



Full wwPDB X-ray Structure Validation Report ⓘ

Sep 7, 2020 – 02:21 PM BST

PDB ID : 4JUJ
Title : Crystal structure of 1918 pandemic influenza virus hemagglutinin mutant D225G complexed with human receptor analogue LSTc
Authors : Zhang, W.; Shi, Y.; Qi, J.; Gao, F.; Li, Q.; Fan, Z.; Yan, J.; Gao, G.F.
Deposited on : 2013-03-25
Resolution : 3.01 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.14.2
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.14.2

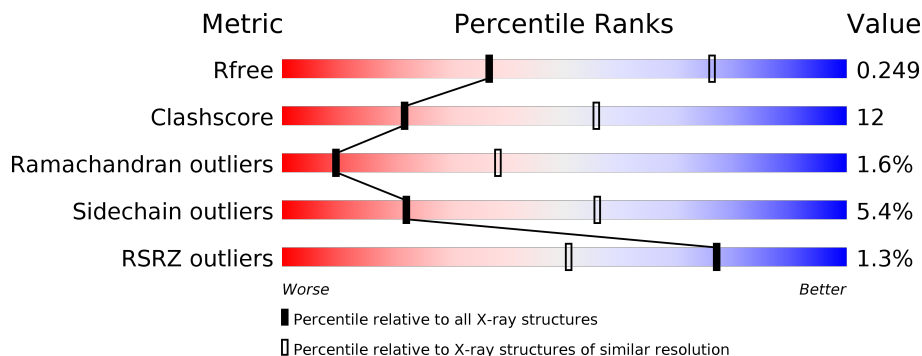
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	2399 (3.04-3.00)
Clashscore	141614	2734 (3.04-3.00)
Ramachandran outliers	138981	2640 (3.04-3.00)
Sidechain outliers	138945	2643 (3.04-3.00)
RSRZ outliers	127900	2287 (3.04-3.00)

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

Mol	Chain	Length	Quality of chain
1	A	324	76% 23% .
1	C	324	77% 22% .
1	E	324	76% 23% .
2	B	170	72% 24% .
2	D	170	3% 71% 20% 5% . . .
2	F	170	4% 70% 22% . . .

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Mol	Chain	Length	Quality of chain
3	G	2	 50% 50%
3	I	2	 50% 50%
3	K	2	 50% 50%
4	H	3	 33% 67%
4	J	3	 67% 33%

2 Entry composition i

There are 5 unique types of molecules in this entry. The entry contains 11784 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	324	2512	1584	434	483	11	0	0	0
1	C	324	2512	1584	434	483	11	0	0	0
1	E	324	2512	1584	434	483	11	0	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	225	GLY	ASP	engineered mutation	UNP Q9WFX3
A	326	ALA	-	expression tag	UNP Q9WFX3
A	327	ARG	-	expression tag	UNP Q9WFX3
C	225	GLY	ASP	engineered mutation	UNP Q9WFX3
C	326	ALA	-	expression tag	UNP Q9WFX3
C	327	ARG	-	expression tag	UNP Q9WFX3
E	225	GLY	ASP	engineered mutation	UNP Q9WFX3
E	326	ALA	-	expression tag	UNP Q9WFX3
E	327	ARG	-	expression tag	UNP Q9WFX3

- Molecule 2 is a protein called Hemagglutinin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	170	1368	855	236	271	6	0	0	0
2	D	164	1317	825	225	261	6	0	0	0
2	F	165	1326	830	226	264	6	0	0	0

- 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			
3	I	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	K	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
			45	25	2	18			
4	J	3	Total	C	N	O	0	0	0
			45	25	2	18			

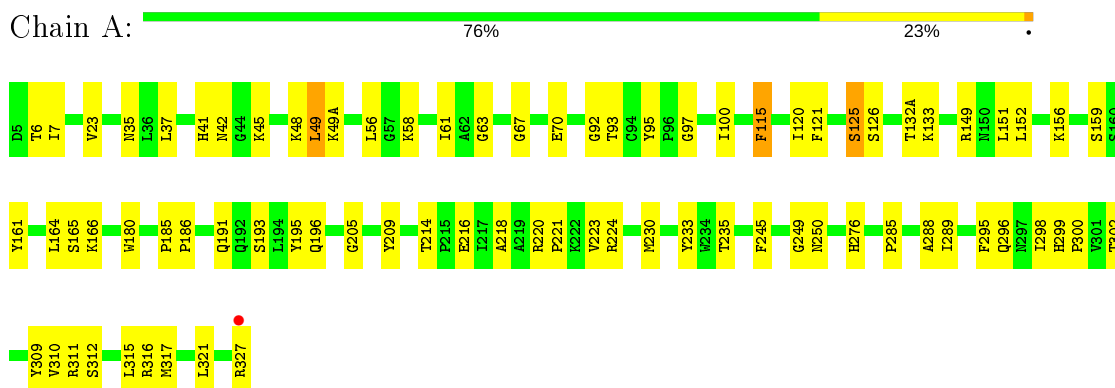
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	21	Total	O	0	0
			21	21		
5	B	6	Total	O	0	0
			6	6		
5	C	7	Total	O	0	0
			7	7		
5	D	2	Total	O	0	0
			2	2		
5	E	23	Total	O	0	0
			23	23		
5	F	4	Total	O	0	0
			4	4		

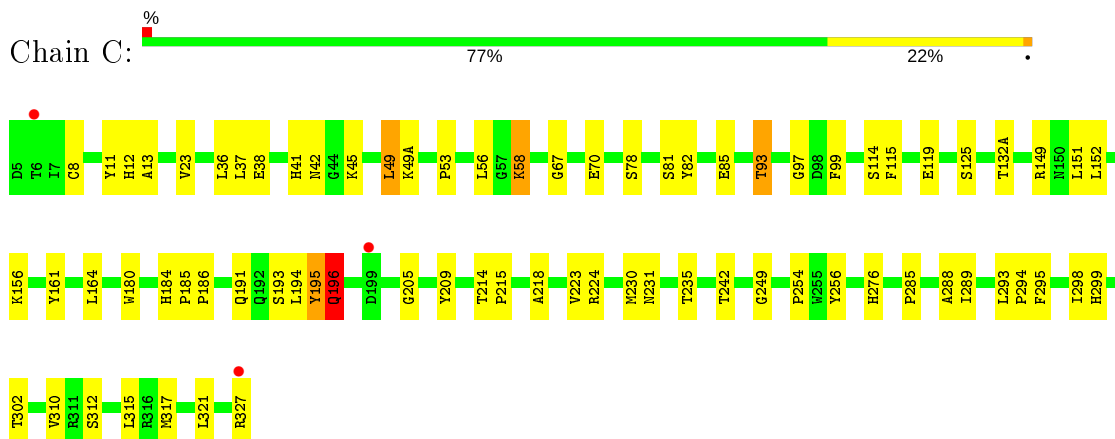
3 Residue-property plots [i](#)

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.

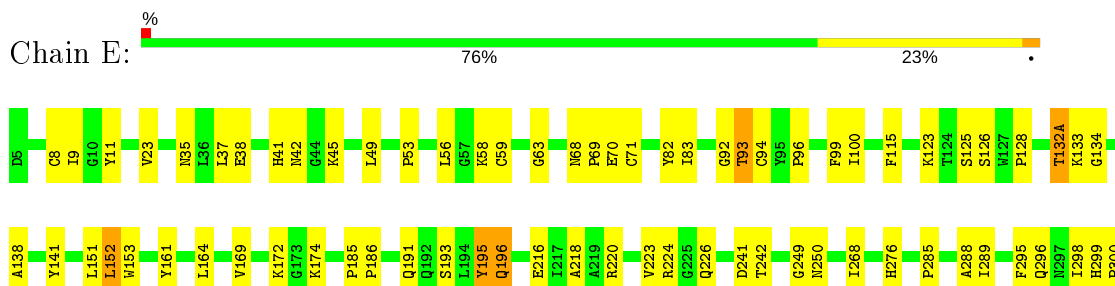
- Molecule 1: Hemagglutinin



- Molecule 1: Hemagglutinin

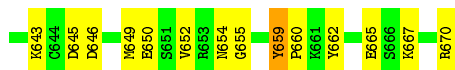
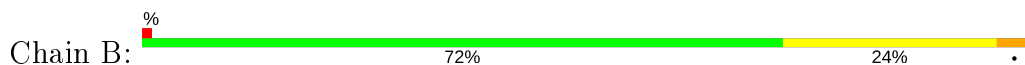


- Molecule 1: Hemagglutinin

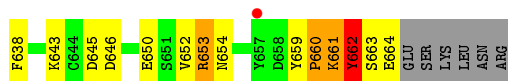
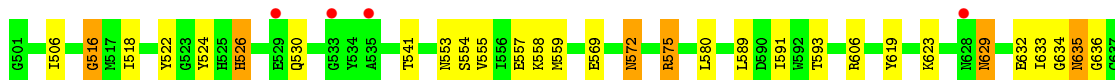




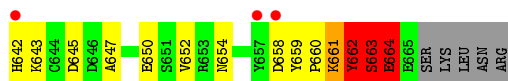
• Molecule 2: Hemagglutinin



• Molecule 2: Hemagglutinin



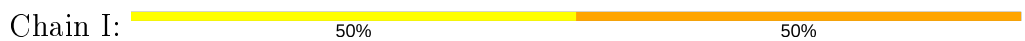
• Molecule 2: Hemagglutinin



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



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



MAG1
MAG2

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

Chain K:
50% 50%

MAG1
MAG2

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

Chain H:
33% 67%

MAG1
GAL2
STA3

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

Chain J:
67% 33%

MAG1
GAL2
STA3

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	71.91Å 242.15Å 71.75Å 90.00° 119.66° 90.00°	Depositor
Resolution (Å)	38.25 – 3.01 43.48 – 3.01	Depositor EDS
% Data completeness (in resolution range)	88.7 (38.25-3.01) 96.3 (43.48-3.01)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.38 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.5_2)	Depositor
R, R_{free}	0.202 , 0.247 0.205 , 0.249	Depositor DCC
R_{free} test set	2044 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	56.8	Xtriage
Anisotropy	0.483	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 43.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.000 for l,k,-h-l 0.000 for -h-l,k,h 0.017 for -h-l,-k,l 0.019 for h,-k,-h-l 0.015 for l,-k,h	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	11784	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.67% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality i

5.1 Standard geometry i

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

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.31	1/2576 (0.0%)	0.43	0/3508
1	C	0.23	0/2576	0.46	2/3508 (0.1%)
1	E	0.24	0/2576	0.45	2/3508 (0.1%)
2	B	0.28	0/1395	0.43	0/1878
2	D	0.27	0/1344	0.44	0/1811
2	F	0.23	0/1353	0.49	1/1823 (0.1%)
All	All	0.26	1/11820 (0.0%)	0.45	5/16036 (0.0%)

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
2	D	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	195	TYR	CE1-CZ	-5.29	1.31	1.38

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	664	GLU	N-CA-CB	10.56	129.61	110.60
1	C	195	TYR	CB-CA-C	8.68	127.76	110.40
1	C	196	GLN	CB-CA-C	-8.43	93.53	110.40
1	E	195	TYR	CB-CA-C	8.08	126.57	110.40
1	E	196	GLN	CB-CA-C	-6.39	97.62	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	662	TYR	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	2512	0	2440	59	0
1	C	2512	0	2440	56	0
1	E	2512	0	2440	51	0
2	B	1368	0	1285	43	0
2	D	1317	0	1231	42	0
2	F	1326	0	1237	41	0
3	G	28	0	25	1	0
3	I	28	0	25	1	0
3	K	28	0	25	1	0
4	H	45	0	38	2	0
4	J	45	0	38	3	0
5	A	21	0	0	0	0
5	B	6	0	0	0	0
5	C	7	0	0	0	0
5	D	2	0	0	0	0
5	E	23	0	0	0	0
5	F	4	0	0	0	0
All	All	11784	0	11224	273	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 12.

All (273) 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:C:196:GLN:OE1	1:C:196:GLN:O	1.71	1.08
2:D:661:LYS:HD3	2:D:662:TYR:N	1.74	1.02
2:F:662:TYR:O	2:F:663:SER:HB3	1.62	0.97

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:663:SER:O	2:F:664:GLU:HG2	1.72	0.88
1:E:310:VAL:HG12	1:E:312:SER:H	1.36	0.88
1:C:195:TYR:O	1:C:196:GLN:HB2	1.71	0.88
2:F:663:SER:OG	2:F:664:GLU:N	2.06	0.85
2:D:661:LYS:HD3	2:D:661:LYS:C	1.98	0.84
1:A:156:LYS:HD2	1:A:196:GLN:HB2	1.58	0.84
2:D:661:LYS:O	2:D:662:TYR:CD2	2.30	0.84
2:D:661:LYS:C	2:D:662:TYR:CD2	2.52	0.82
2:F:642:HIS:HB2	2:F:662:TYR:OH	1.79	0.81
2:D:663:SER:HA	2:D:664:GLU:HB3	1.62	0.81
1:C:41:HIS:HB3	1:C:298:ILE:HD13	1.64	0.80
2:F:642:HIS:HB2	2:F:662:TYR:CZ	2.17	0.78
1:A:156:LYS:CD	1:A:196:GLN:HB2	2.13	0.78
2:D:575:ARG:HH11	2:D:575:ARG:HB2	1.49	0.76
2:B:650:GLU:O	2:B:654:ASN:HB2	1.86	0.76
2:F:662:TYR:O	2:F:663:SER:CB	2.34	0.75
1:A:321:LEU:H	1:A:321:LEU:HD23	1.51	0.75
2:D:661:LYS:O	2:D:662:TYR:HD2	1.69	0.74
2:F:530:GLN:HE22	2:F:645:ASP:HB2	1.51	0.74
1:E:195:TYR:O	1:E:196:GLN:HB3	1.87	0.74
1:A:159:SER:O	1:A:196:GLN:HG2	1.87	0.73
2:F:663:SER:O	2:F:664:GLU:CG	2.37	0.73
1:C:119:GLU:HG3	1:C:256:TYR:CE2	2.24	0.72
1:C:161:TYR:HB2	1:C:195:TYR:O	1.90	0.71
2:B:667:LYS:HA	2:B:667:LYS:HE2	1.73	0.70
1:A:37:LEU:HB2	1:A:315:LEU:HB2	1.74	0.69
1:A:289:ILE:HD11	1:A:298:ILE:HD12	1.74	0.68
1:E:41:HIS:HB3	1:E:298:ILE:HD13	1.74	0.68
1:E:185:PRO:HG2	1:E:191:GLN:NE2	2.09	0.68
2:D:629:ASN:HD22	2:D:629:ASN:N	1.90	0.68
2:B:530:GLN:HE22	2:B:645:ASP:HB2	1.59	0.68
1:C:156:LYS:HD2	1:C:196:GLN:HG2	1.75	0.67
1:A:41:HIS:HB3	1:A:298:ILE:HD13	1.76	0.66
1:C:193:SER:HB2	4:J:1:NAG:H82	1.76	0.66
1:C:93:THR:O	1:C:93:THR:HG23	1.96	0.65
2:B:659:TYR:HB3	2:B:660:PRO:HD3	1.78	0.65
2:B:525:HIS:ND1	2:B:533:GLY:O	2.30	0.65
2:D:572:ASN:HD22	2:D:572:ASN:N	1.93	0.64
2:D:650:GLU:O	2:D:654:ASN:HB2	1.97	0.64
1:C:196:GLN:OE1	1:C:196:GLN:C	2.36	0.64
1:C:285:PRO:HG2	1:C:299:HIS:CE1	2.32	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:310:VAL:HG13	1:C:312:SER:H	1.61	0.64
2:D:516:GLY:O	2:D:518:ILE:HG23	1.98	0.64
2:F:575:ARG:HB2	2:F:575:ARG:HH11	1.63	0.64
2:B:591:ILE:HD13	2:D:591:ILE:HG21	1.80	0.63
1:C:186:PRO:HA	1:C:218:ALA:O	1.99	0.63
1:C:195:TYR:O	1:C:196:GLN:CB	2.40	0.63
1:A:310:VAL:HG22	2:B:593:THR:HA	1.81	0.62
1:E:161:TYR:HB2	1:E:195:TYR:O	1.99	0.62
2:B:575:ARG:HB2	2:B:575:ARG:HH11	1.64	0.62
1:C:327:ARG:HD2	1:C:327:ARG:O	2.00	0.62
2:B:629:ASN:N	2:B:629:ASN:HD22	1.96	0.62
1:E:285:PRO:HG2	1:E:299:HIS:CE1	2.35	0.61
2:F:661:LYS:HB2	2:F:661:LYS:HZ3	1.64	0.61
2:D:591:ILE:HD13	2:F:591:ILE:HG21	1.82	0.61
1:E:185:PRO:HG2	1:E:191:GLN:HE21	1.64	0.61
2:D:661:LYS:CD	2:D:661:LYS:C	2.66	0.60
1:E:42:ASN:HB3	1:E:288:ALA:H	1.66	0.60
1:A:42:ASN:HB3	1:A:288:ALA:H	1.66	0.60
2:D:530:GLN:NE2	2:D:645:ASP:HB2	2.17	0.60
1:E:289:ILE:HD11	1:E:298:ILE:HD12	1.82	0.60
2:D:661:LYS:C	2:D:662:TYR:CG	2.75	0.59
2:F:526:HIS:O	2:F:526:HIS:CG	2.56	0.59
1:C:321:LEU:H	1:C:321:LEU:HD23	1.67	0.59
2:F:659:TYR:N	2:F:660:PRO:HD3	2.17	0.59
1:E:223:VAL:HG12	1:E:224:ARG:HG3	1.85	0.59
1:A:156:LYS:HE2	1:A:193:SER:O	2.02	0.58
1:C:164:LEU:O	1:C:164:LEU:HD12	2.04	0.58
2:B:591:ILE:HG21	2:F:591:ILE:HD13	1.86	0.58
2:D:652:VAL:C	2:D:654:ASN:H	2.07	0.58
2:F:629:ASN:HD22	2:F:629:ASN:N	2.00	0.58
1:A:310:VAL:HG13	1:A:312:SER:H	1.68	0.58
2:D:660:PRO:O	2:D:660:PRO:HG2	2.04	0.57
2:F:530:GLN:NE2	2:F:645:ASP:HB2	2.19	0.57
1:A:133:LYS:O	4:H:3:SIA:H113	2.04	0.57
2:F:658:ASP:C	2:F:660:PRO:HD3	2.24	0.57
1:E:195:TYR:O	1:E:196:GLN:CB	2.52	0.57
1:A:193:SER:HB2	4:H:1:NAG:H82	1.86	0.56
1:E:96:PRO:HG3	1:E:226:GLN:HB2	1.87	0.56
1:C:196:GLN:CA	1:C:196:GLN:OE1	2.53	0.56
2:F:619:TYR:CE1	2:F:636:GLY:HA2	2.39	0.56
1:C:56:LEU:HD12	1:C:85:GLU:HG2	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:37:LEU:HB2	1:E:315:LEU:HB2	1.87	0.56
2:D:629:ASN:ND2	2:D:629:ASN:N	2.55	0.55
1:E:132(A):THR:HG22	1:E:133:LYS:H	1.71	0.55
2:B:629:ASN:ND2	2:B:665:GLU:HG2	2.22	0.55
2:D:575:ARG:NH1	2:D:575:ARG:HB2	2.19	0.55
1:A:295:PHE:CZ	2:B:559:MET:HE2	2.41	0.55
2:D:572:ASN:HD22	2:D:572:ASN:H	1.55	0.55
1:E:123:LYS:HG3	1:E:152:LEU:HD11	1.88	0.54
1:A:45:LYS:HD2	1:A:276:HIS:CG	2.43	0.54
1:C:185:PRO:HG2	1:C:191:GLN:NE2	2.22	0.54
1:A:223:VAL:HG12	1:A:224:ARG:HG3	1.89	0.54
2:B:619:TYR:CE1	2:B:636:GLY:HA2	2.42	0.54
1:C:289:ILE:HD11	1:C:298:ILE:HD12	1.89	0.54
1:E:134:GLY:HA3	1:E:153:TRP:HB3	1.90	0.54
2:F:650:GLU:O	2:F:654:ASN:HB2	2.08	0.54
1:A:186:PRO:HA	1:A:218:ALA:O	2.07	0.53
2:F:554:SER:O	2:F:558:LYS:HG2	2.08	0.53
2:F:623:LYS:HB2	2:F:638:PHE:HZ	1.73	0.53
2:D:572:ASN:ND2	2:D:572:ASN:H	2.06	0.53
1:A:196:GLN:HG3	1:A:196:GLN:O	2.08	0.53
1:E:68:ASN:HB3	1:E:71:CYS:SG	2.49	0.53
1:A:161:TYR:CZ	1:A:249:GLY:HA2	2.44	0.53
1:A:159:SER:C	1:A:196:GLN:HG2	2.29	0.52
1:E:69:PRO:HB3	1:E:141:TYR:HB2	1.90	0.52
1:C:45:LYS:HD2	1:C:276:HIS:CG	2.44	0.52
1:A:48:LYS:O	1:A:49:LEU:HD12	2.09	0.52
1:C:37:LEU:HB2	1:C:315:LEU:HB2	1.91	0.52
1:A:327:ARG:HD2	1:A:327:ARG:O	2.10	0.52
1:C:310:VAL:HG22	2:D:593:THR:HA	1.92	0.52
2:B:660:PRO:C	2:B:662:TYR:H	2.13	0.52
1:E:299:HIS:ND1	1:E:300:PRO:HD2	2.25	0.52
2:D:660:PRO:O	2:D:660:PRO:CG	2.56	0.52
1:A:289:ILE:CD1	1:A:298:ILE:HD12	2.39	0.52
2:B:607:THR:O	2:B:610:PHE:HB3	2.10	0.51
2:F:517:MET:SD	2:F:523:GLY:HA3	2.50	0.51
1:A:156:LYS:HD3	1:A:196:GLN:HB2	1.90	0.51
1:A:285:PRO:HG2	1:A:299:HIS:CE1	2.45	0.51
1:C:93:THR:O	1:C:93:THR:CG2	2.58	0.51
1:E:327:ARG:HD2	1:E:327:ARG:O	2.11	0.51
2:D:572:ASN:ND2	2:D:572:ASN:N	2.58	0.51
1:E:321:LEU:HD23	1:E:321:LEU:H	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:93:THR:O	1:E:93:THR:CG2	2.58	0.51
2:D:606:ARG:HH11	2:F:606:ARG:NH1	2.08	0.51
1:C:161:TYR:CB	1:C:195:TYR:O	2.57	0.51
1:E:196:GLN:O	1:E:196:GLN:HG3	2.10	0.51
4:J:3:SIA:O1A	4:J:3:SIA:H6	2.11	0.51
1:C:8:CYS:O	2:D:524:TYR:HA	2.11	0.50
2:B:652:VAL:C	2:B:654:ASN:H	2.14	0.50
1:A:67:GLY:O	1:A:149:ARG:HG2	2.12	0.50
2:F:652:VAL:C	2:F:654:ASN:H	2.16	0.50
1:C:12:HIS:CD2	1:C:13:ALA:N	2.80	0.49
2:D:526:HIS:O	2:D:526:HIS:CG	2.66	0.49
2:D:553:ASN:O	2:D:557:GLU:HG3	2.13	0.49
1:C:67:GLY:O	1:C:149:ARG:HG2	2.13	0.49
2:B:629:ASN:N	2:B:629:ASN:ND2	2.61	0.49
1:A:56:LEU:HD11	1:A:61:ILE:HD13	1.94	0.49
1:C:184:HIS:HE2	1:C:231:ASN:HD22	1.61	0.48
1:E:327:ARG:HD2	1:E:327:ARG:C	2.34	0.48
1:C:289:ILE:CD1	1:C:298:ILE:HD12	2.43	0.48
1:A:165:SER:HA	1:A:245:PHE:O	2.14	0.48
1:C:180:TRP:HZ3	1:C:235:THR:HG22	1.78	0.48
2:B:633:ILE:HD11	2:B:639:GLU:HB2	1.95	0.48
1:C:327:ARG:HD2	1:C:327:ARG:C	2.33	0.48
1:A:180:TRP:HZ3	1:A:235:THR:HG22	1.79	0.47
1:E:63:GLY:CA	1:E:92:GLY:HA2	2.44	0.47
2:B:530:GLN:HE22	2:B:646:ASP:H	1.62	0.47
1:E:35:ASN:ND2	1:E:37:LEU:H	2.13	0.47
2:B:505:ALA:HA	2:B:509:PHE:CE2	2.50	0.47
1:E:289:ILE:CD1	1:E:298:ILE:HD12	2.45	0.47
1:A:63:GLY:CA	1:A:92:GLY:HA2	2.44	0.47
1:C:11:TYR:CZ	2:D:506:ILE:HG23	2.50	0.47
1:C:97:GLY:HA3	1:C:230:MET:O	2.14	0.47
1:C:185:PRO:HG2	1:C:191:GLN:HE21	1.79	0.47
1:C:36:LEU:HD22	1:C:294:PRO:HG2	1.96	0.47
1:C:42:ASN:HB3	1:C:288:ALA:H	1.79	0.46
1:E:216:GLU:O	1:E:220:ARG:NH2	2.47	0.46
1:A:115:PHE:CD1	1:A:115:PHE:C	2.89	0.46
2:B:625:GLN:HE22	2:B:655:GLY:C	2.18	0.46
1:C:58:LYS:HE2	1:C:70:GLU:HB3	1.97	0.46
1:E:70:GLU:HG2	3:K:1:NAG:H82	1.98	0.46
1:E:100:ILE:HD12	1:E:100:ILE:N	2.30	0.46
2:D:661:LYS:HD3	2:D:662:TYR:CA	2.43	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:169:VAL:HG22	1:E:242:THR:OG1	2.16	0.46
1:A:115:PHE:HD1	1:A:115:PHE:C	2.18	0.46
2:F:555:VAL:O	2:F:559:MET:HG2	2.16	0.46
1:A:49:LEU:O	1:A:49(A):LYS:HB2	2.16	0.45
1:A:327:ARG:HD2	1:A:327:ARG:C	2.36	0.45
2:B:504:GLY:O	2:B:508:GLY:HA3	2.17	0.45
2:B:530:GLN:NE2	2:B:645:ASP:HB2	2.28	0.45
1:A:35:ASN:HB2	1:A:316:ARG:NH2	2.31	0.45
2:B:528:ASN:HD21	2:B:530:GLN:NE2	2.14	0.45
1:C:224:ARG:NH2	3:I:1:NAG:O3	2.48	0.45
1:E:38:GLU:O	1:E:296:GLN:HA	2.16	0.45
2:F:659:TYR:N	2:F:660:PRO:CD	2.80	0.45
1:A:97:GLY:HA3	1:A:230:MET:O	2.17	0.45
1:A:295:PHE:HZ	2:B:559:MET:HE2	1.81	0.45
1:C:78:SER:O	1:C:114:SER:HB2	2.17	0.45
1:A:249:GLY:C	1:A:250:ASN:HD22	2.20	0.45
1:E:11:TYR:CZ	2:F:506:ILE:HG23	2.52	0.45
2:F:659:TYR:O	2:F:661:LYS:HG3	2.16	0.45
1:C:53:PRO:HB3	1:C:82:TYR:CZ	2.52	0.45
1:A:205:GLY:HA2	1:A:209:TYR:O	2.17	0.45
2:D:530:GLN:NE2	2:D:646:ASP:H	2.15	0.45
1:A:100:ILE:HD12	1:A:100:ILE:N	2.32	0.44
2:B:606:ARG:HH21	2:F:605:GLU:HG2	1.80	0.44
1:C:205:GLY:HA2	1:C:209:TYR:O	2.17	0.44
1:C:49:LEU:O	1:C:49(A):LYS:HB2	2.18	0.44
1:E:42:ASN:HD21	1:E:45:LYS:HE3	1.81	0.44
1:C:194:LEU:HG	4:J:1:NAG:C8	2.47	0.44
2:F:528:ASN:HD21	2:F:645:ASP:HA	1.81	0.44
1:A:310:VAL:CG1	1:A:312:SER:H	2.30	0.44
2:D:632:GLU:HG2	2:D:638:PHE:CE2	2.53	0.44
1:A:321:LEU:N	1:A:321:LEU:HD23	2.26	0.44
2:B:524:TYR:CD1	2:B:524:TYR:N	2.84	0.44
2:D:662:TYR:HB2	2:D:663:SER:H	1.43	0.44
2:F:629:ASN:ND2	2:F:629:ASN:N	2.64	0.44
1:A:126:SER:HB3	1:A:166:LYS:HE2	1.98	0.44
1:E:138:ALA:O	1:E:224:ARG:NH1	2.50	0.44
1:A:95:TYR:CD2	1:A:230:MET:HB2	2.52	0.44
1:A:311:ARG:NH2	2:B:589:LEU:HD23	2.32	0.44
2:B:572:ASN:HD22	2:B:572:ASN:N	2.16	0.44
2:B:526:HIS:ND1	2:B:526:HIS:O	2.51	0.44
2:B:528:ASN:HD22	2:B:530:GLN:HB2	1.83	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:GLU:HB2	1:C:293:LEU:HD12	2.00	0.44
1:C:42:ASN:ND2	1:C:45:LYS:HB2	2.33	0.44
1:C:161:TYR:CZ	1:C:249:GLY:HA2	2.53	0.43
2:D:619:TYR:CE1	2:D:636:GLY:HA2	2.53	0.43
1:A:156:LYS:HD3	1:A:196:GLN:N	2.33	0.43
1:A:216:GLU:O	1:A:220:ARG:NH2	2.51	0.43
1:A:100:ILE:HG12	1:A:233:TYR:CE2	2.53	0.43
2:F:634:GLY:O	2:F:636:GLY:N	2.47	0.43
1:E:295:PHE:CZ	2:F:559:MET:HE2	2.54	0.43
2:F:663:SER:O	2:F:664:GLU:CD	2.56	0.43
2:B:572:ASN:ND2	2:B:572:ASN:N	2.66	0.43
2:D:659:TYR:N	2:D:660:PRO:HD3	2.33	0.43
1:E:241:ASP:OD2	1:E:242:THR:N	2.51	0.43
1:A:120:ILE:HG23	1:A:121:PHE:N	2.33	0.43
1:E:83:ILE:HB	1:E:268:ILE:HG12	2.01	0.43
1:A:296:GLN:NE2	1:A:309:TYR:HB2	2.34	0.43
2:B:631:LYS:HG3	2:B:631:LYS:O	2.19	0.43
2:D:524:TYR:CE1	2:D:653:ARG:HB2	2.54	0.43
2:D:555:VAL:O	2:D:559:MET:HG2	2.19	0.42
1:E:9:ILE:N	1:E:9:ILE:HD12	2.33	0.42
1:E:310:VAL:HG13	2:F:593:THR:HA	2.01	0.42
1:E:53:PRO:HB3	1:E:82:TYR:CZ	2.54	0.42
2:B:528:ASN:ND2	2:B:530:GLN:NE2	2.67	0.42
1:C:53:PRO:HB3	1:C:82:TYR:CE2	2.53	0.42
1:E:8:CYS:O	2:F:524:TYR:HA	2.19	0.42
1:E:249:GLY:C	1:E:250:ASN:HD22	2.23	0.42
1:A:164:LEU:C	1:A:164:LEU:HD12	2.40	0.42
1:C:214:THR:HA	1:C:215:PRO:HD3	1.91	0.42
1:E:94:CYS:O	1:E:224:ARG:HD3	2.20	0.42
2:F:506:ILE:HG13	2:F:612:ASP:HA	2.01	0.42
2:F:661:LYS:HD3	2:F:661:LYS:H	1.84	0.42
2:B:526:HIS:HD2	2:B:649:MET:HG3	1.85	0.42
2:B:575:ARG:NH1	2:B:575:ARG:HB2	2.33	0.42
1:C:119:GLU:HG3	1:C:256:TYR:CZ	2.54	0.42
1:E:164:LEU:O	1:E:164:LEU:HD12	2.20	0.42
1:E:96:PRO:CG	1:E:226:GLN:HB2	2.50	0.41
1:A:70:GLU:HG2	3:G:1:NAG:H82	2.02	0.41
1:C:81:SER:HG	1:C:82:TYR:HD2	1.68	0.41
1:A:185:PRO:HG2	1:A:191:GLN:NE2	2.35	0.41
1:A:93:THR:HG22	1:A:93:THR:O	2.21	0.41
1:C:180:TRP:HB3	1:C:254:PRO:HG3	2.01	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:295:PHE:CZ	2:D:559:MET:HE2	2.55	0.41
2:B:509:PHE:CD1	2:B:510:ILE:HG13	2.56	0.41
1:E:56:LEU:HB3	1:E:59:CYS:O	2.21	0.41
1:A:185:PRO:HA	1:A:186:PRO:HD3	1.89	0.41
2:B:530:GLN:NE2	2:B:646:ASP:H	2.18	0.41
1:A:221:PRO:HG3	1:C:242:THR:HB	2.03	0.41
2:D:554:SER:O	2:D:558:LYS:HG2	2.21	0.41
2:B:625:GLN:NE2	2:B:655:GLY:C	2.74	0.41
2:B:660:PRO:C	2:B:662:TYR:N	2.74	0.41
2:F:645:ASP:OD1	2:F:647:ALA:HB3	2.20	0.41
2:B:502:LEU:HD22	2:B:609:ASP:OD2	2.21	0.41
1:C:56:LEU:HD12	1:C:85:GLU:CG	2.49	0.41
1:C:223:VAL:HG12	1:C:224:ARG:HG3	2.03	0.41
1:E:45:LYS:HD2	1:E:276:HIS:CD2	2.56	0.41
2:D:623:LYS:HB2	2:D:638:PHE:HZ	1.85	0.40
1:E:126:SER:C	1:E:128:PRO:HD3	2.41	0.40
1:A:7:ILE:HA	2:B:526:HIS:HA	2.04	0.40
1:E:186:PRO:HA	1:E:218:ALA:O	2.22	0.40
2:F:632:GLU:HG2	2:F:638:PHE:CE2	2.56	0.40
1:A:299:HIS:HA	1:A:300:PRO:HD3	1.86	0.40
1:E:172:LYS:O	1:E:174:LYS:HG3	2.21	0.40
2:F:572:ASN:HD22	2:F:572:ASN:N	2.19	0.40
1:A:311:ARG:CZ	2:B:589:LEU:HD23	2.52	0.40
2:D:589:LEU:HA	2:D:589:LEU:HD12	1.94	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	322/324 (99%)	300 (93%)	21 (6%)	1 (0%)	41 75

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	322/324 (99%)	300 (93%)	21 (6%)	1 (0%)	41	75
1	E	322/324 (99%)	303 (94%)	18 (6%)	1 (0%)	41	75
2	B	168/170 (99%)	150 (89%)	13 (8%)	5 (3%)	4	22
2	D	162/170 (95%)	141 (87%)	13 (8%)	8 (5%)	2	12
2	F	163/170 (96%)	139 (85%)	17 (10%)	7 (4%)	2	14
All	All	1459/1482 (98%)	1333 (91%)	103 (7%)	23 (2%)	9	38

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	663	SER
2	F	664	GLU
2	B	561	THR
2	B	635	ASN
2	D	643	LYS
2	D	660	PRO
2	D	662	TYR
2	F	633	ILE
2	F	643	LYS
1	A	125	SER
2	B	643	LYS
1	E	125	SER
2	F	635	ASN
2	B	659	TYR
2	D	516	GLY
2	D	634	GLY
2	F	634	GLY
2	F	662	TYR
2	D	653	ARG
1	C	125	SER
2	D	633	ILE
2	D	635	ASN
2	B	634	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	278/278 (100%)	266 (96%)	12 (4%)	29	64
1	C	278/278 (100%)	266 (96%)	12 (4%)	29	64
1	E	278/278 (100%)	267 (96%)	11 (4%)	31	67
2	B	145/145 (100%)	136 (94%)	9 (6%)	18	50
2	D	139/145 (96%)	128 (92%)	11 (8%)	12	39
2	F	140/145 (97%)	127 (91%)	13 (9%)	9	32
All	All	1258/1269 (99%)	1190 (95%)	68 (5%)	22	56

All (68) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	THR
1	A	23	VAL
1	A	49	LEU
1	A	58	LYS
1	A	115	PHE
1	A	125	SER
1	A	132(A)	THR
1	A	151	LEU
1	A	152	LEU
1	A	214	THR
1	A	302	THR
1	A	317	MET
2	B	522	TYR
2	B	524	TYR
2	B	526	HIS
2	B	561	THR
2	B	572	ASN
2	B	575	ARG
2	B	580	LEU
2	B	635	ASN
2	B	670	ARG
1	C	23	VAL
1	C	49	LEU
1	C	58	LYS
1	C	93	THR
1	C	99	PHE
1	C	115	PHE

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Mol	Chain	Res	Type
1	C	132(A)	THR
1	C	151	LEU
1	C	152	LEU
1	C	196	GLN
1	C	302	THR
1	C	317	MET
2	D	522	TYR
2	D	526	HIS
2	D	541	THR
2	D	569	GLU
2	D	572	ASN
2	D	575	ARG
2	D	580	LEU
2	D	629	ASN
2	D	635	ASN
2	D	661	LYS
2	D	662	TYR
1	E	23	VAL
1	E	49	LEU
1	E	58	LYS
1	E	93	THR
1	E	99	PHE
1	E	115	PHE
1	E	132(A)	THR
1	E	151	LEU
1	E	152	LEU
1	E	193	SER
1	E	302	THR
2	F	522	TYR
2	F	526	HIS
2	F	529	GLU
2	F	541	THR
2	F	550	ASN
2	F	572	ASN
2	F	575	ARG
2	F	580	LEU
2	F	635	ASN
2	F	661	LYS
2	F	662	TYR
2	F	663	SER
2	F	664	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (37) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	35	ASN
1	A	42	ASN
1	A	88	ASN
1	A	130	HIS
1	A	191	GLN
1	A	250	ASN
2	B	528	ASN
2	B	530	GLN
2	B	572	ASN
2	B	625	GLN
2	B	629	ASN
2	B	654	ASN
1	C	12	HIS
1	C	35	ASN
1	C	42	ASN
1	C	191	GLN
1	C	250	ASN
2	D	530	GLN
2	D	572	ASN
2	D	625	GLN
2	D	629	ASN
2	D	654	ASN
1	E	27	ASN
1	E	35	ASN
1	E	42	ASN
1	E	130	HIS
1	E	191	GLN
1	E	196	GLN
1	E	250	ASN
1	E	276	HIS
2	F	528	ASN
2	F	530	GLN
2	F	550	ASN
2	F	572	ASN
2	F	625	GLN
2	F	629	ASN
2	F	654	ASN

5.3.3 RNA

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](#)

12 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	0.49	0	17,19,21	0.80	0
3	NAG	G	2	3	14,14,15	0.52	0	17,19,21	0.69	0
4	NAG	H	1	4	14,14,15	0.54	0	17,19,21	1.01	1 (5%)
4	GAL	H	2	4	11,11,12	0.63	0	15,15,17	0.92	0
4	SIA	H	3	4	17,20,21	0.40	0	21,28,31	3.44	4 (19%)
3	NAG	I	1	1,3	14,14,15	0.49	0	17,19,21	0.91	1 (5%)
3	NAG	I	2	3	14,14,15	0.49	0	17,19,21	0.81	1 (5%)
4	NAG	J	1	4	14,14,15	0.57	0	17,19,21	0.81	0
4	GAL	J	2	4	11,11,12	0.65	0	15,15,17	1.18	2 (13%)
4	SIA	J	3	4	17,20,21	0.35	0	21,28,31	3.77	5 (23%)
3	NAG	K	1	1,3	14,14,15	0.58	0	17,19,21	0.63	0
3	NAG	K	2	3	14,14,15	0.56	0	17,19,21	0.83	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	-	0/6/23/26	0/1/1/1
4	NAG	H	1	4	-	0/6/23/26	0/1/1/1
4	GAL	H	2	4	-	0/2/19/22	0/1/1/1
4	SIA	H	3	4	-	0/14/34/38	0/1/1/1
3	NAG	I	1	1,3	-	0/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	I	2	3	-	0/6/23/26	0/1/1/1
4	NAG	J	1	4	-	0/6/23/26	0/1/1/1
4	GAL	J	2	4	-	0/2/19/22	0/1/1/1
4	SIA	J	3	4	-	2/14/34/38	0/1/1/1
3	NAG	K	1	1,3	-	0/6/23/26	0/1/1/1
3	NAG	K	2	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	J	3	SIA	C3-C2-C1	-12.09	85.54	111.93
4	H	3	SIA	C3-C2-C1	-10.97	87.97	111.93
4	J	3	SIA	O6-C2-C3	-10.78	90.97	109.87
4	H	3	SIA	O6-C2-C3	-9.73	92.81	109.87
4	H	3	SIA	C4-C3-C2	4.41	117.72	109.81
4	J	3	SIA	C4-C3-C2	4.36	117.62	109.81
4	J	2	GAL	O5-C5-C6	3.01	111.92	107.20
4	H	3	SIA	C6-O6-C2	2.51	116.71	111.34
4	J	2	GAL	C1-C2-C3	2.41	112.62	109.67
3	I	2	NAG	O5-C5-C6	2.35	110.89	107.20
3	I	1	NAG	C1-O5-C5	2.30	115.31	112.19
4	J	3	SIA	C6-O6-C2	2.22	116.09	111.34
4	H	1	NAG	O5-C5-C6	2.22	110.68	107.20
4	J	3	SIA	C4-C5-N5	2.18	114.68	110.38

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	J	3	SIA	C11-C10-N5-C5
4	J	3	SIA	O10-C10-N5-C5

There are no ring outliers.

7 monomers are involved in 8 short contacts:

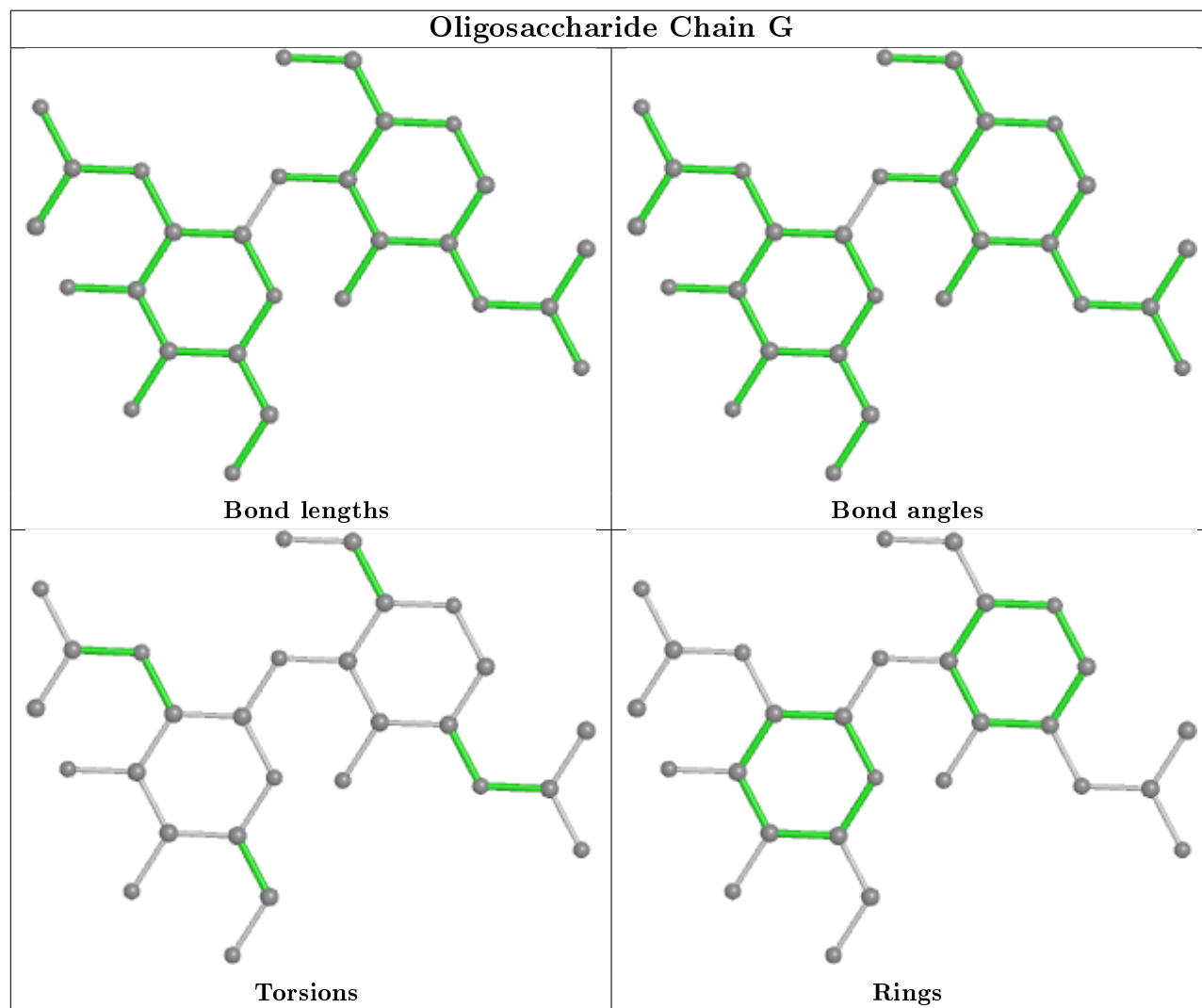
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	1	NAG	1	0
4	H	1	NAG	1	0
4	J	3	SIA	1	0

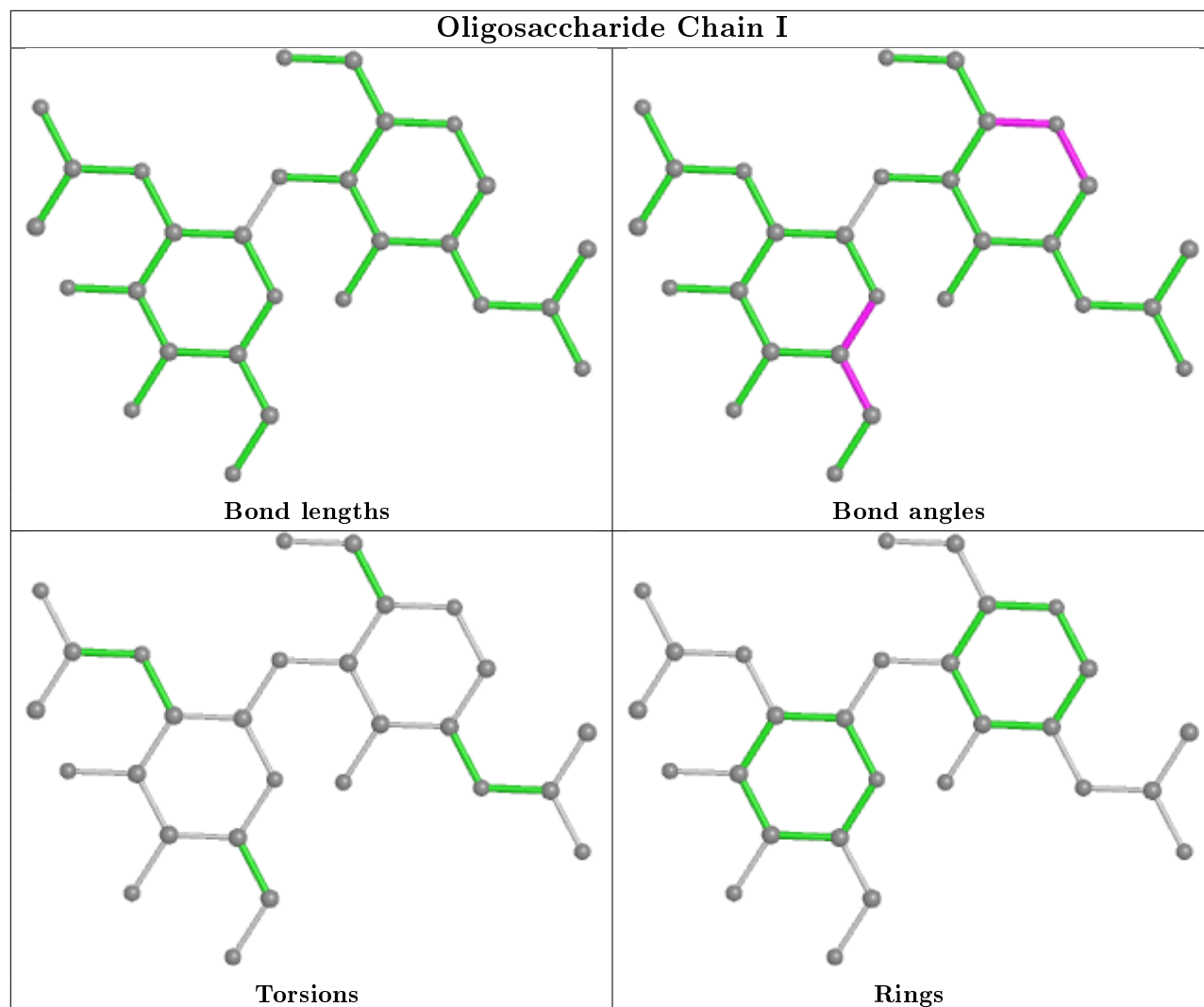
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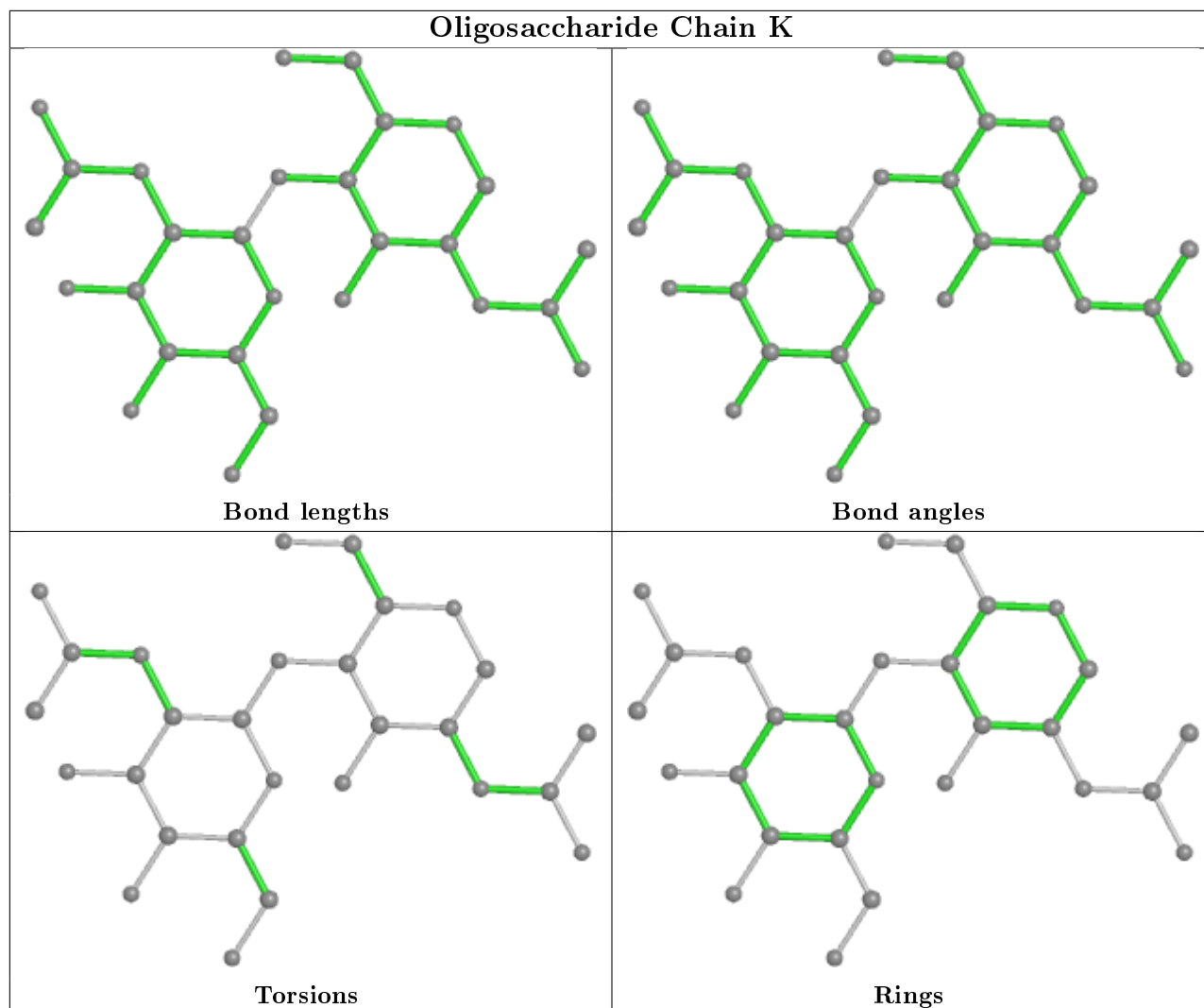
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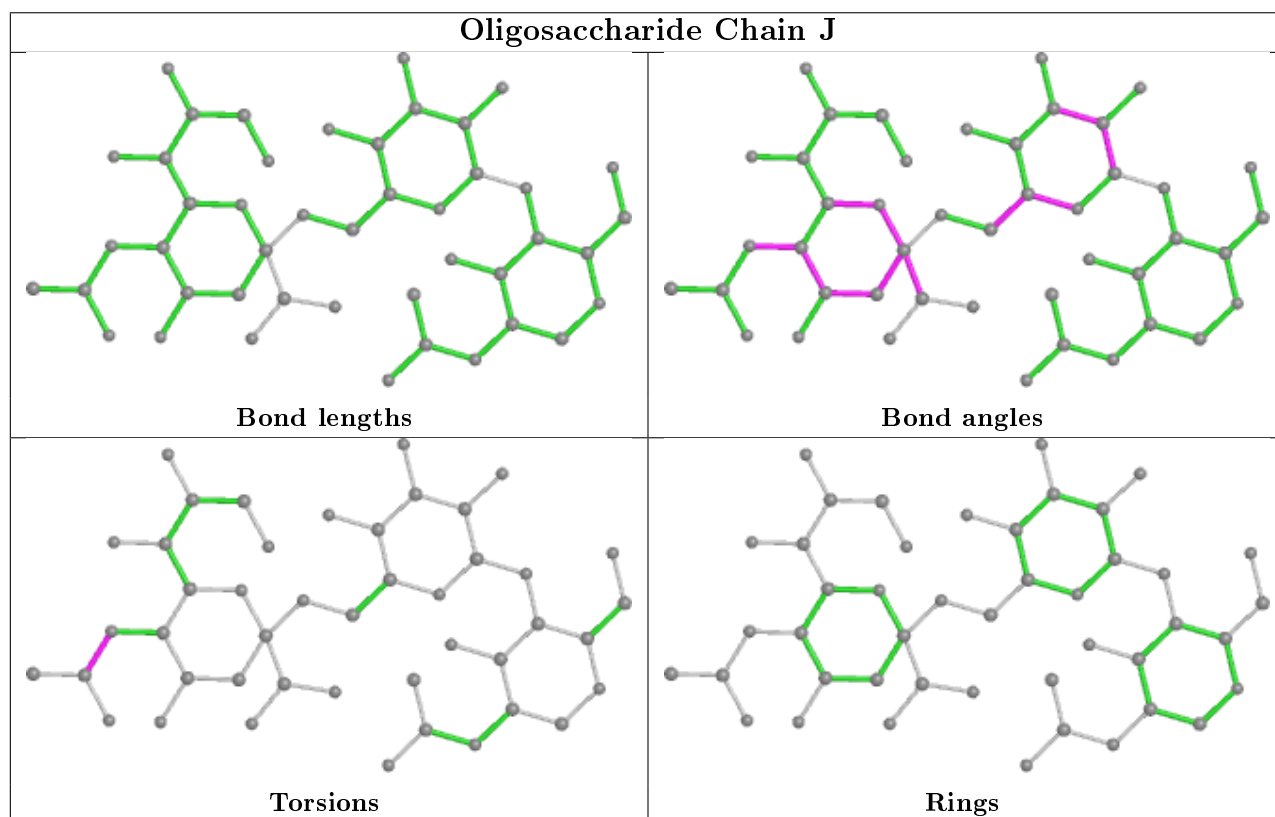
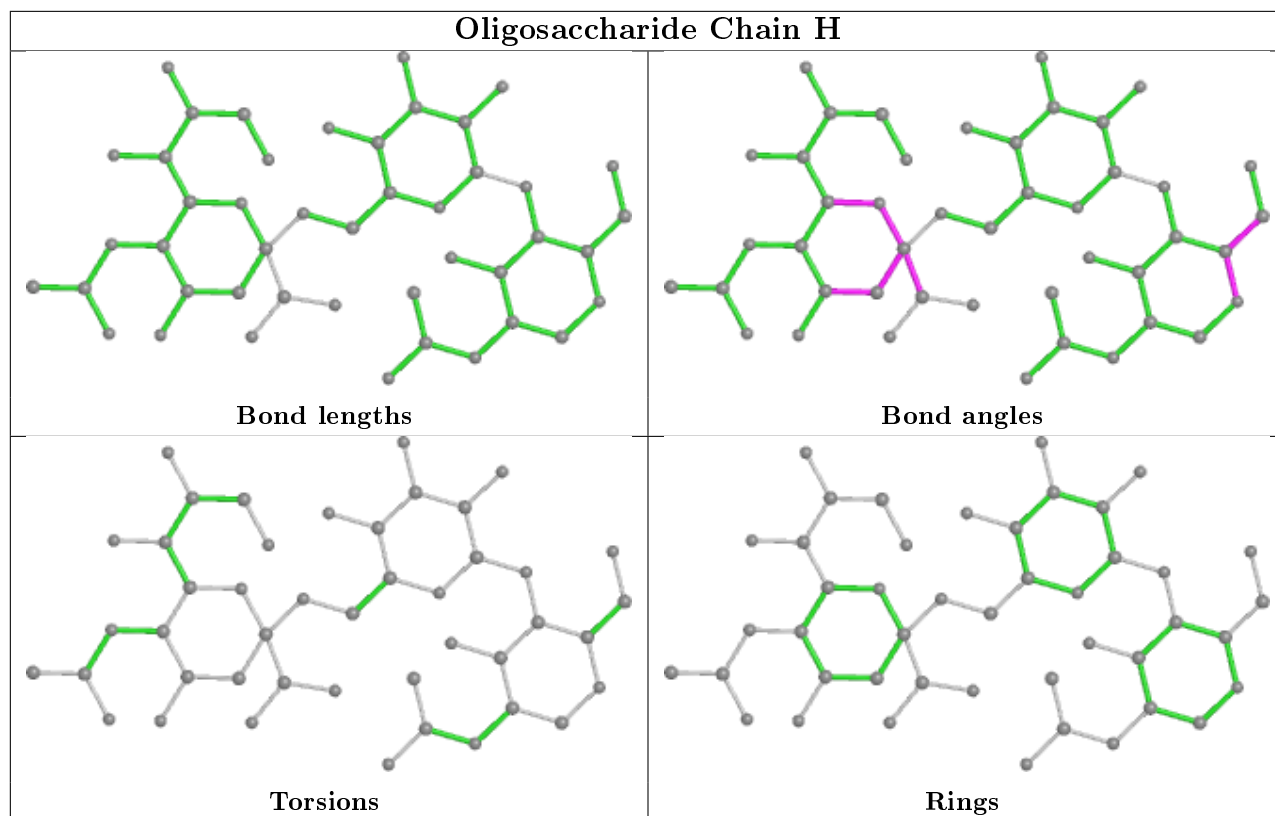
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	H	3	SIA	1	0
3	K	1	NAG	1	0
3	I	1	NAG	1	0
4	J	1	NAG	2	0

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](#)

There are no ligands in this entry.

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	324/324 (100%)	-0.29	1 (0%) 94 83	29, 54, 85, 187	0
1	C	324/324 (100%)	-0.12	3 (0%) 84 62	40, 69, 101, 176	0
1	E	324/324 (100%)	-0.36	2 (0%) 89 72	30, 56, 91, 168	0
2	B	170/170 (100%)	-0.12	1 (0%) 89 72	42, 65, 94, 134	0
2	D	164/170 (96%)	0.16	5 (3%) 50 22	43, 90, 144, 174	0
2	F	165/170 (97%)	0.16	7 (4%) 36 14	40, 81, 141, 161	0
All	All	1471/1482 (99%)	-0.15	19 (1%) 77 51	29, 64, 118, 187	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	327	ARG	6.5
1	E	327	ARG	5.0
1	A	327	ARG	3.8
2	D	533	GLY	3.2
2	F	518	ILE	2.9
2	F	531	GLY	2.9
2	F	642	HIS	2.9
2	D	657	TYR	2.9
2	F	658	ASP	2.8
1	C	6	THR	2.7
1	C	199	ASP	2.7
2	F	640	PHE	2.5
2	D	628	ASN	2.4
2	D	529	GLU	2.4
2	F	519	ASP	2.4
1	E	326	ALA	2.4
2	F	657	TYR	2.3
2	D	535	ALA	2.3
2	B	564	THR	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](#)

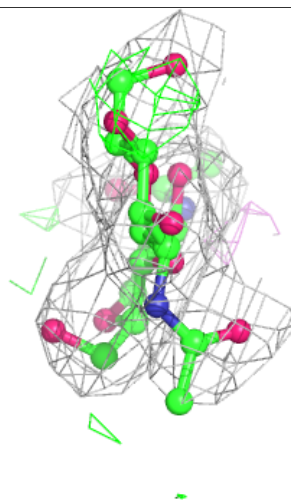
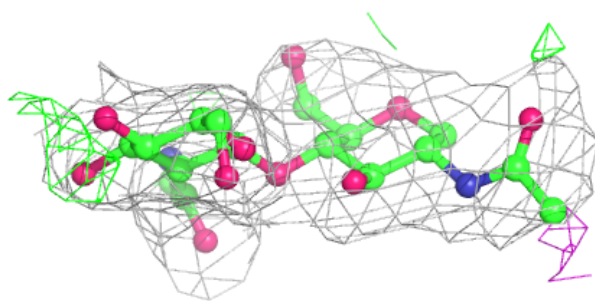
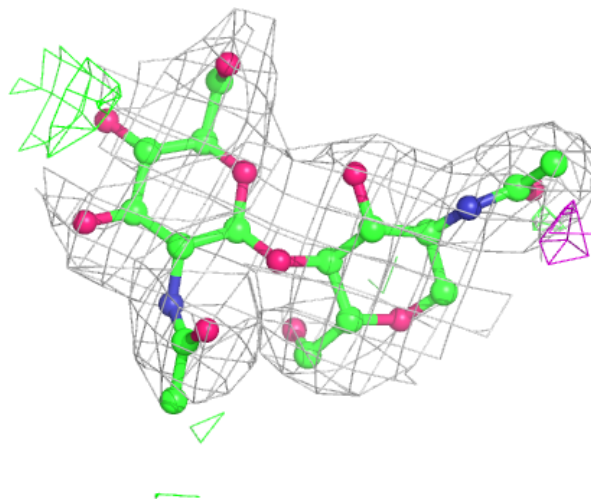
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
4	NAG	J	1	14/15	0.81	0.29	107,114,121,122	0
4	NAG	H	1	14/15	0.83	0.32	101,118,124,128	0
4	GAL	J	2	11/12	0.88	0.19	84,94,105,108	0
4	GAL	H	2	11/12	0.90	0.21	75,95,108,109	0
3	NAG	I	2	14/15	0.91	0.15	70,79,96,96	0
3	NAG	G	2	14/15	0.91	0.20	63,70,85,87	0
3	NAG	K	2	14/15	0.92	0.18	70,87,94,95	0
4	SIA	J	3	20/21	0.93	0.17	62,78,84,85	0
3	NAG	G	1	14/15	0.96	0.14	43,50,59,60	0
4	SIA	H	3	20/21	0.96	0.17	38,55,67,68	0
3	NAG	I	1	14/15	0.96	0.12	64,71,75,81	0
3	NAG	K	1	14/15	0.97	0.14	50,54,60,62	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.

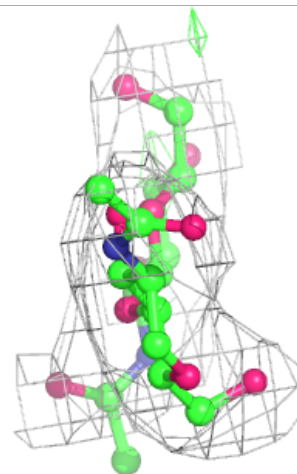
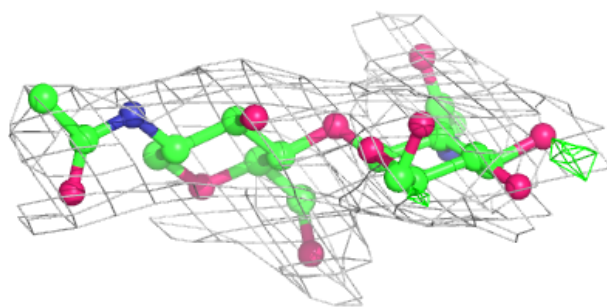
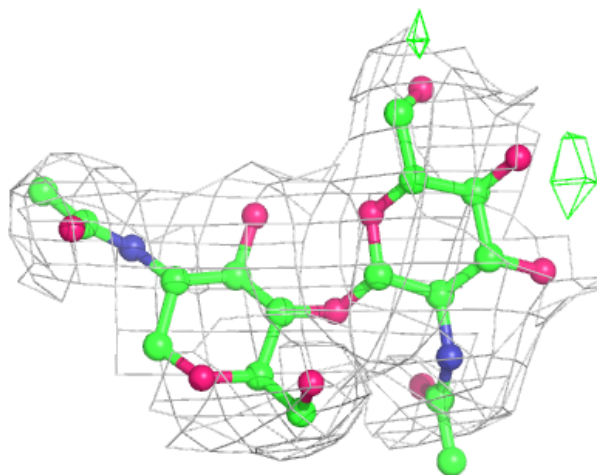
Electron density around Chain G:

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



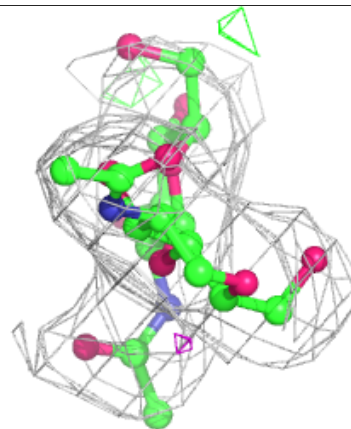
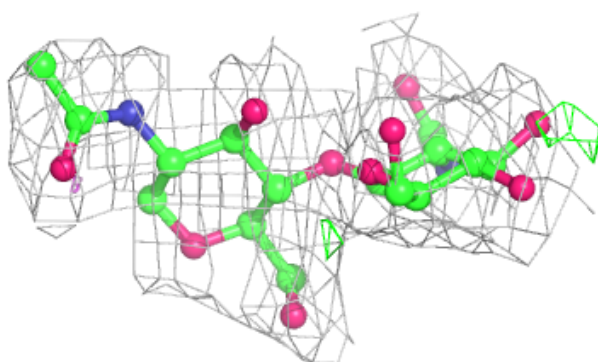
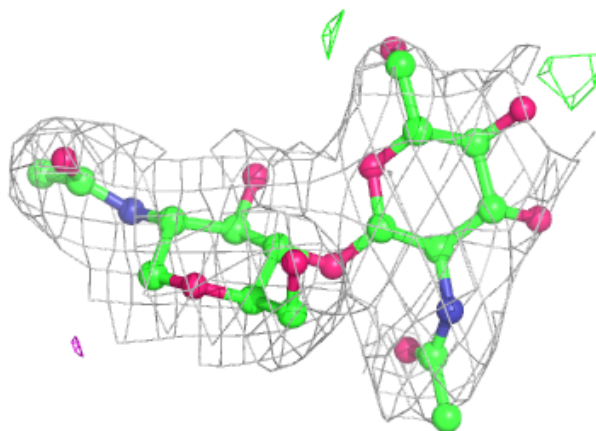
Electron density around Chain I:

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



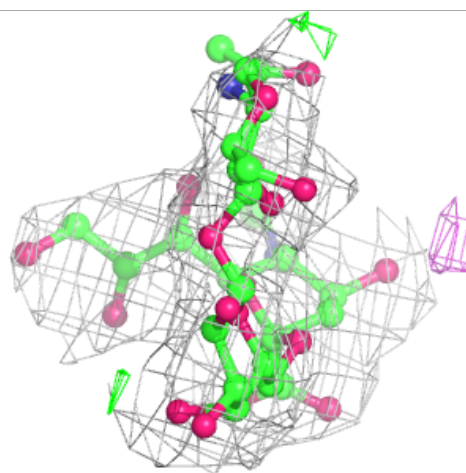
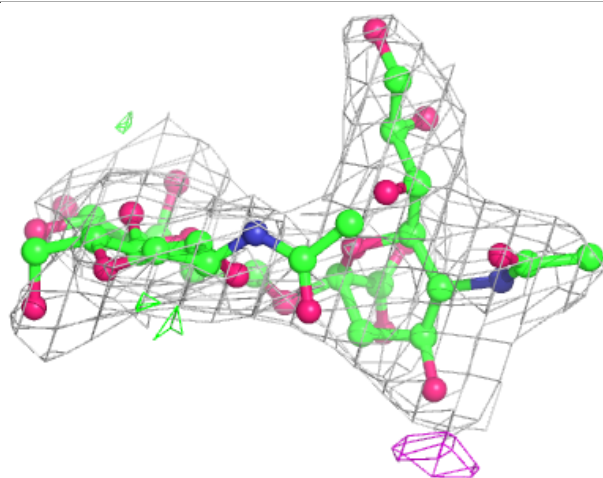
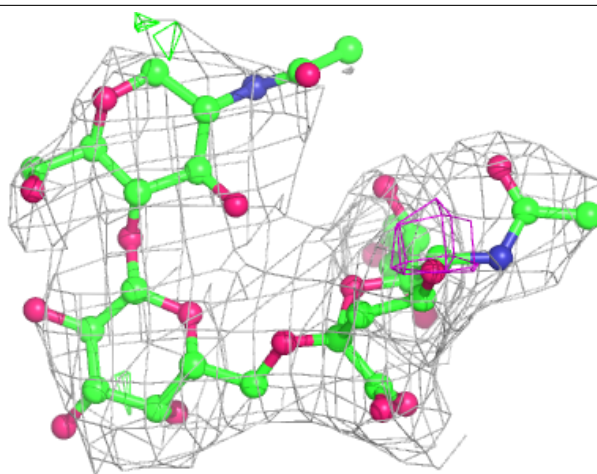
Electron density around Chain K:

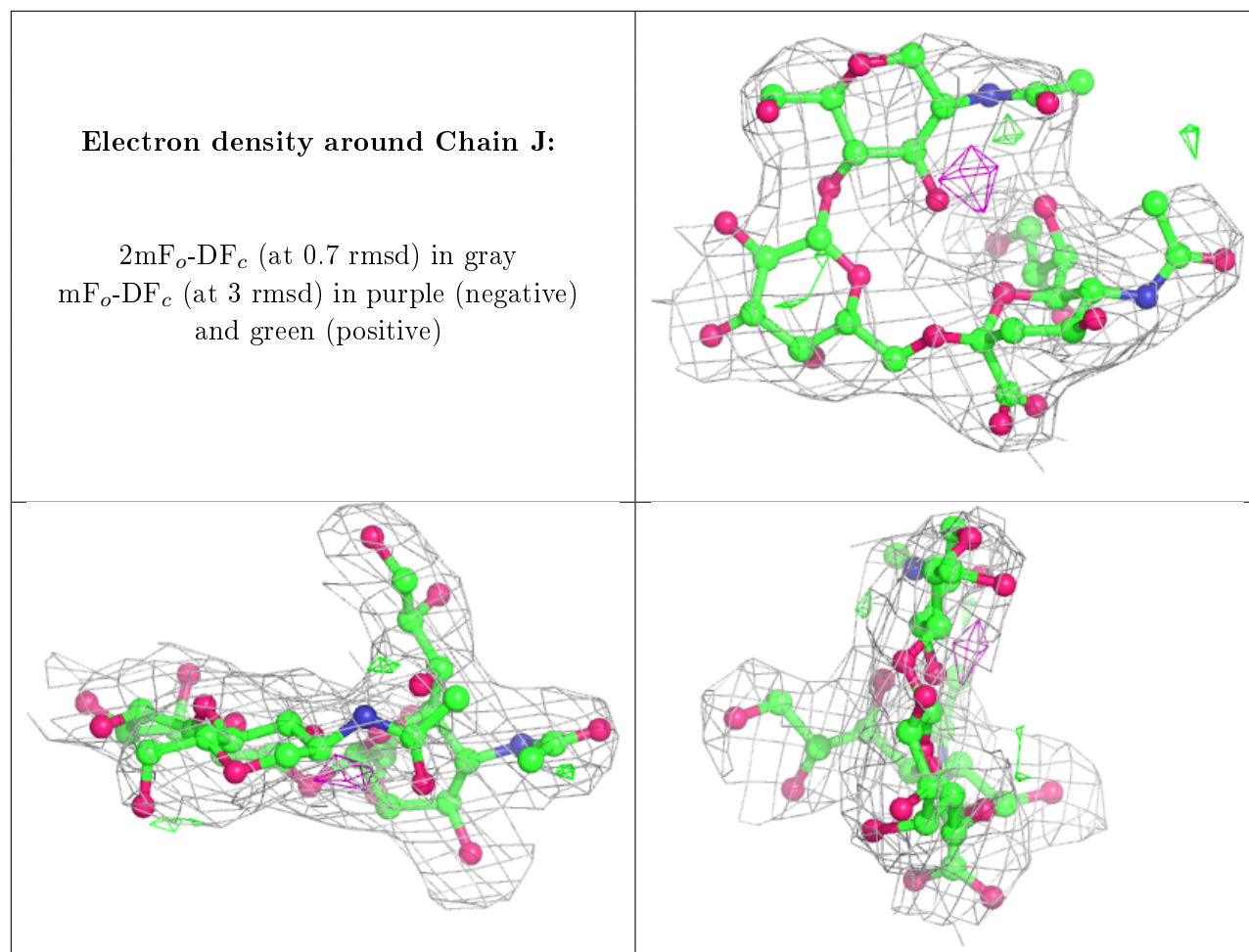
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.