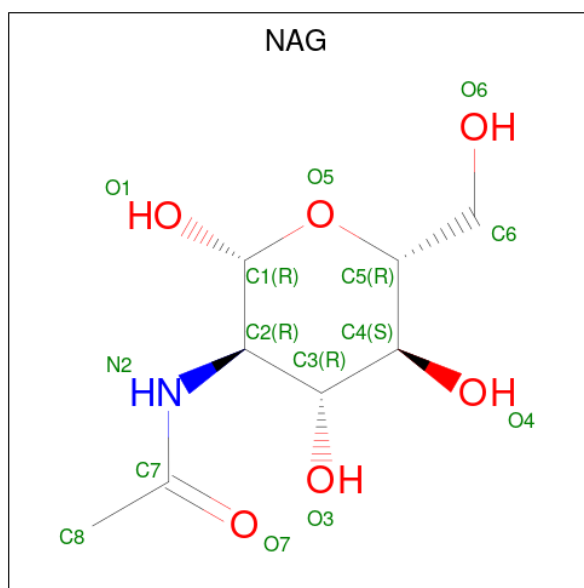
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	3	Total	C	N	O	0	0	0
			46	25	2	19			

- Molecule 5 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.



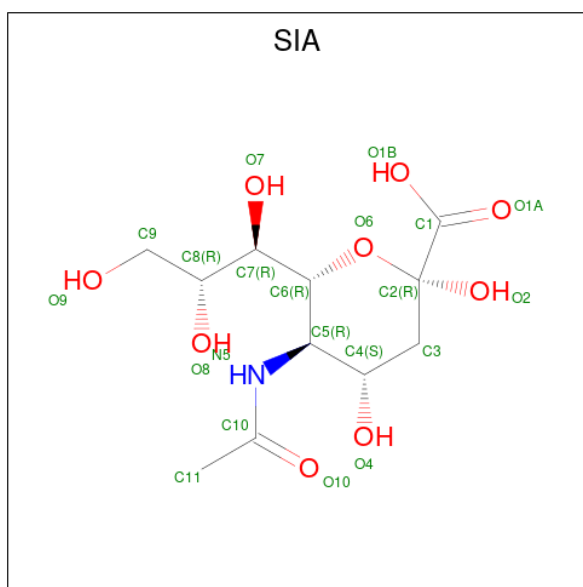
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
5	I	4	Total	C	N	O	0	0	0
			50	28	2	20			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

- Molecule 7 is N-acetyl-alpha-neuraminic acid (CCD ID: SIA) (formula: C₁₁H₁₉NO₉).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
7	C	1	Total	C	N	O	0	0
			21	11	1	9		

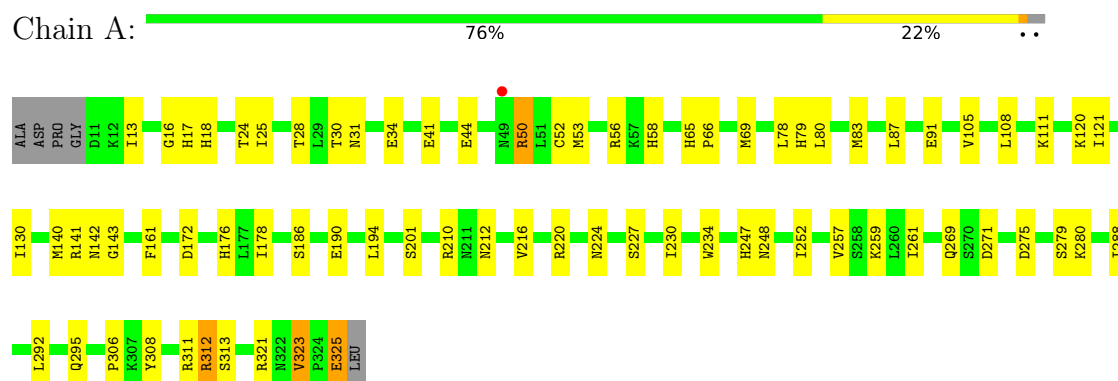
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	32	Total	O	0	0
			32	32		
8	B	11	Total	O	0	0
			11	11		
8	C	47	Total	O	0	0
			47	47		
8	D	16	Total	O	0	0
			16	16		
8	E	43	Total	O	0	0
			43	43		
8	F	19	Total	O	0	0
			19	19		

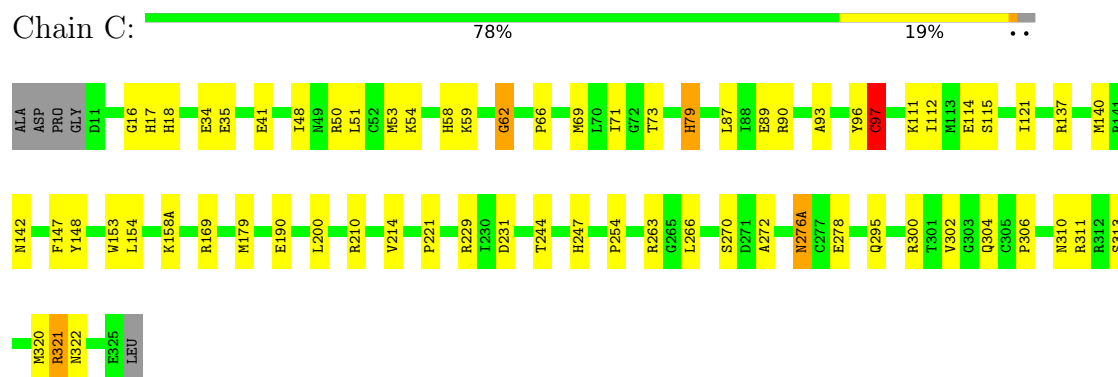
3 Residue-property plots [i](#)

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.

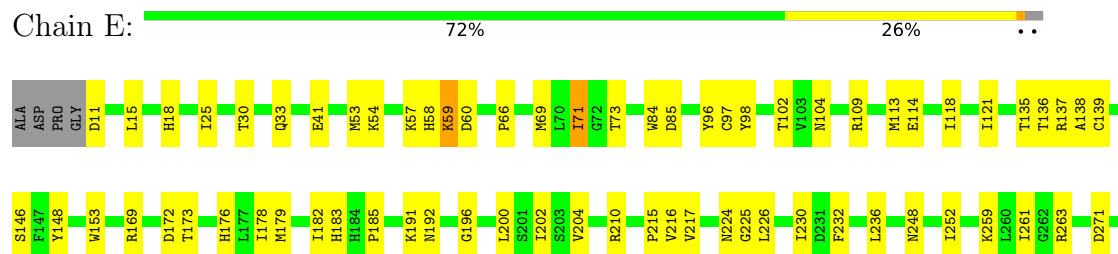
- Molecule 1: Hemagglutinin HA1 chain



- Molecule 1: Hemagglutinin HA1 chain



- Molecule 1: Hemagglutinin HA1 chain





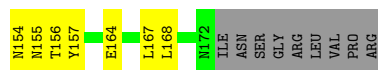
• Molecule 2: Hemagglutinin HA2 chain

Chain B: 68% 26% 5%



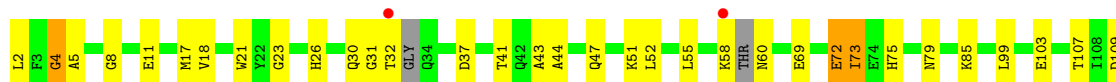
• Molecule 2: Hemagglutinin HA2 chain

Chain D: 73% 22% 5%



• Molecule 2: Hemagglutinin HA2 chain

Chain F: 71% 21% 6%



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

Chain G: 100%



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

Chain H: 33% 33% 33%



- Molecule 5: 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: 25% 50% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	64.19Å 244.45Å 70.76Å 90.00° 111.67° 90.00°	Depositor
Resolution (Å)	37.78 – 2.79 37.78 – 2.79	Depositor EDS
% Data completeness (in resolution range)	96.9 (37.78-2.79) 93.9 (37.78-2.79)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.19	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.80 (at 2.81Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.191 , 0.248 0.192 , 0.246	Depositor DCC
R_{free} test set	2461 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	38.1	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 19.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	11801	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% 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: GAL, BMA, NAG, SIA, MAN

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	0.85	0/2486	1.13	12/3369 (0.4%)
1	C	0.89	1/2486 (0.0%)	1.07	9/3369 (0.3%)
1	E	0.82	0/2486	1.02	1/3369 (0.0%)
2	B	0.71	0/1407	1.00	1/1899 (0.1%)
2	D	0.77	0/1407	0.87	1/1899 (0.1%)
2	F	0.70	0/1394	0.98	4/1878 (0.2%)
All	All	0.81	1/11666 (0.0%)	1.03	28/15783 (0.2%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	93	ALA	CA-CB	-5.34	1.44	1.53

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	276(A)	ASN	N-CA-C	8.39	120.50	111.36
2	F	5	ALA	N-CA-C	7.74	119.35	111.07
1	A	323	VAL	CA-C-N	7.63	127.56	119.85
1	A	323	VAL	C-N-CA	7.63	127.56	119.85
1	A	220	ARG	CA-C-N	-7.29	112.47	119.76
1	A	220	ARG	C-N-CA	-7.29	112.47	119.76
1	A	247	HIS	N-CA-C	7.27	120.40	108.99
1	A	105	VAL	N-CA-C	7.10	117.89	110.72
1	C	97	CYS	CB-CA-C	6.95	122.46	110.79
1	A	91	GLU	N-CA-C	6.55	118.08	111.07
1	A	143	GLY	N-CA-C	-6.37	106.45	115.43
2	F	31	GLY	N-CA-C	6.33	121.62	110.80
1	C	62	GLY	CA-C-N	6.33	129.39	120.28
1	C	62	GLY	C-N-CA	6.33	129.39	120.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	4	GLY	N-CA-C	6.08	121.78	113.99
1	A	194	LEU	N-CA-C	6.03	117.93	111.36
1	A	130	ILE	N-CA-C	6.02	116.97	108.36
1	A	220	ARG	N-CA-C	5.79	121.38	113.16
1	C	79	HIS	N-CA-C	5.79	119.94	112.41
1	C	97	CYS	N-CA-CB	-5.68	101.72	110.13
1	C	270	SER	N-CA-C	5.53	117.67	108.99
2	B	38	TYR	N-CA-C	5.41	116.98	111.14
1	E	285	GLY	N-CA-C	-5.38	103.76	112.31
2	D	23	GLY	N-CA-C	5.34	120.87	110.73
2	F	18	VAL	N-CA-C	-5.30	107.13	112.17
1	C	96	TYR	CA-C-N	5.19	129.37	120.71
1	C	96	TYR	C-N-CA	5.19	129.37	120.71
1	A	261	ILE	N-CA-C	5.13	115.29	108.11

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	2437	0	2394	53	0
1	C	2437	0	2393	55	0
1	E	2437	0	2394	59	0
2	B	1382	0	1279	41	0
2	D	1382	0	1278	28	0
2	F	1371	0	1267	27	0
3	G	28	0	25	0	0
4	H	46	0	40	6	0
5	I	50	0	43	2	0
6	A	14	0	13	0	0
6	C	14	0	13	0	0
6	D	14	0	13	0	0
7	C	21	0	17	7	0
8	A	32	0	0	5	0
8	B	11	0	0	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
8	C	47	0	0	6	0
8	D	16	0	0	1	0
8	E	43	0	0	5	0
8	F	19	0	0	5	0
All	All	11801	0	11169	227	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 (227) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:VAL:O	1:C:210:ARG:NH1	2.03	0.92
2:B:144:CYS:HB3	2:B:149:MET:HE2	1.56	0.88
2:B:105:GLN:NE2	8:B:201:HOH:O	2.07	0.88
2:F:79:ASN:OD1	8:F:201:HOH:O	1.91	0.87
2:F:44:ALA:HA	2:F:110:MET:HE1	1.59	0.85
1:C:73:THR:HG23	1:C:97:CYS:HB3	1.58	0.84
1:E:104:ASN:ND2	8:E:501:HOH:O	2.11	0.82
1:C:231:ASP:OD2	1:E:210:ARG:NH1	2.14	0.80
1:C:41:GLU:OE2	1:C:313:SER:OG	1.99	0.78
2:B:134:GLY:HA2	2:D:124:LYS:HD3	1.65	0.77
1:A:50:ARG:NE	1:A:275:ASP:OD2	2.19	0.76
1:A:78:LEU:O	8:A:501:HOH:O	2.04	0.75
2:D:22:TYR:HD1	2:D:115:MET:HE2	1.52	0.75
1:C:66:PRO:HA	1:C:69:MET:HE3	1.68	0.74
2:B:44:ALA:HB2	2:B:114:GLU:HG3	1.69	0.74
1:E:169:ARG:HD3	5:I:1:NAG:H62	1.72	0.72
1:A:227:SER:O	1:C:210:ARG:NH2	2.25	0.70
2:F:120:GLU:OE1	2:F:123:ARG:NH2	2.25	0.70
1:C:158(A):LYS:NZ	8:C:503:HOH:O	2.15	0.70
1:E:200:LEU:HA	1:E:248:ASN:HD22	1.56	0.70
2:B:22:TYR:H	2:B:41:THR:HG22	1.58	0.69
1:E:41:GLU:OE1	1:E:313:SER:OG	2.07	0.68
1:A:56:ARG:NH2	8:A:503:HOH:O	2.19	0.68
1:C:190:GLU:OE2	7:C:401:SIA:O9	2.11	0.67
1:E:11:ASP:N	8:E:503:HOH:O	2.27	0.67
1:C:300:ARG:NH1	2:D:69:GLU:OE1	2.29	0.66
1:C:114:GLU:OE2	8:C:502:HOH:O	2.13	0.66
1:A:325:GLU:OE1	2:B:15:GLU:HG3	1.96	0.65
1:E:66:PRO:HA	1:E:69:MET:HE2	1.79	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:53:MET:HE2	1:A:58:HIS:HB3	1.78	0.65
2:F:109:ASP:OD2	8:F:202:HOH:O	2.14	0.65
2:F:43:ALA:O	2:F:47:GLN:NE2	2.30	0.64
1:C:112:ILE:O	1:C:115:SER:OG	2.13	0.64
1:A:17:HIS:HB3	2:B:115:MET:CE	2.27	0.64
2:B:151:SER:HA	2:B:156:THR:HB	1.79	0.64
1:C:59:LYS:HB3	1:C:87:LEU:HD12	1.79	0.64
2:F:58:LYS:O	2:F:60:ASN:N	2.31	0.63
1:C:304:GLN:OE1	2:D:60:ASN:HB3	1.99	0.63
2:F:4:GLY:O	2:F:8:GLY:HA3	1.99	0.63
1:E:153:TRP:CE3	4:H:3:SIA:H112	2.34	0.62
2:D:22:TYR:CD1	2:D:115:MET:HE2	2.33	0.60
1:A:172:ASP:OD2	1:A:259:LYS:NZ	2.22	0.59
2:B:37:ASP:O	2:B:41:THR:HG23	2.03	0.59
1:A:311:ARG:NE	2:B:97:GLU:OE1	2.28	0.59
1:C:153:TRP:CZ3	7:C:401:SIA:H112	2.38	0.59
1:A:41:GLU:OE2	1:A:313:SER:OG	2.15	0.59
2:D:154:ASN:HB3	2:D:156:THR:HG23	1.85	0.59
1:E:98:TYR:CE2	1:E:226:LEU:HD13	2.39	0.58
1:C:190:GLU:CD	7:C:401:SIA:HO9	2.10	0.58
1:E:57:LYS:N	1:E:85:ASP:OD1	2.36	0.58
2:D:51:LYS:NZ	2:D:107:THR:OG1	2.36	0.57
1:C:69:MET:HG2	1:C:79:HIS:O	2.04	0.57
1:A:323:VAL:HG23	2:B:13:GLY:H	1.68	0.57
1:E:169:ARG:CD	5:I:1:NAG:H62	2.35	0.57
1:C:114:GLU:O	1:C:263:ARG:HD3	2.05	0.57
1:A:111:LYS:HG2	2:F:75:HIS:CE1	2.40	0.56
1:A:295:GLN:HG2	1:A:306:PRO:HG2	1.87	0.56
2:B:117:ASN:HD21	2:F:2:LEU:HD22	1.71	0.56
1:E:25:ILE:HG23	1:E:33:GLN:HA	1.87	0.56
1:E:97:CYS:O	1:E:224:ASN:ND2	2.38	0.56
1:A:186:SER:OG	1:A:190:GLU:OE1	2.18	0.56
1:E:319:GLY:HA2	2:F:21:TRP:CH2	2.41	0.56
1:A:69:MET:HE1	1:A:87:LEU:HD21	1.88	0.55
1:A:80:LEU:O	1:A:120:LYS:NZ	2.39	0.55
1:C:54:LYS:HD2	1:C:276(A):ASN:O	2.06	0.55
1:C:71:ILE:O	1:C:148:TYR:HB3	2.06	0.55
2:B:51:LYS:HG3	1:E:30:THR:HG22	1.88	0.55
2:B:157:TYR:HE1	2:B:159:HIS:HA	1.70	0.55
2:B:157:TYR:CE1	2:B:159:HIS:HA	2.41	0.55
2:D:121:ARG:HD3	2:D:155:ASN:OD1	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:30:GLN:NE2	2:B:145:ASP:OD1	2.39	0.54
1:C:66:PRO:HA	1:C:69:MET:CE	2.37	0.54
1:C:48:ILE:HG22	1:C:50:ARG:H	1.73	0.54
1:E:172:ASP:OD1	1:E:173:THR:N	2.40	0.54
1:C:62:GLY:O	8:C:504:HOH:O	2.18	0.53
1:C:66:PRO:HD3	1:C:89:GLU:HG3	1.91	0.53
1:A:161:PHE:HB3	1:A:248:ASN:O	2.09	0.53
1:C:53:MET:HE2	1:C:58:HIS:CB	2.38	0.53
1:A:121:ILE:HD13	1:A:176:HIS:CE1	2.44	0.53
1:E:286:GLY:C	8:E:504:HOH:O	2.52	0.53
1:A:69:MET:HG2	1:A:79:HIS:O	2.08	0.52
2:D:164:GLU:O	2:D:168:LEU:HG	2.10	0.52
1:C:17:HIS:HB3	2:D:115:MET:HE3	1.92	0.52
2:D:83:TRP:CH2	2:F:85:LYS:HG2	2.44	0.52
1:C:295:GLN:HG2	1:C:306:PRO:HG2	1.92	0.52
1:C:169:ARG:HG3	1:C:169:ARG:HH11	1.73	0.51
1:A:323:VAL:HG23	2:B:13:GLY:N	2.25	0.51
1:A:66:PRO:HA	1:A:69:MET:HE3	1.93	0.51
2:B:75:HIS:CE1	1:C:111:LYS:HG3	2.45	0.51
1:A:53:MET:HE3	1:A:56:ARG:HB2	1.92	0.51
1:C:54:LYS:HB3	1:C:278:GLU:HB3	1.93	0.51
1:E:183:HIS:O	1:E:215:PRO:HB3	2.11	0.51
4:H:3:SIA:O10	4:H:3:SIA:O7	2.25	0.50
2:B:14:TRP:CH2	2:B:25:ARG:HG3	2.46	0.50
1:A:230:ILE:HD13	1:A:252:ILE:HG13	1.91	0.50
2:D:37:ASP:O	2:D:41:THR:HG23	2.11	0.50
2:F:44:ALA:HA	2:F:110:MET:CE	2.37	0.50
1:C:90:ARG:NH1	1:C:272:ALA:O	2.38	0.50
1:E:182:ILE:HD12	1:E:202:ILE:HD13	1.92	0.50
1:A:53:MET:HE2	1:A:58:HIS:CB	2.41	0.50
1:C:142:ASN:N	8:C:512:HOH:O	2.45	0.49
2:D:145:ASP:O	2:D:148:CYS:N	2.40	0.49
1:E:300:ARG:HD2	2:F:69:GLU:OE2	2.13	0.49
1:A:16:GLY:C	2:B:115:MET:HE1	2.36	0.49
1:C:190:GLU:CD	7:C:401:SIA:O9	2.55	0.49
1:A:17:HIS:CD2	8:A:508:HOH:O	2.65	0.49
1:C:320:MET:HB3	2:D:111:ALA:HB1	1.94	0.49
1:E:135:THR:O	4:H:3:SIA:H4	2.12	0.49
1:A:52:CYS:HB2	1:A:279:SER:HB3	1.94	0.49
1:E:153:TRP:CH2	4:H:3:SIA:H7	2.48	0.49
1:A:178:ILE:HG13	1:A:257:VAL:HG12	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:30:GLN:OE1	2:D:145:ASP:HB2	2.13	0.48
2:D:151:SER:O	2:D:157:TYR:N	2.46	0.48
1:E:114:GLU:CD	1:E:263:ARG:HH22	2.21	0.48
1:E:139:CYS:O	1:E:146:SER:HB3	2.13	0.48
1:A:17:HIS:HB3	2:B:115:MET:HE2	1.93	0.48
1:E:191:LYS:NZ	1:E:196:GLY:O	2.40	0.48
2:F:26:HIS:HB2	2:F:32:THR:HA	1.95	0.48
1:E:109:ARG:NH1	1:E:113:MET:HG3	2.27	0.48
2:B:21:TRP:HB2	2:B:41:THR:HB	1.96	0.48
1:A:279:SER:OG	1:A:280:LYS:N	2.46	0.48
1:C:35:GLU:HG2	1:C:322:ASN:HB3	1.96	0.47
1:E:320:MET:HG3	1:E:321:ARG:O	2.13	0.47
1:E:53:MET:HE2	1:E:58:HIS:CB	2.43	0.47
1:E:137:ARG:N	4:H:3:SIA:O1B	2.41	0.47
1:C:17:HIS:CE1	2:D:6:ILE:HG23	2.49	0.47
1:C:137:ARG:HB2	1:C:140:MET:HE2	1.97	0.47
2:B:148:CYS:O	2:B:151:SER:OG	2.25	0.47
1:C:17:HIS:HB3	2:D:115:MET:CE	2.44	0.47
1:C:266:LEU:HD11	1:C:302:VAL:HG13	1.97	0.47
1:E:54:LYS:HB3	1:E:278:GLU:HB2	1.97	0.47
1:A:16:GLY:O	2:B:115:MET:HE1	2.15	0.47
2:B:22:TYR:H	2:B:41:THR:CG2	2.25	0.47
2:B:53:ASN:O	2:B:57:GLU:HG2	2.15	0.47
2:B:117:ASN:ND2	2:F:2:LEU:HD22	2.28	0.47
1:C:153:TRP:CE3	7:C:401:SIA:H112	2.49	0.47
1:E:316:LEU:HD11	2:F:103:GLU:HG2	1.97	0.47
2:F:21:TRP:HB2	2:F:41:THR:HG22	1.96	0.47
1:A:212:ASN:HB2	1:E:216:VAL:HG11	1.96	0.47
2:B:14:TRP:CZ3	2:B:25:ARG:HG3	2.49	0.46
1:A:28:THR:HG22	1:A:30:THR:H	1.79	0.46
1:E:54:LYS:HD2	1:E:276(A):ASN:O	2.15	0.46
1:A:24:THR:HG22	1:A:25:ILE:N	2.31	0.46
2:F:17:MET:SD	2:F:23:GLY:HA3	2.56	0.46
2:F:172:ASN:ND2	8:F:209:HOH:O	2.49	0.46
1:A:13:ILE:HG22	2:B:140:ILE:HD11	1.98	0.46
1:E:15:LEU:HD13	2:F:119:TYR:HA	1.98	0.46
2:F:44:ALA:CA	2:F:110:MET:HE1	2.39	0.46
1:C:34:GLU:HG2	2:D:108:ILE:HD11	1.97	0.45
1:C:200:LEU:O	1:C:214:VAL:HG13	2.16	0.45
1:E:59:LYS:HZ3	1:E:59:LYS:HG3	1.71	0.45
1:E:225:GLY:N	8:E:514:HOH:O	2.48	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:37:ASP:O	2:F:41:THR:HG23	2.15	0.45
2:F:133:ASP:CG	2:F:137:CYS:HB2	2.40	0.45
1:C:310:ASN:O	1:C:311:ARG:NH1	2.46	0.45
1:E:98:TYR:HD2	1:E:136:THR:HG21	1.82	0.45
2:F:99:LEU:O	2:F:103:GLU:HB2	2.16	0.45
1:A:271:ASP:OD1	8:A:504:HOH:O	2.21	0.45
1:E:121:ILE:HD13	1:E:176:HIS:CE1	2.51	0.45
1:E:204:VAL:O	1:E:210:ARG:HA	2.17	0.45
1:E:288:ILE:HG12	1:E:297:LEU:HD13	1.99	0.44
1:A:17:HIS:HD2	8:A:508:HOH:O	2.00	0.44
1:C:153:TRP:CH2	7:C:401:SIA:H7	2.53	0.44
1:C:221:PRO:O	1:C:229:ARG:NH2	2.50	0.44
1:C:16:GLY:O	2:D:115:MET:HE1	2.17	0.44
1:E:294:PHE:HB3	1:E:309:VAL:HG22	2.00	0.44
1:C:121:ILE:HD12	1:C:121:ILE:N	2.33	0.43
2:D:22:TYR:H	2:D:41:THR:HG22	1.83	0.43
1:E:192:ASN:OD1	8:E:502:HOH:O	2.20	0.43
2:F:51:LYS:HE3	2:F:107:THR:OG1	2.18	0.43
2:B:76:GLN:O	2:B:80:VAL:HG23	2.18	0.43
1:C:51:LEU:O	1:C:53:MET:HG2	2.18	0.43
1:E:71:ILE:HD12	1:E:71:ILE:HA	1.60	0.43
1:E:320:MET:HE2	1:E:320:MET:HB3	1.89	0.43
2:F:172:ASN:O	8:F:203:HOH:O	2.21	0.43
1:E:172:ASP:CG	1:E:173:THR:H	2.26	0.43
1:A:269:GLN:HB3	2:B:70:PHE:CE2	2.52	0.43
1:C:311:ARG:NE	2:D:97:GLU:OE2	2.44	0.43
1:E:236:LEU:HD23	1:E:236:LEU:HA	1.72	0.43
2:F:72:GLU:HG2	2:F:73:ILE:N	2.32	0.43
1:C:53:MET:HE2	1:C:58:HIS:HB2	2.00	0.43
1:A:111:LYS:HE3	1:A:111:LYS:HB3	1.69	0.43
2:B:4:GLY:O	2:B:8:GLY:HA3	2.19	0.43
1:C:247:HIS:ND1	8:C:501:HOH:O	2.10	0.43
1:A:108:LEU:HD22	1:A:234:TRP:CD1	2.54	0.43
1:E:153:TRP:CD2	4:H:3:SIA:H112	2.54	0.43
1:E:185:PRO:HB2	1:E:217:VAL:HG22	2.00	0.43
1:A:44:GLU:N	1:A:292:LEU:HD12	2.34	0.43
1:A:312:ARG:NH1	1:A:313:SER:HB2	2.34	0.43
1:C:321:ARG:NH1	8:C:518:HOH:O	2.53	0.42
2:D:126:LEU:HD23	2:D:126:LEU:HA	1.90	0.42
1:E:96:TYR:OH	1:E:102:THR:OG1	2.34	0.42
1:A:31:ASN:HD22	1:A:34:GLU:CD	2.27	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:21:TRP:H	2:D:41:THR:CG2	2.32	0.42
1:E:84:TRP:CH2	1:E:118:ILE:HG13	2.54	0.42
1:E:179:MET:HE2	1:E:232:PHE:HB3	2.02	0.42
1:E:230:ILE:HD13	1:E:252:ILE:HG13	2.01	0.42
2:B:126:LEU:HA	2:B:126:LEU:HD23	1.85	0.42
7:C:401:SIA:C7	7:C:401:SIA:C10	2.98	0.42
1:E:71:ILE:O	1:E:148:TYR:HB3	2.20	0.42
2:B:22:TYR:HD2	2:B:40:SER:HG	1.67	0.42
2:D:110:MET:C	2:D:110:MET:SD	3.03	0.42
1:A:83:MET:HB3	1:A:83:MET:HE3	1.79	0.41
2:D:53:ASN:HA	2:D:56:VAL:HG22	2.02	0.41
1:C:147:PHE:O	1:C:148:TYR:C	2.63	0.41
1:E:259:LYS:HD2	1:E:261:ILE:HD11	2.01	0.41
1:C:179:MET:O	1:C:254:PRO:HB3	2.21	0.41
1:E:283:TRP:CD1	1:E:286:GLY:O	2.73	0.41
2:B:125:GLN:OE1	2:B:155:ASN:HA	2.19	0.41
2:B:141:TYR:CD1	2:B:170:ARG:HG2	2.55	0.41
1:E:53:MET:HE2	1:E:58:HIS:HB3	2.03	0.41
1:E:60:ASP:HB2	1:E:274:ILE:HD11	2.02	0.41
2:B:75:HIS:NE2	1:C:114:GLU:OE1	2.50	0.41
1:A:30:THR:O	2:D:50:GLY:HA3	2.21	0.41
2:D:147:SER:HB3	8:D:301:HOH:O	2.21	0.41
1:A:140:MET:HE3	1:A:140:MET:HB3	1.98	0.41
1:A:201:SER:H	1:A:248:ASN:HB2	1.85	0.41
1:A:325:GLU:H	1:A:325:GLU:HG3	1.66	0.41
2:B:26:HIS:HB2	2:B:149:MET:SD	2.61	0.41
2:B:151:SER:CA	2:B:156:THR:HB	2.49	0.41
1:E:97:CYS:HB2	1:E:138:ALA:O	2.20	0.41
1:E:283:TRP:CE3	1:E:298:SER:HB2	2.56	0.41
1:A:321:ARG:HG3	2:B:108:ILE:HG23	2.03	0.40
1:A:65:HIS:ND1	1:A:66:PRO:HD2	2.35	0.40
1:A:224:ASN:ND2	1:C:244:THR:HG21	2.36	0.40
1:A:295:GLN:O	1:A:308:TYR:HA	2.22	0.40
2:B:64:GLU:OE1	8:F:201:HOH:O	2.21	0.40

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	A	316/323 (98%)	308 (98%)	8 (2%)	0	100	100
1	C	316/323 (98%)	307 (97%)	9 (3%)	0	100	100
1	E	316/323 (98%)	299 (95%)	17 (5%)	0	100	100
2	B	169/180 (94%)	168 (99%)	1 (1%)	0	100	100
2	D	169/180 (94%)	165 (98%)	4 (2%)	0	100	100
2	F	163/180 (91%)	158 (97%)	5 (3%)	0	100	100
All	All	1449/1509 (96%)	1405 (97%)	44 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	270/273 (99%)	262 (97%)	8 (3%)	36	72
1	C	270/273 (99%)	266 (98%)	4 (2%)	57	84
1	E	270/273 (99%)	264 (98%)	6 (2%)	45	78
2	B	146/154 (95%)	140 (96%)	6 (4%)	27	62
2	D	146/154 (95%)	140 (96%)	6 (4%)	27	62
2	F	145/154 (94%)	138 (95%)	7 (5%)	23	56
All	All	1247/1281 (97%)	1210 (97%)	37 (3%)	36	72

All (37) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	HIS
1	A	50	ARG
1	A	141	ARG
1	A	142	ASN
1	A	210	ARG
1	A	288	ILE
1	A	312	ARG
1	A	325	GLU
2	B	32	THR
2	B	73	ILE
2	B	89	THR
2	B	100	VAL
2	B	140	ILE
2	B	155	ASN
1	C	18	HIS
1	C	97	CYS
1	C	154	LEU
1	C	321	ARG
2	D	39	LYS
2	D	58	LYS
2	D	62	GLU
2	D	100	VAL
2	D	102	MET
2	D	167	LEU
1	E	18	HIS
1	E	59	LYS
1	E	71	ILE
1	E	73	THR
1	E	178	ILE
1	E	271	ASP
2	F	11	GLU
2	F	30	GLN
2	F	52	LEU
2	F	55	LEU
2	F	72	GLU
2	F	73	ILE
2	F	168	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	142	ASN
1	A	166	ASN
2	B	95	GLN
1	C	276	ASN
2	D	105	GLN
1	E	18	HIS
1	E	104	ASN
1	E	192	ASN
1	E	248	ASN
2	F	47	GLN
2	F	79	ASN
2	F	105	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 ⓘ

9 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	NAG	G	1	1,3	14,14,15	1.20	2 (14%)	17,19,21	1.45	3 (17%)
3	NAG	G	2	3	14,14,15	0.73	1 (7%)	17,19,21	1.60	4 (23%)
4	NAG	H	1	4	15,15,15	0.56	0	21,21,21	1.09	2 (9%)
4	GAL	H	2	4	11,11,12	0.63	0	15,15,17	0.63	0
4	SIA	H	3	4	20,20,21	1.03	2 (10%)	21,28,31	1.78	5 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	I	1	1,5	14,14,15	0.74	1 (7%)	17,19,21	0.81	0
5	NAG	I	2	5	14,14,15	0.20	0	17,19,21	0.54	0
5	BMA	I	3	5	11,11,12	1.44	2 (18%)	15,15,17	0.87	1 (6%)
5	MAN	I	4	5	11,11,12	1.55	3 (27%)	15,15,17	1.20	2 (13%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	G	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	2/6/23/26	0/1/1/1
4	NAG	H	1	4	-	3/6/26/26	0/1/1/1
4	GAL	H	2	4	-	0/2/19/22	0/1/1/1
4	SIA	H	3	4	-	1/18/34/38	0/1/1/1
5	NAG	I	1	1,5	-	0/6/23/26	0/1/1/1
5	NAG	I	2	5	-	2/6/23/26	0/1/1/1
5	BMA	I	3	5	-	0/2/19/22	0/1/1/1
5	MAN	I	4	5	-	0/2/19/22	0/1/1/1

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	I	4	MAN	O5-C1	3.13	1.49	1.43
5	I	3	BMA	O5-C1	2.89	1.48	1.43
3	G	1	NAG	C2-N2	-2.84	1.41	1.46
5	I	3	BMA	O5-C5	2.75	1.48	1.43
4	H	3	SIA	O1B-C1	-2.62	1.22	1.30
5	I	4	MAN	O5-C5	2.53	1.48	1.43
3	G	1	NAG	O5-C1	-2.32	1.39	1.43
5	I	1	NAG	O5-C1	-2.32	1.39	1.43
4	H	3	SIA	C2-C1	-2.19	1.49	1.52
5	I	4	MAN	O3-C3	2.15	1.48	1.43
3	G	2	NAG	C2-N2	-2.00	1.42	1.46

All (17) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	3	SIA	O9-C9-C8	-4.67	101.35	111.16
4	H	3	SIA	C6-C5-N5	-3.49	105.34	110.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	3	SIA	O1B-C1-C2	3.45	121.69	112.71
3	G	2	NAG	O5-C5-C6	-3.37	101.11	107.66
5	I	4	MAN	C1-O5-C5	2.77	115.90	112.19
4	H	3	SIA	O1A-C1-C2	-2.76	116.89	122.85
4	H	3	SIA	C9-C8-C7	2.76	117.79	112.17
3	G	1	NAG	C1-C2-N2	2.71	114.71	110.43
3	G	1	NAG	C2-N2-C7	-2.67	119.32	122.90
3	G	2	NAG	C1-C2-N2	-2.62	106.30	110.43
3	G	2	NAG	C2-N2-C7	-2.47	119.59	122.90
3	G	2	NAG	O3-C3-C2	2.42	114.42	109.40
3	G	1	NAG	O5-C5-C6	2.28	112.10	107.66
4	H	1	NAG	C1-C2-N2	-2.27	108.09	110.73
5	I	4	MAN	C1-C2-C3	2.27	112.95	109.64
5	I	3	BMA	C1-O5-C5	2.23	115.17	112.19
4	H	1	NAG	O5-C1-C2	2.20	111.73	109.52

There are no chirality outliers.

All (8) torsion outliers are listed below:

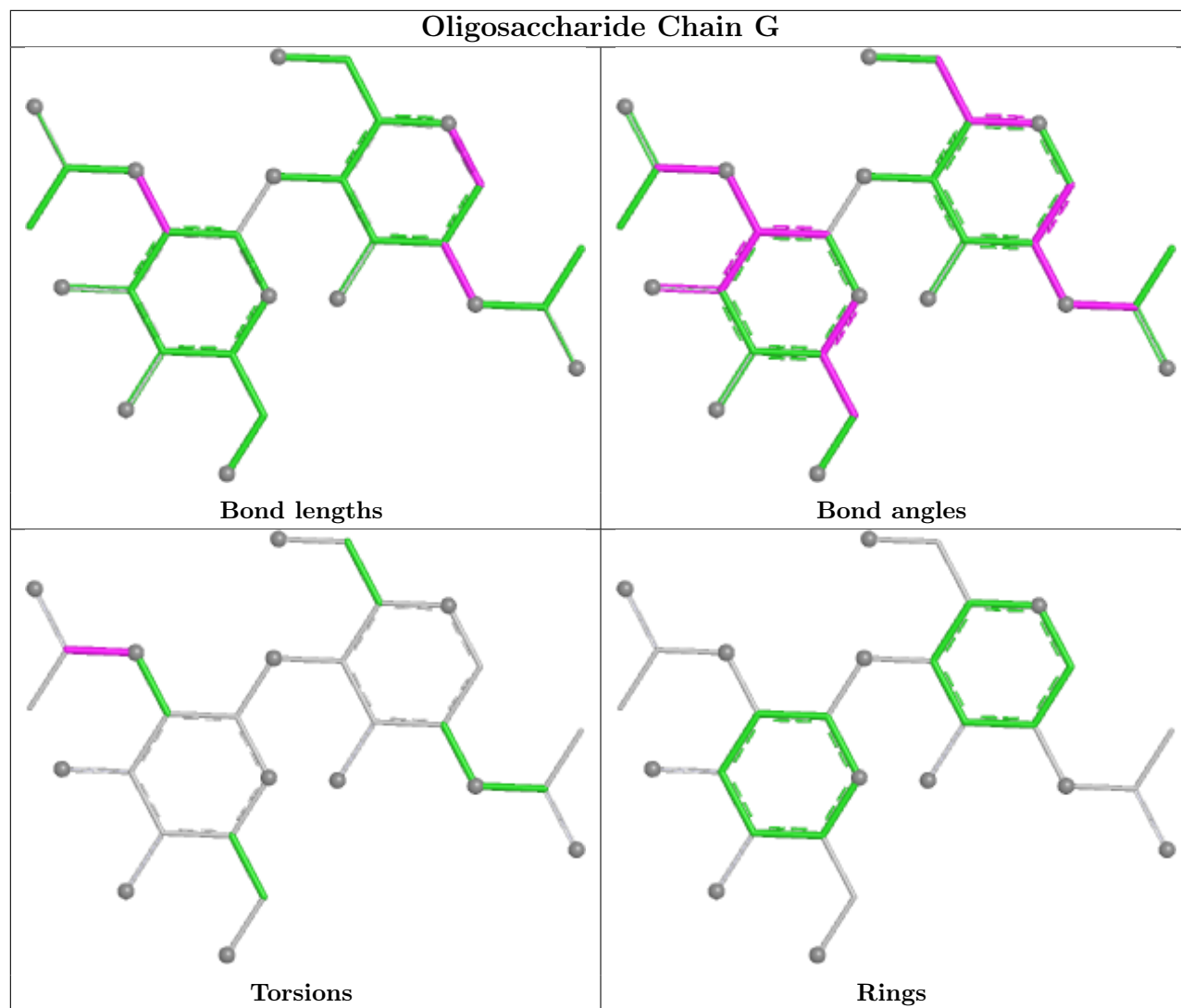
Mol	Chain	Res	Type	Atoms
4	H	1	NAG	C3-C2-N2-C7
4	H	1	NAG	C8-C7-N2-C2
4	H	1	NAG	O7-C7-N2-C2
3	G	2	NAG	C8-C7-N2-C2
5	I	2	NAG	O5-C5-C6-O6
5	I	2	NAG	C4-C5-C6-O6
3	G	2	NAG	O7-C7-N2-C2
4	H	3	SIA	C6-C5-N5-C10

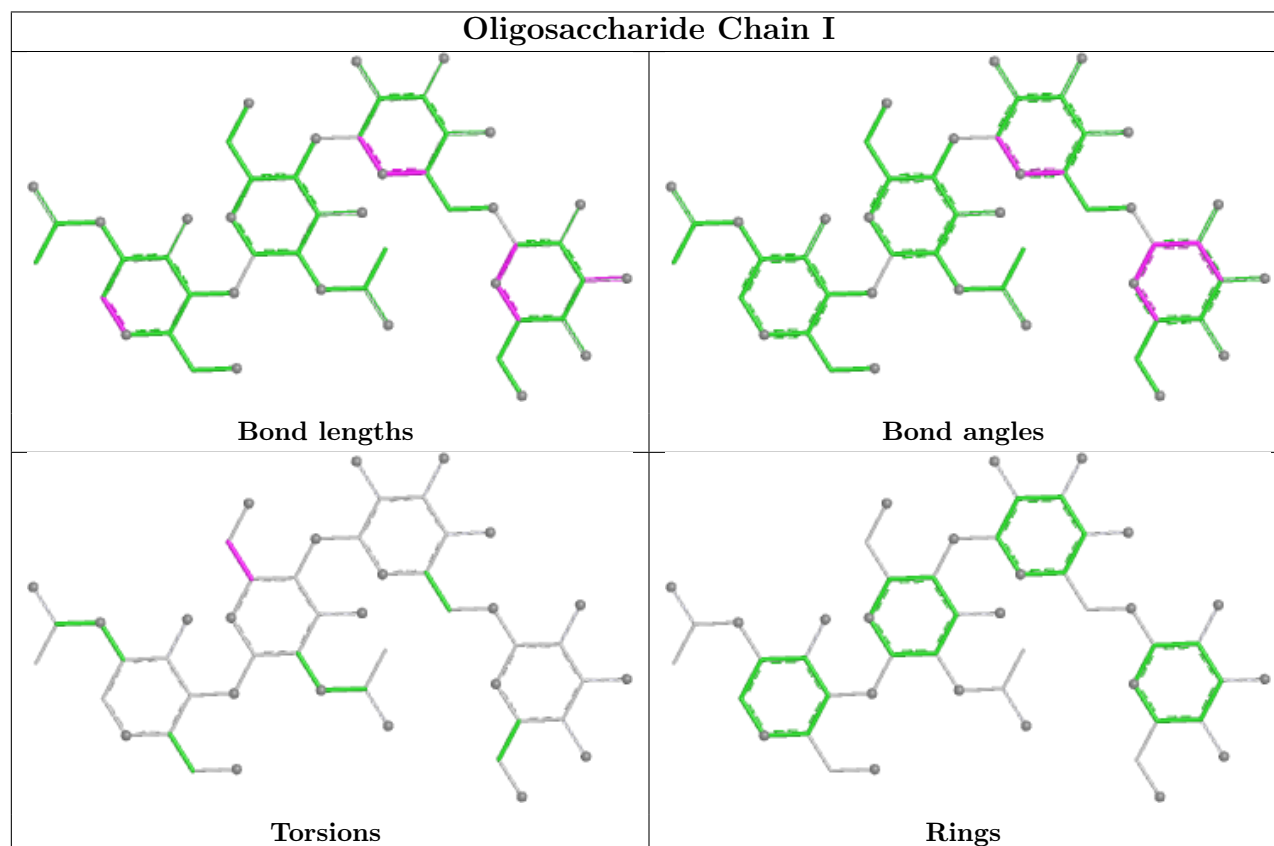
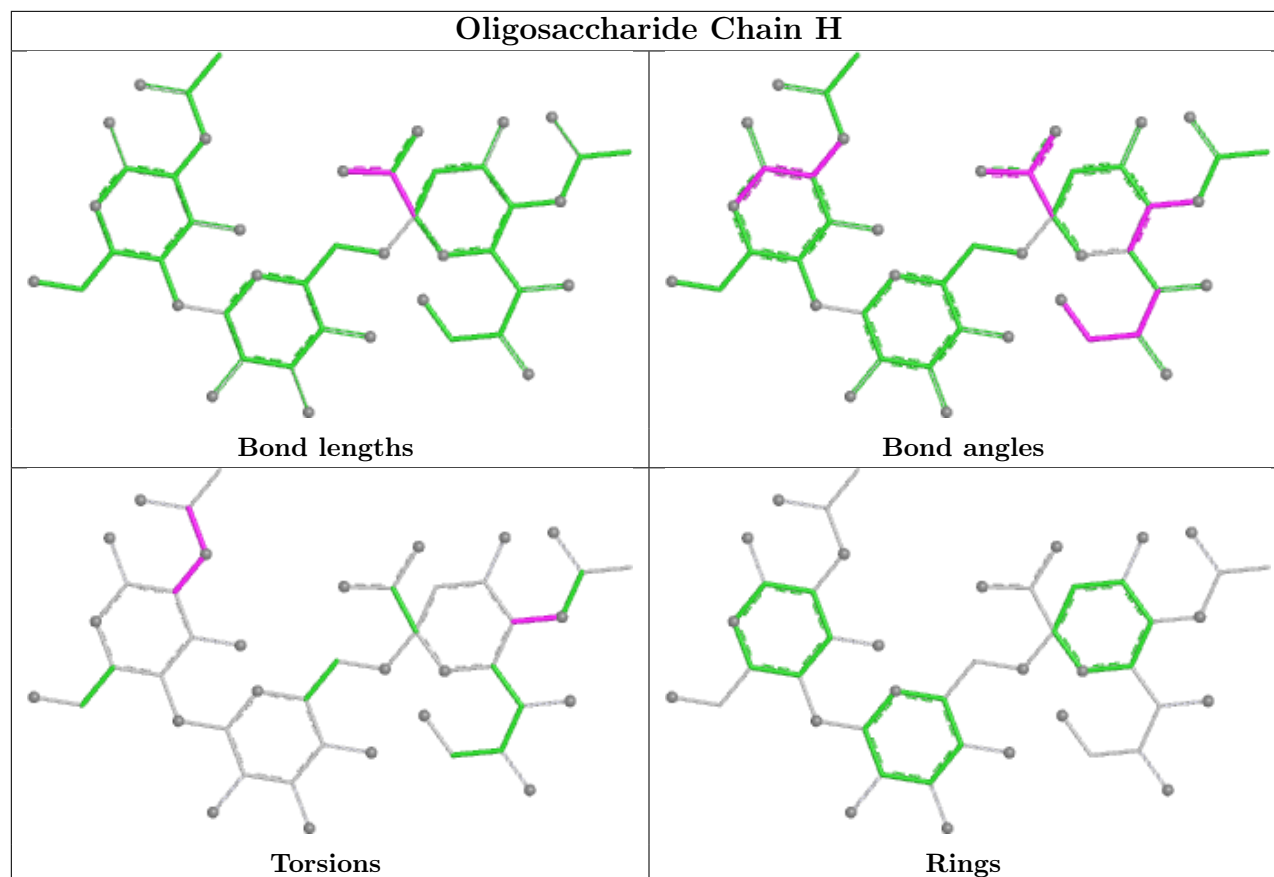
There are no ring outliers.

2 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	I	1	NAG	2	0
4	H	3	SIA	6	0

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





5.6 Ligand geometry

4 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	D	201	2	14,14,15	0.68	1 (7%)	17,19,21	2.14	5 (29%)
6	NAG	C	402	1	14,14,15	1.03	2 (14%)	17,19,21	2.34	4 (23%)
7	SIA	C	401	-	21,21,21	1.14	2 (9%)	24,31,31	1.24	3 (12%)
6	NAG	A	401	1	14,14,15	0.75	1 (7%)	17,19,21	0.88	1 (5%)

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	D	201	2	-	2/6/23/26	0/1/1/1
6	NAG	C	402	1	-	3/6/23/26	0/1/1/1
7	SIA	C	401	-	-	3/20/38/38	0/1/1/1
6	NAG	A	401	1	-	0/6/23/26	0/1/1/1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	C	401	SIA	O2-C2	-3.76	1.34	1.39
6	C	402	NAG	C2-N2	-2.48	1.42	1.46
6	A	401	NAG	O5-C1	-2.37	1.39	1.43
6	C	402	NAG	O5-C1	-2.30	1.39	1.43
7	C	401	SIA	C2-C1	-2.15	1.50	1.53
6	D	201	NAG	O5-C1	-2.11	1.40	1.43

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	C	402	NAG	C1-C2-N2	-6.59	100.05	110.43
6	C	402	NAG	C6-C5-C4	-5.31	99.97	113.02
6	D	201	NAG	C1-C2-N2	-4.40	103.50	110.43
6	D	201	NAG	C2-N2-C7	4.22	128.56	122.90
6	D	201	NAG	O4-C4-C3	3.52	118.66	110.38
7	C	401	SIA	O1A-C1-C2	-3.31	118.33	123.85
7	C	401	SIA	O2-C2-O6	-3.00	101.95	109.51
6	C	402	NAG	O5-C5-C6	-2.94	101.94	107.66
7	C	401	SIA	O6-C6-C5	2.87	112.47	109.84
6	D	201	NAG	O6-C6-C5	-2.65	102.32	111.33
6	D	201	NAG	C6-C5-C4	-2.48	106.93	113.02
6	C	402	NAG	O3-C3-C2	2.10	113.76	109.40
6	A	401	NAG	C1-C2-N2	-2.04	107.22	110.43

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	C	402	NAG	C8-C7-N2-C2
6	C	402	NAG	O7-C7-N2-C2
6	D	201	NAG	O5-C5-C6-O6
7	C	401	SIA	O8-C8-C9-O9
6	D	201	NAG	C4-C5-C6-O6
7	C	401	SIA	C7-C8-C9-O9
6	C	402	NAG	O5-C5-C6-O6
7	C	401	SIA	C6-C5-N5-C10

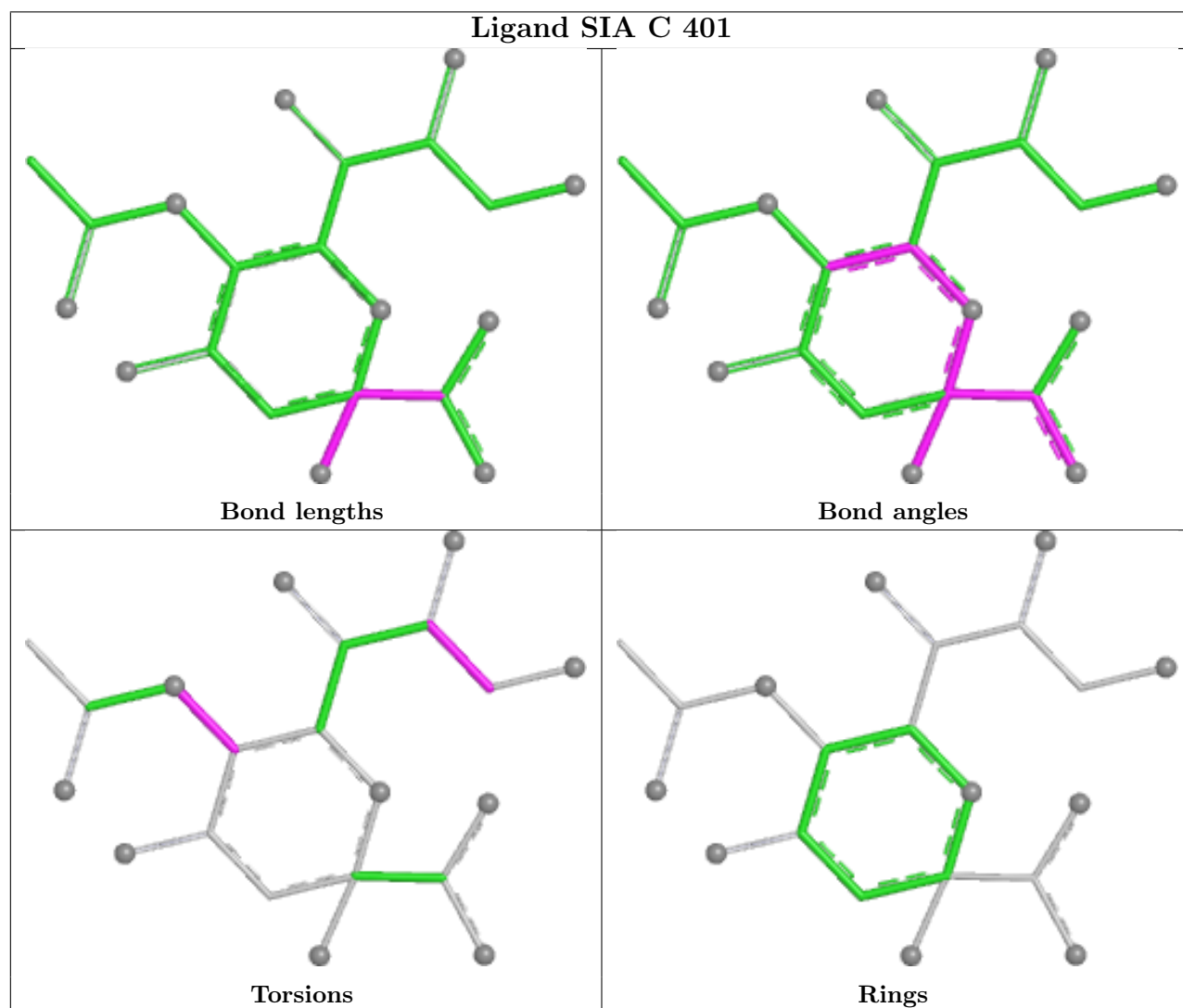
There are no ring outliers.

1 monomer is involved in 7 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
7	C	401	SIA	7	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier.

The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	318/323 (98%)	-0.22	1 (0%) 90 86	15, 28, 53, 73	0
1	C	318/323 (98%)	-0.45	0 100 100	14, 23, 39, 55	0
1	E	318/323 (98%)	-0.39	0 100 100	16, 25, 45, 68	0
2	B	171/180 (95%)	0.17	2 (1%) 76 68	17, 43, 67, 79	0
2	D	171/180 (95%)	-0.26	0 100 100	17, 35, 50, 60	0
2	F	169/180 (93%)	-0.01	2 (1%) 76 68	21, 41, 55, 69	0
All	All	1465/1509 (97%)	-0.24	5 (0%) 90 86	14, 30, 55, 79	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	23	GLY	3.0
2	B	57	GLU	2.7
2	F	58	LYS	2.4
1	A	49	ASN	2.3
2	F	32	THR	2.1

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](#)

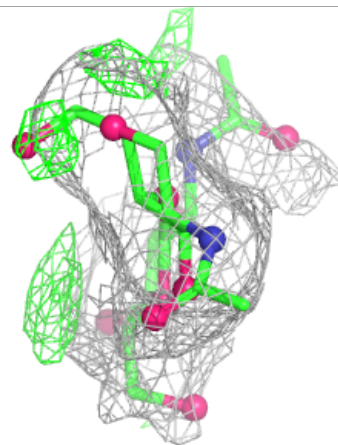
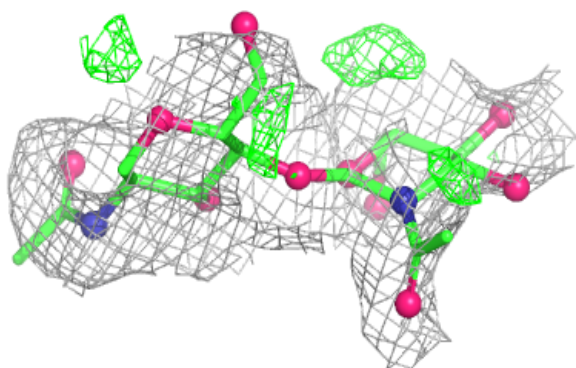
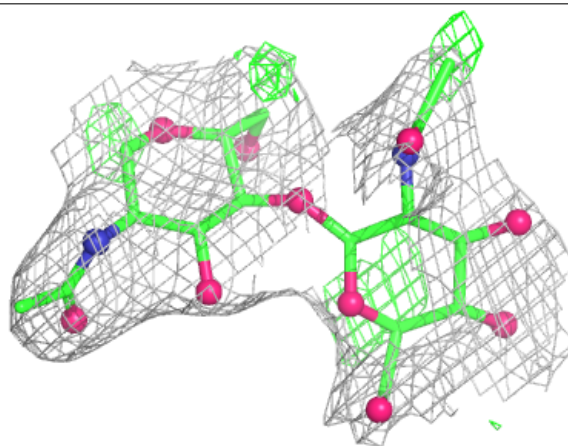
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
5	MAN	I	4	11/12	0.38	0.17	87,103,109,109	0
4	NAG	H	1	15/15	0.50	0.19	79,94,106,107	0
3	NAG	G	2	14/15	0.57	0.16	71,79,81,81	0
5	BMA	I	3	11/12	0.78	0.11	71,76,89,95	0
5	NAG	I	2	14/15	0.80	0.13	50,60,65,70	0
3	NAG	G	1	14/15	0.81	0.13	39,48,66,66	0
4	GAL	H	2	11/12	0.86	0.11	63,84,90,94	0
4	SIA	H	3	20/21	0.88	0.10	34,45,56,58	0
5	NAG	I	1	14/15	0.91	0.09	35,40,46,50	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.

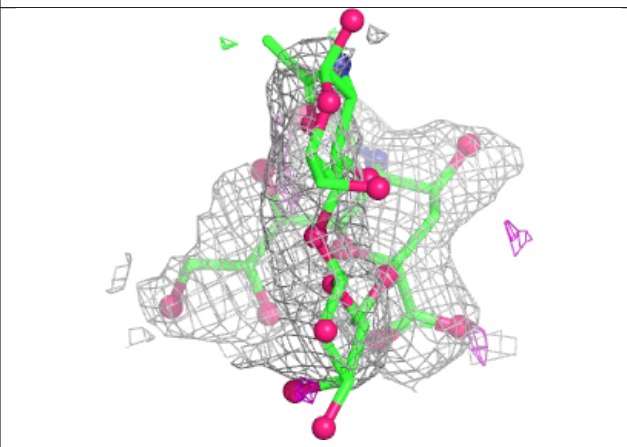
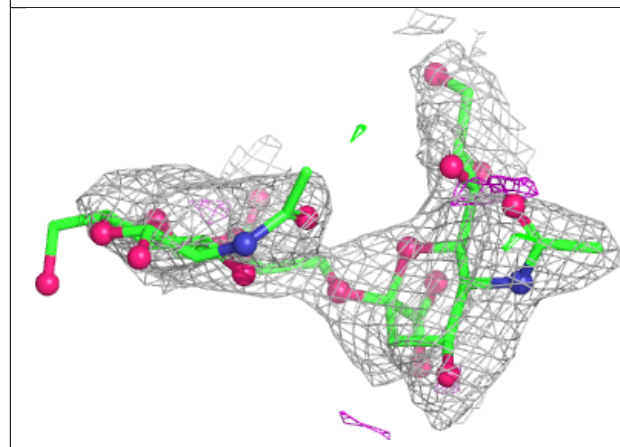
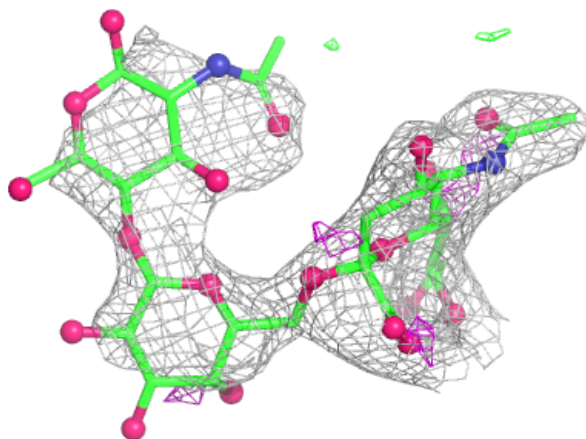
Electron density around Chain G:

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

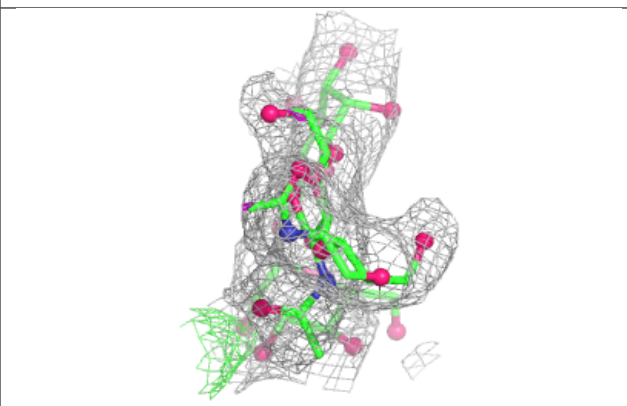
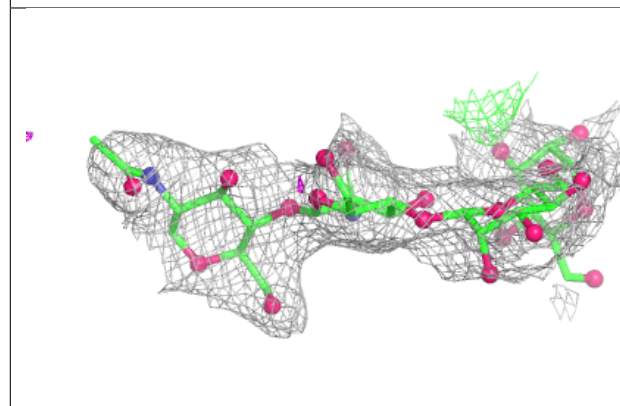
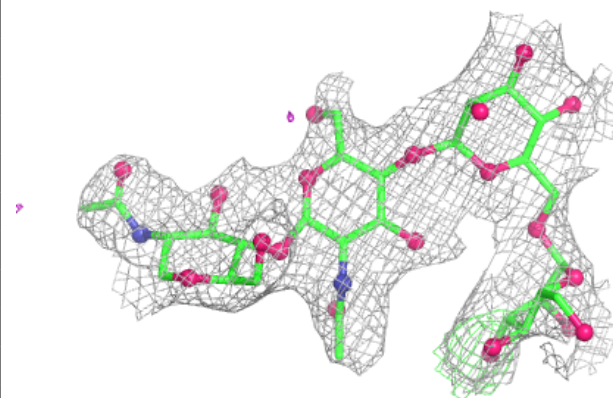


Electron density around Chain H:

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

**Electron density around Chain I:**

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



6.4 Ligands

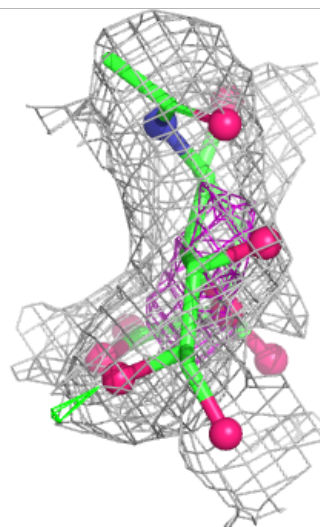
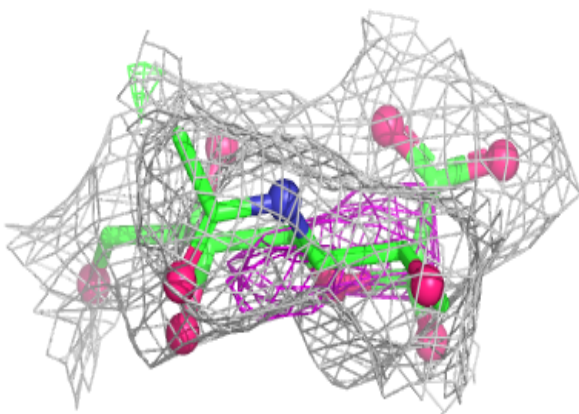
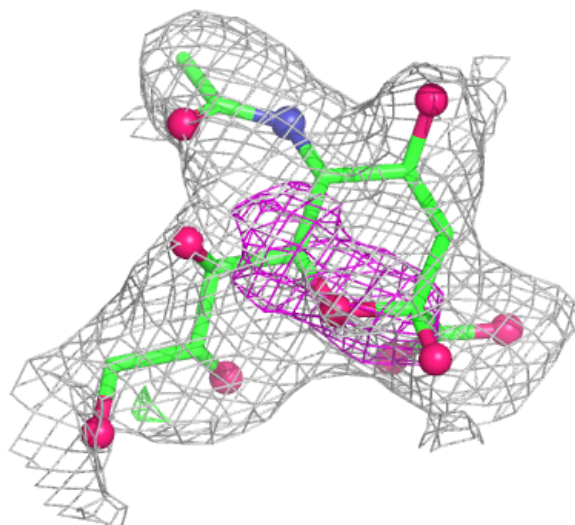
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	C	402	14/15	0.57	0.16	76,82,88,89	0
6	NAG	D	201	14/15	0.66	0.15	31,52,59,64	0
6	NAG	A	401	14/15	0.67	0.15	45,52,62,67	0
7	SIA	C	401	21/21	0.71	0.14	37,53,68,68	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around SIA C 401:

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



6.5 Other polymers [i](#)

There are no such residues in this entry.