Full wwPDB X-ray Structure Validation Report (i)

Jun 23, 2024 – 03:45 PM EDT

PDB ID	:	5J67
Title	:	Structure of Astrotactin-2, a conserved vertebrate-specific and perforin-like membrane protein involved in neuronal development
Authors	:	Ni, T.; Harlos, K.; Gilbert, R.J.C.
Deposited on	:	2016-04-04
Resolution	:	3.16 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:

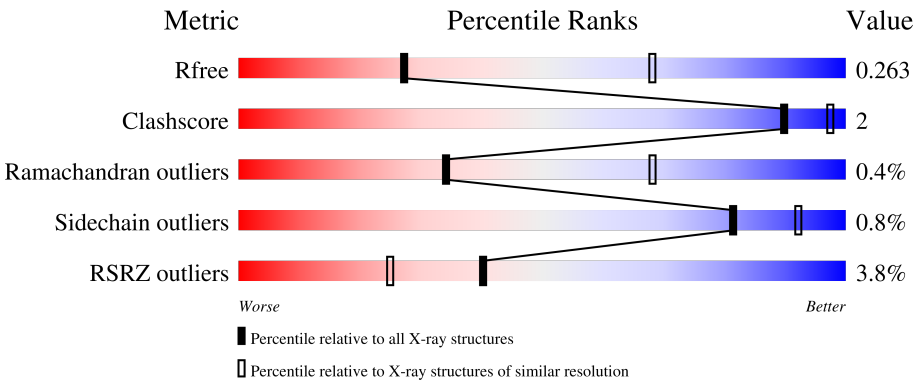
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & 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 3.16 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	1665 (3.20-3.12)
Clashscore	141614	1804 (3.20-3.12)
Ramachandran outliers	138981	1770 (3.20-3.12)
Sidechain outliers	138945	1769 (3.20-3.12)
RSRZ outliers	127900	1616 (3.20-3.12)

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	572	%, 92%, 5% •
1	B	572	6%, 91%, 7% •
1	C	572	2%, 92%, 6% •
1	D	572	6%, 92%, 5% •
2	E	10	20%, 80%

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Mol	Chain	Length	Quality of chain
2	F	10	30% 70%
2	G	10	40% 60%
2	H	10	30% 70%

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	MAN	E	8	X	-	-	-
2	MAN	F	6	-	-	-	X
2	MAN	F	8	X	-	-	-
2	MAN	G	8	X	-	-	-
2	MAN	H	8	X	-	-	-
3	NAG	D	1310	-	-	-	X
4	I3P	B	1311	-	-	-	X
4	I3P	D	1311	-	-	-	X

2 Entry composition (i)

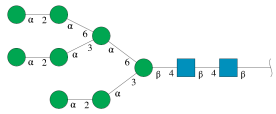
There are 4 unique types of molecules in this entry. The entry contains 18450 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Astrotactin-2.

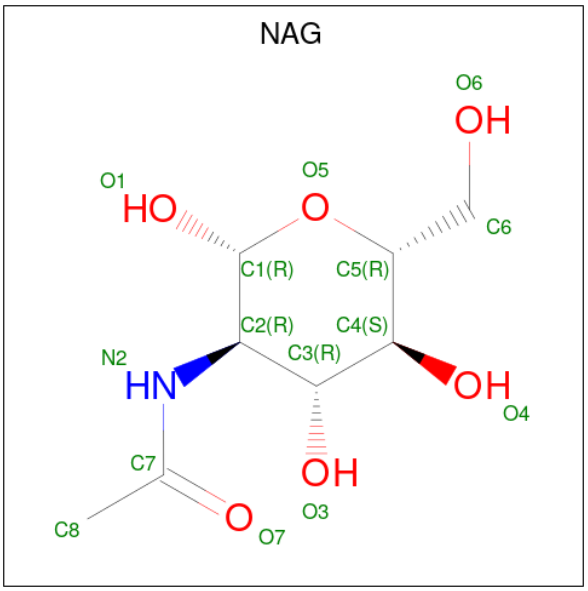
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	559	4456	2823	758	846	29	0	0	0
1	B	558	4445	2817	754	845	29	0	0	0
1	C	563	4492	2845	762	856	29	0	0	0
1	D	557	4441	2814	752	846	29	0	0	0

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



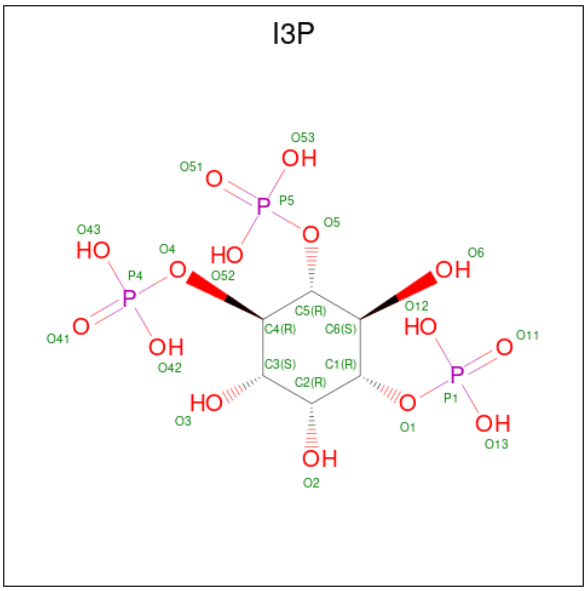
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
			Total	C	N	O			
2	E	10	116	64	2	50	0	0	0
2	F	10	116	64	2	50	0	0	0
2	G	10	116	64	2	50	0	0	0
2	H	10	116	64	2	50	0	0	0

- Molecule 3 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		
3	A	1	14	8	1	5	0	0
			Total	C	N	O		
3	B	1	14	8	1	5	0	0
			Total	C	N	O		
3	C	1	14	8	1	5	0	0
			Total	C	N	O		
3	D	1	14	8	1	5	0	0

- Molecule 4 is D-MYO-INOSITOL-1,4,5-TRIPHOSPHATE (three-letter code: I3P) (formula: $C_6H_{15}O_{15}P_3$).



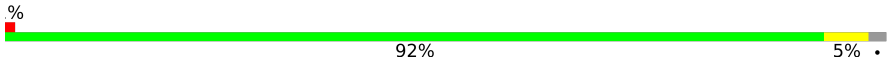
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
4	A	1	24	6	15	3	0	0
			Total	C	O	P		
4	B	1	24	6	15	3	0	0
			Total	C	O	P		
4	C	1	24	6	15	3	0	0
			Total	C	O	P		
4	D	1	24	6	15	3	0	0

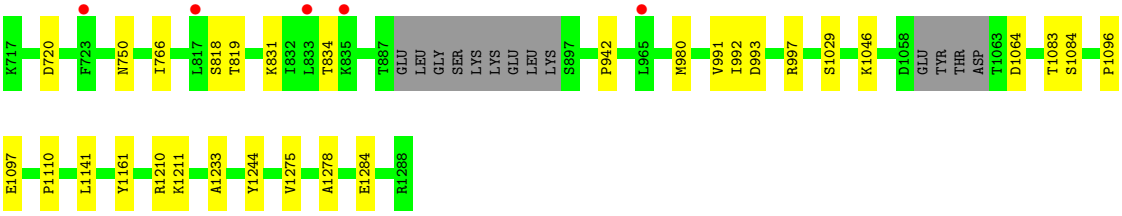
3 Residue-property plots (i)

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: Astrotactin-2

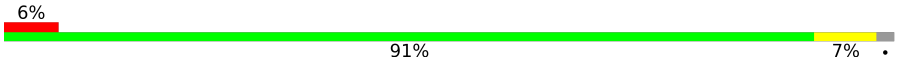
Chain A:

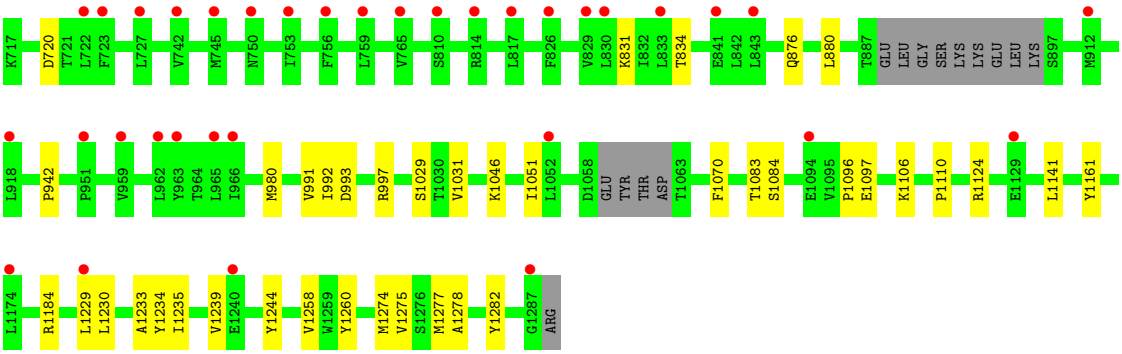




• Molecule 1: Astrotactin-2

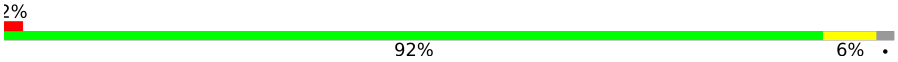
Chain B:

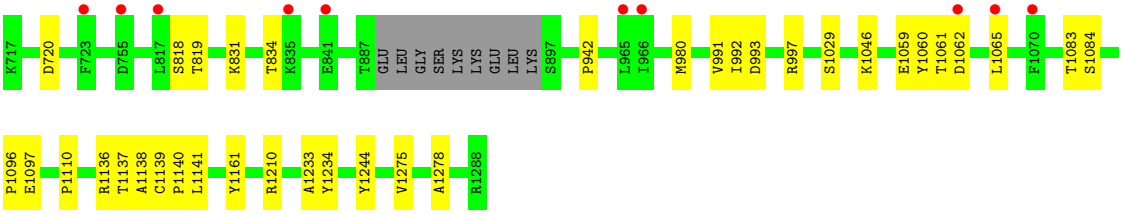




• Molecule 1: Astrotactin-2

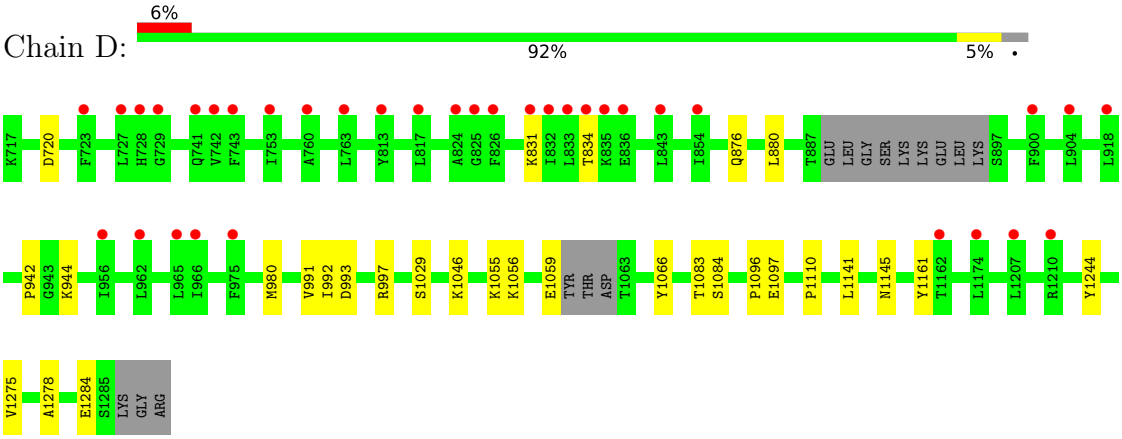
Chain C:





• Molecule 1: Astrotactin-2

Chain D:



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

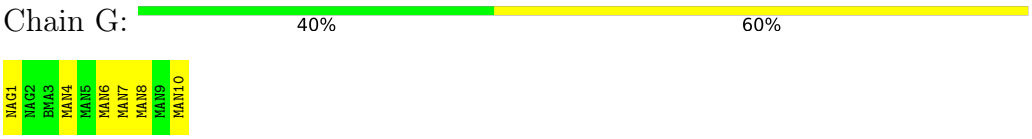
Chain E:

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

Chain F:

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

Chain G:



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

Chain H:

NAG1
NAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α, β, γ	101.24Å 108.19Å 111.36Å 87.18° 84.50° 65.67°	Depositor
Resolution (Å)	98.58 – 3.16 98.58 – 3.16	Depositor EDS
% Data completeness (in resolution range)	99.9 (98.58-3.16) 99.9 (98.58-3.16)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.32 (at 3.13Å)	Xtriage
Refinement program	REFMAC 5.8.0135	Depositor
R, R_{free}	0.233 , 0.259 0.235 , 0.263	Depositor DCC
R_{free} test set	3722 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å^2)	109.3	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.30 , 77.0	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	0.94	EDS
Total number of atoms	18450	wwPDB-VP
Average B, all atoms (Å^2)	136.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.94% 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: BMA, NAG, MAN, I3P

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	0.39	0/4552	0.65	0/6166
1	B	0.39	0/4540	0.65	0/6149
1	C	0.40	0/4588	0.69	5/6214 (0.1%)
1	D	0.38	0/4537	0.65	0/6148
All	All	0.39	0/18217	0.66	5/24677 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	1210	ARG	NE-CZ-NH2	9.94	125.27	120.30
1	C	1210	ARG	NE-CZ-NH1	-6.12	117.24	120.30
1	C	1136	ARG	NE-CZ-NH1	-5.58	117.51	120.30
1	C	1062	ASP	CB-CG-OD2	5.21	122.99	118.30
1	C	1136	ARG	NE-CZ-NH2	5.02	122.81	120.30

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	4456	0	4386	14	1
1	B	4445	0	4372	29	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	4492	0	4411	19	0
1	D	4441	0	4363	11	1
2	E	116	0	97	0	0
2	F	116	0	97	0	0
2	G	116	0	97	0	0
2	H	116	0	97	0	0
3	A	14	0	13	0	0
3	B	14	0	13	0	0
3	C	14	0	13	0	0
3	D	14	0	13	0	0
4	A	24	0	9	0	0
4	B	24	0	9	0	0
4	C	24	0	9	0	0
4	D	24	0	9	0	0
All	All	18450	0	18008	66	1

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 2.

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:C:1138:ALA:C	1:C:1139:CYS:N	2.21	0.94
1:B:880:LEU:HD22	1:C:1065:LEU:HD23	1.56	0.88
1:B:880:LEU:HD22	1:C:1065:LEU:CD2	2.07	0.85
1:B:1051:ILE:HG23	1:B:1070:PHE:HE1	1.46	0.80
1:B:1051:ILE:HG23	1:B:1070:PHE:CE1	2.19	0.78
1:C:1060:TYR:O	1:C:1061:THR:OG1	2.08	0.69
1:C:1233:ALA:O	1:C:1234:TYR:N	2.28	0.66
1:A:766:ILE:HG23	1:B:1274:MET:HE3	1.85	0.59
1:B:1106:LYS:NZ	1:B:1244:TYR:OH	2.34	0.59
1:B:1051:ILE:CG2	1:B:1070:PHE:HE1	2.13	0.59
1:D:980:MET:HB3	1:D:991:VAL:HG11	1.85	0.58
1:C:980:MET:HB3	1:C:991:VAL:HG11	1.84	0.58
1:B:980:MET:HB3	1:B:991:VAL:HG11	1.86	0.58
1:B:1141:LEU:HD21	1:B:1278:ALA:HB1	1.87	0.57
1:B:1260:TYR:CD1	1:D:944:LYS:HE2	2.39	0.57
1:A:980:MET:HB3	1:A:991:VAL:HG11	1.86	0.56
1:B:1051:ILE:CG2	1:B:1070:PHE:CE1	2.88	0.55
1:C:1029:SER:N	1:C:1138:ALA:HB3	2.22	0.55
1:C:1083:THR:HB	1:C:1084:SER:HA	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1083:THR:HB	1:A:1084:SER:HA	1.89	0.54
1:D:1083:THR:HB	1:D:1084:SER:HA	1.89	0.54
1:C:1141:LEU:HD21	1:C:1278:ALA:HB1	1.90	0.54
1:B:1083:THR:HB	1:B:1084:SER:HA	1.89	0.53
1:B:1233:ALA:O	1:B:1234:TYR:N	2.43	0.52
1:D:1029:SER:HB2	1:D:1110:PRO:HB3	1.93	0.51
1:B:1051:ILE:HD12	1:B:1070:PHE:CE1	2.46	0.51
1:B:1124:ARG:NH1	1:B:1258:VAL:HG12	2.26	0.51
1:A:1210:ARG:HG3	1:A:1211:LYS:N	2.26	0.50
1:B:1029:SER:HB2	1:B:1110:PRO:HB3	1.92	0.50
1:A:1029:SER:HB2	1:A:1110:PRO:HB3	1.93	0.50
1:C:1029:SER:HB2	1:C:1110:PRO:HB3	1.93	0.49
1:A:766:ILE:HG23	1:B:1274:MET:CE	2.42	0.48
1:B:880:LEU:HB3	1:C:1065:LEU:HD22	1.96	0.48
1:A:750:ASN:O	1:B:1239:VAL:HB	2.14	0.48
1:C:1244:TYR:CD1	1:C:1275:VAL:HG23	2.51	0.46
1:B:1244:TYR:CD1	1:B:1275:VAL:HG23	2.52	0.45
1:B:1051:ILE:HD12	1:B:1070:PHE:HE1	1.81	0.44
1:C:1137:THR:O	1:C:1138:ALA:C	2.55	0.44
1:A:831:LYS:O	1:A:834:THR:HG22	2.18	0.44
1:B:1031:VAL:HG11	1:B:1277:MET:CE	2.48	0.44
1:A:1096:PRO:HA	1:A:1097:GLU:HA	1.82	0.43
1:B:831:LYS:O	1:B:834:THR:HG22	2.18	0.43
1:C:831:LYS:O	1:C:834:THR:HG22	2.18	0.43
1:C:818:SER:HA	1:C:819:THR:HA	1.86	0.43
1:A:818:SER:HA	1:A:819:THR:HA	1.86	0.43
1:D:831:LYS:O	1:D:834:THR:HG22	2.18	0.43
1:C:1110:PRO:HB2	1:C:1140:PRO:HG3	2.01	0.42
1:B:992:ILE:HD13	1:B:997:ARG:HB2	2.01	0.42
1:D:1244:TYR:CD1	1:D:1275:VAL:HG23	2.53	0.42
1:A:1244:TYR:CD1	1:A:1275:VAL:HG23	2.55	0.42
1:C:992:ILE:HD13	1:C:997:ARG:HB2	2.02	0.42
1:C:1096:PRO:HA	1:C:1097:GLU:HA	1.81	0.42
1:D:992:ILE:HD13	1:D:997:ARG:HB2	2.01	0.42
1:B:1230:LEU:HA	1:B:1235:ILE:HD11	2.02	0.41
1:D:1096:PRO:HA	1:D:1097:GLU:HA	1.82	0.41
1:A:992:ILE:HD13	1:A:997:ARG:HB2	2.02	0.41
1:B:1096:PRO:HA	1:B:1097:GLU:HA	1.82	0.41
1:B:1184:ARG:CZ	1:B:1235:ILE:HG13	2.51	0.41
1:B:876:GLN:O	1:B:880:LEU:HG	2.21	0.41
1:C:1141:LEU:CD2	1:C:1278:ALA:HB1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:1141:LEU:HD21	1:D:1278:ALA:HB1	2.02	0.41
1:A:1141:LEU:HD21	1:A:1278:ALA:HB1	2.03	0.40
1:D:876:GLN:O	1:D:880:LEU:HG	2.21	0.40
1:B:1229:LEU:HB3	1:B:1282:TYR:CE1	2.56	0.40
1:D:1055:LYS:CD	1:D:1066:TYR:HD1	2.35	0.40
1:A:1233:ALA:HB3	1:A:1284:GLU:HA	2.02	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1210:ARG:NH1	1:D:1145:ASN:OD1[1_554]	2.03	0.17

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	553/572 (97%)	521 (94%)	29 (5%)	3 (0%)	29	65
1	B	550/572 (96%)	520 (94%)	28 (5%)	2 (0%)	34	68
1	C	555/572 (97%)	524 (94%)	29 (5%)	2 (0%)	34	68
1	D	551/572 (96%)	521 (95%)	28 (5%)	2 (0%)	34	68
All	All	2209/2288 (96%)	2086 (94%)	114 (5%)	9 (0%)	34	68

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1064	ASP
1	A	1161	TYR
1	B	1161	TYR
1	C	942	PRO

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Mol	Chain	Res	Type
1	C	1161	TYR
1	D	1161	TYR
1	A	942	PRO
1	B	942	PRO
1	D	942	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	501/513 (98%)	498 (99%)	3 (1%)	86	94
1	B	500/513 (98%)	497 (99%)	3 (1%)	86	94
1	C	505/513 (98%)	501 (99%)	4 (1%)	81	92
1	D	500/513 (98%)	494 (99%)	6 (1%)	71	87
All	All	2006/2052 (98%)	1990 (99%)	16 (1%)	81	92

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

Mol	Chain	Res	Type
1	A	720	ASP
1	A	993	ASP
1	A	1046	LYS
1	B	720	ASP
1	B	993	ASP
1	B	1046	LYS
1	C	720	ASP
1	C	993	ASP
1	C	1046	LYS
1	C	1059	GLU
1	D	720	ASP
1	D	993	ASP
1	D	1046	LYS
1	D	1056	LYS
1	D	1059	GLU
1	D	1284	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

40 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	NAG	E	1	1,2	14,14,15	0.59	0	17,19,21	2.59	8 (47%)
2	MAN	E	10	2	11,11,12	0.38	0	15,15,17	0.94	1 (6%)
2	NAG	E	2	2	14,14,15	0.31	0	17,19,21	0.85	0
2	BMA	E	3	2	11,11,12	0.20	0	15,15,17	0.85	0
2	MAN	E	4	2	11,11,12	0.43	0	15,15,17	1.20	2 (13%)
2	MAN	E	5	2	11,11,12	0.46	0	15,15,17	1.09	1 (6%)
2	MAN	E	6	2	11,11,12	0.31	0	15,15,17	0.99	1 (6%)
2	MAN	E	7	2	11,11,12	0.63	0	15,15,17	1.24	2 (13%)
2	MAN	E	8	2	11,11,12	0.60	0	15,15,17	2.12	2 (13%)
2	MAN	E	9	2	11,11,12	0.30	0	15,15,17	0.66	1 (6%)
2	NAG	F	1	1,2	14,14,15	0.46	0	17,19,21	2.45	5 (29%)
2	MAN	F	10	2	11,11,12	0.43	0	15,15,17	0.88	0
2	NAG	F	2	2	14,14,15	0.38	0	17,19,21	0.74	0
2	BMA	F	3	2	11,11,12	0.30	0	15,15,17	0.71	0
2	MAN	F	4	2	11,11,12	0.46	0	15,15,17	0.99	2 (13%)
2	MAN	F	5	2	11,11,12	0.43	0	15,15,17	1.39	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	MAN	F	6	2	11,11,12	0.29	0	15,15,17	0.84	1 (6%)
2	MAN	F	7	2	11,11,12	0.46	0	15,15,17	1.40	3 (20%)
2	MAN	F	8	2	11,11,12	0.46	0	15,15,17	1.84	2 (13%)
2	MAN	F	9	2	11,11,12	0.33	0	15,15,17	1.03	1 (6%)
2	NAG	G	1	1,2	14,14,15	0.44	0	17,19,21	2.13	3 (17%)
2	MAN	G	10	2	11,11,12	0.35	0	15,15,17	0.87	1 (6%)
2	NAG	G	2	2	14,14,15	0.43	0	17,19,21	0.79	0
2	BMA	G	3	2	11,11,12	0.27	0	15,15,17	0.74	0
2	MAN	G	4	2	11,11,12	0.29	0	15,15,17	0.76	1 (6%)
2	MAN	G	5	2	11,11,12	0.30	0	15,15,17	0.76	0
2	MAN	G	6	2	11,11,12	0.37	0	15,15,17	1.00	1 (6%)
2	MAN	G	7	2	11,11,12	0.46	0	15,15,17	1.10	1 (6%)
2	MAN	G	8	2	11,11,12	0.44	0	15,15,17	1.62	2 (13%)
2	MAN	G	9	2	11,11,12	0.25	0	15,15,17	0.61	0
2	NAG	H	1	1,2	14,14,15	0.43	0	17,19,21	2.49	5 (29%)
2	MAN	H	10	2	11,11,12	0.44	0	15,15,17	0.98	1 (6%)
2	NAG	H	2	2	14,14,15	0.35	0	17,19,21	0.87	0
2	BMA	H	3	2	11,11,12	0.29	0	15,15,17	1.07	0
2	MAN	H	4	2	11,11,12	0.40	0	15,15,17	0.60	0
2	MAN	H	5	2	11,11,12	0.41	0	15,15,17	1.04	1 (6%)
2	MAN	H	6	2	11,11,12	0.38	0	15,15,17	0.90	1 (6%)
2	MAN	H	7	2	11,11,12	0.35	0	15,15,17	0.79	1 (6%)
2	MAN	H	8	2	11,11,12	0.38	0	15,15,17	1.57	3 (20%)
2	MAN	H	9	2	11,11,12	0.35	0	15,15,17	0.96	1 (6%)

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	NAG	E	1	1,2	-	3/6/23/26	0/1/1/1
2	MAN	E	10	2	-	2/2/19/22	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	BMA	E	3	2	-	0/2/19/22	0/1/1/1
2	MAN	E	4	2	-	2/2/19/22	0/1/1/1
2	MAN	E	5	2	-	0/2/19/22	0/1/1/1
2	MAN	E	6	2	-	0/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MAN	E	7	2	-	2/2/19/22	0/1/1/1
2	MAN	E	8	2	1/1/4/5	1/2/19/22	0/1/1/1
2	MAN	E	9	2	-	0/2/19/22	0/1/1/1
2	NAG	F	1	1,2	-	3/6/23/26	0/1/1/1
2	MAN	F	10	2	-	0/2/19/22	0/1/1/1
2	NAG	F	2	2	-	0/6/23/26	0/1/1/1
2	BMA	F	3	2	-	0/2/19/22	0/1/1/1
2	MAN	F	4	2	-	1/2/19/22	0/1/1/1
2	MAN	F	5	2	-	0/2/19/22	0/1/1/1
2	MAN	F	6	2	-	2/2/19/22	0/1/1/1
2	MAN	F	7	2	-	1/2/19/22	0/1/1/1
2	MAN	F	8	2	1/1/4/5	2/2/19/22	0/1/1/1
2	MAN	F	9	2	-	1/2/19/22	0/1/1/1
2	NAG	G	1	1,2	-	4/6/23/26	0/1/1/1
2	MAN	G	10	2	-	0/2/19/22	0/1/1/1
2	NAG	G	2	2	-	0/6/23/26	0/1/1/1
2	BMA	G	3	2	-	0/2/19/22	0/1/1/1
2	MAN	G	4	2	-	2/2/19/22	0/1/1/1
2	MAN	G	5	2	-	0/2/19/22	0/1/1/1
2	MAN	G	6	2	-	2/2/19/22	0/1/1/1
2	MAN	G	7	2	-	1/2/19/22	0/1/1/1
2	MAN	G	8	2	1/1/4/5	2/2/19/22	0/1/1/1
2	MAN	G	9	2	-	0/2/19/22	0/1/1/1
2	NAG	H	1	1,2	-	3/6/23/26	0/1/1/1
2	MAN	H	10	2	-	0/2/19/22	0/1/1/1
2	NAG	H	2	2	-	0/6/23/26	0/1/1/1
2	BMA	H	3	2	-	0/2/19/22	0/1/1/1
2	MAN	H	4	2	-	2/2/19/22	0/1/1/1
2	MAN	H	5	2	-	0/2/19/22	0/1/1/1
2	MAN	H	6	2	-	2/2/19/22	0/1/1/1
2	MAN	H	7	2	-	1/2/19/22	0/1/1/1
2	MAN	H	8	2	1/1/4/5	2/2/19/22	0/1/1/1
2	MAN	H	9	2	-	2/2/19/22	0/1/1/1

There are no bond length outliers.

All (55) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	1	NAG	C1-O5-C5	8.30	123.43	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	1	NAG	C1-O5-C5	8.09	123.16	112.19
2	E	1	NAG	C1-O5-C5	7.77	122.72	112.19
2	E	8	MAN	C1-O5-C5	6.72	121.30	112.19
2	G	1	NAG	C1-O5-C5	6.29	120.71	112.19
2	F	8	MAN	C1-O5-C5	6.10	120.45	112.19
2	G	8	MAN	C1-O5-C5	5.02	118.99	112.19
2	F	5	MAN	C1-O5-C5	4.41	118.17	112.19
2	H	8	MAN	C1-O5-C5	4.40	118.15	112.19
2	G	1	NAG	C8-C7-N2	3.90	122.71	116.10
2	E	8	MAN	C1-C2-C3	3.52	113.99	109.67
2	E	1	NAG	C8-C7-N2	3.37	121.80	116.10
2	F	7	MAN	C3-C4-C5	3.10	115.78	110.24
2	E	6	MAN	C1-O5-C5	3.02	116.29	112.19
2	H	5	MAN	C1-O5-C5	2.99	116.24	112.19
2	E	1	NAG	O5-C5-C4	2.93	117.96	110.83
2	F	1	NAG	C8-C7-N2	2.89	120.99	116.10
2	H	1	NAG	C8-C7-N2	2.87	120.95	116.10
2	H	1	NAG	O5-C5-C4	2.79	117.60	110.83
2	F	1	NAG	O5-C5-C4	2.78	117.59	110.83
2	E	1	NAG	C1-C2-N2	-2.77	105.75	110.49
2	H	10	MAN	C1-O5-C5	2.75	115.92	112.19
2	F	7	MAN	C1-O5-C5	2.72	115.88	112.19
2	E	1	NAG	O5-C1-C2	2.68	115.52	111.29
2	F	4	MAN	C1-O5-C5	2.65	115.79	112.19
2	H	8	MAN	C3-C4-C5	2.65	114.97	110.24
2	E	4	MAN	C3-C4-C5	2.64	114.95	110.24
2	G	8	MAN	C3-C4-C5	2.63	114.93	110.24
2	F	9	MAN	C1-O5-C5	2.60	115.71	112.19
2	G	6	MAN	C1-O5-C5	2.59	115.70	112.19
2	H	7	MAN	C1-O5-C5	2.50	115.58	112.19
2	F	6	MAN	C1-O5-C5	2.49	115.57	112.19
2	H	6	MAN	C1-O5-C5	2.49	115.56	112.19
2	E	5	MAN	C1-O5-C5	2.47	115.55	112.19
2	H	8	MAN	O5-C5-C6	2.46	111.06	107.20
2	E	7	MAN	C2-C3-C4	2.45	115.13	110.89
2	F	7	MAN	O5-C1-C2	-2.39	107.08	110.77
2	G	7	MAN	C3-C4-C5	2.39	114.50	110.24
2	H	1	NAG	C4-C3-C2	-2.32	107.61	111.02
2	H	9	MAN	C1-C2-C3	2.28	112.46	109.67
2	E	1	NAG	C2-N2-C7	2.26	126.12	122.90
2	E	10	MAN	C1-O5-C5	2.25	115.24	112.19
2	F	1	NAG	C2-N2-C7	2.23	126.08	122.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	E	1	NAG	C4-C3-C2	-2.22	107.76	111.02
2	E	1	NAG	O7-C7-N2	-2.21	117.88	121.95
2	F	1	NAG	C4-C3-C2	-2.17	107.84	111.02
2	E	7	MAN	C3-C4-C5	2.14	114.05	110.24
2	F	4	MAN	C1-C2-C3	2.11	112.26	109.67
2	G	4	MAN	C1-O5-C5	2.11	115.05	112.19
2	E	4	MAN	C1-O5-C5	2.09	115.02	112.19
2	E	9	MAN	C1-O5-C5	2.06	114.98	112.19
2	F	8	MAN	O5-C5-C6	2.06	110.43	107.20
2	G	10	MAN	C1-O5-C5	2.04	114.96	112.19
2	H	1	NAG	C2-N2-C7	2.03	125.79	122.90
2	G	1	NAG	O5-C5-C4	2.01	115.71	110.83

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	E	8	MAN	C1
2	F	8	MAN	C1
2	G	8	MAN	C1
2	H	8	MAN	C1

All (43) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	10	MAN	C4-C5-C6-O6
2	H	4	MAN	C4-C5-C6-O6
2	E	10	MAN	O5-C5-C6-O6
2	G	1	NAG	O5-C5-C6-O6
2	E	7	MAN	O5-C5-C6-O6
2	F	8	MAN	O5-C5-C6-O6
2	G	4	MAN	C4-C5-C6-O6
2	E	4	MAN	C4-C5-C6-O6
2	G	8	MAN	O5-C5-C6-O6
2	H	8	MAN	O5-C5-C6-O6
2	H	4	MAN	O5-C5-C6-O6
2	G	1	NAG	C4-C5-C6-O6
2	E	7	MAN	C4-C5-C6-O6
2	E	4	MAN	O5-C5-C6-O6
2	G	4	MAN	O5-C5-C6-O6
2	F	6	MAN	C4-C5-C6-O6
2	F	8	MAN	C4-C5-C6-O6
2	H	9	MAN	C4-C5-C6-O6

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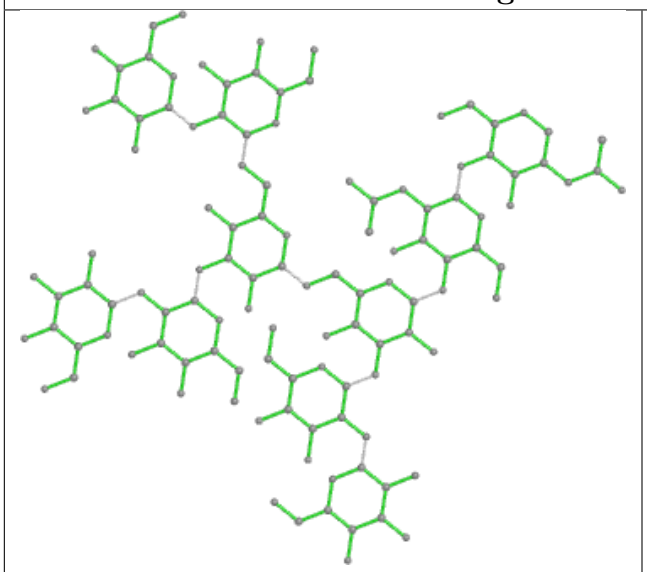
Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	F	1	NAG	C8-C7-N2-C2
2	F	1	NAG	O7-C7-N2-C2
2	G	1	NAG	C8-C7-N2-C2
2	G	1	NAG	O7-C7-N2-C2
2	H	1	NAG	C8-C7-N2-C2
2	H	1	NAG	O7-C7-N2-C2
2	F	6	MAN	O5-C5-C6-O6
2	H	6	MAN	C4-C5-C6-O6
2	H	6	MAN	O5-C5-C6-O6
2	H	8	MAN	C4-C5-C6-O6
2	H	9	MAN	O5-C5-C6-O6
2	G	8	MAN	C4-C5-C6-O6
2	F	4	MAN	O5-C5-C6-O6
2	F	9	MAN	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	F	7	MAN	O5-C5-C6-O6
2	G	6	MAN	C4-C5-C6-O6
2	H	7	MAN	O5-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
2	G	7	MAN	O5-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
2	E	8	MAN	C4-C5-C6-O6
2	G	6	MAN	O5-C5-C6-O6

There are no ring outliers.

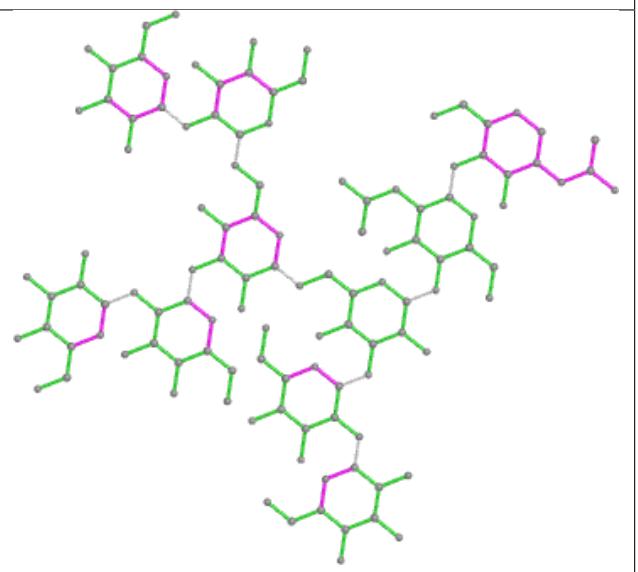
No monomer is involved in short contacts.

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

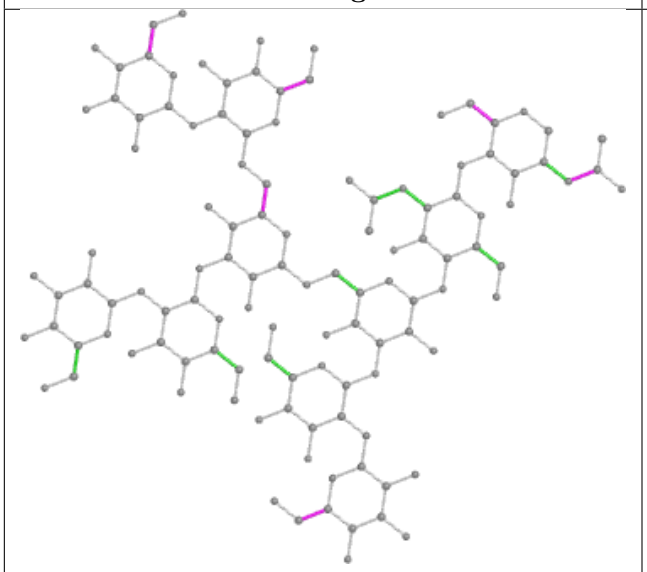
Oligosaccharide Chain E



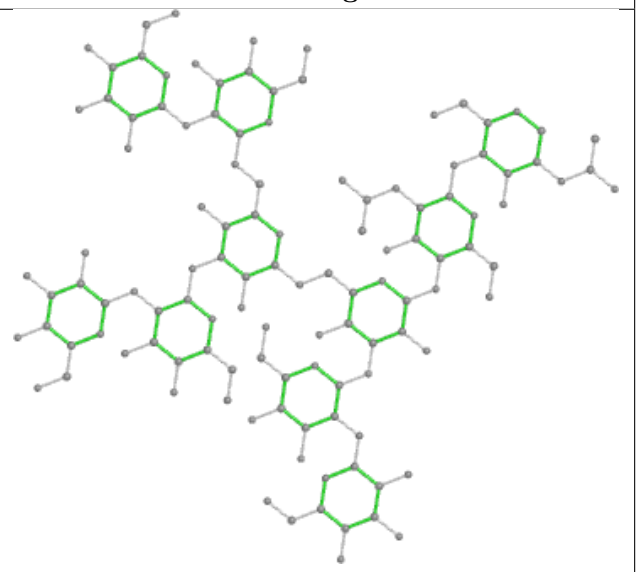
Bond lengths



Bond angles

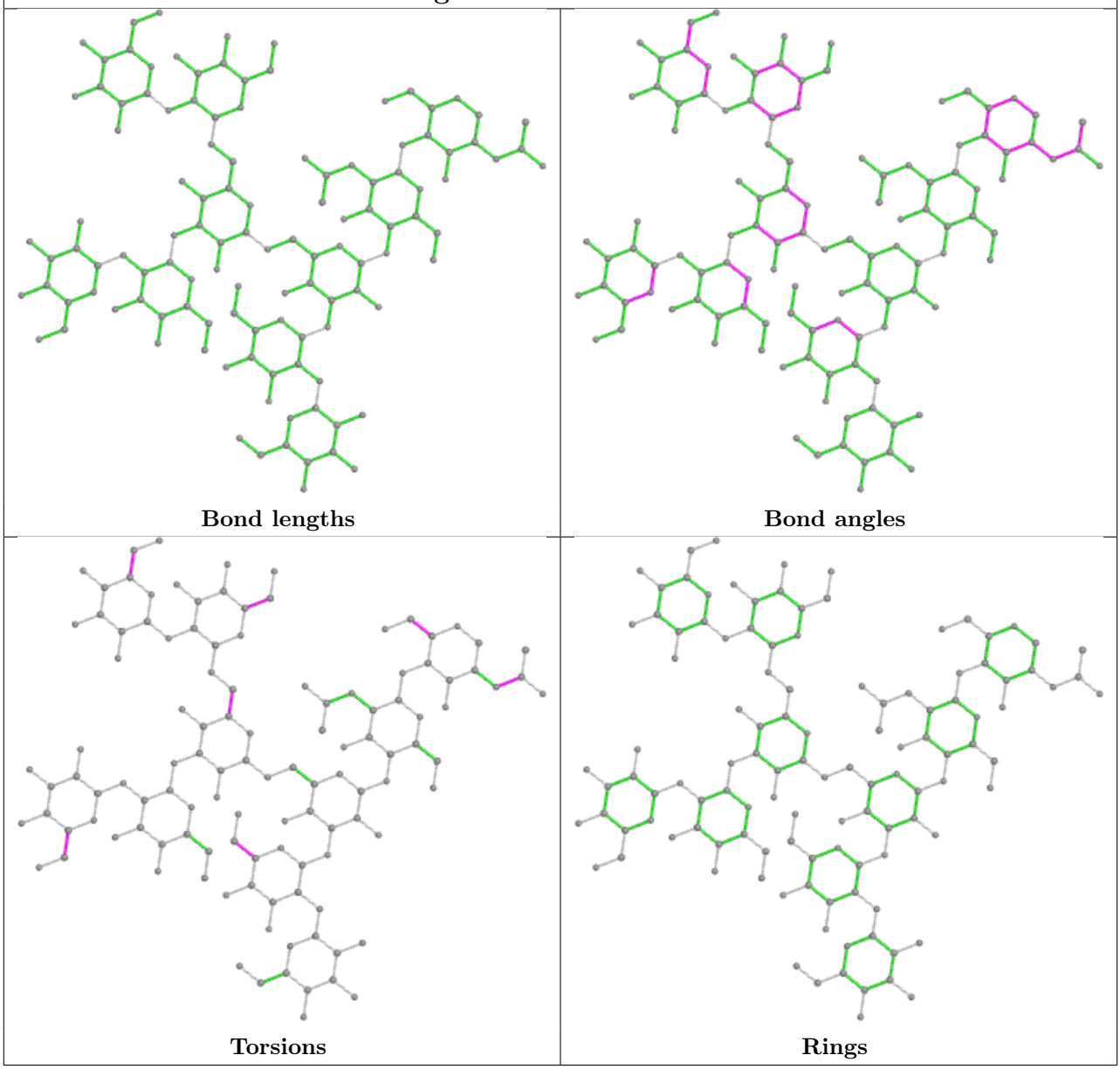


Torsions

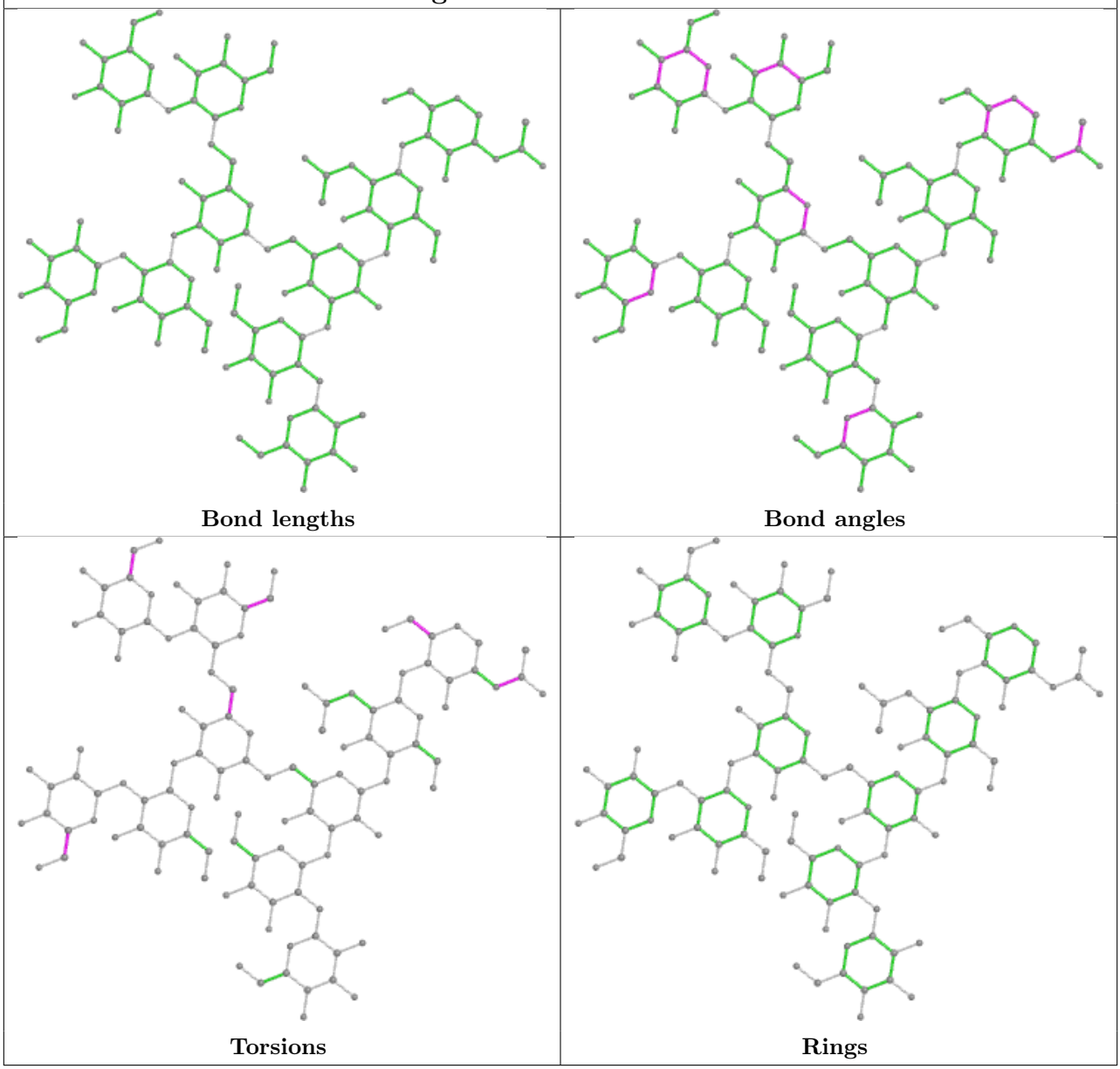


Rings

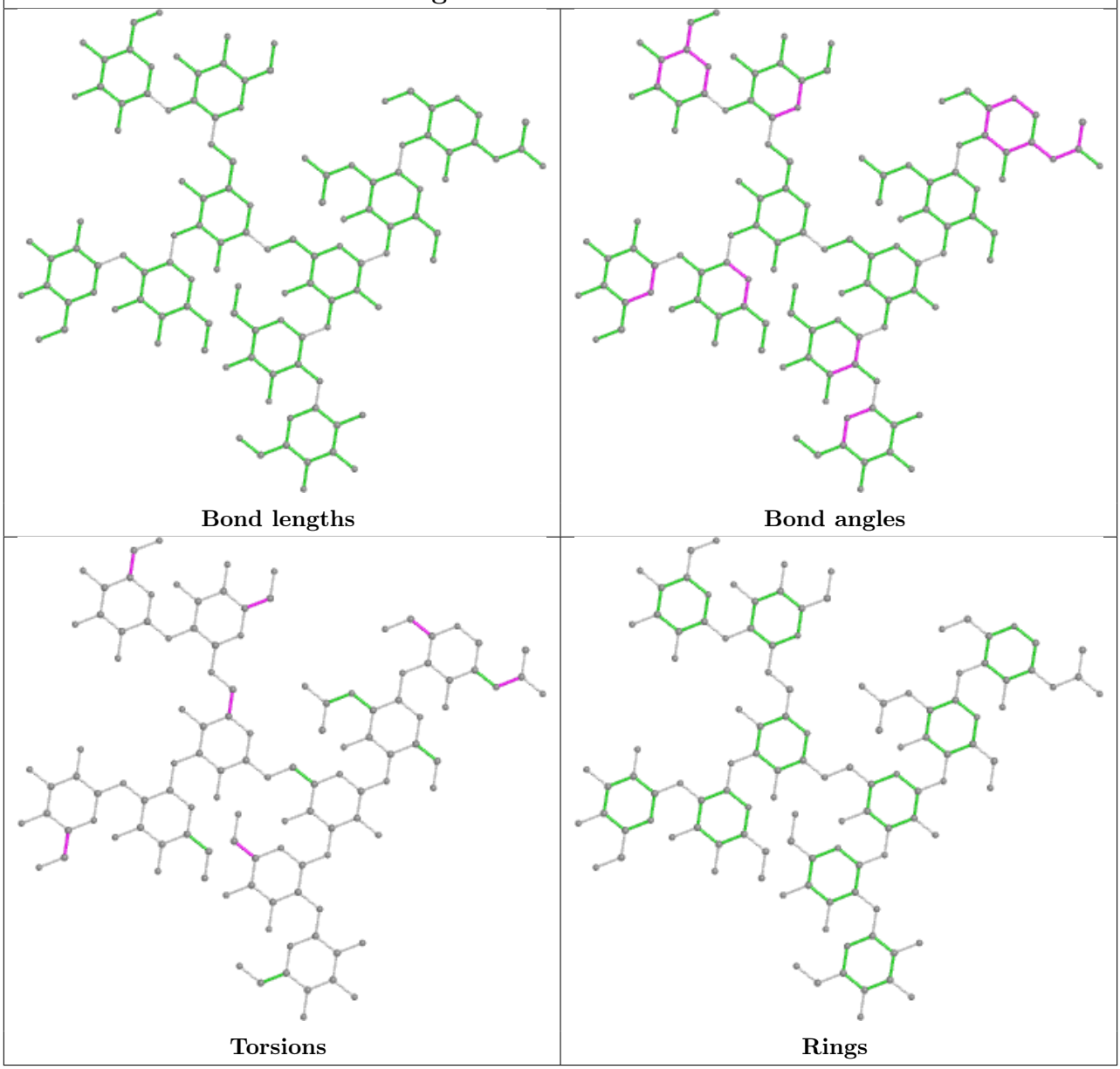
Oligosaccharide Chain F



Oligosaccharide Chain G



Oligosaccharide Chain H



5.6 Ligand geometry (i)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	I3P	D	1311	-	24,24,24	0.80	0	36,39,39	1.19	4 (11%)
4	I3P	B	1311	-	24,24,24	0.72	0	36,39,39	1.04	0
3	NAG	A	1310	1	14,14,15	0.29	0	17,19,21	1.08	2 (11%)
3	NAG	D	1310	1	14,14,15	0.36	0	17,19,21	0.81	0
3	NAG	B	1310	1	14,14,15	0.39	0	17,19,21	0.87	0
3	NAG	C	1310	1	14,14,15	0.36	0	17,19,21	1.58	3 (17%)
4	I3P	A	1311	-	24,24,24	0.79	1 (4%)	36,39,39	0.97	0
4	I3P	C	1311	-	24,24,24	0.87	1 (4%)	36,39,39	1.31	3 (8%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	I3P	D	1311	-	-	7/15/39/39	0/1/1/1
4	I3P	B	1311	-	-	0/15/39/39	0/1/1/1
3	NAG	A	1310	1	-	2/6/23/26	0/1/1/1
3	NAG	D	1310	1	-	2/6/23/26	0/1/1/1
3	NAG	B	1310	1	-	2/6/23/26	0/1/1/1
3	NAG	C	1310	1	-	2/6/23/26	0/1/1/1
4	I3P	A	1311	-	-	3/15/39/39	0/1/1/1
4	I3P	C	1311	-	-	4/15/39/39	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1311	I3P	P5-O5	2.13	1.63	1.59
4	A	1311	I3P	P5-O5	2.04	1.63	1.59

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	C	1310	NAG	C1-O5-C5	4.89	118.82	112.19
4	C	1311	I3P	C6-C1-C2	3.52	115.92	110.85
3	C	1310	NAG	O5-C1-C2	-2.82	106.84	111.29
4	C	1311	I3P	C5-C6-C1	2.64	114.44	108.96
4	D	1311	I3P	O4-C4-C5	2.48	114.54	108.69
4	D	1311	I3P	C6-C1-C2	2.43	114.35	110.85

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	1310	NAG	C1-O5-C5	2.37	115.40	112.19
3	A	1310	NAG	O5-C1-C2	-2.20	107.81	111.29
4	D	1311	I3P	C6-C5-C4	-2.14	106.77	111.66
3	C	1310	NAG	C3-C4-C5	2.12	114.02	110.24
4	C	1311	I3P	O4-C4-C5	2.10	113.64	108.69
4	D	1311	I3P	O5-C5-C4	2.02	113.44	108.69

There are no chirality outliers.

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	1311	I3P	C1-O1-P1-O11
4	C	1311	I3P	C2-C1-O1-P1
4	C	1311	I3P	C6-C1-O1-P1
4	C	1311	I3P	C1-O1-P1-O11
4	D	1311	I3P	C2-C1-O1-P1
4	D	1311	I3P	C6-C1-O1-P1
3	B	1310	NAG	O5-C5-C6-O6
3	B	1310	NAG	C4-C5-C6-O6
3	C	1310	NAG	O5-C5-C6-O6
3	C	1310	NAG	C4-C5-C6-O6
3	A	1310	NAG	C4-C5-C6-O6
3	D	1310	NAG	C4-C5-C6-O6
3	A	1310	NAG	O5-C5-C6-O6
4	C	1311	I3P	C5-O5-P5-O53
4	D	1311	I3P	C4-O4-P4-O43
3	D	1310	NAG	O5-C5-C6-O6
4	A	1311	I3P	C4-O4-P4-O41
4	D	1311	I3P	C1-O1-P1-O11
4	D	1311	I3P	C5-O5-P5-O51
4	A	1311	I3P	C5-O5-P5-O52
4	D	1311	I3P	C1-O1-P1-O12
4	D	1311	I3P	C5-O5-P5-O52

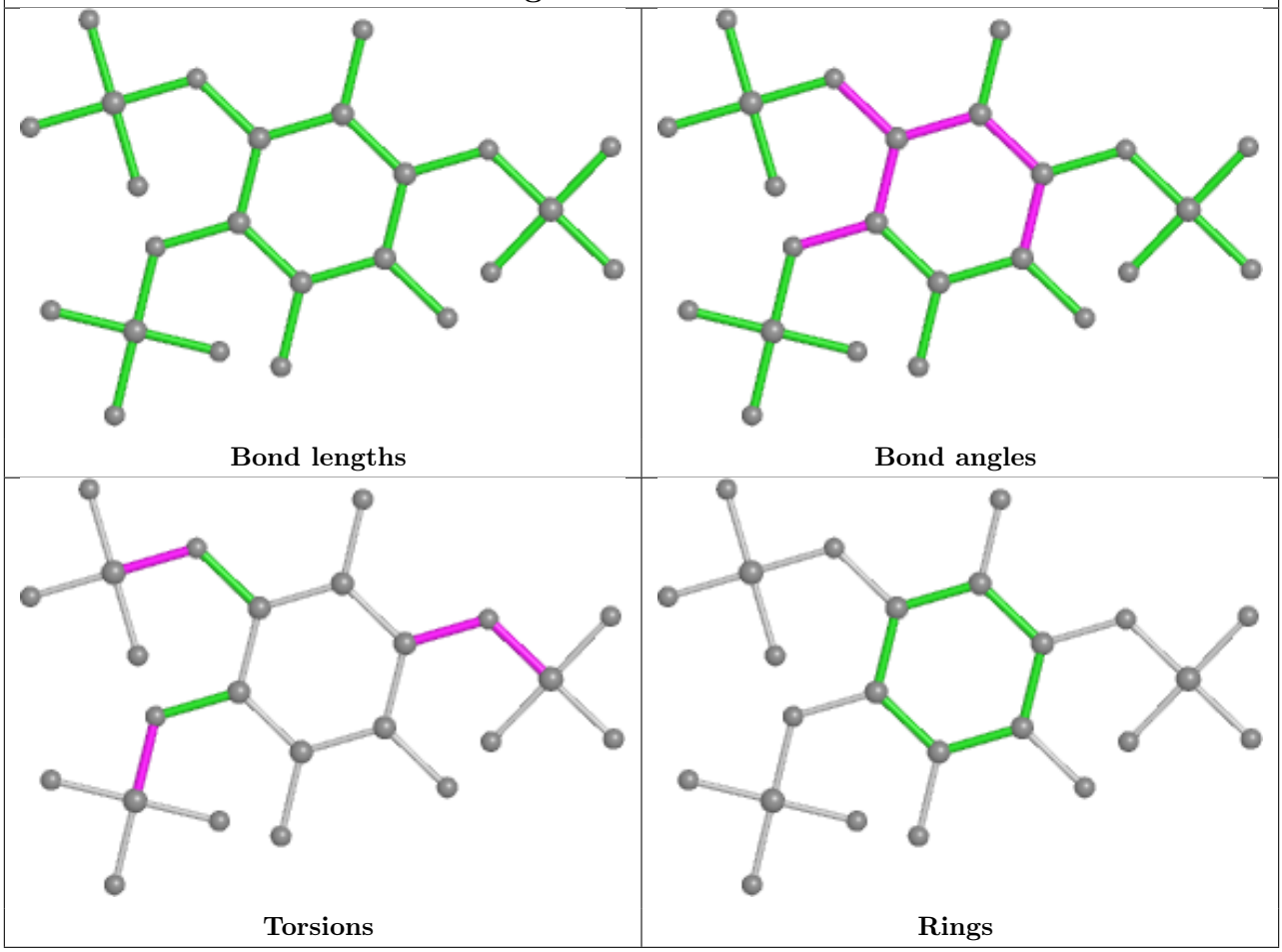
There are no ring outliers.

No monomer is involved in short contacts.

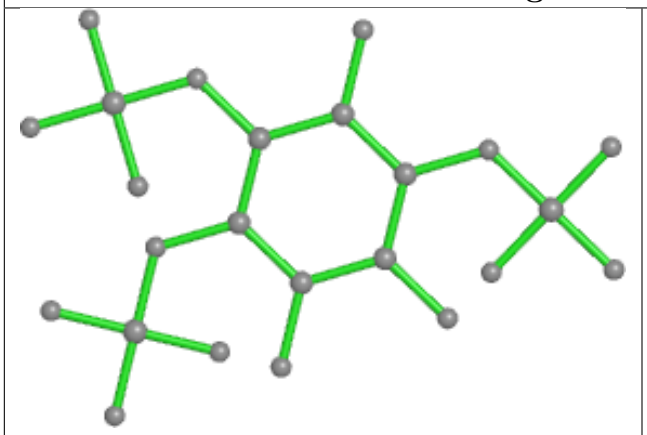
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

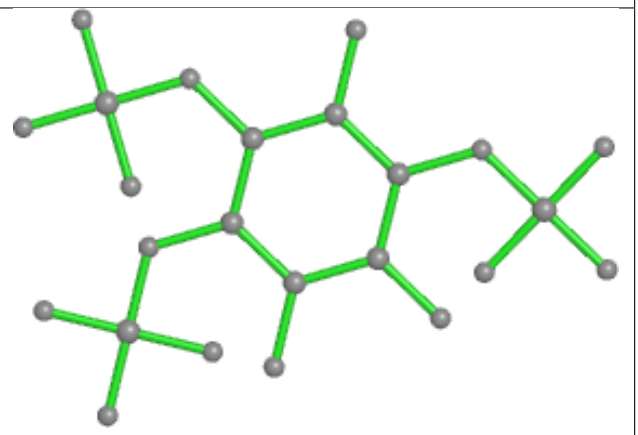
Ligand I3P D 1311



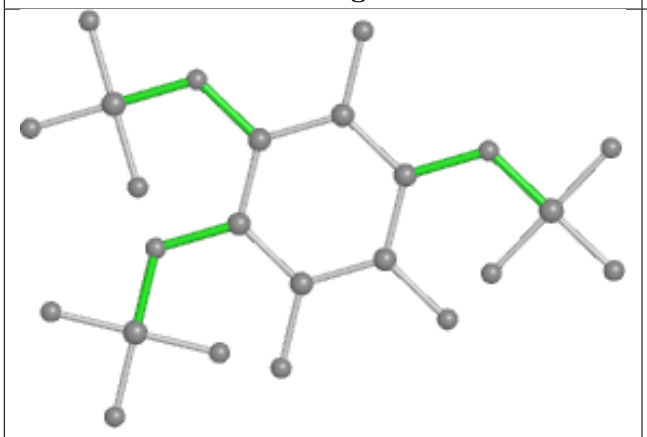
Ligand I3P B 1311



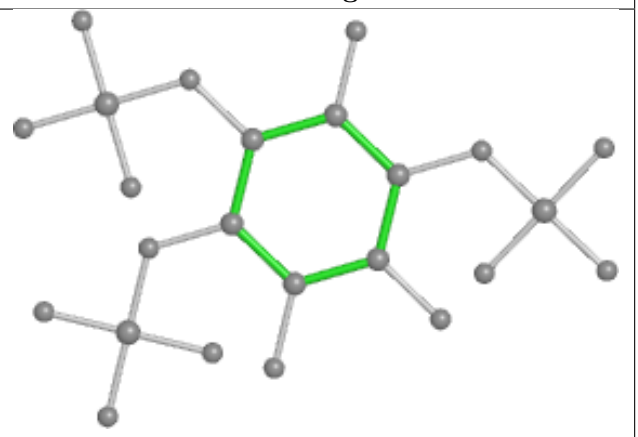
Bond lengths



Bond angles

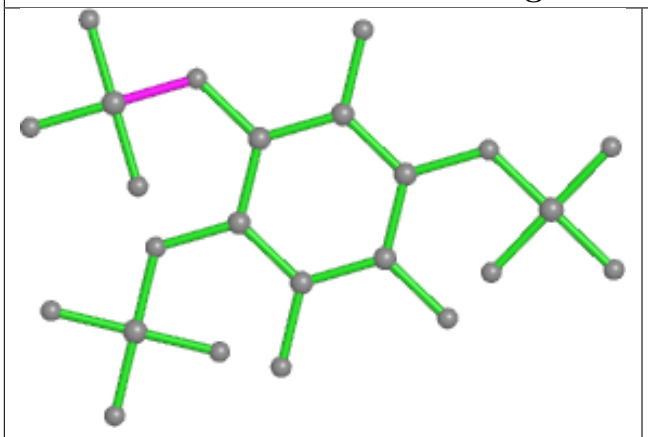


Torsions

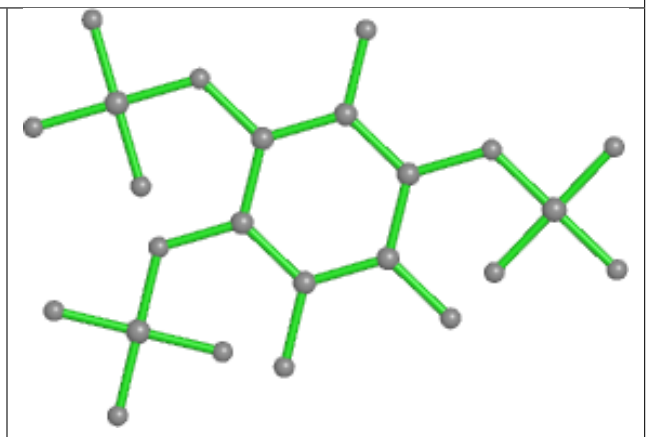


Rings

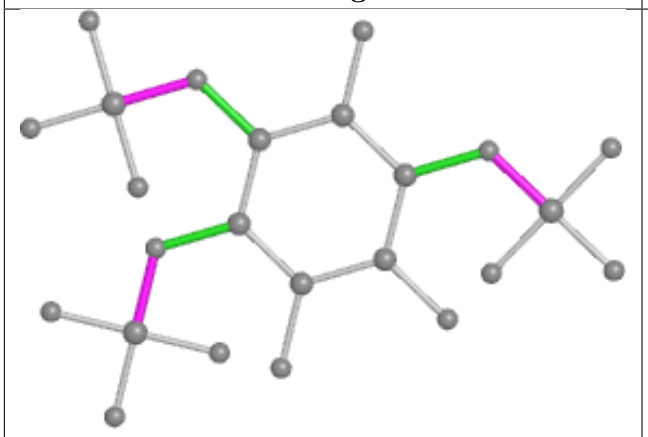
Ligand I3P A 1311



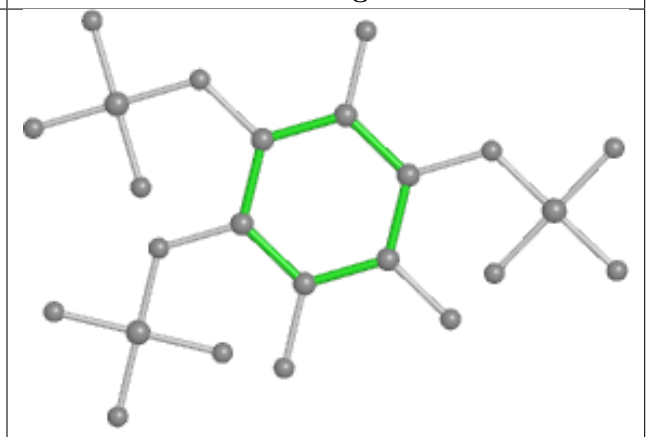
Bond lengths



Bond angles

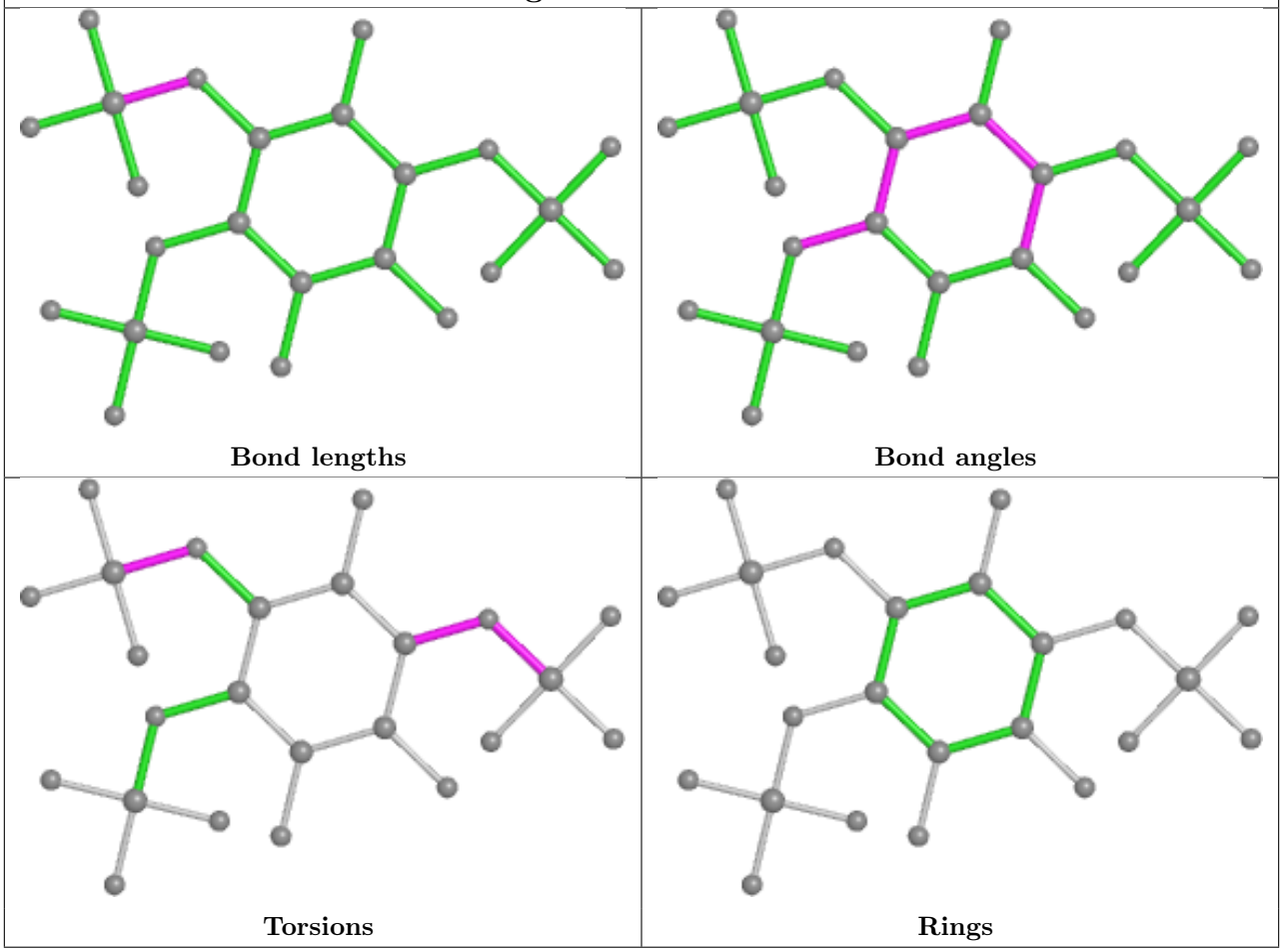


Torsions



Rings

Ligand I3P C 1311



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	C	2
1	B	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	B	1233:ALA	C	1234:TYR	N	3.09
1	C	1233:ALA	C	1234:TYR	N	3.08
1	C	1138:ALA	C	1139:CYS	N	2.21

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	559/572 (97%)	0.22	5 (0%)	84	75	74, 109, 161, 198	0
1	B	558/572 (97%)	0.38	34 (6%)	21	11	92, 151, 226, 251	0
1	C	563/572 (98%)	0.28	10 (1%)	68	55	72, 113, 169, 215	0
1	D	557/572 (97%)	0.48	35 (6%)	20	10	99, 153, 222, 243	0
All	All	2237/2288 (97%)	0.34	84 (3%)	40	25	72, 126, 211, 251	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	833	LEU	9.3
1	D	965	LEU	8.7
1	D	826	PHE	6.2
1	D	1174	LEU	5.7
1	D	966	ILE	5.7
1	D	835	LYS	5.4
1	D	962	LEU	5.1
1	D	763	LEU	5.1
1	D	729	GLY	4.5
1	D	832	ILE	4.4
1	B	965	LEU	4.0
1	B	727	LEU	3.8
1	B	843	LEU	3.7
1	B	742	VAL	3.7
1	B	841	GLU	3.6
1	C	965	LEU	3.6
1	B	753	ILE	3.5
1	C	817	LEU	3.5
1	B	833	LEU	3.5
1	B	826	PHE	3.4
1	A	965	LEU	3.4

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Mol	Chain	Res	Type	RSRZ
1	D	824	ALA	3.4
1	B	1094	GLU	3.2
1	B	723	PHE	3.1
1	B	951	PRO	3.1
1	D	843	LEU	3.1
1	B	814	ARG	3.1
1	B	745	MET	3.0
1	B	1174	LEU	3.0
1	D	1207	LEU	3.0
1	D	760	ALA	2.9
1	D	825	GLY	2.9
1	B	817	LEU	2.9
1	D	836	GLU	2.9
1	C	755	ASP	2.9
1	D	1162	THR	2.8
1	B	830	LEU	2.7
1	D	900	PHE	2.7
1	C	723	PHE	2.7
1	B	963	TYR	2.7
1	B	966	ILE	2.7
1	B	912	MET	2.7
1	D	741	GLN	2.7
1	D	975	PHE	2.6
1	D	834	THR	2.6
1	B	810	SER	2.5
1	C	1065	LEU	2.5
1	B	829	VAL	2.5
1	A	833	LEU	2.5
1	D	743	PHE	2.5
1	D	918	LEU	2.4
1	B	1287	GLY	2.4
1	D	723	PHE	2.4
1	B	1229	LEU	2.4
1	D	904	LEU	2.4
1	C	1070	PHE	2.4
1	C	1062	ASP	2.4
1	D	742	VAL	2.4
1	D	727	LEU	2.3
1	B	918	LEU	2.3
1	B	1052	LEU	2.3
1	D	813	TYR	2.3
1	D	817	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	722	LEU	2.2
1	B	756	PHE	2.2
1	C	835	LYS	2.2
1	B	765	VAL	2.2
1	D	753	ILE	2.2
1	D	854	ILE	2.2
1	D	831	LYS	2.2
1	B	959	VAL	2.1
1	A	835	LYS	2.1
1	C	966	ILE	2.1
1	B	759	LEU	2.1
1	D	956	ILE	2.1
1	B	1240	GLU	2.1
1	A	817	LEU	2.1
1	A	723	PHE	2.1
1	D	728	HIS	2.1
1	B	1129	GLU	2.1
1	B	962	LEU	2.1
1	C	841	GLU	2.0
1	D	1210	ARG	2.0
1	B	750	ASN	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
2	MAN	F	8	11/12	0.58	0.23	145,156,159,159	0
2	MAN	E	8	11/12	0.67	0.26	152,156,159,159	0
2	MAN	F	6	11/12	0.72	0.41	148,149,152,153	0
2	MAN	G	10	11/12	0.79	0.26	127,133,140,143	0
2	MAN	G	6	11/12	0.80	0.16	138,145,149,150	0
2	MAN	E	5	11/12	0.81	0.15	147,150,152,153	0
2	MAN	H	6	11/12	0.82	0.26	142,147,150,150	0

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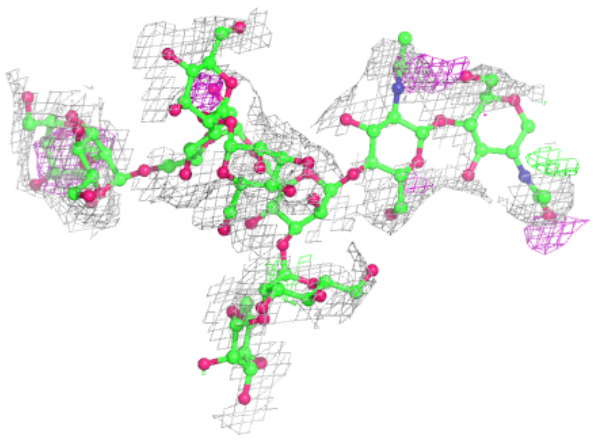
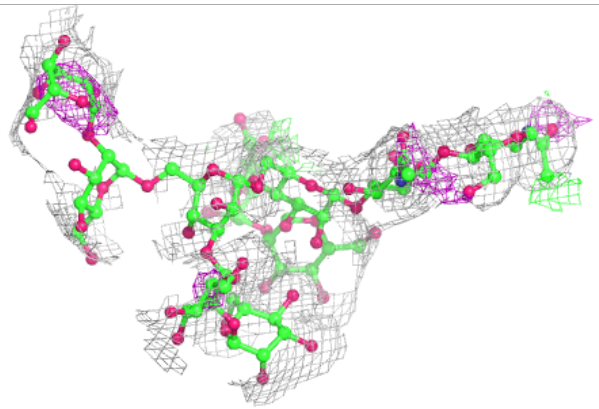
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	MAN	F	5	11/12	0.83	0.19	138,140,143,146	0
2	MAN	E	6	11/12	0.83	0.20	149,154,160,162	0
2	MAN	H	7	11/12	0.83	0.18	141,143,144,146	0
2	MAN	G	8	11/12	0.85	0.18	135,143,146,150	0
2	MAN	F	10	11/12	0.85	0.17	120,125,127,129	0
2	NAG	H	1	14/15	0.86	0.21	111,113,119,121	0
2	MAN	F	4	11/12	0.86	0.19	131,134,140,140	0
2	MAN	E	7	11/12	0.86	0.16	131,134,143,151	0
2	MAN	E	10	11/12	0.87	0.17	138,146,150,150	0
2	MAN	H	8	11/12	0.87	0.19	132,138,140,141	0
2	MAN	H	5	11/12	0.88	0.16	124,135,139,144	0
2	MAN	G	7	11/12	0.88	0.12	126,129,131,135	0
2	NAG	F	2	14/15	0.89	0.22	122,122,123,124	0
2	MAN	H	10	11/12	0.89	0.12	133,136,138,138	0
2	MAN	F	7	11/12	0.90	0.18	136,140,146,152	0
2	BMA	F	3	11/12	0.90	0.15	118,122,125,128	0
2	MAN	H	4	11/12	0.90	0.08	126,129,134,138	0
2	NAG	F	1	14/15	0.90	0.12	109,115,121,122	0
2	MAN	E	4	11/12	0.91	0.10	119,122,130,136	0
2	NAG	E	2	14/15	0.91	0.20	92,95,101,102	0
2	NAG	H	2	14/15	0.91	0.18	108,117,123,127	0
2	BMA	H	3	11/12	0.92	0.09	113,118,120,122	0
2	MAN	G	5	11/12	0.93	0.11	133,137,141,142	0
2	MAN	H	9	11/12	0.93	0.12	121,123,125,131	0
2	MAN	F	9	11/12	0.93	0.15	124,126,128,128	0
2	NAG	E	1	14/15	0.94	0.20	94,96,99,99	0
2	NAG	G	1	14/15	0.94	0.19	78,83,89,94	0
2	MAN	G	4	11/12	0.94	0.13	106,110,120,123	0
2	MAN	E	9	11/12	0.95	0.14	120,127,130,134	0
2	NAG	G	2	14/15	0.95	0.17	75,77,80,82	0
2	BMA	E	3	11/12	0.96	0.16	103,105,113,115	0
2	BMA	G	3	11/12	0.96	0.16	86,89,99,100	0
2	MAN	G	9	11/12	0.97	0.12	107,110,114,123	0

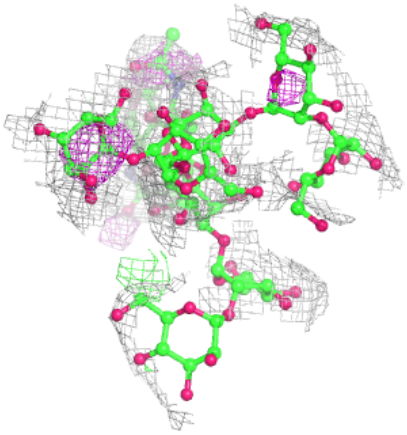
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

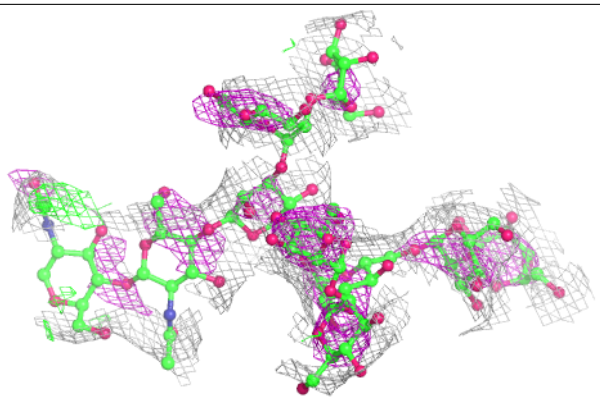


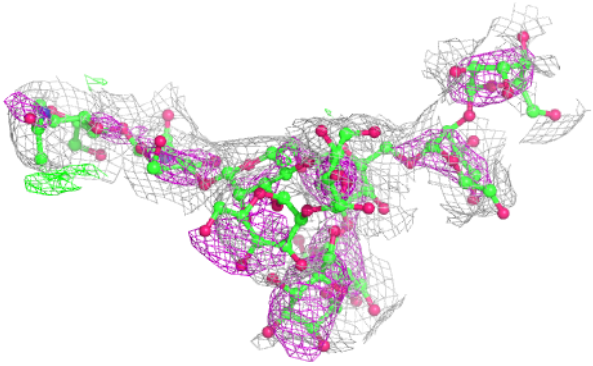


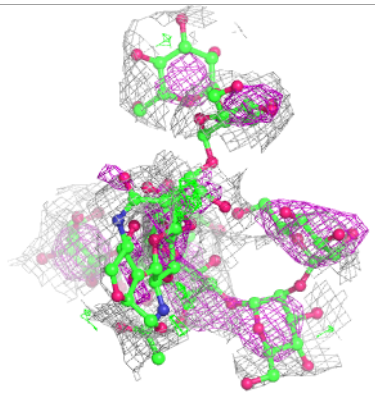


Electron density around Chain F:

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

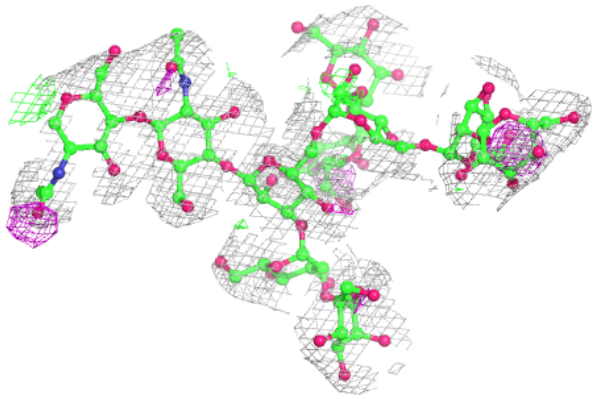


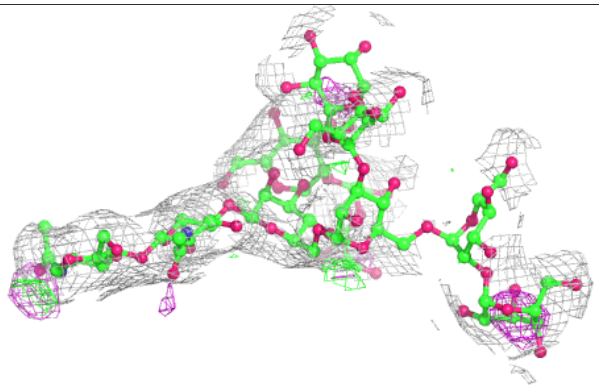


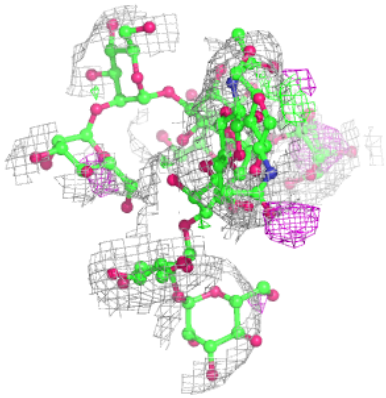


Electron density around Chain G:

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

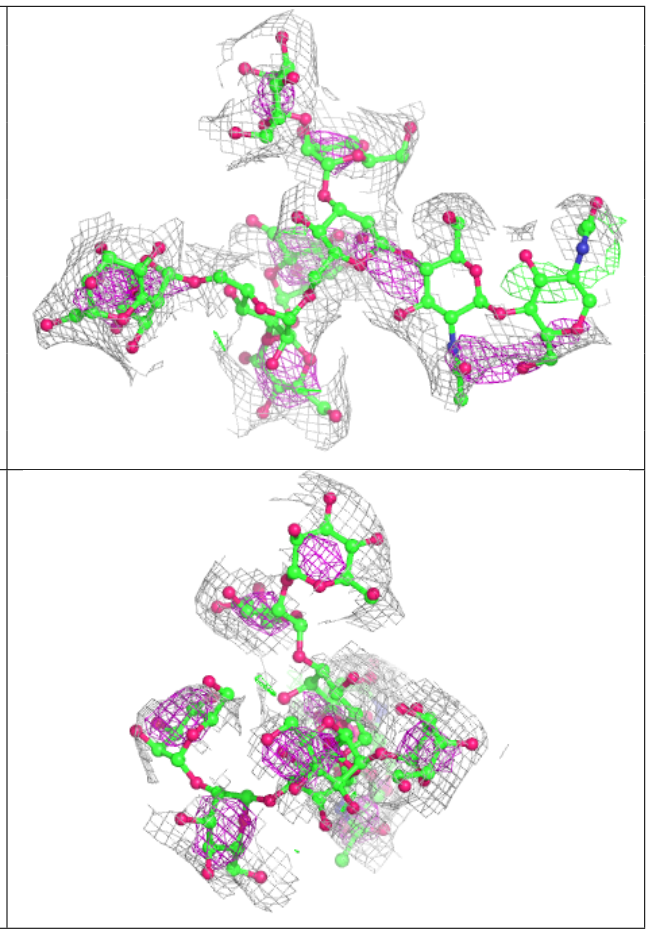






Electron density around Chain H:

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



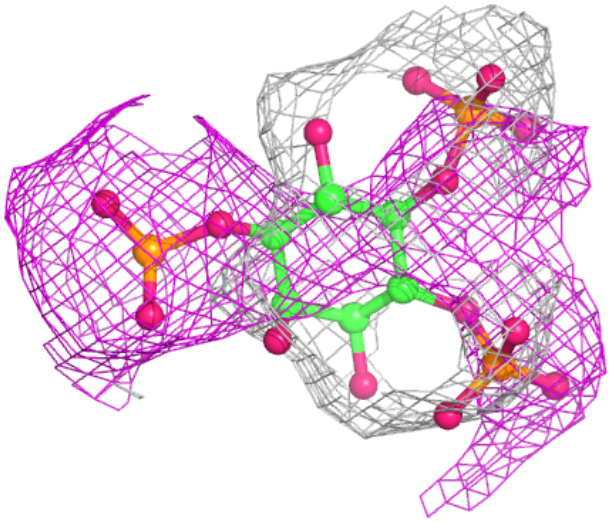
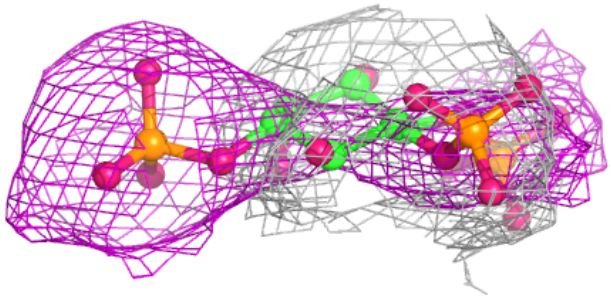
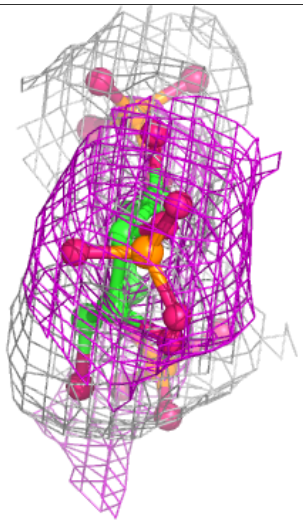
6.4 Ligands (i)

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	I3P	D	1311	24/24	0.62	0.46	190,201,218,219	0
4	I3P	C	1311	24/24	0.69	0.35	166,178,188,190	0
3	NAG	D	1310	14/15	0.69	0.44	100,104,106,106	0
4	I3P	B	1311	24/24	0.70	0.41	204,209,217,218	0
4	I3P	A	1311	24/24	0.72	0.21	185,204,218,219	0
3	NAG	B	1310	14/15	0.84	0.40	82,86,86,87	0
3	NAG	A	1310	14/15	0.90	0.21	115,121,126,130	0
3	NAG	C	1310	14/15	0.92	0.10	96,104,107,108	0

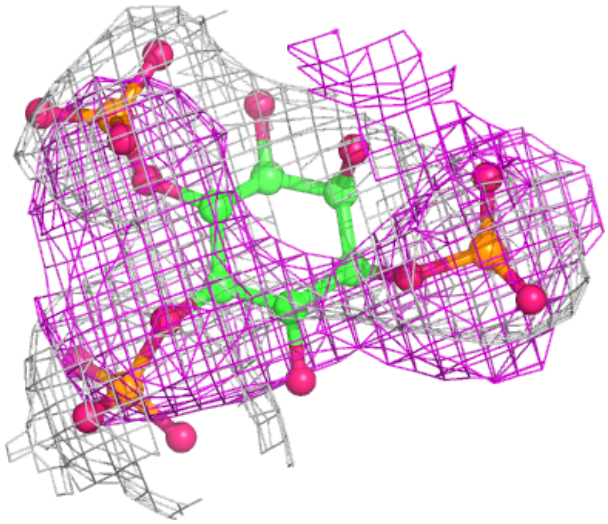
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different

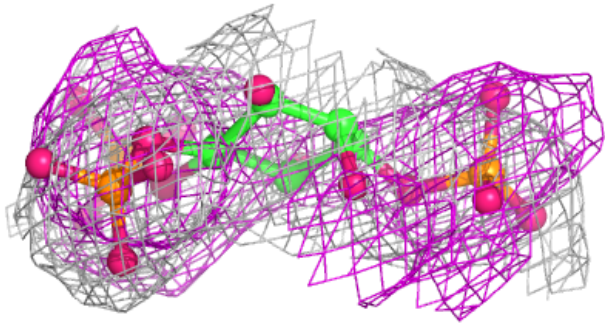
orientation to approximate a three-dimensional view.

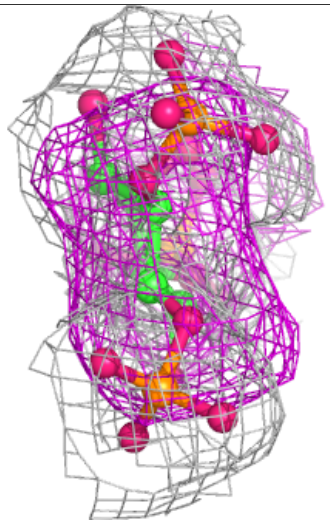
Electron density around I3P D 1311: $2mF_o$-DF_c (at 0.7 rmsd) in gray mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)	
	

Electron density around I3P C 1311:

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

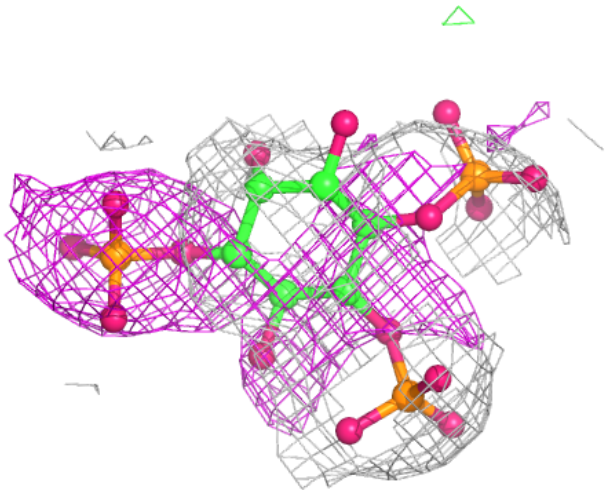


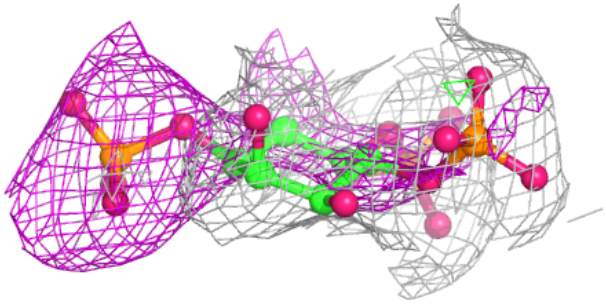


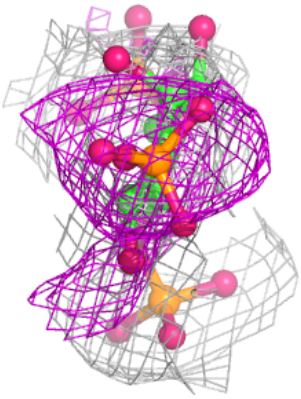


Electron density around I3P B 1311:

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

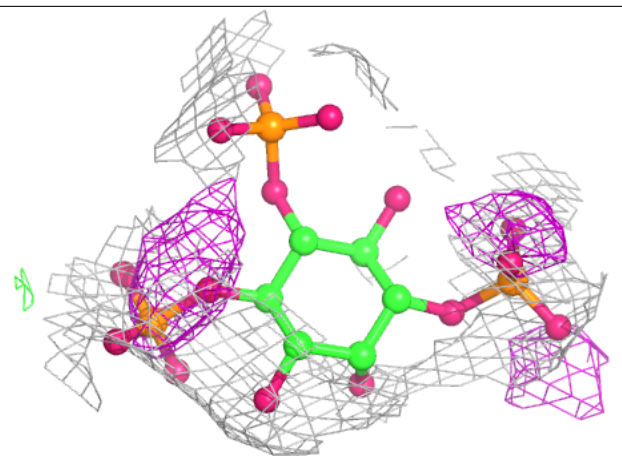


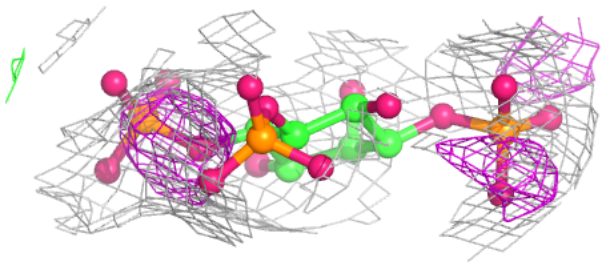


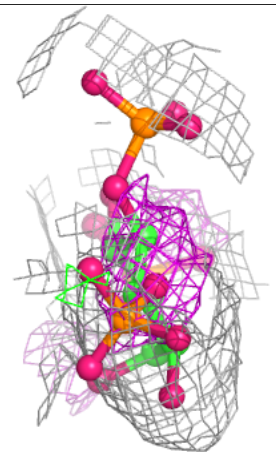


Electron density around I3P A 1311:

$2mF_o$-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.