



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 12:44 AM UTC

PDB ID : 5THB / pdb_00005thb
Title : Crystal structure of H10 hemagglutinin mutant (T193D-Q226L-G228S) from Jiangxi-Donghu (2013) H10N8 influenza virus
Authors : Tzarum, N.; Wilson, I.A.
Deposited on : 2016-09-29
Resolution : 2.41 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

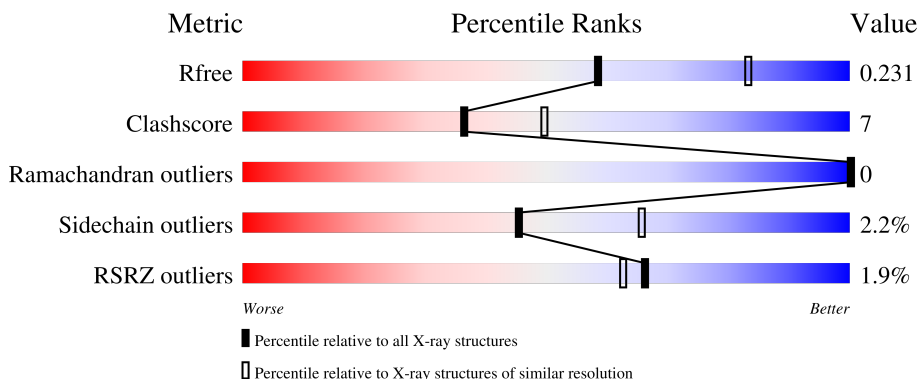
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.41 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	6062 (2.44-2.40)
Clashscore	190562	6562 (2.44-2.40)
Ramachandran outliers	187476	6481 (2.44-2.40)
Sidechain outliers	187428	6482 (2.44-2.40)
RSRZ outliers	180081	6066 (2.44-2.40)

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	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 85%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 12%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">85% 12% ..</p>
1	C	323	<div style="display: flex; align-items: center;"> <div style="width: 82%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 3%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">82% 15% ..</p>
1	E	323	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 83%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">3% 83% 15% .</p>
2	B	180	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 74%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 21%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 2%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">3% 74% 21% . 5%</p>
2	D	180	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 77%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 16%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 3%; height: 10px; background-color: orange; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">3% 77% 16% . 6%</p>

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Mol	Chain	Length	Quality of chain
2	F	180	
3	G	4	
4	H	2	

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 12224 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
			Total	C	N	O	S			
1	A	318	2437	1508	448	464	17	0	0	0
1	C	319	2445	1514	449	465	17	0	0	0
1	E	318	2437	1508	448	464	17	0	0	0

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
			Total	C	N	O	S			
2	B	171	1382	854	239	281	8	0	0	0
2	D	169	1367	846	236	277	8	0	0	0
2	F	169	1371	848	237	278	8	0	0	0

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 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.



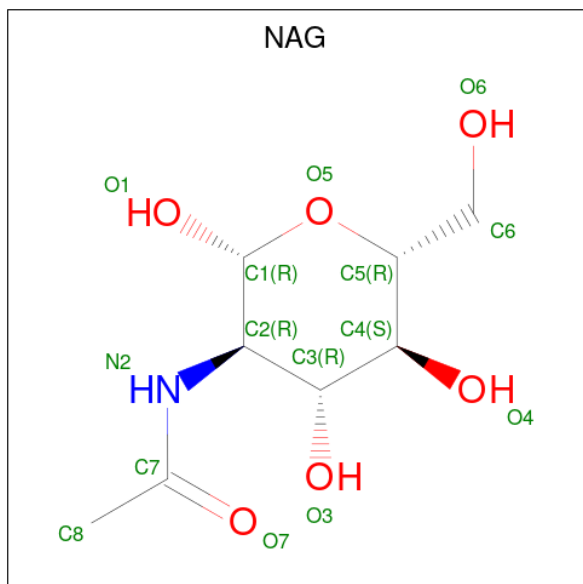
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace	
			Total	C	N				O
3	G	4	50	28	2	20	0	0	0

- Molecule 4 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
			Total	C	N	O			
4	H	2	28	16	2	10	0	0	0

- Molecule 5 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	138	Total	O	0	0
			138	138		
6	B	64	Total	O	0	0
			64	64		
6	C	150	Total	O	0	0
			150	150		

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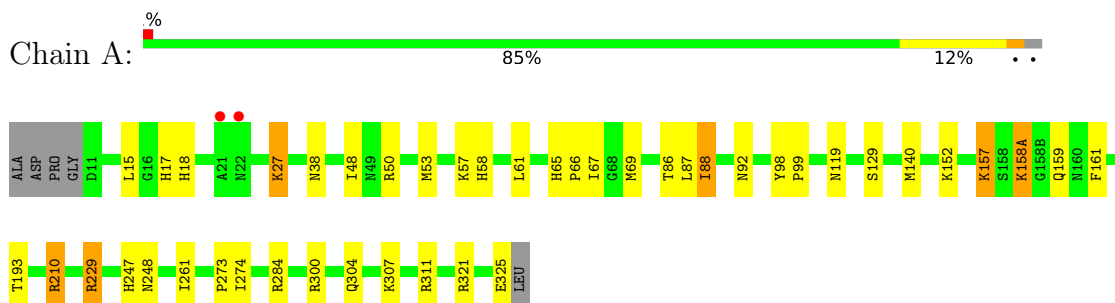
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	D	72	Total 72	O 72	0	0
6	E	135	Total 135	O 135	0	0
6	F	78	Total 78	O 78	0	0

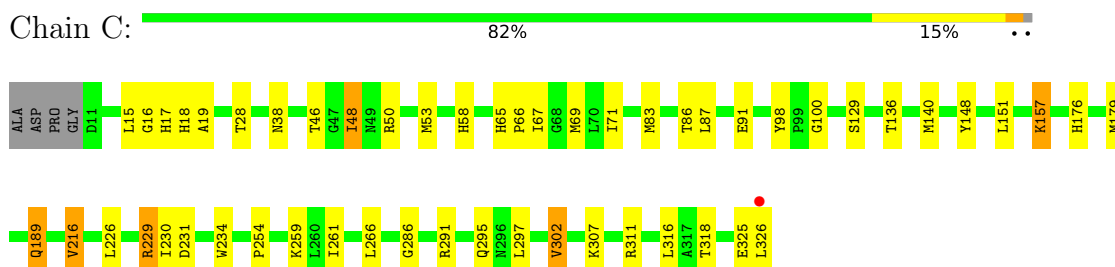
3 Residue-property plots [i](#)

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.

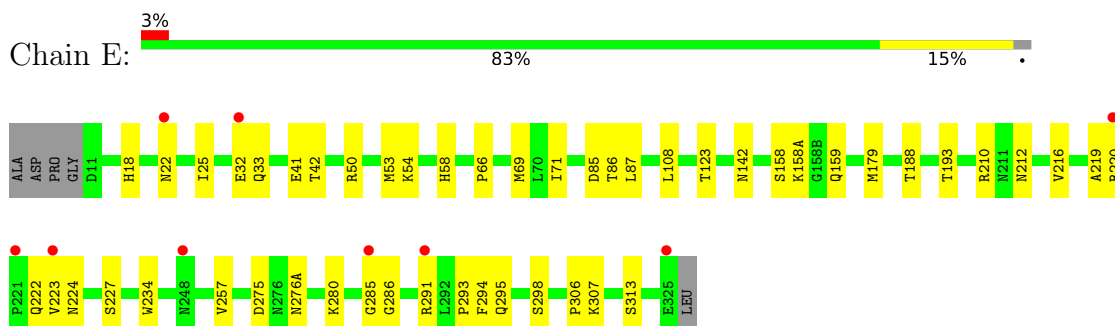
- Molecule 1: Hemagglutinin HA1 chain



- Molecule 1: Hemagglutinin HA1 chain

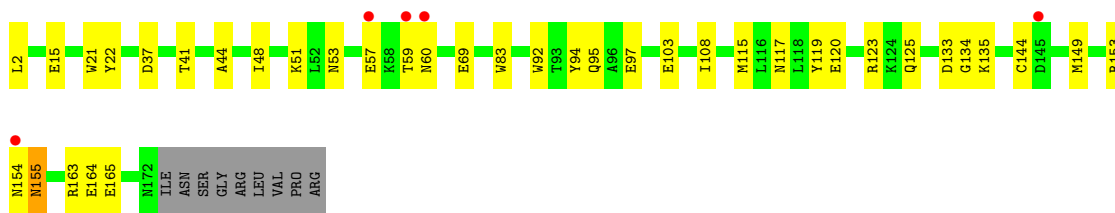


- Molecule 1: Hemagglutinin HA1 chain

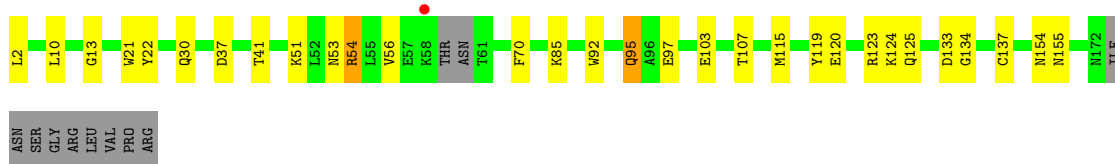
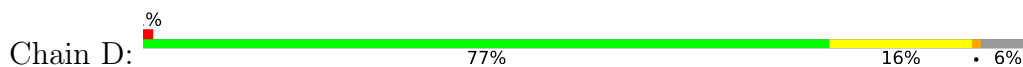


- Molecule 2: Hemagglutinin HA2 chain

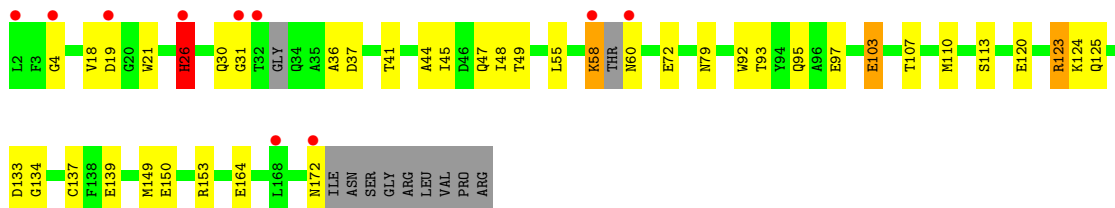




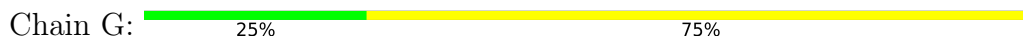
- Molecule 2: Hemagglutinin HA2 chain



- Molecule 2: Hemagglutinin HA2 chain



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



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



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	63.83Å 241.29Å 70.06Å 90.00° 111.43° 90.00°	Depositor
Resolution (Å)	33.05 – 2.41 33.05 – 2.41	Depositor EDS
% Data completeness (in resolution range)	96.8 (33.05-2.41) 95.7 (33.05-2.41)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.93 (at 2.39Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.180 , 0.230 0.184 , 0.231	Depositor DCC
R_{free} test set	3707 reflections (4.85%)	wwPDB-VP
Wilson B-factor (Å ²)	21.9	Xtrriage
Anisotropy	0.259	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12224	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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: NAG, BMA, 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	0.68	0/2486	0.94	7/3369 (0.2%)
1	C	0.67	0/2494	0.91	2/3380 (0.1%)
1	E	0.68	0/2486	1.03	15/3369 (0.4%)
2	B	0.65	0/1407	0.91	0/1899
2	D	0.62	0/1391	0.88	0/1875
2	F	0.61	0/1394	0.92	4/1878 (0.2%)
All	All	0.66	0/11658	0.94	28/15770 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	1
2	B	0	1
2	F	0	2
All	All	0	4

There are no bond length outliers.

All (28) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	220	ARG	CA-C-N	-9.65	110.11	119.76
1	E	220	ARG	C-N-CA	-9.65	110.11	119.76
1	E	220	ARG	N-CA-C	7.80	123.40	113.25
1	E	298	SER	CA-C-N	7.72	128.38	120.04
1	E	298	SER	C-N-CA	7.72	128.38	120.04
1	E	285	GLY	N-CA-C	-7.61	102.32	112.14
1	C	48	ILE	N-CA-C	7.58	119.22	109.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	286	GLY	N-CA-C	7.56	131.09	113.18
2	F	26	HIS	N-CA-C	7.44	121.95	110.42
2	F	31	GLY	N-CA-C	7.36	130.61	113.18
1	E	224	ASN	N-CA-C	-6.24	100.00	108.24
1	A	210	ARG	NE-CZ-NH2	6.02	124.62	119.20
1	A	247	HIS	N-CA-C	5.88	118.22	108.99
1	A	158(A)	LYS	N-CA-C	5.79	118.33	111.33
1	A	38	ASN	N-CA-C	5.76	117.61	108.79
1	A	210	ARG	NE-CZ-NH1	-5.68	115.82	121.50
1	C	286	GLY	N-CA-C	5.49	118.62	110.60
2	F	18	VAL	N-CA-C	-5.29	107.32	112.29
1	E	291	ARG	N-CA-C	-5.28	107.01	113.50
2	F	134	GLY	N-CA-C	5.24	122.33	115.47
1	E	32	GLU	N-CA-C	-5.16	102.85	110.48
1	A	159	GLN	N-CA-C	5.10	118.03	110.48
1	E	85	ASP	N-CA-C	-5.08	107.11	113.72
1	E	142	ASN	CA-C-N	-5.04	111.54	121.41
1	E	142	ASN	C-N-CA	-5.04	111.54	121.41
1	A	88	ILE	N-CA-C	5.03	115.34	107.99
1	E	222	GLN	CA-C-N	-5.01	116.45	122.22
1	E	222	GLN	C-N-CA	-5.01	116.45	122.22

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	59	THR	Peptide
1	E	219	ALA	Peptide
2	F	30	GLN	Peptide
2	F	4	GLY	Peptide

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	2393	38	0
1	C	2445	0	2405	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2437	0	2394	27	0
2	B	1382	0	1278	34	0
2	D	1367	0	1264	35	0
2	F	1371	0	1266	27	0
3	G	50	0	43	0	0
4	H	28	0	25	0	0
5	A	28	0	26	0	0
5	B	14	0	13	0	0
5	D	14	0	13	0	0
5	F	14	0	13	1	0
6	A	138	0	0	4	0
6	B	64	0	0	4	0
6	C	150	0	0	2	1
6	D	72	0	0	6	0
6	E	135	0	0	3	1
6	F	78	0	0	8	0
All	All	12224	0	11133	164	1

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 7.

All (164) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:44:ALA:HA	2:F:110:MET:HE1	1.41	1.01
2:D:54:ARG:NH1	6:D:301:HOH:O	2.01	0.92
2:B:133:ASP:OD2	2:B:135:LYS:HE3	1.78	0.83
2:B:53:ASN:O	2:B:57:GLU:HG2	1.81	0.80
1:C:53:MET:HE2	1:C:58:HIS:HB3	1.69	0.75
2:F:153:ARG:NH1	6:F:304:HOH:O	2.19	0.74
1:C:28:THR:O	6:C:601:HOH:O	2.04	0.73
1:A:210:ARG:NH1	1:E:227:SER:O	2.22	0.72
1:A:53:MET:HE2	1:A:58:HIS:HB3	1.73	0.70
2:D:30:GLN:OE1	6:D:302:HOH:O	2.10	0.70
1:A:53:MET:HE2	1:A:58:HIS:CB	2.21	0.70
2:F:58:LYS:O	2:F:60:ASN:N	2.25	0.69
2:D:53:ASN:OD1	6:D:303:HOH:O	2.11	0.68
2:B:163:ARG:NH1	6:B:304:HOH:O	2.27	0.68
2:D:10:LEU:HD13	2:D:115:MET:HE3	1.76	0.67
2:F:120:GLU:OE1	2:F:123:ARG:NH2	2.27	0.67
1:A:69:MET:HE1	1:A:87:LEU:HD21	1.77	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:266:LEU:HD11	1:C:302:VAL:HG13	1.78	0.66
2:B:95:GLN:NE2	2:F:95:GLN:OE1	2.29	0.66
2:B:165:GLU:OE2	6:B:301:HOH:O	2.13	0.66
2:B:117:ASN:O	6:B:302:HOH:O	2.14	0.65
2:F:103:GLU:OE2	6:F:301:HOH:O	2.14	0.65
2:D:154:ASN:ND2	6:D:304:HOH:O	2.28	0.64
2:D:134:GLY:HA2	2:F:124:LYS:HD3	1.79	0.64
1:C:129:SER:O	1:C:157:LYS:HE3	1.97	0.64
2:D:125:GLN:HE22	2:D:155:ASN:HA	1.64	0.63
1:C:48:ILE:HG22	1:C:50:ARG:H	1.64	0.62
2:F:139:GLU:OE1	6:F:302:HOH:O	2.16	0.62
1:A:17:HIS:HB3	2:B:115:MET:CE	2.31	0.61
2:D:10:LEU:CD1	2:D:115:MET:HE3	2.30	0.61
1:C:16:GLY:C	2:D:115:MET:HE1	2.25	0.60
1:C:179:MET:HG2	1:C:234:TRP:HB3	1.82	0.60
1:C:231:ASP:OD2	1:E:210:ARG:NH1	2.35	0.59
1:C:189:GLN:NE2	6:C:603:HOH:O	2.27	0.59
2:B:37:ASP:O	2:B:41:THR:HG23	2.02	0.59
1:C:53:MET:HE2	1:C:58:HIS:CB	2.32	0.59
1:A:284:ARG:NH1	2:B:69:GLU:HG3	2.18	0.58
1:E:123:THR:HG22	1:E:257:VAL:HG13	1.85	0.58
2:D:120:GLU:OE1	2:D:123:ARG:NH2	2.22	0.58
2:F:26:HIS:CD2	2:F:149:MET:HE2	2.39	0.58
2:D:21:TRP:H	2:D:41:THR:CG2	2.17	0.57
2:B:154:ASN:O	2:B:155:ASN:HB2	2.03	0.57
1:A:88:ILE:HD13	1:A:274:ILE:HD11	1.85	0.57
2:F:79:ASN:OD1	5:F:201:NAG:C8	2.52	0.57
1:C:216:VAL:HG11	1:E:212:ASN:HB2	1.87	0.57
2:F:47:GLN:CD	2:F:110:MET:HE3	2.30	0.57
2:F:120:GLU:OE2	6:F:303:HOH:O	2.17	0.55
1:A:92:ASN:ND2	6:A:501:HOH:O	2.19	0.55
1:A:53:MET:CE	1:A:58:HIS:HB3	2.37	0.55
1:C:15:LEU:HD13	2:D:119:TYR:HA	1.89	0.55
1:C:229:ARG:HD2	6:E:816:HOH:O	2.06	0.55
2:D:22:TYR:H	2:D:41:THR:HG22	1.71	0.55
2:F:164:GLU:HG2	6:F:330:HOH:O	2.07	0.54
2:D:133:ASP:OD2	2:D:137:CYS:HB2	2.08	0.54
2:F:60:ASN:ND2	6:F:305:HOH:O	2.20	0.53
1:E:53:MET:HE2	1:E:58:HIS:CB	2.38	0.53
1:C:66:PRO:HA	1:C:69:MET:HE2	1.91	0.53
1:C:53:MET:HE1	1:C:86:THR:HG21	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:83:MET:HE1	1:C:261:ILE:HD12	1.91	0.53
1:A:321:ARG:HG3	2:B:108:ILE:HG23	1.92	0.52
1:C:17:HIS:HB3	2:D:115:MET:HE2	1.92	0.52
2:B:21:TRP:H	2:B:41:THR:CG2	2.22	0.52
2:F:19:ASP:HB3	2:F:36:ALA:HB2	1.91	0.52
2:F:93:THR:O	2:F:97:GLU:HG3	2.09	0.52
1:A:284:ARG:HG3	2:B:69:GLU:HG2	1.92	0.52
1:A:304:GLN:HB3	2:B:60:ASN:CB	2.41	0.51
1:A:50:ARG:NH2	1:A:273:PRO:HG2	2.26	0.51
2:D:37:ASP:O	2:D:41:THR:HG23	2.10	0.51
1:C:17:HIS:HB3	2:D:115:MET:CE	2.41	0.51
1:A:300:ARG:NE	6:A:510:HOH:O	2.44	0.50
1:C:151:LEU:HD23	1:C:254:PRO:HA	1.93	0.50
1:A:152:LYS:NZ	6:A:511:HOH:O	2.44	0.50
1:C:307:LYS:HG3	2:D:92:TRP:CE2	2.47	0.50
2:B:22:TYR:H	2:B:41:THR:HG22	1.75	0.50
2:B:125:GLN:OE1	2:B:155:ASN:HA	2.12	0.50
2:F:21:TRP:H	2:F:41:THR:CG2	2.24	0.50
1:A:229:ARG:HD2	6:A:595:HOH:O	2.11	0.50
1:A:158(A):LYS:O	1:A:193:THR:O	2.28	0.50
1:A:53:MET:HE2	1:A:58:HIS:HB2	1.91	0.49
2:F:47:GLN:NE2	2:F:110:MET:HE3	2.27	0.49
1:A:15:LEU:HD13	2:B:119:TYR:HA	1.94	0.49
2:B:21:TRP:H	2:B:41:THR:HG21	1.78	0.48
1:C:176:HIS:CD2	1:C:259:LYS:HG3	2.48	0.48
1:C:91:GLU:HB2	2:D:70:PHE:CE1	2.47	0.48
2:F:45:ILE:O	2:F:49:THR:HG23	2.13	0.48
1:A:65:HIS:CE1	1:A:67:ILE:HG12	2.48	0.48
2:F:110:MET:O	2:F:113:SER:HB3	2.13	0.48
1:E:42:THR:HG22	2:F:55:LEU:HD21	1.96	0.48
1:C:98:TYR:HD2	1:C:136:THR:HG21	1.79	0.48
1:E:53:MET:HE2	1:E:58:HIS:CG	2.49	0.48
1:A:304:GLN:HB3	2:B:60:ASN:HB2	1.96	0.47
1:E:22:ASN:OD1	1:E:22:ASN:N	2.41	0.47
2:F:133:ASP:CG	2:F:137:CYS:HB2	2.39	0.47
1:E:295:GLN:HG2	1:E:306:PRO:HG2	1.96	0.47
2:D:2:LEU:HB3	6:F:320:HOH:O	2.14	0.47
1:A:311:ARG:NE	2:B:97:GLU:OE1	2.36	0.47
1:A:66:PRO:HA	1:A:69:MET:CE	2.44	0.47
1:A:119:ASN:ND2	1:A:261:ILE:HD11	2.30	0.47
2:B:94:TYR:HE1	2:D:95:GLN:NE2	2.13	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:125:GLN:HG2	6:F:317:HOH:O	2.15	0.47
1:E:188:THR:HG22	6:E:812:HOH:O	2.15	0.46
2:B:44:ALA:O	2:B:48:ILE:HG12	2.15	0.46
1:E:69:MET:HE1	1:E:87:LEU:HD21	1.97	0.46
1:E:179:MET:HE2	1:E:179:MET:HB3	1.71	0.46
1:C:311:ARG:HE	2:D:97:GLU:CD	2.23	0.46
1:A:210:ARG:NH2	1:E:216:VAL:O	2.47	0.46
1:C:69:MET:HE1	1:C:87:LEU:HD21	1.97	0.45
2:D:54:ARG:NE	6:D:305:HOH:O	2.33	0.45
1:E:41:GLU:OE1	1:E:313:SER:OG	2.33	0.45
1:E:22:ASN:ND2	6:E:713:HOH:O	2.49	0.45
2:B:134:GLY:HA2	2:D:124:LYS:HD3	1.97	0.45
2:B:153:ARG:NH2	6:B:307:HOH:O	2.50	0.45
2:F:21:TRP:H	2:F:41:THR:HG22	1.82	0.44
1:C:17:HIS:N	2:D:115:MET:HE1	2.31	0.44
1:E:307:LYS:HG3	2:F:92:TRP:CE2	2.53	0.44
1:C:226:LEU:O	1:C:229:ARG:NH2	2.51	0.44
1:E:158:SER:HB2	1:E:159:GLN:HG3	2.00	0.44
1:A:129:SER:O	1:A:157:LYS:HE2	2.17	0.44
2:D:2:LEU:C	2:D:2:LEU:HD23	2.42	0.44
2:D:22:TYR:H	2:D:41:THR:CG2	2.31	0.44
2:F:48:ILE:HD11	2:F:107:THR:HG23	1.99	0.44
1:A:17:HIS:HB3	2:B:115:MET:HE1	2.00	0.44
1:A:284:ARG:HH11	2:B:69:GLU:HG3	1.82	0.44
1:A:50:ARG:CZ	1:A:273:PRO:HG2	2.47	0.44
2:B:144:CYS:HB3	2:B:149:MET:HE2	2.00	0.44
2:D:54:ARG:NH2	6:D:305:HOH:O	2.40	0.44
1:C:38:ASN:C	1:C:318:THR:HG22	2.43	0.43
1:E:50:ARG:NE	1:E:275:ASP:OD2	2.39	0.43
1:C:316:LEU:HD11	2:D:103:GLU:HG2	1.99	0.43
1:E:54:LYS:HD2	1:E:276(A):ASN:O	2.17	0.43
1:E:66:PRO:HA	1:E:69:MET:HE3	2.00	0.43
1:C:100:GLY:HA3	1:C:230:ILE:O	2.18	0.43
1:C:176:HIS:HD2	1:C:259:LYS:HG3	1.84	0.43
1:A:27:LYS:HB2	1:A:27:LYS:HE3	1.75	0.43
2:D:21:TRP:H	2:D:41:THR:HG21	1.84	0.43
1:A:140:MET:HB3	1:A:140:MET:HE3	1.79	0.43
2:B:53:ASN:C	2:B:57:GLU:HG2	2.41	0.43
1:C:71:ILE:O	1:C:148:TYR:HB3	2.19	0.43
1:E:223:VAL:HG23	1:E:223:VAL:O	2.19	0.43
2:B:51:LYS:HD3	2:B:103:GLU:HG3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:66:PRO:HA	1:E:69:MET:CE	2.49	0.43
1:A:161:PHE:HB3	1:A:248:ASN:O	2.19	0.42
2:D:51:LYS:NZ	2:D:107:THR:OG1	2.31	0.42
1:E:293:PRO:HG2	1:E:294:PHE:CD2	2.55	0.42
1:E:158(A):LYS:HG2	1:E:193:THR:HG23	2.02	0.42
1:C:295:GLN:OE1	1:C:297:LEU:N	2.53	0.42
1:C:291:ARG:HG3	2:D:56:VAL:HB	2.02	0.42
2:D:133:ASP:CG	2:D:137:CYS:HB2	2.43	0.42
1:A:325:GLU:HG3	2:B:15:GLU:HG3	2.01	0.42
1:E:25:ILE:HG23	1:E:33:GLN:HA	2.02	0.42
1:E:53:MET:HE1	1:E:86:THR:HG21	2.02	0.42
1:C:19:ALA:HB2	2:D:13:GLY:HA3	2.02	0.41
1:C:65:HIS:CE1	1:C:67:ILE:HG12	2.55	0.41
1:A:53:MET:HE1	1:A:86:THR:HG21	2.02	0.41
1:C:325:GLU:OE2	1:C:326:LEU:N	2.52	0.41
1:E:108:LEU:HB2	1:E:234:TRP:CE2	2.56	0.41
1:A:61:LEU:HD11	1:A:69:MET:HE2	2.02	0.41
2:B:2:LEU:HD23	2:B:2:LEU:C	2.46	0.41
1:A:50:ARG:HD3	1:A:50:ARG:HA	1.93	0.41
2:F:37:ASP:O	2:F:41:THR:HG23	2.20	0.41
2:B:83:TRP:CH2	2:D:85:LYS:HG2	2.55	0.41
2:B:120:GLU:OE1	2:B:123:ARG:NH1	2.54	0.40
1:A:307:LYS:HG3	2:B:92:TRP:CE2	2.57	0.40
1:A:98:TYR:CD1	1:A:99:PRO:HD2	2.57	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:C:740:HOH:O	6:E:721:HOH:O[1_454]	2.17	0.03

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	316/323 (98%)	309 (98%)	7 (2%)	0	100	100
1	C	317/323 (98%)	308 (97%)	9 (3%)	0	100	100
1	E	316/323 (98%)	310 (98%)	6 (2%)	0	100	100
2	B	169/180 (94%)	162 (96%)	7 (4%)	0	100	100
2	D	165/180 (92%)	162 (98%)	3 (2%)	0	100	100
2	F	163/180 (91%)	157 (96%)	6 (4%)	0	100	100
All	All	1446/1509 (96%)	1408 (97%)	38 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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%)	264 (98%)	6 (2%)	45	66
1	C	271/273 (99%)	263 (97%)	8 (3%)	36	56
1	E	270/273 (99%)	267 (99%)	3 (1%)	65	81
2	B	146/154 (95%)	144 (99%)	2 (1%)	59	77
2	D	144/154 (94%)	142 (99%)	2 (1%)	59	77
2	F	145/154 (94%)	138 (95%)	7 (5%)	23	38
All	All	1246/1281 (97%)	1218 (98%)	28 (2%)	45	66

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

Mol	Chain	Res	Type
1	A	18	HIS
1	A	27	LYS
1	A	48	ILE
1	A	57	LYS
1	A	157	LYS

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Mol	Chain	Res	Type
1	A	229	ARG
2	B	155	ASN
2	B	164	GLU
1	C	18	HIS
1	C	46	THR
1	C	140	MET
1	C	157	LYS
1	C	189	GLN
1	C	216	VAL
1	C	229	ARG
1	C	302	VAL
2	D	54	ARG
2	D	95	GLN
1	E	18	HIS
1	E	71	ILE
1	E	280	LYS
2	F	26	HIS
2	F	58	LYS
2	F	72	GLU
2	F	103	GLU
2	F	123	ARG
2	F	150	GLU
2	F	172	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	17	HIS
1	A	119	ASN
1	A	212	ASN
1	A	238	GLN
1	A	276	ASN
1	A	310	ASN
2	B	30	GLN
2	B	79	ASN
2	B	95	GLN
2	B	105	GLN
1	C	18	HIS
1	C	110	GLN
1	C	310	ASN
2	D	79	ASN
2	D	95	GLN

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Mol	Chain	Res	Type
2	D	125	GLN
2	D	161	GLN
1	E	110	GLN
1	E	163	GLN
1	E	166	ASN
1	E	222	GLN
1	E	238	GLN
2	F	12	ASN
2	F	26	HIS
2	F	30	GLN
2	F	47	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

6 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	1.06	2 (14%)	17,19,21	1.46	3 (17%)
3	NAG	G	2	3	14,14,15	0.21	0	17,19,21	0.55	0
3	BMA	G	3	3	11,11,12	1.44	2 (18%)	15,15,17	0.85	1 (6%)
3	MAN	G	4	3	11,11,12	1.54	3 (27%)	15,15,17	1.18	2 (13%)
4	NAG	H	1	1,4	14,14,15	1.17	2 (14%)	17,19,21	1.20	2 (11%)
4	NAG	H	2	4	14,14,15	0.33	0	17,19,21	0.77	0

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

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

All (9) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	G	4	MAN	O5-C1	3.09	1.48	1.43
3	G	3	BMA	O5-C1	2.87	1.48	1.43
3	G	3	BMA	O5-C5	2.78	1.48	1.43
3	G	4	MAN	O5-C5	2.49	1.48	1.43
3	G	1	NAG	O5-C1	-2.33	1.39	1.43
4	H	1	NAG	O5-C1	-2.27	1.39	1.43
4	H	1	NAG	C2-N2	-2.23	1.42	1.46
3	G	1	NAG	O7-C7	-2.15	1.18	1.23
3	G	4	MAN	O3-C3	2.14	1.48	1.43

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	1	NAG	C1-C2-N2	-3.59	104.77	110.43
3	G	1	NAG	C6-C5-C4	-2.90	105.91	113.02
3	G	4	MAN	C1-O5-C5	2.75	115.87	112.19
3	G	1	NAG	O3-C3-C2	2.50	114.59	109.40
3	G	4	MAN	C1-C2-C3	2.20	112.85	109.64
3	G	3	BMA	C1-O5-C5	2.15	115.07	112.19
4	H	1	NAG	O5-C5-C6	-2.03	103.71	107.66
4	H	1	NAG	O3-C3-C2	2.02	113.60	109.40

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	H	1	NAG	O5-C5-C6-O6

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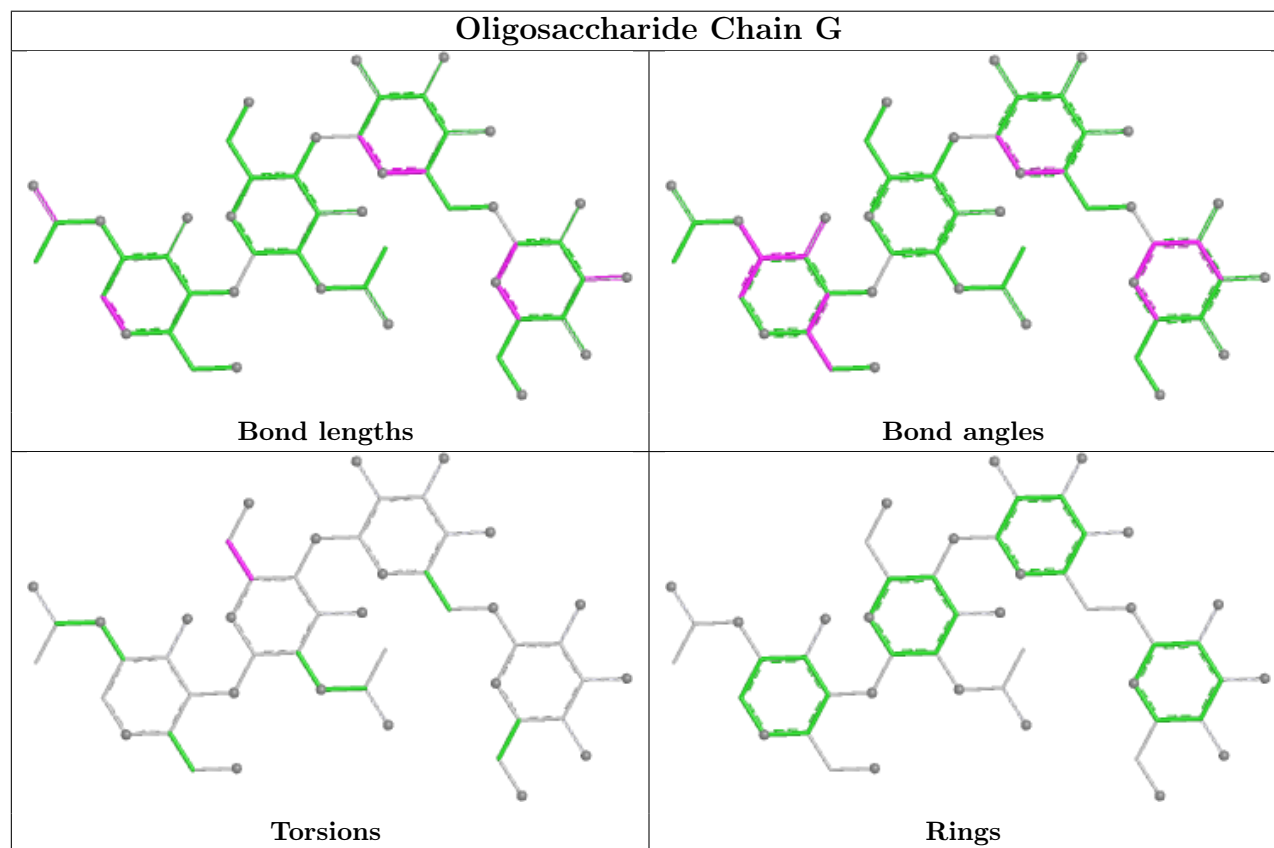
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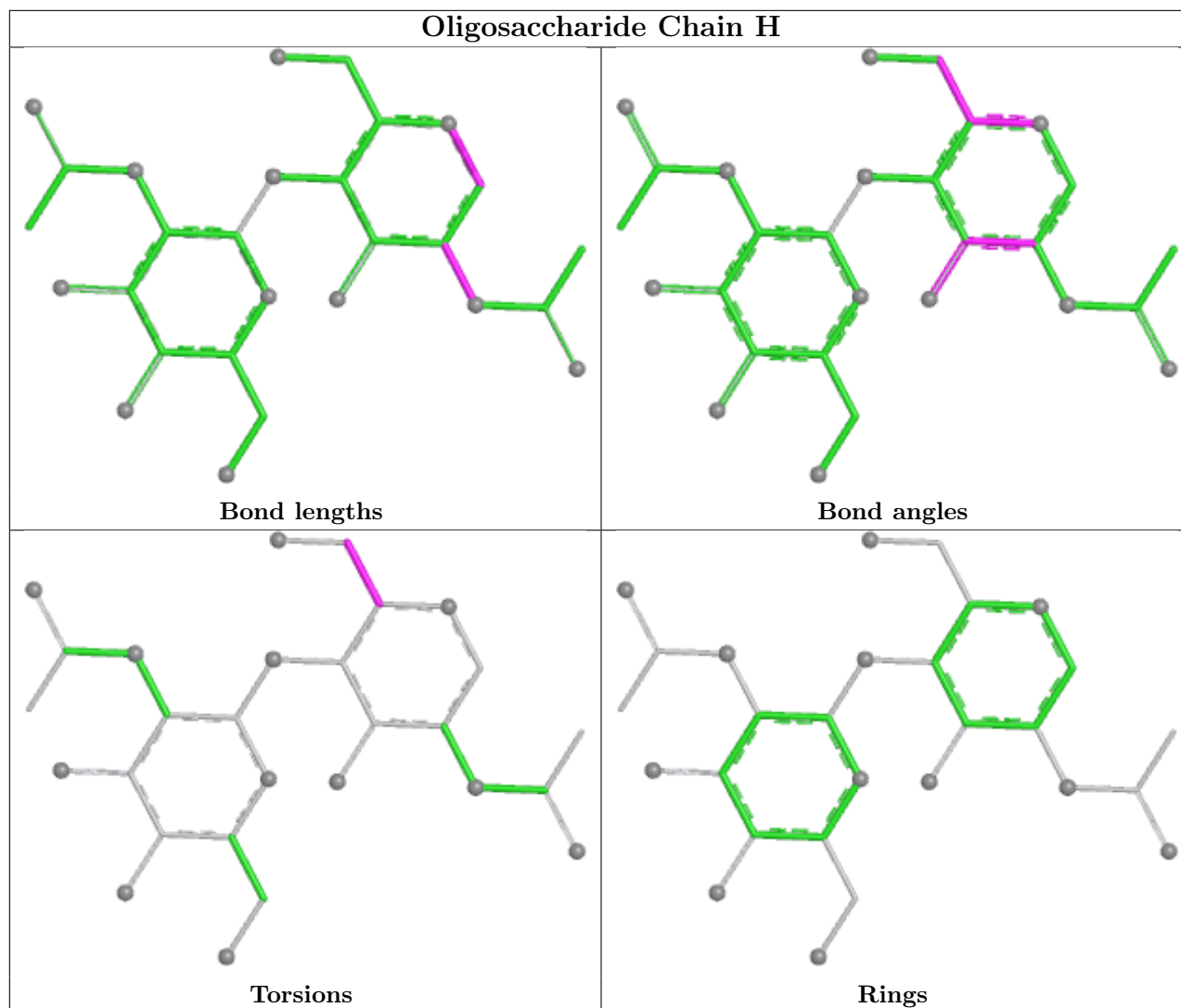
Mol	Chain	Res	Type	Atoms
4	H	1	NAG	C4-C5-C6-O6
3	G	2	NAG	O5-C5-C6-O6
3	G	2	NAG	C4-C5-C6-O6

There are no ring outliers.

No monomer is involved in short contacts.

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





5.6 Ligand geometry [i](#)

5 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
5	NAG	A	400	1	14,14,15	0.64	1 (7%)	17,19,21	1.34	1 (5%)
5	NAG	A	401	1	14,14,15	0.74	1 (7%)	17,19,21	0.88	1 (5%)
5	NAG	D	201	2	14,14,15	0.79	1 (7%)	17,19,21	3.01	5 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	NAG	F	201	2	14,14,15	0.79	1 (7%)	17,19,21	3.02	5 (29%)
5	NAG	B	201	2	14,14,15	0.64	1 (7%)	17,19,21	1.35	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
5	NAG	A	400	1	-	4/6/23/26	0/1/1/1
5	NAG	A	401	1	-	0/6/23/26	0/1/1/1
5	NAG	D	201	2	-	0/6/23/26	0/1/1/1
5	NAG	F	201	2	-	0/6/23/26	0/1/1/1
5	NAG	B	201	2	-	4/6/23/26	0/1/1/1

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	401	NAG	O5-C1	-2.29	1.39	1.43
5	B	201	NAG	O5-C1	-2.14	1.40	1.43
5	A	400	NAG	O5-C1	-2.12	1.40	1.43
5	D	201	NAG	O5-C1	-2.12	1.40	1.43
5	F	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(°)
5	F	201	NAG	O3-C3-C2	-8.18	92.42	109.40
5	D	201	NAG	O3-C3-C2	-8.16	92.44	109.40
5	F	201	NAG	C1-C2-N2	-6.74	99.81	110.43
5	D	201	NAG	C1-C2-N2	-6.73	99.83	110.43
5	D	201	NAG	O4-C4-C5	4.55	120.52	109.32
5	F	201	NAG	O4-C4-C5	4.54	120.50	109.32
5	A	400	NAG	O4-C4-C3	-3.19	102.85	110.38
5	B	201	NAG	O4-C4-C3	-3.19	102.86	110.38
5	D	201	NAG	C6-C5-C4	2.15	118.30	113.02
5	F	201	NAG	C6-C5-C4	2.15	118.29	113.02
5	D	201	NAG	C2-N2-C7	-2.07	120.13	122.90
5	F	201	NAG	C2-N2-C7	-2.07	120.13	122.90
5	A	401	NAG	C1-C2-N2	-2.02	107.26	110.43

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
5	A	400	NAG	C8-C7-N2-C2
5	A	400	NAG	O7-C7-N2-C2
5	B	201	NAG	C8-C7-N2-C2
5	B	201	NAG	O7-C7-N2-C2
5	A	400	NAG	C4-C5-C6-O6
5	B	201	NAG	C4-C5-C6-O6
5	A	400	NAG	O5-C5-C6-O6
5	B	201	NAG	O5-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	F	201	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	318/323 (98%)	-0.46	2 (0%) 85 84	11, 21, 36, 58	0
1	C	319/323 (98%)	-0.48	1 (0%) 90 88	12, 21, 36, 59	0
1	E	318/323 (98%)	-0.29	9 (2%) 55 51	12, 22, 41, 58	0
2	B	171/180 (95%)	-0.03	5 (2%) 53 50	14, 30, 49, 55	0
2	D	169/180 (93%)	-0.26	1 (0%) 85 84	15, 27, 41, 56	0
2	F	169/180 (93%)	-0.02	10 (5%) 28 24	14, 29, 47, 59	0
All	All	1464/1509 (97%)	-0.30	28 (1%) 66 63	11, 24, 44, 59	0

All (28) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	220	ARG	5.4
2	D	58	LYS	4.2
2	F	26	HIS	3.4
2	F	168	LEU	3.1
1	C	326	LEU	3.1
2	F	172	ASN	3.0
1	E	223	VAL	3.0
2	B	154	ASN	2.9
2	F	32	THR	2.8
2	B	57	GLU	2.7
1	E	22	ASN	2.6
2	F	2	LEU	2.5
2	F	4	GLY	2.5
2	F	31	GLY	2.4
1	E	221	PRO	2.4
2	F	58	LYS	2.4
2	F	60	ASN	2.4
2	F	19	ASP	2.3
1	E	248	ASN	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	21	ALA	2.2
2	B	60	ASN	2.2
1	E	32	GLU	2.2
1	E	325	GLU	2.2
2	B	59	THR	2.2
1	E	285	GLY	2.2
2	B	145	ASP	2.1
1	A	22	ASN	2.1
1	E	291	ARG	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 [\(i\)](#)

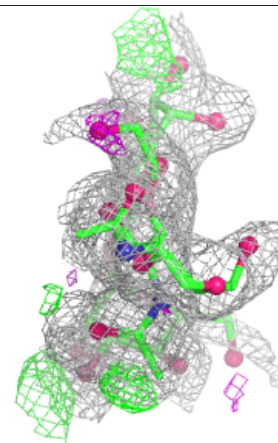
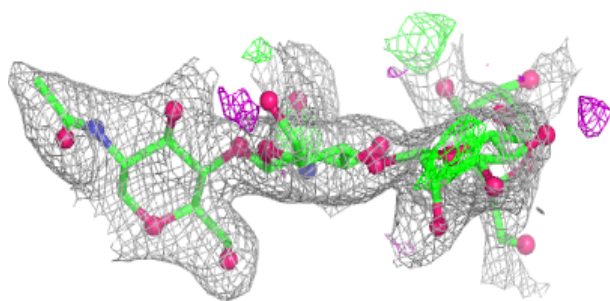
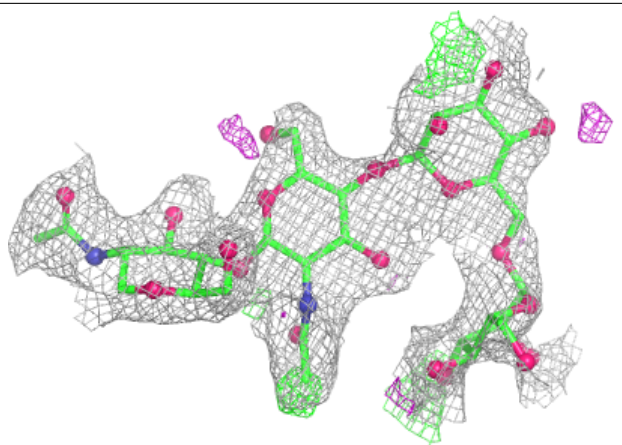
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

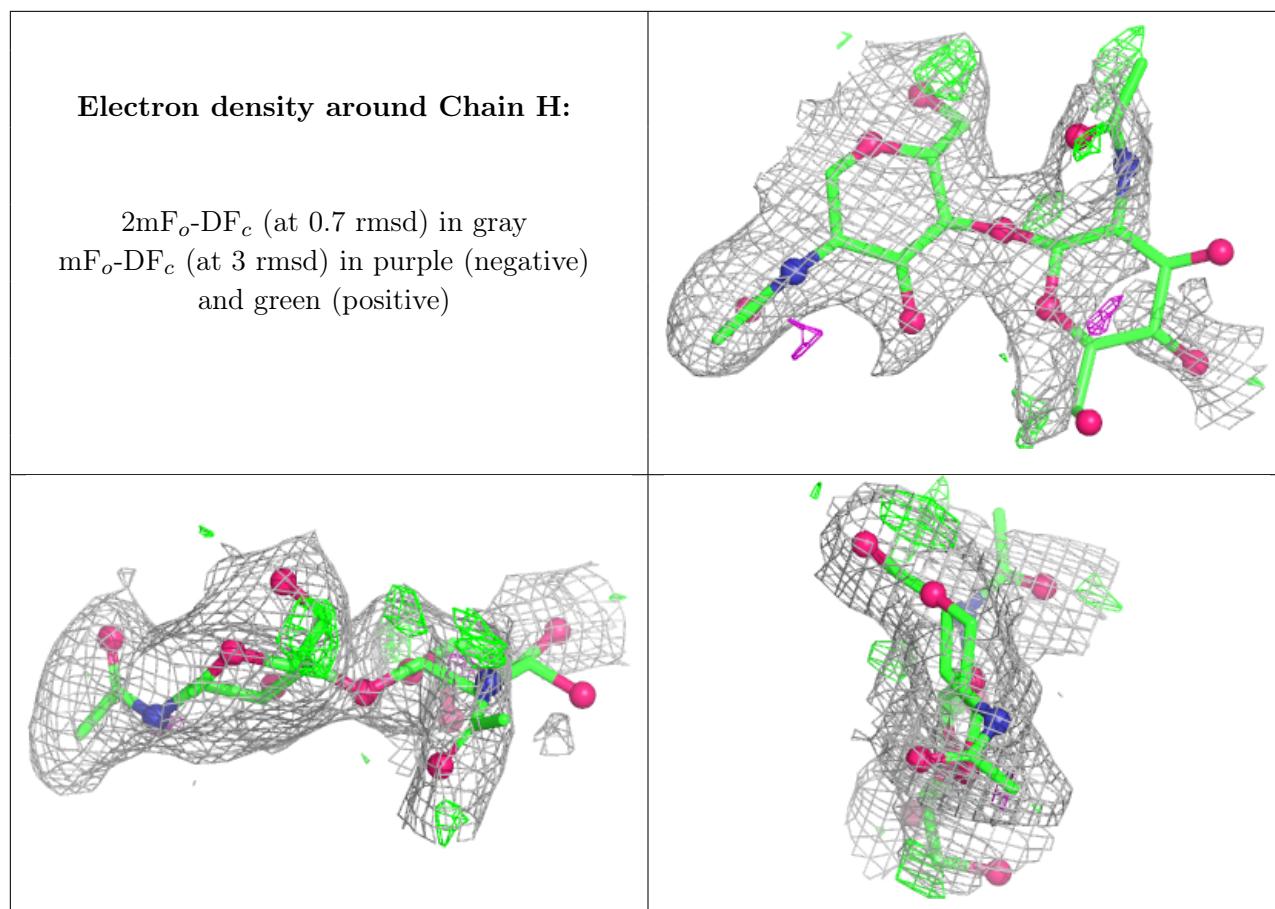
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	H	2	14/15	0.52	0.25	63,75,91,91	0
3	MAN	G	4	11/12	0.57	0.22	70,83,87,88	0
3	BMA	G	3	11/12	0.77	0.17	56,63,71,77	0
3	NAG	G	2	14/15	0.79	0.16	34,43,57,60	0
4	NAG	H	1	14/15	0.81	0.14	26,44,59,67	0
3	NAG	G	1	14/15	0.90	0.10	28,34,41,43	0

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)





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NAG	A	400	14/15	0.54	0.23	51,60,68,72	0
5	NAG	B	201	14/15	0.60	0.25	47,58,68,81	0
5	NAG	F	201	14/15	0.64	0.21	47,57,63,70	0
5	NAG	A	401	14/15	0.65	0.21	41,50,61,67	0
5	NAG	D	201	14/15	0.72	0.17	47,54,60,63	0

6.5 Other polymers [i](#)

There are no such residues in this entry.