



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 20, 2026 – 06:09 AM UTC

PDB ID : 2QJR / pdb_00002qjr
Title : dipeptidyl peptidase IV in complex with inhibitor PZF
Authors : Shenping, L.
Deposited on : 2007-07-09
Resolution : 2.20 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

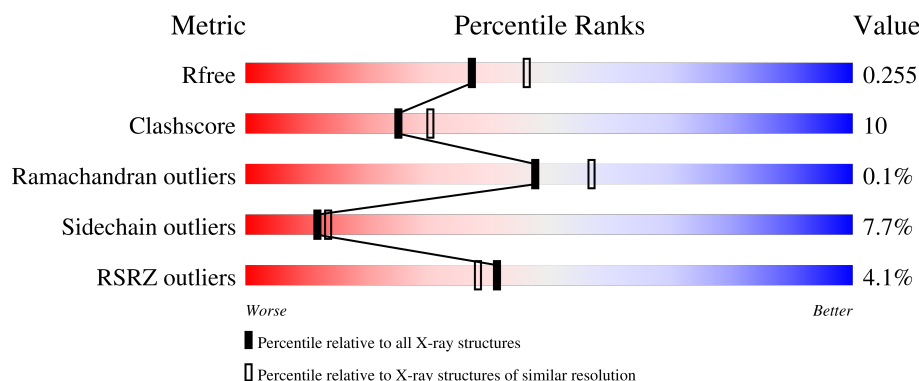
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	748	3% 73% 21% • •
1	B	748	5% 68% 24% 5% •
2	C	2	100%
2	E	2	100%
2	F	2	50% 50%

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Mol	Chain	Length	Quality of chain
2	G	2	 100%
2	H	2	 50% 50%
3	D	4	 100%

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	NAG	G	1	X	-	-	-
2	NAG	H	1	X	-	-	-
3	LGU	D	3	X	-	-	-
6	LGU	A	803	X	-	-	-

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 12532 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 membrane form.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	0	0
			5957	3825	980	1126	26			
1	B	728	Total	C	N	O	S	0	0	0
			5957	3825	980	1126	26			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	767	LEU	-	expression tag	UNP P27487
A	768	VAL	-	expression tag	UNP P27487
A	769	PRO	-	expression tag	UNP P27487
A	770	ARG	-	expression tag	UNP P27487
A	771	GLY	-	expression tag	UNP P27487
A	772	SER	-	expression tag	UNP P27487
A	773	HIS	-	expression tag	UNP P27487
A	774	HIS	-	expression tag	UNP P27487
A	775	HIS	-	expression tag	UNP P27487
A	776	HIS	-	expression tag	UNP P27487
A	777	HIS	-	expression tag	UNP P27487
A	778	HIS	-	expression tag	UNP P27487
B	767	LEU	-	expression tag	UNP P27487
B	768	VAL	-	expression tag	UNP P27487
B	769	PRO	-	expression tag	UNP P27487
B	770	ARG	-	expression tag	UNP P27487
B	771	GLY	-	expression tag	UNP P27487
B	772	SER	-	expression tag	UNP P27487
B	773	HIS	-	expression tag	UNP P27487
B	774	HIS	-	expression tag	UNP P27487
B	775	HIS	-	expression tag	UNP P27487
B	776	HIS	-	expression tag	UNP P27487
B	777	HIS	-	expression tag	UNP P27487
B	778	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			
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			
2	H	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-L-gulopyranuronic acid-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



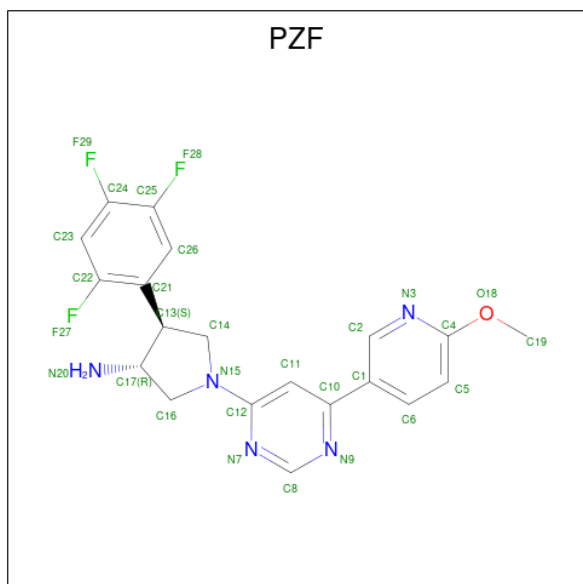
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	D	4	Total	C	N	O	0	0	0
			50	28	2	20			

- 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	B	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is (3R,4S)-1-[6-(6-METHOXYPYRIDIN-3-YL)PYRIMIDIN-4-YL]-4-(2,4,5-TRIFLUOROPHENYL)PYRROLIDIN-3-AMINE (CCD ID: PZF) (formula: $C_{20}H_{18}F_3N_5O$).



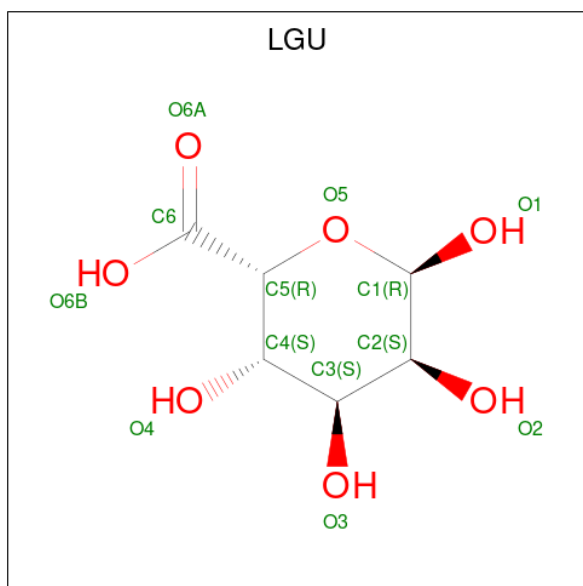
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	F	N	O	0	0
			29	20	3	5	1		

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	B	1	Total	C	F	N	O	0	0
			29	20	3	5	1		

- Molecule 6 is alpha-L-gulopyranuronic acid (CCD ID: LGU) (formula: $C_6H_{10}O_7$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			11	6	5		

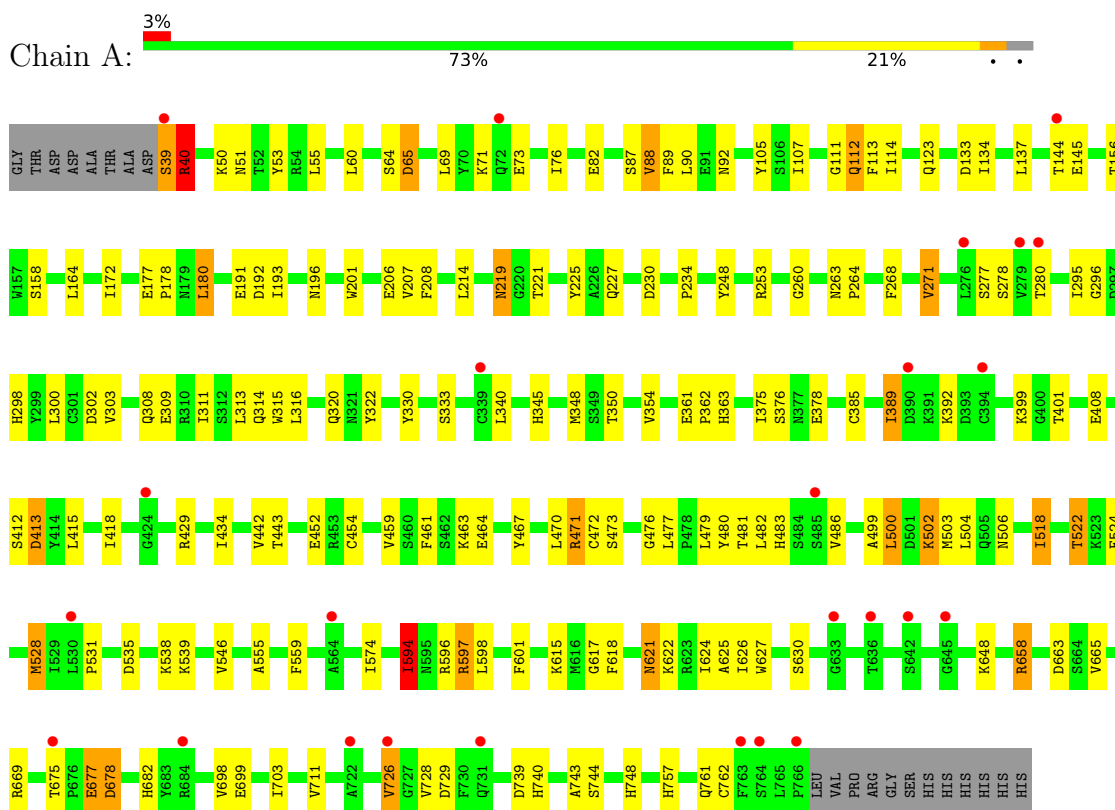
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	105	Total	O	0	0
			105	105		
7	B	226	Total	O	0	0
			226	226		

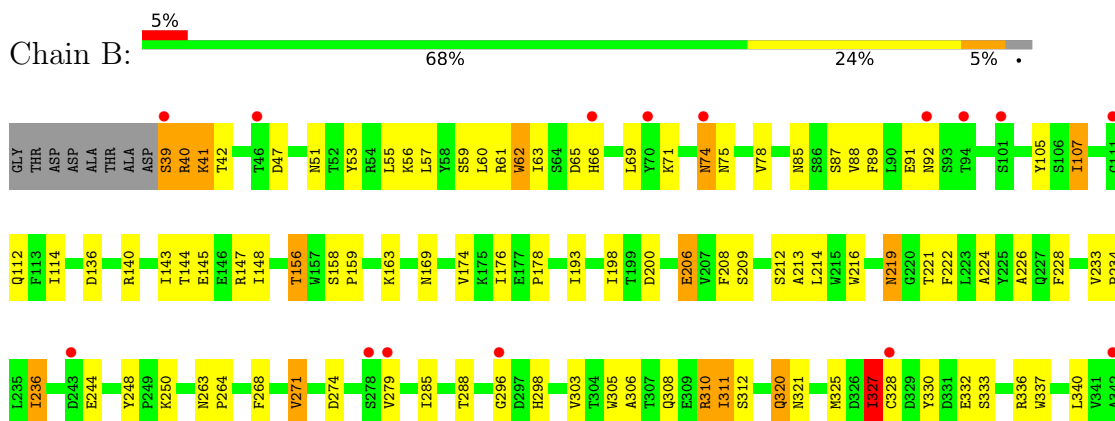
3 Residue-property plots

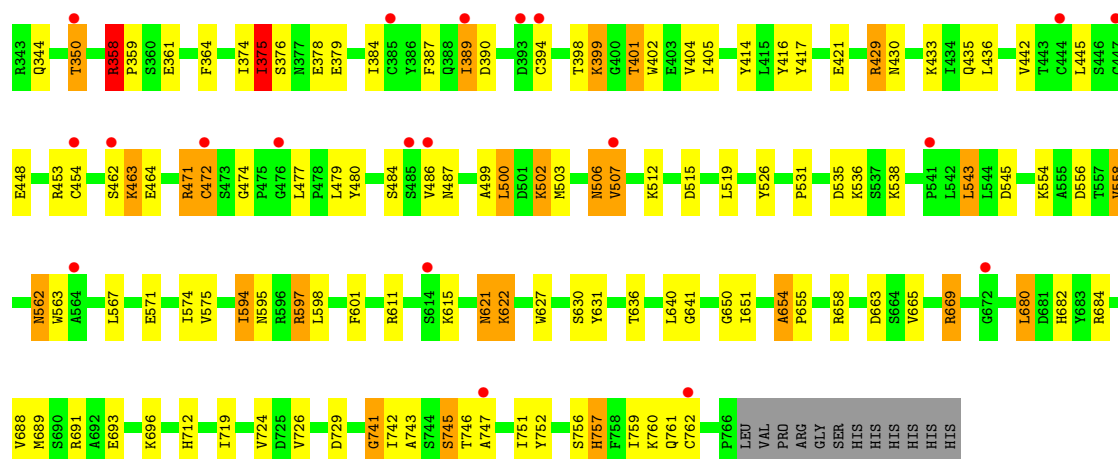
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Dipeptidyl peptidase 4 membrane form



- Molecule 1: Dipeptidyl peptidase 4 membrane form





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

Chain C: 100%



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

Chain E: 100%



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

Chain F: 50% 50%



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

Chain G: 100%



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

Chain H: 50% 50%

MAG1
MAG2

- Molecule 3: α -D-mannopyranose-(1-3)- α -L-gulopyranuronic acid-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain D:

100%

MAG1
MAG2
LGD3
MAN4

4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.47Å 68.80Å 422.61Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	41.90 – 2.20 41.90 – 2.20	Depositor EDS
% Data completeness (in resolution range)	85.7 (41.90-2.20) 85.7 (41.90-2.20)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.65 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.191 , 0.253 0.200 , 0.255	Depositor DCC
R_{free} test set	4280 reflections (2.83%)	wwPDB-VP
Wilson B-factor (Å ²)	39.1	Xtriage
Anisotropy	0.929	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 64.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.055 for k,h,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12532	wwPDB-VP
Average B, all atoms (Å ²)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.04% 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: LGU, NAG, MAN, PZF

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	1.35	22/6129 (0.4%)	1.21	13/8336 (0.2%)
1	B	1.67	57/6129 (0.9%)	1.39	38/8336 (0.5%)
All	All	1.52	79/12258 (0.6%)	1.31	51/16672 (0.3%)

All (79) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	47	ASP	CA-C	-9.46	1.40	1.52
1	B	198	ILE	C-O	-7.89	1.15	1.23
1	B	74	ASN	CA-C	7.76	1.63	1.53
1	A	658	ARG	NE-CZ	7.57	1.41	1.33
1	B	669	ARG	CD-NE	-7.51	1.35	1.46
1	B	515	ASP	C-O	-7.33	1.17	1.24
1	B	429	ARG	NE-CZ	7.13	1.40	1.33
1	B	62	TRP	C-O	-7.08	1.15	1.23
1	B	417	TYR	C-O	-6.74	1.15	1.23
1	A	88	VAL	CA-CB	6.70	1.61	1.54
1	A	315	TRP	N-CA	-6.70	1.37	1.45
1	B	486	VAL	CA-C	6.69	1.60	1.52
1	B	745	SER	N-CA	-6.49	1.38	1.46
1	B	226	ALA	C-O	6.46	1.31	1.23
1	A	739	ASP	C-O	-6.43	1.17	1.24
1	B	85	ASN	C-O	-6.35	1.16	1.23
1	B	474	GLY	C-O	-6.19	1.15	1.24
1	B	384	ILE	C-O	-6.16	1.16	1.24
1	A	665	VAL	C-O	-6.10	1.17	1.24
1	B	654	ALA	C-O	-6.08	1.20	1.23
1	B	321	ASN	CA-CB	6.06	1.64	1.53
1	B	390	ASP	N-CA	5.97	1.52	1.46
1	B	747	ALA	CA-CB	5.93	1.62	1.53
1	B	176	ILE	C-O	-5.92	1.16	1.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	430	ASN	C-O	-5.92	1.16	1.23
1	A	253	ARG	C-O	-5.88	1.17	1.24
1	A	123	GLN	CA-C	5.86	1.55	1.52
1	A	303	VAL	C-O	-5.83	1.17	1.24
1	A	678	ASP	CA-C	5.82	1.55	1.52
1	A	518	ILE	CA-CB	-5.74	1.47	1.54
1	B	350	THR	CA-CB	5.72	1.64	1.53
1	B	107	ILE	C-O	-5.66	1.16	1.24
1	A	555	ALA	CA-CB	-5.63	1.45	1.53
1	B	416	TYR	C-O	-5.61	1.17	1.24
1	B	404	VAL	CA-CB	5.60	1.60	1.54
1	A	625	ALA	CA-CB	-5.58	1.45	1.53
1	A	698	VAL	C-O	-5.57	1.17	1.23
1	B	305	TRP	C-O	-5.55	1.17	1.23
1	B	174	VAL	CA-CB	5.54	1.61	1.54
1	B	563	TRP	N-CA	5.50	1.53	1.46
1	B	574	ILE	CA-C	-5.44	1.46	1.52
1	B	436	LEU	C-O	-5.43	1.17	1.24
1	A	443	THR	CA-CB	5.39	1.60	1.53
1	B	688	VAL	CA-C	5.38	1.59	1.52
1	B	402	TRP	C-O	-5.36	1.17	1.23
1	B	472	CYS	CA-CB	5.36	1.61	1.53
1	B	471	ARG	CA-C	5.35	1.59	1.52
1	B	310	ARG	C-O	-5.35	1.18	1.23
1	B	375	ILE	CA-CB	-5.34	1.50	1.55
1	A	64	SER	N-CA	-5.34	1.39	1.46
1	B	743	ALA	C-O	-5.33	1.17	1.24
1	B	398	THR	N-CA	5.33	1.52	1.45
1	B	442	VAL	N-CA	-5.31	1.39	1.46
1	A	260	GLY	C-O	-5.30	1.16	1.24
1	B	558	VAL	CA-CB	5.29	1.60	1.54
1	A	461	PHE	C-O	-5.25	1.17	1.23
1	B	250	LYS	CD-CE	5.23	1.68	1.52
1	A	201	TRP	CA-CB	5.22	1.61	1.53
1	B	463	LYS	CE-NZ	5.20	1.65	1.49
1	A	748	HIS	C-O	-5.20	1.18	1.24
1	B	193	ILE	N-CA	5.19	1.52	1.46
1	A	476	GLY	N-CA	5.16	1.51	1.45
1	B	212	SER	C-O	-5.12	1.17	1.23
1	B	42	THR	N-CA	-5.12	1.40	1.45
1	B	228	PHE	C-O	-5.12	1.17	1.23
1	B	595	ASN	C-O	-5.12	1.17	1.23

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	531	PRO	CA-C	5.11	1.54	1.51
1	A	574	ILE	C-O	-5.10	1.18	1.24
1	B	484	SER	N-CA	5.10	1.52	1.46
1	B	631	TYR	CA-C	-5.10	1.46	1.52
1	B	311	ILE	CA-CB	5.08	1.61	1.54
1	B	421	GLU	N-CA	-5.07	1.40	1.46
1	B	222	PHE	C-O	-5.07	1.17	1.23
1	B	746	THR	C-O	-5.06	1.17	1.24
1	B	445	LEU	C-O	-5.03	1.17	1.24
1	B	399	LYS	CA-CB	-5.02	1.45	1.53
1	B	303	VAL	CA-CB	-5.02	1.47	1.54
1	B	669	ARG	NE-CZ	-5.01	1.27	1.33
1	B	507	VAL	C-O	-5.00	1.19	1.24

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	47	ASP	N-CA-C	-9.29	100.27	111.69
1	B	271	VAL	CB-CA-C	-8.39	98.06	110.81
1	B	148	ILE	N-CA-C	-8.11	101.48	109.02
1	A	389	ILE	N-CA-C	7.89	118.67	110.62
1	B	233	VAL	N-CA-C	-7.76	101.95	108.63
1	A	401	THR	OG1-CB-CG2	-7.75	93.80	109.30
1	B	394	CYS	N-CA-C	-7.37	96.61	108.55
1	B	506	ASN	N-CA-C	7.34	122.09	113.20
1	B	219	ASN	N-CA-C	7.25	120.23	111.82
1	B	321	ASN	CB-CA-C	-7.05	97.37	109.65
1	B	74	ASN	N-CA-CB	-6.82	101.68	111.84
1	B	78	VAL	CB-CA-C	-6.74	101.68	110.84
1	B	453	ARG	NE-CZ-NH2	6.73	125.26	119.20
1	A	443	THR	OG1-CB-CG2	-6.65	96.00	109.30
1	B	453	ARG	NE-CZ-NH1	-6.61	114.89	121.50
1	A	354	VAL	N-CA-C	6.53	117.99	108.58
1	B	374	ILE	CB-CA-C	-6.49	102.02	111.68
1	B	320	GLN	CA-C-N	6.41	133.35	122.26
1	B	320	GLN	C-N-CA	6.41	133.35	122.26
1	B	719	ILE	N-CA-C	6.22	116.33	110.30
1	B	321	ASN	N-CA-CB	-6.16	101.35	110.53
1	A	271	VAL	CB-CA-C	-6.12	101.50	110.62
1	B	169	ASN	CB-CA-C	-6.06	102.95	111.80
1	B	358	ARG	N-CA-C	-5.96	102.76	108.07
1	A	454	CYS	N-CA-C	5.91	118.04	108.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	39	SER	CA-CB-OG	5.80	122.69	111.10
1	B	669	ARG	NE-CZ-NH2	-5.70	114.07	119.20
1	B	724	VAL	N-CA-C	5.69	115.88	110.42
1	A	486	VAL	N-CA-C	5.59	116.05	110.23
1	B	112	GLN	N-CA-C	5.56	118.27	111.82
1	B	206	GLU	CB-CA-C	-5.46	99.90	109.07
1	B	156	THR	CB-CA-C	-5.45	98.56	109.35
1	B	741	GLY	N-CA-C	-5.37	107.28	115.66
1	B	454	CYS	N-CA-C	5.36	117.56	108.02
1	A	597	ARG	N-CA-CB	-5.34	105.38	111.79
1	B	213	ALA	CA-C-N	5.31	130.92	122.74
1	B	213	ALA	C-N-CA	5.31	130.92	122.74
1	B	327	ILE	CA-C-N	-5.31	115.51	122.99
1	B	327	ILE	C-N-CA	-5.31	115.51	122.99
1	A	219	ASN	CB-CA-C	-5.26	100.57	110.67
1	A	112	GLN	N-CA-C	5.23	117.88	111.82
1	B	311	ILE	CA-C-N	-5.22	115.36	122.93
1	B	311	ILE	C-N-CA	-5.22	115.36	122.93
1	B	463	LYS	CA-C-N	-5.14	116.15	125.66
1	B	463	LYS	C-N-CA	-5.14	116.15	125.66
1	B	575	VAL	N-CA-C	5.13	115.48	107.99
1	B	219	ASN	CB-CA-C	-5.10	100.88	110.67
1	A	40	ARG	N-CA-CB	5.09	119.10	110.49
1	B	757	HIS	N-CA-C	-5.01	105.89	111.36
1	A	594	ILE	N-CA-C	-5.01	107.87	112.83
1	B	472	CYS	N-CA-CB	5.01	118.59	110.43

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	5957	0	5672	101	0
1	B	5957	0	5672	130	0
2	C	28	0	25	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	28	0	25	0	0
2	F	28	0	25	1	0
2	G	28	0	25	2	0
2	H	28	0	25	1	0
3	D	50	0	40	0	0
4	A	14	0	13	0	0
4	B	14	0	13	1	0
5	A	29	0	18	1	0
5	B	29	0	18	0	0
6	A	11	0	7	0	0
7	A	105	0	0	10	0
7	B	226	0	0	22	0
All	All	12532	0	11578	231	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 10.

All (231) 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:39:SER:O	1:A:40:ARG:CZ	1.79	1.29
1:B:40:ARG:HD3	1:B:40:ARG:N	1.64	1.12
1:B:39:SER:HB2	1:B:40:ARG:HD2	1.31	1.08
1:A:621:ASN:C	1:A:621:ASN:HD22	1.64	1.04
1:B:597:ARG:HD3	7:B:992:HOH:O	1.56	1.02
1:B:39:SER:HB2	1:B:40:ARG:CD	1.91	0.99
1:B:74:ASN:HB2	7:B:1034:HOH:O	1.63	0.94
1:A:630:SER:HG	1:A:740:HIS:HE2	1.05	0.90
1:B:401:THR:HG22	1:B:401:THR:O	1.74	0.87
1:B:39:SER:O	7:B:901:HOH:O	1.93	0.87
1:B:40:ARG:HD3	1:B:40:ARG:H	1.29	0.87
1:B:147:ARG:HD3	7:B:905:HOH:O	1.75	0.87
1:A:434:ILE:HD12	1:A:442:VAL:HG22	1.59	0.85
1:B:499:ALA:O	1:B:502:LYS:HG2	1.78	0.84
1:B:65:ASP:OD1	1:B:463:LYS:O	1.95	0.83
1:B:236:ILE:HG12	1:B:712:HIS:CE1	2.13	0.82
1:A:39:SER:O	1:A:40:ARG:NH1	2.14	0.80
1:A:39:SER:O	1:A:40:ARG:NH2	2.16	0.79
1:B:598:LEU:O	1:B:682:HIS:HE1	1.65	0.79
1:B:325:MET:HE2	1:B:327:ILE:HD11	1.66	0.77
1:B:219:ASN:HB2	1:B:308:GLN:OE1	1.85	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:61:ARG:HD3	7:B:1068:HOH:O	1.85	0.75
1:A:345:HIS:HD2	7:A:969:HOH:O	1.70	0.74
1:A:630:SER:OG	1:A:740:HIS:NE2	2.09	0.74
1:B:66:HIS:HD2	7:B:1105:HOH:O	1.71	0.74
1:B:306:ALA:HB3	1:B:310:ARG:HG2	1.70	0.73
1:B:389:ILE:O	7:B:902:HOH:O	2.06	0.73
1:A:658:ARG:NE	7:A:901:HOH:O	1.93	0.72
1:B:66:HIS:CD2	7:B:1105:HOH:O	2.42	0.72
1:B:477:LEU:HD22	1:B:500:LEU:HD13	1.73	0.71
1:A:463:LYS:O	1:A:464:GLU:HB2	1.91	0.70
1:A:177:GLU:HB2	1:A:180:LEU:HD22	1.73	0.70
1:A:71:LYS:HZ1	1:A:92:ASN:HD21	1.40	0.69
1:A:499:ALA:O	1:A:502:LYS:HG3	1.91	0.69
1:B:74:ASN:O	1:B:75:ASN:OD1	2.10	0.69
1:A:206:GLU:OE2	1:A:663:ASP:OD2	2.11	0.68
1:B:55:LEU:HD23	1:B:500:LEU:CD1	2.23	0.68
1:A:89:PHE:CE2	1:A:107:ILE:HD13	2.29	0.68
1:B:57:LEU:HD23	7:B:958:HOH:O	1.94	0.68
1:A:71:LYS:NZ	1:A:92:ASN:HD21	1.92	0.68
1:B:89:PHE:CE2	1:B:107:ILE:HD13	2.30	0.67
1:B:71:LYS:NZ	1:B:74:ASN:HA	2.09	0.67
1:B:621:ASN:C	1:B:621:ASN:HD22	2.02	0.66
1:A:761:GLN:NE2	7:A:903:HOH:O	2.26	0.66
1:B:39:SER:CB	1:B:40:ARG:HH11	2.09	0.66
1:A:621:ASN:HD22	1:A:622:LYS:N	1.93	0.65
1:B:503:MET:O	1:B:506:ASN:HB2	1.97	0.65
1:B:147:ARG:NH1	7:B:905:HOH:O	2.29	0.64
1:A:53:TYR:CD1	1:A:503:MET:HE2	2.32	0.64
1:A:345:HIS:CD2	7:A:969:HOH:O	2.46	0.64
1:B:680:LEU:HD22	1:B:684:ARG:HD3	1.80	0.63
1:B:89:PHE:HE2	1:B:107:ILE:HD13	1.64	0.63
1:B:429:ARG:NE	7:B:903:HOH:O	2.18	0.63
1:A:298:HIS:HE1	7:A:970:HOH:O	1.81	0.63
1:B:414:TYR:CD1	1:B:433:LYS:HD2	2.32	0.63
1:B:598:LEU:O	1:B:682:HIS:CE1	2.49	0.63
1:B:71:LYS:HZ1	1:B:74:ASN:HA	1.64	0.63
1:B:658:ARG:HB2	1:B:689:MET:HE1	1.80	0.62
1:A:111:GLY:O	1:A:137:LEU:HD12	2.00	0.62
1:A:621:ASN:C	1:A:621:ASN:ND2	2.42	0.62
1:B:364:PHE:HE2	1:B:389:ILE:HD11	1.65	0.62
1:B:358:ARG:HB2	1:B:359:PRO:HD2	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:LEU:HD22	1:A:504:LEU:HG	1.82	0.62
1:B:640:LEU:HD11	1:B:650:GLY:HA3	1.80	0.62
1:A:361:GLU:OE2	1:A:363:HIS:HE1	1.83	0.61
1:B:40:ARG:N	1:B:40:ARG:CD	2.48	0.61
1:B:163:LYS:NZ	1:B:274:ASP:OD1	2.31	0.61
1:B:562:ASN:C	1:B:562:ASN:HD22	2.08	0.61
1:B:39:SER:HB2	1:B:40:ARG:HD3	1.80	0.60
4:B:801:NAG:C4	7:B:910:HOH:O	2.49	0.60
1:B:401:THR:O	1:B:401:THR:CG2	2.44	0.60
1:B:40:ARG:CB	1:B:506:ASN:O	2.50	0.59
1:A:39:SER:O	1:A:40:ARG:NE	2.32	0.59
1:A:434:ILE:HD12	1:A:442:VAL:CG2	2.32	0.58
1:B:680:LEU:HD22	1:B:684:ARG:CD	2.32	0.58
1:A:76:ILE:HD12	1:A:90:LEU:HB3	1.85	0.58
1:B:405:ILE:HG13	1:B:429:ARG:CD	2.33	0.58
1:A:528:MET:HE1	1:A:618:PHE:CZ	2.38	0.58
1:B:325:MET:HE2	1:B:327:ILE:CD1	2.33	0.57
1:B:55:LEU:HD23	1:B:500:LEU:HD12	1.85	0.57
1:A:470:LEU:HD12	1:A:483:HIS:CE1	2.40	0.57
1:A:481:THR:OG1	1:A:483:HIS:HE1	1.88	0.57
1:B:535:ASP:OD2	1:B:538:LYS:NZ	2.38	0.57
1:B:55:LEU:HD23	1:B:500:LEU:HD11	1.87	0.56
1:B:39:SER:OG	1:B:40:ARG:NH1	2.39	0.56
1:A:89:PHE:HE2	1:A:107:ILE:HD13	1.71	0.55
1:B:693:GLU:CD	1:B:726:VAL:HG21	2.32	0.55
1:B:597:ARG:CD	7:B:992:HOH:O	2.30	0.55
1:A:207:VAL:HG12	1:A:208:PHE:CD1	2.42	0.55
1:A:134:ILE:HD11	1:A:164:LEU:CD1	2.37	0.54
1:B:40:ARG:HB2	1:B:506:ASN:O	2.07	0.54
1:B:296:GLY:O	1:B:298:HIS:HD2	1.90	0.54
7:B:1032:HOH:O	2:G:1:NAG:H82	2.07	0.54
1:A:65:ASP:OD1	1:A:65:ASP:N	2.27	0.53
1:B:399:LYS:NZ	7:B:915:HOH:O	2.42	0.53
1:A:314:GLN:HE22	1:A:362:PRO:HD3	1.73	0.53
1:A:726:VAL:CG1	1:A:728:VAL:HG23	2.39	0.53
1:A:219:ASN:HB2	1:A:308:GLN:OE1	2.08	0.53
1:A:320:GLN:OE1	1:A:669:ARG:HD3	2.10	0.52
1:A:624:ILE:HD12	7:A:916:HOH:O	2.07	0.52
1:B:39:SER:C	1:B:40:ARG:HD3	2.32	0.52
1:A:675:THR:HB	1:A:677:GLU:OE1	2.09	0.52
1:A:594:ILE:CD1	1:A:601:PHE:HB2	2.40	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:545:ASP:OD1	1:B:554:LYS:NZ	2.40	0.52
1:A:528:MET:HE1	1:A:618:PHE:HZ	1.74	0.51
1:A:191:GLU:O	1:A:192:ASP:HB2	2.10	0.51
1:A:230:ASP:OD1	1:A:264:PRO:HB3	2.09	0.51
1:A:378:GLU:CG	7:A:906:HOH:O	2.58	0.51
1:A:546:VAL:HG21	1:A:626:ILE:HD11	1.93	0.51
1:A:302:ASP:HB3	1:A:314:GLN:HB2	1.92	0.51
1:A:206:GLU:CD	5:A:802:PZF:H202	2.20	0.50
1:A:296:GLY:O	1:A:298:HIS:HD2	1.94	0.50
1:A:596:ARG:NH2	1:A:678:ASP:OD1	2.42	0.50
1:B:463:LYS:O	1:B:464:GLU:HB2	2.12	0.50
1:B:87:SER:OG	2:F:1:NAG:O7	2.29	0.50
1:B:621:ASN:HD22	1:B:622:LYS:N	2.09	0.50
1:B:594:ILE:CD1	1:B:601:PHE:HB2	2.42	0.49
1:A:415:LEU:HD23	1:A:415:LEU:C	2.37	0.49
1:A:40:ARG:NE	1:A:40:ARG:HA	2.24	0.49
7:B:1124:HOH:O	2:H:1:NAG:H83	2.12	0.49
1:B:477:LEU:CD2	1:B:500:LEU:HD13	2.42	0.49
1:B:611:ARG:O	1:B:615:LYS:HG2	2.13	0.49
1:B:200:ASP:OD2	1:B:264:PRO:HG3	2.12	0.49
1:B:206:GLU:HG3	1:B:665:VAL:HB	1.95	0.49
1:A:594:ILE:HG23	1:A:594:ILE:O	2.13	0.49
1:A:322:TYR:CE2	1:A:348:MET:HE2	2.47	0.49
1:B:463:LYS:O	1:B:464:GLU:CB	2.58	0.48
1:B:41:LYS:HE2	1:B:53:TYR:OH	2.13	0.48
1:A:473:SER:OG	7:A:902:HOH:O	2.20	0.48
1:B:224:ALA:HB1	1:B:268:PHE:CZ	2.48	0.48
1:A:192:ASP:C	1:A:193:ILE:HG13	2.39	0.48
1:A:757:HIS:HD2	1:B:729:ASP:OD2	1.97	0.48
1:A:726:VAL:HG13	1:A:728:VAL:HG23	1.94	0.48
1:A:477:LEU:HD22	1:A:500:LEU:HD13	1.95	0.48
1:A:522:THR:HG22	1:A:524:PHE:CE2	2.50	0.47
1:A:502:LYS:HE2	1:A:503:MET:HG3	1.95	0.47
1:A:408:GLU:HG3	1:A:418:ILE:HD12	1.97	0.47
1:B:39:SER:CB	1:B:40:ARG:CD	2.81	0.46
1:A:69:LEU:HD13	1:A:107:ILE:HD12	1.98	0.46
1:A:630:SER:HG	1:A:740:HIS:CE1	2.27	0.46
1:B:39:SER:N	1:B:40:ARG:HH11	2.14	0.46
1:B:310:ARG:HA	1:B:328:CYS:O	2.15	0.46
1:B:571:GLU:CD	1:B:760:LYS:HD3	2.40	0.46
1:A:105:TYR:HB2	1:A:114:ILE:HD11	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:413:ASP:OD2	1:A:413:ASP:N	2.49	0.46
1:A:361:GLU:OE2	1:A:363:HIS:CE1	2.68	0.46
1:A:55:LEU:HD21	1:A:559:PHE:HE2	1.80	0.46
1:B:693:GLU:O	1:B:696:LYS:HG3	2.15	0.46
1:B:761:GLN:C	1:B:761:GLN:OE1	2.59	0.46
1:A:429:ARG:NE	7:A:911:HOH:O	2.43	0.46
1:A:134:ILE:HD11	1:A:164:LEU:HD11	1.97	0.46
1:B:358:ARG:HB2	1:B:359:PRO:CD	2.46	0.45
1:B:414:TYR:CE1	1:B:433:LYS:HD2	2.50	0.45
1:B:751:ILE:HG23	1:B:752:TYR:N	2.32	0.45
1:B:636:THR:HG21	1:B:651:ILE:O	2.15	0.45
1:B:41:LYS:CE	1:B:53:TYR:OH	2.65	0.45
1:A:375:ILE:HG23	7:A:994:HOH:O	2.17	0.45
1:B:158:SER:HB2	1:B:159:PRO:CD	2.47	0.45
1:B:562:ASN:C	1:B:562:ASN:ND2	2.73	0.45
1:B:143:ILE:HD13	1:B:178:PRO:HB2	2.00	0.44
1:B:756:SER:O	1:B:760:LYS:HG3	2.18	0.44
1:B:144:THR:HG22	7:B:1120:HOH:O	2.17	0.44
1:B:526:TYR:CD2	1:B:526:TYR:C	2.95	0.44
1:B:306:ALA:CB	1:B:310:ARG:HG2	2.44	0.44
1:A:277:SER:HG	1:A:280:THR:H	1.62	0.44
1:A:648:LYS:NZ	1:A:699:GLU:OE1	2.50	0.44
1:B:325:MET:CE	1:B:327:ILE:HD11	2.43	0.44
1:A:214:LEU:HD23	1:A:225:TYR:HB3	2.00	0.44
1:A:471:ARG:HD2	1:A:480:TYR:CE2	2.53	0.44
1:B:325:MET:O	1:B:344:GLN:HA	2.18	0.44
1:B:359:PRO:HA	7:B:908:HOH:O	2.18	0.44
1:B:105:TYR:HB2	1:B:114:ILE:HD11	1.98	0.44
1:A:268:PHE:CZ	1:A:313:LEU:HD21	2.54	0.43
1:B:206:GLU:OE2	1:B:663:ASP:OD2	2.36	0.43
1:A:528:MET:HE3	1:A:618:PHE:CE1	2.53	0.43
1:B:654:ALA:N	1:B:655:PRO:CD	2.81	0.43
1:A:375:ILE:HG22	1:A:376:SER:N	2.33	0.43
1:A:535:ASP:OD2	1:A:538:LYS:HG2	2.18	0.43
1:A:598:LEU:O	1:A:682:HIS:NE2	2.39	0.43
1:B:63:ILE:HG21	1:B:69:LEU:CD1	2.48	0.43
1:B:144:THR:CG2	7:B:1120:HOH:O	2.66	0.43
1:B:641:GLY:O	1:B:691:ARG:HB3	2.17	0.43
1:B:759:ILE:HG23	1:B:759:ILE:HD12	1.66	0.43
1:A:134:ILE:CD1	1:A:164:LEU:HD11	2.49	0.43
1:A:248:TYR:CZ	1:B:234:PRO:HB2	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:158:SER:HA	1:B:216:TRP:CD1	2.53	0.43
1:B:332:GLU:OE1	2:G:2:NAG:O3	2.36	0.43
1:A:729:ASP:OD2	1:B:757:HIS:HD2	2.02	0.43
1:B:74:ASN:HB3	1:B:92:ASN:CG	2.44	0.43
1:A:309:GLU:HB3	1:A:330:TYR:HB3	2.00	0.42
1:B:136:ASP:OD1	1:B:136:ASP:C	2.60	0.42
1:A:648:LYS:NZ	1:A:699:GLU:OE2	2.52	0.42
2:C:1:NAG:H61	2:C:2:NAG:C7	2.49	0.42
1:A:172:ILE:HD13	1:A:214:LEU:HD21	2.01	0.42
1:A:418:ILE:HD11	1:A:459:VAL:HG12	2.01	0.42
1:B:556:ASP:C	1:B:556:ASP:OD1	2.63	0.42
1:B:208:PHE:O	1:B:209:SER:C	2.58	0.42
1:B:312:SER:HB2	1:B:325:MET:HE3	2.01	0.42
1:B:627:TRP:HB2	1:B:651:ILE:HB	2.02	0.42
1:A:69:LEU:CD1	1:A:107:ILE:HD12	2.50	0.42
1:A:82:GLU:HB2	1:A:467:TYR:OH	2.19	0.41
1:B:364:PHE:CE2	1:B:389:ILE:HD11	2.51	0.41
1:A:711:VAL:CG2	1:A:740:HIS:CE1	3.03	0.41
1:B:448:GLU:OE1	1:B:448:GLU:HA	2.20	0.41
1:B:320:GLN:OE1	1:B:669:ARG:HD3	2.21	0.41
1:B:375:ILE:CD1	1:B:387:PHE:HZ	2.34	0.41
1:B:543:LEU:HD12	1:B:567:LEU:HD13	2.01	0.41
1:A:112:GLN:HB3	1:A:113:PHE:CE2	2.55	0.41
1:A:234:PRO:HB2	1:B:248:TYR:CZ	2.56	0.41
1:A:477:LEU:HD22	1:A:500:LEU:CD1	2.51	0.41
1:B:330:TYR:HB2	1:B:337:TRP:CH2	2.55	0.41
1:B:405:ILE:HG13	1:B:429:ARG:HD3	2.03	0.41
1:B:487:ASN:N	7:B:907:HOH:O	2.34	0.41
1:A:196:ASN:OD1	1:A:227:GLN:HG3	2.21	0.41
1:A:703:ILE:HD13	1:A:703:ILE:HG21	1.78	0.41
1:B:512:LYS:NZ	1:B:556:ASP:O	2.49	0.41
1:B:741:GLY:O	1:B:742:ILE:C	2.63	0.41
1:B:742:ILE:O	1:B:742:ILE:HG22	2.20	0.41
1:A:177:GLU:HA	1:A:178:PRO:HD3	1.94	0.41
1:A:539:LYS:HE2	1:A:617:GLY:O	2.21	0.41
1:B:62:TRP:CG	1:B:462:SER:HA	2.56	0.41
1:B:39:SER:CB	1:B:40:ARG:HD3	2.49	0.41
1:B:236:ILE:HG12	1:B:712:HIS:ND1	2.34	0.41
1:B:471:ARG:HG3	1:B:480:TYR:CE2	2.56	0.41
1:A:481:THR:OG1	1:A:483:HIS:CE1	2.71	0.40
1:A:743:ALA:O	1:A:744:SER:C	2.59	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:435:GLN:NE2	7:B:934:HOH:O	2.51	0.40
1:A:677:GLU:CD	1:A:677:GLU:H	2.29	0.40
1:B:285:ILE:HG23	1:B:336:ARG:HH11	1.85	0.40
1:B:669:ARG:HD2	7:B:1085:HOH:O	2.20	0.40
1:B:40:ARG:HG2	1:B:506:ASN:C	2.46	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	726/748 (97%)	694 (96%)	31 (4%)	1 (0%)	48	57
1	B	726/748 (97%)	689 (95%)	36 (5%)	1 (0%)	48	57
All	All	1452/1496 (97%)	1383 (95%)	67 (5%)	2 (0%)	48	57

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	73	GLU
1	B	630	SER

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	651/669 (97%)	602 (92%)	49 (8%)	12	14
1	B	651/669 (97%)	600 (92%)	51 (8%)	11	13
All	All	1302/1338 (97%)	1202 (92%)	100 (8%)	12	13

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	40	ARG
1	A	50	LYS
1	A	51	ASN
1	A	60	LEU
1	A	65	ASP
1	A	87	SER
1	A	88	VAL
1	A	133	ASP
1	A	144	THR
1	A	145	GLU
1	A	156	THR
1	A	158	SER
1	A	180	LEU
1	A	221	THR
1	A	263	ASN
1	A	271	VAL
1	A	278	SER
1	A	295	ILE
1	A	300	LEU
1	A	311	ILE
1	A	316	LEU
1	A	333	SER
1	A	340	LEU
1	A	350	THR
1	A	385	CYS
1	A	389	ILE
1	A	392	LYS
1	A	399	LYS
1	A	412	SER
1	A	413	ASP
1	A	452	GLU
1	A	471	ARG
1	A	472	CYS
1	A	479	LEU
1	A	482	LEU

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Mol	Chain	Res	Type
1	A	500	LEU
1	A	502	LYS
1	A	506	ASN
1	A	518	ILE
1	A	522	THR
1	A	528	MET
1	A	594	ILE
1	A	597	ARG
1	A	615	LYS
1	A	621	ASN
1	A	627	TRP
1	A	677	GLU
1	A	726	VAL
1	A	762	CYS
1	B	39	SER
1	B	40	ARG
1	B	41	LYS
1	B	51	ASN
1	B	56	LYS
1	B	59	SER
1	B	60	LEU
1	B	88	VAL
1	B	91	GLU
1	B	140	ARG
1	B	145	GLU
1	B	156	THR
1	B	214	LEU
1	B	221	THR
1	B	236	ILE
1	B	244	GLU
1	B	263	ASN
1	B	271	VAL
1	B	279	VAL
1	B	288	THR
1	B	311	ILE
1	B	327	ILE
1	B	333	SER
1	B	340	LEU
1	B	350	THR
1	B	358	ARG
1	B	361	GLU
1	B	375	ILE

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Mol	Chain	Res	Type
1	B	376	SER
1	B	378	GLU
1	B	379	GLU
1	B	389	ILE
1	B	401	THR
1	B	472	CYS
1	B	479	LEU
1	B	500	LEU
1	B	502	LYS
1	B	507	VAL
1	B	519	LEU
1	B	531	PRO
1	B	536	LYS
1	B	543	LEU
1	B	558	VAL
1	B	562	ASN
1	B	594	ILE
1	B	597	ARG
1	B	621	ASN
1	B	622	LYS
1	B	680	LEU
1	B	745	SER
1	B	762	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	51	ASN
1	A	80	ASN
1	A	92	ASN
1	A	169	ASN
1	A	227	GLN
1	A	247	GLN
1	A	263	ASN
1	A	298	HIS
1	A	314	GLN
1	A	338	ASN
1	A	363	HIS
1	A	430	ASN
1	A	483	HIS
1	A	533	HIS
1	A	595	ASN

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Mol	Chain	Res	Type
1	A	612	GLN
1	A	621	ASN
1	A	748	HIS
1	A	757	HIS
1	B	75	ASN
1	B	123	GLN
1	B	169	ASN
1	B	227	GLN
1	B	263	ASN
1	B	298	HIS
1	B	383	HIS
1	B	430	ASN
1	B	450	ASN
1	B	487	ASN
1	B	506	ASN
1	B	562	ASN
1	B	572	ASN
1	B	621	ASN
1	B	682	HIS
1	B	697	GLN
1	B	757	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 ⓘ

14 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	1,2	14,14,15	1.50	3 (21%)	17,19,21	1.63	3 (17%)
2	NAG	C	2	2	14,14,15	0.87	0	17,19,21	1.97	6 (35%)
3	NAG	D	1	1,3	14,14,15	1.29	2 (14%)	17,19,21	2.22	9 (52%)
3	NAG	D	2	3	14,14,15	0.92	0	17,19,21	1.11	1 (5%)
3	LGU	D	3	6,3	11,11,13	0.82	1 (9%)	15,15,19	2.25	5 (33%)
3	MAN	D	4	3	11,11,12	1.06	1 (9%)	15,15,17	1.29	1 (6%)
2	NAG	E	1	1,2	14,14,15	0.74	0	17,19,21	2.56	6 (35%)
2	NAG	E	2	2	14,14,15	0.84	1 (7%)	17,19,21	1.27	2 (11%)
2	NAG	F	1	1,2	14,14,15	1.21	2 (14%)	17,19,21	2.26	4 (23%)
2	NAG	F	2	2	14,14,15	1.17	1 (7%)	17,19,21	1.74	5 (29%)
2	NAG	G	1	1,2	14,14,15	1.16	1 (7%)	17,19,21	1.91	4 (23%)
2	NAG	G	2	2	14,14,15	0.52	0	17,19,21	1.55	2 (11%)
2	NAG	H	1	1,2	14,14,15	0.87	1 (7%)	17,19,21	1.79	3 (17%)
2	NAG	H	2	2	14,14,15	0.65	0	17,19,21	1.85	5 (29%)

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	1,2	-	0/6/23/26	0/1/1/1
2	NAG	C	2	2	-	2/6/23/26	0/1/1/1
3	NAG	D	1	1,3	-	2/6/23/26	0/1/1/1
3	NAG	D	2	3	-	0/6/23/26	0/1/1/1
3	LGU	D	3	6,3	1/1/4/6	2/2/19/24	0/1/1/1
3	MAN	D	4	3	-	2/2/19/22	0/1/1/1
2	NAG	E	1	1,2	-	2/6/23/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	4/6/23/26	0/1/1/1
2	NAG	G	1	1,2	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	G	2	2	-	4/6/23/26	0/1/1/1
2	NAG	H	1	1,2	1/1/5/7	0/6/23/26	0/1/1/1
2	NAG	H	2	2	-	2/6/23/26	0/1/1/1

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	2	NAG	O5-C1	-3.41	1.38	1.43
3	D	1	NAG	O5-C1	-3.36	1.38	1.43
2	C	1	NAG	C2-N2	-3.24	1.40	1.46
2	C	1	NAG	O5-C1	-3.02	1.38	1.43
2	G	1	NAG	C1-C2	2.69	1.56	1.52
2	H	1	NAG	O5-C1	-2.59	1.39	1.43
3	D	4	MAN	O5-C1	-2.58	1.39	1.43
3	D	1	NAG	O5-C5	-2.42	1.38	1.43
2	F	1	NAG	C2-N2	-2.40	1.42	1.46
2	E	2	NAG	C1-C2	2.29	1.55	1.52
2	C	1	NAG	C1-C2	2.22	1.55	1.52
3	D	3	LGU	C4-C5	2.14	1.57	1.53
2	F	1	NAG	O5-C5	-2.04	1.39	1.43

All (56) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	NAG	C1-O5-C5	7.03	121.60	112.19
2	F	1	NAG	O5-C1-C2	-6.97	100.50	111.29
3	D	3	LGU	C1-O5-C5	5.28	119.27	112.19
2	E	1	NAG	O5-C1-C2	5.21	119.35	111.29
2	G	1	NAG	O5-C1-C2	5.12	119.22	111.29
2	C	2	NAG	C1-O5-C5	5.11	119.04	112.19
2	G	2	NAG	C1-O5-C5	4.99	118.87	112.19
2	H	1	NAG	O5-C1-C2	4.88	118.84	111.29
2	H	2	NAG	C1-O5-C5	4.51	118.24	112.19
2	C	1	NAG	C2-N2-C7	4.32	128.69	122.90
3	D	1	NAG	C1-C2-N2	4.26	117.15	110.43
3	D	3	LGU	C1-C2-C3	-4.01	103.80	109.64
3	D	4	MAN	O5-C1-C2	-3.71	101.93	110.79
2	F	2	NAG	O5-C1-C2	3.12	116.11	111.29
2	H	2	NAG	C1-C2-N2	-3.12	105.52	110.43
2	H	1	NAG	O7-C7-N2	3.11	127.48	121.98
2	G	1	NAG	O4-C4-C3	3.07	117.62	110.38
2	F	2	NAG	O7-C7-N2	3.04	127.36	121.98
3	D	1	NAG	O6-C6-C5	-3.04	100.97	111.33
3	D	1	NAG	O3-C3-C4	-3.03	103.22	110.38
2	F	2	NAG	C3-C4-C5	3.00	115.67	110.23
3	D	3	LGU	O2-C2-C1	2.98	116.05	109.22
2	H	2	NAG	O5-C1-C2	2.95	115.86	111.29
2	E	1	NAG	C3-C4-C5	2.95	115.58	110.23
3	D	1	NAG	C3-C4-C5	2.95	115.58	110.23
2	E	1	NAG	O5-C5-C4	2.82	117.68	110.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	H	2	NAG	O4-C4-C5	2.74	116.08	109.32
3	D	1	NAG	O4-C4-C3	-2.63	104.18	110.38
3	D	2	NAG	O4-C4-C3	-2.59	104.27	110.38
2	H	2	NAG	C3-C4-C5	-2.57	105.58	110.23
2	F	2	NAG	C1-O5-C5	2.56	115.62	112.19
3	D	1	NAG	O5-C5-C6	-2.52	102.76	107.66
2	F	1	NAG	O3-C3-C2	-2.52	104.17	109.40
2	E	1	NAG	O5-C5-C6	2.45	112.43	107.66
2	F	1	NAG	O7-C7-C8	-2.44	117.72	122.05
2	F	2	NAG	O4-C4-C3	-2.39	104.73	110.38
2	C	2	NAG	C2-N2-C7	2.36	126.07	122.90
2	E	2	NAG	O5-C5-C6	2.36	112.26	107.66
2	C	2	NAG	O6-C6-C5	-2.35	103.34	111.33
3	D	1	NAG	O5-C5-C4	2.28	116.37	110.83
2	C	2	NAG	O3-C3-C2	2.27	114.12	109.40
2	C	2	NAG	C6-C5-C4	-2.27	107.45	113.02
2	E	2	NAG	C4-C3-C2	2.23	114.29	111.02
2	C	1	NAG	O5-C5-C4	2.20	116.17	110.83
3	D	3	LGU	C2-C3-C4	-2.18	107.02	110.86
3	D	1	NAG	C2-N2-C7	-2.18	119.98	122.90
2	F	1	NAG	C1-O5-C5	2.17	115.10	112.19
2	G	1	NAG	O7-C7-C8	-2.15	118.23	122.05
3	D	3	LGU	O5-C5-C4	2.13	116.01	110.83
2	C	1	NAG	O5-C1-C2	-2.11	108.02	111.29
2	G	1	NAG	C2-N2-C7	2.11	125.73	122.90
3	D	1	NAG	C1-O5-C5	2.10	115.00	112.19
2	G	2	NAG	O4-C4-C3	2.09	115.30	110.38
2	E	1	NAG	O3-C3-C4	-2.09	105.46	110.38
2	C	2	NAG	C3-C4-C5	-2.07	106.47	110.23
2	H	1	NAG	O7-C7-C8	-2.07	118.37	122.05

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	G	1	NAG	C1
2	H	1	NAG	C1
3	D	3	LGU	C5

All (22) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	3	LGU	C4-C5-C6-O6B

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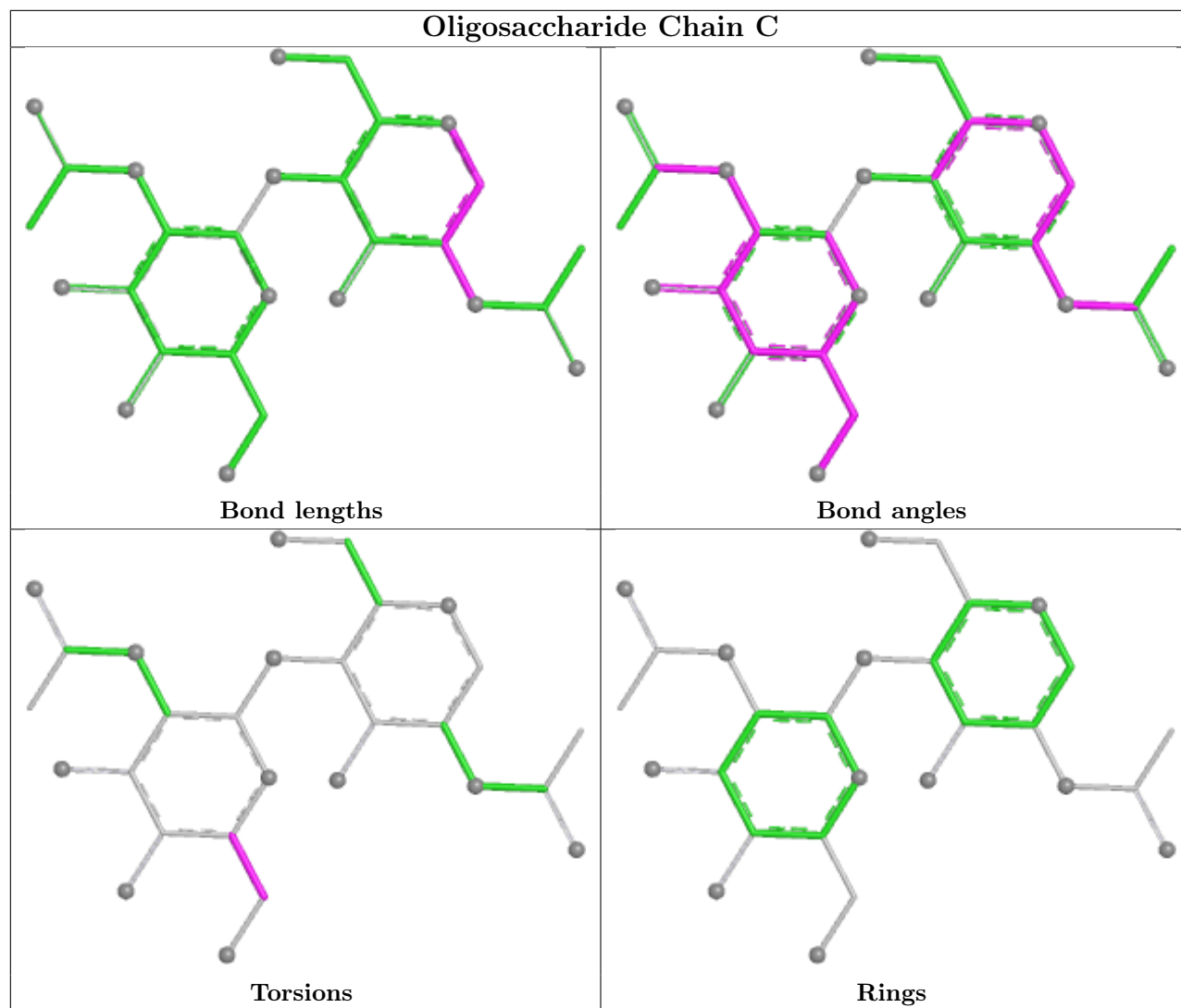
Mol	Chain	Res	Type	Atoms
2	E	2	NAG	O5-C5-C6-O6
2	E	2	NAG	C4-C5-C6-O6
3	D	3	LGU	O5-C5-C6-O6B
2	C	2	NAG	O5-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6
3	D	1	NAG	C4-C5-C6-O6
2	C	2	NAG	C4-C5-C6-O6
2	G	2	NAG	C4-C5-C6-O6
2	G	2	NAG	C8-C7-N2-C2
2	F	2	NAG	O5-C5-C6-O6
3	D	4	MAN	C4-C5-C6-O6
2	G	2	NAG	O5-C5-C6-O6
3	D	4	MAN	O5-C5-C6-O6
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	G	2	NAG	O7-C7-N2-C2
2	F	2	NAG	C8-C7-N2-C2
2	F	2	NAG	O7-C7-N2-C2
2	H	2	NAG	O5-C5-C6-O6
3	D	1	NAG	O5-C5-C6-O6
2	H	2	NAG	C4-C5-C6-O6

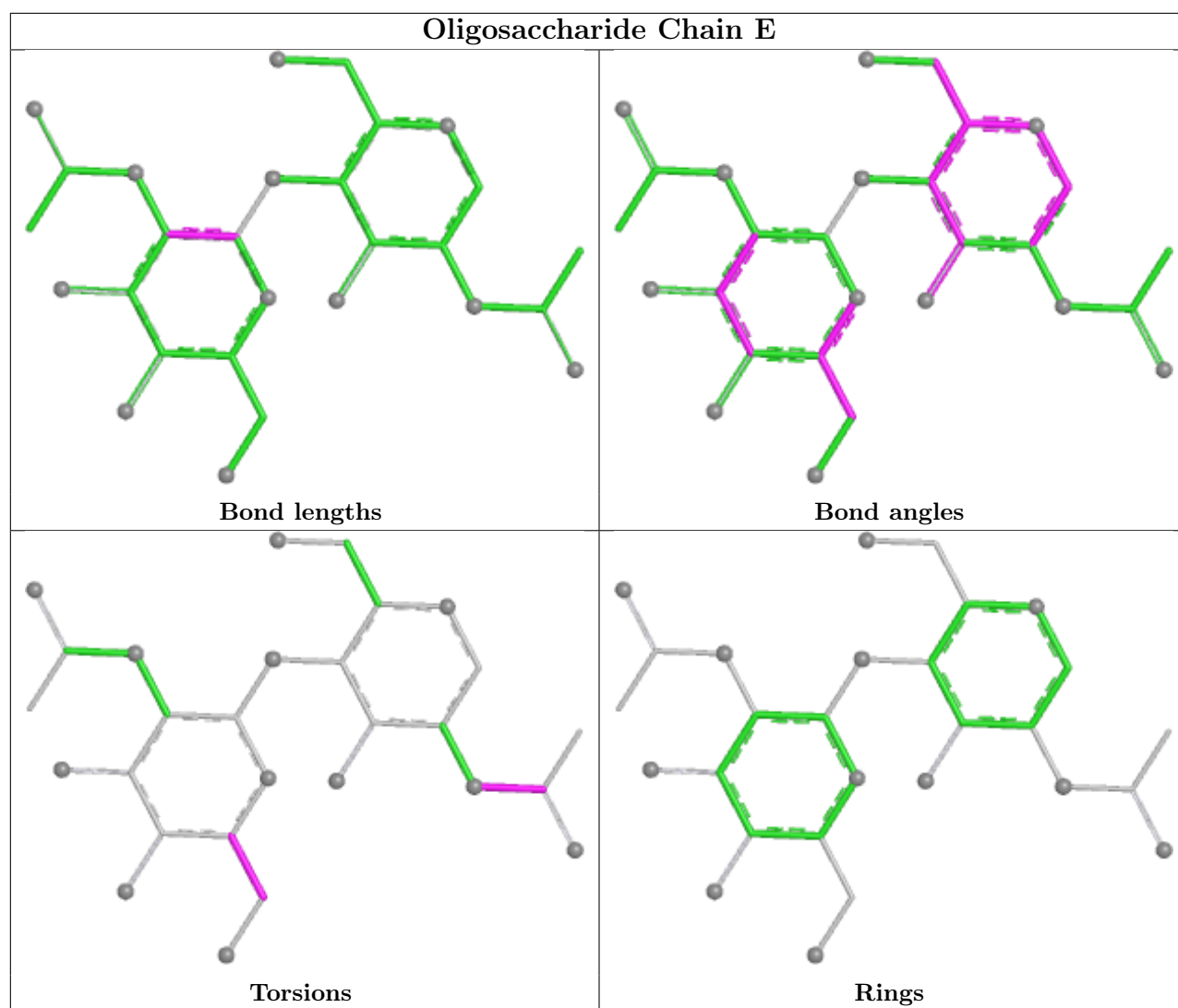
There are no ring outliers.

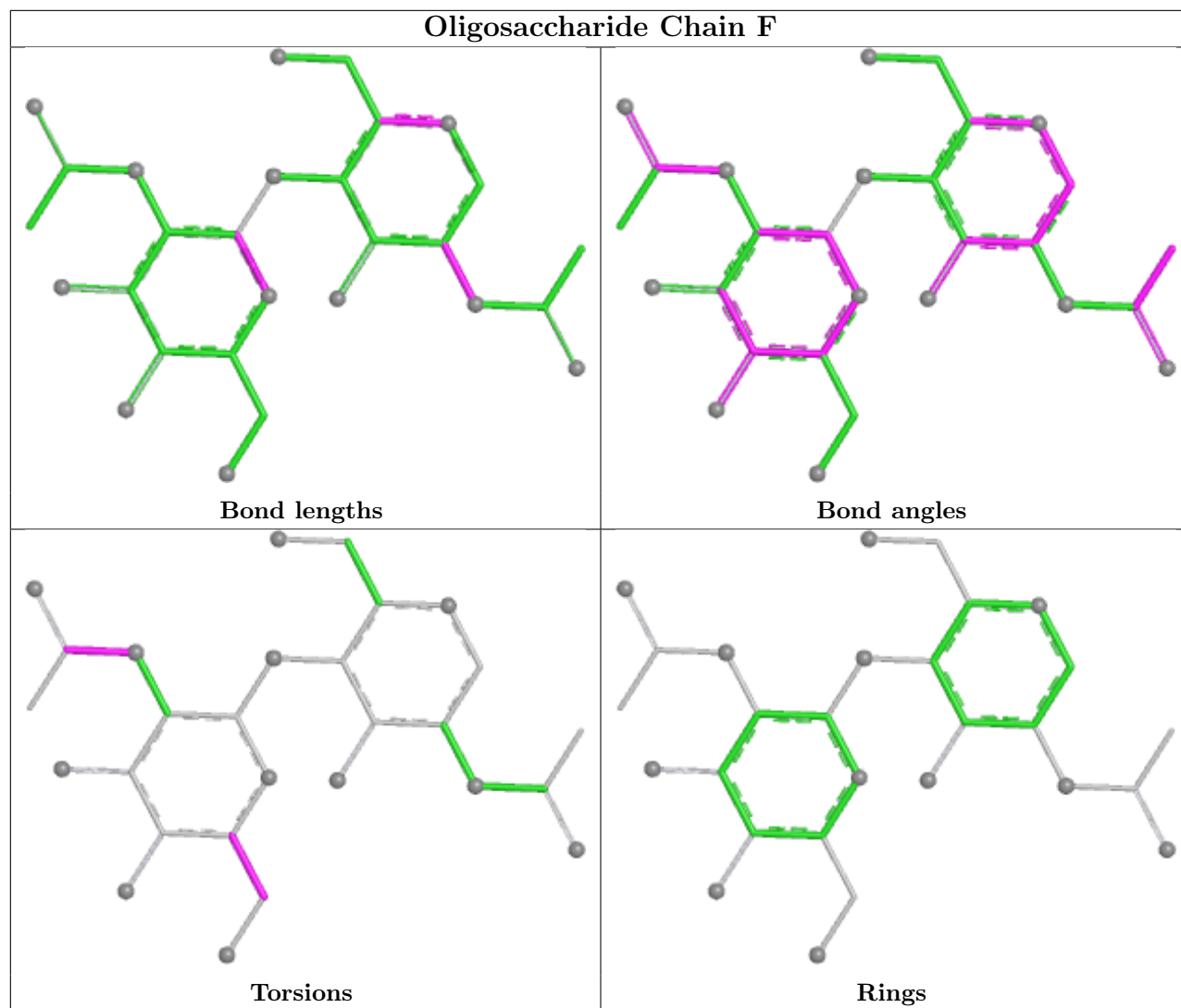
6 monomers are involved in 5 short contacts:

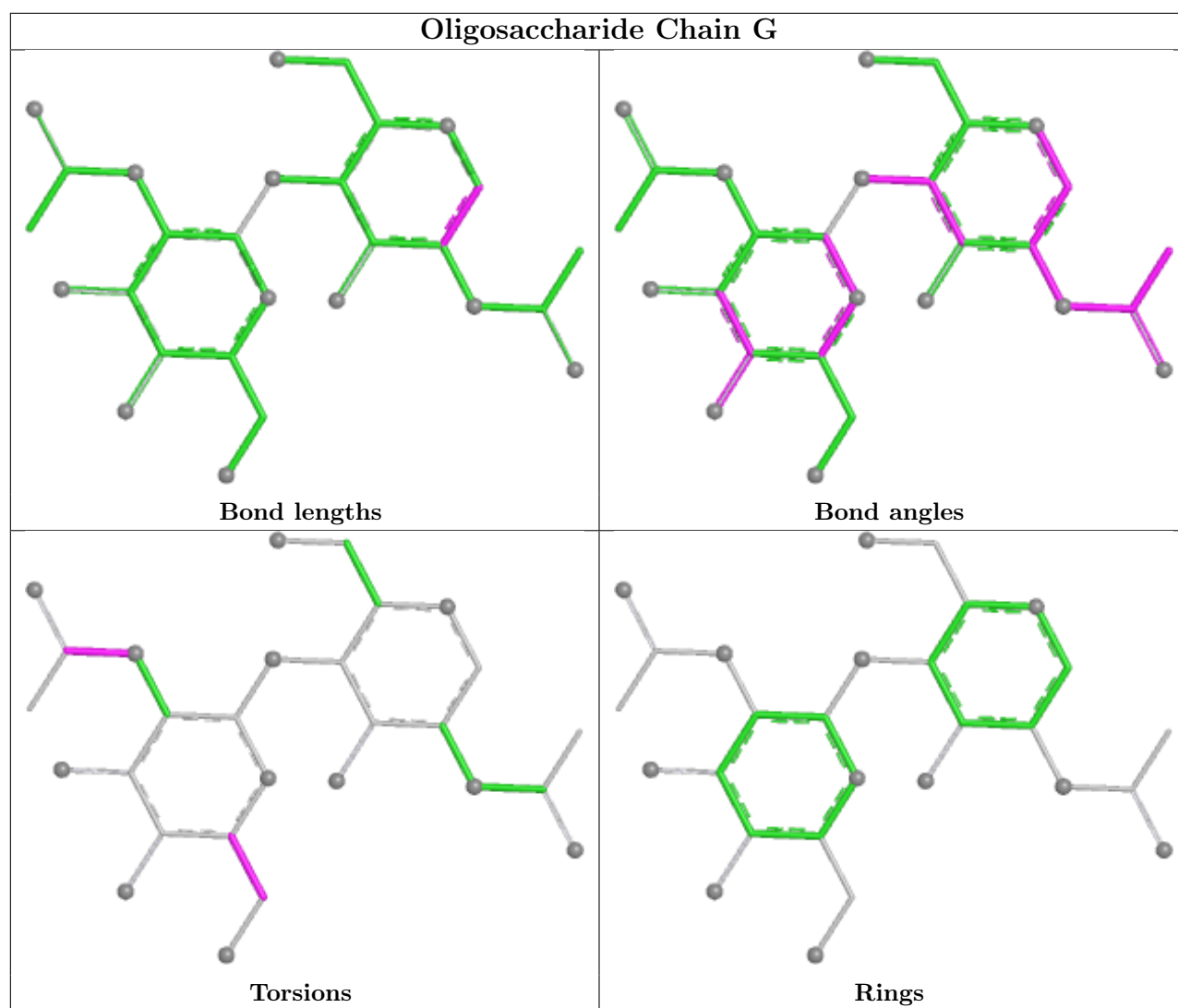
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	1	NAG	1	0
2	F	1	NAG	1	0
2	C	2	NAG	1	0
2	G	2	NAG	1	0
2	H	1	NAG	1	0
2	G	1	NAG	1	0

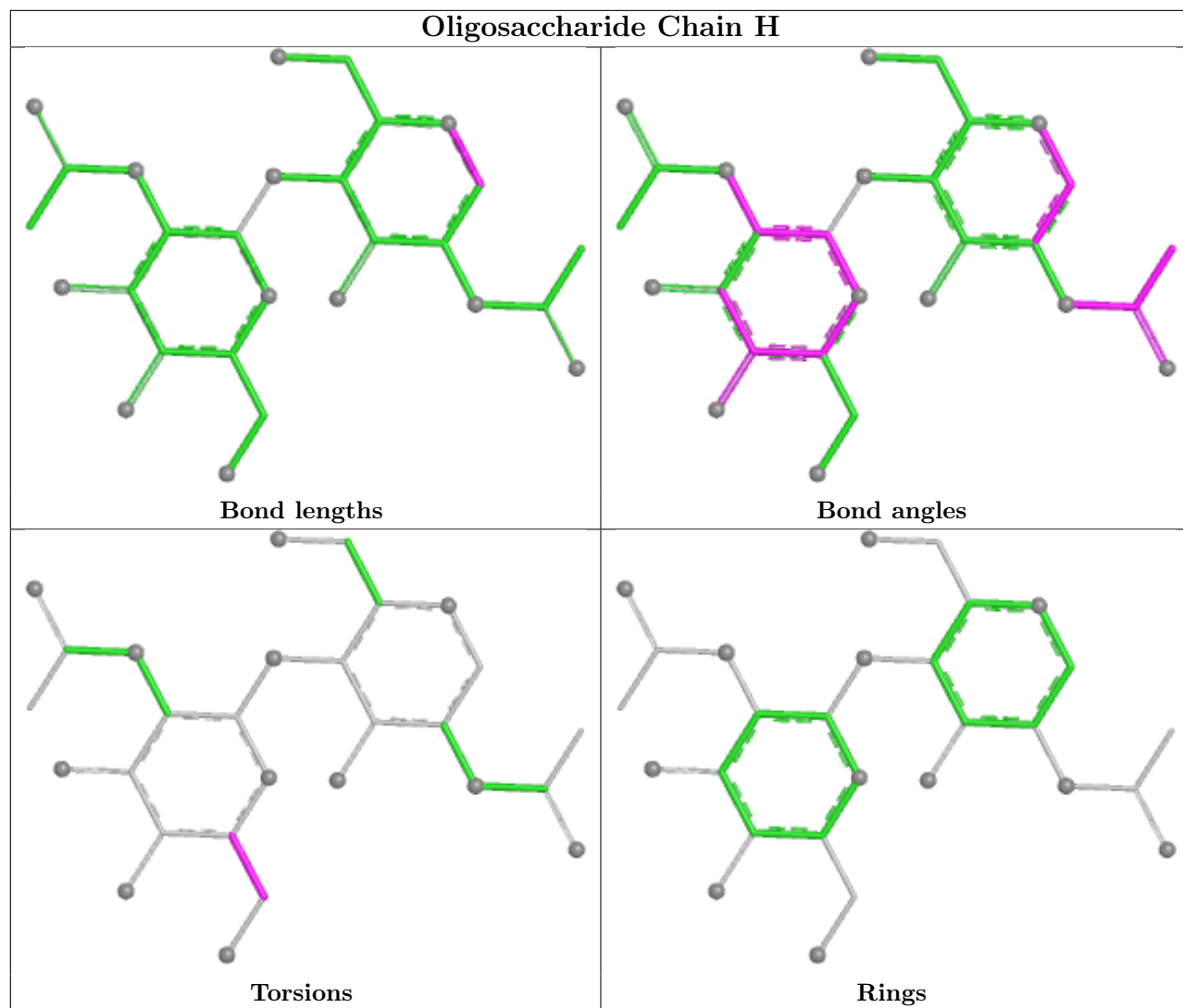
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

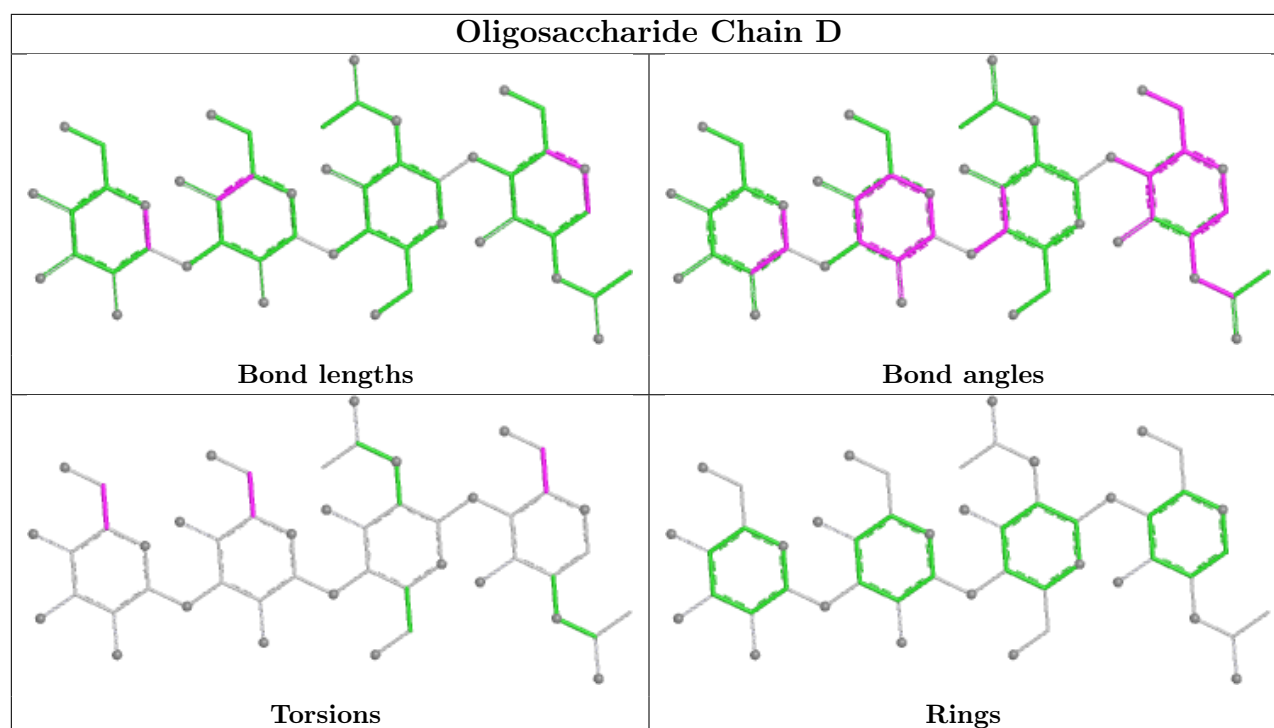












5.6 Ligand geometry [i](#)

5 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	801	1	14,14,15	1.13	1 (7%)	17,19,21	2.33	6 (35%)
5	PZF	B	802	-	32,32,32	1.48	8 (25%)	37,46,46	2.46	13 (35%)
4	NAG	A	801	1	14,14,15	0.91	0	17,19,21	1.54	4 (23%)
6	LGU	A	803	3	11,11,13	0.85	0	15,15,19	1.70	4 (26%)
5	PZF	A	802	-	32,32,32	0.98	2 (6%)	37,46,46	2.35	14 (37%)

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	801	1	-	2/6/23/26	0/1/1/1
5	PZF	B	802	-	-	0/14/26/26	0/4/4/4
4	NAG	A	801	1	-	4/6/23/26	0/1/1/1
6	LGU	A	803	3	1/1/4/6	2/2/19/24	0/1/1/1
5	PZF	A	802	-	-	0/14/26/26	0/4/4/4

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	802	PZF	C26-C21	-3.44	1.34	1.39
4	B	801	NAG	O5-C1	-3.41	1.38	1.43
5	B	802	PZF	C4-N3	2.66	1.36	1.32
5	B	802	PZF	F29-C24	-2.62	1.28	1.35
5	B	802	PZF	C8-N9	2.35	1.38	1.33
5	A	802	PZF	C2-N3	-2.31	1.29	1.34
5	B	802	PZF	C21-C13	-2.15	1.47	1.52
5	B	802	PZF	O18-C4	2.09	1.38	1.35
5	B	802	PZF	C11-C10	2.07	1.42	1.39
5	B	802	PZF	C14-N15	2.02	1.49	1.46
5	A	802	PZF	C13-C17	-2.01	1.49	1.54

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	802	PZF	C8-N7-C12	8.47	122.06	114.95
5	A	802	PZF	C8-N7-C12	6.13	120.09	114.95
5	A	802	PZF	C8-N9-C10	5.55	123.63	115.89
5	A	802	PZF	N9-C8-N7	-5.51	120.25	128.58
4	B	801	NAG	O4-C4-C5	-5.17	96.60	109.32
5	B	802	PZF	N9-C8-N7	-4.62	121.59	128.58
5	B	802	PZF	O18-C4-C5	4.53	124.78	116.77
4	B	801	NAG	C1-O5-C5	3.99	117.53	112.19
5	A	802	PZF	C11-C10-N9	-3.97	117.11	122.13
4	B	801	NAG	C3-C4-C5	3.91	117.32	110.23
5	B	802	PZF	C2-N3-C4	3.87	120.24	116.72
5	B	802	PZF	C11-C12-N7	-3.74	116.60	122.82
6	A	803	LGU	C1-O5-C5	-3.69	107.24	112.19
5	A	802	PZF	C14-N15-C16	3.59	116.84	111.64
5	B	802	PZF	C11-C10-N9	-3.30	117.96	122.13
5	B	802	PZF	C8-N9-C10	3.09	120.19	115.89
4	A	801	NAG	C1-O5-C5	3.05	116.28	112.19
5	A	802	PZF	C2-C1-C10	-3.02	116.83	121.29

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	803	LGU	O5-C1-C2	-2.94	103.78	110.79
5	B	802	PZF	C2-C1-C10	-2.89	117.01	121.29
5	B	802	PZF	O18-C4-N3	-2.78	114.61	119.68
5	A	802	PZF	C2-N3-C4	2.76	119.23	116.72
4	A	801	NAG	C1-C2-N2	2.70	114.68	110.43
4	B	801	NAG	O5-C5-C4	2.69	117.37	110.83
5	B	802	PZF	C5-C4-N3	-2.61	120.58	124.65
5	A	802	PZF	C14-N15-C12	-2.60	120.43	123.58
5	A	802	PZF	C23-C22-C21	-2.56	121.09	123.93
4	B	801	NAG	C6-C5-C4	-2.49	106.90	113.02
6	A	803	LGU	C2-C3-C4	2.39	115.07	110.86
6	A	803	LGU	C3-C4-C5	2.38	114.54	110.23
5	A	802	PZF	C11-C12-N7	-2.28	119.02	122.82
4	B	801	NAG	O5-C5-C6	2.27	112.08	107.66
5	A	802	PZF	C21-C26-C25	2.27	122.48	119.33
5	B	802	PZF	C11-C10-C1	2.25	124.87	121.82
4	A	801	NAG	C2-N2-C7	-2.23	119.91	122.90
5	A	802	PZF	N7-C12-N15	2.21	119.92	116.76
5	B	802	PZF	C21-C26-C25	2.19	122.37	119.33
5	A	802	PZF	F28-C25-C24	2.17	124.24	118.45
5	A	802	PZF	C26-C21-C22	2.09	118.17	116.42
5	B	802	PZF	C14-N15-C12	2.04	126.06	123.58
4	A	801	NAG	O7-C7-C8	-2.02	118.45	122.05

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
6	A	803	LGU	C5

All (8) torsion outliers are listed below:

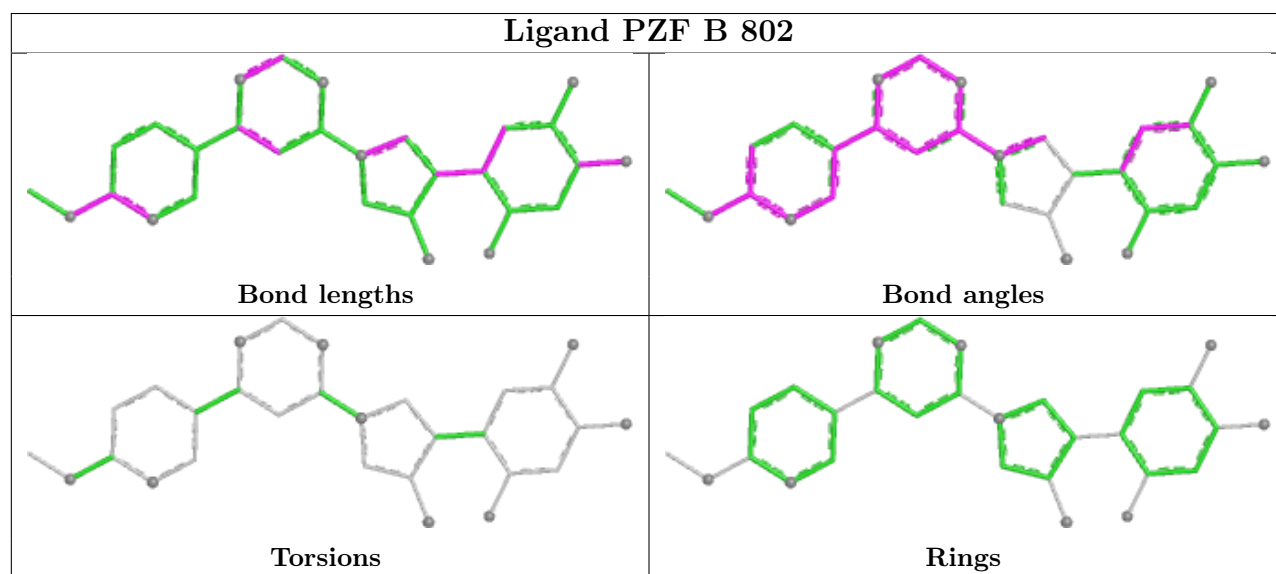
Mol	Chain	Res	Type	Atoms
4	A	801	NAG	C8-C7-N2-C2
4	A	801	NAG	O7-C7-N2-C2
6	A	803	LGU	O5-C5-C6-O6B
6	A	803	LGU	C4-C5-C6-O6B
4	B	801	NAG	C4-C5-C6-O6
4	B	801	NAG	O5-C5-C6-O6
4	A	801	NAG	C4-C5-C6-O6
4	A	801	NAG	O5-C5-C6-O6

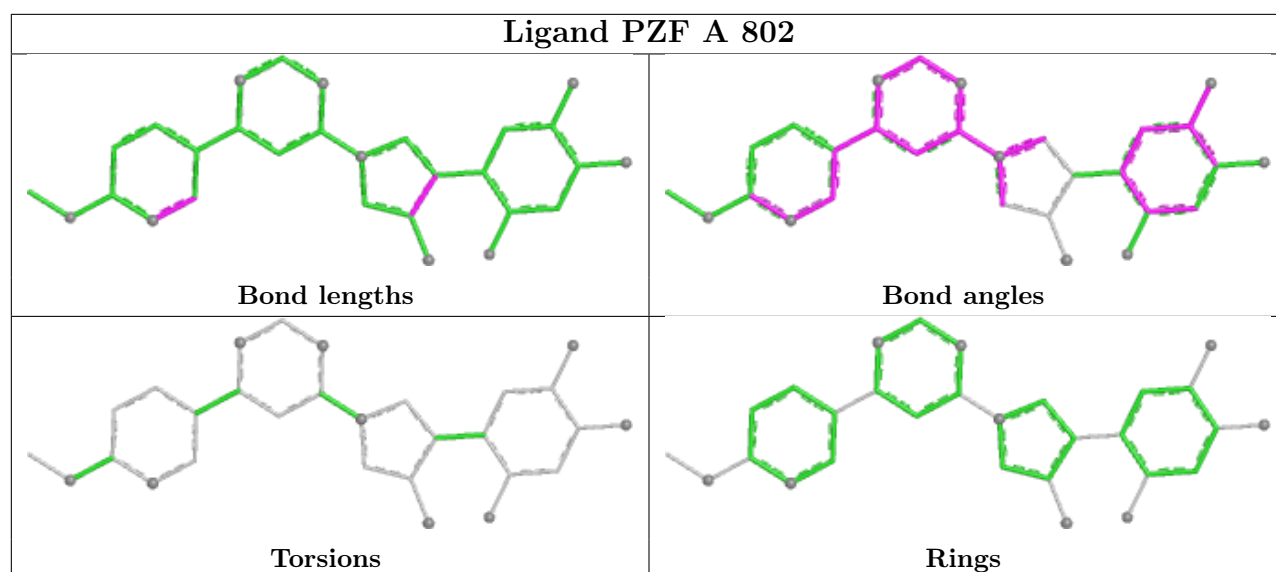
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	801	NAG	1	0
5	A	802	PZF	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

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	728/748 (97%)	0.85	25 (3%)	48 45	46, 55, 70, 79	0
1	B	728/748 (97%)	0.77	35 (4%)	35 32	44, 55, 71, 79	0
All	All	1456/1496 (97%)	0.81	60 (4%)	41 38	44, 55, 71, 79	0

All (60) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	39	SER	3.7
1	B	472	CYS	3.6
1	B	447	CYS	3.1
1	A	39	SER	2.9
1	B	92	ASN	2.9
1	A	279	VAL	2.8
1	B	350	THR	2.8
1	A	390	ASP	2.8
1	B	393	ASP	2.7
1	A	642	SER	2.7
1	A	722	ALA	2.6
1	A	280	THR	2.6
1	B	394	CYS	2.6
1	A	766	PRO	2.6
1	B	279	VAL	2.5
1	B	762	CYS	2.5
1	B	342	ALA	2.5
1	A	144	THR	2.4
1	A	764	SER	2.4
1	B	485	SER	2.4
1	A	645	GLY	2.4
1	B	111	GLY	2.3
1	A	424	GLY	2.3
1	B	328	CYS	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	296	GLY	2.3
1	B	74	ASN	2.2
1	A	72	GLN	2.2
1	A	636	THR	2.2
1	A	731	GLN	2.2
1	B	94	THR	2.2
1	B	278	SER	2.2
1	B	462	SER	2.2
1	A	633	GLY	2.2
1	B	672	GLY	2.2
1	B	614	SER	2.2
1	B	507	VAL	2.2
1	A	276	LEU	2.2
1	B	243	ASP	2.2
1	A	394	CYS	2.2
1	B	46	THR	2.2
1	A	530	LEU	2.1
1	A	763	PHE	2.1
1	B	66	HIS	2.1
1	B	70	TYR	2.1
1	B	389	ILE	2.1
1	B	541	PRO	2.1
1	A	684	ARG	2.1
1	B	444	CYS	2.1
1	A	485	SER	2.1
1	A	675	THR	2.1
1	B	101	SER	2.1
1	B	454	CYS	2.1
1	B	486	VAL	2.1
1	B	476	GLY	2.1
1	B	747	ALA	2.0
1	A	726	VAL	2.0
1	A	564	ALA	2.0
1	B	564	ALA	2.0
1	A	339	CYS	2.0
1	B	385	CYS	2.0

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

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

6.3 Carbohydrates

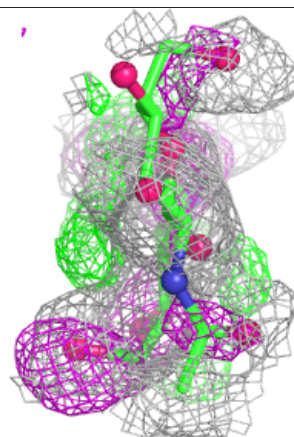
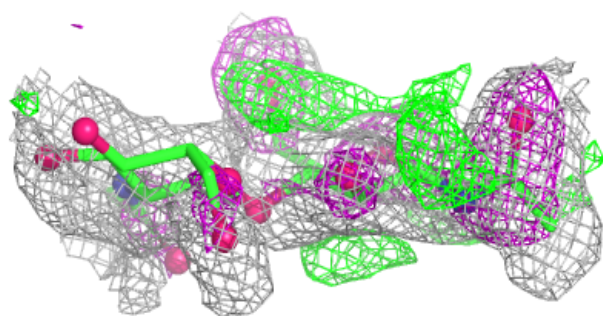
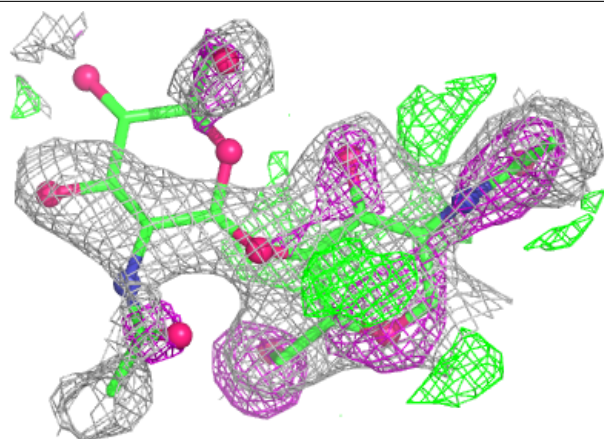
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists 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	C	2	14/15	0.61	0.19	75,83,85,86	0
2	NAG	E	2	14/15	0.62	0.15	99,102,104,105	0
3	LGU	D	3	11/13	0.69	0.19	88,93,97,98	0
2	NAG	G	2	14/15	0.71	0.16	87,91,94,95	0
2	NAG	H	2	14/15	0.77	0.17	96,100,102,103	0
2	NAG	G	1	14/15	0.78	0.17	79,85,92,93	0
3	MAN	D	4	11/12	0.79	0.16	89,91,93,96	0
2	NAG	E	1	14/15	0.80	0.15	81,87,90,94	0
2	NAG	F	2	14/15	0.81	0.18	63,68,78,80	0
2	NAG	H	1	14/15	0.84	0.15	75,84,87,91	0
3	NAG	D	2	14/15	0.84	0.18	65,76,79,83	0
3	NAG	D	1	14/15	0.86	0.18	51,59,64,69	0
2	NAG	C	1	14/15	0.89	0.20	44,48,53,63	0
2	NAG	F	1	14/15	0.90	0.17	42,47,56,61	0

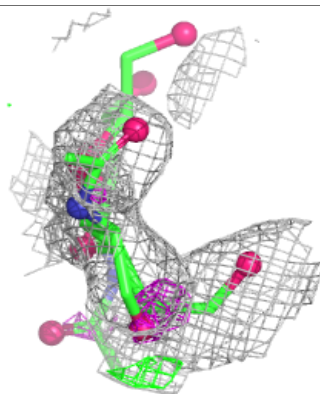
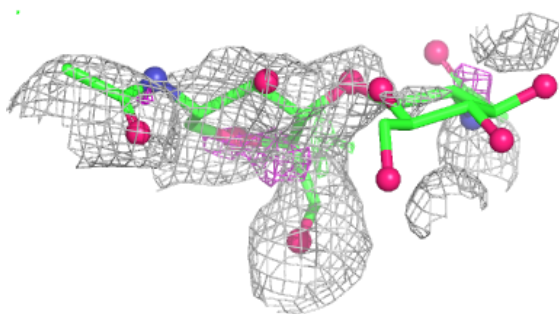
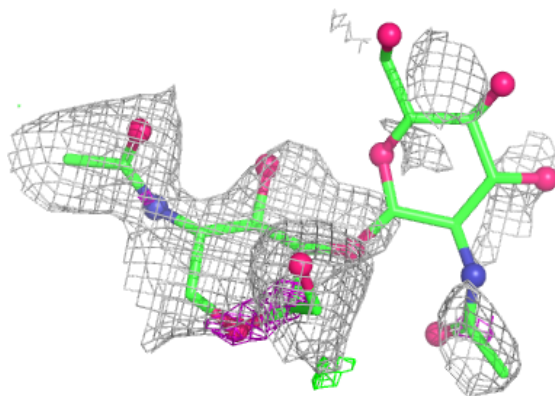
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain C:

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

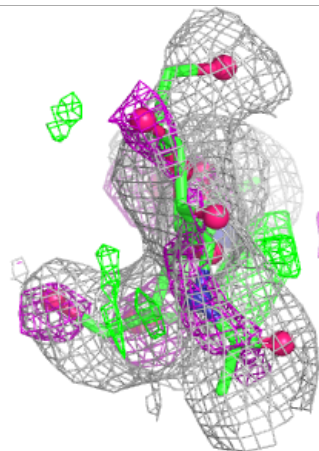
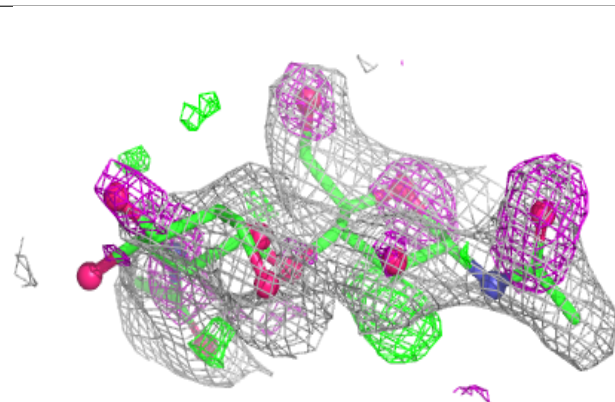
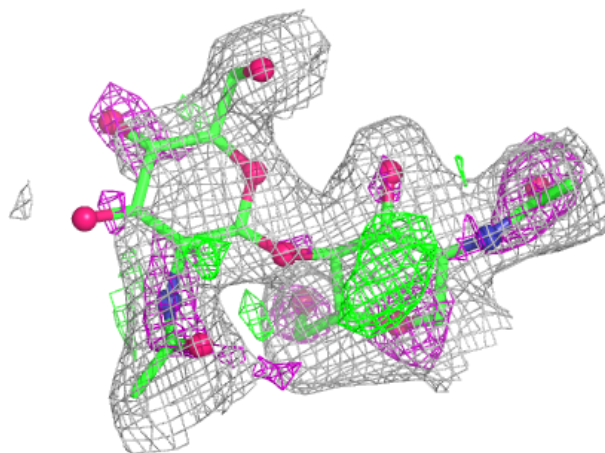
**Electron density around Chain E:**

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



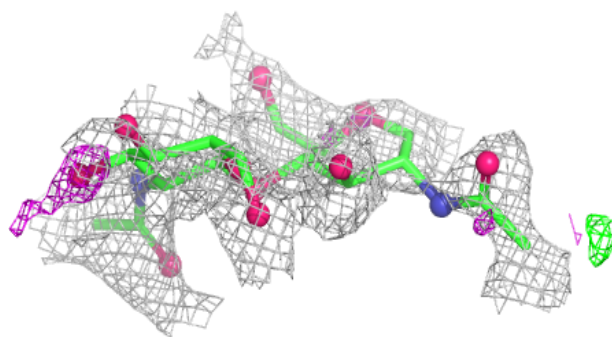
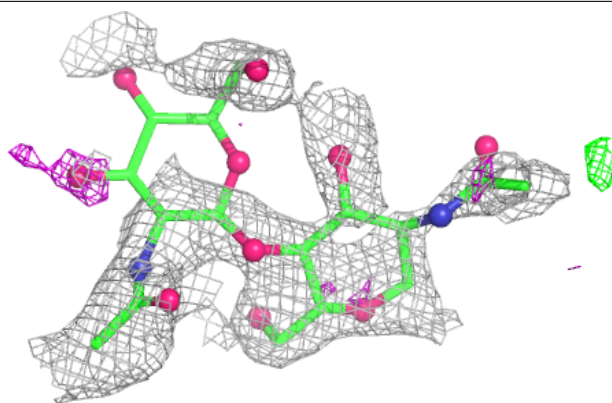
Electron density around Chain F:

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



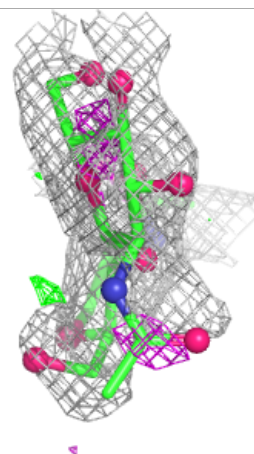
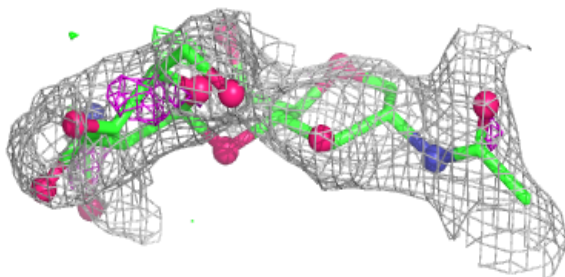
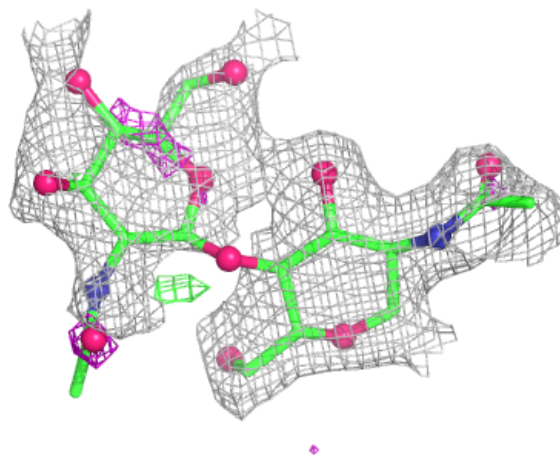
Electron density around Chain G:

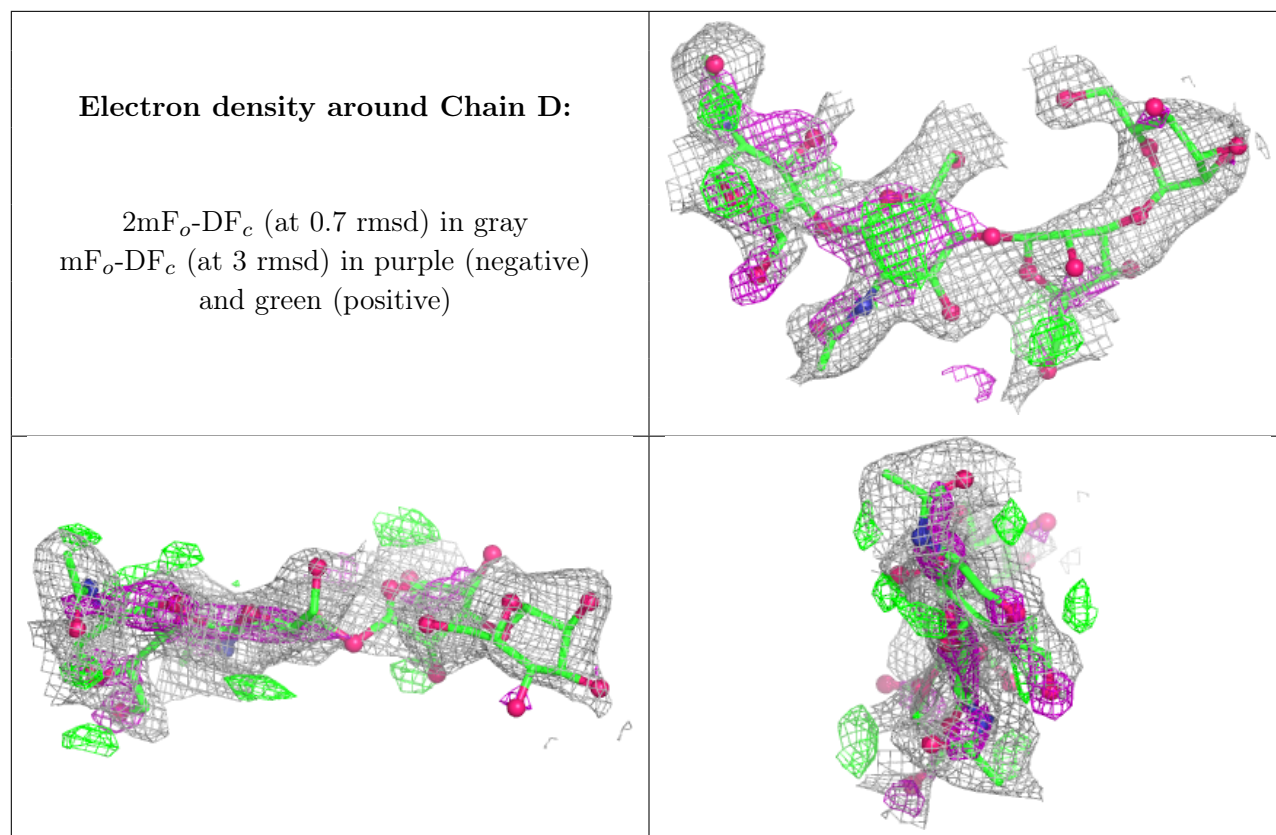
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



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](#)

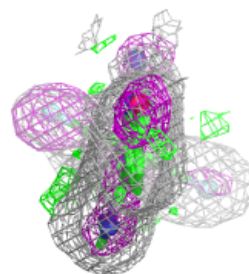
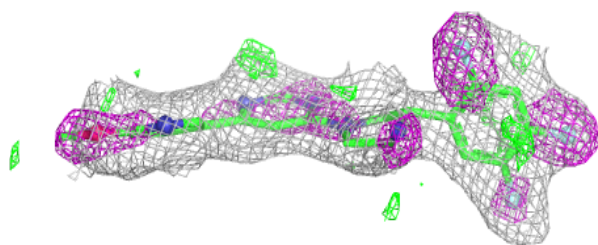
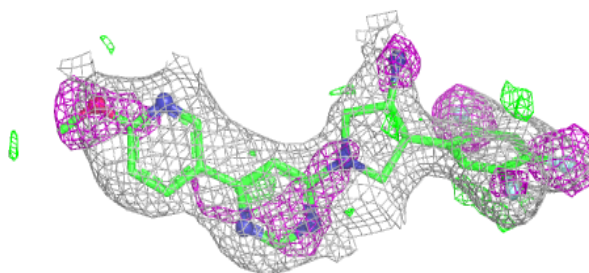
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	LGU	A	803	11/13	0.74	0.18	87,93,96,98	0
4	NAG	A	801	14/15	0.84	0.16	77,82,87,88	0
4	NAG	B	801	14/15	0.86	0.15	54,59,68,72	0
5	PZF	A	802	29/29	0.88	0.17	44,57,71,80	0
5	PZF	B	802	29/29	0.93	0.14	35,44,62,72	0

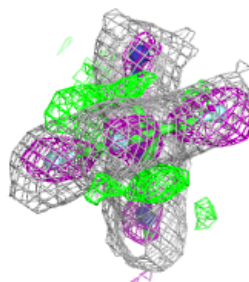
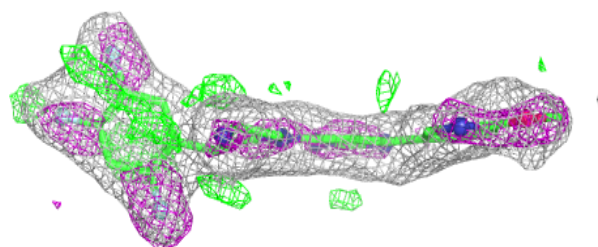
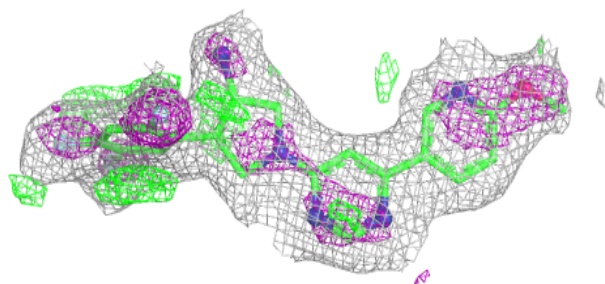
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around PZF A 802:

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

**Electron density around PZF B 802:**

$2mF_o-DF_c$ (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.