



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 12:13 AM UTC

PDB ID : 5J67 / pdb_00005j67
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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

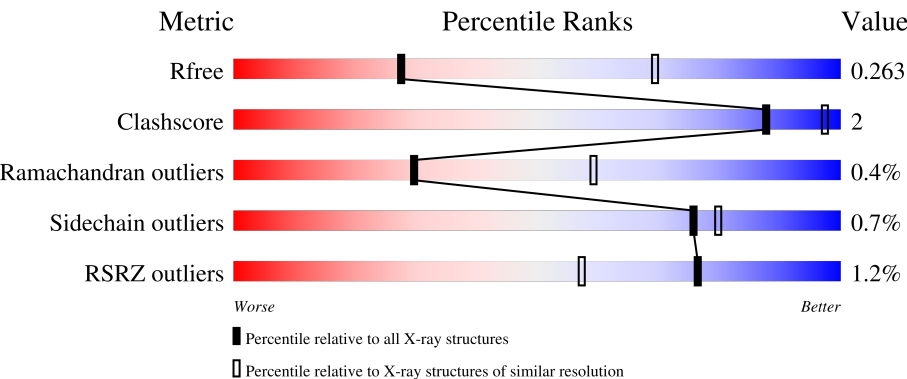
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance 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.



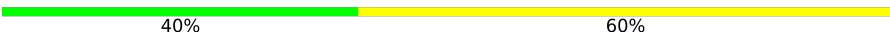
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R _{free}	180053	2361 (3.20-3.12)
Clashscore	190562	2486 (3.20-3.12)
Ramachandran outliers	187476	2405 (3.20-3.12)
Sidechain outliers	187428	2404 (3.20-3.12)
RSRZ outliers	180081	2361 (3.20-3.12)

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 $\leq 5\%$. 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	572	92%6%
1	B	572	% 90%7%
1	C	572	% 92%7%
1	D	572	3% 92%6%
2	E	10	20%80%

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Mol	Chain	Length	Quality of chain
2	F	10	
2	G	10	
2	H	10	

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	8	X	-	-	-
2	MAN	G	8	X	-	-	-
2	MAN	H	8	X	-	-	-

2 Entry composition [i](#)

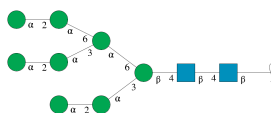
There are 4 unique types of molecules in this entry. The entry contains 18450 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- 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.



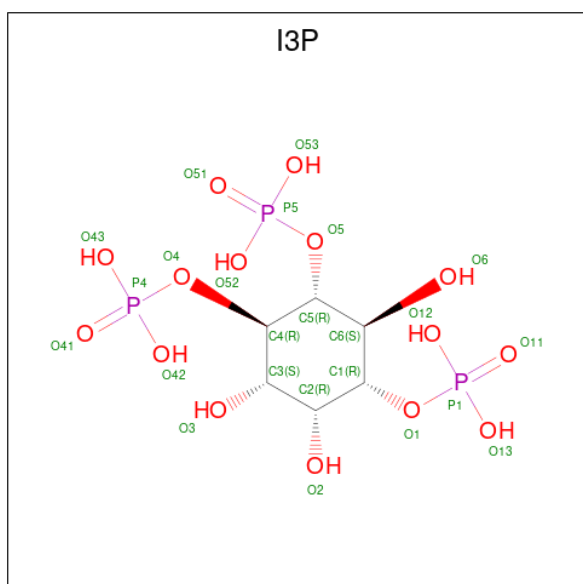
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	10	Total	C	N	O	0	0	0
			116	64	2	50			
2	F	10	Total	C	N	O	0	0	0
			116	64	2	50			
2	G	10	Total	C	N	O	0	0	0
			116	64	2	50			
2	H	10	Total	C	N	O	0	0	0
			116	64	2	50			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

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

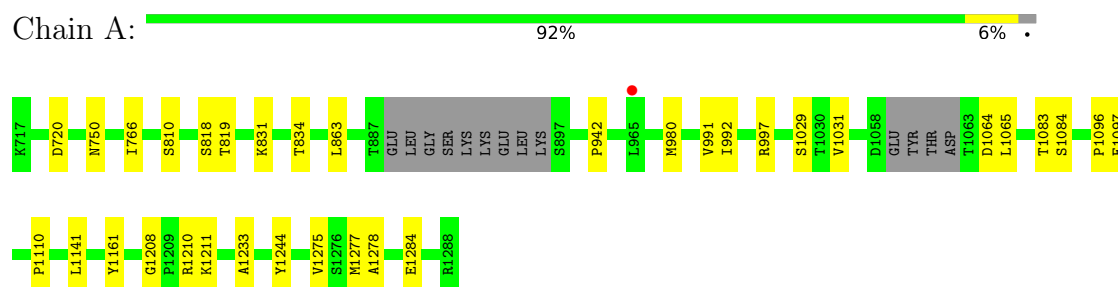


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total 24	C 6	O 15	P 3	0	0
4	B	1	Total 24	C 6	O 15	P 3	0	0
4	C	1	Total 24	C 6	O 15	P 3	0	0
4	D	1	Total 24	C 6	O 15	P 3	0	0

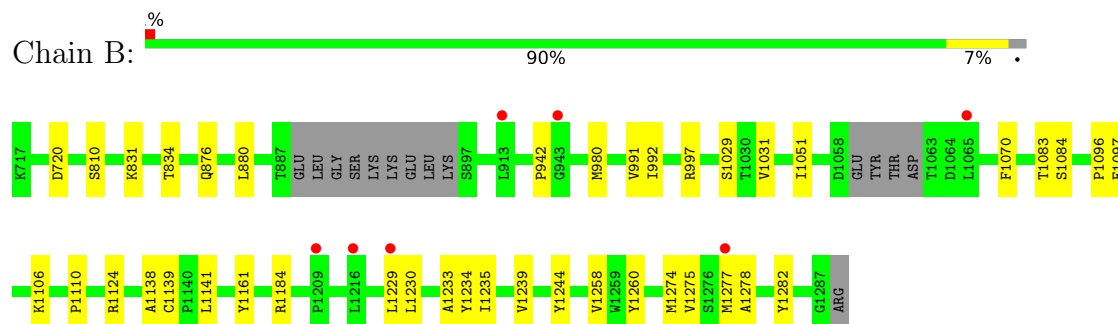
3 Residue-property plots

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.

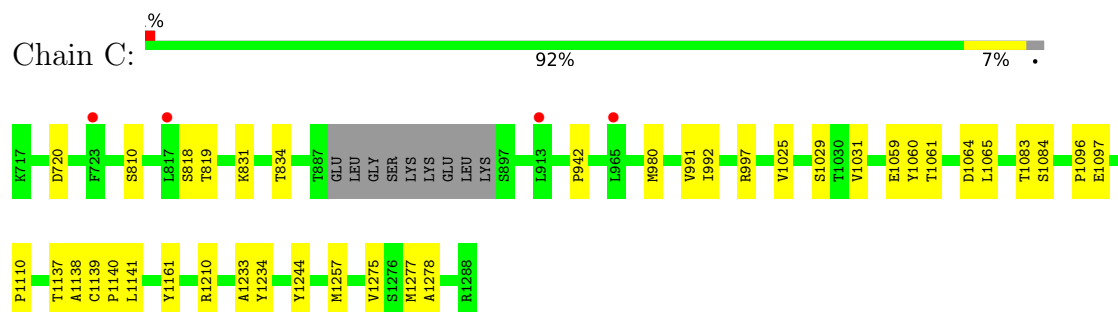
• Molecule 1: Astrotactin-2



• Molecule 1: Astrotactin-2

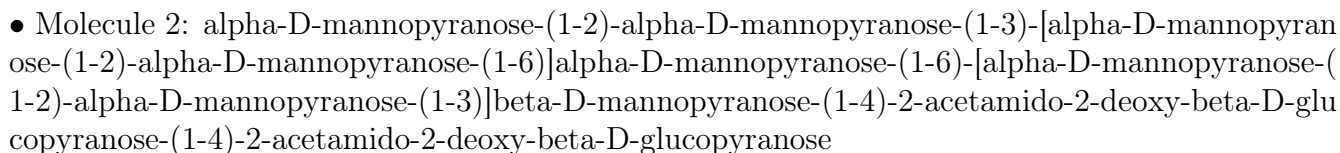


• Molecule 1: Astrotactin-2



• Molecule 1: Astrotactin-2





NAG1
NAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

• 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

NAG1
NAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

- 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

NAG1
NAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

- 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

NAG1
NAG2
BMA3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

4 Data and refinement statistics

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
$\langle I/\sigma(I) \rangle$ 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)	0.31 , 75.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	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.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, BMA, I3P, MAN

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	0.56	0/4552	0.82	2/6166 (0.0%)
1	B	0.56	1/4540 (0.0%)	0.83	1/6149 (0.0%)
1	C	0.57	1/4588 (0.0%)	0.84	1/6214 (0.0%)
1	D	0.55	0/4537	0.82	0/6148
All	All	0.56	2/18217 (0.0%)	0.83	4/24677 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	1064	ASP	N-CA	5.84	1.53	1.46
1	B	1139	CYS	N-CA	5.06	1.53	1.45

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	1210	ARG	NE-CZ-NH2	6.75	125.27	119.20
1	B	1138	ALA	CA-C-O	-5.39	115.84	121.99
1	A	1208	GLY	CA-C-N	5.03	124.47	119.24
1	A	1208	GLY	C-N-CA	5.03	124.47	119.24

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	16	1
1	B	4445	0	4372	29	0
1	C	4492	0	4411	24	0
1	D	4441	0	4363	12	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	73	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 (73) 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	0.98
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:1031:VAL:HG11	1:B:1277:MET:HE2	1.64	0.79
1:A:766:ILE:HG23	1:B:1274:MET:HE3	1.64	0.79
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:D:1031:VAL:HG11	1:D:1277:MET:HE2	1.81	0.62
1:A:1031:VAL:HG11	1:A:1277:MET:HE2	1.81	0.62
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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:C:1031:VAL:HG11	1:C:1277:MET:HE2	1.90	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:1137:THR:O	1:C:1138:ALA:C	2.55	0.49
1:C:1029:SER:HB2	1:C:1110:PRO:HB3	1.93	0.49
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:A:766:ILE:CG2	1:B:1274:MET:HE3	2.39	0.47
1:C:1244:TYR:CD1	1:C:1275:VAL:HG23	2.51	0.46
1:A:863:LEU:HD21	1:C:1257:MET:HE2	1.98	0.46
1:B:1244:TYR:CD1	1:B:1275:VAL:HG23	2.52	0.45
1:C:1138:ALA:C	1:C:1139:CYS:CA	2.90	0.45
1:B:1051:ILE:HD12	1:B:1070:PHE:HE1	1.81	0.44
1:A:831:LYS:O	1:A:834:THR:HG22	2.18	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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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:C:1060:TYR:C	1:C:1061:THR:HG1	2.22	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
1:D:1141:LEU:HD21	1:D:1278:ALA:HB1	2.02	0.41
1:C:1025:VAL:CG1	1:C:1277:MET:HE1	2.51	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%)	24	55
1	B	550/572 (96%)	520 (94%)	28 (5%)	2 (0%)	30	59
1	C	555/572 (97%)	524 (94%)	29 (5%)	2 (0%)	30	59
1	D	551/572 (96%)	521 (95%)	28 (5%)	2 (0%)	30	59
All	All	2209/2288 (96%)	2086 (94%)	114 (5%)	9 (0%)	30	59

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
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 ⓘ

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%)	78	81
1	B	500/513 (98%)	498 (100%)	2 (0%)	84	84
1	C	505/513 (98%)	502 (99%)	3 (1%)	78	81
1	D	500/513 (98%)	494 (99%)	6 (1%)	63	74
All	All	2006/2052 (98%)	1992 (99%)	14 (1%)	76	80

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	720	ASP
1	A	810	SER
1	A	1065	LEU
1	B	720	ASP
1	B	810	SER
1	C	720	ASP
1	C	810	SER
1	C	1059	GLU
1	D	720	ASP
1	D	810	SER
1	D	1056	LYS

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Mol	Chain	Res	Type
1	D	1059	GLU
1	D	1065	LEU
1	D	1284	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	828	ASN
1	A	848	HIS
1	A	881	GLN
1	A	883	GLN
1	A	946	GLN
1	A	1054	HIS
1	A	1172	ASN
1	B	736	HIS
1	B	828	ASN
1	B	848	HIS
1	B	881	GLN
1	B	883	GLN
1	B	946	GLN
1	B	1192	HIS
1	C	828	ASN
1	C	848	HIS
1	C	881	GLN
1	C	883	GLN
1	C	946	GLN
1	C	1054	HIS
1	C	1172	ASN
1	D	828	ASN
1	D	848	HIS
1	D	881	GLN
1	D	883	GLN
1	D	946	GLN
1	D	1054	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	2.64	8 (47%)
2	MAN	E	10	2	11,11,12	0.38	0	15,15,17	0.85	1 (6%)
2	NAG	E	2	2	14,14,15	0.31	0	17,19,21	0.80	0
2	BMA	E	3	2	11,11,12	0.20	0	15,15,17	0.79	0
2	MAN	E	4	2	11,11,12	0.42	0	15,15,17	1.17	2 (13%)
2	MAN	E	5	2	11,11,12	0.46	0	15,15,17	1.01	1 (6%)
2	MAN	E	6	2	11,11,12	0.31	0	15,15,17	0.95	1 (6%)
2	MAN	E	7	2	11,11,12	0.63	0	15,15,17	1.12	2 (13%)
2	MAN	E	8	2	11,11,12	0.59	0	15,15,17	2.03	2 (13%)
2	MAN	E	9	2	11,11,12	0.29	0	15,15,17	0.65	1 (6%)
2	NAG	F	1	1,2	14,14,15	0.47	0	17,19,21	2.48	5 (29%)
2	MAN	F	10	2	11,11,12	0.43	0	15,15,17	0.80	0
2	NAG	F	2	2	14,14,15	0.38	0	17,19,21	0.73	0
2	BMA	F	3	2	11,11,12	0.30	0	15,15,17	0.62	0
2	MAN	F	4	2	11,11,12	0.45	0	15,15,17	0.96	1 (6%)
2	MAN	F	5	2	11,11,12	0.42	0	15,15,17	1.33	1 (6%)
2	MAN	F	6	2	11,11,12	0.29	0	15,15,17	0.78	1 (6%)
2	MAN	F	7	2	11,11,12	0.46	0	15,15,17	1.32	2 (13%)
2	MAN	F	8	2	11,11,12	0.47	0	15,15,17	1.79	1 (6%)
2	MAN	F	9	2	11,11,12	0.33	0	15,15,17	0.99	1 (6%)
2	NAG	G	1	1,2	14,14,15	0.46	0	17,19,21	2.17	5 (29%)
2	MAN	G	10	2	11,11,12	0.35	0	15,15,17	0.80	1 (6%)
2	NAG	G	2	2	14,14,15	0.45	0	17,19,21	0.75	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	BMA	G	3	2	11,11,12	0.26	0	15,15,17	0.69	0
2	MAN	G	4	2	11,11,12	0.29	0	15,15,17	0.70	1 (6%)
2	MAN	G	5	2	11,11,12	0.30	0	15,15,17	0.72	0
2	MAN	G	6	2	11,11,12	0.37	0	15,15,17	0.93	1 (6%)
2	MAN	G	7	2	11,11,12	0.46	0	15,15,17	1.05	1 (6%)
2	MAN	G	8	2	11,11,12	0.44	0	15,15,17	1.59	2 (13%)
2	MAN	G	9	2	11,11,12	0.25	0	15,15,17	0.59	0
2	NAG	H	1	1,2	14,14,15	0.44	0	17,19,21	2.52	5 (29%)
2	MAN	H	10	2	11,11,12	0.44	0	15,15,17	0.91	1 (6%)
2	NAG	H	2	2	14,14,15	0.36	0	17,19,21	0.84	0
2	BMA	H	3	2	11,11,12	0.28	0	15,15,17	0.97	0
2	MAN	H	4	2	11,11,12	0.39	0	15,15,17	0.56	0
2	MAN	H	5	2	11,11,12	0.41	0	15,15,17	1.03	1 (6%)
2	MAN	H	6	2	11,11,12	0.38	0	15,15,17	0.80	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	1.49	2 (13%)
2	MAN	H	9	2	11,11,12	0.35	0	15,15,17	0.88	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
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	-	1/2/19/22	0/1/1/1
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

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
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 (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	1	NAG	C1-O5-C5	8.39	123.43	112.19
2	F	1	NAG	C1-O5-C5	8.19	123.16	112.19
2	E	1	NAG	C1-O5-C5	7.86	122.72	112.19
2	E	8	MAN	C1-O5-C5	6.80	121.30	112.19
2	G	1	NAG	C1-O5-C5	6.36	120.71	112.19
2	F	8	MAN	C1-O5-C5	6.17	120.45	112.19
2	G	8	MAN	C1-O5-C5	5.08	118.99	112.19
2	F	5	MAN	C1-O5-C5	4.47	118.17	112.19
2	H	8	MAN	C1-O5-C5	4.45	118.15	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	1	NAG	C8-C7-N2	3.97	122.71	116.12
2	E	1	NAG	C8-C7-N2	3.42	121.80	116.12
2	E	6	MAN	C1-O5-C5	3.06	116.29	112.19
2	F	7	MAN	C3-C4-C5	3.06	115.78	110.23
2	H	5	MAN	C1-O5-C5	3.03	116.24	112.19
2	E	8	MAN	C1-C2-C3	2.98	113.99	109.64
2	E	1	NAG	C1-C2-N2	-2.97	105.75	110.43
2	F	1	NAG	C8-C7-N2	2.94	120.99	116.12
2	E	1	NAG	O5-C5-C4	2.93	117.96	110.83
2	H	1	NAG	C8-C7-N2	2.91	120.95	116.12
2	H	1	NAG	O5-C5-C4	2.79	117.60	110.83
2	H	10	MAN	C1-O5-C5	2.78	115.92	112.19
2	F	1	NAG	O5-C5-C4	2.78	117.59	110.83
2	F	7	MAN	C1-O5-C5	2.75	115.88	112.19
2	E	1	NAG	O5-C1-C2	2.73	115.52	111.29
2	F	4	MAN	C1-O5-C5	2.69	115.79	112.19
2	F	9	MAN	C1-O5-C5	2.63	115.71	112.19
2	G	6	MAN	C1-O5-C5	2.62	115.70	112.19
2	H	8	MAN	C3-C4-C5	2.61	114.97	110.23
2	E	4	MAN	C3-C4-C5	2.60	114.95	110.23
2	G	8	MAN	C3-C4-C5	2.59	114.93	110.23
2	H	7	MAN	C1-O5-C5	2.54	115.58	112.19
2	F	6	MAN	C1-O5-C5	2.52	115.57	112.19
2	H	6	MAN	C1-O5-C5	2.52	115.56	112.19
2	E	5	MAN	C1-O5-C5	2.51	115.55	112.19
2	E	7	MAN	C2-C3-C4	2.43	115.13	110.86
2	E	1	NAG	C2-N2-C7	2.40	126.12	122.90
2	F	1	NAG	C2-N2-C7	2.37	126.08	122.90
2	G	7	MAN	C3-C4-C5	2.35	114.50	110.23
2	H	1	NAG	C4-C3-C2	-2.32	107.61	111.02
2	E	1	NAG	O7-C7-N2	-2.32	117.88	121.98
2	E	10	MAN	C1-O5-C5	2.28	115.24	112.19
2	E	1	NAG	C4-C3-C2	-2.22	107.76	111.02
2	F	1	NAG	C4-C3-C2	-2.17	107.84	111.02
2	H	1	NAG	C2-N2-C7	2.16	125.79	122.90
2	G	4	MAN	C1-O5-C5	2.14	115.05	112.19
2	G	1	NAG	C2-N2-C7	2.12	125.74	122.90
2	E	4	MAN	C1-O5-C5	2.11	115.02	112.19
2	E	7	MAN	C3-C4-C5	2.11	114.05	110.23
2	E	9	MAN	C1-O5-C5	2.09	114.98	112.19
2	G	10	MAN	C1-O5-C5	2.07	114.96	112.19
2	G	1	NAG	O7-C7-C8	-2.02	118.46	122.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
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 (44) 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	G	8	MAN	O5-C5-C6-O6
2	H	8	MAN	O5-C5-C6-O6
2	E	4	MAN	C4-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	8	MAN	C4-C5-C6-O6
2	F	6	MAN	C4-C5-C6-O6
2	H	9	MAN	C4-C5-C6-O6
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	8	MAN	C4-C5-C6-O6

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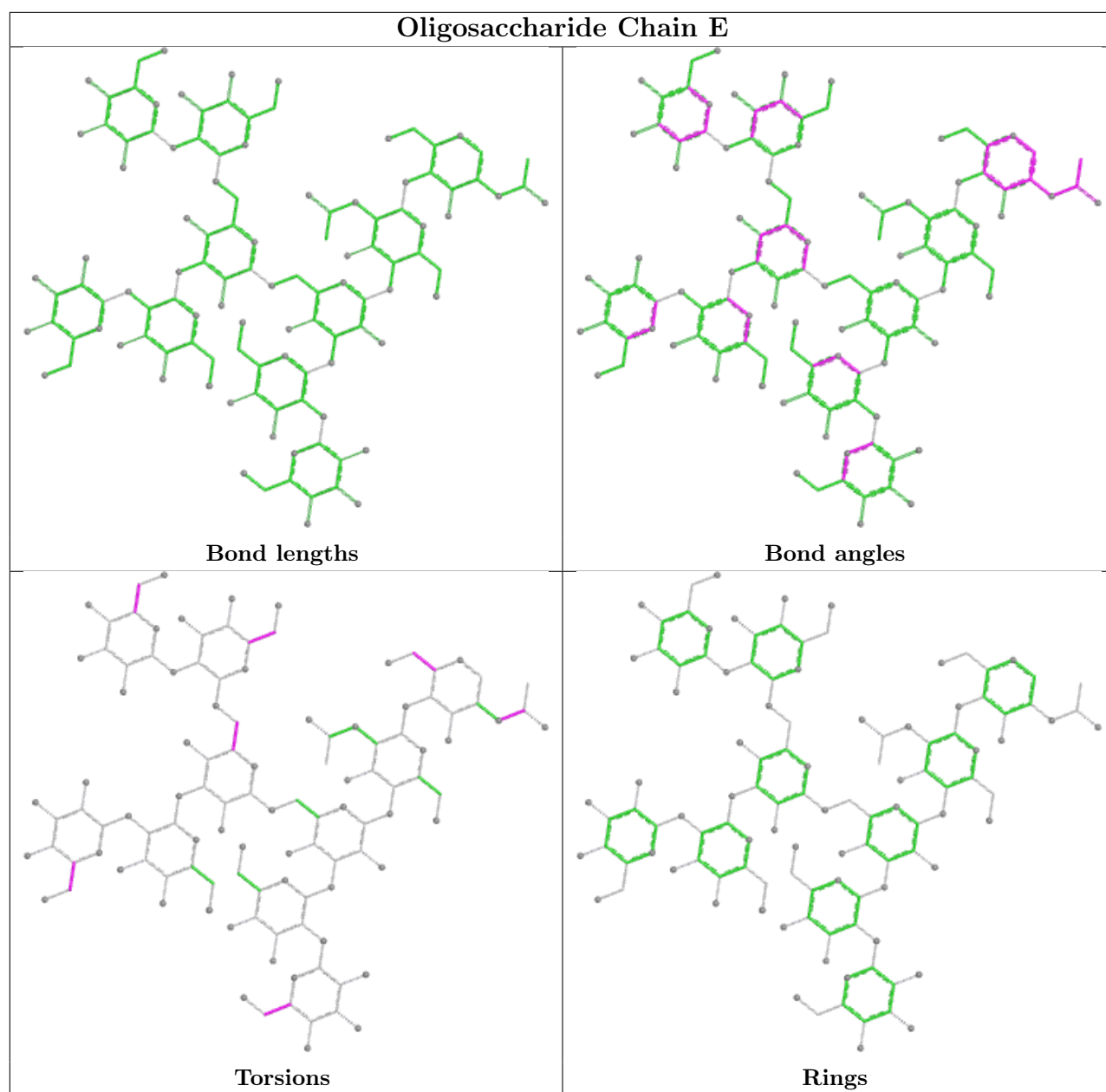
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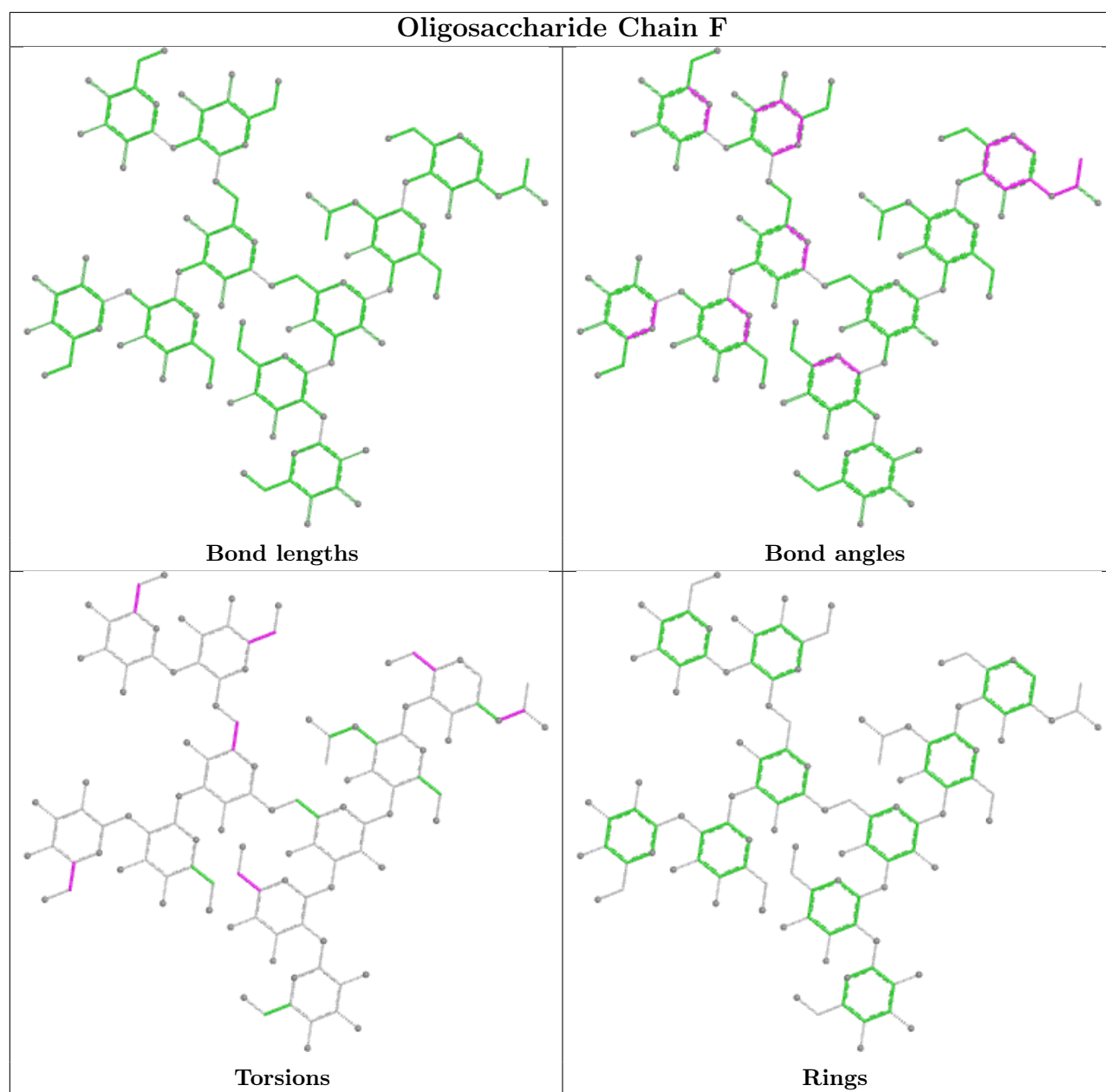
Mol	Chain	Res	Type	Atoms
2	H	9	MAN	O5-C5-C6-O6
2	G	8	MAN	C4-C5-C6-O6
2	H	6	MAN	O5-C5-C6-O6
2	F	4	MAN	O5-C5-C6-O6
2	F	9	MAN	O5-C5-C6-O6
2	E	1	NAG	O5-C5-C6-O6
2	G	6	MAN	C4-C5-C6-O6
2	F	7	MAN	O5-C5-C6-O6
2	H	7	MAN	O5-C5-C6-O6
2	H	1	NAG	O5-C5-C6-O6
2	G	7	MAN	O5-C5-C6-O6
2	F	1	NAG	O5-C5-C6-O6
2	E	8	MAN	C4-C5-C6-O6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

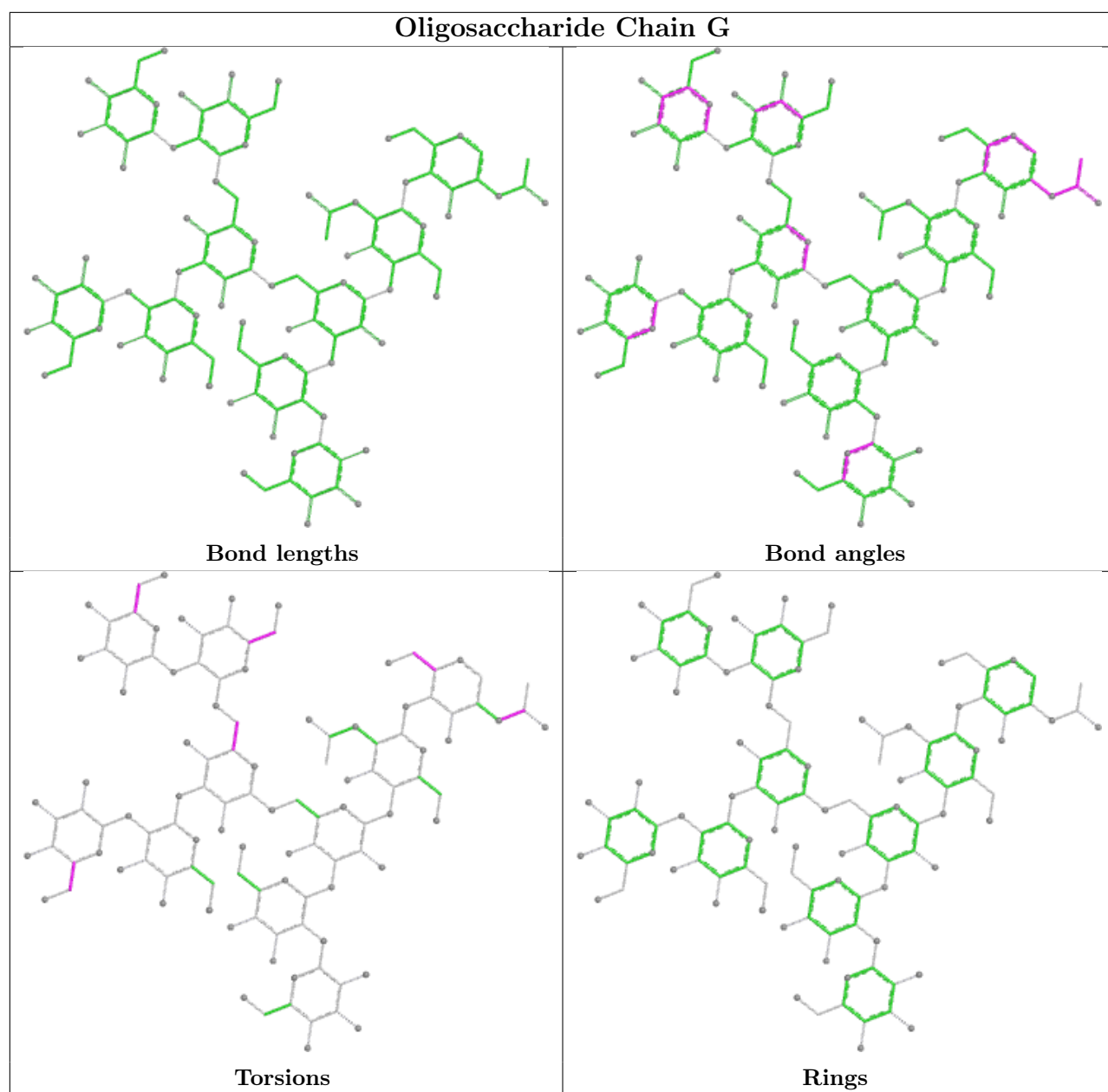
There are no ring outliers.

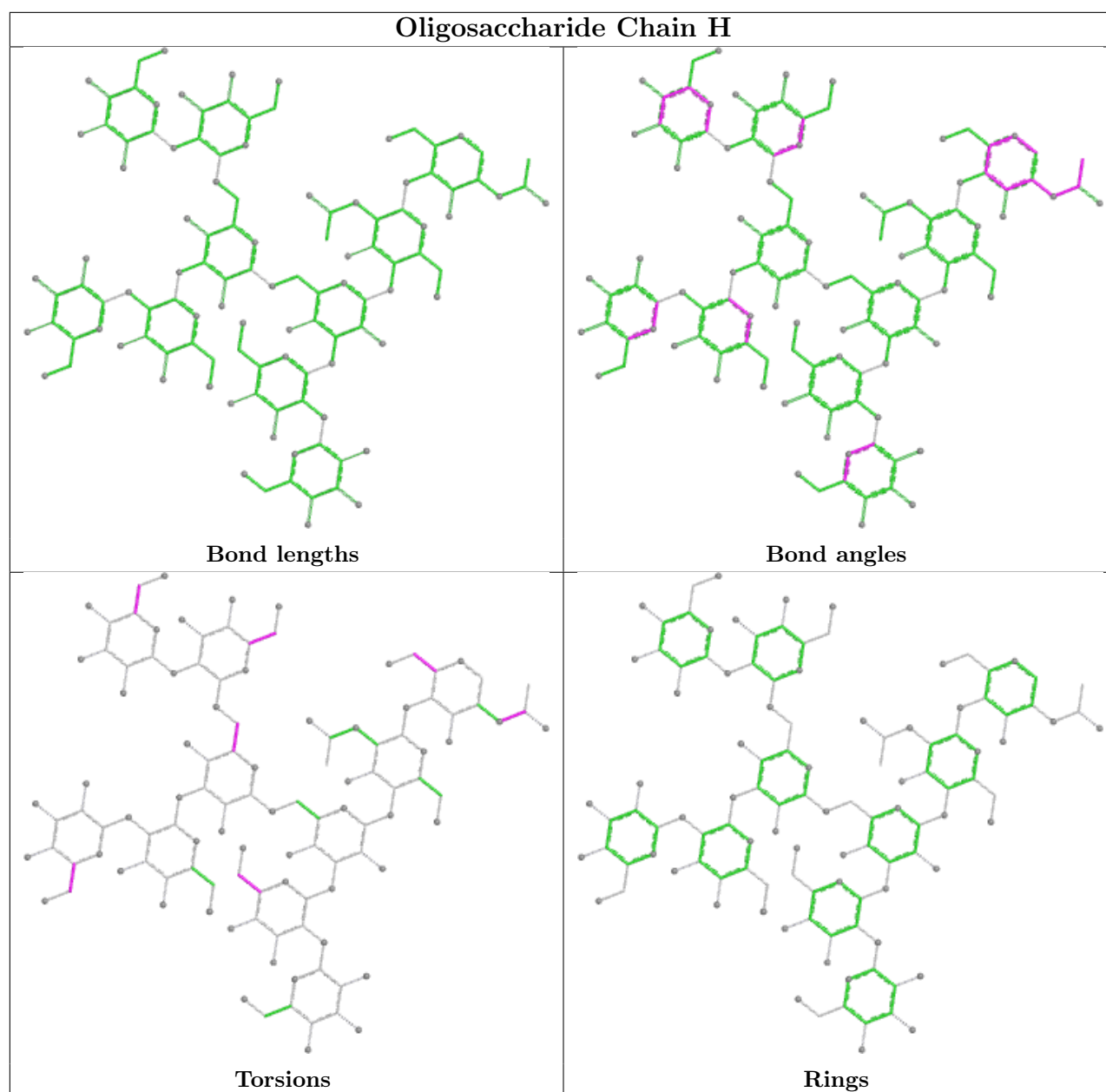
No monomer is involved in short contacts.

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









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	B	1311	-	24,24,24	0.73	0	39,39,39	1.04	0
4	I3P	A	1311	-	24,24,24	0.80	1 (4%)	39,39,39	0.97	0
3	NAG	C	1310	1	14,14,15	0.36	0	17,19,21	1.60	3 (17%)
4	I3P	D	1311	-	24,24,24	0.81	0	39,39,39	1.20	4 (10%)
3	NAG	A	1310	1	14,14,15	0.30	0	17,19,21	1.09	2 (11%)
3	NAG	D	1310	1	14,14,15	0.38	0	17,19,21	0.80	0
3	NAG	B	1310	1	14,14,15	0.40	0	17,19,21	0.82	0
4	I3P	C	1311	-	24,24,24	0.89	1 (4%)	39,39,39	1.30	6 (15%)

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	B	1311	-	-	0/15/39/39	0/1/1/1
4	I3P	A	1311	-	-	2/15/39/39	0/1/1/1
3	NAG	C	1310	1	-	2/6/23/26	0/1/1/1
4	I3P	D	1311	-	-	6/15/39/39	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
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	2.17	1.63	1.59
4	A	1311	I3P	P5-O5	2.07	1.63	1.59

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	1310	NAG	C1-O5-C5	4.95	118.82	112.19
4	C	1311	I3P	C6-C1-C2	3.65	115.92	110.86
3	C	1310	NAG	O5-C1-C2	-2.88	106.84	111.29
4	D	1311	I3P	O4-C4-C5	2.71	114.54	108.76
4	C	1311	I3P	C5-C6-C1	2.63	114.44	109.11
4	D	1311	I3P	C6-C1-C2	2.52	114.35	110.86

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1310	NAG	C1-O5-C5	2.40	115.40	112.19
4	C	1311	I3P	O4-C4-C5	2.29	113.64	108.76
3	A	1310	NAG	O5-C1-C2	-2.25	107.81	111.29
4	D	1311	I3P	C6-C5-C4	-2.20	106.77	111.72
4	D	1311	I3P	O5-C5-C4	2.20	113.44	108.76
4	C	1311	I3P	O5-C5-C4	2.13	113.29	108.76
3	C	1310	NAG	C3-C4-C5	2.09	114.02	110.23
4	C	1311	I3P	C6-C5-C4	-2.03	107.15	111.72
4	C	1311	I3P	C3-C4-C5	-2.02	107.17	111.72

There are no chirality outliers.

All (20) torsion outliers are listed below:

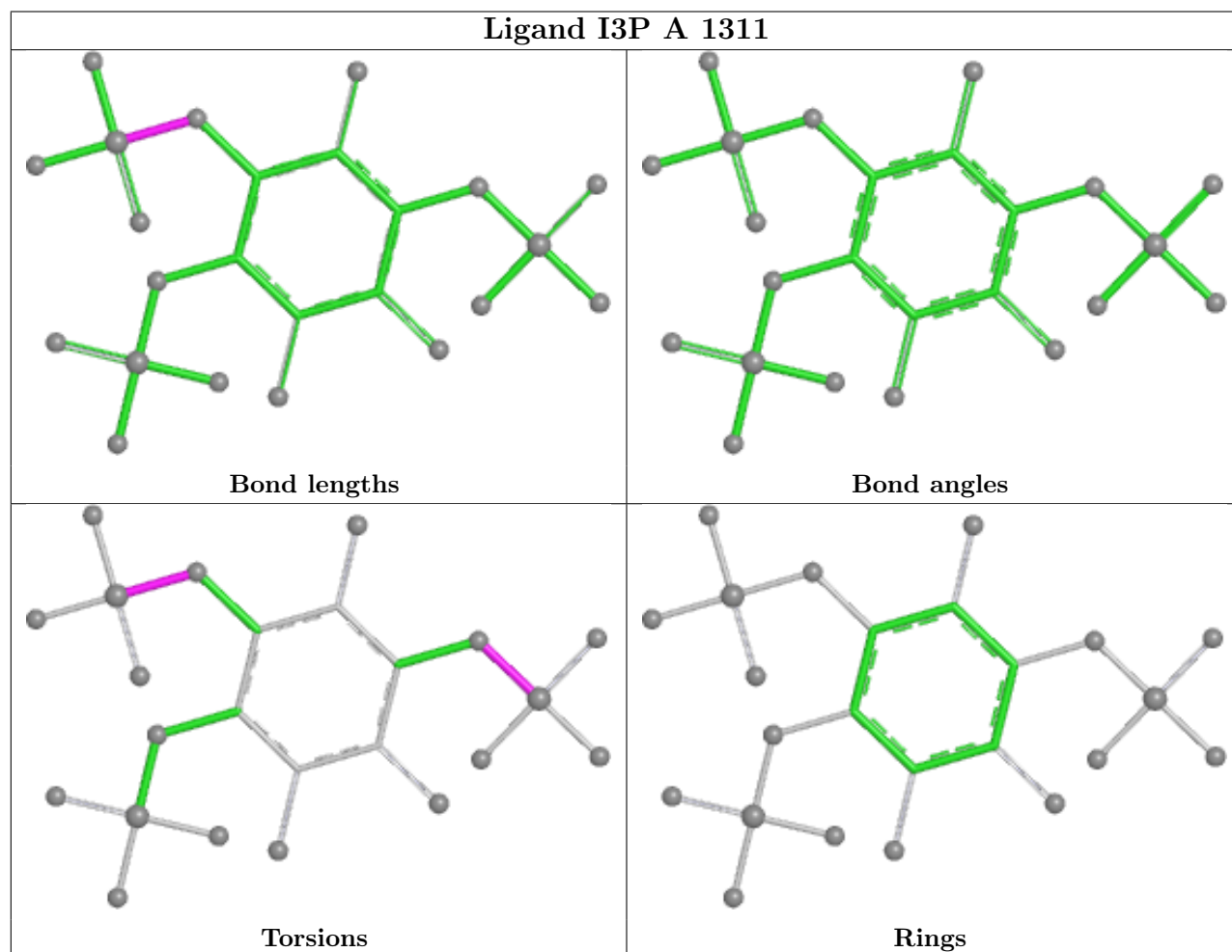
Mol	Chain	Res	Type	Atoms
4	C	1311	I3P	C2-C1-O1-P1
4	C	1311	I3P	C6-C1-O1-P1
4	D	1311	I3P	C2-C1-O1-P1
4	D	1311	I3P	C6-C1-O1-P1
3	B	1310	NAG	O5-C5-C6-O6
3	B	1310	NAG	C4-C5-C6-O6
3	C	1310	NAG	C4-C5-C6-O6
3	C	1310	NAG	O5-C5-C6-O6
3	A	1310	NAG	C4-C5-C6-O6
3	D	1310	NAG	C4-C5-C6-O6
3	A	1310	NAG	O5-C5-C6-O6
3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

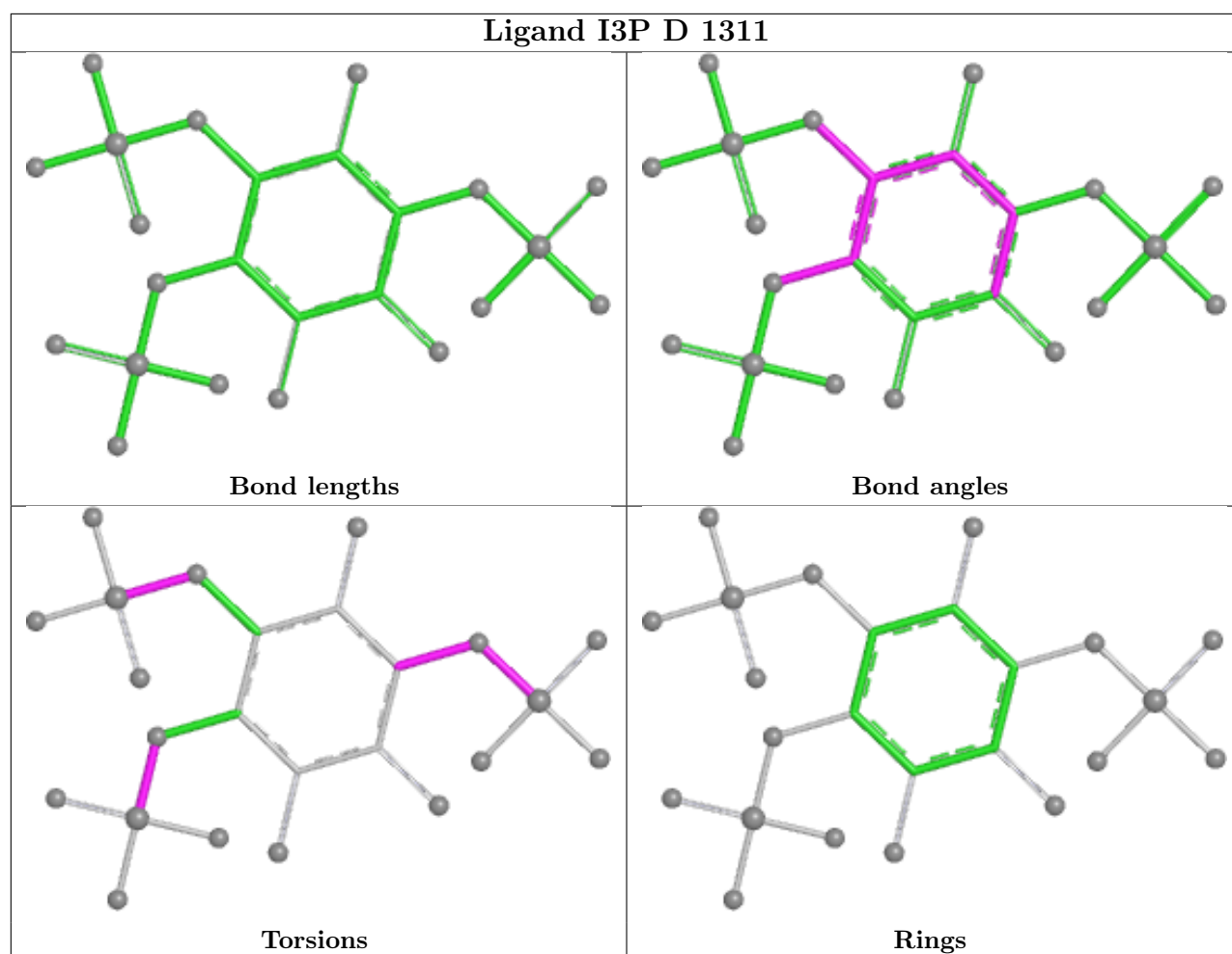
There are no ring outliers.

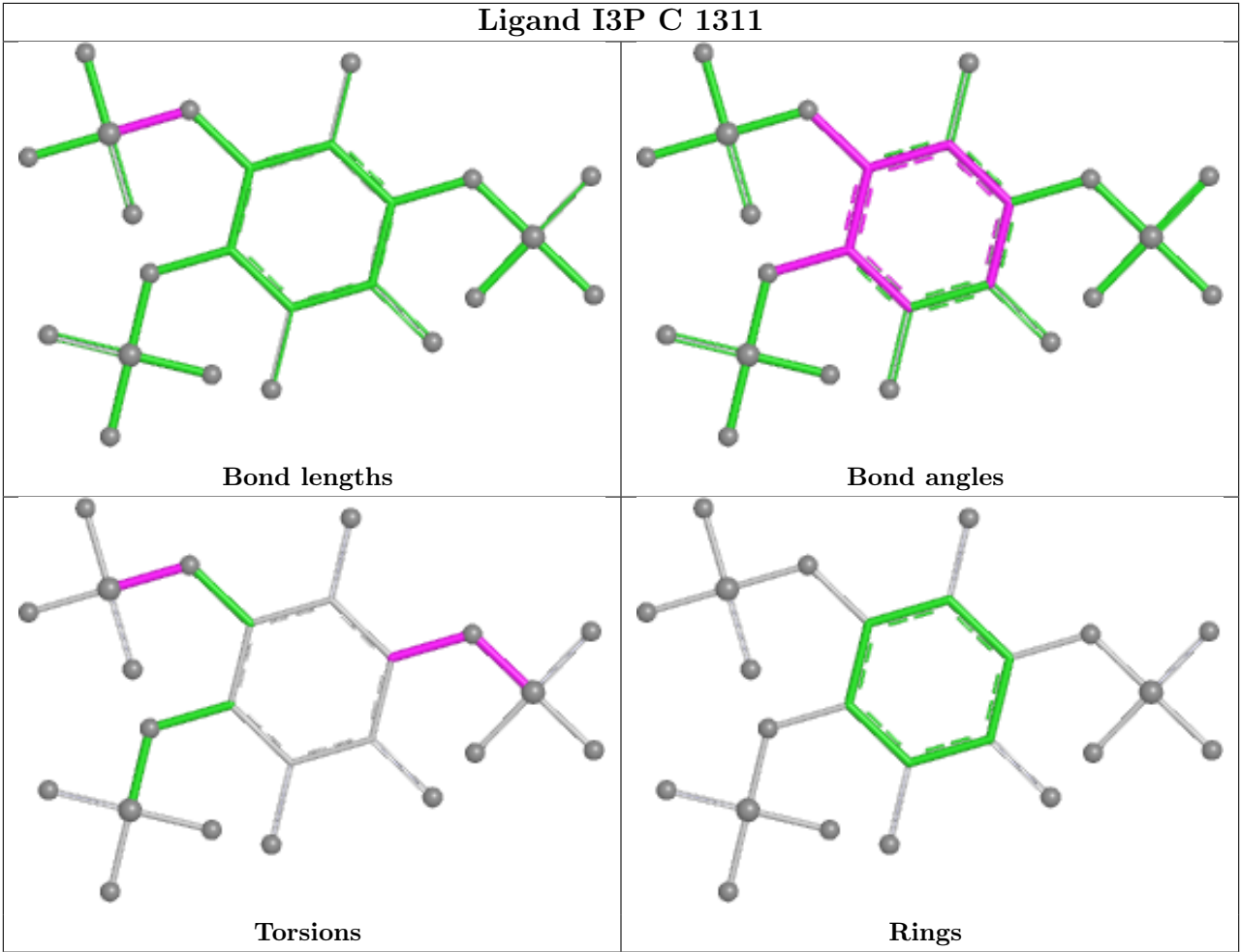
No monomer is involved in short contacts.

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 than 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.







5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	559/572 (97%)	-0.13	1 (0%) 91 84	74, 109, 161, 198	0
1	B	558/572 (97%)	0.16	7 (1%) 75 55	92, 151, 226, 251	0
1	C	563/572 (98%)	0.03	4 (0%) 84 69	72, 113, 169, 215	0
1	D	557/572 (97%)	0.15	15 (2%) 56 35	99, 153, 222, 243	0
All	All	2237/2288 (97%)	0.05	27 (1%) 76 57	72, 126, 211, 251	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	833	LEU	4.6
1	D	965	LEU	4.3
1	D	1174	LEU	3.5
1	D	826	PHE	3.2
1	D	760	ALA	2.8
1	B	1229	LEU	2.8
1	B	1216	LEU	2.8
1	D	880	LEU	2.6
1	D	1201	TRP	2.5
1	C	965	LEU	2.5
1	D	1057	VAL	2.4
1	B	913	LEU	2.4
1	B	1065	LEU	2.3
1	B	943	GLY	2.3
1	D	1209	PRO	2.2
1	D	950	THR	2.2
1	D	832	ILE	2.2
1	A	965	LEU	2.2
1	C	723	PHE	2.2
1	B	1277	MET	2.1
1	D	824	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	836	GLU	2.1
1	C	913	LEU	2.1
1	C	817	LEU	2.1
1	D	763	LEU	2.0
1	D	913	LEU	2.0
1	B	1209	PRO	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	0.21	0.15	145,156,159,159	0
2	MAN	F	5	11/12	0.25	0.14	138,140,143,146	0
2	MAN	H	6	11/12	0.37	0.13	142,147,150,150	0
2	MAN	H	5	11/12	0.44	0.12	124,135,139,144	0
2	MAN	F	6	11/12	0.47	0.14	148,149,152,153	0
2	MAN	E	8	11/12	0.48	0.13	152,156,159,159	0
2	MAN	H	10	11/12	0.49	0.12	133,136,138,138	0
2	MAN	H	7	11/12	0.55	0.11	141,143,144,146	0
2	MAN	G	5	11/12	0.55	0.10	133,137,141,142	0
2	MAN	E	5	11/12	0.56	0.11	147,150,152,153	0
2	MAN	G	6	11/12	0.56	0.10	138,145,149,150	0
2	MAN	H	8	11/12	0.58	0.10	132,138,140,141	0
2	NAG	F	1	14/15	0.60	0.14	109,115,121,122	0
2	MAN	G	8	11/12	0.64	0.10	135,143,146,150	0
2	MAN	H	4	11/12	0.65	0.11	126,129,134,138	0
2	MAN	F	7	11/12	0.65	0.11	136,140,146,152	0
2	MAN	G	7	11/12	0.66	0.10	126,129,131,135	0
2	MAN	F	10	11/12	0.66	0.12	120,125,127,129	0
2	MAN	E	10	11/12	0.68	0.11	138,146,150,150	0
2	MAN	F	9	11/12	0.71	0.10	124,126,128,128	0
2	MAN	F	4	11/12	0.72	0.10	131,134,140,140	0
2	MAN	H	9	11/12	0.75	0.09	121,123,125,131	0

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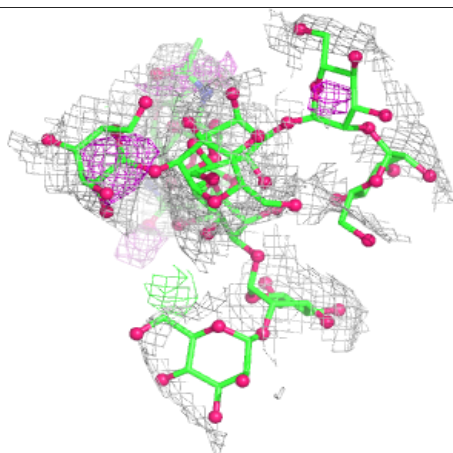
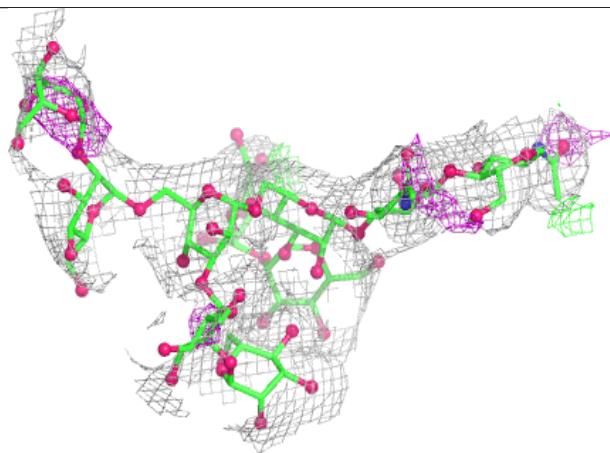
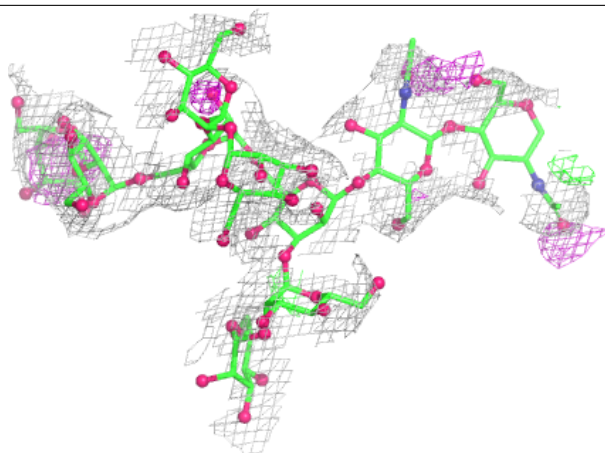
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	MAN	E	6	11/12	0.77	0.09	149,154,160,162	0
2	NAG	F	2	14/15	0.77	0.13	122,122,123,124	0
2	NAG	H	1	14/15	0.78	0.12	111,113,119,121	0
2	MAN	G	10	11/12	0.80	0.11	127,133,140,143	0
2	MAN	E	7	11/12	0.81	0.09	131,134,143,151	0
2	BMA	H	3	11/12	0.83	0.08	113,118,120,122	0
2	NAG	E	1	14/15	0.84	0.11	94,96,99,99	0
2	NAG	H	2	14/15	0.85	0.12	108,117,123,127	0
2	MAN	E	9	11/12	0.85	0.07	120,127,130,134	0
2	MAN	G	4	11/12	0.85	0.09	106,110,120,123	0
2	MAN	G	9	11/12	0.86	0.07	107,110,114,123	0
2	NAG	G	1	14/15	0.90	0.09	78,83,89,94	0
2	BMA	F	3	11/12	0.90	0.09	118,122,125,128	0
2	BMA	E	3	11/12	0.92	0.07	103,105,113,115	0
2	MAN	E	4	11/12	0.92	0.06	119,122,130,136	0
2	BMA	G	3	11/12	0.95	0.06	86,89,99,100	0
2	NAG	E	2	14/15	0.95	0.09	92,95,101,102	0
2	NAG	G	2	14/15	0.96	0.07	75,77,80,82	0

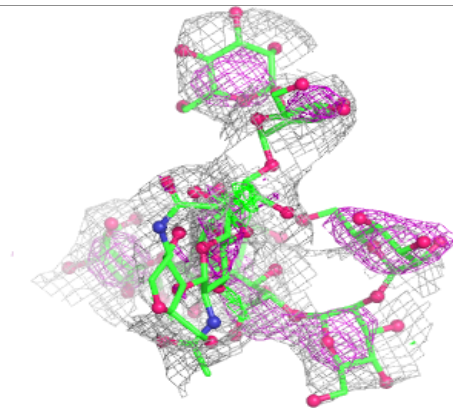
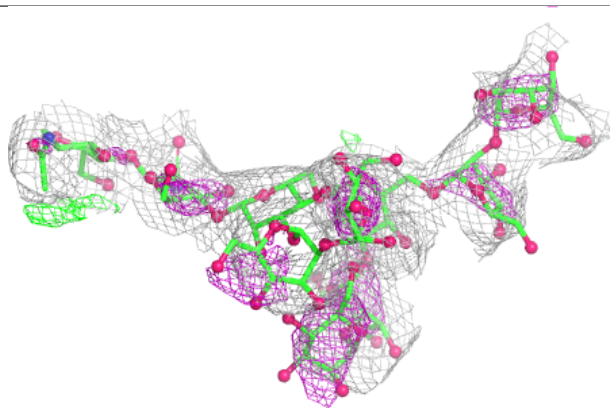
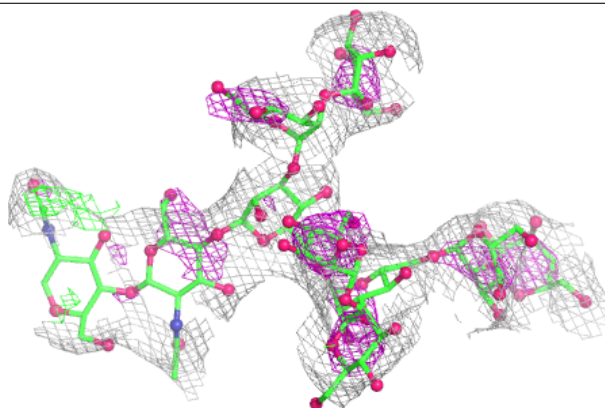
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

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)

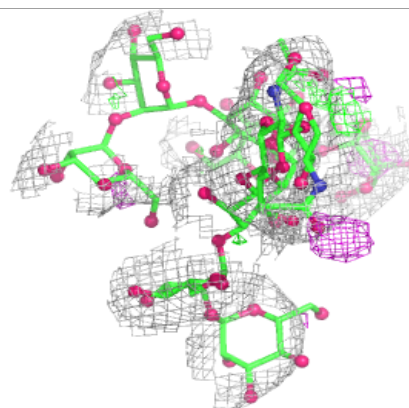
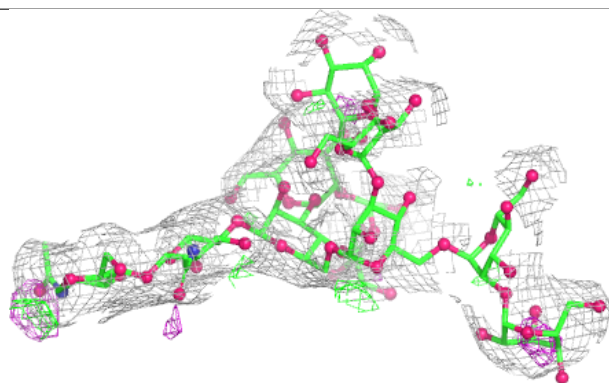
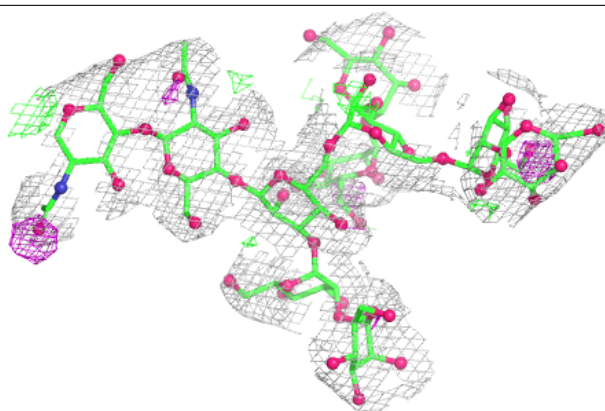
**Electron density around Chain F:**

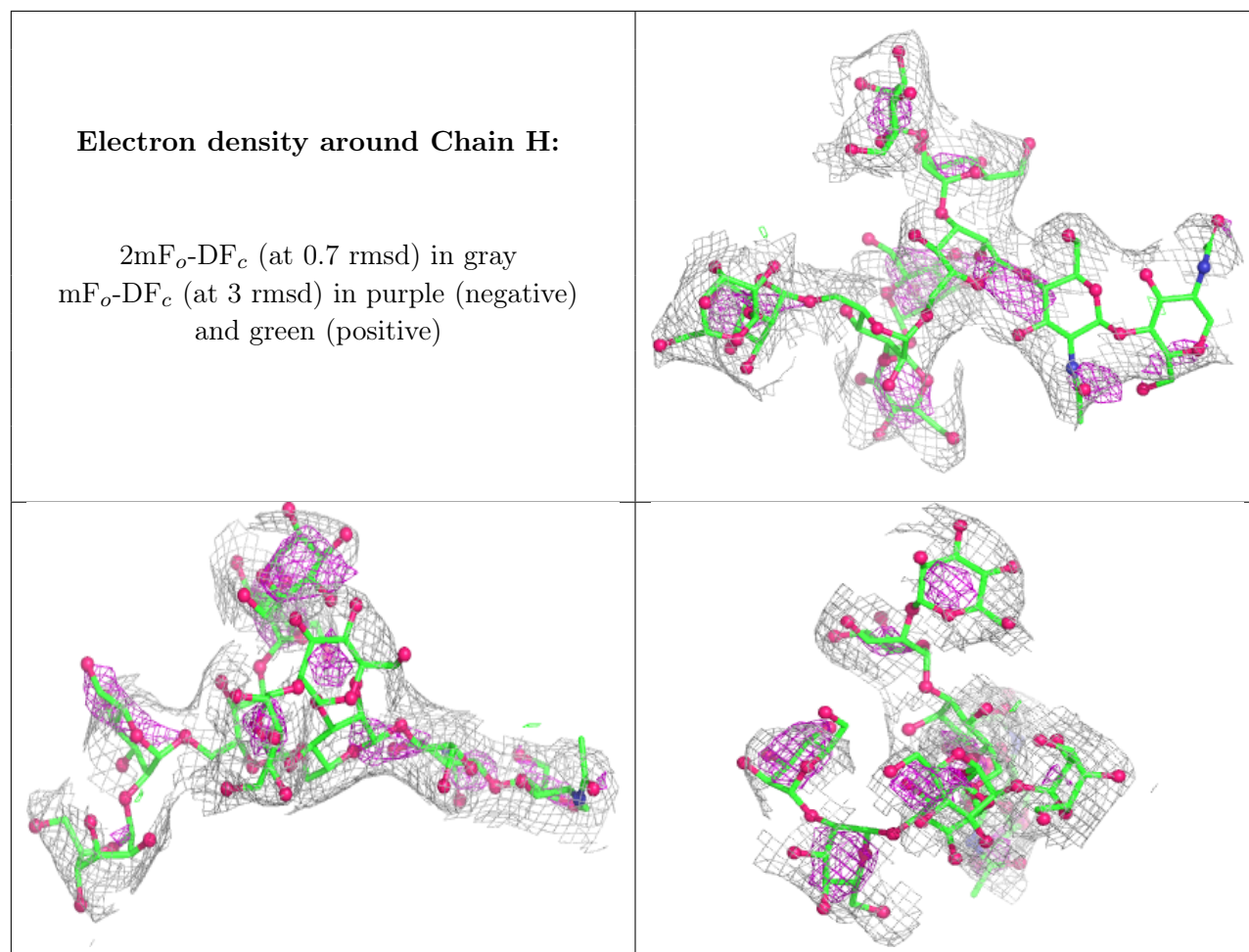
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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)





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.

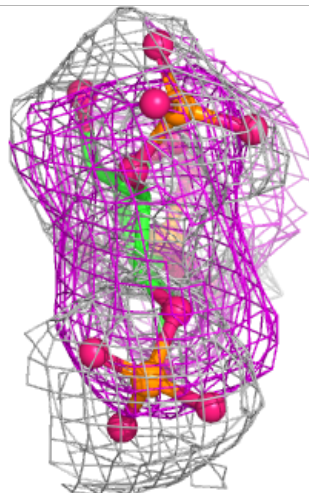
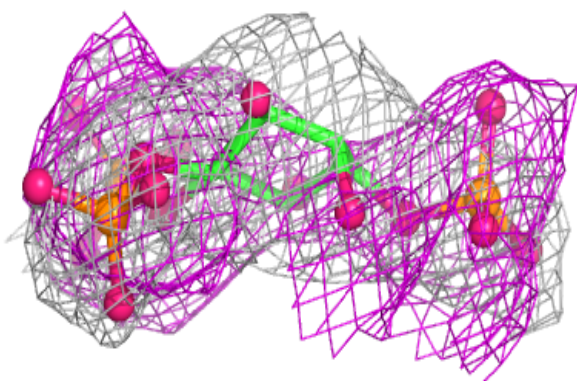
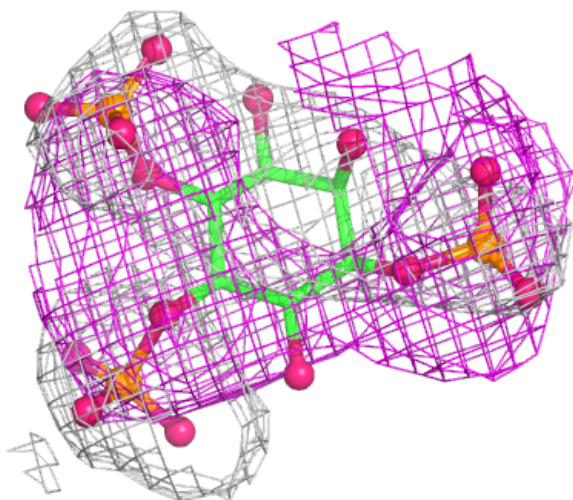
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	D	1310	14/15	0.18	0.17	100,104,106,106	0
3	NAG	B	1310	14/15	0.32	0.17	82,86,86,87	0
4	I3P	B	1311	24/24	0.44	0.11	204,209,217,218	0
4	I3P	C	1311	24/24	0.56	0.13	166,178,188,190	0
4	I3P	D	1311	24/24	0.59	0.12	190,201,218,219	0
3	NAG	A	1310	14/15	0.63	0.12	115,121,126,130	0
4	I3P	A	1311	24/24	0.65	0.12	185,204,218,219	0
3	NAG	C	1310	14/15	0.80	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.

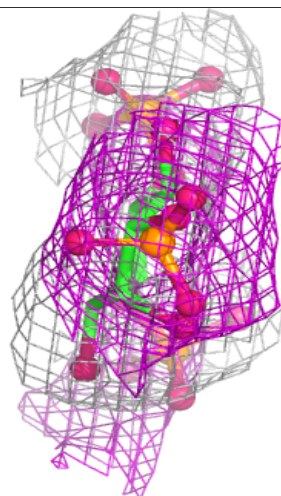
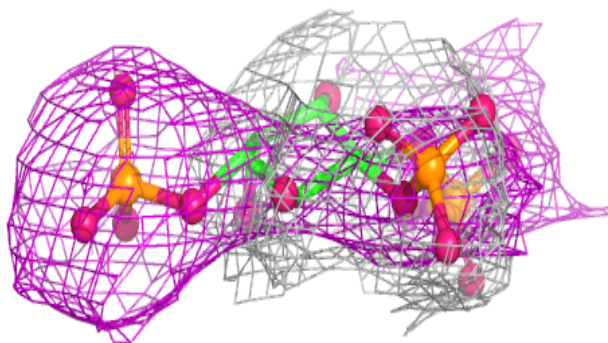
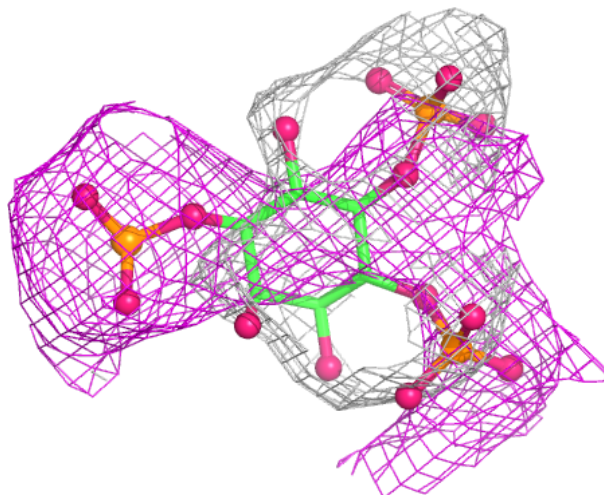
Electron density around I3P C 1311:

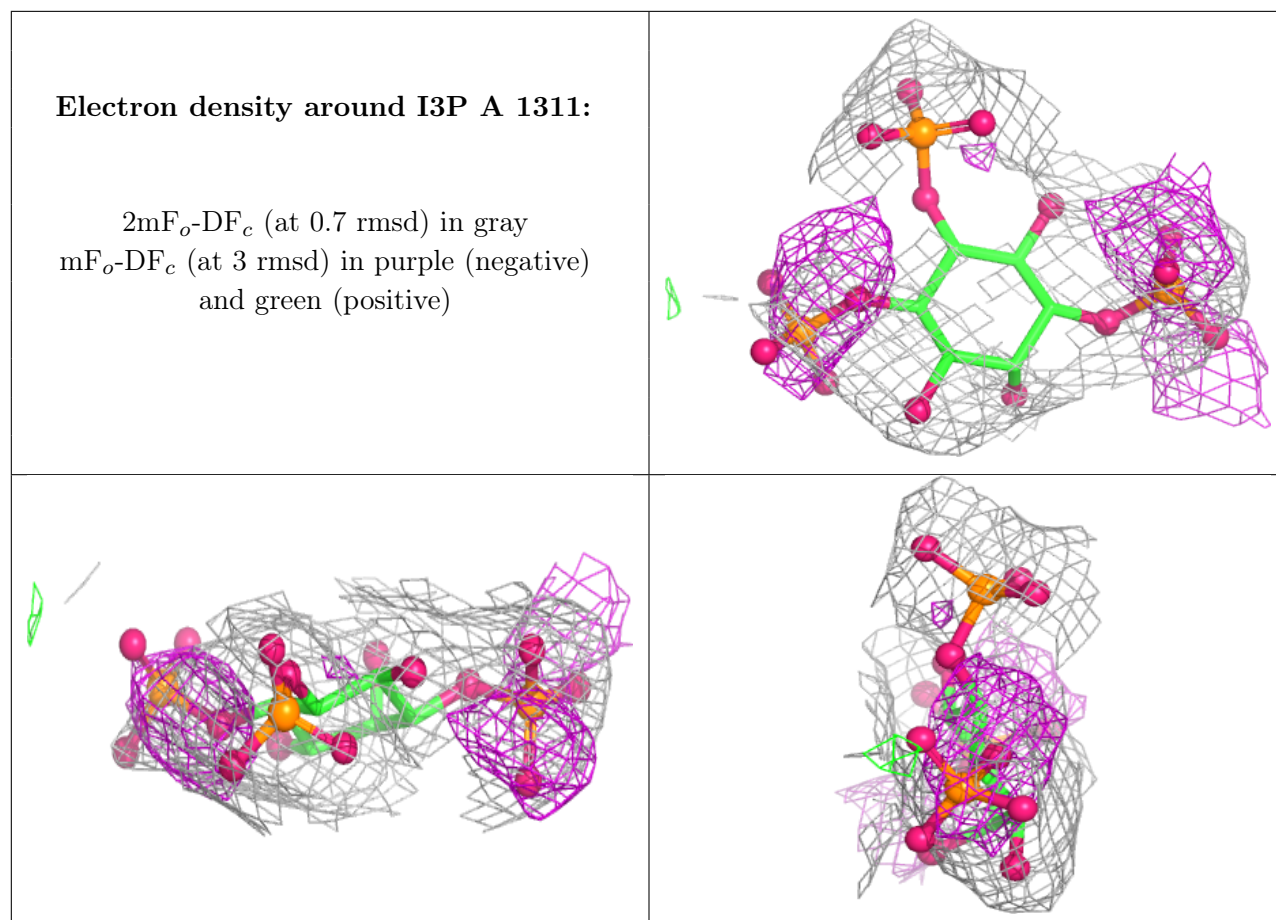
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around I3P 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)





6.5 Other polymers ⓘ

There are no such residues in this entry.