Full wwPDB X-ray Structure Validation Report (i)

Jun 18, 2024 – 01:47 PM EDT

PDB ID	:	4IRS
Title	:	Structure of the mouse CD1d-PyrC-alpha-GalCer-iNKT TCR complex
Authors	:	Nemcovic, M.; Zajonc, D.M.
Deposited on	:	2013-01-15
Resolution	:	2.80 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 (i) 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 (i)) were used in the production of this report:

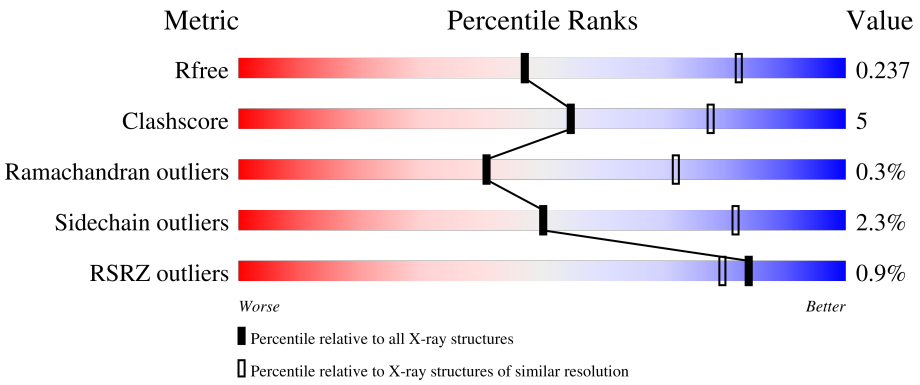
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 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	3140 (2.80-2.80)
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	285	2%, 76%, 17%, • 6%
2	B	99	%, 84%, 15%, •
3	C	209	86%, 10%, •
4	D	241	87%, 10%, • •
5	E	2	50%, 50%

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Mol	Chain	Length	Quality of chain
6	F	3	67% 33%

2 Entry composition (i)

There are 9 unique types of molecules in this entry. The entry contains 6604 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 Antigen-presenting glycoprotein CD1d1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	268	2139	1367	365	394	13	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	HIS	ASP	conflict	UNP P11609
A	280	HIS	-	expression tag	UNP P11609
A	281	HIS	-	expression tag	UNP P11609
A	282	HIS	-	expression tag	UNP P11609
A	283	HIS	-	expression tag	UNP P11609
A	284	HIS	-	expression tag	UNP P11609
A	285	HIS	-	expression tag	UNP P11609

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	98	788	502	133	146	7	0	0	0

- Molecule 3 is a protein called Valpha14 (mouse variable domain, human constant domain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	C	202	1537	952	260	317	8	0	0	0

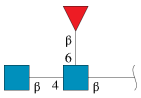
- Molecule 4 is a protein called Vbeta8.2 (mouse variable domain, human constant domain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
4	D	239	1879	1178	334	361	6	0	0	0

- Molecule 5 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.

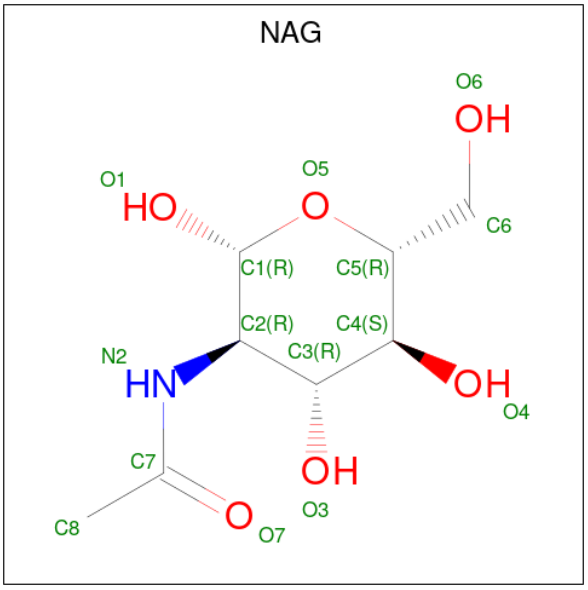
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
5	E	2	28	16	2	10	0	0	0

- Molecule 6 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[beta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose.



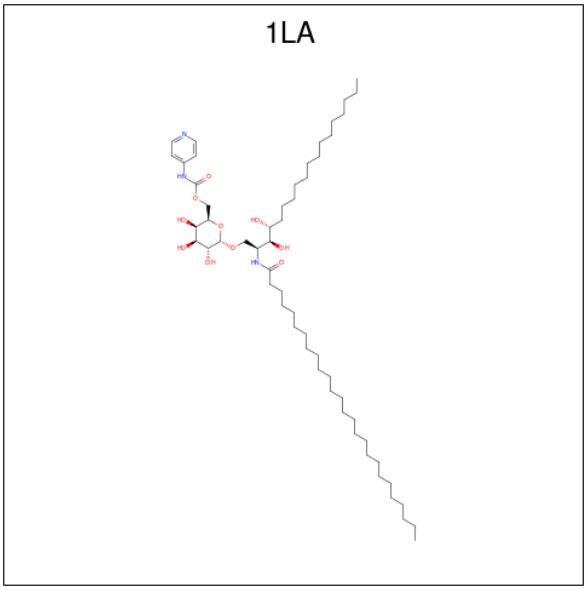
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
6	F	3	38	22	2	14	0	0	0

- Molecule 7 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



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

- Molecule 8 is N-[(2S,3S,4R)-3,4-dihydroxy-1-{[6-O-(pyridin-4-ylcarbamoyl)-alpha-D-galactopyranosyl]oxy}octadecan-2-yl]hexacosanamide (three-letter code: 1LA) (formula: $C_{56}H_{103}N_3O_{10}$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
8	A	1	69	56	3	10	0	0

- Molecule 9 is water.

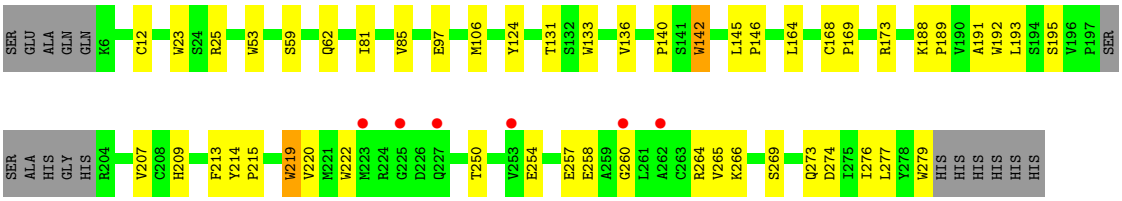
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
9	A	41	41	41	0	0
			Total	O		
9	B	14	14	14	0	0
			Total	O		
9	C	25	25	25	0	0
			Total	O		
9	D	32	32	32	0	0

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.

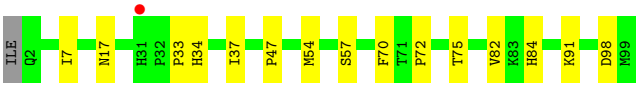
• Molecule 1: Antigen-presenting glycoprotein CD1d1

Chain A:



• Molecule 2: Beta-2-microglobulin

Chain B:

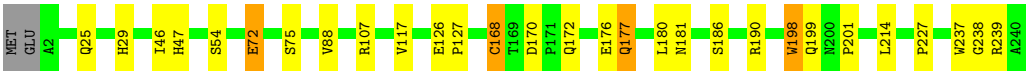


• Molecule 3: Valpha14 (mouse variable domain, human constant domain)

Chain C:

• Molecule 4: Vbeta8.2 (mouse variable domain, human constant domain)

Chain D:

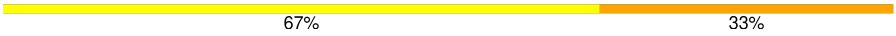


• Molecule 5: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain E:

• Molecule 6: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-[beta-L-fucopyranose-(1-6)]2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	78.97Å 191.40Å 151.22Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.33 – 2.80 59.33 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.6 (59.33-2.80) 97.6 (59.33-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.78 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.0	Depositor
R, R_{free}	0.190 , 0.239 0.190 , 0.237	Depositor DCC
R_{free} test set	1415 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å^2)	41.7	Xtriage
Anisotropy	0.087	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 36.3	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	6604	wwPDB-VP
Average B, all atoms (Å^2)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.40% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: 1LA, NAG, FUL

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	0.72	7/2201 (0.3%)	0.62	0/2993
2	B	0.62	0/814	0.62	0/1110
3	C	0.58	1/1564 (0.1%)	0.62	0/2130
4	D	0.63	2/1930 (0.1%)	0.61	0/2631
All	All	0.65	10/6509 (0.2%)	0.62	0/8864

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	TRP	CD2-CE2	5.46	1.48	1.41
1	A	53	TRP	CD2-CE2	5.46	1.48	1.41
3	C	178	TRP	CD2-CE2	5.38	1.47	1.41
1	A	133	TRP	CD2-CE2	5.33	1.47	1.41
4	D	198	TRP	CD2-CE2	5.28	1.47	1.41
1	A	23	TRP	CD2-CE2	5.25	1.47	1.41
1	A	192	TRP	CD2-CE2	5.19	1.47	1.41
4	D	237	TRP	CD2-CE2	5.19	1.47	1.41
1	A	219	TRP	CD2-CE2	5.11	1.47	1.41
1	A	222	TRP	CD2-CE2	5.11	1.47	1.41

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen

atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2139	0	2030	29	0
2	B	788	0	734	9	0
3	C	1537	0	1447	11	0
4	D	1879	0	1782	19	0
5	E	28	0	25	1	0
6	F	38	0	34	2	0
7	A	14	0	13	0	0
8	A	69	0	103	3	0
9	A	41	0	0	2	0
9	B	14	0	0	0	0
9	C	25	0	0	0	0
9	D	32	0	0	0	0
All	All	6604	0	6168	66	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 5.

All (66) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:GLU:HG2	9:A:435:HOH:O	1.76	0.86
2:B:17:ASN:HA	2:B:72:PRO:O	1.86	0.75
1:A:59:SER:H	1:A:62:GLN:NE2	1.86	0.73
3:C:155:VAL:HA	3:C:179:SER:HB2	1.74	0.69
2:B:33:PRO:HB2	2:B:54:MET:HE1	1.76	0.68
8:A:307:1LA:OCK	8:A:307:1LA:H100	1.92	0.67
1:A:12:CYS:HB3	8:A:307:1LA:H28	1.79	0.65
4:D:25:GLN:NE2	4:D:29:HIS:H	1.94	0.64
4:D:201:PRO:HA	4:D:238:GLY:O	1.98	0.64
1:A:193:LEU:HG	1:A:277:LEU:HD23	1.80	0.64
1:A:254:GLU:O	1:A:257:GLU:HG2	1.96	0.64
1:A:168:CYS:HB3	1:A:169:PRO:HD3	1.81	0.62
1:A:264:ARG:HH11	1:A:276:ILE:HD13	1.66	0.61
2:B:33:PRO:CB	2:B:54:MET:HE1	2.30	0.61
4:D:25:GLN:HE22	4:D:29:HIS:H	1.49	0.60
4:D:176:GLU:OE2	4:D:186:SER:OG	2.20	0.59
4:D:170:ASP:OD1	4:D:190:ARG:NH2	2.35	0.59
1:A:258:GLU:HB3	1:A:279:TRP:HD1	1.68	0.58
1:A:191:ALA:HA	1:A:209:HIS:O	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:TYR:CZ	1:A:136:VAL:HG11	2.40	0.57
1:A:219:TRP:HB3	1:A:266:LYS:HB2	1.86	0.56
2:B:33:PRO:CB	2:B:54:MET:CE	2.84	0.56
1:A:220:VAL:HG22	1:A:265:VAL:HG22	1.89	0.54
2:B:34:HIS:O	2:B:84:HIS:HD2	1.92	0.53
4:D:46:ILE:HG22	4:D:47:HIS:CD2	2.43	0.53
2:B:7:ILE:HD12	2:B:91:LYS:HD2	1.91	0.53
1:A:189:PRO:HB3	1:A:213:PHE:HB3	1.90	0.52
8:A:307:1LA:OCK	8:A:307:1LA:CCL	2.58	0.52
4:D:72:GLU:H	4:D:72:GLU:CD	2.14	0.51
4:D:127:PRO:HD2	4:D:198:TRP:CZ2	2.48	0.49
2:B:33:PRO:HB3	2:B:54:MET:CE	2.43	0.48
2:B:37:ILE:HG12	2:B:82:VAL:HG22	1.94	0.48
3:C:29:ASP:OD1	3:C:29:ASP:N	2.46	0.48
3:C:40:GLY:CA	4:D:107:ARG:HH22	2.27	0.48
4:D:177:GLN:HG2	4:D:180:LEU:HD12	1.95	0.47
1:A:207:VAL:HG22	1:A:250:THR:HG22	1.97	0.47
3:C:163:LEU:HB3	4:D:168:CYS:HB3	1.97	0.47
1:A:81:ILE:O	1:A:85:VAL:HG23	2.16	0.46
1:A:140:PRO:HB2	1:A:142:TRP:CD1	2.51	0.46
4:D:72:GLU:CD	4:D:72:GLU:N	2.70	0.45
3:C:55:LYS:HG2	3:C:64:THR:HG22	1.98	0.45
1:A:25:ARG:HB3	5:E:1:NAG:H82	1.99	0.44
1:A:195:SER:O	2:B:98:ASP:HB3	2.17	0.44
1:A:266:LYS:HG2	1:A:274:ASP:OD2	2.18	0.44
3:C:156:TYR:O	3:C:177:ALA:HA	2.17	0.44
1:A:59:SER:H	1:A:62:GLN:HE21	1.64	0.43
3:C:40:GLY:HA2	4:D:107:ARG:HH22	1.83	0.43
1:A:145:LEU:HB3	1:A:146:PRO:HD3	2.01	0.42
1:A:260:GLY:HA2	9:A:408:HOH:O	2.18	0.42
4:D:117:VAL:HG12	4:D:227:PRO:HB2	2.01	0.42
3:C:21:GLN:OE1	3:C:73:THR:HG22	2.18	0.42
3:C:36:LYS:HB2	3:C:46:LEU:HD11	2.00	0.42
1:A:214:TYR:CG	1:A:215:PRO:HA	2.55	0.42
3:C:202:PRO:O	3:C:203:SER:C	2.57	0.42
1:A:164:LEU:CD1	6:F:3:FUL:H62	2.50	0.41
1:A:106:MET:CE	1:A:173:ARG:HG2	2.51	0.41
1:A:145:LEU:HB3	1:A:146:PRO:CD	2.51	0.41
1:A:188:LYS:HA	1:A:269:SER:OG	2.20	0.41
4:D:25:GLN:HE22	4:D:29:HIS:N	2.18	0.41
4:D:214:LEU:HD13	4:D:227:PRO:HG2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:164:LEU:HD12	6:F:3:FUL:H62	2.02	0.40
1:A:59:SER:CB	1:A:62:GLN:HE21	2.34	0.40
3:C:138:PHE:HB2	3:C:190:PHE:CE1	2.56	0.40
4:D:199:GLN:HA	4:D:239:ARG:O	2.21	0.40
4:D:88:VAL:HG22	4:D:107:ARG:HG2	2.03	0.40
4:D:126:GLU:HA	4:D:127:PRO:HD3	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	264/285 (93%)	253 (96%)	11 (4%)	0	100	100
2	B	96/99 (97%)	94 (98%)	1 (1%)	1 (1%)	15	44
3	C	198/209 (95%)	186 (94%)	11 (6%)	1 (0%)	29	61
4	D	237/241 (98%)	231 (98%)	6 (2%)	0	100	100
All	All	795/834 (95%)	764 (96%)	29 (4%)	2 (0%)	41	72

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	179	SER
2	B	47	PRO

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	229/249 (92%)	227 (99%)	2 (1%)	78	94
2	B	87/93 (94%)	84 (97%)	3 (3%)	37	71
3	C	175/188 (93%)	171 (98%)	4 (2%)	50	82
4	D	204/208 (98%)	197 (97%)	7 (3%)	37	71
All	All	695/738 (94%)	679 (98%)	16 (2%)	50	82

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	131	THR
1	A	273	GLN
2	B	57	SER
2	B	70	PHE
2	B	75	THR
3	C	10	LEU
3	C	49	LEU
3	C	132	ASP
3	C	143	SER
4	D	54	SER
4	D	72	GLU
4	D	75	SER
4	D	168	CYS
4	D	172	GLN
4	D	177	GLN
4	D	181	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	62	GLN
1	A	110	ASN
1	A	273	GLN
2	B	17	ASN
3	C	30	ASN
4	D	25	GLN
4	D	136	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

5 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	$\#\|Z\| > 2$
5	NAG	E	1	1,5	14,14,15	0.59	0	17,19,21	0.99	1 (5%)
5	NAG	E	2	5	14,14,15	0.67	0	17,19,21	0.86	1 (5%)
6	NAG	F	1	1,6	14,14,15	0.53	0	17,19,21	1.39	2 (11%)
6	NAG	F	2	6	14,14,15	0.55	0	17,19,21	1.19	2 (11%)
6	FUL	F	3	6	10,10,11	0.57	0	14,14,16	1.53	4 (28%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	NAG	E	1	1,5	-	1/6/23/26	0/1/1/1
5	NAG	E	2	5	-	1/6/23/26	0/1/1/1
6	NAG	F	1	1,6	-	2/6/23/26	0/1/1/1
6	NAG	F	2	6	-	2/6/23/26	0/1/1/1
6	FUL	F	3	6	-	-	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	F	1	NAG	C1-O5-C5	4.41	118.10	112.19
6	F	3	FUL	C1-C2-C3	-3.17	105.02	109.64
5	E	2	NAG	C4-C3-C2	2.57	114.78	111.02
6	F	2	NAG	C1-O5-C5	2.45	115.47	112.19
6	F	3	FUL	O5-C1-C2	-2.31	105.29	110.79
6	F	1	NAG	C2-N2-C7	2.28	125.96	122.90
5	E	1	NAG	C1-O5-C5	2.28	115.24	112.19
6	F	3	FUL	O3-C3-C2	2.26	114.66	110.05
6	F	3	FUL	O5-C5-C4	2.22	113.56	109.55
6	F	2	NAG	O7-C7-N2	2.15	125.79	121.98

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	F	2	NAG	O5-C5-C6-O6
6	F	1	NAG	O5-C5-C6-O6
6	F	1	NAG	C4-C5-C6-O6
6	F	2	NAG	C4-C5-C6-O6
5	E	1	NAG	C4-C5-C6-O6
5	E	2	NAG	C1-C2-N2-C7

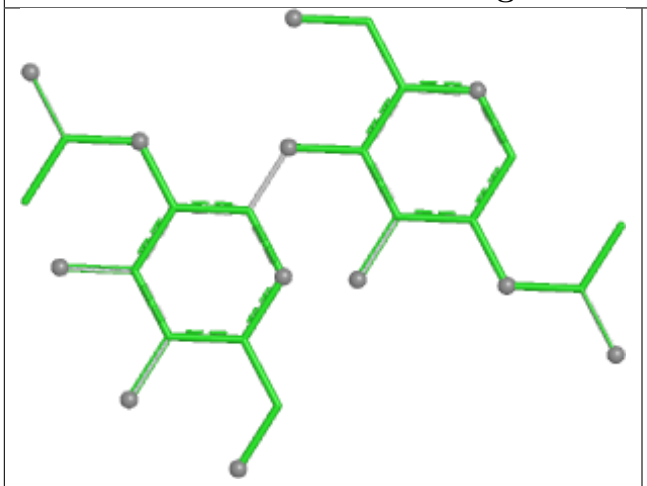
There are no ring outliers.

2 monomers are involved in 3 short contacts:

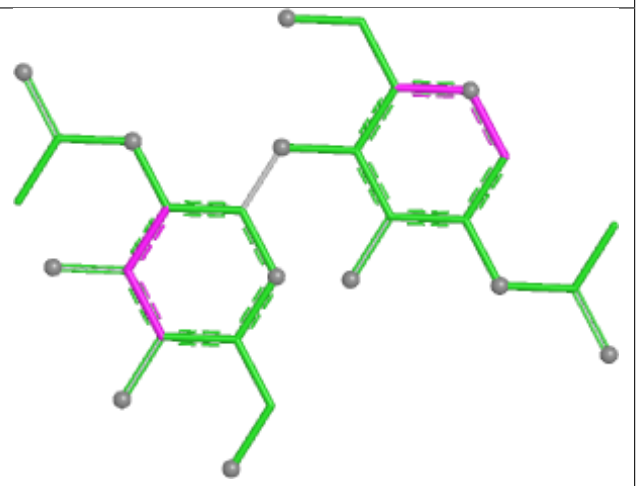
Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	F	3	FUL	2	0
5	E	1	NAG	1	0

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

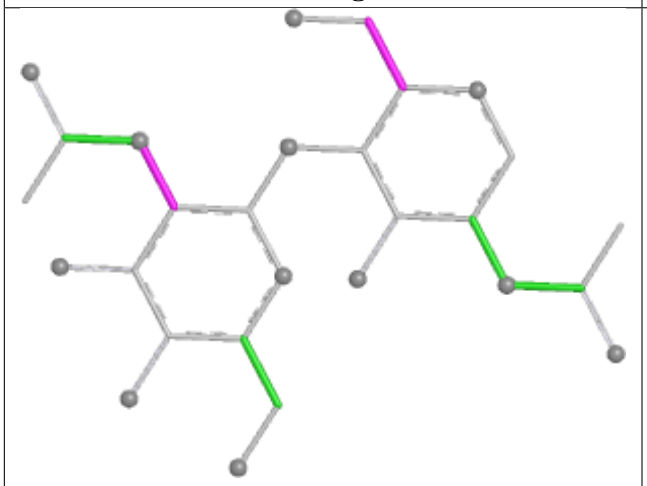
Oligosaccharide Chain E



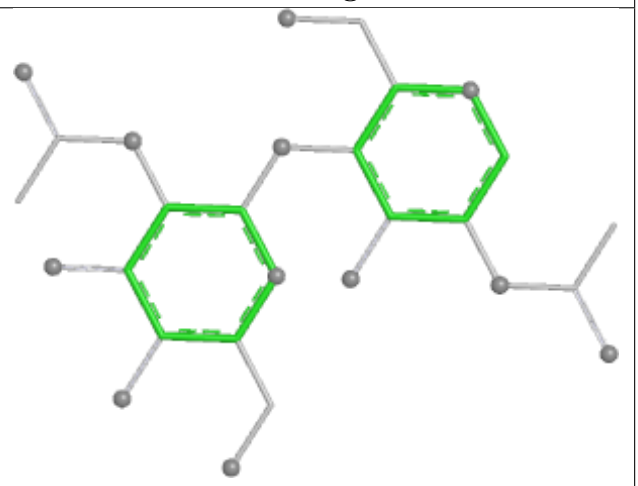
Bond lengths



Bond angles

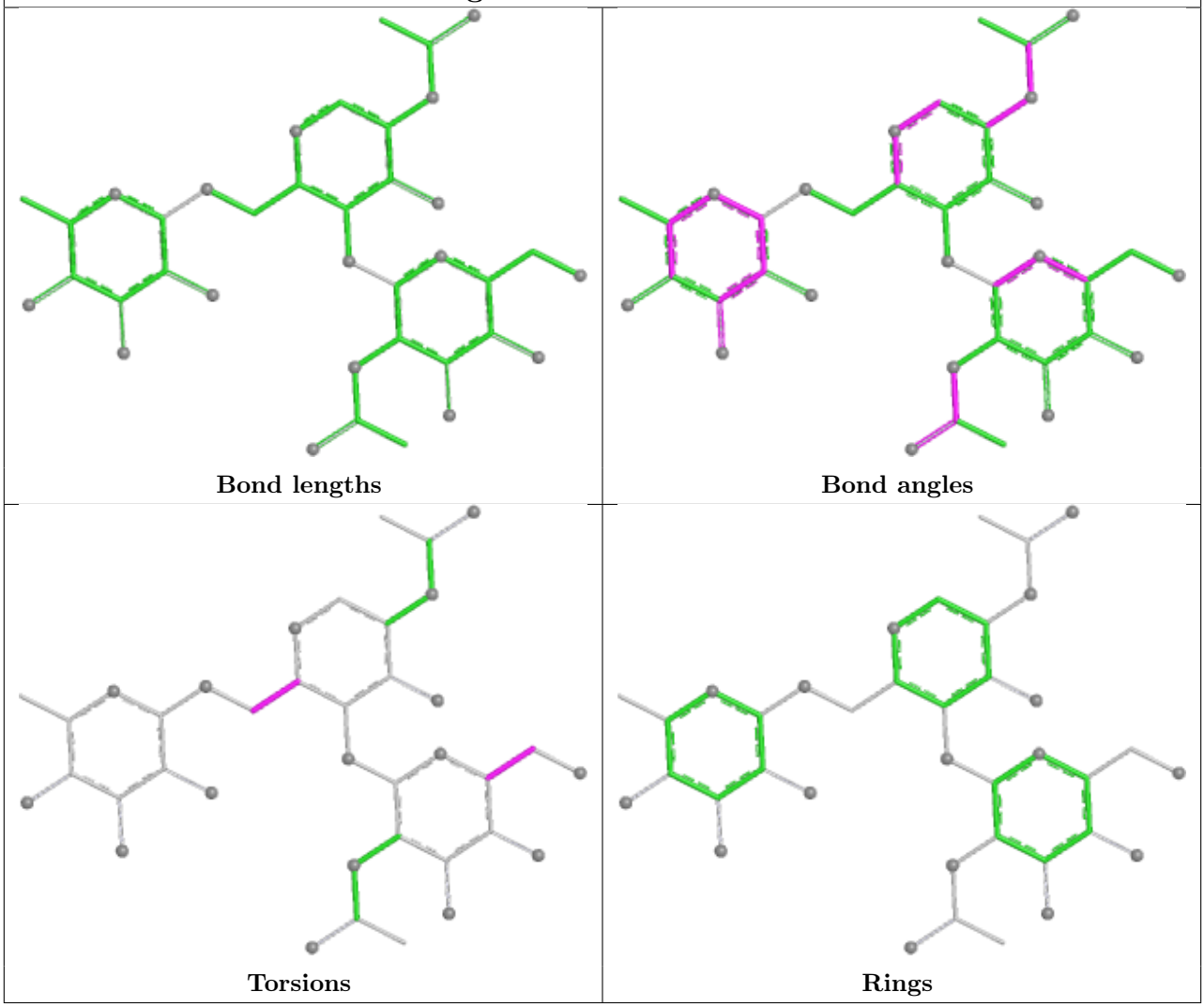


Torsions



Rings

Oligosaccharide Chain F



5.6 Ligand geometry (i)

2 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	$\#\|Z\| > 2$
8	1LA	A	307	-	70,70,70	0.91	4 (5%)	79,82,82	1.47	11 (13%)
7	NAG	A	301	1	14,14,15	0.66	0	17,19,21	1.21	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
8	1LA	A	307	-	-	21/65/85/85	0/2/2/2
7	NAG	A	301	1	-	0/6/23/26	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
8	A	307	1LA	CCQ-N	3.09	1.42	1.33
8	A	307	1LA	CCS-N	2.59	1.41	1.33
8	A	307	1LA	O6A-C1A	2.57	1.48	1.41
8	A	307	1LA	CCU-NCJ	-2.13	1.37	1.41

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
8	A	307	1LA	OAZ-CCI-NCJ	6.11	118.94	109.35
8	A	307	1LA	O6A-C5M-C6A	5.61	117.81	106.69
8	A	307	1LA	C1A-O6A-C5M	-2.74	108.37	113.72
7	A	301	NAG	C1-O5-C5	2.70	115.80	112.19
8	A	307	1LA	C6A-OAZ-CCI	-2.62	109.64	115.91
7	A	301	NAG	O5-C5-C6	2.52	112.57	107.66
8	A	307	1LA	OAZ-C6A-C5M	2.39	113.46	108.41
8	A	307	1LA	CCU-NCJ-CCI	-2.31	122.73	126.39
8	A	307	1LA	OCK-CCI-NCJ	-2.31	121.21	126.12
8	A	307	1LA	OAZ-CCI-OCK	-2.30	119.84	124.26
8	A	307	1LA	CCT-CCS-N	-2.18	119.88	123.60
8	A	307	1LA	C6-C5-C4	-2.09	110.69	114.11
8	A	307	1LA	CCS-CCT-CCU	2.03	121.55	118.88

There are no chirality outliers.

All (21) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
8	A	307	1LA	OCK-CCI-OAZ-C6A
8	A	307	1LA	NCJ-CCI-OAZ-C6A
8	A	307	1LA	CAJ-CAK-CAL-CAM
8	A	307	1LA	C14-C15-C16-C17
8	A	307	1LA	CAP-CAQ-CAR-CAS

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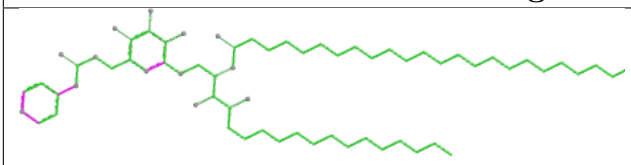
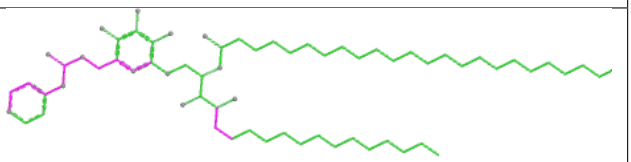
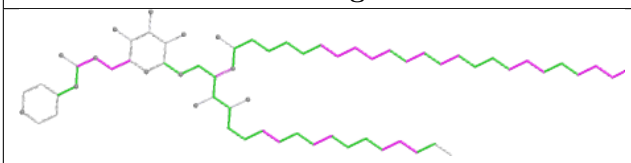
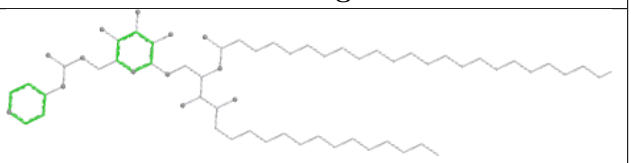
Mol	Chain	Res	Type	Atoms
8	A	307	1LA	CAF-CAG-CAH-CAI
8	A	307	1LA	C9-C10-C11-C12
8	A	307	1LA	CAT-CAU-CAV-CAW
8	A	307	1LA	C6-C7-C8-C9
8	A	307	1LA	CAV-CAW-CAX-CAY
8	A	307	1LA	CAG-CAH-CAI-CAJ
8	A	307	1LA	CAQ-CAR-CAS-CAT
8	A	307	1LA	CAW-CAX-CAY-CAZ
8	A	307	1LA	C13-C14-C15-C16
8	A	307	1LA	CAL-CAM-CAN-CAO
8	A	307	1LA	C5M-C6A-OAZ-CCI
8	A	307	1LA	CAU-CAV-CAW-CAX
8	A	307	1LA	CAH-CAI-CAJ-CAK
8	A	307	1LA	C4A-C5M-C6A-OAZ
8	A	307	1LA	C3-C2-N2-CAA
8	A	307	1LA	CAE-CAF-CAG-CAH

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
8	A	307	1LA	3	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.

Ligand 1LA A 307	
 Bond lengths	 Bond angles
 Torsions	 Rings

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	268/285 (94%)	0.03	6 (2%)	62	52	18, 31, 72, 84	0
2	B	98/99 (98%)	-0.13	1 (1%)	82	77	21, 37, 60, 76	1 (1%)
3	C	202/209 (96%)	0.00	0	100	100	20, 34, 73, 96	0
4	D	239/241 (99%)	-0.15	0	100	100	19, 32, 54, 83	0
All	All	807/834 (96%)	-0.05	7 (0%)	84	80	18, 33, 70, 96	1 (0%)

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	262	ALA	3.1
1	A	260	GLY	2.5
2	B	31	HIS	2.4
1	A	223	MET	2.2
1	A	227	GLN	2.1
1	A	225	GLY	2.0
1	A	253	VAL	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)

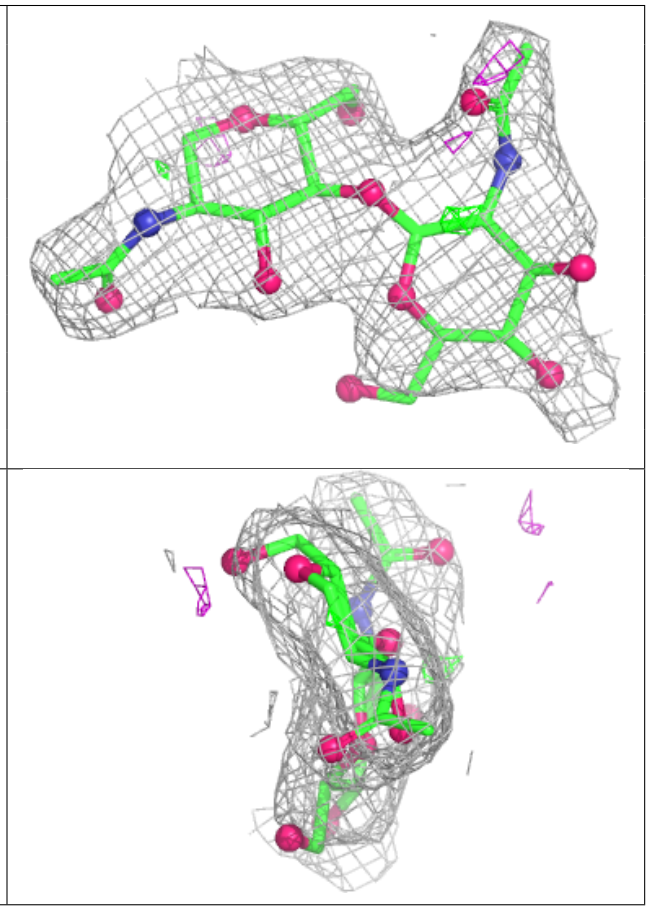
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	NAG	E	2	14/15	0.87	0.25	63,68,75,75	0
6	FUL	F	3	10/11	0.92	0.27	51,55,57,60	0
6	NAG	F	2	14/15	0.94	0.23	37,43,45,46	0
5	NAG	E	1	14/15	0.95	0.20	36,41,48,55	0
6	NAG	F	1	14/15	0.97	0.18	32,35,43,51	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.

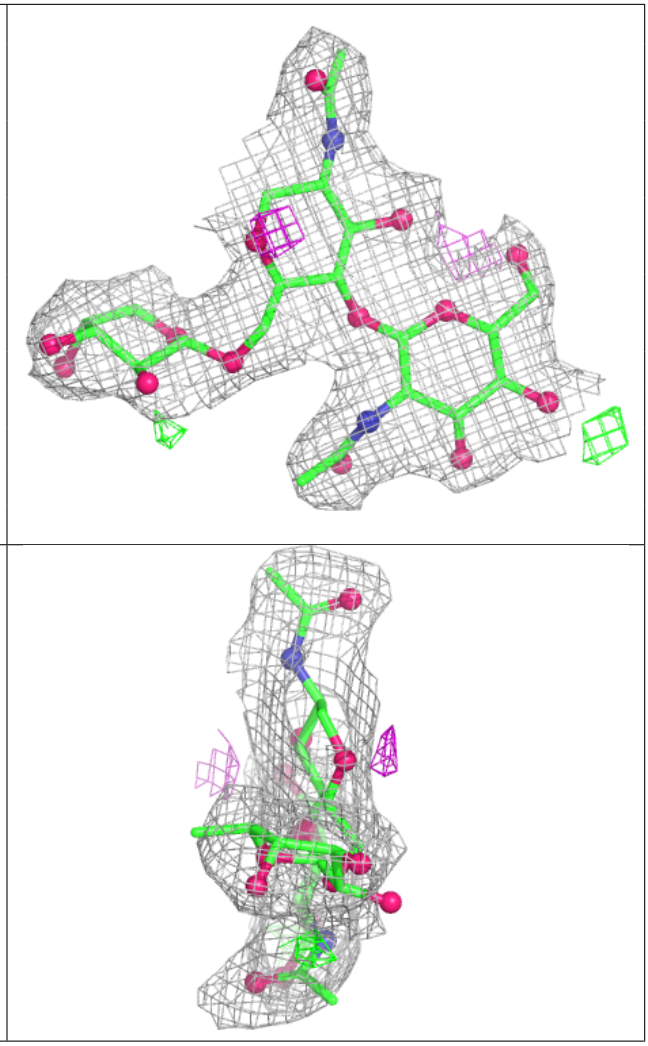
Electron density around Chain E:

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 F:

$2mF_o$-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)

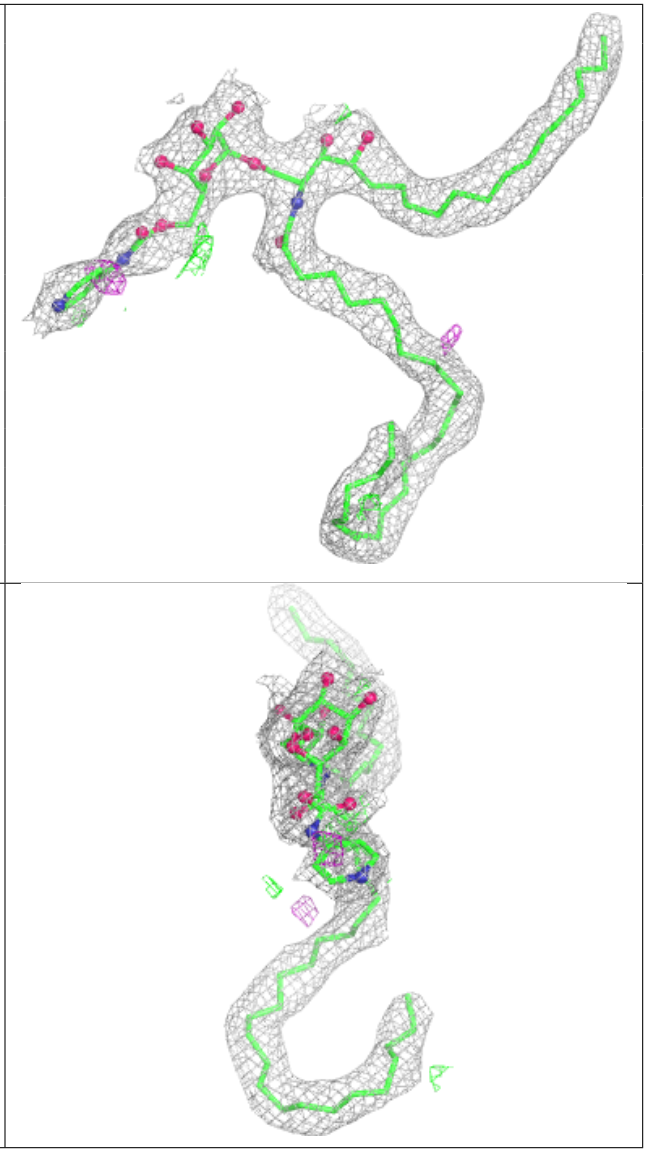
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
7	NAG	A	301	14/15	0.85	0.22	53,67,71,74	0
8	1LA	A	307	69/69	0.95	0.21	22,26,43,44	0

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

Electron density around 1LA A 307:

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



6.5 Other polymers (i)

There are no such residues in this entry.