



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 12:08 PM UTC

PDB ID : 3VJL / pdb_00003vjL
Title : Crystal structure of human deipeptidyl peptidase IV (DPP-4) in complex with a prolylthiazolidine inhibitor #2
Authors : Akahoshi, F.; Kishida, H.; Miyaguchi, I.; Yoshida, T.; Ishii, S.
Deposited on : 2011-10-24
Resolution : 2.39 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

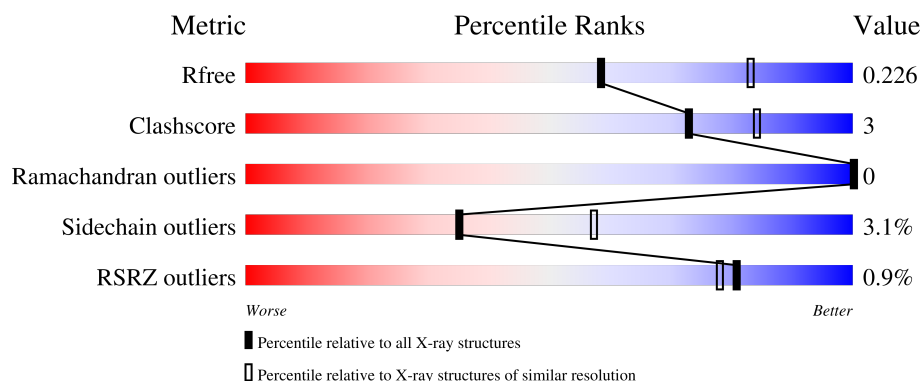
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.39 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.



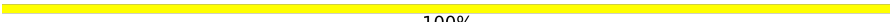
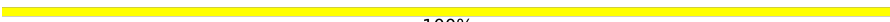
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	4912 (2.40-2.40)
Clashscore	190562	5391 (2.40-2.40)
Ramachandran outliers	187476	5320 (2.40-2.40)
Sidechain outliers	187428	5321 (2.40-2.40)
RSRZ outliers	180081	4916 (2.40-2.40)

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	740	88% 10% ..
1	B	740	89% 9% ..
2	C	2	50% 50%
2	D	2	100%
2	E	2	50% 50%

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

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 13126 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 Dipeptidyl peptidase 4.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	729	Total	C	N	O	S	0	0	0
			5971	3831	983	1131	26			
1	B	729	Total	C	N	O	S	0	0	0
			5971	3831	983	1131	26			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	767	HIS	-	expression tag	UNP P27487
A	768	HIS	-	expression tag	UNP P27487
A	769	HIS	-	expression tag	UNP P27487
A	770	HIS	-	expression tag	UNP P27487
A	771	HIS	-	expression tag	UNP P27487
A	772	HIS	-	expression tag	UNP P27487
B	767	HIS	-	expression tag	UNP P27487
B	768	HIS	-	expression tag	UNP P27487
B	769	HIS	-	expression tag	UNP P27487
B	770	HIS	-	expression tag	UNP P27487
B	771	HIS	-	expression tag	UNP P27487
B	772	HIS	-	expression tag	UNP P27487

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



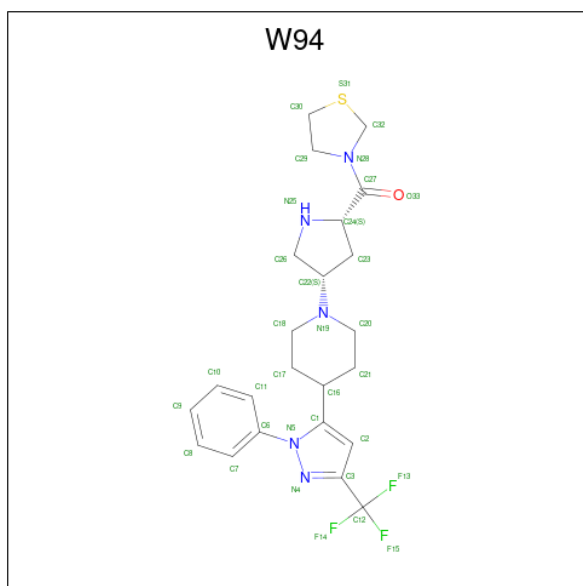
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	C	2	Total	C	N	O	0	0	0
			28	16	2	10			

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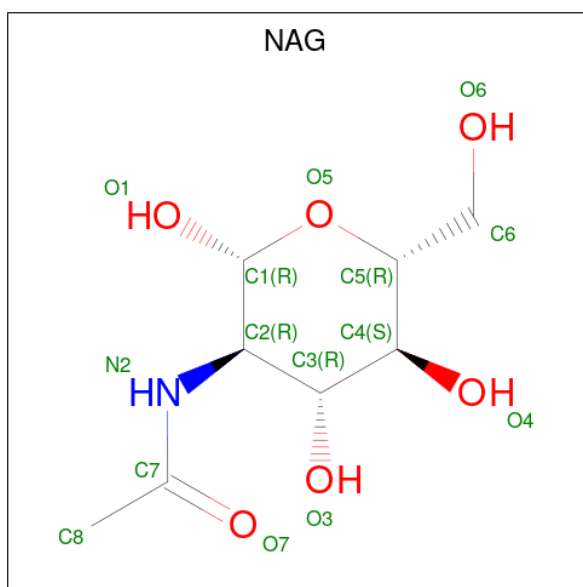
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	E	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	F	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	G	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is [(2S,4S)-4-{4-[1-phenyl-3-(trifluoromethyl)-1H-pyrazol-5-yl]piperidin-1-yl}pyrrolidin-2-yl](1,3-thiazolidin-3-yl)methanone (CCD ID: W94) (formula: C₂₃H₂₈F₃N₅OS).



Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	S	0	0
			33	23	3	5	1	1		
3	B	1	Total	C	F	N	O	S	0	0
			33	23	3	5	1	1		

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).



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

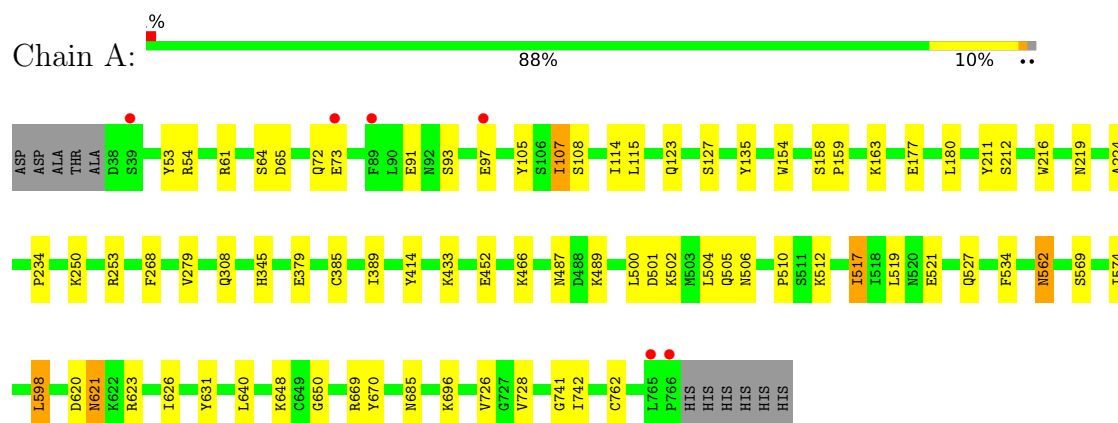
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	420	Total	O	0	0
			420	420		
5	B	474	Total	O	0	0
			474	474		

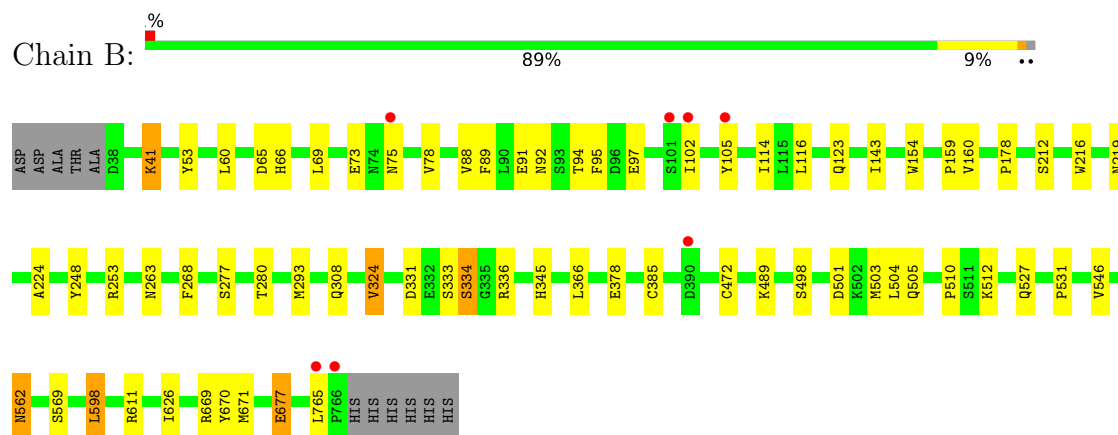
3 Residue-property plots [i](#)

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: Dipeptidyl peptidase 4



- Molecule 1: Dipeptidyl peptidase 4



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



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

Chain D:  100%

MAG1
MAG2

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

Chain E:  50% 50%

MAG1
MAG2

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

Chain F:  100%

MAG1
MAG2

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

Chain G:  100%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	117.81Å 125.83Å 137.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.39 30.00 – 2.39	Depositor EDS
% Data completeness (in resolution range)	99.1 (30.00-2.39) 99.0 (30.00-2.39)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	10.44 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0110	Depositor
R, R_{free}	0.184 , 0.230 0.181 , 0.226	Depositor DCC
R_{free} test set	4025 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	27.8	Xtriage
Anisotropy	0.039	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 36.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.51$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	13126	wwPDB-VP
Average B, all atoms (Å ²)	25.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.77% 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, W94

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	0.58	1/6143 (0.0%)	0.77	0/8355
1	B	0.60	2/6143 (0.0%)	0.78	0/8355
All	All	0.59	3/12286 (0.0%)	0.77	0/16710

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	531	PRO	CA-C	5.71	1.55	1.51
1	B	123	GLN	CA-C	5.30	1.55	1.52
1	A	123	GLN	CA-C	5.06	1.54	1.52

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5971	0	5684	40	0
1	B	5971	0	5683	37	0
2	C	28	0	25	0	0
2	D	28	0	25	0	0
2	E	28	0	25	0	0
2	F	28	0	25	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	G	28	0	25	0	0
3	A	33	0	28	0	0
3	B	33	0	28	0	0
4	A	42	0	39	1	0
4	B	42	0	39	0	0
5	A	420	0	0	3	0
5	B	474	0	0	5	0
All	All	13126	0	11626	76	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 3.

All (76) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:253:ARG:HH21	1:B:253:ARG:HH21	1.22	0.87
1:A:640:LEU:HD11	1:A:650:GLY:HA3	1.62	0.81
4:A:1229:NAG:H81	5:A:2134:HOH:O	1.81	0.80
1:A:219:ASN:H	1:A:308:GLN:HE22	1.38	0.71
1:B:345:HIS:HD2	5:B:2628:HOH:O	1.74	0.71
1:B:219:ASN:H	1:B:308:GLN:HE22	1.46	0.62
1:A:105:TYR:HB2	1:A:114:ILE:HD11	1.83	0.59
1:B:334:SER:HB3	1:B:336:ARG:H	1.69	0.57
1:B:331:ASP:OD1	1:B:334:SER:HB2	2.05	0.56
1:B:546:VAL:HG21	1:B:626:ILE:HD11	1.88	0.56
1:B:512:LYS:HE3	1:B:527:GLN:OE1	2.06	0.56
1:A:510:PRO:HD3	1:A:569:SER:HB2	1.89	0.54
1:A:620:ASP:OD2	1:A:623:ARG:HD3	2.08	0.54
1:B:489:LYS:HE2	5:B:2597:HOH:O	2.07	0.54
1:A:114:ILE:HG23	1:A:135:TYR:HB3	1.89	0.54
1:B:510:PRO:HD3	1:B:569:SER:HB2	1.89	0.54
1:A:219:ASN:N	1:A:308:GLN:HE22	2.04	0.53
1:B:677:GLU:H	1:B:677:GLU:CD	2.18	0.51
1:A:72:GLN:HG3	1:A:73:GLU:HG3	1.92	0.50
1:B:75:ASN:ND2	1:B:92:ASN:H	2.09	0.50
1:A:648:LYS:HE2	1:A:762:CYS:O	2.12	0.50
1:A:696:LYS:HG3	1:A:728:VAL:HG22	1.93	0.50
1:B:41:LYS:HE3	1:B:53:TYR:OH	2.12	0.50
1:A:621:ASN:HD22	1:A:621:ASN:H	1.57	0.50
1:B:562:ASN:C	1:B:562:ASN:HD22	2.19	0.49
1:B:78:VAL:HG23	1:B:89:PHE:HB2	1.95	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:219:ASN:N	1:B:308:GLN:HE22	2.11	0.48
1:B:277:SER:HB3	1:B:280:THR:HG22	1.96	0.47
1:A:234:PRO:HB2	1:B:248:TYR:CZ	2.50	0.47
1:A:726:VAL:HG23	1:A:728:VAL:HG23	1.96	0.47
1:B:512:LYS:HE3	1:B:527:GLN:CD	2.39	0.47
1:B:501:ASP:O	1:B:505:GLN:HG2	2.13	0.47
1:A:626:ILE:O	1:A:650:GLY:HA2	2.15	0.47
1:B:69:LEU:HD23	1:B:78:VAL:HG22	1.97	0.47
1:A:127:SER:HB3	1:A:211:TYR:CD1	2.50	0.47
1:B:105:TYR:HB3	1:B:114:ILE:HD11	1.96	0.46
1:B:60:LEU:HD12	1:B:60:LEU:C	2.40	0.46
1:B:75:ASN:HD22	1:B:91:GLU:HA	1.80	0.46
1:A:534:PHE:HE1	1:A:574:ILE:HD11	1.82	0.45
1:A:741:GLY:O	1:A:742:ILE:C	2.60	0.45
1:B:611:ARG:NH2	5:B:2664:HOH:O	2.50	0.45
1:A:107:ILE:HG22	1:A:108:SER:O	2.17	0.45
1:B:95:PHE:CE2	1:B:116:LEU:HD11	2.53	0.44
1:A:65:ASP:OD2	1:A:466:LYS:HB2	2.18	0.44
1:A:224:ALA:HB1	1:A:268:PHE:CZ	2.52	0.44
1:B:345:HIS:CD2	5:B:2628:HOH:O	2.59	0.44
1:A:64:SER:HB2	5:A:2385:HOH:O	2.18	0.44
1:B:159:PRO:HD3	1:B:216:TRP:CB	2.47	0.43
1:A:414:TYR:CE1	1:A:433:LYS:HE2	2.53	0.43
1:A:501:ASP:O	1:A:505:GLN:HG2	2.19	0.43
1:A:154:TRP:CE2	1:A:212:SER:HB3	2.54	0.43
1:A:562:ASN:C	1:A:562:ASN:HD22	2.25	0.43
1:A:91:GLU:C	1:A:93:SER:H	2.26	0.42
1:A:219:ASN:OD1	1:A:219:ASN:C	2.62	0.42
1:A:345:HIS:HE1	1:A:389:ILE:O	2.02	0.42
1:B:598:LEU:HD22	1:B:671:MET:HG2	2.01	0.42
1:A:669:ARG:HD2	1:A:670:TYR:CZ	2.54	0.42
1:B:143:ILE:HD13	1:B:178:PRO:HB2	2.02	0.42
1:A:53:TYR:HB3	1:A:500:LEU:HD11	2.02	0.42
1:B:41:LYS:HG2	5:B:2813:HOH:O	2.20	0.41
1:B:293:MET:HE3	1:B:324:VAL:HG13	2.01	0.41
1:A:114:ILE:CG2	1:A:135:TYR:HB3	2.49	0.41
1:B:669:ARG:HD2	1:B:670:TYR:CZ	2.56	0.41
1:A:345:HIS:HD2	5:A:2077:HOH:O	2.04	0.41
1:B:154:TRP:CE2	1:B:212:SER:HB2	2.56	0.41
1:B:224:ALA:HB1	1:B:268:PHE:CZ	2.56	0.41
1:A:598:LEU:HG	1:A:631:TYR:OH	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:105:TYR:HA	1:A:115:LEU:O	2.21	0.41
1:A:512:LYS:HE3	1:A:527:GLN:OE1	2.20	0.41
1:A:158:SER:OG	1:A:163:LYS:HB2	2.21	0.40
1:A:177:GLU:HG3	1:A:180:LEU:HD12	2.02	0.40
1:A:517:ILE:HD12	1:A:519:LEU:HG	2.03	0.40
1:B:159:PRO:HD3	1:B:216:TRP:HB3	2.01	0.40
1:B:88:VAL:HG11	1:B:91:GLU:HG2	2.03	0.40
1:A:159:PRO:HD3	1:A:216:TRP:HB3	2.03	0.40
1:B:65:ASP:HB3	1:B:66:HIS:CE1	2.57	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	727/740 (98%)	700 (96%)	27 (4%)	0	100	100
1	B	727/740 (98%)	703 (97%)	24 (3%)	0	100	100
All	All	1454/1480 (98%)	1403 (96%)	51 (4%)	0	100	100

There are no Ramachandran outliers to report.

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	654/663 (99%)	634 (97%)	20 (3%)	35	57
1	B	654/663 (99%)	633 (97%)	21 (3%)	34	56
All	All	1308/1326 (99%)	1267 (97%)	41 (3%)	35	57

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ARG
1	A	61	ARG
1	A	97	GLU
1	A	107	ILE
1	A	250	LYS
1	A	279	VAL
1	A	379	GLU
1	A	385	CYS
1	A	452	GLU
1	A	487	ASN
1	A	489	LYS
1	A	502	LYS
1	A	504	LEU
1	A	506	ASN
1	A	517	ILE
1	A	521	GLU
1	A	562	ASN
1	A	598	LEU
1	A	621	ASN
1	A	685	ASN
1	B	41	LYS
1	B	73	GLU
1	B	94	THR
1	B	97	GLU
1	B	102	ILE
1	B	160	VAL
1	B	263	ASN
1	B	324	VAL
1	B	333	SER
1	B	334	SER
1	B	366	LEU
1	B	378	GLU
1	B	385	CYS
1	B	472	CYS
1	B	498	SER

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Mol	Chain	Res	Type
1	B	503	MET
1	B	504	LEU
1	B	562	ASN
1	B	598	LEU
1	B	677	GLU
1	B	765	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (33) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	75	ASN
1	A	112	GLN
1	A	138	ASN
1	A	169	ASN
1	A	227	GLN
1	A	263	ASN
1	A	308	GLN
1	A	345	HIS
1	A	430	ASN
1	A	487	ASN
1	A	505	GLN
1	A	506	ASN
1	A	562	ASN
1	A	572	ASN
1	A	621	ASN
1	A	679	ASN
1	A	694	ASN
1	B	66	HIS
1	B	75	ASN
1	B	112	GLN
1	B	123	GLN
1	B	169	ASN
1	B	263	ASN
1	B	286	GLN
1	B	308	GLN
1	B	345	HIS
1	B	505	GLN
1	B	562	ASN
1	B	572	ASN
1	B	606	GLN
1	B	679	ASN

Continued on next page...

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Mol	Chain	Res	Type
1	B	694	ASN

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 ⓘ

10 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	C	1	2,1	14,14,15	0.58	0	17,19,21	0.98	0
2	NAG	C	2	2	14,14,15	0.51	0	17,19,21	1.26	1 (5%)
2	NAG	D	1	2,1	14,14,15	0.74	0	17,19,21	1.05	1 (5%)
2	NAG	D	2	2	14,14,15	0.50	0	17,19,21	0.97	1 (5%)
2	NAG	E	1	2,1	14,14,15	0.55	0	17,19,21	0.77	0
2	NAG	E	2	2	14,14,15	0.59	0	17,19,21	1.11	1 (5%)
2	NAG	F	1	2,1	14,14,15	0.61	0	17,19,21	1.33	1 (5%)
2	NAG	F	2	2	14,14,15	0.60	0	17,19,21	1.26	1 (5%)
2	NAG	G	1	2,1	14,14,15	0.53	0	17,19,21	0.98	1 (5%)
2	NAG	G	2	2	14,14,15	0.52	0	17,19,21	1.88	3 (17%)

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	C	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
2	NAG	D	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	1/6/23/26	0/1/1/1
2	NAG	E	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	0/6/23/26	0/1/1/1
2	NAG	F	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	1/6/23/26	0/1/1/1
2	NAG	G	1	2,1	-	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	1/6/23/26	0/1/1/1

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	G	2	NAG	C1-O5-C5	6.13	120.40	112.19
2	F	1	NAG	C1-O5-C5	4.74	118.53	112.19
2	C	2	NAG	C1-O5-C5	3.27	116.57	112.19
2	F	2	NAG	C1-O5-C5	3.18	116.44	112.19
2	D	1	NAG	O5-C1-C2	-3.00	106.65	111.29
2	G	2	NAG	C3-C4-C5	2.60	114.94	110.23
2	D	2	NAG	C1-O5-C5	2.57	115.63	112.19
2	G	2	NAG	C4-C3-C2	2.22	114.27	111.02
2	G	1	NAG	O5-C1-C2	-2.13	107.99	111.29
2	E	2	NAG	O5-C5-C6	2.12	111.79	107.66

There are no chirality outliers.

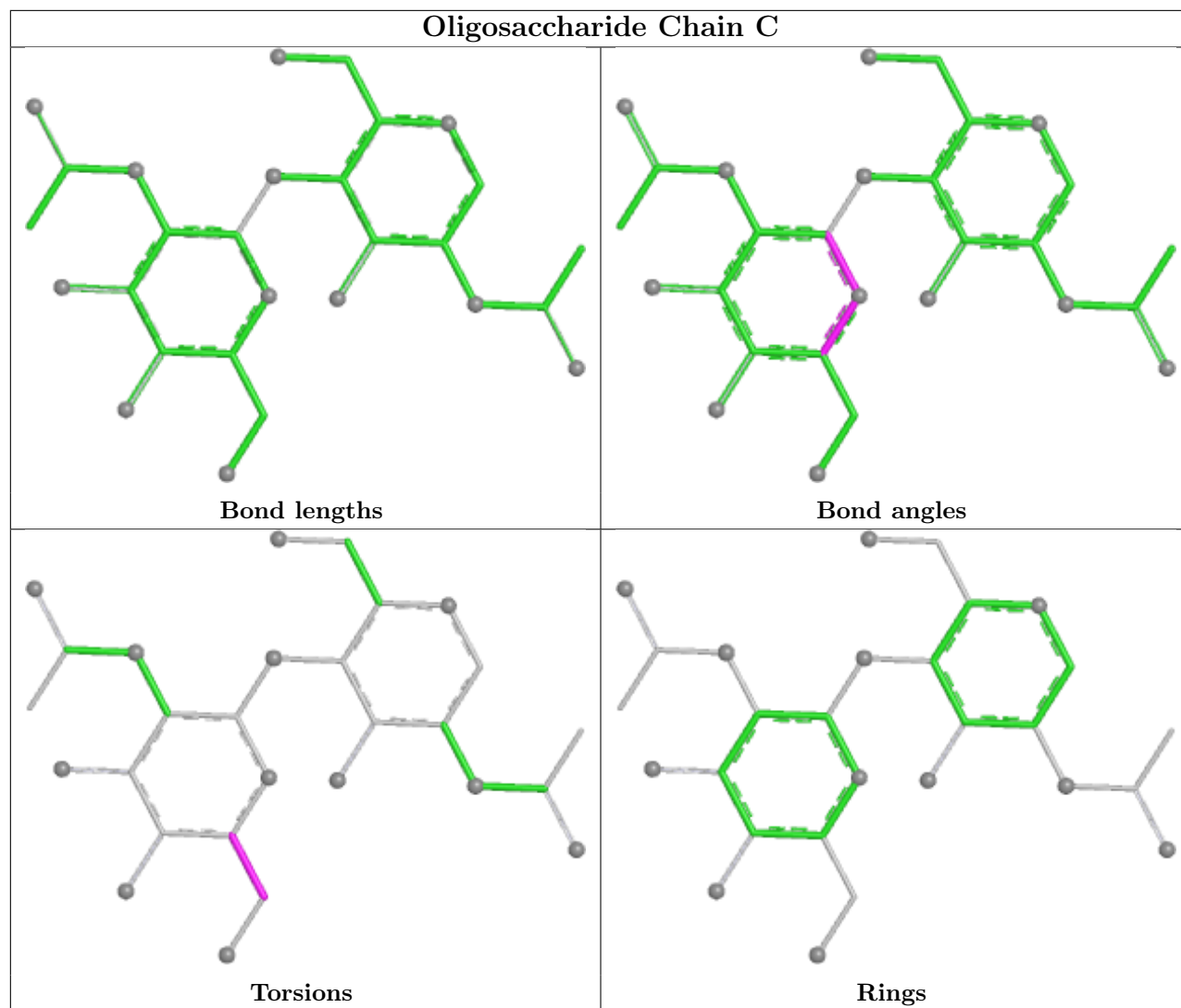
All (5) torsion outliers are listed below:

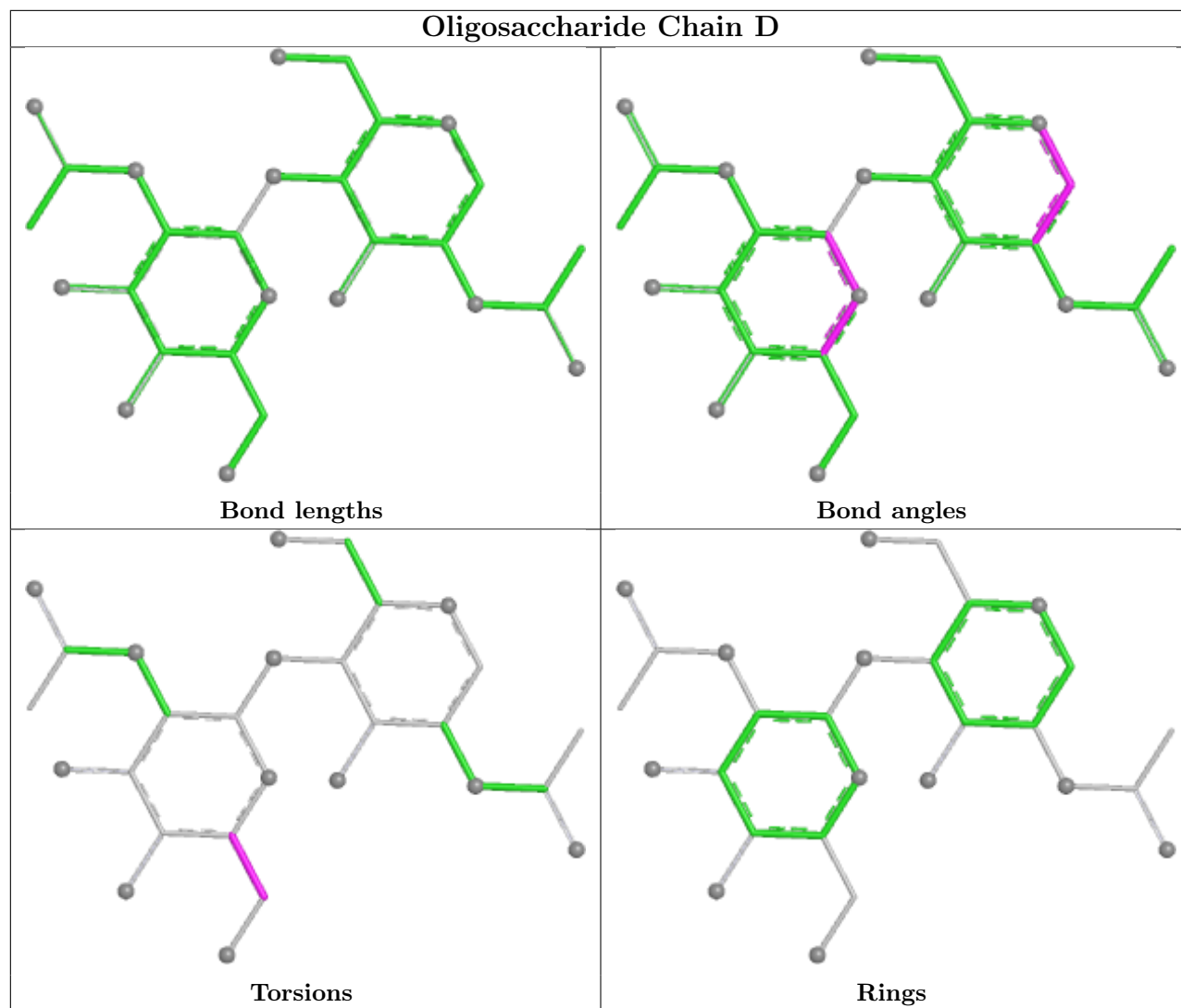
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C4-C5-C6-O6
2	C	2	NAG	O5-C5-C6-O6
2	D	2	NAG	C4-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6

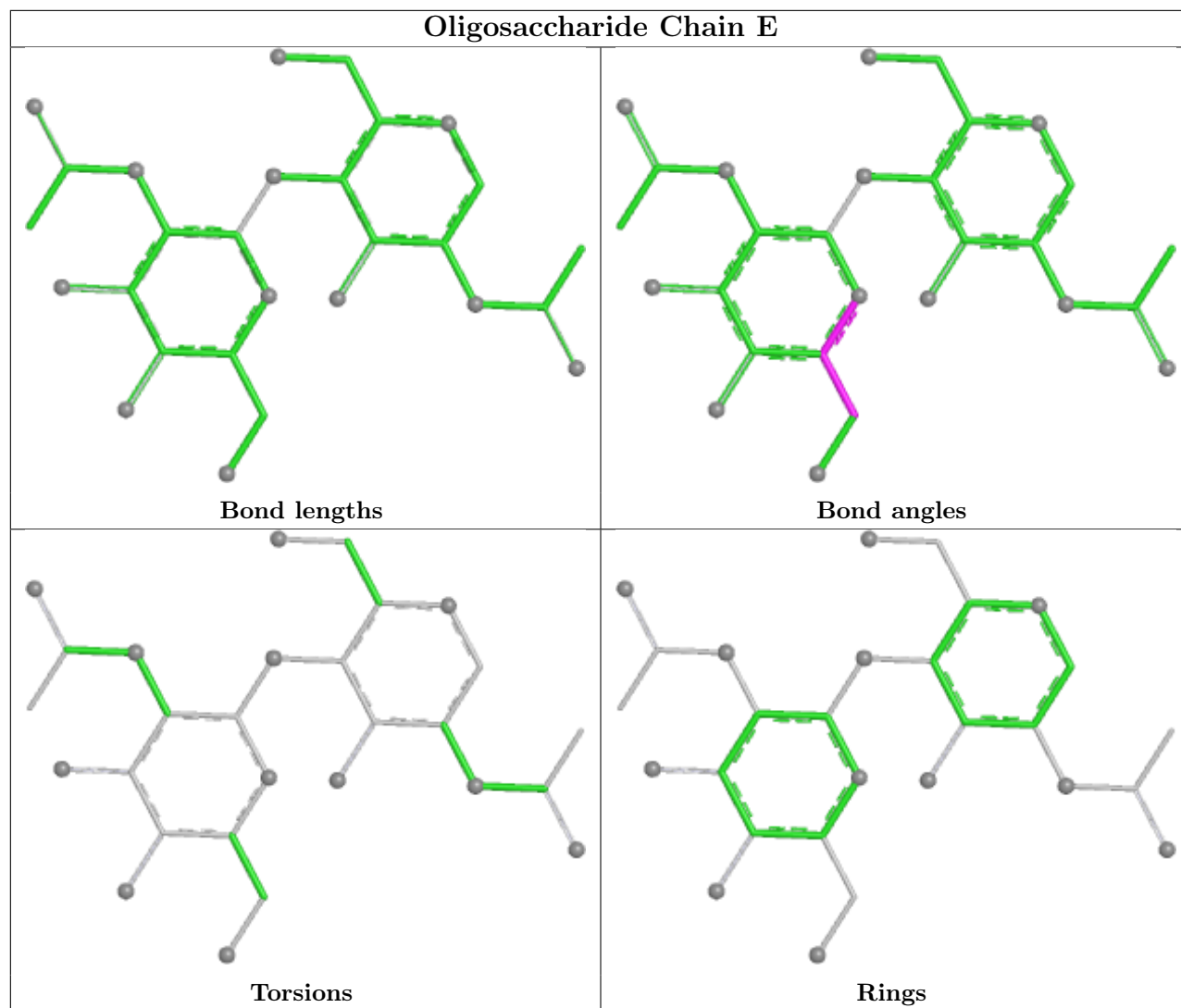
There are no ring outliers.

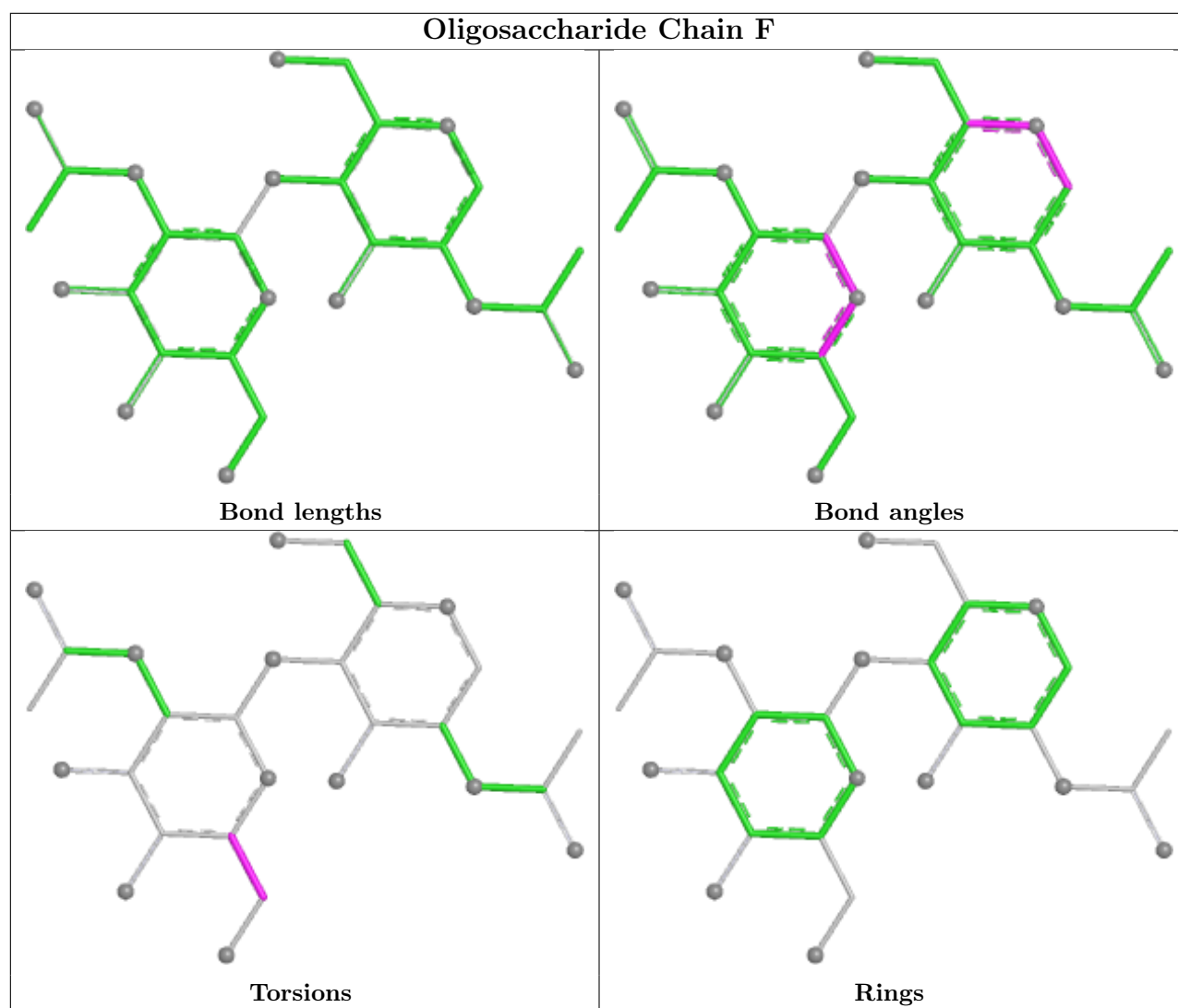
No monomer is involved in short contacts.

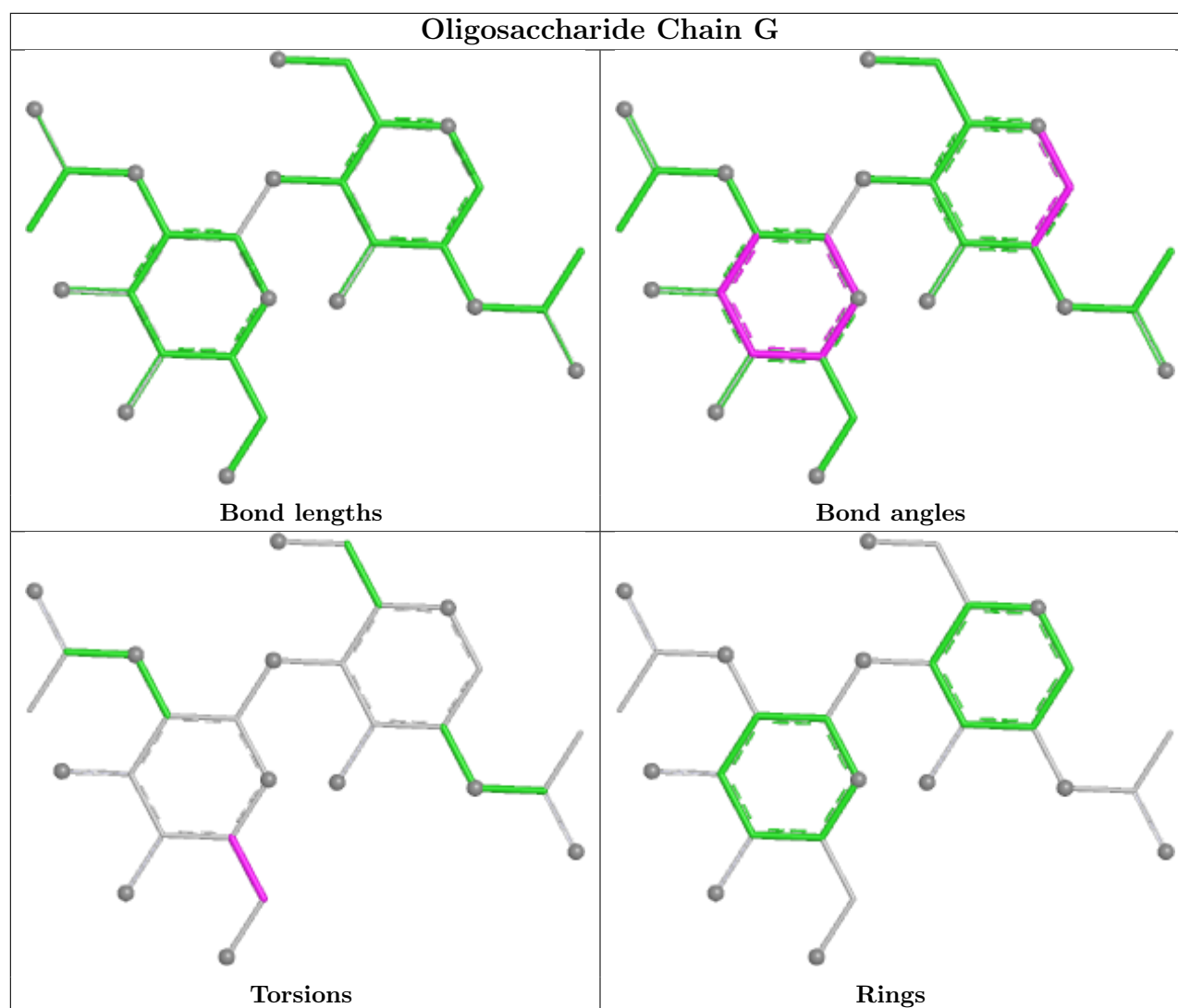
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.











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	NAG	B	1150	1	14,14,15	0.51	0	17,19,21	1.30	1 (5%)
3	W94	A	1000	-	36,37,37	1.09	1 (2%)	44,54,54	2.12	12 (27%)
4	NAG	A	1085	1	14,14,15	0.42	0	17,19,21	1.63	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	1229	1	14,14,15	0.53	0	17,19,21	1.59	3 (17%)
4	NAG	B	1085	1	14,14,15	0.51	0	17,19,21	1.61	1 (5%)
4	NAG	B	1281	1	14,14,15	0.61	0	17,19,21	1.50	3 (17%)
4	NAG	A	1321	1	14,14,15	0.58	0	17,19,21	1.68	1 (5%)
3	W94	B	1000	-	36,37,37	1.14	2 (5%)	44,54,54	2.03	10 (22%)

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	NAG	B	1150	1	-	0/6/23/26	0/1/1/1
3	W94	A	1000	-	-	0/26/52/52	0/5/5/5
4	NAG	A	1085	1	-	1/6/23/26	0/1/1/1
4	NAG	A	1229	1	-	3/6/23/26	0/1/1/1
4	NAG	B	1085	1	-	0/6/23/26	0/1/1/1
4	NAG	B	1281	1	-	2/6/23/26	0/1/1/1
4	NAG	A	1321	1	-	0/6/23/26	0/1/1/1
3	W94	B	1000	-	-	0/26/52/52	0/5/5/5

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1000	W94	C6-N5	-5.31	1.33	1.43
3	B	1000	W94	C6-N5	-4.91	1.34	1.43
3	B	1000	W94	C2-C1	2.39	1.41	1.37

All (32) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	1000	W94	C16-C1-N5	7.23	136.04	123.63
3	B	1000	W94	C16-C1-N5	6.97	135.59	123.63
4	A	1085	NAG	C1-O5-C5	6.11	120.37	112.19
4	B	1085	NAG	C1-O5-C5	5.71	119.84	112.19
4	A	1321	NAG	C1-O5-C5	5.61	119.70	112.19
3	A	1000	W94	C3-N4-N5	-4.98	100.09	104.27
3	B	1000	W94	C20-N19-C18	4.87	118.42	109.13
3	A	1000	W94	C1-C2-C3	-4.40	102.87	106.16
4	B	1150	NAG	C1-O5-C5	4.36	118.03	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1000	W94	C3-N4-N5	-4.27	100.69	104.27
3	A	1000	W94	C20-N19-C18	3.91	116.58	109.13
3	A	1000	W94	C2-C3-N4	3.68	116.21	111.86
4	A	1229	NAG	O5-C1-C2	-3.56	105.78	111.29
4	B	1281	NAG	C1-O5-C5	3.40	116.75	112.19
4	A	1229	NAG	C1-C2-N2	3.37	115.74	110.43
3	B	1000	W94	C7-C6-N5	3.19	124.53	119.72
3	B	1000	W94	C1-C2-C3	-3.19	103.78	106.16
4	A	1229	NAG	C1-O5-C5	3.09	116.32	112.19
3	B	1000	W94	C10-C11-C6	3.06	123.31	119.68
3	A	1000	W94	C8-C7-C6	2.97	123.21	119.68
3	A	1000	W94	C11-C6-N5	2.94	124.15	119.72
4	B	1281	NAG	O5-C5-C6	2.84	113.19	107.66
4	B	1281	NAG	C1-C2-N2	2.77	114.80	110.43
3	B	1000	W94	C2-C3-N4	2.73	115.08	111.86
3	A	1000	W94	F15-C12-C3	-2.52	107.70	112.86
3	A	1000	W94	F13-C12-C3	-2.49	107.75	112.86
3	B	1000	W94	C30-C29-N28	2.48	112.07	107.13
3	A	1000	W94	C12-C3-C2	-2.44	124.90	128.19
3	B	1000	W94	F13-C12-C3	-2.30	108.13	112.86
3	A	1000	W94	C30-C29-N28	2.19	111.49	107.13
3	B	1000	W94	C12-C3-C2	-2.16	125.28	128.19
3	A	1000	W94	C17-C18-N19	-2.04	108.29	111.42

There are no chirality outliers.

All (6) torsion outliers are listed below:

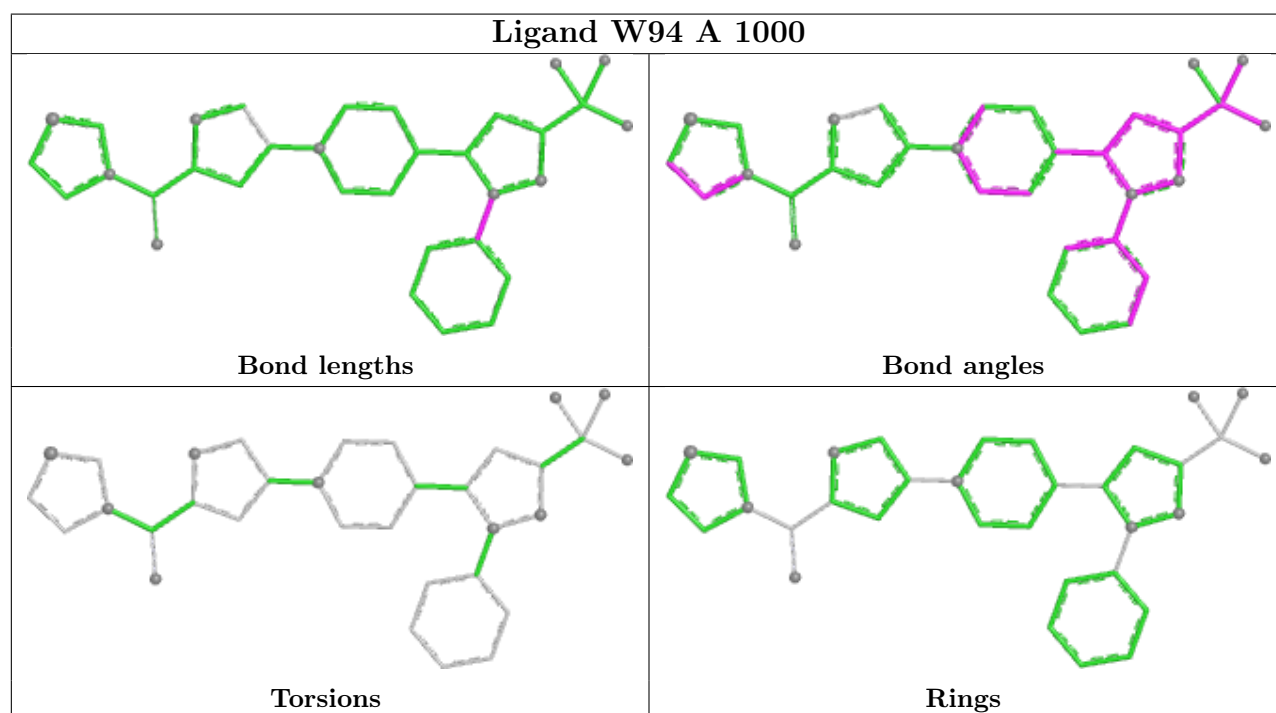
Mol	Chain	Res	Type	Atoms
4	B	1281	NAG	C4-C5-C6-O6
4	B	1281	NAG	O5-C5-C6-O6
4	A	1229	NAG	C8-C7-N2-C2
4	A	1229	NAG	O7-C7-N2-C2
4	A	1085	NAG	C4-C5-C6-O6
4	A	1229	NAG	C4-C5-C6-O6

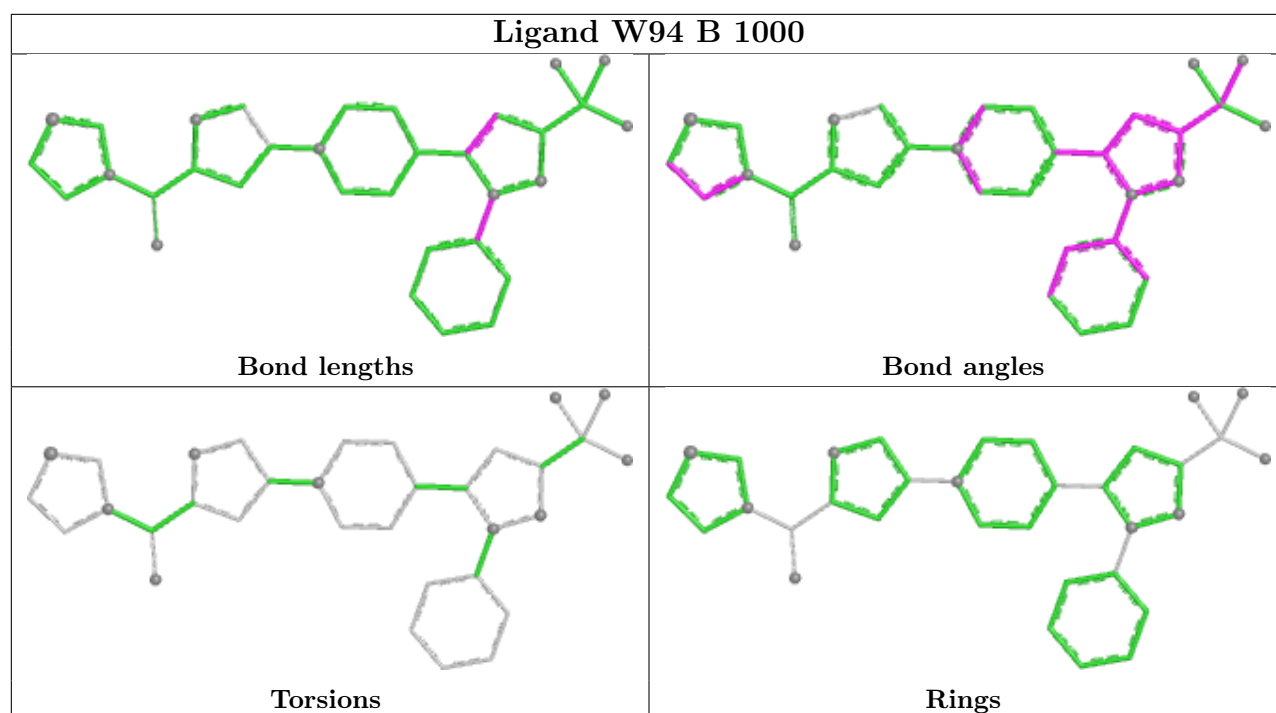
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	1229	NAG	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for 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 [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	729/740 (98%)	-0.24	6 (0%) 82 80	14, 23, 46, 66	0
1	B	729/740 (98%)	-0.32	7 (0%) 79 76	13, 22, 42, 56	0
All	All	1458/1480 (98%)	-0.28	13 (0%) 81 78	13, 23, 44, 66	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	766	PRO	6.6
1	B	766	PRO	5.6
1	A	765	LEU	5.3
1	B	105	TYR	2.8
1	B	765	LEU	2.6
1	A	89	PHE	2.4
1	A	73	GLU	2.3
1	A	39	SER	2.3
1	B	101	SER	2.3
1	B	390	ASP	2.2
1	B	102	ILE	2.2
1	B	75	ASN	2.1
1	A	97	GLU	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.

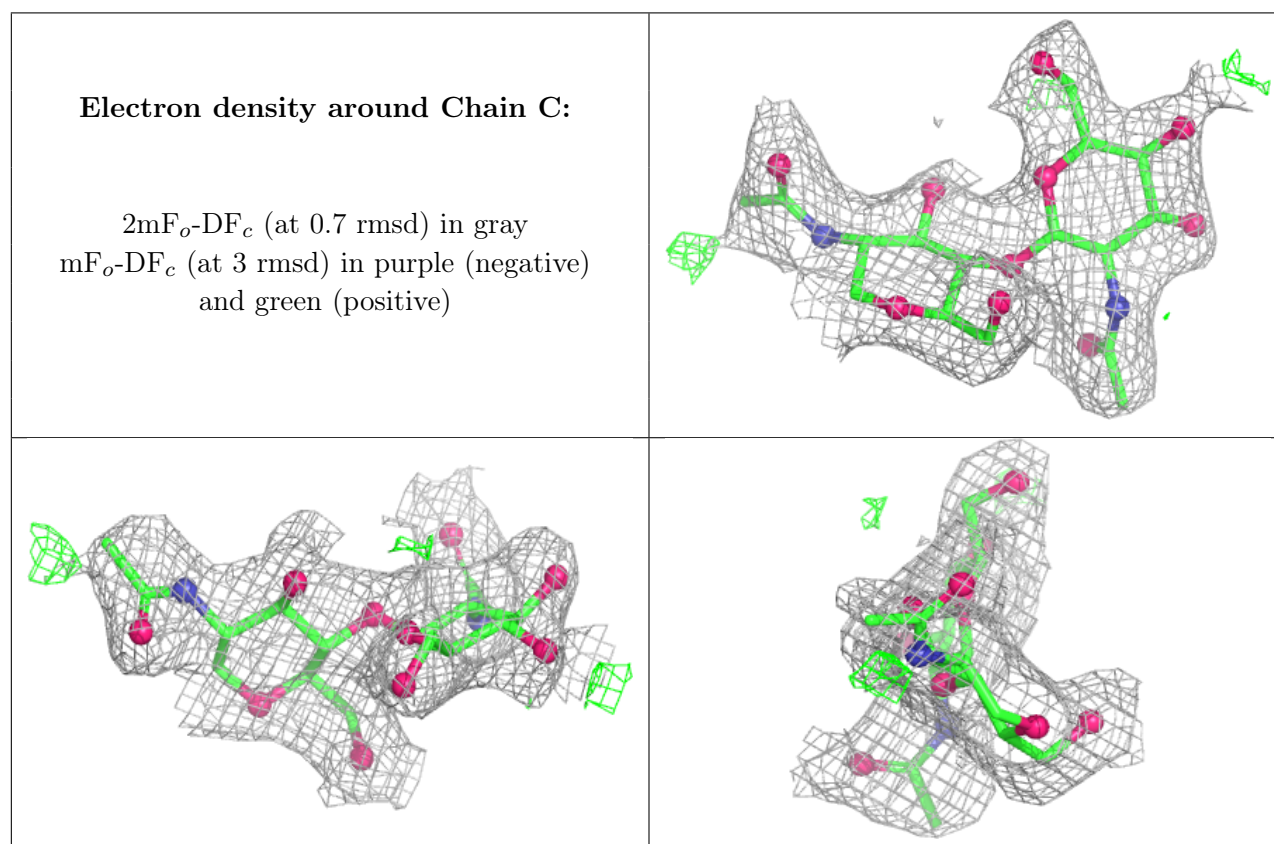
6.3 Carbohydrates [i](#)

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.

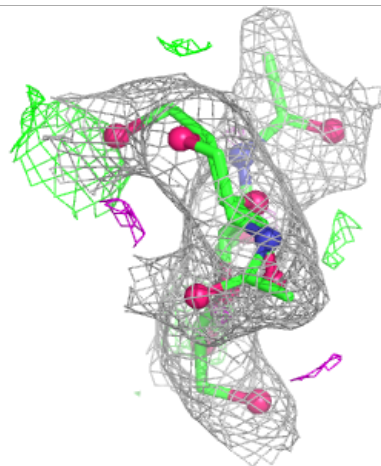
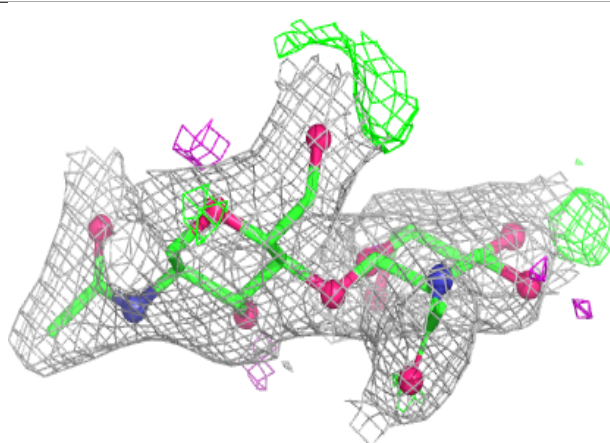
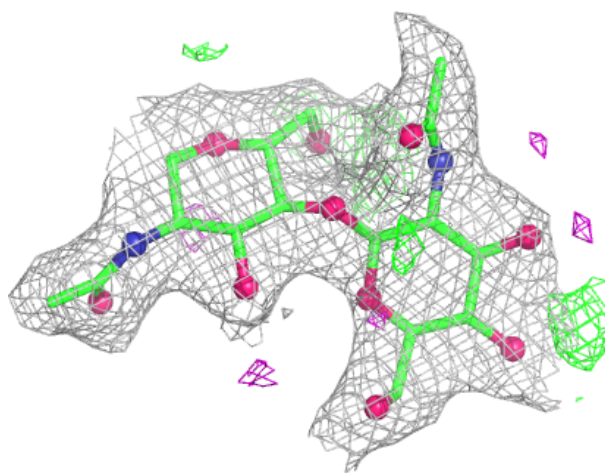
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	F	2	14/15	0.74	0.15	50,53,56,57	0
2	NAG	G	2	14/15	0.77	0.15	53,55,58,59	0
2	NAG	E	2	14/15	0.79	0.14	53,54,57,57	0
2	NAG	E	1	14/15	0.86	0.11	40,45,49,50	0
2	NAG	C	2	14/15	0.86	0.11	50,52,55,55	0
2	NAG	G	1	14/15	0.87	0.10	40,45,48,51	0
2	NAG	D	2	14/15	0.88	0.11	27,33,35,35	0
2	NAG	C	1	14/15	0.92	0.09	33,42,47,47	0
2	NAG	F	1	14/15	0.93	0.07	34,38,42,46	0
2	NAG	D	1	14/15	0.96	0.06	23,24,29,31	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.



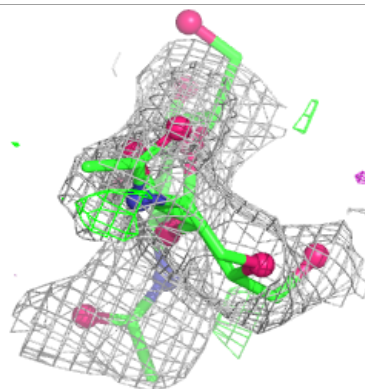
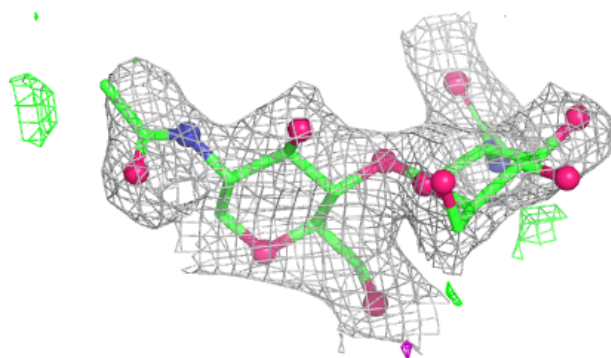
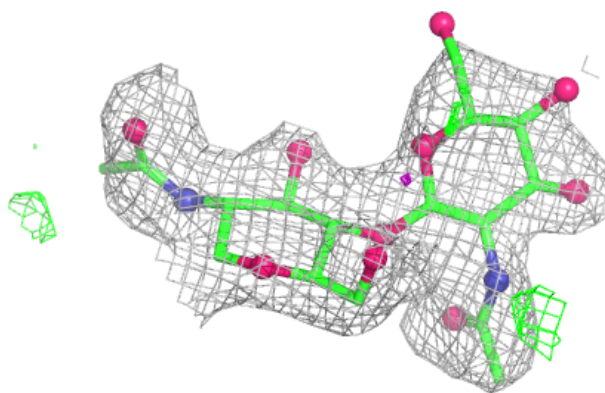
Electron density around Chain D:

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



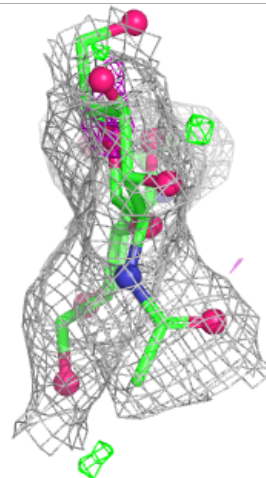
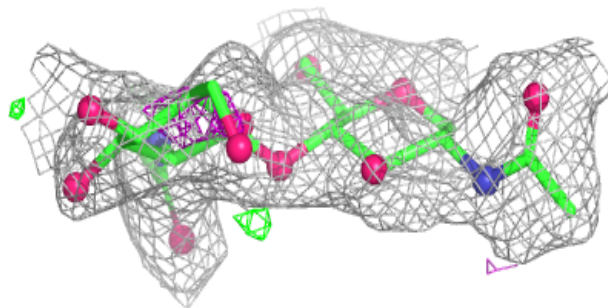
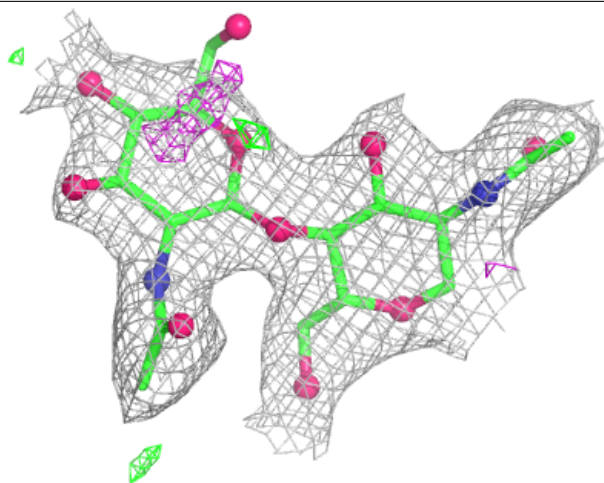
Electron density around Chain E:

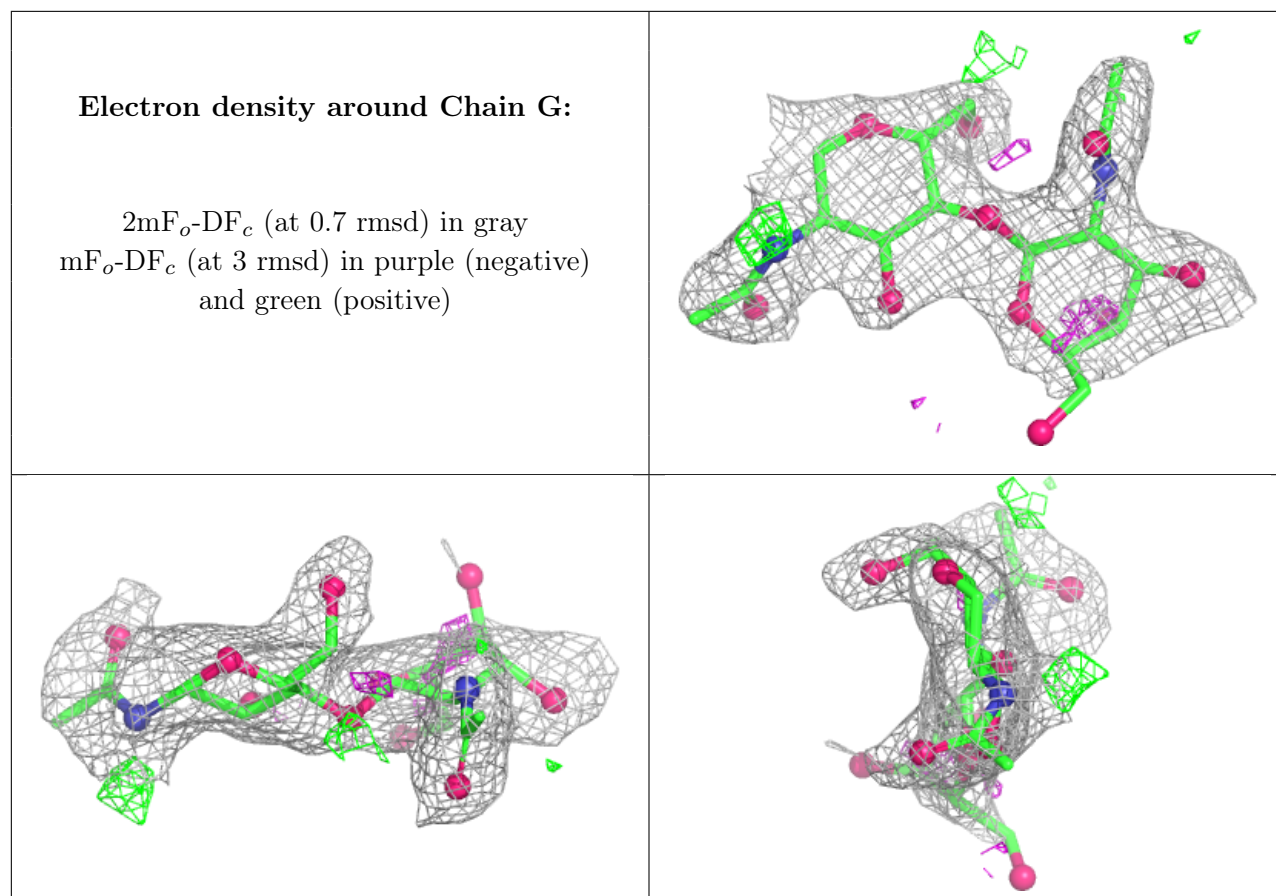
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain F:

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





6.4 Ligands [i](#)

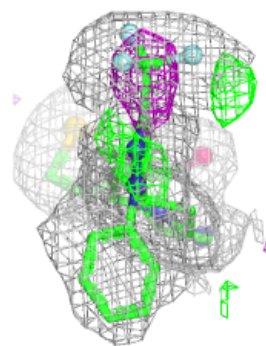
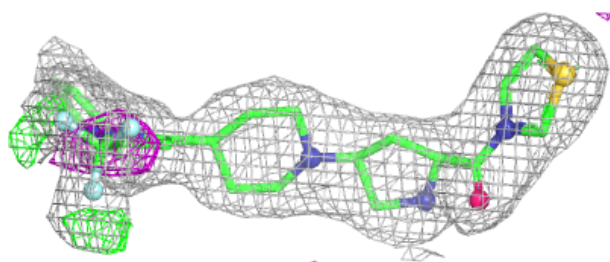
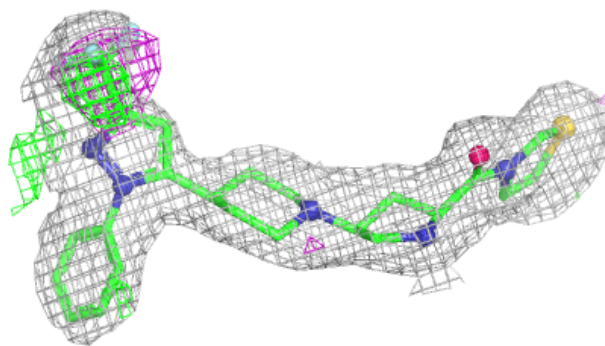
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	NAG	B	1150	14/15	0.79	0.13	45,49,51,51	0
4	NAG	A	1085	14/15	0.80	0.13	53,55,57,58	0
4	NAG	A	1321	14/15	0.82	0.11	35,38,40,41	0
4	NAG	A	1229	14/15	0.83	0.13	29,33,35,38	0
4	NAG	B	1281	14/15	0.85	0.11	46,48,49,50	0
4	NAG	B	1085	14/15	0.89	0.09	39,41,44,44	0
3	W94	B	1000	33/33	0.91	0.13	19,34,53,54	0
3	W94	A	1000	33/33	0.93	0.10	16,29,52,53	0

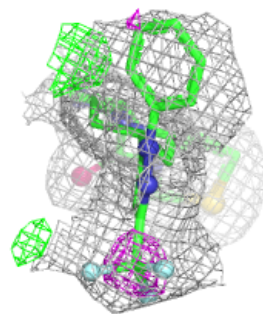
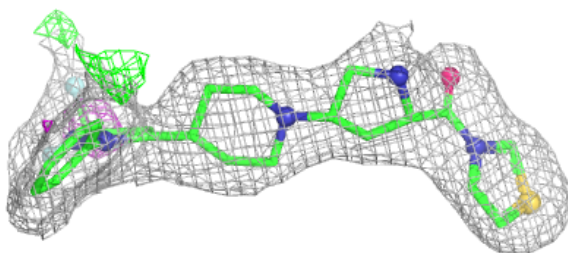
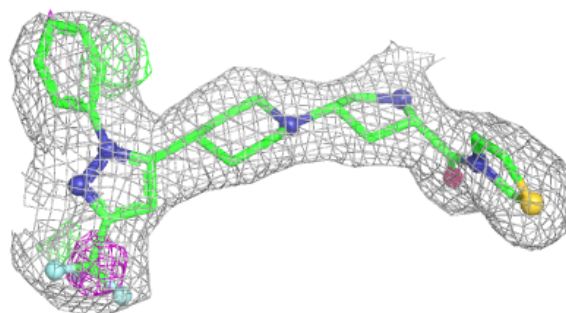
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around W94 B 1000:

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

**Electron density around W94 A 1000:**

$2mF_o-DF_c$ (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.