



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 12:04 PM UTC

PDB ID : 4KDM / pdb_00004kdm
Title : Crystal structure of the hemagglutinin of ferret-transmissible H5N1 virus
Authors : Lu, X.; Shi, Y.; Zhang, W.; Zhang, Y.; Qi, J.; Gao, G.F.
Deposited on : 2013-04-25
Resolution : 2.50 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)
Xtrriage (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

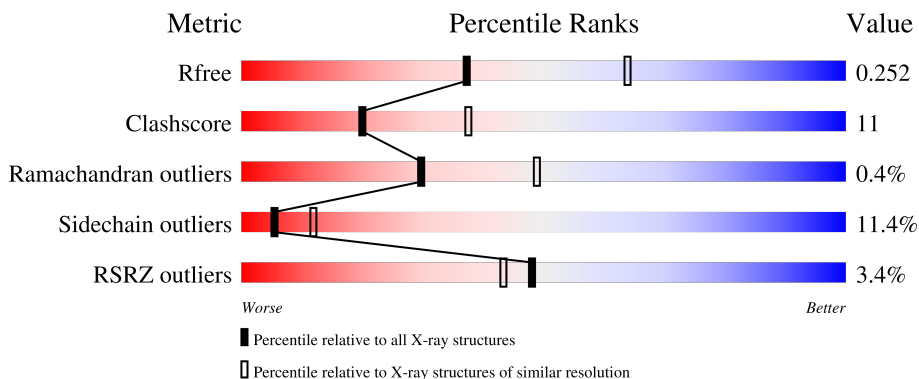
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION



The reported resolution of this entry is 2.50 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.



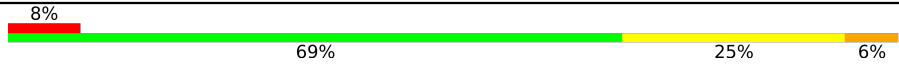

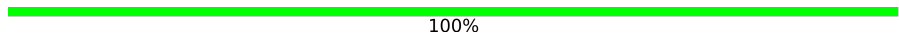
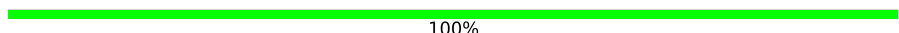
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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 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	322	
1	C	322	
1	E	322	
2	B	175	
2	D	175	

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Mol	Chain	Length	Quality of chain
2	F	175	 8% 69% 25% 6%
3	G	2	 50% 50%
3	H	2	 100%
3	I	2	 100%

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 12221 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.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	322	2559	1621	440	483	15	0	0	0
1	C	322	2559	1621	440	483	15	0	0	0
1	E	322	2559	1621	440	483	15	0	0	0

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	4	GLN	-	expression tag	UNP Q6DQ33
A	158	ASP	ASN	engineered mutation	UNP Q6DQ33
A	224	LYS	ASN	engineered mutation	UNP Q6DQ33
A	226	LEU	GLN	engineered mutation	UNP Q6DQ33
A	319	ILE	THR	engineered mutation	UNP Q6DQ33
C	4	GLN	-	expression tag	UNP Q6DQ33
C	158	ASP	ASN	engineered mutation	UNP Q6DQ33
C	224	LYS	ASN	engineered mutation	UNP Q6DQ33
C	226	LEU	GLN	engineered mutation	UNP Q6DQ33
C	319	ILE	THR	engineered mutation	UNP Q6DQ33
E	4	GLN	-	expression tag	UNP Q6DQ33
E	158	ASP	ASN	engineered mutation	UNP Q6DQ33
E	224	LYS	ASN	engineered mutation	UNP Q6DQ33
E	226	LEU	GLN	engineered mutation	UNP Q6DQ33
E	319	ILE	THR	engineered mutation	UNP Q6DQ33

- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	175	1416	880	246	282	8	0	0	0

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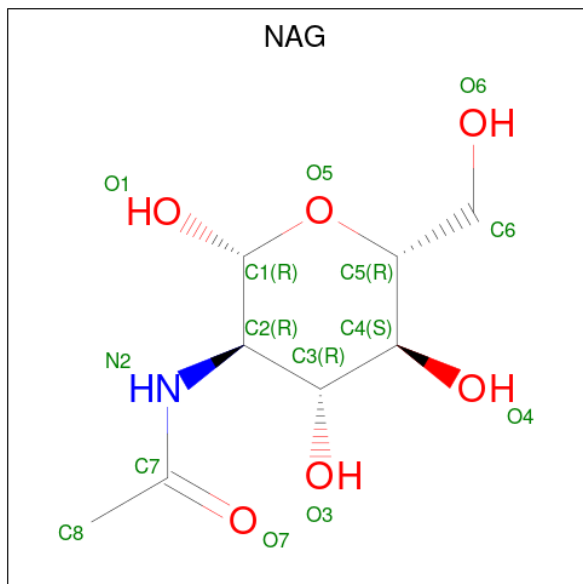
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	D	175	Total	C	N	O	S	0	0	0
			1416	880	246	282	8			
2	F	175	Total	C	N	O	S	0	0	0
			1416	880	246	282	8			

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

- Molecule 4 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		
4	A	1	Total	C	N	O	0	0
			14	8	1	5		

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

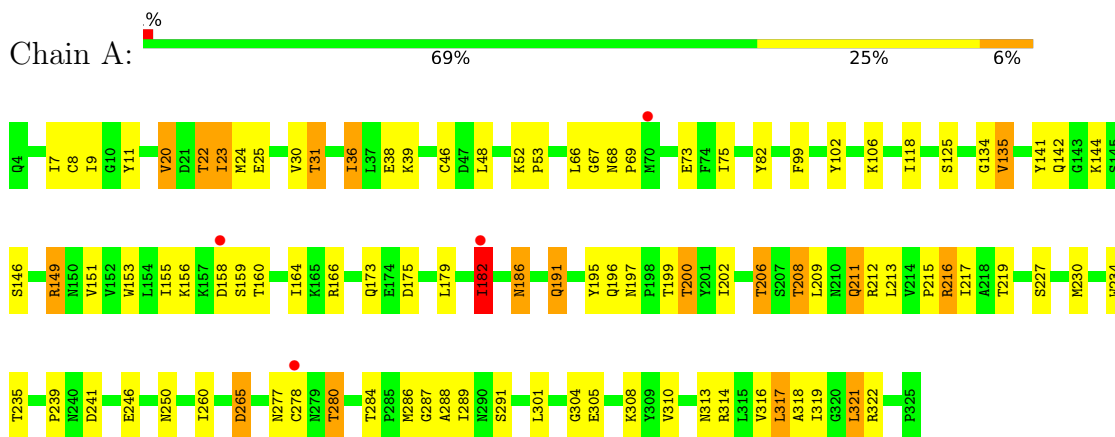
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	53	Total	O	0	0
			53	53		
5	B	13	Total	O	0	0
			13	13		
5	C	42	Total	O	0	0
			42	42		
5	D	17	Total	O	0	0
			17	17		
5	E	33	Total	O	0	0
			33	33		
5	F	12	Total	O	0	0
			12	12		

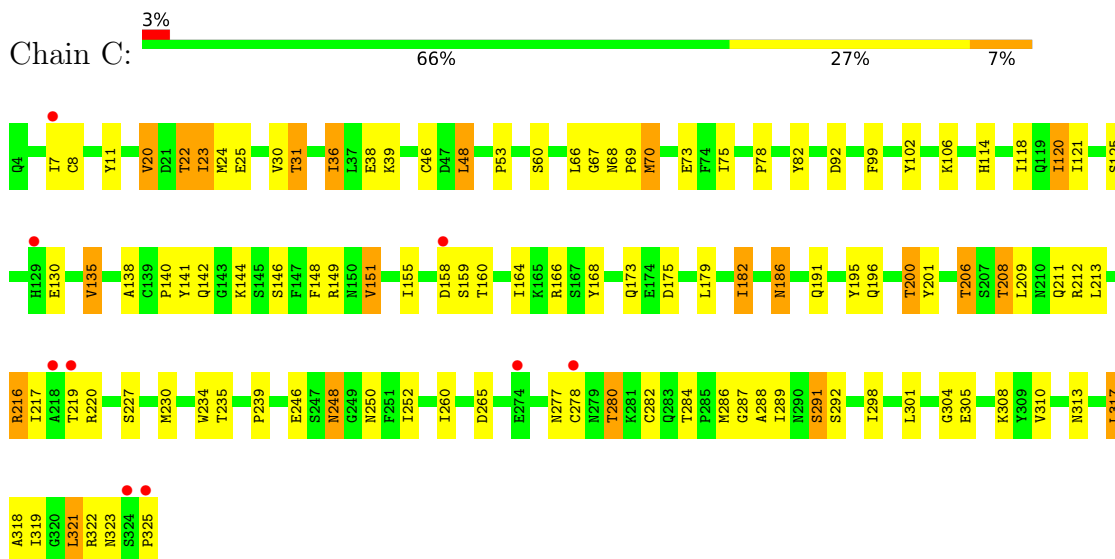
3 Residue-property plots i

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

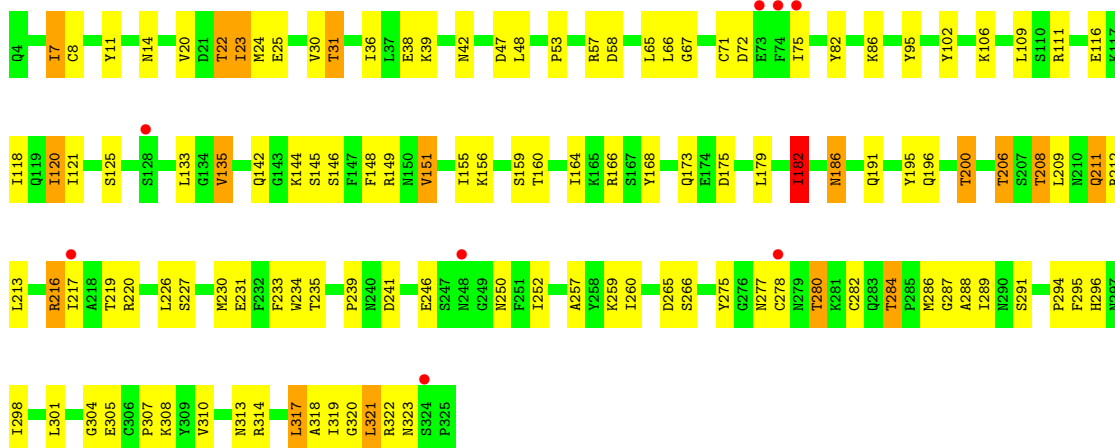


- Molecule 1: Hemagglutinin

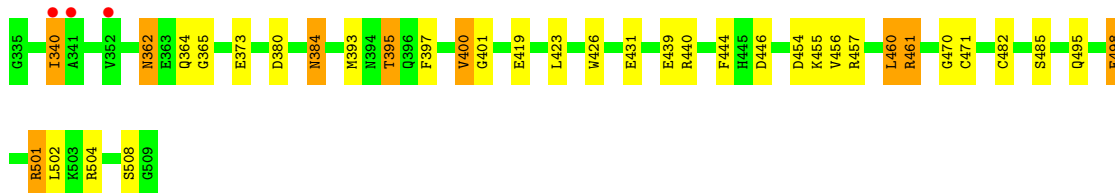
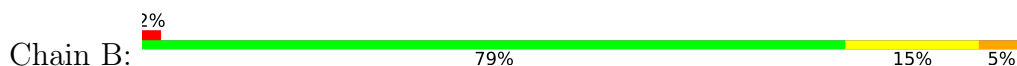


- Molecule 1: Hemagglutinin

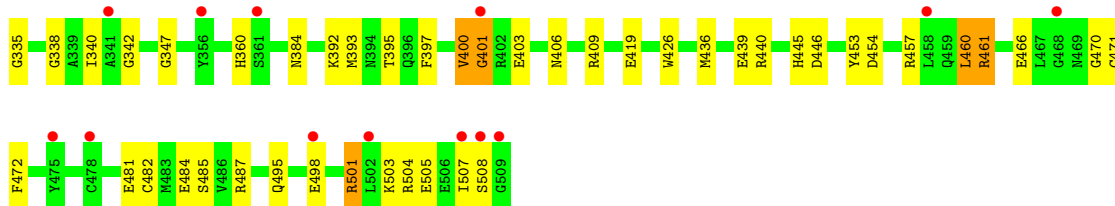
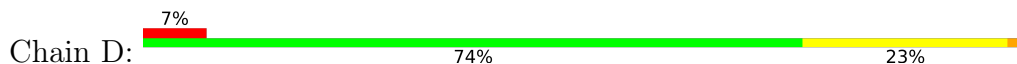




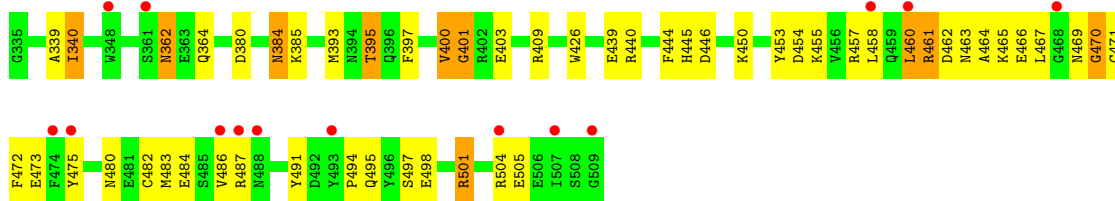
- Molecule 2: Hemagglutinin



- Molecule 2: Hemagglutinin



- Molecule 2: Hemagglutinin



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

Chain G:  50% 50%

HA61
HA62

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

Chain H:  100%

HA61
HA62

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

Chain I:  100%

HA61
HA62

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.76Å 243.85Å 71.61Å 90.00° 110.01° 90.00°	Depositor
Resolution (Å)	35.07 – 2.50 35.07 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.4 (35.07-2.50) 99.4 (35.07-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.60 (at 2.51Å)	Xtrriage
Refinement program	PHENIX (phenix.refine: 1.8_1069)	Depositor
R, R_{free}	0.215 , 0.251 0.217 , 0.252	Depositor DCC
R_{free} test set	3723 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	54.8	Xtrriage
Anisotropy	0.293	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 46.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	12221	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.41% 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

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	0.34	0/2621	0.83	3/3558 (0.1%)
1	C	0.33	0/2621	0.82	5/3558 (0.1%)
1	E	0.34	0/2621	0.84	3/3558 (0.1%)
2	B	0.42	0/1443	0.83	4/1939 (0.2%)
2	D	0.41	0/1443	0.80	2/1939 (0.1%)
2	F	0.39	0/1443	0.82	3/1939 (0.2%)
All	All	0.36	0/12192	0.83	20/16491 (0.1%)

There are no bond length outliers.

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	461	ARG	CB-CA-C	-8.07	107.26	116.54
2	F	470	GLY	N-CA-C	-7.58	104.74	115.43
1	E	71	CYS	N-CA-C	-7.13	100.45	110.50
2	B	470	GLY	N-CA-C	-7.04	105.50	115.43
2	B	460	LEU	CB-CA-C	7.01	123.75	109.67
1	E	66	LEU	N-CA-C	-6.88	101.94	110.41
2	F	461	ARG	CB-CA-C	-6.60	108.95	116.54
1	E	182	ILE	CB-CA-C	-6.56	100.72	110.82
2	B	461	ARG	CB-CA-C	-6.53	109.05	116.63
1	A	66	LEU	N-CA-C	-6.38	102.57	110.41
2	F	340	ILE	N-CA-C	5.97	116.75	110.72
1	C	248	ASN	N-CA-C	-5.87	106.46	113.97
1	C	292	SER	N-CA-C	-5.69	106.46	113.41
1	C	158	ASP	N-CA-C	-5.68	99.68	108.31
2	B	340	ILE	N-CA-C	5.62	116.39	110.72
1	A	158	ASP	CB-CA-C	-5.60	110.13	116.63
1	C	66	LEU	N-CA-C	-5.55	103.58	110.41
2	D	460	LEU	CB-CA-C	5.19	120.09	109.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	182	ILE	CB-CA-C	-5.17	102.86	110.82
1	C	291	SER	N-CA-C	5.11	116.75	108.41

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2559	0	2513	66	0
1	C	2559	0	2513	63	1
1	E	2559	0	2513	73	1
2	B	1416	0	1319	28	0
2	D	1416	0	1319	35	0
2	F	1416	0	1319	46	0
3	G	28	0	25	1	0
3	H	28	0	25	0	0
3	I	28	0	25	0	0
4	A	14	0	13	0	0
4	C	14	0	13	0	0
4	E	14	0	13	0	0
5	A	53	0	0	8	0
5	B	13	0	0	2	0
5	C	42	0	0	4	0
5	D	17	0	0	6	0
5	E	33	0	0	6	0
5	F	12	0	0	2	0
All	All	12221	0	11610	270	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 11.

All (270) 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:206:THR:HG22	1:E:208:THR:H	1.37	0.89
5:A:753:HOH:O	3:G:1:NAG:O3	1.88	0.89
1:E:58:ASP:OD1	5:E:726:HOH:O	1.90	0.88
1:E:38:GLU:OE1	5:E:725:HOH:O	1.92	0.87
1:C:138:ALA:O	5:C:705:HOH:O	1.94	0.84
2:D:461:ARG:NH2	2:F:466:GLU:O	2.10	0.84
1:A:206:THR:HG22	1:A:208:THR:H	1.42	0.83
1:A:284:THR:HG22	1:A:286:MET:H	1.44	0.82
1:A:182:ILE:HD11	1:A:213:LEU:HD13	1.63	0.81
1:E:318:ALA:O	5:E:730:HOH:O	1.98	0.81
1:A:38:GLU:OE1	5:A:727:HOH:O	1.98	0.80
1:A:199:THR:OG1	5:A:752:HOH:O	1.99	0.80
1:E:57:ARG:NH1	1:E:72:ASP:OD1	2.15	0.79
1:C:206:THR:HG22	1:C:208:THR:H	1.48	0.78
1:E:67:GLY:HA3	1:E:149:ARG:HG2	1.66	0.77
1:C:135:VAL:HG22	1:C:146:SER:HA	1.67	0.77
1:E:284:THR:HG22	1:E:286:MET:H	1.50	0.76
1:C:38:GLU:OE1	5:C:734:HOH:O	2.03	0.76
2:B:419:GLU:O	5:B:608:HOH:O	2.04	0.75
1:C:186:ASN:OD1	1:C:186:ASN:N	2.20	0.74
1:C:284:THR:HG22	1:C:286:MET:H	1.50	0.74
1:E:280:THR:HG21	1:E:288:ALA:HB1	1.69	0.74
2:F:403:GLU:OE2	5:F:604:HOH:O	2.05	0.74
1:E:186:ASN:OD1	1:E:186:ASN:N	2.21	0.74
1:A:22:THR:HG22	1:A:24:MET:H	1.53	0.73
1:A:135:VAL:HG22	1:A:146:SER:HA	1.70	0.73
1:A:186:ASN:N	1:A:186:ASN:OD1	2.19	0.72
1:C:280:THR:HG21	1:C:288:ALA:HB1	1.72	0.72
2:D:347:GLY:O	5:D:607:HOH:O	2.08	0.71
1:A:39:LYS:NZ	1:A:313:ASN:O	2.21	0.70
1:E:25:GLU:OE2	1:E:322:ARG:NH2	2.22	0.70
2:D:453:TYR:HE1	2:D:470:GLY:HA2	1.56	0.70
1:E:22:THR:HG22	1:E:24:MET:H	1.58	0.69
1:C:39:LYS:NZ	1:C:313:ASN:O	2.22	0.69
1:C:130:GLU:O	5:C:703:HOH:O	2.11	0.69
1:E:156:LYS:HD2	1:E:196:GLN:HG2	1.75	0.69
1:A:125:SER:OG	1:A:166:ARG:NH2	2.26	0.68
2:F:484:GLU:HG2	2:F:487:ARG:HH12	1.58	0.67
1:A:265:ASP:OD2	2:D:409:ARG:NH2	2.27	0.67
1:C:25:GLU:OE2	1:C:322:ARG:NH2	2.28	0.67
2:F:494:PRO:HA	2:F:497:SER:HB3	1.75	0.67
1:C:182:ILE:HD11	1:C:213:LEU:HD13	1.77	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:301:LEU:HA	2:B:400:VAL:HG22	1.77	0.66
1:C:22:THR:HG22	1:C:24:MET:H	1.61	0.65
1:C:216:ARG:O	1:C:220:ARG:NH2	2.30	0.65
2:B:362:ASN:OD1	2:B:364:GLN:NE2	2.30	0.65
1:A:25:GLU:OE2	1:A:322:ARG:NH2	2.29	0.64
1:C:284:THR:HB	1:C:287:GLY:O	1.97	0.64
1:E:39:LYS:NZ	1:E:313:ASN:O	2.27	0.64
1:E:179:LEU:HD23	1:E:234:TRP:HB3	1.79	0.64
2:D:453:TYR:CE1	2:D:470:GLY:HA2	2.32	0.64
2:F:466:GLU:HG2	2:F:472:PHE:HE2	1.63	0.63
1:E:8:CYS:HA	2:F:471:CYS:HA	1.79	0.62
1:A:212:ARG:NH1	5:A:732:HOH:O	2.14	0.62
1:A:289:ILE:HG22	1:A:291:SER:HB3	1.81	0.62
2:B:508:SER:OG	2:F:501:ARG:NE	2.28	0.61
1:E:145:SER:OG	5:E:732:HOH:O	2.16	0.61
2:D:440:ARG:NH1	2:F:439:GLU:OE2	2.34	0.61
1:C:67:GLY:HA3	1:C:149:ARG:HB2	1.81	0.60
2:B:431:GLU:OE1	5:B:605:HOH:O	2.17	0.60
2:D:393:MET:O	5:D:617:HOH:O	2.17	0.60
1:C:120:ILE:HG23	1:C:121:ILE:HG13	1.84	0.60
1:C:289:ILE:HG22	1:C:291:SER:HB3	1.83	0.59
2:D:360:HIS:HD2	2:D:487:ARG:HH21	1.50	0.59
1:A:20:VAL:HG11	1:A:318:ALA:HB2	1.82	0.59
1:A:280:THR:HG21	1:A:288:ALA:HB1	1.84	0.59
1:E:125:SER:OG	1:E:166:ARG:NH2	2.35	0.59
2:F:364:GLN:HE22	2:F:480:ASN:H	1.50	0.59
1:A:191:GLN:OE1	1:A:250:ASN:ND2	2.34	0.59
1:A:314:ARG:HG3	1:A:316:VAL:HG23	1.85	0.59
1:A:31:THR:HB	1:A:321:LEU:H	1.68	0.58
1:C:23:ILE:HG22	1:C:24:MET:HG3	1.86	0.58
1:A:179:LEU:HD23	1:A:234:TRP:HB3	1.86	0.57
1:A:206:THR:HG21	5:A:708:HOH:O	2.03	0.57
2:D:335:GLY:N	5:D:601:HOH:O	2.38	0.57
2:D:466:GLU:HG2	2:D:472:PHE:HE2	1.68	0.57
1:A:156:LYS:HD2	1:A:196:GLN:HG2	1.87	0.56
1:E:308:LYS:HZ3	2:F:393:MET:HE2	1.71	0.56
1:C:22:THR:HG23	2:D:439:GLU:HB2	1.86	0.56
2:D:482:CYS:O	2:D:485:SER:OG	2.20	0.56
1:E:23:ILE:HG22	1:E:24:MET:HG3	1.86	0.56
1:A:36:ILE:HG12	1:A:317:LEU:HD22	1.88	0.56
1:A:11:TYR:CZ	2:B:340:ILE:HG23	2.41	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:284:THR:HB	1:A:287:GLY:O	2.06	0.56
2:B:440:ARG:HH21	2:F:440:ARG:HH21	1.54	0.56
1:E:182:ILE:CD1	1:E:213:LEU:HD13	2.36	0.55
1:A:308:LYS:HZ3	2:B:393:MET:HE2	1.71	0.55
1:E:206:THR:HB	1:E:209:LEU:HB3	1.88	0.55
1:A:23:ILE:HG22	1:A:24:MET:HG3	1.89	0.55
2:F:362:ASN:H	2:F:362:ASN:HD22	1.55	0.55
1:E:284:THR:HB	1:E:287:GLY:O	2.06	0.55
2:F:501:ARG:HD3	2:F:505:GLU:HG3	1.88	0.55
1:A:164:ILE:O	1:A:246:GLU:HA	2.07	0.55
1:E:120:ILE:HD13	1:E:257:ALA:HB3	1.89	0.54
1:E:289:ILE:HD11	1:E:298:ILE:HG13	1.89	0.54
1:C:179:LEU:HD23	1:C:234:TRP:HB3	1.90	0.54
1:E:289:ILE:HG22	1:E:291:SER:HB3	1.89	0.54
1:C:301:LEU:HA	2:D:400:VAL:HG22	1.89	0.54
1:E:116:GLU:HB3	1:E:259:LYS:HB2	1.90	0.54
2:F:466:GLU:HG2	2:F:472:PHE:CE2	2.41	0.54
1:C:304:GLY:HA2	2:D:397:PHE:CD1	2.43	0.53
1:A:22:THR:HG23	2:B:439:GLU:HB2	1.90	0.53
1:C:200:THR:HG21	1:C:250:ASN:OD1	2.08	0.53
1:E:11:TYR:CZ	2:F:340:ILE:HG23	2.44	0.53
1:A:73:GLU:OE2	1:A:141:TYR:OH	2.24	0.53
1:A:200:THR:HG21	1:A:250:ASN:OD1	2.09	0.53
1:E:25:GLU:CD	1:E:322:ARG:HH22	2.17	0.53
1:C:200:THR:HA	1:C:248:ASN:HD22	1.74	0.53
1:E:182:ILE:HD11	1:E:213:LEU:HD13	1.91	0.53
2:D:466:GLU:HG2	2:D:472:PHE:CE2	2.44	0.52
2:D:454:ASP:OD1	2:D:457:ARG:NH1	2.42	0.52
2:B:362:ASN:ND2	2:B:365:GLY:O	2.43	0.52
2:B:393:MET:HE1	2:B:426:TRP:CH2	2.43	0.52
1:A:197:ASN:ND2	5:A:712:HOH:O	1.92	0.52
1:E:148:PHE:HB2	1:E:151:VAL:HG12	1.92	0.52
1:E:182:ILE:HD11	1:E:233:PHE:CE1	2.45	0.52
2:D:403:GLU:OE2	5:D:612:HOH:O	2.18	0.52
1:A:102:TYR:CE2	1:A:106:LYS:HD2	2.45	0.51
1:C:102:TYR:CZ	1:C:106:LYS:HD2	2.46	0.51
1:C:280:THR:HB	1:C:282:CYS:H	1.75	0.51
2:B:362:ASN:HD22	2:B:362:ASN:H	1.59	0.51
2:F:482:CYS:O	2:F:486:VAL:HG23	2.11	0.51
1:E:320:GLY:O	2:F:445:HIS:NE2	2.44	0.51
1:E:301:LEU:HA	2:F:400:VAL:HG22	1.93	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:135:VAL:HG22	1:E:146:SER:HA	1.92	0.50
1:E:200:THR:HG21	1:E:250:ASN:OD1	2.10	0.50
1:A:52:LYS:HG2	1:A:53:PRO:HD2	1.93	0.50
1:E:125:SER:C	1:E:166:ARG:HH22	2.20	0.50
1:E:53:PRO:HB3	1:E:82:TYR:CE2	2.47	0.50
1:E:133:LEU:N	5:E:714:HOH:O	2.44	0.50
2:F:463:ASN:ND2	2:F:491:TYR:OH	2.26	0.50
1:E:164:ILE:O	1:E:246:GLU:HA	2.12	0.50
1:E:156:LYS:NZ	1:E:196:GLN:OE1	2.42	0.49
1:A:206:THR:HB	1:A:209:LEU:HB3	1.93	0.49
1:A:212:ARG:HB2	1:C:216:ARG:HG2	1.93	0.49
1:E:216:ARG:O	1:E:220:ARG:NH2	2.45	0.49
2:B:508:SER:HG	2:F:501:ARG:HE	1.58	0.49
1:C:125:SER:CB	1:C:166:ARG:HH22	2.25	0.49
2:F:453:TYR:CE1	2:F:470:GLY:HA2	2.47	0.49
1:A:159:SER:O	1:A:196:GLN:HG3	2.12	0.49
1:C:206:THR:HB	1:C:209:LEU:HB3	1.94	0.49
1:C:164:ILE:O	1:C:246:GLU:HA	2.12	0.49
2:F:364:GLN:HE22	2:F:480:ASN:N	2.09	0.49
2:D:481:GLU:O	2:D:484:GLU:HB3	2.13	0.49
1:C:73:GLU:OE2	1:C:141:TYR:OH	2.19	0.48
2:F:339:ALA:HB2	2:F:450:LYS:HB2	1.94	0.48
2:B:454:ASP:OD1	2:B:457:ARG:NH1	2.46	0.48
2:F:380:ASP:O	2:F:384:ASN:HB2	2.14	0.48
1:A:25:GLU:CD	1:A:322:ARG:HH22	2.20	0.48
1:E:142:GLN:C	1:E:144:LYS:H	2.22	0.48
2:F:454:ASP:OD1	2:F:457:ARG:NH1	2.45	0.48
2:B:482:CYS:O	2:B:485:SER:OG	2.21	0.48
1:C:125:SER:HB2	1:C:166:ARG:HH22	1.79	0.48
1:E:211:GLN:OE1	1:E:213:LEU:HD11	2.14	0.48
1:C:308:LYS:HD2	2:D:426:TRP:CE2	2.49	0.48
2:F:465:LYS:HB2	2:F:475:TYR:CZ	2.48	0.48
1:E:14:ASN:O	1:E:323:ASN:ND2	2.42	0.48
2:F:393:MET:HE1	2:F:426:TRP:CH2	2.48	0.48
2:D:505:GLU:HA	2:D:508:SER:HB3	1.95	0.47
1:E:47:ASP:OD1	1:E:275:TYR:OH	2.24	0.47
1:C:60:SER:OG	1:C:92:ASP:OD1	2.27	0.47
1:C:323:ASN:O	1:C:325:PRO:HD3	2.15	0.47
2:D:501:ARG:HD3	2:D:505:GLU:HG3	1.97	0.47
1:A:200:THR:HG22	1:A:215:PRO:HG3	1.97	0.47
1:C:53:PRO:HB3	1:C:82:TYR:CE2	2.50	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:102:TYR:CE2	1:C:106:LYS:HD2	2.50	0.47
1:E:111:ARG:HH11	1:E:266:SER:HB3	1.79	0.47
1:E:120:ILE:HG13	1:E:168:TYR:CE1	2.50	0.47
1:E:321:LEU:HB3	2:F:445:HIS:CD2	2.50	0.47
1:C:70:MET:SD	1:C:140:PRO:HD2	2.55	0.47
2:D:393:MET:HE1	2:D:426:TRP:CH2	2.50	0.47
2:D:503:LYS:O	2:D:507:ILE:HG12	2.15	0.47
2:F:462:ASP:O	2:F:504:ARG:NH1	2.48	0.47
1:A:99:PHE:HB3	1:A:102:TYR:HB2	1.97	0.46
1:C:308:LYS:HZ3	2:D:393:MET:HE2	1.80	0.46
2:F:460:LEU:HD13	2:F:464:ALA:HB3	1.98	0.46
1:A:211:GLN:NE2	5:A:725:HOH:O	2.37	0.46
1:A:308:LYS:HD2	2:B:426:TRP:CE2	2.51	0.46
1:E:296:HIS:CD2	1:E:307:PRO:HG2	2.51	0.46
1:A:211:GLN:OE1	1:A:213:LEU:HD11	2.14	0.46
1:E:57:ARG:O	1:E:86:LYS:HG3	2.14	0.46
2:F:362:ASN:OD1	2:F:364:GLN:NE2	2.48	0.46
1:E:7:ILE:HG12	2:F:483:MET:HE2	1.98	0.46
1:C:99:PHE:HB3	1:C:102:TYR:HB2	1.98	0.46
1:C:182:ILE:HD11	1:C:213:LEU:CD1	2.44	0.46
1:C:291:SER:HA	5:C:734:HOH:O	2.16	0.46
1:A:102:TYR:CZ	1:A:106:LYS:HD2	2.50	0.46
1:C:36:ILE:HG12	1:C:317:LEU:HD22	1.97	0.45
2:D:360:HIS:CD2	2:D:487:ARG:HH21	2.31	0.45
1:A:159:SER:C	1:A:196:GLN:HG3	2.42	0.45
1:C:120:ILE:HD11	1:C:168:TYR:CZ	2.50	0.45
1:A:23:ILE:HG23	2:F:385:LYS:HG3	1.98	0.45
1:E:120:ILE:HG23	1:E:121:ILE:HG13	1.99	0.45
1:E:280:THR:HB	1:E:282:CYS:H	1.82	0.45
2:F:454:ASP:O	2:F:458:LEU:HG	2.17	0.45
1:A:53:PRO:HB3	1:A:82:TYR:CE2	2.52	0.45
2:B:498:GLU:O	2:B:502:LEU:HG	2.16	0.45
1:E:294:PRO:HG2	1:E:295:PHE:CD1	2.51	0.45
1:A:289:ILE:CG2	1:A:291:SER:HB3	2.46	0.45
1:A:308:LYS:NZ	2:B:393:MET:HE2	2.32	0.45
1:C:175:ASP:OD1	1:C:239:PRO:HD3	2.16	0.45
1:E:31:THR:HB	1:E:321:LEU:H	1.82	0.45
1:E:159:SER:C	1:E:196:GLN:HG3	2.41	0.45
2:F:457:ARG:HD2	2:F:466:GLU:OE2	2.17	0.45
1:C:289:ILE:HD11	1:C:298:ILE:HG13	1.98	0.44
1:A:11:TYR:HB2	1:A:321:LEU:HD11	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:182:ILE:HG23	1:A:202:ILE:HD12	1.99	0.44
1:A:216:ARG:HG2	1:E:212:ARG:HB2	1.99	0.44
1:A:304:GLY:HA2	2:B:397:PHE:CE1	2.53	0.44
2:B:397:PHE:C	2:B:397:PHE:CD2	2.95	0.44
1:E:65:LEU:HD11	1:E:109:LEU:HD11	1.98	0.44
1:E:102:TYR:CE2	1:E:106:LYS:HD2	2.52	0.44
1:E:308:LYS:HD2	2:F:426:TRP:CE2	2.53	0.44
1:A:8:CYS:HA	2:B:471:CYS:HA	2.00	0.44
1:C:67:GLY:HA3	1:C:149:ARG:H	1.82	0.44
1:C:67:GLY:CA	1:C:149:ARG:H	2.30	0.44
2:D:406:ASN:O	5:D:615:HOH:O	2.21	0.44
1:E:175:ASP:OD1	1:E:239:PRO:HD3	2.18	0.44
2:F:484:GLU:HG2	2:F:487:ARG:NH1	2.30	0.44
2:F:467:LEU:HD21	2:F:473:GLU:HB2	2.00	0.44
1:C:20:VAL:HG21	1:C:318:ALA:HB2	2.00	0.43
1:C:142:GLN:C	1:C:144:LYS:H	2.26	0.43
2:F:469:ASN:HB2	2:F:471:CYS:HB2	2.00	0.43
1:E:65:LEU:HD11	1:E:109:LEU:CD1	2.48	0.43
1:A:142:GLN:N	5:A:704:HOH:O	2.20	0.43
1:C:148:PHE:HB2	1:C:151:VAL:HG12	2.00	0.42
1:E:304:GLY:HA2	2:F:397:PHE:CD1	2.53	0.42
2:B:444:PHE:CD1	1:C:24:MET:HE1	2.55	0.42
1:E:22:THR:HG23	2:F:439:GLU:HB2	1.99	0.42
1:A:67:GLY:HA3	1:A:149:ARG:HB2	2.02	0.42
1:C:78:PRO:O	1:C:114:HIS:HA	2.19	0.42
1:E:308:LYS:HD2	2:F:426:TRP:NE1	2.35	0.42
2:F:400:VAL:HA	2:F:401:GLY:HA3	1.68	0.42
1:C:11:TYR:CZ	2:D:340:ILE:HG23	2.54	0.42
1:C:48:LEU:HD23	1:C:48:LEU:HA	1.92	0.42
1:A:9:ILE:HD11	2:B:456:VAL:HG21	2.02	0.42
2:B:440:ARG:HH22	2:D:436:MET:HB3	1.84	0.42
1:C:201:TYR:OH	1:C:246:GLU:OE1	2.26	0.42
1:A:24:MET:HE1	2:F:444:PHE:CE1	2.54	0.42
1:C:68:ASN:HA	1:C:69:PRO:HD3	1.85	0.42
2:B:395:THR:O	2:B:395:THR:OG1	2.38	0.42
1:C:321:LEU:HB3	2:D:445:HIS:CD2	2.54	0.42
2:D:392:LYS:HD3	2:D:392:LYS:HA	1.91	0.42
1:E:206:THR:HG23	1:E:241:ASP:OD2	2.20	0.42
1:C:31:THR:HB	1:C:321:LEU:H	1.85	0.41
2:D:338:GLY:O	2:D:342:GLY:HA3	2.20	0.41
2:D:400:VAL:HA	2:D:401:GLY:HA3	1.70	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:395:THR:O	2:F:395:THR:OG1	2.38	0.41
1:A:134:GLY:HA3	1:A:153:TRP:HB3	2.01	0.41
1:A:206:THR:HG23	1:A:241:ASP:OD2	2.21	0.41
1:C:212:ARG:HB2	1:E:216:ARG:HG2	2.01	0.41
1:E:314:ARG:HE	1:E:314:ARG:HB3	1.57	0.41
1:A:304:GLY:HA2	2:B:397:PHE:CD1	2.55	0.41
1:C:46:CYS:HB3	1:C:278:CYS:C	2.46	0.41
1:E:42:ASN:ND2	1:E:278:CYS:SG	2.94	0.41
2:D:419:GLU:OE2	5:D:613:HOH:O	2.22	0.41
1:E:212:ARG:NE	5:E:703:HOH:O	1.92	0.41
1:A:46:CYS:HB3	1:A:278:CYS:C	2.45	0.41
2:B:501:ARG:HG2	2:B:504:ARG:NH2	2.36	0.41
2:D:501:ARG:HG2	2:D:504:ARG:NH2	2.36	0.41
1:E:95:TYR:CE1	1:E:226:LEU:HD13	2.56	0.41
1:E:182:ILE:HG12	1:E:231:GLU:HB3	2.03	0.40
1:A:142:GLN:C	1:A:144:LYS:H	2.30	0.40
1:C:8:CYS:HA	2:D:471:CYS:HA	2.03	0.40
1:C:159:SER:O	1:C:196:GLN:HG3	2.21	0.40
1:A:175:ASP:OD1	1:A:239:PRO:HD3	2.22	0.40
1:E:317:LEU:HD12	1:E:317:LEU:HA	1.83	0.40
1:A:68:ASN:HA	1:A:69:PRO:HD3	1.91	0.40
2:B:380:ASP:O	2:B:384:ASN:HB2	2.22	0.40
2:F:409:ARG:NH1	5:F:611:HOH:O	2.55	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)
1:C:141:TYR:O	1:E:144:LYS:NZ[1_655]	2.14	0.06

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	320/322 (99%)	297 (93%)	22 (7%)	1 (0%)	36	55
1	C	320/322 (99%)	301 (94%)	18 (6%)	1 (0%)	36	55
1	E	320/322 (99%)	299 (93%)	20 (6%)	1 (0%)	36	55
2	B	173/175 (99%)	164 (95%)	8 (5%)	1 (1%)	21	38
2	D	173/175 (99%)	165 (95%)	7 (4%)	1 (1%)	21	38
2	F	173/175 (99%)	166 (96%)	6 (4%)	1 (1%)	21	38
All	All	1479/1491 (99%)	1392 (94%)	81 (6%)	6 (0%)	30	49

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	277	ASN
1	C	277	ASN
1	E	277	ASN
2	B	401	GLY
2	D	401	GLY
2	F	401	GLY

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	290/290 (100%)	252 (87%)	38 (13%)	4	8
1	C	290/290 (100%)	250 (86%)	40 (14%)	3	7
1	E	290/290 (100%)	250 (86%)	40 (14%)	3	7
2	B	149/149 (100%)	136 (91%)	13 (9%)	9	21
2	D	149/149 (100%)	141 (95%)	8 (5%)	20	41
2	F	149/149 (100%)	138 (93%)	11 (7%)	13	27
All	All	1317/1317 (100%)	1167 (89%)	150 (11%)	5	12

All (150) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	ILE
1	A	20	VAL
1	A	22	THR
1	A	23	ILE
1	A	30	VAL
1	A	31	THR
1	A	36	ILE
1	A	48	LEU
1	A	75	ILE
1	A	118	ILE
1	A	135	VAL
1	A	149	ARG
1	A	151	VAL
1	A	155	ILE
1	A	160	THR
1	A	173	GLN
1	A	182	ILE
1	A	186	ASN
1	A	191	GLN
1	A	195	TYR
1	A	200	THR
1	A	206	THR
1	A	208	THR
1	A	211	GLN
1	A	216	ARG
1	A	217	ILE
1	A	219	THR
1	A	227	SER
1	A	230	MET
1	A	235	THR
1	A	260	ILE
1	A	265	ASP
1	A	280	THR
1	A	305	GLU
1	A	310	VAL
1	A	317	LEU
1	A	319	ILE
1	A	321	LEU
2	B	362	ASN
2	B	373	GLU
2	B	384	ASN
2	B	395	THR
2	B	400	VAL

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Mol	Chain	Res	Type
2	B	423	LEU
2	B	446	ASP
2	B	455	LYS
2	B	460	LEU
2	B	461	ARG
2	B	495	GLN
2	B	498	GLU
2	B	501	ARG
1	C	7	ILE
1	C	20	VAL
1	C	22	THR
1	C	23	ILE
1	C	30	VAL
1	C	31	THR
1	C	36	ILE
1	C	48	LEU
1	C	70	MET
1	C	75	ILE
1	C	118	ILE
1	C	120	ILE
1	C	135	VAL
1	C	151	VAL
1	C	155	ILE
1	C	160	THR
1	C	173	GLN
1	C	182	ILE
1	C	186	ASN
1	C	191	GLN
1	C	195	TYR
1	C	200	THR
1	C	206	THR
1	C	208	THR
1	C	211	GLN
1	C	216	ARG
1	C	217	ILE
1	C	219	THR
1	C	227	SER
1	C	230	MET
1	C	235	THR
1	C	252	ILE
1	C	260	ILE
1	C	265	ASP

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Mol	Chain	Res	Type
1	C	280	THR
1	C	305	GLU
1	C	310	VAL
1	C	317	LEU
1	C	319	ILE
1	C	321	LEU
2	D	384	ASN
2	D	395	THR
2	D	400	VAL
2	D	446	ASP
2	D	460	LEU
2	D	495	GLN
2	D	498	GLU
2	D	501	ARG
1	E	7	ILE
1	E	20	VAL
1	E	22	THR
1	E	23	ILE
1	E	30	VAL
1	E	31	THR
1	E	36	ILE
1	E	48	LEU
1	E	75	ILE
1	E	118	ILE
1	E	120	ILE
1	E	135	VAL
1	E	151	VAL
1	E	155	ILE
1	E	160	THR
1	E	173	GLN
1	E	182	ILE
1	E	186	ASN
1	E	191	GLN
1	E	195	TYR
1	E	200	THR
1	E	206	THR
1	E	208	THR
1	E	211	GLN
1	E	216	ARG
1	E	217	ILE
1	E	219	THR
1	E	227	SER

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Mol	Chain	Res	Type
1	E	230	MET
1	E	235	THR
1	E	252	ILE
1	E	260	ILE
1	E	265	ASP
1	E	280	THR
1	E	284	THR
1	E	305	GLU
1	E	310	VAL
1	E	317	LEU
1	E	319	ILE
1	E	321	LEU
2	F	362	ASN
2	F	384	ASN
2	F	395	THR
2	F	400	VAL
2	F	446	ASP
2	F	455	LYS
2	F	460	LEU
2	F	461	ARG
2	F	495	GLN
2	F	498	GLU
2	F	501	ARG

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

Mol	Chain	Res	Type
1	A	19	GLN
1	A	211	GLN
1	A	323	ASN
2	B	364	GLN
1	C	76	ASN
1	C	248	ASN
1	C	323	ASN
2	D	360	HIS
2	D	376	GLN
1	E	150	ASN
1	E	197	ASN
2	F	364	GLN
2	F	429	ASN
2	F	469	ASN
2	F	480	ASN

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	3,1	14,14,15	0.52	0	17,19,21	0.66	0
3	NAG	G	2	3	14,14,15	0.44	0	17,19,21	0.88	0
3	NAG	H	1	3,1	14,14,15	0.54	0	17,19,21	0.66	0
3	NAG	H	2	3	14,14,15	0.55	0	17,19,21	0.68	0
3	NAG	I	1	3,1	14,14,15	0.52	0	17,19,21	0.80	0
3	NAG	I	2	3	14,14,15	0.42	0	17,19,21	0.88	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	3,1	-	0/6/23/26	0/1/1/1
3	NAG	G	2	3	-	0/6/23/26	0/1/1/1
3	NAG	H	1	3,1	-	0/6/23/26	0/1/1/1
3	NAG	H	2	3	-	1/6/23/26	0/1/1/1
3	NAG	I	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	I	2	3	-	3/6/23/26	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (6) torsion outliers are listed below:

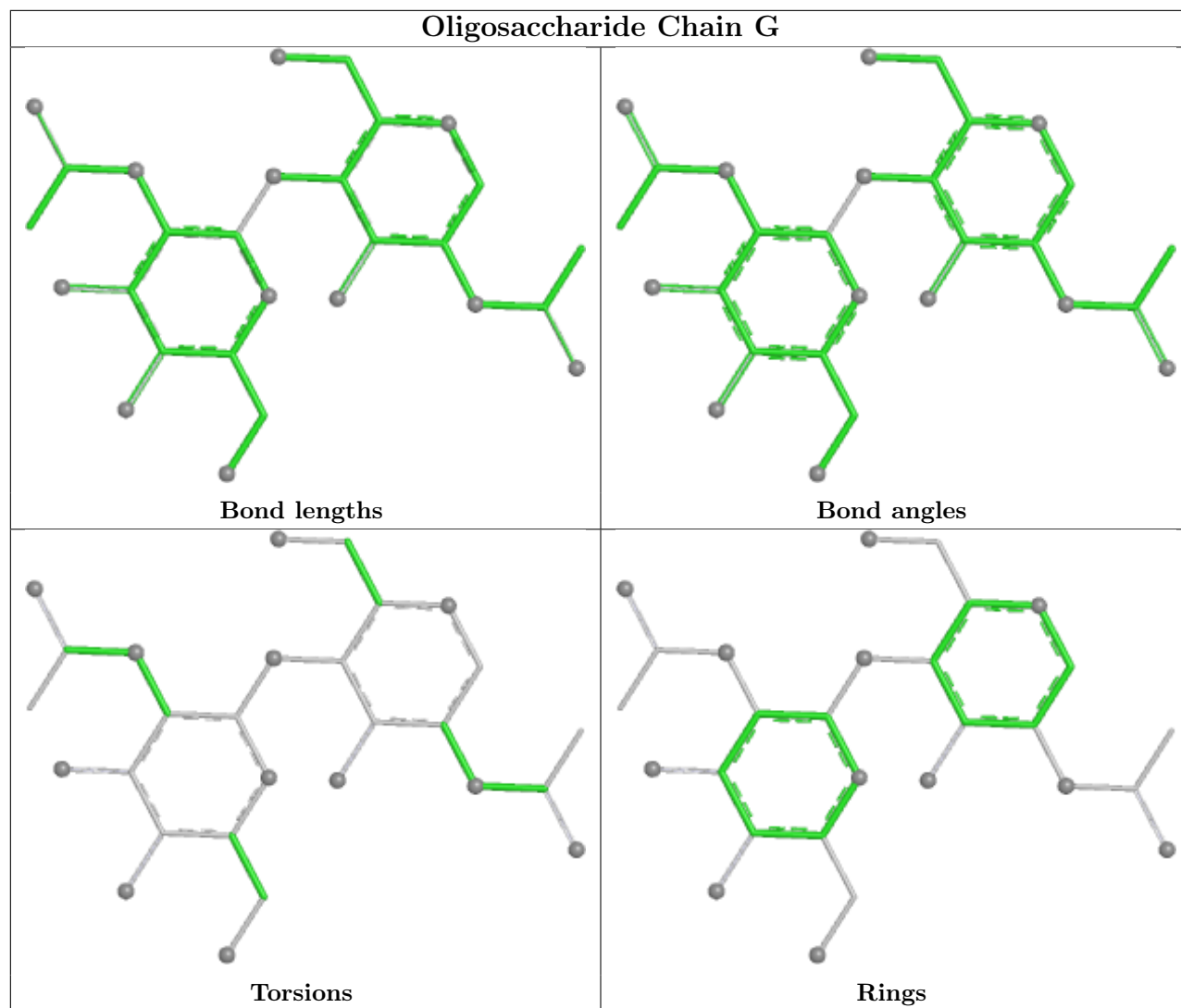
Mol	Chain	Res	Type	Atoms
3	I	1	NAG	C8-C7-N2-C2
3	I	1	NAG	O7-C7-N2-C2
3	I	2	NAG	C8-C7-N2-C2
3	I	2	NAG	O7-C7-N2-C2
3	H	2	NAG	O5-C5-C6-O6
3	I	2	NAG	O5-C5-C6-O6

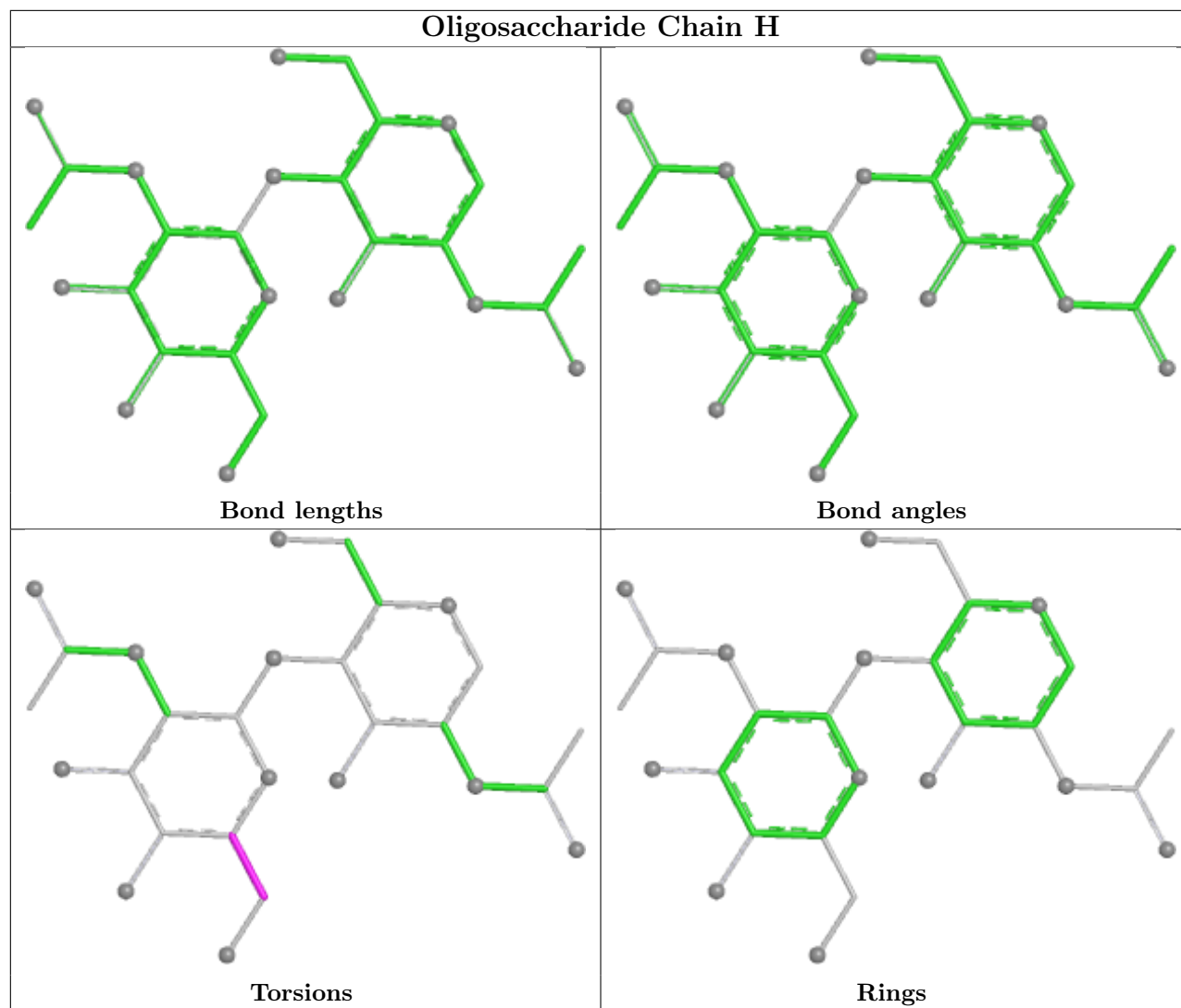
There are no ring outliers.

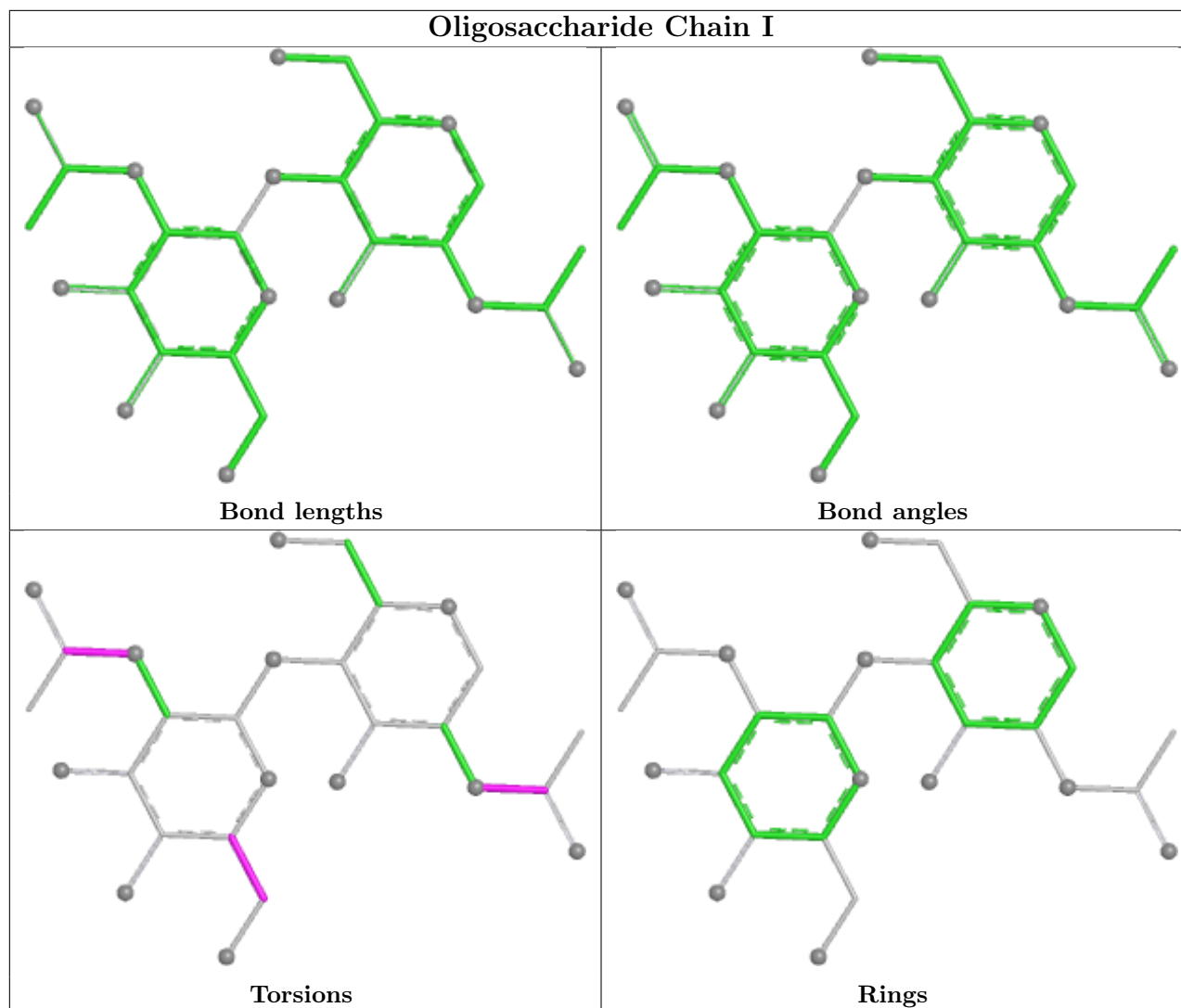
1 monomer is involved in 1 short contact:

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

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







5.6 Ligand geometry [i](#)

3 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	C	601	1	14,14,15	0.49	0	17,19,21	0.79	0
4	NAG	A	601	1	14,14,15	0.47	0	17,19,21	0.75	0
4	NAG	E	601	1	14,14,15	0.42	0	17,19,21	1.39	2 (11%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	C	601	1	-	2/6/23/26	0/1/1/1
4	NAG	A	601	1	-	0/6/23/26	0/1/1/1
4	NAG	E	601	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(°)
4	E	601	NAG	C1-O5-C5	4.25	117.88	112.19
4	E	601	NAG	C4-C3-C2	-2.29	107.66	111.02

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	E	601	NAG	C8-C7-N2-C2
4	E	601	NAG	O7-C7-N2-C2
4	C	601	NAG	C8-C7-N2-C2
4	C	601	NAG	O7-C7-N2-C2

There are no ring outliers.

No monomer is involved in short contacts.

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	322/322 (100%)	-0.11	4 (1%) 76 73	20, 44, 72, 119	0
1	C	322/322 (100%)	0.06	9 (2%) 55 50	23, 51, 89, 121	0
1	E	322/322 (100%)	0.15	8 (2%) 58 54	30, 56, 97, 143	0
2	B	175/175 (100%)	0.05	3 (1%) 69 65	30, 57, 88, 140	0
2	D	175/175 (100%)	0.38	13 (7%) 20 18	25, 75, 123, 191	0
2	F	175/175 (100%)	0.53	14 (8%) 18 16	29, 84, 147, 196	0
All	All	1491/1491 (100%)	0.13	51 (3%) 48 43	20, 54, 115, 196	0

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	75	ILE	4.2
2	F	507	ILE	4.0
2	F	460	LEU	3.3
1	E	73	GLU	3.2
2	F	361	SER	3.1
1	C	129	HIS	3.1
2	F	475	TYR	3.0
2	F	468	GLY	3.0
2	D	507	ILE	2.9
1	C	218	ALA	2.8
2	D	498	GLU	2.8
1	C	325	PRO	2.8
2	D	341	ALA	2.8
1	E	74	PHE	2.8
1	A	70	MET	2.7
1	C	278	CYS	2.7
2	B	352	VAL	2.6
2	F	486	VAL	2.6
2	D	401	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	278	CYS	2.5
2	F	487	ARG	2.5
1	E	128	SER	2.5
1	E	278	CYS	2.4
2	D	361	SER	2.4
2	F	509	GLY	2.4
1	C	324	SER	2.4
2	D	468	GLY	2.4
1	E	217	ILE	2.3
1	A	182	ILE	2.3
1	C	7	ILE	2.3
2	F	474	PHE	2.3
2	D	475	TYR	2.3
2	D	458	LEU	2.3
1	C	158	ASP	2.3
2	D	509	GLY	2.3
2	D	356	TYR	2.3
2	F	488	ASN	2.2
2	D	508	SER	2.2
2	F	458	LEU	2.2
1	C	274	GLU	2.2
2	F	504	ARG	2.2
1	E	324	SER	2.1
1	E	248	ASN	2.1
2	D	502	LEU	2.1
2	F	493	TYR	2.1
2	F	348	TRP	2.1
1	C	219	THR	2.1
2	B	340	ILE	2.1
2	D	478	CYS	2.1
2	B	341	ALA	2.0
1	A	158	ASP	2.0

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

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

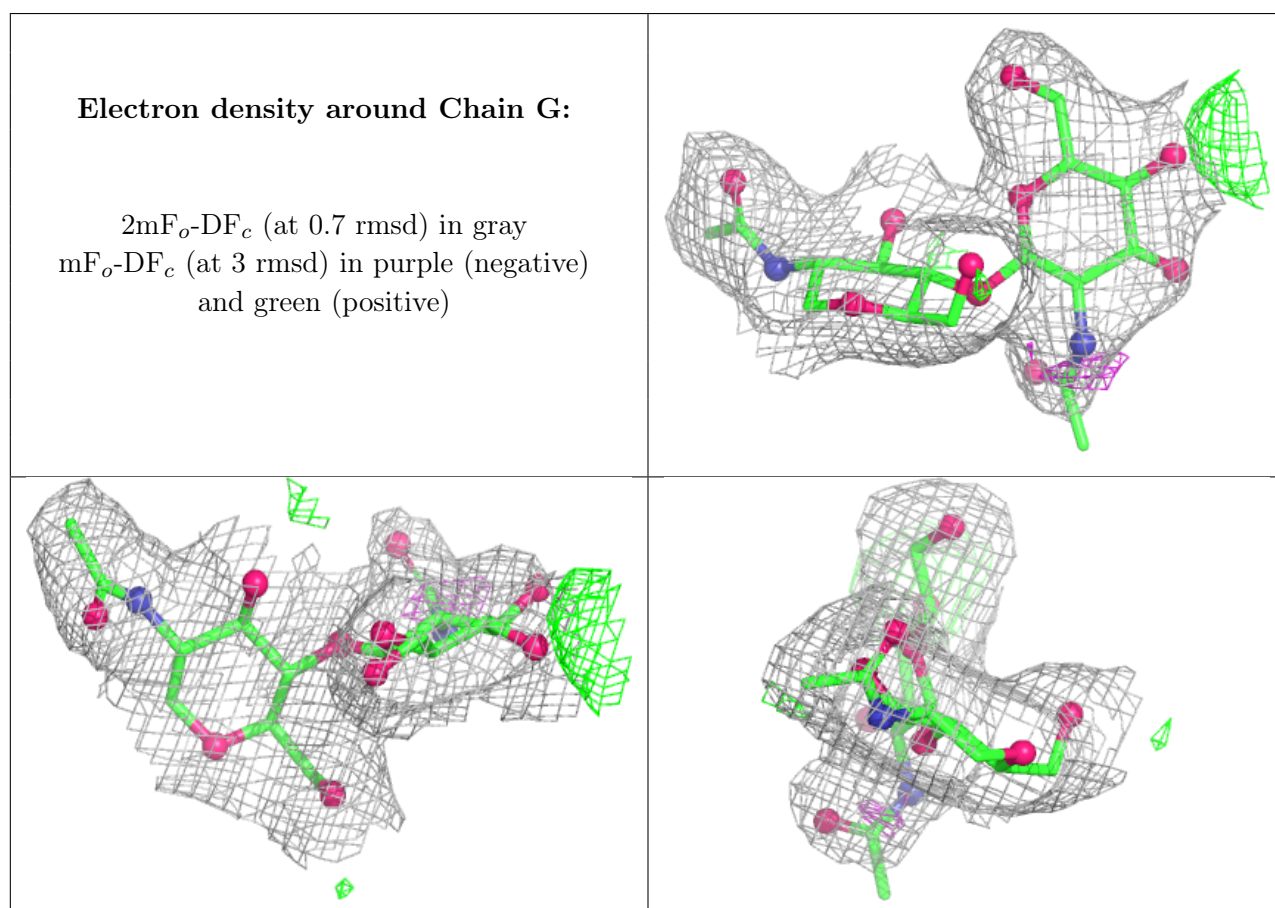
6.3 Carbohydrates [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum,

median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

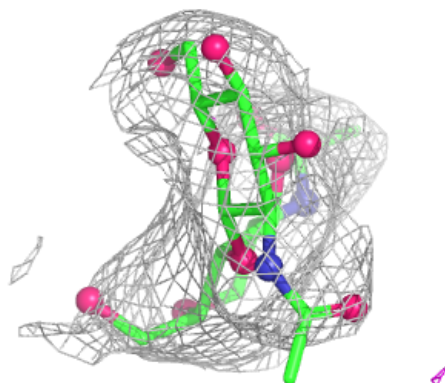
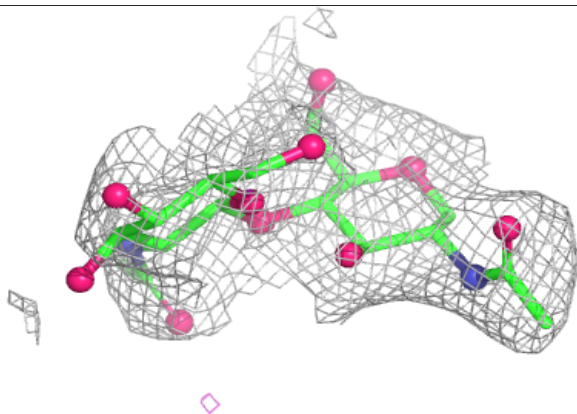
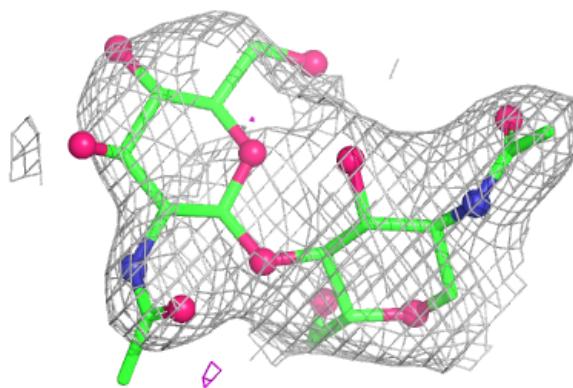
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	I	2	14/15	0.52	0.12	73,89,97,99	0
3	NAG	G	2	14/15	0.71	0.12	52,56,66,72	0
3	NAG	H	2	14/15	0.72	0.13	80,89,92,96	0
3	NAG	I	1	14/15	0.87	0.09	46,62,71,78	0
3	NAG	H	1	14/15	0.88	0.08	47,71,85,85	0
3	NAG	G	1	14/15	0.93	0.06	28,43,53,55	0

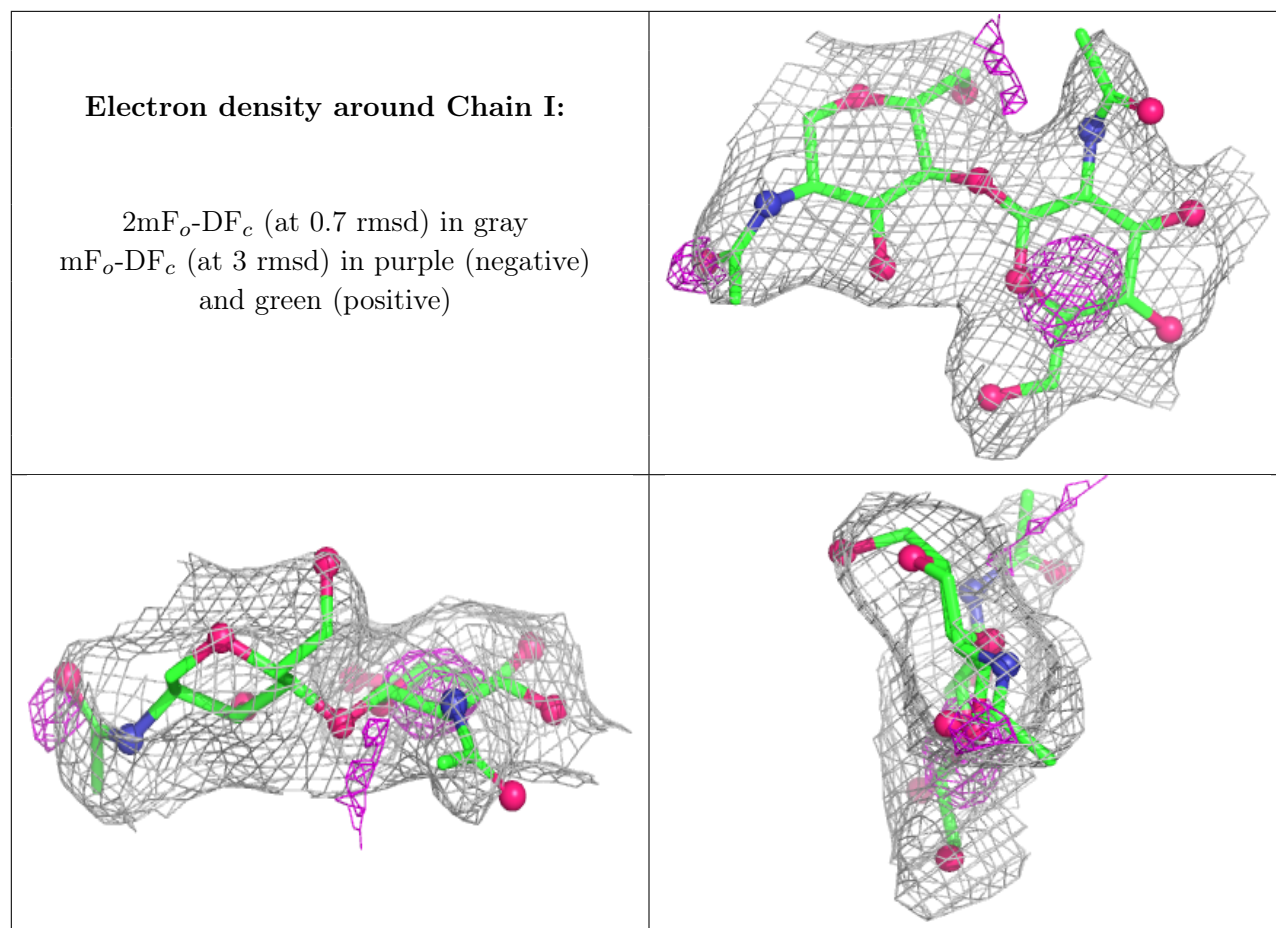
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.



Electron density around Chain H:

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





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
4	NAG	E	601	14/15	0.45	0.15	85,92,96,99	0
4	NAG	C	601	14/15	0.64	0.16	79,85,90,90	0
4	NAG	A	601	14/15	0.64	0.13	78,82,87,87	0

6.5 Other polymers [i](#)

There are no such residues in this entry.