



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 10, 2024 – 09:42 AM EST

PDB ID : 2P37
Title : Crystal structure of a lectin from Canavalia maritima seeds (CML) in complex with man1-3man-OMe
Authors : Moreno, F.B.M.B.; Bezerra, G.A.; Oliveira, T.M.; Souza, E.P.; Rocha, B.A.M.; Benevides, R.G.; Delatorre, P.; Cavada, B.S.; de Azevedo Jr., W.
Deposited on : 2007-03-08
Resolution : 2.10 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 : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36
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) : 2.36

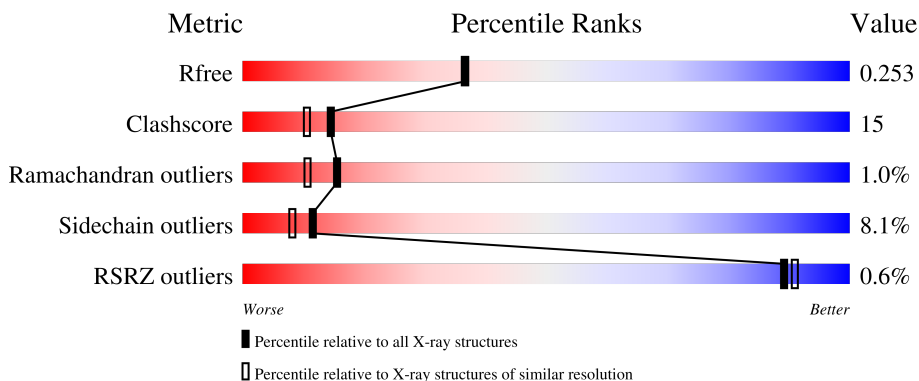
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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.



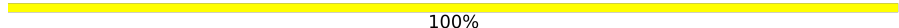
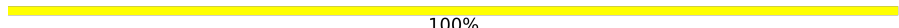

Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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	237	
1	B	237	
1	C	237	
1	D	237	
2	E	2	

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	G	2	 100%
2	H	2	 50% 50%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ZEL	E	1	X	-	-	-
2	ZEL	F	1	X	-	-	-
2	ZEL	G	1	X	-	-	-

2 Entry composition [i](#)

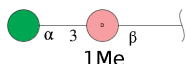
There are 5 unique types of molecules in this entry. The entry contains 7544 atoms, of which 6 are hydrogens and 0 are deuteriums.

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 Concanavalin A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace	
1	A	234	Total	C	N	O	S	0	0	0	
			1775	1120	297	357	1				
1	B	234	Total	C	N	O	S	0	0	0	
			1778	1122	297	358	1				
1	C	232	Total	C	H	N	O	S	0	0	0
			1767	1113	6	294	353	1			
1	D	233	Total	C	N	O	S	0	0	0	
			1770	1118	295	356	1				

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-methyl beta-D-altropyranoside.



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace
2	E	2	Total	C	O	0	0	0
			24	13	11			
2	F	2	Total	C	O	0	0	0
			24	13	11			
2	G	2	Total	C	O	0	0	0
			24	13	11			
2	H	2	Total	C	O	0	0	0
			24	13	11			

- Molecule 3 is MANGANESE (II) ION (three-letter code: MN) (formula: Mn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	Mn	0	0
			1	1		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total Mn 1 1	0	0
3	C	1	Total Mn 1 1	0	0
3	D	1	Total Mn 1 1	0	0

- Molecule 4 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total Ca 1 1	0	0
4	B	1	Total Ca 1 1	0	0
4	C	1	Total Ca 1 1	0	0
4	D	1	Total Ca 1 1	0	0

- Molecule 5 is water.

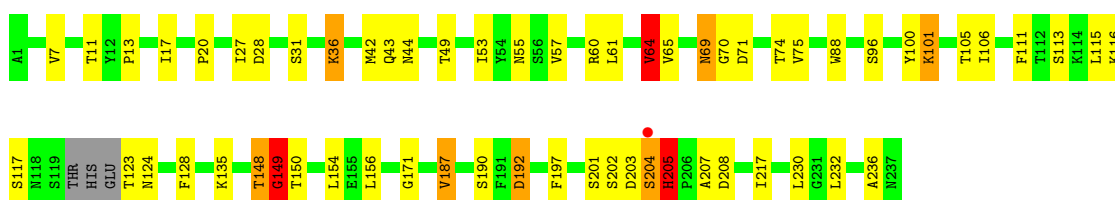
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	105	Total O 105 105	0	0
5	B	114	Total O 114 114	0	0
5	C	64	Total O 64 64	0	0
5	D	67	Total O 67 67	0	0

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.

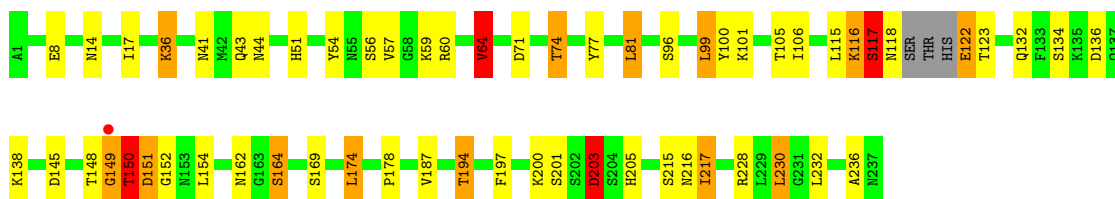
- Molecule 1: Concanavalin A

Chain A: 



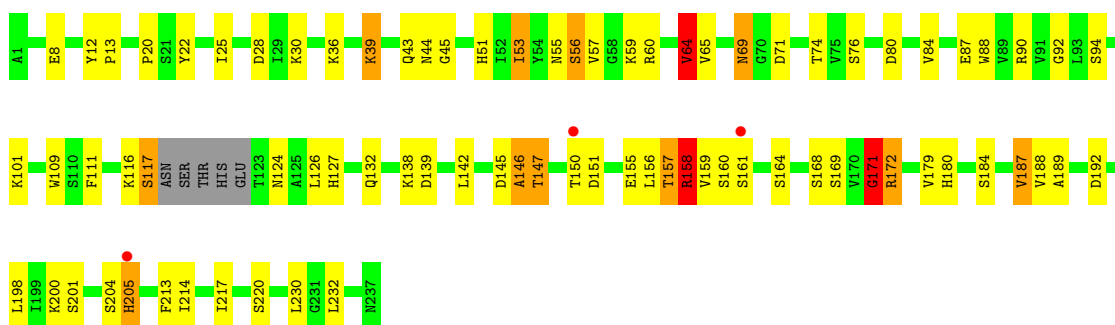
- Molecule 1: Concanavalin A

Chain B: 



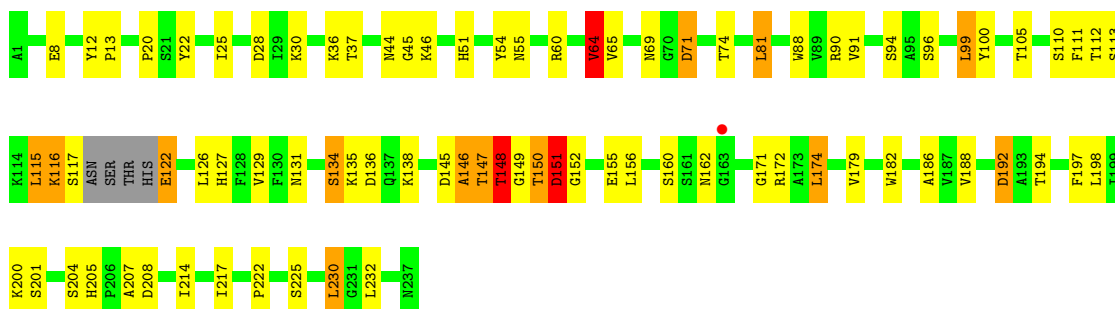
- Molecule 1: Concanavalin A

Chain C: 



- Molecule 1: Concanavalin A

Chain D: 



- Molecule 2: alpha-D-mannopyranose-(1-3)-methyl beta-D-altropyranoside

Chain E: 100%

ZEL1
MAN2

- Molecule 2: alpha-D-mannopyranose-(1-3)-methyl beta-D-altropyranoside

Chain F: 100%

ZEL1
MAN2

- Molecule 2: alpha-D-mannopyranose-(1-3)-methyl beta-D-altropyranoside

Chain G: 100%

ZEL1
MAN2

- Molecule 2: alpha-D-mannopyranose-(1-3)-methyl beta-D-altropyranoside

Chain H: 50%

ZEL1
MAN2

4 Data and refinement statistics i

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	69.39Å 69.39Å 161.30Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	33.92 – 2.10 33.92 – 2.10	Depositor EDS
% Data completeness (in resolution range)	98.9 (33.92-2.10) 98.9 (33.92-2.10)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.10Å)	Xtrriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.187 , 0.255 0.186 , 0.253	Depositor DCC
R_{free} test set	2524 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	26.2	Xtrriage
Anisotropy	0.009	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 18.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	0.022 for -h,-k,l 0.477 for h,-h-k,-l 0.022 for -k,-h,-l	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	7544	wwPDB-VP
Average B, all atoms (Å ²)	22.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.07% 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: CA, MAN, ZEL, MN

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	1.03	1/1814 (0.1%)	1.01	3/2469 (0.1%)
1	B	1.02	2/1817 (0.1%)	1.04	6/2473 (0.2%)
1	C	0.90	0/1800	1.06	12/2450 (0.5%)
1	D	0.92	0/1809	1.02	8/2462 (0.3%)
All	All	0.97	3/7240 (0.0%)	1.03	29/9854 (0.3%)

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	A	0	3
1	B	0	3
1	C	0	4
1	D	0	2
All	All	0	12

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	203	ASP	CB-CG	6.03	1.64	1.51
1	B	8	GLU	CD-OE2	-5.20	1.20	1.25
1	A	149	GLY	N-CA	5.05	1.53	1.46

All (29) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	172	ARG	O-C-N	-10.94	105.20	122.70
1	C	158	ARG	N-CA-C	-9.26	86.01	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	172	ARG	CA-C-N	9.19	137.41	117.20
1	A	205	HIS	N-CA-C	9.03	135.38	111.00
1	C	205	HIS	N-CA-C	8.35	133.55	111.00
1	D	148	THR	N-CA-C	8.34	133.52	111.00
1	B	151	ASP	N-CA-C	8.06	132.76	111.00
1	C	172	ARG	C-N-CA	8.01	141.72	121.70
1	D	64	VAL	CB-CA-C	-7.93	96.34	111.40
1	B	174	LEU	CA-CB-CG	7.88	133.44	115.30
1	D	146	ALA	C-N-CA	7.83	141.28	121.70
1	B	64	VAL	CB-CA-C	-7.75	96.68	111.40
1	C	64	VAL	CB-CA-C	-6.58	98.89	111.40
1	C	161	SER	N-CA-C	-6.49	93.47	111.00
1	C	172	ARG	CB-CA-C	-6.42	97.56	110.40
1	D	205	HIS	N-CA-CB	-6.18	99.47	110.60
1	A	61	LEU	CB-CG-CD2	-6.13	100.59	111.00
1	D	146	ALA	CA-C-N	5.98	130.36	117.20
1	D	174	LEU	CA-CB-CG	5.98	129.05	115.30
1	D	205	HIS	N-CA-C	5.95	127.06	111.00
1	A	64	VAL	CB-CA-C	-5.79	100.40	111.40
1	C	171	GLY	O-C-N	-5.76	113.48	122.70
1	D	115	LEU	CB-CG-CD1	-5.61	101.47	111.00
1	B	203	ASP	CB-CG-OD1	5.60	123.34	118.30
1	C	146	ALA	C-N-CA	5.48	135.41	121.70
1	C	205	HIS	N-CA-CB	-5.22	101.20	110.60
1	C	139	ASP	CB-CG-OD1	5.14	122.93	118.30
1	B	174	LEU	CB-CG-CD2	5.13	119.72	111.00
1	B	64	VAL	CG1-CB-CG2	5.06	118.99	110.90

There are no chirality outliers.

All (12) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	148	THR	Peptide
1	A	149	GLY	Peptide
1	A	204	SER	Peptide
1	B	116	LYS	Peptide
1	B	117	SER	Peptide
1	B	150	THR	Peptide
1	C	146	ALA	Peptide
1	C	157	THR	Peptide
1	C	171	GLY	Mainchain
1	C	204	SER	Peptide

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Mol	Chain	Res	Type	Group
1	D	147	THR	Peptide
1	D	204	SER	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	1775	0	1727	55	0
1	B	1778	0	1728	54	0
1	C	1761	6	1716	65	0
1	D	1770	0	1722	65	0
2	E	24	0	10	0	0
2	F	24	0	10	0	0
2	G	24	0	10	0	0
2	H	24	0	10	1	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	A	105	0	0	2	0
5	B	114	0	0	10	0
5	C	64	0	0	5	0
5	D	67	0	0	6	0
All	All	7538	6	6933	218	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 15.

All (218) 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:B:60:ARG:HH11	1:D:60:ARG:NH1	1.09	1.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:ARG:NH1	1:D:60:ARG:NH1	1.89	1.20
1:B:215:SER:HB2	5:B:310:HOH:O	1.43	1.18
1:C:232:LEU:HG	5:C:275:HOH:O	1.45	1.16
1:B:59:LYS:HE2	5:B:285:HOH:O	1.48	1.13
1:B:117:SER:HB2	1:B:118:ASN:HB2	1.26	1.12
1:D:150:THR:HA	1:D:152:GLY:N	1.67	1.08
1:B:117:SER:HB2	1:B:118:ASN:CB	1.86	1.06
1:B:60:ARG:NH1	1:D:60:ARG:HH11	1.50	1.05
1:A:203:ASP:OD2	1:A:205:HIS:CG	2.12	1.03
1:A:60:ARG:HH21	1:C:60:ARG:NH2	1.59	1.01
1:D:150:THR:HA	1:D:152:GLY:H	0.85	1.00
1:C:88:TRP:HB3	1:C:217:ILE:HD11	1.44	0.99
1:C:44:ASN:HD21	1:C:201:SER:H	1.11	0.99
1:A:201:SER:OG	1:A:205:HIS:HD2	1.51	0.94
1:C:117:SER:HA	1:C:187:VAL:HG12	1.47	0.93
1:A:60:ARG:HH21	1:C:60:ARG:HH21	1.11	0.93
1:A:44:ASN:HD21	1:A:201:SER:H	1.11	0.89
1:D:150:THR:CA	1:D:152:GLY:H	1.81	0.89
1:C:117:SER:HA	1:C:187:VAL:CG1	2.05	0.86
1:B:117:SER:CB	1:B:118:ASN:HB2	2.05	0.85
1:A:203:ASP:OD2	1:A:205:HIS:ND1	2.09	0.84
1:D:44:ASN:HD21	1:D:201:SER:H	1.24	0.84
1:B:44:ASN:HD21	1:B:201:SER:H	1.26	0.83
1:A:201:SER:OG	1:A:205:HIS:CD2	2.33	0.81
1:A:203:ASP:OD2	1:A:205:HIS:HB2	1.82	0.80
1:D:90:ARG:NH1	1:D:217:ILE:HG23	1.97	0.80
1:A:101:LYS:O	1:A:205:HIS:NE2	2.16	0.79
1:A:60:ARG:NH2	1:C:60:ARG:HH21	1.83	0.77
1:A:124:ASN:HD21	1:B:132:GLN:H	1.30	0.77
1:A:44:ASN:HD21	1:A:201:SER:N	1.83	0.77
1:C:172:ARG:NH1	1:C:213:PHE:CE2	2.52	0.77
1:A:203:ASP:OD2	1:A:205:HIS:CB	2.33	0.76
1:C:90:ARG:NH1	1:C:217:ILE:HG23	2.02	0.75
1:A:60:ARG:NH2	1:C:60:ARG:NH2	2.32	0.74
1:A:88:TRP:HB3	1:A:217:ILE:HD11	1.68	0.74
1:B:150:THR:HA	1:B:152:GLY:H	1.53	0.73
1:C:44:ASN:HD21	1:C:201:SER:N	1.87	0.73
1:A:203:ASP:CG	1:A:205:HIS:HB2	2.09	0.72
1:B:117:SER:HB2	1:B:118:ASN:CA	2.19	0.72
1:B:194:THR:CG2	5:B:303:HOH:O	2.36	0.72
1:B:60:ARG:HH11	1:D:60:ARG:HH11	0.74	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:194:THR:HG23	5:B:303:HOH:O	1.89	0.72
1:C:142:LEU:HD13	1:C:147:THR:H	1.55	0.72
1:B:71:ASP:HB2	5:B:325:HOH:O	1.89	0.71
1:C:147:THR:HB	5:C:271:HOH:O	1.91	0.71
1:C:69:ASN:HD21	1:C:71:ASP:HB2	1.58	0.69
1:A:117:SER:HA	1:A:187:VAL:HG13	1.76	0.68
1:D:149:GLY:CA	1:D:150:THR:O	2.42	0.68
1:C:51:HIS:HB2	1:C:64:VAL:HG23	1.75	0.67
1:B:56:SER:O	1:B:59:LYS:HE3	1.94	0.66
1:B:178:PRO:HB3	1:B:217:ILE:HD11	1.78	0.66
1:B:215:SER:CB	5:B:310:HOH:O	2.17	0.65
1:D:145:ASP:O	5:D:265:HOH:O	2.14	0.65
1:D:156:LEU:O	1:D:171:GLY:HA3	1.96	0.65
1:C:156:LEU:O	1:C:171:GLY:HA3	1.97	0.65
1:B:118:ASN:HB3	1:B:187:VAL:HG21	1.80	0.64
1:A:71:ASP:HB2	5:A:296:HOH:O	1.98	0.64
1:D:149:GLY:HA3	1:D:150:THR:C	2.18	0.64
1:A:43:GLN:HE22	1:A:69:ASN:HD21	1.47	0.63
1:A:101:LYS:O	1:A:205:HIS:CD2	2.50	0.63
1:C:230:LEU:O	1:C:232:LEU:HD13	1.99	0.63
1:C:53:ILE:HD13	1:C:53:ILE:O	1.98	0.63
1:A:43:GLN:HE22	1:A:69:ASN:ND2	1.97	0.63
1:A:57:VAL:HG21	1:C:64:VAL:HG22	1.80	0.62
1:A:101:LYS:H	1:A:205:HIS:CE1	2.17	0.62
1:C:45:GLY:HA2	1:C:198:LEU:HD21	1.81	0.62
1:A:106:ILE:HB	1:A:154:LEU:HB3	1.83	0.61
1:D:149:GLY:HA2	1:D:150:THR:O	2.01	0.61
1:C:53:ILE:HD13	1:C:53:ILE:C	2.22	0.60
1:C:55:ASN:HA	5:C:268:HOH:O	2.00	0.60
1:B:203:ASP:C	1:B:205:HIS:H	2.05	0.59
1:D:54:TYR:CE1	1:D:81:LEU:HD22	2.37	0.59
1:D:36:LYS:HE2	5:D:251:HOH:O	2.03	0.58
1:D:150:THR:HB	1:D:151:ASP:HA	1.85	0.58
1:B:162:ASN:OD1	1:B:164:SER:HB2	2.04	0.58
1:A:74:THR:HG23	1:C:57:VAL:HG13	1.85	0.58
1:A:69:ASN:HD22	1:A:70:GLY:N	2.02	0.58
1:C:172:ARG:HD2	1:C:213:PHE:CZ	2.39	0.57
1:C:60:ARG:HD3	1:C:76:SER:HB3	1.87	0.57
1:B:64:VAL:HG13	1:B:74:THR:HG22	1.86	0.57
1:D:116:LYS:HG3	1:D:188:VAL:HB	1.86	0.57
1:C:56:SER:CB	1:C:189:ALA:H	2.18	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:57:VAL:CG2	1:D:64:VAL:HG22	2.34	0.57
1:C:43:GLN:HE22	1:C:69:ASN:ND2	2.03	0.57
1:C:69:ASN:ND2	1:C:71:ASP:H	2.02	0.57
1:D:45:GLY:HA2	1:D:198:LEU:HD21	1.86	0.56
1:B:136:ASP:OD2	1:B:138:LYS:HE2	2.06	0.56
1:C:124:ASN:ND2	1:D:131:ASN:H	2.03	0.56
1:C:94:SER:HB3	1:C:232:LEU:HD11	1.88	0.56
1:B:115:LEU:O	1:B:123:THR:HA	2.06	0.56
1:B:118:ASN:HB3	1:B:187:VAL:CG2	2.36	0.56
1:A:36:LYS:HD2	1:A:75:VAL:HG23	1.88	0.55
1:B:43:GLN:NE2	1:B:71:ASP:HB3	2.22	0.55
1:D:146:ALA:HA	1:D:155:GLU:O	2.07	0.55
1:D:51:HIS:HB2	1:D:64:VAL:HG23	1.87	0.55
1:C:90:ARG:HH12	1:C:217:ILE:HG23	1.71	0.55
1:C:147:THR:HG22	1:C:155:GLU:HB2	1.89	0.55
1:D:126:LEU:CD2	1:D:179:VAL:HG22	2.36	0.55
1:A:111:PHE:CE2	1:A:113:SER:HB2	2.44	0.53
1:D:117:SER:H	1:D:186:ALA:HA	1.73	0.53
1:B:44:ASN:HD21	1:B:201:SER:N	2.01	0.53
1:D:110:SER:HB3	1:D:129:VAL:HG13	1.90	0.53
1:D:149:GLY:HA3	1:D:150:THR:O	2.08	0.53
1:D:91:VAL:HG12	1:D:214:ILE:HG12	1.91	0.53
1:A:57:VAL:CG2	1:C:64:VAL:HG22	2.39	0.53
1:C:30:LYS:HD2	1:C:84:VAL:HG13	1.91	0.52
1:C:92:GLY:HA2	1:C:109:TRP:CH2	2.45	0.52
1:C:44:ASN:ND2	1:C:201:SER:H	1.94	0.52
1:C:172:ARG:HD2	1:C:213:PHE:CE1	2.45	0.52
1:D:96:SER:OG	1:D:230:LEU:HA	2.10	0.52
1:B:36:LYS:HE2	1:B:77:TYR:CD1	2.46	0.51
1:C:88:TRP:CB	1:C:217:ILE:HD11	2.29	0.51
1:B:106:ILE:HB	1:B:154:LEU:HB3	1.93	0.51
1:D:30:LYS:HA	5:D:243:HOH:O	2.10	0.51
1:B:57:VAL:HG21	1:D:64:VAL:HG22	1.92	0.51
1:A:7:VAL:HG22	1:A:27:ILE:CD1	2.41	0.51
1:C:36:LYS:HE3	1:C:76:SER:O	2.11	0.51
1:B:122:GLU:OE1	1:B:123:THR:OG1	2.29	0.51
1:C:20:PRO:HB2	1:C:22:TYR:CZ	2.46	0.51
1:D:8:GLU:OE1	1:D:28:ASP:OD2	2.29	0.51
1:A:111:PHE:HB3	1:A:128:PHE:CZ	2.47	0.50
1:C:145:ASP:HB3	1:C:158:ARG:HG2	1.93	0.50
1:D:54:TYR:CD1	1:D:81:LEU:HD22	2.46	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:HIS:CG	1:A:205:HIS:O	2.65	0.49
1:A:74:THR:HG23	1:C:57:VAL:CG1	2.42	0.49
1:B:96:SER:OG	1:B:230:LEU:HA	2.12	0.49
1:D:222:PRO:HD2	1:D:225:SER:OG	2.11	0.49
1:D:69:ASN:OD1	1:D:71:ASP:HB2	2.12	0.49
1:B:215:SER:CA	5:B:310:HOH:O	2.53	0.49
1:C:116:LYS:HG3	1:C:188:VAL:HB	1.94	0.49
1:B:60:ARG:HH11	1:D:60:ARG:CZ	2.06	0.49
1:C:8:GLU:OE1	1:C:28:ASP:OD2	2.31	0.49
1:B:99:LEU:HD13	1:B:100:TYR:CZ	2.47	0.49
1:A:100:TYR:HB3	1:A:205:HIS:O	2.13	0.48
1:B:105:THR:O	1:B:197:PHE:HA	2.13	0.48
1:B:51:HIS:HB2	1:B:64:VAL:HG23	1.94	0.48
1:D:99:LEU:HB2	2:H:2:MAN:C1	2.42	0.48
1:D:149:GLY:CA	1:D:150:THR:C	2.80	0.48
1:D:111:PHE:CE2	1:D:113:SER:HB2	2.49	0.48
1:D:117:SER:HA	1:D:122:GLU:O	2.12	0.48
1:C:43:GLN:HE22	1:C:69:ASN:HD21	1.60	0.47
1:B:203:ASP:C	1:B:205:HIS:N	2.67	0.47
1:C:74:THR:O	5:C:259:HOH:O	2.20	0.47
1:D:20:PRO:HB2	1:D:22:TYR:CZ	2.49	0.47
1:A:192:ASP:HB2	5:A:309:HOH:O	2.14	0.47
1:A:148:THR:O	1:A:149:GLY:O	2.33	0.47
1:A:203:ASP:C	1:A:205:HIS:HB2	2.35	0.47
1:B:54:TYR:CE1	1:B:81:LEU:HD22	2.50	0.47
1:D:136:ASP:OD2	1:D:138:LYS:NZ	2.44	0.47
1:A:115:LEU:O	1:A:123:THR:HA	2.15	0.46
1:C:25:ILE:HG21	1:C:65:VAL:HG21	1.97	0.46
1:D:100:TYR:HB2	1:D:207:ALA:HB3	1.97	0.46
1:B:145:ASP:OD2	1:B:169:SER:OG	2.31	0.46
1:C:87:GLU:HG3	1:C:180:HIS:CD2	2.50	0.46
1:D:147:THR:HB	5:D:265:HOH:O	2.14	0.46
1:D:90:ARG:HH12	1:D:217:ILE:HG23	1.75	0.46
1:A:105:THR:O	1:A:197:PHE:HA	2.15	0.46
1:B:148:THR:O	1:B:149:GLY:O	2.32	0.46
1:B:203:ASP:CG	5:B:269:HOH:O	2.53	0.46
1:D:192:ASP:HB2	5:D:287:HOH:O	2.15	0.46
1:D:105:THR:O	1:D:197:PHE:HA	2.16	0.46
1:A:17:ILE:HG21	1:A:236:ALA:HB1	1.98	0.45
1:D:160:SER:C	1:D:162:ASN:H	2.19	0.45
1:B:116:LYS:HG2	1:B:123:THR:HG23	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:126:LEU:CD2	1:C:179:VAL:HG22	2.47	0.45
1:A:17:ILE:HG21	1:A:236:ALA:CB	2.46	0.45
1:B:118:ASN:CB	1:B:187:VAL:HG21	2.45	0.45
1:D:160:SER:C	1:D:162:ASN:N	2.69	0.45
1:A:88:TRP:HB3	1:A:217:ILE:CD1	2.44	0.45
1:C:138:LYS:HB2	1:D:88:TRP:CE3	2.52	0.45
1:A:36:LYS:HD2	1:A:75:VAL:CG2	2.47	0.45
1:A:96:SER:OG	1:A:230:LEU:HA	2.17	0.44
1:D:74:THR:HG23	5:D:244:HOH:O	2.16	0.44
1:D:134:SER:O	1:D:148:THR:HB	2.18	0.44
1:D:207:ALA:HA	1:D:208:ASP:HA	1.83	0.44
1:B:149:GLY:HA3	1:B:150:THR:C	2.39	0.43
1:D:45:GLY:HA3	1:D:200:LYS:HE3	2.00	0.43
1:C:30:LYS:HA	5:C:250:HOH:O	2.18	0.43
1:D:44:ASN:ND2	1:D:201:SER:H	2.04	0.43
1:B:17:ILE:HG21	1:B:236:ALA:CB	2.49	0.43
1:D:46:LYS:HD3	1:D:46:LYS:HA	1.83	0.43
1:C:59:LYS:HD3	1:C:80:ASP:HB2	2.00	0.43
1:A:116:LYS:HG2	1:A:123:THR:HG23	2.00	0.42
1:A:156:LEU:O	1:A:171:GLY:HA3	2.19	0.42
1:C:12:TYR:HA	1:C:13:PRO:HD3	1.88	0.42
1:D:25:ILE:O	1:D:37:THR:HA	2.19	0.42
1:B:44:ASN:ND2	1:B:200:LYS:HA	2.34	0.42
1:A:28:ASP:HB3	1:A:31:SER:O	2.20	0.42
1:A:207:ALA:HA	1:A:208:ASP:HA	1.87	0.42
1:C:44:ASN:ND2	1:C:200:LYS:HA	2.34	0.42
1:C:111:PHE:O	1:C:127:HIS:HA	2.20	0.42
1:D:94:SER:OG	1:D:172:ARG:HG2	2.20	0.42
1:A:111:PHE:HB3	1:A:128:PHE:CE2	2.55	0.42
1:D:12:TYR:HA	1:D:13:PRO:HD3	1.93	0.42
1:A:64:VAL:HG22	1:C:57:VAL:HG21	2.02	0.41
1:C:124:ASN:HD22	1:C:124:ASN:N	2.18	0.41
1:C:39:LYS:HE2	1:C:39:LYS:HB2	1.47	0.41
1:A:13:PRO:HB3	1:A:20:PRO:O	2.19	0.41
1:D:81:LEU:HD12	1:D:81:LEU:HA	1.97	0.41
1:A:11:THR:HG22	1:A:42:MET:HG3	2.03	0.41
1:A:57:VAL:HG21	1:C:64:VAL:CG2	2.50	0.41
1:B:36:LYS:HE2	1:B:77:TYR:HD1	1.85	0.41
1:C:160:SER:O	1:C:164:SER:HB3	2.21	0.41
1:A:49:THR:O	1:A:65:VAL:HA	2.21	0.41
1:B:215:SER:HA	5:B:310:HOH:O	2.18	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:157:THR:HB	1:C:169:SER:OG	2.21	0.41
1:D:25:ILE:HG21	1:D:65:VAL:HG21	2.03	0.41
1:C:94:SER:HA	1:C:172:ARG:HA	2.03	0.41
1:C:158:ARG:HD3	1:C:158:ARG:HA	1.88	0.41
1:D:115:LEU:HD21	1:D:182:TRP:HA	2.03	0.40
1:A:55:ASN:HA	1:A:190:SER:HA	2.03	0.40
1:D:126:LEU:HD23	1:D:179:VAL:HG22	2.03	0.40
1:B:194:THR:HG22	5:B:303:HOH:O	2.15	0.40
1:C:138:LYS:HD2	1:D:217:ILE:CD1	2.51	0.40
1:D:112:THR:OG1	1:D:127:HIS:ND1	2.53	0.40
1:B:14:ASN:ND2	1:B:228:ARG:HB3	2.37	0.40
1:B:117:SER:H	1:B:187:VAL:HB	1.86	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	230/237 (97%)	217 (94%)	11 (5%)	2 (1%)	17 12
1	B	230/237 (97%)	219 (95%)	8 (4%)	3 (1%)	12 7
1	C	228/237 (96%)	216 (95%)	11 (5%)	1 (0%)	34 32
1	D	229/237 (97%)	218 (95%)	8 (4%)	3 (1%)	12 7
All	All	917/948 (97%)	870 (95%)	38 (4%)	9 (1%)	15 11

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	150	THR
1	B	151	ASP
1	C	147	THR

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Mol	Chain	Res	Type
1	A	149	GLY
1	A	205	HIS
1	B	149	GLY
1	D	150	THR
1	D	148	THR
1	D	151	ASP

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	199/202 (98%)	187 (94%)	12 (6%)	19	16
1	B	199/202 (98%)	181 (91%)	18 (9%)	9	6
1	C	197/202 (98%)	178 (90%)	19 (10%)	8	5
1	D	198/202 (98%)	183 (92%)	15 (8%)	13	10
All	All	793/808 (98%)	729 (92%)	64 (8%)	11	8

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	LYS
1	A	53	ILE
1	A	64	VAL
1	A	69	ASN
1	A	101	LYS
1	A	135	LYS
1	A	150	THR
1	A	187	VAL
1	A	192	ASP
1	A	202	SER
1	A	204	SER
1	A	232	LEU
1	B	36	LYS
1	B	41	ASN
1	B	64	VAL

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Mol	Chain	Res	Type
1	B	74	THR
1	B	81	LEU
1	B	99	LEU
1	B	101	LYS
1	B	117	SER
1	B	122	GLU
1	B	134	SER
1	B	164	SER
1	B	174	LEU
1	B	194	THR
1	B	203	ASP
1	B	216	ASN
1	B	217	ILE
1	B	230	LEU
1	B	232	LEU
1	C	39	LYS
1	C	53	ILE
1	C	56	SER
1	C	64	VAL
1	C	69	ASN
1	C	101	LYS
1	C	117	SER
1	C	132	GLN
1	C	150	THR
1	C	151	ASP
1	C	158	ARG
1	C	159	VAL
1	C	168	SER
1	C	184	SER
1	C	187	VAL
1	C	192	ASP
1	C	205	HIS
1	C	214	ILE
1	C	220	SER
1	D	55	ASN
1	D	64	VAL
1	D	71	ASP
1	D	81	LEU
1	D	99	LEU
1	D	116	LYS
1	D	122	GLU
1	D	134	SER

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Mol	Chain	Res	Type
1	D	135	LYS
1	D	151	ASP
1	D	174	LEU
1	D	192	ASP
1	D	194	THR
1	D	230	LEU
1	D	232	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (21) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	44	ASN
1	A	69	ASN
1	A	83	ASN
1	A	104	ASN
1	A	124	ASN
1	A	166	GLN
1	A	205	HIS
1	B	41	ASN
1	B	44	ASN
1	B	51	HIS
1	B	104	ASN
1	B	132	GLN
1	B	166	GLN
1	C	44	ASN
1	C	69	ASN
1	C	104	ASN
1	C	124	ASN
1	C	132	GLN
1	D	44	ASN
1	D	104	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

8 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
2	ZEL	E	1	2	13,13,13	0.61	0	18,18,18	2.14	5 (27%)
2	MAN	E	2	2	11,11,12	0.61	0	15,15,17	1.10	1 (6%)
2	ZEL	F	1	2	13,13,13	0.86	0	18,18,18	3.24	7 (38%)
2	MAN	F	2	2	11,11,12	1.04	1 (9%)	15,15,17	1.83	3 (20%)
2	ZEL	G	1	2	13,13,13	0.64	0	18,18,18	3.19	7 (38%)
2	MAN	G	2	2	11,11,12	0.79	0	15,15,17	1.47	3 (20%)
2	ZEL	H	1	2	13,13,13	0.72	0	18,18,18	2.92	9 (50%)
2	MAN	H	2	2	11,11,12	1.08	1 (9%)	15,15,17	1.37	4 (26%)

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
2	ZEL	E	1	2	1/1/5/5	2/4/24/24	0/1/1/1
2	MAN	E	2	2	-	0/2/19/22	0/1/1/1
2	ZEL	F	1	2	1/1/5/5	2/4/24/24	0/1/1/1
2	MAN	F	2	2	-	0/2/19/22	0/1/1/1
2	ZEL	G	1	2	1/1/5/5	0/4/24/24	0/1/1/1
2	MAN	G	2	2	-	0/2/19/22	0/1/1/1
2	ZEL	H	1	2	-	4/4/24/24	0/1/1/1
2	MAN	H	2	2	-	1/2/19/22	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	MAN	C2-C3	2.38	1.56	1.52
2	H	2	MAN	C2-C3	2.13	1.55	1.52

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1	ZEL	O1-C1-C2	9.69	119.51	108.15
2	F	1	ZEL	C1-C2-C3	-7.82	93.72	110.00
2	F	1	ZEL	O1-C1-C2	6.94	116.28	108.15
2	H	1	ZEL	O1-C1-C2	6.79	116.11	108.15
2	G	1	ZEL	C1-C2-C3	-6.07	97.36	110.00
2	H	1	ZEL	C4-C3-C2	5.90	121.12	110.82
2	E	1	ZEL	C1-C2-C3	-5.53	98.48	110.00
2	F	1	ZEL	C1-O5-C5	-4.81	104.26	113.69
2	G	1	ZEL	C1-O5-C5	-4.14	105.56	113.69
2	F	1	ZEL	O5-C5-C6	4.11	116.64	106.44
2	F	1	ZEL	O2-C2-C1	3.97	119.68	110.05
2	E	1	ZEL	O1-C1-C2	3.79	112.60	108.15
2	F	2	MAN	O2-C2-C1	-3.66	101.66	109.15
2	H	1	ZEL	O3-C3-C4	3.47	118.37	110.35
2	H	1	ZEL	C1M-O1-C1	-3.44	107.97	113.27
2	E	1	ZEL	O5-C5-C6	3.43	114.96	106.44
2	G	2	MAN	O3-C3-C4	3.37	118.14	110.35
2	F	1	ZEL	O5-C1-O1	3.32	118.67	110.97
2	F	2	MAN	C1-O5-C5	3.19	116.51	112.19
2	F	2	MAN	O5-C1-C2	-3.17	105.88	110.77
2	E	1	ZEL	O2-C2-C1	3.12	117.63	110.05
2	H	1	ZEL	C1-O5-C5	-2.79	108.20	113.69
2	H	1	ZEL	C3-C4-C5	2.74	115.12	110.24
2	E	2	MAN	C1-O5-C5	2.72	115.87	112.19
2	G	2	MAN	O5-C1-C2	2.69	114.92	110.77
2	F	1	ZEL	C1M-O1-C1	2.68	117.41	113.27
2	G	1	ZEL	C1M-O1-C1	-2.64	109.19	113.27
2	H	1	ZEL	O4-C4-C3	2.62	116.40	110.35
2	G	1	ZEL	O2-C2-C1	2.58	116.30	110.05
2	H	2	MAN	O3-C3-C4	2.46	116.04	110.35
2	E	1	ZEL	C1M-O1-C1	-2.43	109.52	113.27
2	H	1	ZEL	O2-C2-C3	-2.35	104.91	110.35
2	H	2	MAN	C1-O5-C5	2.34	115.36	112.19
2	H	2	MAN	O4-C4-C5	-2.24	103.74	109.30
2	G	1	ZEL	C4-C3-C2	2.18	114.62	110.82
2	H	2	MAN	O5-C5-C6	2.16	110.58	107.20
2	H	1	ZEL	C6-C5-C4	-2.15	107.96	113.00
2	G	1	ZEL	O3-C3-C2	-2.09	105.52	110.35
2	G	2	MAN	O3-C3-C2	-2.08	106.00	109.99

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	E	1	ZEL	C3
2	F	1	ZEL	C3
2	G	1	ZEL	C3

All (9) torsion outliers are listed below:

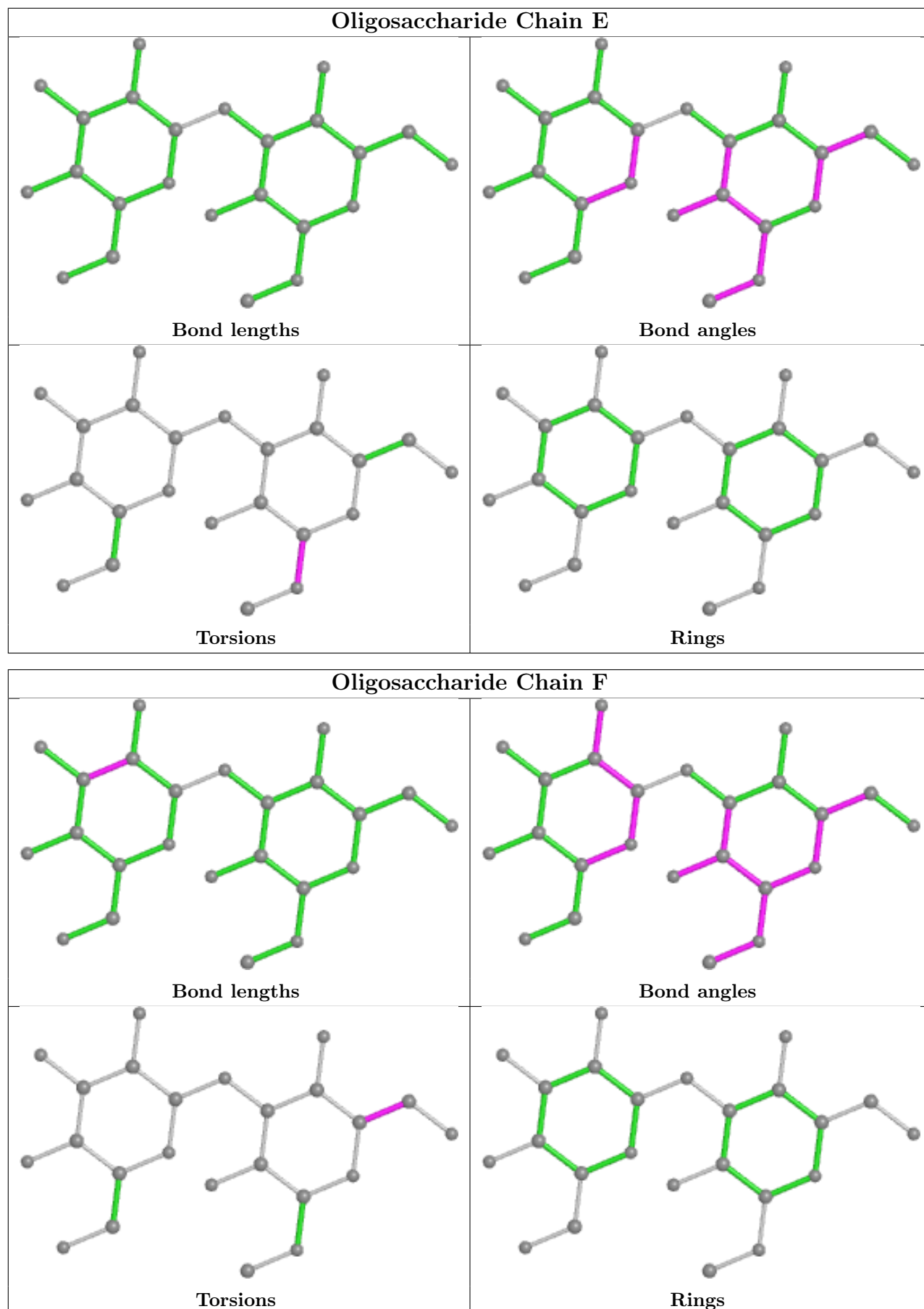
Mol	Chain	Res	Type	Atoms
2	H	1	ZEL	O5-C1-O1-C1M
2	H	1	ZEL	C2-C1-O1-C1M
2	H	1	ZEL	O5-C5-C6-O6
2	H	1	ZEL	C4-C5-C6-O6
2	E	1	ZEL	C2-C1-O1-C1M
2	F	1	ZEL	O5-C5-C6-O6
2	E	1	ZEL	O5-C1-O1-C1M
2	H	2	MAN	C4-C5-C6-O6
2	F	1	ZEL	C4-C5-C6-O6

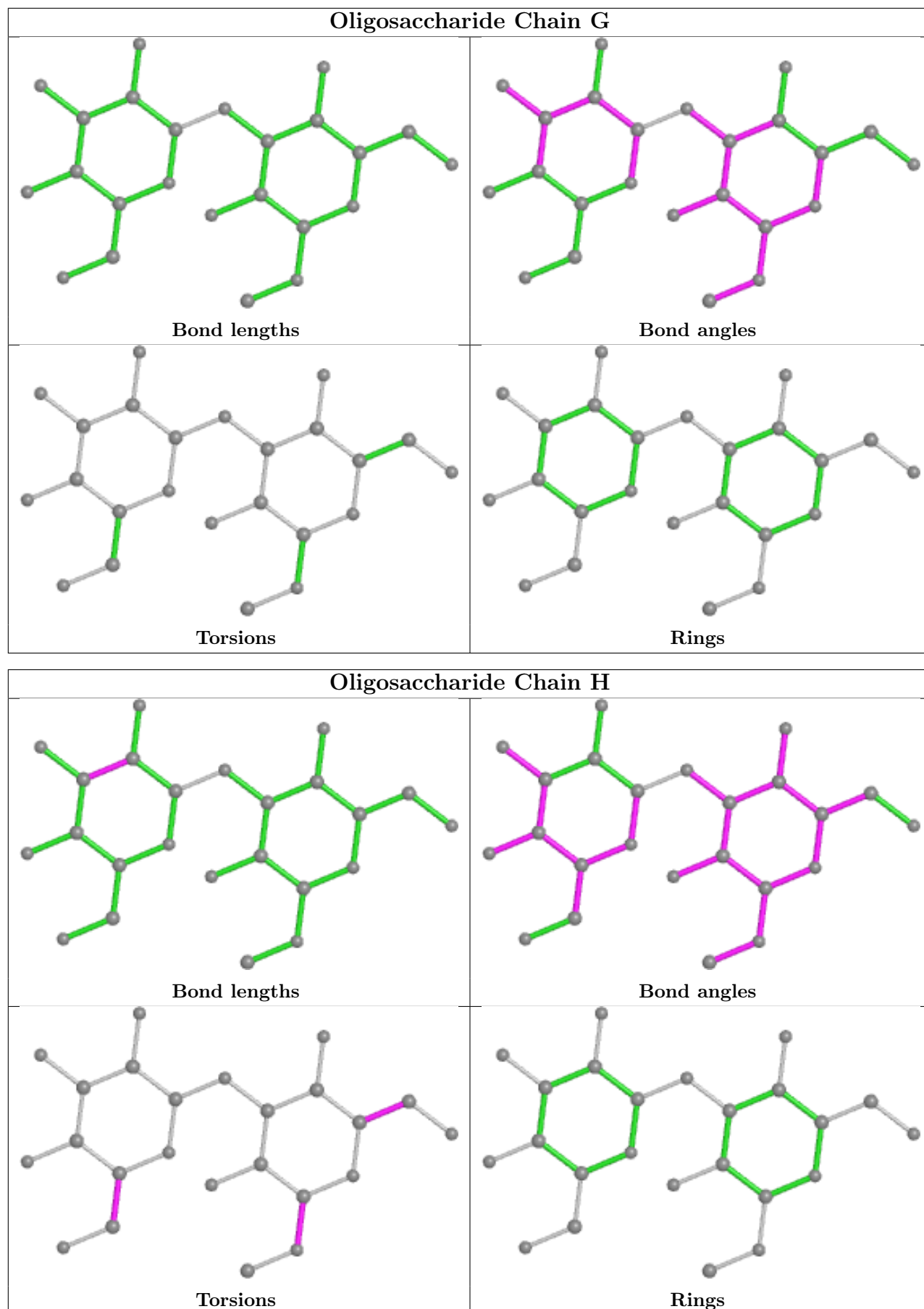
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	2	MAN	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](#)

Of 8 ligands modelled in this entry, 8 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

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 [i](#)

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

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	234/237 (98%)	-0.41	1 (0%) 92 93	12, 18, 31, 47	6 (2%)
1	B	234/237 (98%)	-0.39	1 (0%) 92 93	12, 18, 31, 43	4 (1%)
1	C	232/237 (97%)	-0.20	3 (1%) 77 80	15, 24, 38, 47	6 (2%)
1	D	233/237 (98%)	-0.18	1 (0%) 92 93	15, 24, 37, 47	6 (2%)
All	All	933/948 (98%)	-0.30	6 (0%) 89 91	12, 21, 36, 47	22 (2%)

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	204	SER	3.8
1	D	163	GLY	3.3
1	B	149	GLY	3.2
1	C	205	HIS	2.8
1	C	150	THR	2.5
1	C	161	SER	2.2

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.

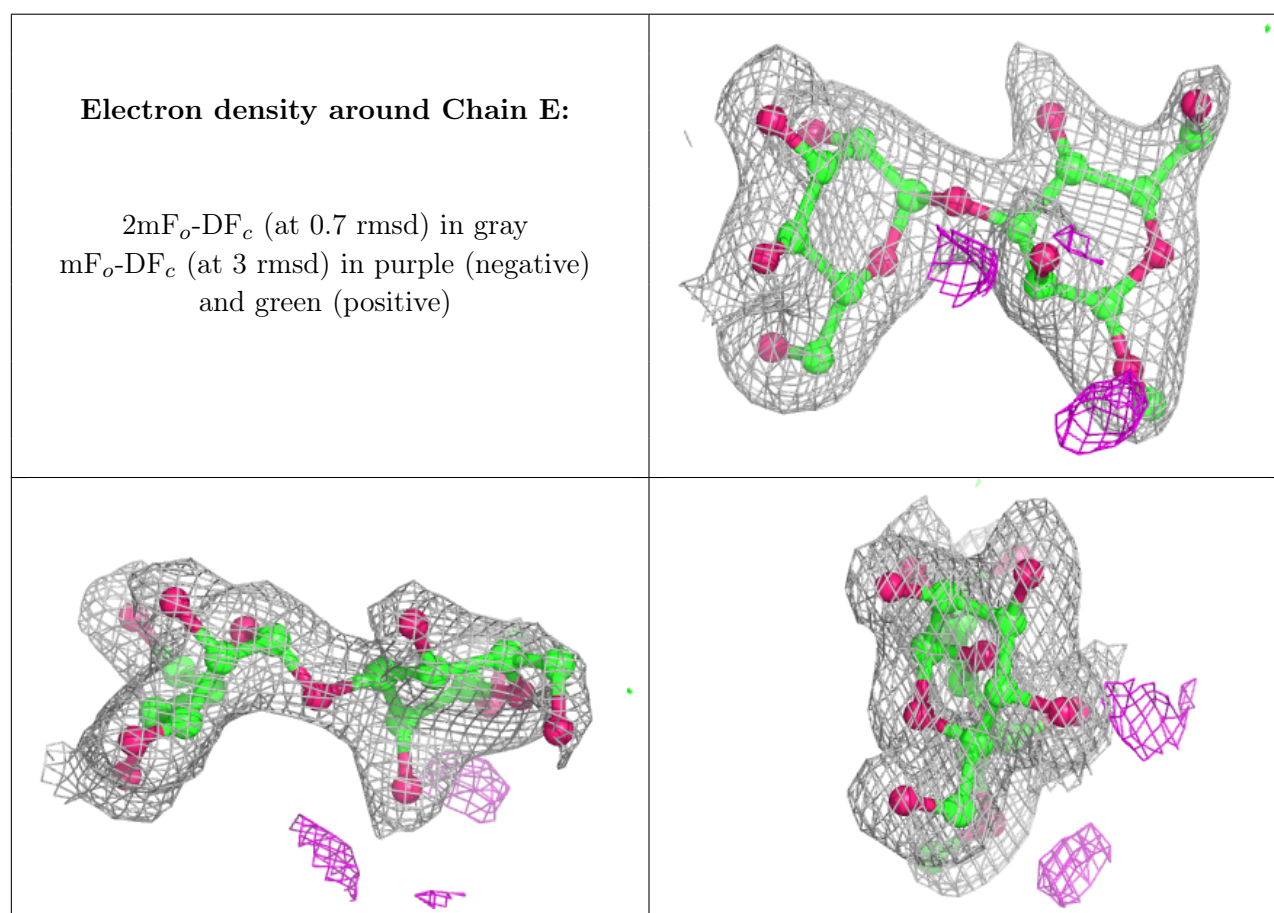
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	ZEL	E	1	13/13	0.91	0.17	25,33,37,40	0

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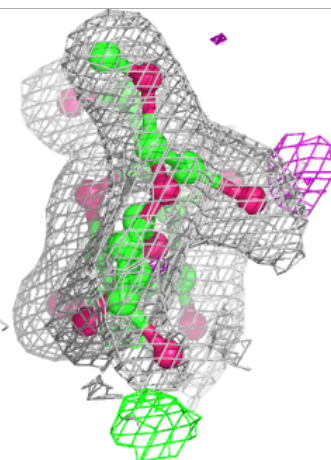
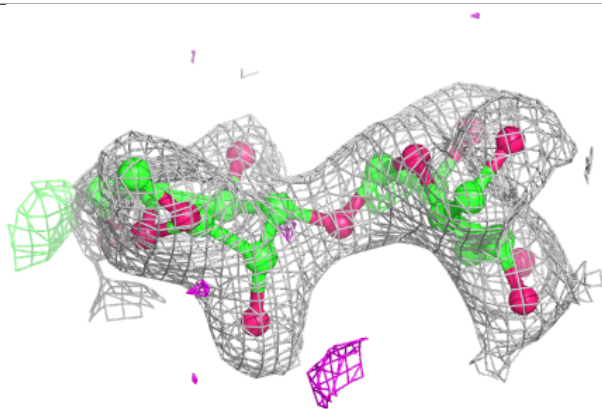
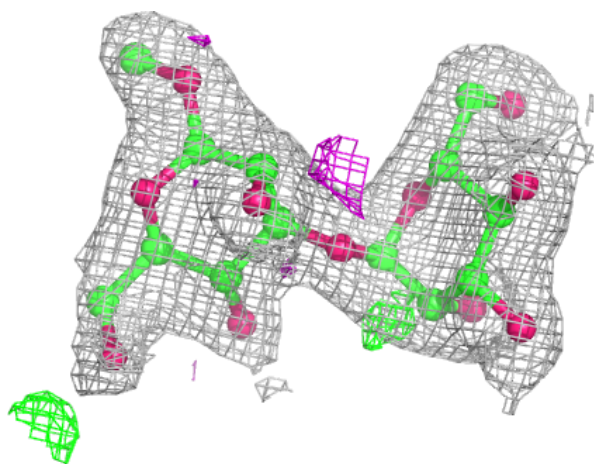
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
2	ZEL	F	1	13/13	0.92	0.13	25,32,39,42	0
2	ZEL	H	1	13/13	0.94	0.14	32,40,43,46	0
2	ZEL	G	1	13/13	0.96	0.10	27,33,36,37	0
2	MAN	G	2	11/12	0.96	0.10	21,27,28,29	0
2	MAN	F	2	11/12	0.96	0.11	14,20,22,23	0
2	MAN	H	2	11/12	0.96	0.12	22,27,29,31	0
2	MAN	E	2	11/12	0.97	0.09	13,18,22,22	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.



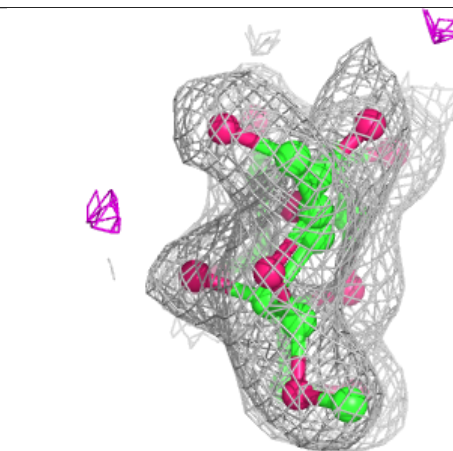
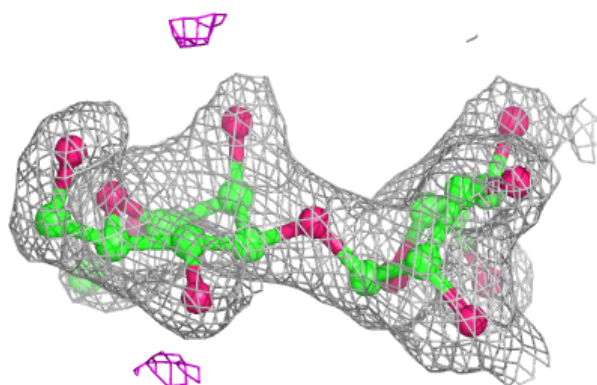
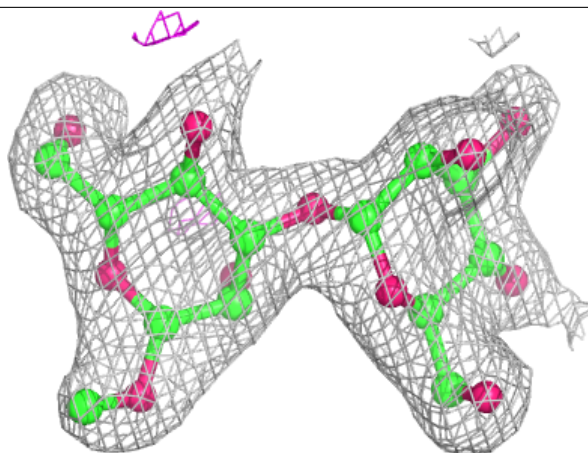
Electron density around Chain F:

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

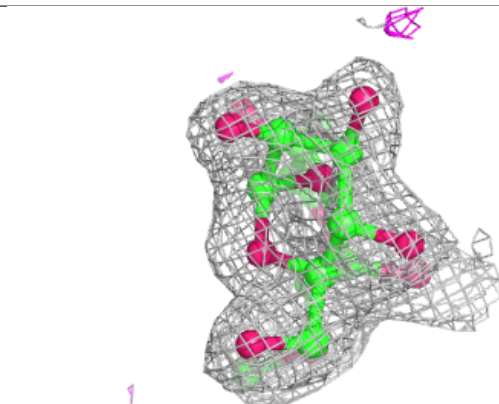
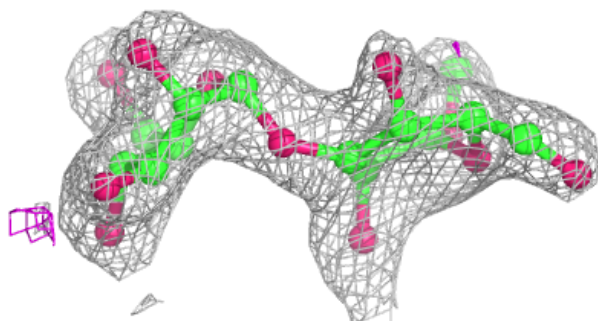
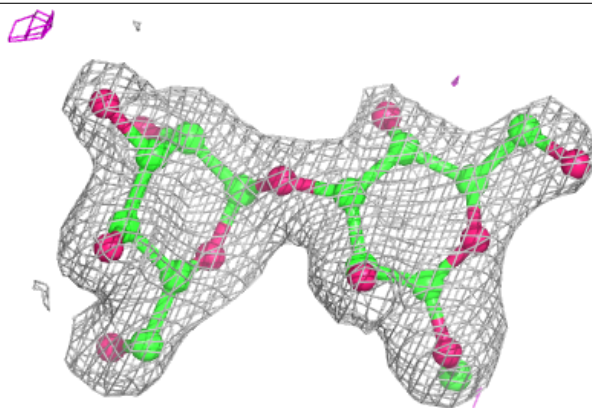


Electron density around Chain G:

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

**Electron density around Chain H:**

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



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	CA	D	239	1/1	0.97	0.07	20,20,20,20	0
4	CA	C	239	1/1	0.98	0.06	20,20,20,20	0
3	MN	B	238	1/1	0.99	0.10	18,18,18,18	0
3	MN	C	238	1/1	0.99	0.08	23,23,23,23	0
4	CA	A	239	1/1	1.00	0.05	18,18,18,18	0
4	CA	B	239	1/1	1.00	0.08	17,17,17,17	0
3	MN	A	238	1/1	1.00	0.12	18,18,18,18	0
3	MN	D	238	1/1	1.00	0.09	23,23,23,23	0

6.5 Other polymers [i](#)

There are no such residues in this entry.