



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 05:50 PM UTC

PDB ID : 2C6F / pdb_00002c6f
Title : Structure of human somatic angiotensin-I converting enzyme N domain
Authors : Corradi, H.R.; Schwager, S.L.U.; Nichinda, A.; Sturrock, E.D.; Acharya, K.R.
Deposited on : 2005-11-09
Resolution : 3.01 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
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

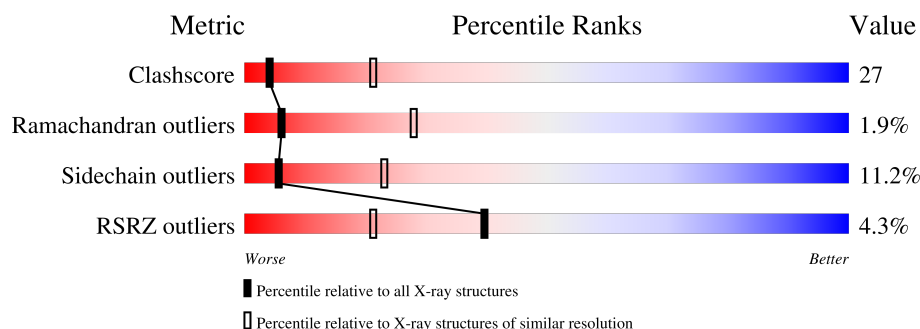
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.01 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))
Clashscore	190562	3444 (3.04-3.00)
Ramachandran outliers	187476	3319 (3.04-3.00)
Sidechain outliers	187428	3322 (3.04-3.00)
RSRZ outliers	180081	3130 (3.04-3.00)

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	612	3% 54% 37% 9%
1	B	612	5% 51% 40% 7%
2	C	2	50% 50%
2	D	2	50% 50%
2	E	2	50% 50%
2	F	2	50% 50%

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
6	GOL	A	707	-	X	-	-
6	GOL	B	706	-	X	-	-

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9970 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 ANGIOTENSIN-CONVERTING ENZYME, SOMATIC ISOFORM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	612	Total	C	N	O	S	6	0	0
			4884	3145	840	880	19			
1	B	611	Total	C	N	O	S	14	0	0
			4852	3124	830	879	19			

- 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	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			

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

- Molecule 4 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Zn	0	0
			1	1		
4	B	1	Total	Zn	0	0
			1	1		

- Molecule 5 is ACETATE ION (CCD ID: ACT) (formula: C₂H₃O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		
5	B	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total	C	O	0	0
			6	3	3		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 7 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	1	Total	Cl	0	0
			1	1		
7	B	1	Total	Cl	0	0
			1	1		

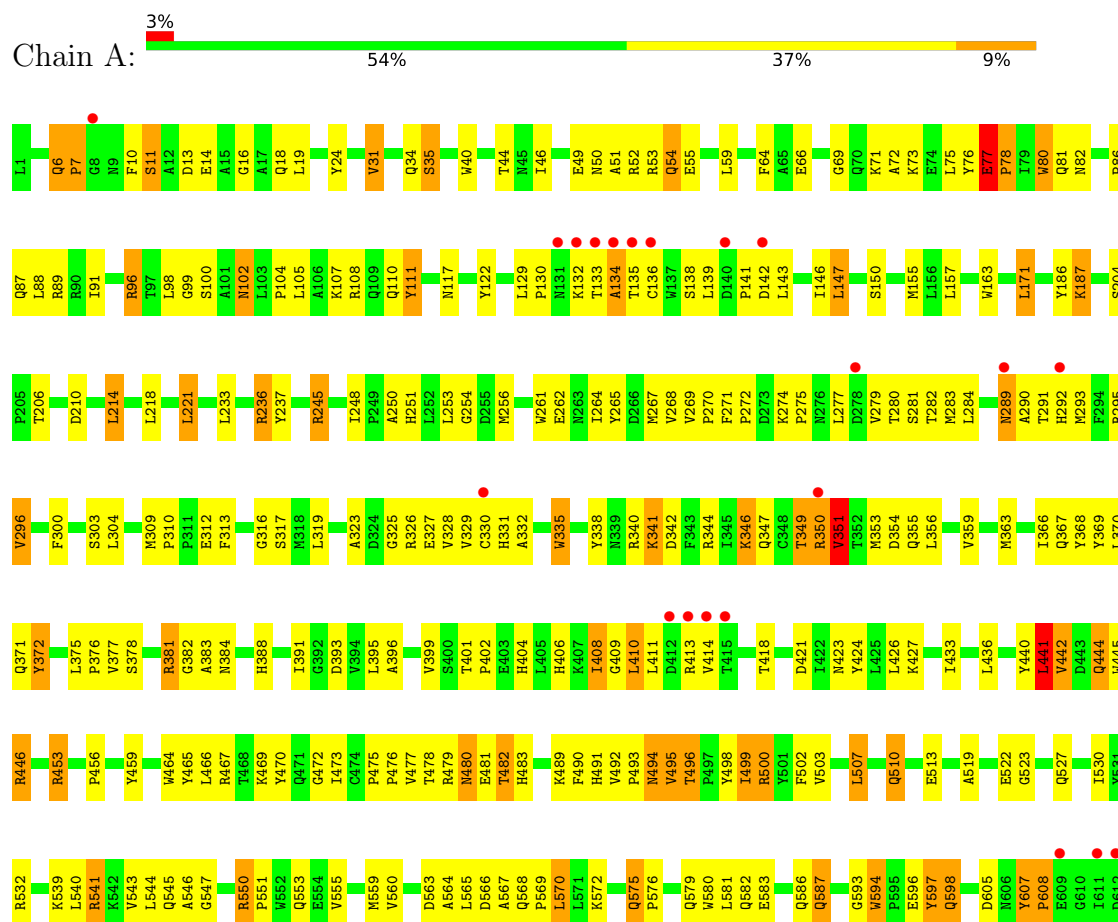
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	11	Total	O	0	0
			11	11		
8	B	13	Total	O	0	0
			13	13		

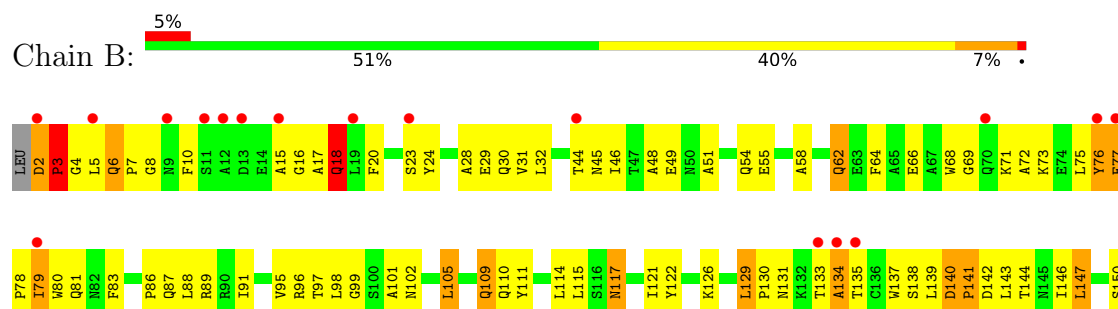
3 Residue-property plots

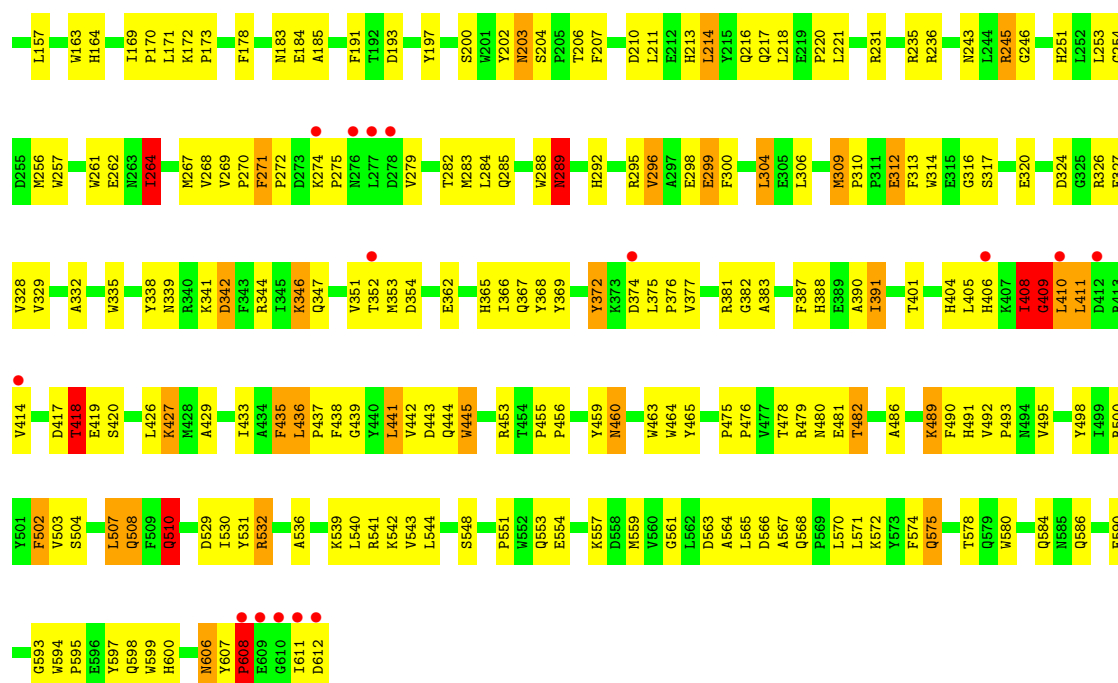
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: ANGIOTENSIN-CONVERTING ENZYME, SOMATIC ISOFORM



• Molecule 1: ANGIOTENSIN-CONVERTING ENZYME, SOMATIC ISOFORM





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

Chain C: 50% 50%



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

Chain D: 50% 50%



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

Chain E: 50% 50%



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

Chain F: 50% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	101.12Å 211.32Å 171.27Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.91 – 3.01 47.91 – 3.01	Depositor EDS
% Data completeness (in resolution range)	96.7 (47.91-3.01) 96.9 (47.91-3.01)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.33 (at 3.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.273 0.219 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	45.6	Xtriage
Anisotropy	0.335	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 49.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	9970	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.90% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, ZN, NAG, ACT, GOL

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	0.63	2/5042 (0.0%)	1.14	46/6885 (0.7%)
1	B	0.62	4/5010 (0.1%)	1.27	70/6845 (1.0%)
All	All	0.62	6/10052 (0.1%)	1.21	116/13730 (0.8%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	3	2
All	All	3	3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	117	ASN	CA-CB	-6.30	1.43	1.53
1	B	436	LEU	CA-CB	-5.73	1.48	1.54
1	B	575	GLN	CA-CB	-5.62	1.49	1.54
1	A	117	ASN	CB-CG	-5.43	1.38	1.52
1	B	264	ILE	CA-CB	-5.19	1.47	1.54
1	B	140	ASP	CG-OD2	5.18	1.35	1.25

All (116) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	3	PRO	CA-N-CD	-23.00	79.80	112.00
1	B	608	PRO	CA-N-CD	-19.25	85.05	112.00
1	A	598	GLN	OE1-CD-NE2	-10.43	112.17	122.60

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	527	GLN	OE1-CD-NE2	-10.28	112.33	122.60
1	B	410	LEU	N-CA-C	-10.26	100.08	112.59
1	B	444	GLN	OE1-CD-NE2	-9.90	112.70	122.60
1	B	109	GLN	OE1-CD-NE2	-9.84	112.76	122.60
1	B	18	GLN	OE1-CD-NE2	-9.76	112.84	122.60
1	B	6	GLN	OE1-CD-NE2	-9.66	112.94	122.60
1	B	62	GLN	OE1-CD-NE2	-9.54	113.06	122.60
1	B	110	GLN	OE1-CD-NE2	-9.49	113.11	122.60
1	B	30	GLN	OE1-CD-NE2	-9.42	113.18	122.60
1	B	87	GLN	OE1-CD-NE2	-9.36	113.24	122.60
1	B	54	GLN	OE1-CD-NE2	-9.32	113.28	122.60
1	B	418	THR	N-CA-C	-9.26	100.61	112.93
1	A	409	GLY	CA-C-N	-8.75	108.89	122.37
1	A	409	GLY	C-N-CA	-8.75	108.89	122.37
1	B	117	ASN	CA-CB-CG	-8.23	104.37	112.60
1	A	465	TYR	N-CA-C	-8.18	102.31	111.14
1	B	465	TYR	N-CA-C	-8.06	102.45	111.07
1	B	445	TRP	N-CA-C	-7.86	101.81	111.40
1	A	444	GLN	N-CA-C	-7.77	102.74	111.14
1	B	309	MET	CA-C-N	7.39	128.00	120.38
1	B	309	MET	C-N-CA	7.39	128.00	120.38
1	A	441	LEU	N-CA-C	7.28	119.84	111.11
1	A	605	ASP	N-CA-C	7.24	119.80	111.11
1	B	435	PHE	N-CA-C	7.24	118.81	111.07
1	B	30	GLN	CG-CD-NE2	7.14	127.11	116.40
1	B	482	THR	N-CA-C	-7.07	104.39	113.16
1	B	594	TRP	CA-C-N	7.07	127.06	119.28
1	B	594	TRP	C-N-CA	7.07	127.06	119.28
1	B	600	HIS	CA-C-N	-7.01	113.16	120.38
1	B	600	HIS	C-N-CA	-7.01	113.16	120.38
1	B	18	GLN	CG-CD-NE2	6.99	126.89	116.40
1	A	133	THR	N-CA-C	-6.97	105.20	113.21
1	A	349	THR	CA-C-N	-6.95	110.58	123.03
1	A	349	THR	C-N-CA	-6.95	110.58	123.03
1	A	598	GLN	CG-CD-NE2	6.91	126.77	116.40
1	A	564	ALA	N-CA-C	6.76	117.16	108.24
1	B	575	GLN	N-CA-C	6.73	122.65	113.77
1	B	87	GLN	CG-CD-NE2	6.70	126.45	116.40
1	B	455	PRO	CA-C-N	-6.70	112.73	119.56
1	B	455	PRO	C-N-CA	-6.70	112.73	119.56
1	B	271	PHE	CA-C-N	6.66	126.61	119.28
1	B	271	PHE	C-N-CA	6.66	126.61	119.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	290	ALA	N-CA-C	-6.64	104.66	112.89
1	B	109	GLN	CG-CD-NE2	6.59	126.28	116.40
1	B	48	ALA	N-CA-C	-6.58	105.83	114.31
1	B	6	GLN	CG-CD-NE2	6.57	126.25	116.40
1	B	54	GLN	CG-CD-NE2	6.54	126.21	116.40
1	B	110	GLN	CG-CD-NE2	6.51	126.17	116.40
1	A	527	GLN	CG-CD-NE2	6.41	126.02	116.40
1	B	444	GLN	CG-CD-NE2	6.34	125.91	116.40
1	A	519	ALA	N-CA-C	-6.33	104.43	111.71
1	A	296	VAL	N-CA-C	-6.32	104.34	110.72
1	B	62	GLN	CG-CD-NE2	6.32	125.88	116.40
1	A	349	THR	N-CA-C	6.30	118.85	109.59
1	A	575	GLN	N-CA-C	6.29	122.08	113.77
1	B	289	ASN	N-CA-C	6.27	119.76	110.48
1	B	2	ASP	CA-C-N	6.25	127.65	119.84
1	B	2	ASP	C-N-CA	6.25	127.65	119.84
1	B	306	LEU	N-CA-C	-6.24	99.73	109.72
1	B	296	VAL	N-CA-C	-6.23	104.97	111.58
1	B	409	GLY	N-CA-C	6.21	127.91	113.18
1	B	463	TRP	N-CA-C	-6.19	104.16	111.03
1	A	446	ARG	N-CA-C	6.14	117.64	111.07
1	B	599	TRP	N-CA-C	6.03	118.74	110.24
1	B	335	TRP	N-CA-C	6.02	118.76	109.07
1	B	510	GLN	N-CA-C	-6.01	104.81	111.36
1	A	502	PHE	N-CA-C	-6.01	104.36	111.03
1	A	384	ASN	N-CA-C	-5.89	102.92	108.22
1	B	436	LEU	N-CA-C	5.86	121.50	113.77
1	A	510	GLN	N-CA-C	-5.84	104.91	111.28
1	B	3	PRO	CA-CB-CG	-5.82	93.45	104.50
1	A	80	TRP	N-CA-C	-5.81	102.83	110.33
1	A	10	PHE	N-CA-C	5.79	117.98	109.24
1	B	164	HIS	N-CA-C	5.77	117.57	111.28
1	B	611	ILE	N-CA-C	5.74	116.88	107.24
1	B	375	LEU	N-CA-C	5.71	117.16	110.31
1	A	78	PRO	N-CA-C	-5.71	106.65	114.98
1	B	578	THR	N-CA-C	-5.64	105.13	111.28
1	B	564	ALA	N-CA-C	5.64	115.28	107.73
1	A	117	ASN	CB-CG-ND2	-5.63	107.95	116.40
1	A	102	ASN	CB-CA-C	-5.62	99.43	110.11
1	A	442	VAL	N-CA-C	5.60	116.05	110.23
1	A	523	GLY	N-CA-C	-5.58	100.96	112.34
1	A	110	GLN	N-CA-C	-5.54	105.14	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	304	LEU	N-CA-C	-5.50	106.42	113.02
1	B	46	ILE	N-CA-C	5.42	115.39	108.12
1	A	335	TRP	N-CA-C	5.41	117.78	109.07
1	B	3	PRO	CB-CA-C	5.41	120.48	111.56
1	B	110	GLN	N-CA-C	-5.41	105.30	111.14
1	B	502	PHE	N-CA-C	-5.38	105.06	111.03
1	A	134	ALA	CA-C-N	5.38	131.81	121.54
1	A	134	ALA	C-N-CA	5.38	131.81	121.54
1	A	135	THR	CA-C-N	5.36	129.93	120.87
1	A	135	THR	C-N-CA	5.36	129.93	120.87
1	B	532	ARG	N-CA-C	5.34	122.17	110.80
1	B	8	GLY	N-CA-C	-5.33	105.11	112.25
1	A	14	GLU	N-CA-C	-5.27	106.36	112.89
1	B	606	ASN	CA-C-N	-5.23	112.66	121.03
1	B	606	ASN	C-N-CA	-5.23	112.66	121.03
1	A	597	TYR	N-CA-C	-5.23	103.58	110.43
1	A	187	LYS	N-CA-C	-5.20	105.61	111.28
1	B	5	LEU	N-CA-C	5.17	120.51	113.37
1	A	328	VAL	N-CA-C	5.15	117.45	108.95
1	A	135	THR	N-CA-C	5.12	121.71	110.80
1	A	136	CYS	N-CA-C	5.12	117.14	110.43
1	A	351	VAL	N-CA-C	5.11	119.97	109.34
1	A	282	THR	N-CA-C	-5.11	106.90	113.23
1	B	401	THR	CA-C-N	5.09	124.88	119.28
1	B	401	THR	C-N-CA	5.09	124.88	119.28
1	B	491	HIS	N-CA-C	5.08	118.58	112.38
1	A	77	GLU	N-CA-C	5.06	121.00	109.81
1	A	330	CYS	N-CA-C	5.06	118.56	112.38
1	B	16	GLY	N-CA-C	-5.03	101.26	113.18

All (3) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	B	79	ILE	CA
1	B	410	LEU	CA
1	B	530	ILE	CB

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	607	TYR	Peptide
1	B	275	PRO	Peptide

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Mol	Chain	Res	Type	Group
1	B	409	GLY	Peptide

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	4884	0	4582	266	1
1	B	4852	0	4514	256	0
2	C	28	0	25	1	1
2	D	28	0	25	6	0
2	E	28	0	25	3	0
2	F	28	0	25	2	0
3	A	42	0	36	4	0
3	B	28	0	23	2	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	8	0	6	0	0
5	B	4	0	3	0	0
6	A	6	0	6	1	0
6	B	6	0	6	1	0
7	A	1	0	0	1	0
7	B	1	0	0	0	0
8	A	11	0	0	0	0
8	B	13	0	0	0	0
All	All	9970	0	9276	526	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 27.

All (526) 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:B:3:PRO:HD3	1:B:6:GLN:NE2	1.36	1.40
1:B:282:THR:HG21	1:B:410:LEU:CB	1.60	1.31
1:A:279:VAL:HG21	1:A:410:LEU:CD1	1.60	1.29
1:B:282:THR:CG2	1:B:410:LEU:CB	2.18	1.20

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:VAL:CG1	1:A:410:LEU:HD13	1.73	1.17
1:B:80:TRP:HA	1:B:83:PHE:CD2	1.78	1.17
1:A:279:VAL:HG21	1:A:410:LEU:HD12	1.17	1.13
1:B:80:TRP:O	1:B:89:ARG:HG2	1.48	1.12
1:A:206:THR:HG23	1:A:210:ASP:OD2	1.51	1.10
1:B:3:PRO:CD	1:B:6:GLN:NE2	2.13	1.10
1:A:279:VAL:HG11	1:A:410:LEU:HD13	1.09	1.09
3:B:702:NAG:H3	3:B:702:NAG:C8	1.48	1.09
1:A:139:LEU:HD22	1:A:163:TRP:CZ2	1.89	1.08
3:B:702:NAG:C8	3:B:702:NAG:C3	2.30	1.08
1:A:6:GLN:HG3	1:A:7:PRO:HD2	1.35	1.07
1:B:80:TRP:HA	1:B:83:PHE:CE2	1.88	1.07
1:A:6:GLN:HG3	1:A:7:PRO:CD	1.85	1.06
1:A:279:VAL:HG11	1:A:410:LEU:CD1	1.85	1.04
1:B:3:PRO:CD	1:B:6:GLN:HE22	1.70	1.03
1:A:340:ARG:NH2	2:D:2:NAG:HN2	1.56	1.03
1:A:279:VAL:CG2	1:A:410:LEU:CD1	2.37	1.01
1:A:279:VAL:CG2	1:A:410:LEU:HD12	1.91	1.00
1:A:279:VAL:HG13	1:A:410:LEU:HB2	1.49	0.95
1:A:340:ARG:HH22	2:D:2:NAG:HN2	1.09	0.95
1:B:482:THR:HG21	2:E:1:NAG:H82	1.49	0.93
1:A:72:ALA:O	1:A:76:TYR:HB2	1.71	0.91
1:B:129:LEU:HD23	1:B:612:ASP:CB	2.00	0.91
1:B:10:PHE:O	1:B:75:LEU:HD21	1.72	0.90
1:A:274:LYS:HB3	1:A:275:PRO:HD2	1.52	0.90
1:A:139:LEU:HD22	1:A:163:TRP:CH2	2.08	0.88
1:A:46:ILE:HG13	1:A:327:GLU:O	1.72	0.88
1:A:279:VAL:HG21	1:A:410:LEU:HD13	1.53	0.88
1:A:279:VAL:CB	1:A:410:LEU:HD13	2.04	0.87
1:B:29:GLU:OE1	2:F:1:NAG:H5	1.74	0.87
1:B:80:TRP:CE2	1:B:81:GLN:HG3	2.10	0.86
1:A:157:LEU:HG	1:A:607:TYR:OH	1.76	0.85
1:A:279:VAL:HG13	1:A:410:LEU:CB	2.07	0.84
1:B:353:MET:HE1	1:B:427:LYS:CE	2.07	0.84
1:A:408:ILE:HG12	1:A:408:ILE:O	1.77	0.84
1:A:279:VAL:CG2	1:A:410:LEU:HD13	2.04	0.84
1:A:270:PRO:HD3	1:A:426:LEU:HD22	1.60	0.84
1:B:3:PRO:HD3	1:B:6:GLN:HE22	0.76	0.83
1:A:566:ASP:OD2	1:A:568:GLN:HB2	1.79	0.83
1:A:441:LEU:HD23	1:A:442:VAL:N	1.94	0.82
1:B:464:TRP:CE3	1:B:475:PRO:HD3	2.16	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:353:MET:HE1	1:B:427:LYS:HE3	1.61	0.80
1:B:282:THR:HG21	1:B:410:LEU:CA	2.10	0.80
1:A:80:TRP:O	1:A:81:GLN:HB2	1.79	0.80
1:B:289:ASN:HB3	1:B:292:HIS:H	1.45	0.80
1:B:77:GLU:H	1:B:78:PRO:HD2	1.46	0.79
1:A:500:ARG:HG2	7:A:708:CL:CL	2.19	0.79
1:B:91:ILE:O	1:B:95:VAL:HG23	1.81	0.79
1:B:406:HIS:HA	1:B:411:LEU:O	1.82	0.79
1:A:597:TYR:O	1:A:598:GLN:HB3	1.83	0.79
1:B:2:ASP:CB	1:B:3:PRO:HD2	2.13	0.79
3:A:701:NAG:O3	3:A:701:NAG:C8	2.31	0.79
1:B:236:ARG:HE	1:B:267:MET:HE2	1.48	0.79
1:B:507:LEU:HA	1:B:510:GLN:HG3	1.66	0.78
1:B:129:LEU:CD2	1:B:612:ASP:CB	2.61	0.78
1:A:480:ASN:OD1	1:A:482:THR:HG23	1.84	0.78
1:A:132:LYS:C	1:A:134:ALA:H	1.93	0.77
1:A:329:VAL:O	1:A:346:LYS:HE3	1.83	0.77
1:B:441:LEU:HD23	1:B:442:VAL:N	1.99	0.77
1:A:129:LEU:HD12	1:A:129:LEU:N	1.99	0.76
1:A:372:TYR:HB2	1:A:375:LEU:HD12	1.68	0.76
1:B:295:ARG:HG2	1:B:314:TRP:CH2	2.20	0.76
1:A:221:LEU:HD23	1:A:433:ILE:HD12	1.67	0.76
1:B:147:LEU:HD22	1:B:256:MET:HA	1.68	0.76
1:B:78:PRO:O	1:B:79:ILE:CB	2.32	0.75
1:A:441:LEU:O	1:A:444:GLN:HB2	1.87	0.74
1:B:80:TRP:CA	1:B:83:PHE:CE2	2.70	0.74
1:B:77:GLU:H	1:B:78:PRO:CD	2.01	0.74
1:B:129:LEU:C	1:B:131:ASN:H	1.96	0.74
1:B:530:ILE:HG23	1:B:530:ILE:O	1.86	0.74
1:A:96:ARG:HH21	1:A:96:ARG:HG3	1.54	0.73
1:A:369:TYR:HA	1:A:372:TYR:CE1	2.24	0.73
1:B:3:PRO:HA	1:B:6:GLN:NE2	2.03	0.73
1:B:405:LEU:O	1:B:408:ILE:HG12	1.87	0.73
1:B:129:LEU:O	1:B:131:ASN:N	2.22	0.72
1:A:279:VAL:CG1	1:A:410:LEU:CD1	2.54	0.72
1:A:69:GLY:HA3	1:A:98:LEU:HD21	1.70	0.72
1:B:540:LEU:O	1:B:543:VAL:HG22	1.89	0.72
1:B:332:ALA:HB2	1:B:347:GLN:HG3	1.71	0.72
1:A:279:VAL:CG1	1:A:410:LEU:CB	2.68	0.71
1:A:279:VAL:CG1	1:A:410:LEU:HB2	2.19	0.71
1:B:79:ILE:O	1:B:80:TRP:CD1	2.43	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:236:ARG:NE	1:B:267:MET:CE	2.53	0.71
1:B:129:LEU:HD12	1:B:129:LEU:N	2.05	0.71
1:A:279:VAL:HG12	1:A:283:MET:HG3	1.71	0.71
1:B:482:THR:CG2	2:E:1:NAG:H82	2.20	0.71
1:A:6:GLN:HG3	1:A:7:PRO:HD3	1.74	0.70
1:B:329:VAL:O	1:B:346:LYS:HE3	1.90	0.70
1:A:292:HIS:O	1:A:296:VAL:HG23	1.92	0.70
1:A:456:PRO:HA	1:A:459:TYR:CD2	2.26	0.69
1:B:282:THR:HG22	1:B:410:LEU:CB	2.19	0.69
1:A:489:LYS:HE3	1:A:491:HIS:HD2	1.56	0.69
1:B:284:LEU:HD23	1:B:351:VAL:HG11	1.73	0.69
1:B:464:TRP:CZ3	1:B:475:PRO:HD3	2.28	0.69
1:B:105:LEU:O	1:B:109:GLN:HG3	1.93	0.69
1:A:291:THR:HB	3:A:703:NAG:H2	1.73	0.69
1:B:441:LEU:HD23	1:B:441:LEU:C	2.17	0.68
1:A:157:LEU:HD11	1:A:477:VAL:HG13	1.73	0.68
1:B:489:LYS:O	1:B:493:PRO:HD2	1.93	0.68
1:A:147:LEU:HD22	1:A:256:MET:HA	1.75	0.68
1:A:279:VAL:HG11	1:A:410:LEU:CG	2.23	0.68
1:A:86:PRO:HA	1:A:89:ARG:NH1	2.09	0.68
1:B:279:VAL:HG12	1:B:279:VAL:O	1.91	0.68
1:A:283:MET:HE1	1:A:356:LEU:HB2	1.76	0.67
1:A:489:LYS:O	1:A:493:PRO:HD2	1.94	0.67
1:B:129:LEU:HD21	1:B:137:TRP:CZ2	2.30	0.67
1:A:464:TRP:CE3	1:A:475:PRO:HD3	2.30	0.67
1:A:6:GLN:CG	1:A:7:PRO:HD2	2.19	0.67
1:B:203:ASN:N	1:B:203:ASN:HD22	1.93	0.66
1:B:129:LEU:C	1:B:131:ASN:N	2.53	0.66
1:B:264:ILE:O	1:B:264:ILE:HG23	1.94	0.66
1:B:171:LEU:HD21	1:B:493:PRO:HB3	1.77	0.66
1:B:58:ALA:O	1:B:62:GLN:HG3	1.95	0.66
1:B:157:LEU:HD13	1:B:476:PRO:HB2	1.77	0.66
1:A:206:THR:CG2	1:A:210:ASP:OD2	2.36	0.65
1:A:66:GLU:HA	1:A:98:LEU:HD22	1.78	0.65
1:B:29:GLU:HG2	1:B:338:TYR:O	1.97	0.65
1:A:440:TYR:O	1:A:444:GLN:HG2	1.97	0.65
1:A:11:SER:C	1:A:13:ASP:H	2.05	0.64
1:A:236:ARG:HG2	1:A:267:MET:HE1	1.79	0.64
1:A:395:LEU:O	1:A:399:VAL:HG23	1.98	0.64
1:B:316:GLY:HA3	1:B:344:ARG:HD3	1.79	0.64
1:B:221:LEU:HD12	1:B:433:ILE:HD13	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:264:ILE:O	1:B:264:ILE:CG2	2.46	0.64
1:A:340:ARG:NH2	2:D:2:NAG:N2	2.37	0.64
1:A:146:ILE:O	1:A:150:SER:HB3	1.98	0.64
1:A:73:LYS:HB2	1:A:77:GLU:HG3	1.78	0.64
1:B:316:GLY:C	1:B:344:ARG:HD3	2.22	0.64
1:B:183:ASN:HD21	1:B:193:ASP:HB2	1.64	0.63
1:A:332:ALA:HB2	1:A:347:GLN:HG3	1.80	0.63
1:B:607:TYR:CG	1:B:608:PRO:HD2	2.34	0.63
1:A:347:GLN:NE2	1:A:349:THR:OG1	2.31	0.63
1:A:31:VAL:HG21	1:A:64:PHE:CD1	2.33	0.63
1:A:236:ARG:NE	1:A:267:MET:HE3	2.14	0.63
1:A:541:ARG:O	1:A:541:ARG:HG3	1.99	0.63
1:A:251:HIS:CE1	1:A:476:PRO:HB3	2.34	0.62
1:B:203:ASN:HD22	1:B:203:ASN:H	1.45	0.62
1:A:340:ARG:NH1	2:D:2:NAG:H2	2.14	0.62
1:B:271:PHE:CD1	1:B:419:GLU:HB3	2.34	0.62
1:B:329:VAL:O	1:B:346:LYS:CE	2.48	0.62
1:A:283:MET:HE3	1:A:356:LEU:HD22	1.81	0.62
1:B:3:PRO:CD	1:B:6:GLN:HE21	2.10	0.62
1:B:10:PHE:O	1:B:75:LEU:CD2	2.47	0.62
1:A:206:THR:HG23	1:A:210:ASP:CG	2.23	0.62
1:B:312:GLU:HG3	1:B:341:LYS:O	2.00	0.62
1:A:157:LEU:HD13	1:A:476:PRO:HB2	1.81	0.61
1:B:324:ASP:OD1	1:B:326:ARG:HG2	2.00	0.61
1:B:3:PRO:CA	1:B:6:GLN:NE2	2.64	0.61
1:A:96:ARG:HG3	1:A:96:ARG:NH2	2.14	0.61
1:B:279:VAL:O	1:B:279:VAL:CG1	2.48	0.61
1:A:270:PRO:HD3	1:A:426:LEU:CD2	2.30	0.61
1:B:404:HIS:HB2	1:B:529:ASP:OD2	2.01	0.61
1:B:279:VAL:HG13	1:B:282:THR:HB	1.83	0.60
1:A:73:LYS:CB	1:A:77:GLU:HG3	2.31	0.60
1:B:86:PRO:HA	1:B:89:ARG:HE	1.66	0.60
1:A:147:LEU:HD22	1:A:256:MET:CA	2.32	0.59
1:B:146:ILE:O	1:B:150:SER:HB3	2.02	0.59
1:A:377:VAL:HG13	1:A:378:SER:N	2.18	0.59
1:A:513:GLU:OE2	1:A:572:LYS:HE3	2.02	0.59
1:B:441:LEU:CD2	1:B:442:VAL:N	2.64	0.59
1:B:530:ILE:O	1:B:530:ILE:CG2	2.51	0.59
1:B:142:ASP:O	1:B:146:ILE:HG13	2.02	0.59
1:A:300:PHE:HE2	1:A:304:LEU:HD11	1.68	0.59
1:B:353:MET:HE1	1:B:427:LYS:NZ	2.17	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:98:LEU:HB2	1:B:102:ASN:ND2	2.18	0.59
1:B:504:SER:O	1:B:508:GLN:HB3	2.03	0.59
1:B:7:PRO:HB2	1:B:71:LYS:HG2	1.85	0.58
1:A:248:ILE:O	1:A:473:ILE:HA	2.03	0.58
1:B:77:GLU:N	1:B:78:PRO:CD	2.66	0.58
1:A:236:ARG:HG2	1:A:267:MET:CE	2.33	0.58
1:B:2:ASP:CB	1:B:3:PRO:CD	2.78	0.58
1:B:139:LEU:HD22	1:B:163:TRP:CZ2	2.39	0.58
1:B:607:TYR:C	1:B:608:PRO:O	2.43	0.58
1:B:580:TRP:O	1:B:584:GLN:HG2	2.04	0.58
1:A:147:LEU:HD22	1:A:256:MET:N	2.19	0.57
1:A:261:TRP:CD1	1:A:261:TRP:N	2.71	0.57
1:B:218:LEU:CD2	1:B:570:LEU:HD23	2.34	0.57
1:A:75:LEU:H	1:A:75:LEU:HD23	1.69	0.57
1:A:489:LYS:O	1:A:491:HIS:N	2.38	0.57
1:A:338:TYR:HD2	1:A:377:VAL:HG23	1.68	0.57
1:A:597:TYR:O	1:A:598:GLN:CB	2.53	0.57
1:B:326:ARG:O	1:B:327:GLU:C	2.48	0.57
1:A:142:ASP:O	1:A:146:ILE:HG13	2.05	0.57
1:A:236:ARG:HE	1:A:267:MET:HE3	1.69	0.57
1:B:541:ARG:O	1:B:541:ARG:HG2	2.04	0.57
1:A:51:ALA:O	1:A:55:GLU:HG3	2.04	0.57
1:B:97:THR:O	1:B:97:THR:HG22	2.04	0.56
1:A:441:LEU:HD23	1:A:441:LEU:C	2.29	0.56
1:B:80:TRP:CG	1:B:81:GLN:N	2.73	0.56
1:A:406:HIS:HD2	1:A:413:ARG:O	1.89	0.56
1:A:446:ARG:NH2	1:A:496:THR:O	2.35	0.56
1:A:404:HIS:O	1:A:408:ILE:HG23	2.05	0.56
1:A:75:LEU:HD23	1:A:75:LEU:N	2.21	0.56
1:A:129:LEU:HD12	1:A:129:LEU:H	1.71	0.55
1:A:77:GLU:N	1:A:78:PRO:HD2	2.21	0.55
1:A:98:LEU:HB2	1:A:102:ASN:HD21	1.71	0.55
1:B:310:PRO:O	1:B:313:PHE:HB3	2.06	0.55
1:A:323:ALA:C	1:A:325:GLY:H	2.14	0.55
1:A:261:TRP:O	1:A:264:ILE:HG12	2.07	0.55
1:A:367:GLN:O	1:A:371:GLN:HG2	2.05	0.55
1:B:213:HIS:O	1:B:216:GLN:HB2	2.07	0.55
1:B:366:ILE:O	1:B:369:TYR:N	2.39	0.55
1:A:129:LEU:N	1:A:129:LEU:CD1	2.67	0.55
1:A:469:LYS:HE2	1:A:470:TYR:CZ	2.42	0.55
1:A:277:LEU:CD1	1:A:424:TYR:HA	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:PRO:O	1:B:272:PRO:HD3	2.08	0.54
1:B:279:VAL:HG21	1:B:405:LEU:HD13	1.89	0.54
1:B:270:PRO:HD3	1:B:426:LEU:HD22	1.89	0.54
1:B:141:PRO:HG2	1:B:142:ASP:N	2.22	0.54
1:A:91:ILE:HG23	1:A:378:SER:OG	2.06	0.54
1:A:268:VAL:O	1:A:269:VAL:C	2.51	0.54
1:A:104:PRO:HG2	1:A:107:LYS:HD2	1.88	0.54
1:B:316:GLY:CA	1:B:344:ARG:HD3	2.38	0.54
1:B:17:ALA:O	1:B:20:PHE:HB3	2.08	0.53
1:B:64:PHE:CE1	1:B:68:TRP:NE1	2.75	0.53
1:B:326:ARG:O	1:B:328:VAL:HG13	2.09	0.53
1:B:478:THR:O	1:B:479:ARG:HD2	2.08	0.53
1:A:492:VAL:N	1:A:493:PRO:HD2	2.23	0.53
1:B:251:HIS:CE1	1:B:476:PRO:HB3	2.43	0.53
1:A:147:LEU:O	1:A:254:GLY:HA2	2.09	0.53
1:B:99:GLY:O	1:B:102:ASN:ND2	2.36	0.53
1:B:147:LEU:O	1:B:254:GLY:HA2	2.08	0.53
1:A:54:GLN:HG3	1:A:55:GLU:N	2.24	0.53
1:A:132:LYS:C	1:A:134:ALA:N	2.63	0.52
1:A:326:ARG:HD2	3:A:701:NAG:HN2	1.72	0.52
1:A:607:TYR:CD1	1:A:608:PRO:HA	2.44	0.52
1:B:218:LEU:O	1:B:221:LEU:HB2	2.09	0.52
1:A:350:ARG:HD2	1:A:351:VAL:H	1.75	0.52
1:B:29:GLU:OE1	2:F:1:NAG:C5	2.54	0.52
1:B:29:GLU:CG	1:B:338:TYR:O	2.57	0.52
1:A:214:LEU:HD11	1:A:565:LEU:HB3	1.90	0.52
1:A:510:GLN:OE1	1:A:566:ASP:N	2.40	0.52
1:B:141:PRO:O	1:B:142:ASP:C	2.51	0.52
1:A:480:ASN:OD1	1:A:482:THR:CG2	2.57	0.52
1:B:426:LEU:O	1:B:429:ALA:HB3	2.10	0.52
1:A:98:LEU:HB2	1:A:102:ASN:ND2	2.24	0.52
1:A:69:GLY:HA3	1:A:98:LEU:CD2	2.38	0.52
1:B:292:HIS:O	1:B:296:VAL:HG23	2.09	0.52
1:A:597:TYR:CD1	1:A:597:TYR:C	2.88	0.51
1:A:147:LEU:CD2	1:A:256:MET:HA	2.41	0.51
1:A:353:MET:HE1	1:A:427:LYS:HE3	1.92	0.51
1:A:489:LYS:O	1:A:490:PHE:C	2.53	0.51
1:B:133:THR:O	1:B:134:ALA:C	2.53	0.51
1:B:408:ILE:CG1	1:B:409:GLY:N	2.70	0.51
1:B:597:TYR:CD1	1:B:597:TYR:C	2.89	0.51
1:A:66:GLU:HA	1:A:98:LEU:CD2	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:269:VAL:O	1:A:269:VAL:HG13	2.09	0.51
1:A:312:GLU:HG2	1:A:341:LYS:HE2	1.93	0.51
1:B:71:LYS:HE3	1:B:75:LEU:HD11	1.91	0.51
1:A:204:SER:OG	1:A:206:THR:HG22	2.11	0.51
1:A:495:VAL:HG12	1:A:495:VAL:O	2.10	0.51
1:A:204:SER:C	1:A:206:THR:H	2.18	0.50
1:B:140:ASP:OD1	1:B:140:ASP:C	2.53	0.50
1:A:445:TRP:HB2	1:A:466:LEU:HD12	1.94	0.50
1:B:3:PRO:N	1:B:6:GLN:HE21	2.10	0.50
1:B:200:SER:O	1:B:203:ASN:ND2	2.44	0.50
1:A:583:GLU:O	1:A:587:GLN:HG3	2.12	0.50
1:B:510:GLN:OE1	1:B:566:ASP:N	2.44	0.49
1:B:563:ASP:OD1	1:B:563:ASP:O	2.30	0.49
1:A:40:TRP:CE2	1:A:44:THR:HG21	2.46	0.49
1:A:139:LEU:CD2	1:A:163:TRP:CZ2	2.79	0.49
1:A:478:THR:O	1:A:479:ARG:HD2	2.13	0.49
1:B:3:PRO:O	1:B:4:GLY:C	2.52	0.49
1:B:138:SER:O	1:B:143:LEU:HG	2.11	0.49
1:B:362:GLU:O	1:B:365:HIS:HB2	2.13	0.49
1:A:265:TYR:CZ	1:A:269:VAL:HG23	2.47	0.49
1:A:583:GLU:O	1:A:587:GLN:CG	2.61	0.49
1:A:218:LEU:HD13	1:A:436:LEU:HD13	1.95	0.49
1:B:2:ASP:O	1:B:3:PRO:HB2	2.12	0.49
1:B:129:LEU:N	1:B:129:LEU:CD1	2.76	0.49
1:B:332:ALA:CB	1:B:347:GLN:HG3	2.39	0.49
1:B:256:MET:HB3	1:B:257:TRP:CE3	2.48	0.48
1:A:499:ILE:HG13	1:A:499:ILE:O	2.14	0.48
1:B:3:PRO:CA	1:B:6:GLN:HE21	2.26	0.48
1:B:593:GLY:HA3	6:B:706:GOL:O1	2.13	0.48
1:A:271:PHE:HB2	1:A:423:ASN:HD21	1.78	0.48
1:A:393:ASP:O	1:A:396:ALA:HB3	2.13	0.48
1:B:79:ILE:O	1:B:81:GLN:N	2.41	0.48
1:A:270:PRO:O	1:A:272:PRO:HD3	2.13	0.48
1:A:300:PHE:CE2	1:A:304:LEU:HD11	2.48	0.48
1:A:49:GLU:HG3	1:A:53:ARG:HE	1.79	0.48
1:A:99:GLY:HA2	1:A:186:TYR:CE2	2.48	0.48
1:B:169:ILE:N	1:B:170:PRO:HD2	2.29	0.48
1:A:293:MET:HE3	1:A:356:LEU:HA	1.96	0.48
1:A:399:VAL:HA	1:A:404:HIS:CD2	2.48	0.48
1:A:551:PRO:O	1:A:555:VAL:HG23	2.14	0.48
1:B:51:ALA:O	1:B:55:GLU:HG3	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:171:LEU:HD21	1:A:493:PRO:HB3	1.95	0.48
1:A:441:LEU:HD23	1:A:442:VAL:CA	2.43	0.48
1:A:507:LEU:HB3	1:A:565:LEU:CD2	2.44	0.48
1:B:489:LYS:O	1:B:490:PHE:C	2.56	0.48
1:B:295:ARG:HG2	1:B:314:TRP:HH2	1.72	0.48
1:B:482:THR:HG21	2:E:1:NAG:C8	2.34	0.48
1:A:139:LEU:HA	1:A:143:LEU:HB2	1.95	0.47
1:A:453:ARG:O	1:A:453:ARG:HG2	2.14	0.47
1:A:539:LYS:HE3	1:A:559:MET:O	2.13	0.47
1:B:44:THR:O	1:B:328:VAL:HG12	2.14	0.47
1:A:274:LYS:HB3	1:A:275:PRO:CD	2.34	0.47
1:A:498:TYR:C	1:A:500:ARG:H	2.22	0.47
1:B:567:ALA:O	1:B:568:GLN:C	2.57	0.47
1:B:606:ASN:O	1:B:608:PRO:O	2.32	0.47
1:A:98:LEU:HD12	1:A:102:ASN:OD1	2.13	0.47
1:B:80:TRP:CZ2	1:B:81:GLN:HG3	2.48	0.47
1:B:77:GLU:CG	1:B:77:GLU:O	2.62	0.47
1:A:141:PRO:O	1:A:142:ASP:C	2.57	0.47
1:A:355:GLN:O	1:A:359:VAL:HG23	2.15	0.47
1:A:482:THR:HG21	2:C:1:NAG:HN2	1.80	0.47
1:B:243:ASN:OD1	1:B:245:ARG:N	2.41	0.47
1:A:274:LYS:CB	1:A:275:PRO:HD2	2.35	0.47
1:A:316:GLY:C	1:A:344:ARG:HD3	2.39	0.47
1:B:98:LEU:HB2	1:B:102:ASN:HD21	1.78	0.47
1:A:490:PHE:C	1:A:490:PHE:CD2	2.92	0.47
1:B:218:LEU:HD22	1:B:570:LEU:HD23	1.95	0.47
1:B:390:ALA:O	1:B:391:ILE:C	2.57	0.47
1:A:122:TYR:CD1	1:A:493:PRO:HG3	2.49	0.47
1:A:277:LEU:HD11	1:A:424:TYR:HB2	1.97	0.47
1:B:236:ARG:HE	1:B:267:MET:CE	2.11	0.47
1:A:490:PHE:O	1:A:493:PRO:HG2	2.15	0.47
1:B:551:PRO:HB2	1:B:554:GLU:HG3	1.96	0.47
1:A:34:GLN:HG3	1:A:35:SER:H	1.80	0.46
1:A:284:LEU:HD23	1:A:351:VAL:HG11	1.97	0.46
1:B:262:GLU:OE1	1:B:262:GLU:N	2.48	0.46
1:B:270:PRO:HD3	1:B:426:LEU:CD2	2.44	0.46
1:B:329:VAL:HB	1:B:346:LYS:HE3	1.97	0.46
1:B:418:THR:HG22	1:B:419:GLU:N	2.29	0.46
1:B:495:VAL:HG12	1:B:495:VAL:O	2.14	0.46
1:A:129:LEU:H	1:A:129:LEU:CD1	2.27	0.46
1:B:283:MET:HE2	1:B:288:TRP:CE3	2.50	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:486:ALA:O	1:B:492:VAL:HG21	2.16	0.46
1:A:77:GLU:N	1:A:78:PRO:CD	2.78	0.46
1:A:456:PRO:HA	1:A:459:TYR:CE2	2.50	0.46
1:B:217:GLN:O	1:B:220:PRO:HD2	2.15	0.46
1:A:80:TRP:CE2	1:A:81:GLN:HG3	2.50	0.46
1:B:75:LEU:HD23	1:B:76:TYR:CZ	2.51	0.46
1:B:492:VAL:HB	1:B:493:PRO:CD	2.46	0.46
1:A:80:TRP:O	1:A:81:GLN:CB	2.57	0.46
1:A:279:VAL:HG11	1:A:410:LEU:HD22	1.98	0.46
1:B:147:LEU:HD13	1:B:256:MET:HE2	1.98	0.46
1:B:382:GLY:O	1:B:383:ALA:C	2.58	0.46
1:B:408:ILE:HG12	1:B:409:GLY:N	2.30	0.46
1:B:437:PRO:O	1:B:441:LEU:HB3	2.15	0.46
1:B:529:ASP:OD1	1:B:531:TYR:HB2	2.16	0.46
1:A:143:LEU:O	1:A:147:LEU:HD12	2.15	0.46
1:A:338:TYR:HD2	1:A:377:VAL:CG2	2.29	0.46
1:A:418:THR:O	1:A:421:ASP:N	2.49	0.46
1:B:204:SER:C	1:B:206:THR:H	2.23	0.46
1:B:268:VAL:O	1:B:269:VAL:C	2.57	0.46
1:A:157:LEU:HG	1:A:607:TYR:HH	1.80	0.46
1:A:510:GLN:HE21	1:A:560:VAL:HG11	1.81	0.46
1:A:340:ARG:CZ	2:D:2:NAG:HN2	2.21	0.46
1:A:375:LEU:O	1:A:376:PRO:C	2.58	0.46
1:B:147:LEU:HD22	1:B:256:MET:CA	2.44	0.46
1:B:607:TYR:CD2	1:B:608:PRO:HD2	2.50	0.46
1:A:138:SER:O	1:A:139:LEU:C	2.59	0.45
1:A:586:GLN:O	1:A:587:GLN:C	2.58	0.45
1:B:44:THR:O	1:B:44:THR:HG22	2.15	0.45
1:B:45:ASN:OD1	1:B:326:ARG:HB2	2.17	0.45
1:B:435:PHE:CE2	1:B:439:GLY:HA3	2.51	0.45
1:A:11:SER:C	1:A:13:ASP:N	2.73	0.45
1:A:481:GLU:OE2	1:A:481:GLU:HA	2.16	0.45
1:A:507:LEU:O	1:A:507:LEU:HG	2.16	0.45
1:B:141:PRO:CG	1:B:142:ASP:N	2.79	0.45
1:B:507:LEU:O	1:B:507:LEU:HG	2.16	0.45
1:B:129:LEU:HD12	1:B:129:LEU:H	1.78	0.45
1:B:536:ALA:O	1:B:539:LYS:N	2.50	0.45
1:B:309:MET:HB3	1:B:313:PHE:CD2	2.52	0.45
1:B:406:HIS:O	1:B:408:ILE:O	2.34	0.45
1:B:98:LEU:HB3	1:B:101:ALA:HB3	1.99	0.45
1:B:191:PHE:CZ	1:B:197:TYR:HD1	2.35	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:438:PHE:O	1:B:442:VAL:HG23	2.17	0.45
1:B:570:LEU:CD1	1:B:574:PHE:HE1	2.30	0.45
1:B:607:TYR:CD2	1:B:608:PRO:CD	3.00	0.45
1:A:369:TYR:HA	1:A:372:TYR:HE1	1.79	0.45
1:A:469:LYS:HE2	1:A:470:TYR:CE1	2.51	0.45
1:A:570:LEU:O	1:A:570:LEU:HD12	2.17	0.45
1:B:32:LEU:CD1	1:B:377:VAL:HG11	2.46	0.45
1:B:441:LEU:C	1:B:441:LEU:CD2	2.89	0.45
1:B:539:LYS:HE2	1:B:559:MET:O	2.16	0.45
1:B:568:GLN:HA	1:B:571:LEU:HD12	1.98	0.45
2:D:1:NAG:H61	2:D:2:NAG:C7	2.47	0.45
1:B:71:LYS:HE3	1:B:75:LEU:CD1	2.47	0.45
1:B:289:ASN:HB2	1:B:292:HIS:HB2	1.99	0.45
1:A:31:VAL:O	1:A:34:GLN:HG3	2.17	0.45
1:A:289:ASN:HB3	1:A:292:HIS:H	1.82	0.45
1:A:377:VAL:CG1	1:A:378:SER:N	2.80	0.45
1:B:71:LYS:O	1:B:72:ALA:C	2.59	0.45
1:B:77:GLU:N	1:B:78:PRO:HD2	2.23	0.44
1:A:279:VAL:O	1:A:281:SER:N	2.50	0.44
1:A:575:GLN:HB3	1:A:576:PRO:HD3	1.99	0.44
1:B:111:TYR:CE2	1:B:115:LEU:HD11	2.53	0.44
1:B:218:LEU:HD21	1:B:570:LEU:HD23	1.99	0.44
1:A:480:ASN:OD1	1:A:483:HIS:CE1	2.71	0.44
1:B:44:THR:C	1:B:328:VAL:HG12	2.42	0.44
1:B:417:ASP:HB3	1:B:420:SER:OG	2.16	0.44
1:B:86:PRO:HB3	1:B:89:ARG:HH21	1.81	0.44
1:A:18:GLN:HG2	1:A:88:LEU:HD22	2.00	0.44
1:A:300:PHE:CD2	1:A:395:LEU:HD23	2.53	0.44
1:B:172:LYS:N	1:B:173:PRO:HD2	2.32	0.44
1:A:77:GLU:H	1:A:78:PRO:HD2	1.82	0.44
1:B:80:TRP:NE1	1:B:81:GLN:HG3	2.32	0.44
1:B:117:ASN:O	1:B:121:ILE:HG13	2.18	0.44
1:A:218:LEU:O	1:A:221:LEU:HB2	2.18	0.44
1:A:279:VAL:HG11	1:A:410:LEU:CD2	2.47	0.44
1:A:44:THR:O	1:A:326:ARG:NH2	2.47	0.44
1:A:221:LEU:HD23	1:A:433:ILE:CD1	2.42	0.44
1:B:69:GLY:HA3	1:B:98:LEU:CD2	2.47	0.44
1:B:171:LEU:HD21	1:B:493:PRO:CB	2.47	0.44
1:B:183:ASN:HD21	1:B:193:ASP:CB	2.30	0.44
1:B:288:TRP:CH2	1:B:296:VAL:HG21	2.51	0.44
1:B:383:ALA:HB3	1:B:387:PHE:CG	2.53	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:LEU:HD22	1:A:237:TYR:HE1	1.83	0.43
1:A:269:VAL:O	1:A:269:VAL:CG1	2.65	0.43
1:A:593:GLY:HA3	6:A:707:GOL:O1	2.18	0.43
1:B:75:LEU:HD23	1:B:76:TYR:OH	2.17	0.43
1:A:24:TYR:CE1	1:A:64:PHE:HE2	2.35	0.43
1:A:283:MET:CE	1:A:356:LEU:HD22	2.47	0.43
1:A:494:ASN:N	1:A:494:ASN:HD22	2.15	0.43
1:B:15:ALA:HA	1:B:18:GLN:HB2	2.00	0.43
1:B:207:PHE:CZ	1:B:211:LEU:HD11	2.53	0.43
1:B:507:LEU:HB3	1:B:565:LEU:CD2	2.49	0.43
1:A:71:LYS:O	1:A:75:LEU:HD23	2.17	0.43
1:A:441:LEU:CD1	1:A:467:ARG:HA	2.48	0.43
1:B:122:TYR:CE1	1:B:163:TRP:HZ2	2.36	0.43
1:B:298:GLU:O	1:B:299:GLU:C	2.62	0.43
1:A:382:GLY:O	1:A:383:ALA:C	2.61	0.43
1:B:352:THR:HG22	1:B:353:MET:N	2.32	0.43
1:B:366:ILE:O	1:B:367:GLN:C	2.62	0.43
1:B:406:HIS:C	1:B:408:ILE:N	2.76	0.43
1:A:147:LEU:CD1	1:A:256:MET:HE2	2.48	0.43
1:A:464:TRP:CZ3	1:A:475:PRO:HD3	2.52	0.43
1:A:580:TRP:CZ3	1:A:581:LEU:HD23	2.54	0.43
1:B:245:ARG:O	1:B:595:PRO:HD2	2.19	0.43
1:B:317:SER:HA	1:B:344:ARG:HB3	2.01	0.43
1:B:442:VAL:O	1:B:445:TRP:HB3	2.19	0.43
1:A:86:PRO:O	1:A:87:GLN:C	2.61	0.43
1:A:366:ILE:O	1:A:370:LEU:HG	2.18	0.43
1:A:503:VAL:HG13	1:A:565:LEU:HD11	2.00	0.43
1:B:83:PHE:CD2	1:B:88:LEU:HD12	2.53	0.43
1:A:245:ARG:O	1:A:596:GLU:HB2	2.18	0.43
1:A:510:GLN:HG2	1:A:569:PRO:HG3	2.00	0.43
1:B:236:ARG:NE	1:B:267:MET:HE2	2.16	0.43
1:A:472:GLY:CA	1:A:594:TRP:CE2	3.02	0.42
1:A:594:TRP:CE3	1:A:594:TRP:N	2.87	0.42
1:A:381:ARG:O	1:A:547:GLY:HA2	2.20	0.42
1:B:502:PHE:O	1:B:503:VAL:C	2.62	0.42
1:A:368:TYR:CE2	1:A:544:LEU:HA	2.54	0.42
1:A:494:ASN:O	1:A:495:VAL:C	2.60	0.42
1:B:115:LEU:HD21	1:B:178:PHE:CE1	2.54	0.42
1:B:140:ASP:HA	1:B:141:PRO:HA	1.86	0.42
1:B:304:LEU:HD21	1:B:530:ILE:O	2.19	0.42
1:A:31:VAL:HG21	1:A:64:PHE:HD1	1.79	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:ILE:HA	1:A:369:TYR:CD1	2.54	0.42
1:A:472:GLY:HA2	1:A:594:TRP:CE2	2.53	0.42
1:B:372:TYR:CD1	1:B:372:TYR:C	2.97	0.42
1:B:299:GLU:O	1:B:300:PHE:C	2.62	0.42
1:B:405:LEU:O	1:B:408:ILE:CG1	2.63	0.42
1:B:460:ASN:N	1:B:481:GLU:OE2	2.34	0.42
1:A:143:LEU:HD13	1:A:163:TRP:HB2	2.01	0.42
1:A:204:SER:C	1:A:206:THR:N	2.77	0.42
1:A:607:TYR:HA	1:A:608:PRO:HA	1.76	0.42
1:B:570:LEU:CD1	1:B:574:PHE:CE1	3.03	0.42
1:A:335:TRP:CD1	1:A:335:TRP:N	2.87	0.42
1:A:510:GLN:HE21	1:A:510:GLN:HB3	1.63	0.42
1:B:83:PHE:HB2	1:B:89:ARG:HG3	2.02	0.42
1:B:436:LEU:N	1:B:437:PRO:HD2	2.35	0.42
1:A:210:ASP:O	1:A:214:LEU:HB2	2.19	0.42
1:A:478:THR:HG22	1:A:479:ARG:N	2.34	0.42
1:A:596:GLU:C	1:A:597:TYR:O	2.61	0.42
1:B:362:GLU:OE1	1:B:362:GLU:HA	2.19	0.42
1:A:108:ARG:O	1:A:111:TYR:HB3	2.20	0.41
1:A:312:GLU:OE1	1:A:312:GLU:N	2.46	0.41
1:B:372:TYR:C	1:B:372:TYR:HD1	2.28	0.41
1:A:309:MET:HA	1:A:310:PRO:HD3	1.93	0.41
1:A:567:ALA:O	1:A:570:LEU:HB3	2.20	0.41
1:B:406:HIS:CA	1:B:411:LEU:O	2.61	0.41
1:B:531:TYR:O	1:B:532:ARG:HB2	2.20	0.41
1:B:23:SER:O	1:B:24:TYR:C	2.62	0.41
1:A:303:SER:HB2	1:A:532:ARG:HG2	2.03	0.41
1:A:498:TYR:C	1:A:500:ARG:N	2.77	0.41
1:A:498:TYR:O	1:A:500:ARG:N	2.54	0.41
1:B:80:TRP:CG	1:B:81:GLN:H	2.37	0.41
1:B:279:VAL:CG2	1:B:405:LEU:HD13	2.49	0.41
1:B:376:PRO:O	1:B:377:VAL:C	2.64	0.41
1:B:541:ARG:O	1:B:541:ARG:CG	2.66	0.41
1:A:157:LEU:HD23	1:A:607:TYR:CZ	2.56	0.41
1:A:326:ARG:HD2	3:A:701:NAG:N2	2.36	0.41
1:B:114:LEU:HD23	1:B:114:LEU:HA	1.93	0.41
1:B:339:ASN:OD1	1:B:342:ASP:N	2.50	0.41
1:B:498:TYR:C	1:B:500:ARG:N	2.77	0.41
1:A:317:SER:HB3	1:A:319:LEU:HD21	2.01	0.41
1:A:350:ARG:HD2	1:A:351:VAL:HG23	2.01	0.41
1:B:202:TYR:O	1:B:203:ASN:C	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:557:LYS:O	1:B:561:GLY:HA2	2.19	0.41
1:A:331:HIS:CB	1:A:490:PHE:CE1	3.04	0.41
1:A:544:LEU:C	1:A:546:ALA:H	2.28	0.41
1:A:104:PRO:CG	1:A:107:LYS:HD2	2.51	0.41
1:A:441:LEU:HD23	1:A:442:VAL:HA	2.03	0.41
1:A:401:THR:HA	1:A:402:PRO:HD3	1.97	0.41
1:B:79:ILE:O	1:B:80:TRP:CG	2.74	0.41
1:B:147:LEU:HD13	1:B:256:MET:CE	2.51	0.41
1:B:210:ASP:O	1:B:214:LEU:HB2	2.21	0.41
1:B:243:ASN:OD1	1:B:246:GLY:N	2.54	0.41
1:A:309:MET:HB3	1:A:313:PHE:CD2	2.56	0.41
1:A:351:VAL:O	1:A:351:VAL:HG12	2.21	0.41
1:A:550:ARG:HG2	1:A:550:ARG:HH11	1.85	0.41
1:B:24:TYR:C	1:B:24:TYR:CD2	2.98	0.41
1:B:261:TRP:CD1	1:B:261:TRP:N	2.88	0.41
1:B:368:TYR:CE2	1:B:544:LEU:HA	2.56	0.41
1:A:157:LEU:CD2	1:A:607:TYR:CZ	3.03	0.40
1:A:279:VAL:HG13	1:A:410:LEU:HB3	1.98	0.40
1:A:500:ARG:H	1:A:500:ARG:HG3	1.54	0.40
1:A:550:ARG:HG2	1:A:550:ARG:NH1	2.36	0.40
1:B:28:ALA:HA	1:B:31:VAL:HG12	2.03	0.40
1:A:6:GLN:CG	1:A:7:PRO:CD	2.77	0.40
1:A:16:GLY:O	1:A:19:LEU:HB2	2.21	0.40
1:A:279:VAL:HG12	1:A:279:VAL:O	2.22	0.40
1:B:253:LEU:HG	1:B:261:TRP:CE2	2.56	0.40
1:A:6:GLN:HE21	1:A:6:GLN:HB2	1.60	0.40
1:A:50:ASN:HA	1:A:53:ARG:HD2	2.02	0.40
1:B:31:VAL:HG11	1:B:64:PHE:CD2	2.56	0.40
1:B:381:ARG:HB2	1:B:548:SER:HB2	2.02	0.40
1:B:456:PRO:HA	1:B:459:TYR:CD2	2.57	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:596:GLU:OE2	2:C:1:NAG:O7[3_554]	2.01	0.19

5.3 Torsion angles

5.3.1 Protein backbone

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	610/612 (100%)	537 (88%)	60 (10%)	13 (2%)	5	25
1	B	609/612 (100%)	526 (86%)	73 (12%)	10 (2%)	7	32
All	All	1219/1224 (100%)	1063 (87%)	133 (11%)	23 (2%)	6	28

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	608	PRO
1	B	3	PRO
1	B	608	PRO
1	A	7	PRO
1	A	250	ALA
1	A	280	THR
1	A	545	GLN
1	B	130	PRO
1	B	134	ALA
1	B	135	THR
1	B	391	ILE
1	A	351	VAL
1	A	111	TYR
1	A	414	VAL
1	B	77	GLU
1	A	391	ILE
1	A	495	VAL
1	B	185	ALA
1	B	408	ILE
1	A	130	PRO
1	B	79	ILE
1	A	499	ILE
1	A	530	ILE

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	496/526 (94%)	440 (89%)	56 (11%)	5	23
1	B	489/526 (93%)	435 (89%)	54 (11%)	6	24
All	All	985/1052 (94%)	875 (89%)	110 (11%)	6	23

All (110) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	11	SER
1	A	31	VAL
1	A	35	SER
1	A	52	ARG
1	A	54	GLN
1	A	59	LEU
1	A	77	GLU
1	A	82	ASN
1	A	96	ARG
1	A	100	SER
1	A	105	LEU
1	A	147	LEU
1	A	155	MET
1	A	171	LEU
1	A	187	LYS
1	A	214	LEU
1	A	221	LEU
1	A	236	ARG
1	A	245	ARG
1	A	253	LEU
1	A	262	GLU
1	A	289	ASN
1	A	295	ARG
1	A	341	LYS
1	A	342	ASP
1	A	346	LYS

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Mol	Chain	Res	Type
1	A	350	ARG
1	A	354	ASP
1	A	363	MET
1	A	372	TYR
1	A	381	ARG
1	A	388	HIS
1	A	408	ILE
1	A	410	LEU
1	A	411	LEU
1	A	441	LEU
1	A	453	ARG
1	A	480	ASN
1	A	482	THR
1	A	494	ASN
1	A	496	THR
1	A	500	ARG
1	A	507	LEU
1	A	522	GLU
1	A	540	LEU
1	A	541	ARG
1	A	543	VAL
1	A	550	ARG
1	A	553	GLN
1	A	563	ASP
1	A	570	LEU
1	A	579	GLN
1	A	582	GLN
1	A	587	GLN
1	A	594	TRP
1	B	3	PRO
1	B	18	GLN
1	B	49	GLU
1	B	66	GLU
1	B	73	LYS
1	B	76	TYR
1	B	96	ARG
1	B	105	LEU
1	B	126	LYS
1	B	129	LEU
1	B	141	PRO
1	B	144	THR
1	B	147	LEU

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Mol	Chain	Res	Type
1	B	184	GLU
1	B	203	ASN
1	B	214	LEU
1	B	231	ARG
1	B	235	ARG
1	B	245	ARG
1	B	264	ILE
1	B	274	LYS
1	B	285	GLN
1	B	289	ASN
1	B	299	GLU
1	B	312	GLU
1	B	320	GLU
1	B	342	ASP
1	B	346	LYS
1	B	354	ASP
1	B	372	TYR
1	B	374	ASP
1	B	388	HIS
1	B	408	ILE
1	B	411	LEU
1	B	414	VAL
1	B	418	THR
1	B	427	LYS
1	B	441	LEU
1	B	443	ASP
1	B	453	ARG
1	B	460	ASN
1	B	480	ASN
1	B	489	LYS
1	B	507	LEU
1	B	508	GLN
1	B	510	GLN
1	B	542	LYS
1	B	553	GLN
1	B	572	LYS
1	B	575	GLN
1	B	586	GLN
1	B	590	GLU
1	B	598	GLN
1	B	608	PRO

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (31)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	50	ASN
1	A	81	GLN
1	A	331	HIS
1	A	347	GLN
1	A	371	GLN
1	A	404	HIS
1	A	406	HIS
1	A	471	GLN
1	A	491	HIS
1	A	494	ASN
1	A	545	GLN
1	A	568	GLN
1	A	575	GLN
1	A	587	GLN
1	A	600	HIS
1	B	6	GLN
1	B	18	GLN
1	B	145	ASN
1	B	183	ASN
1	B	203	ASN
1	B	289	ASN
1	B	347	GLN
1	B	367	GLN
1	B	371	GLN
1	B	444	GLN
1	B	545	GLN
1	B	553	GLN
1	B	575	GLN
1	B	579	GLN
1	B	582	GLN

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 ⓘ

8 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	0.57	0	17,19,21	0.76	1 (5%)
2	NAG	C	2	2	14,14,15	0.52	0	17,19,21	0.69	0
2	NAG	D	1	1,2	14,14,15	0.58	0	17,19,21	0.75	1 (5%)
2	NAG	D	2	2	14,14,15	0.51	0	17,19,21	0.68	0
2	NAG	E	1	1,2	14,14,15	0.57	0	17,19,21	0.75	1 (5%)
2	NAG	E	2	2	14,14,15	0.53	0	17,19,21	0.68	0
2	NAG	F	1	1,2	14,14,15	2.36	1 (7%)	17,19,21	1.06	1 (5%)
2	NAG	F	2	2	14,14,15	2.35	1 (7%)	17,19,21	1.06	1 (5%)

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	-	3/6/23/26	0/1/1/1
2	NAG	D	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	D	2	2	-	3/6/23/26	0/1/1/1
2	NAG	E	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	E	2	2	-	3/6/23/26	0/1/1/1
2	NAG	F	1	1,2	-	0/6/23/26	0/1/1/1
2	NAG	F	2	2	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	F	1	NAG	C8-C7	-8.77	1.32	1.50
2	F	2	NAG	C8-C7	-8.72	1.32	1.50

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	NAG	C1-C2-N2	-3.44	105.01	110.43
2	F	2	NAG	C1-C2-N2	-3.42	105.04	110.43
2	C	1	NAG	C2-N2-C7	-2.17	120.00	122.90
2	D	1	NAG	C2-N2-C7	-2.14	120.04	122.90
2	E	1	NAG	C2-N2-C7	-2.13	120.05	122.90

There are no chirality outliers.

All (11) torsion outliers are listed below:

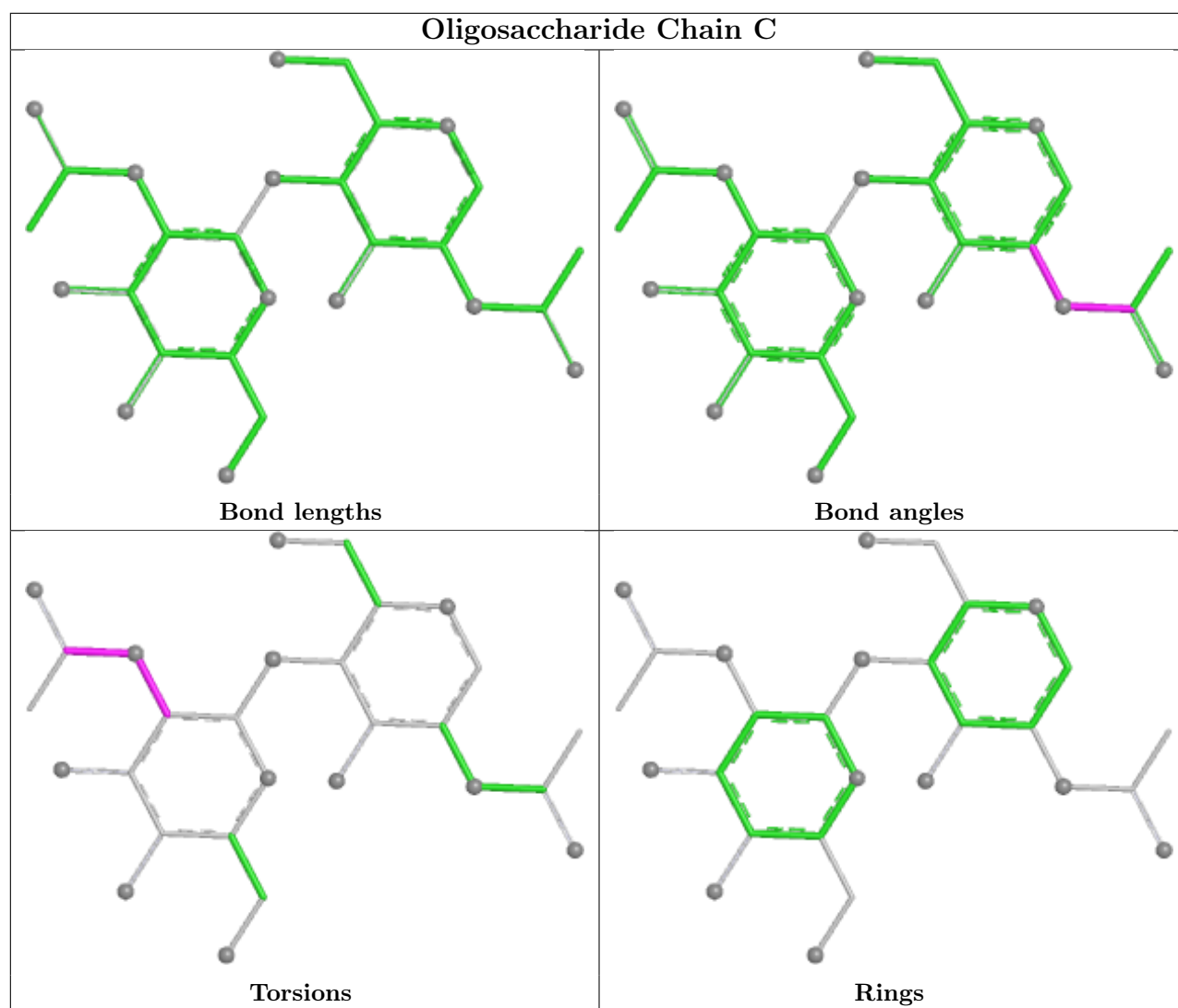
Mol	Chain	Res	Type	Atoms
2	C	2	NAG	C3-C2-N2-C7
2	D	2	NAG	C3-C2-N2-C7
2	E	2	NAG	C3-C2-N2-C7
2	C	2	NAG	C8-C7-N2-C2
2	C	2	NAG	O7-C7-N2-C2
2	D	2	NAG	C8-C7-N2-C2
2	D	2	NAG	O7-C7-N2-C2
2	E	2	NAG	C8-C7-N2-C2
2	E	2	NAG	O7-C7-N2-C2
2	F	2	NAG	O5-C5-C6-O6
2	F	2	NAG	C4-C5-C6-O6

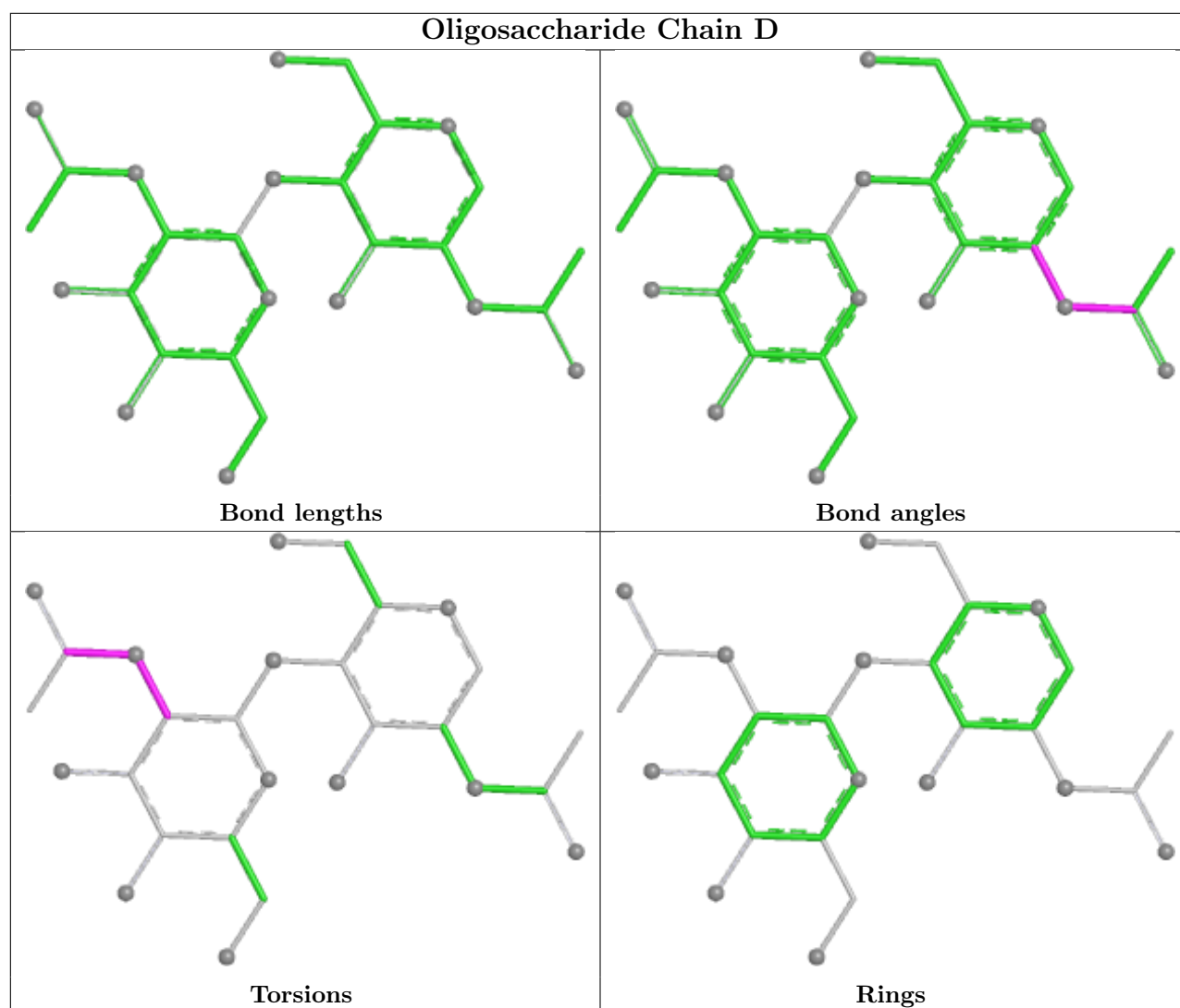
There are no ring outliers.

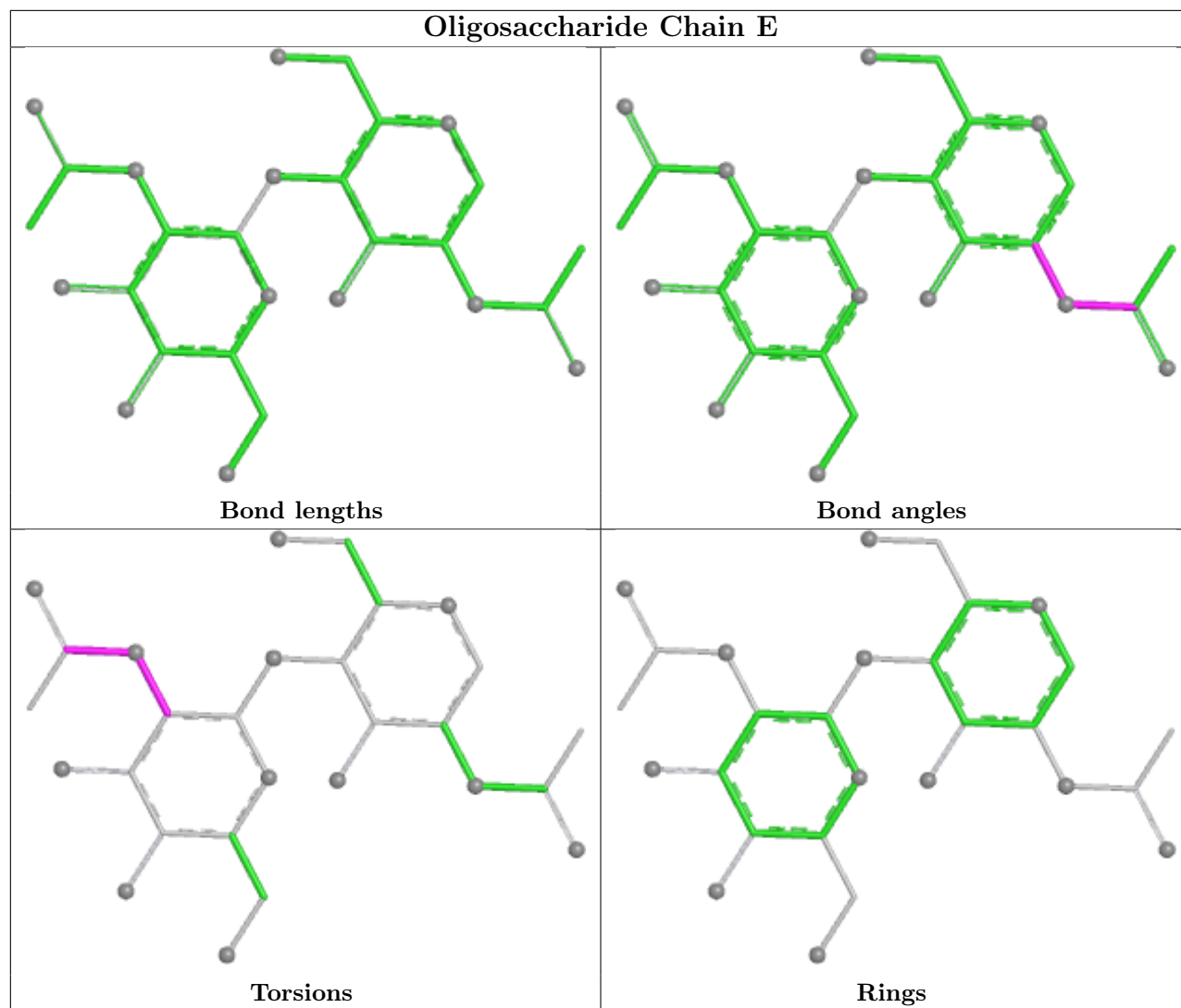
5 monomers are involved in 13 short contacts:

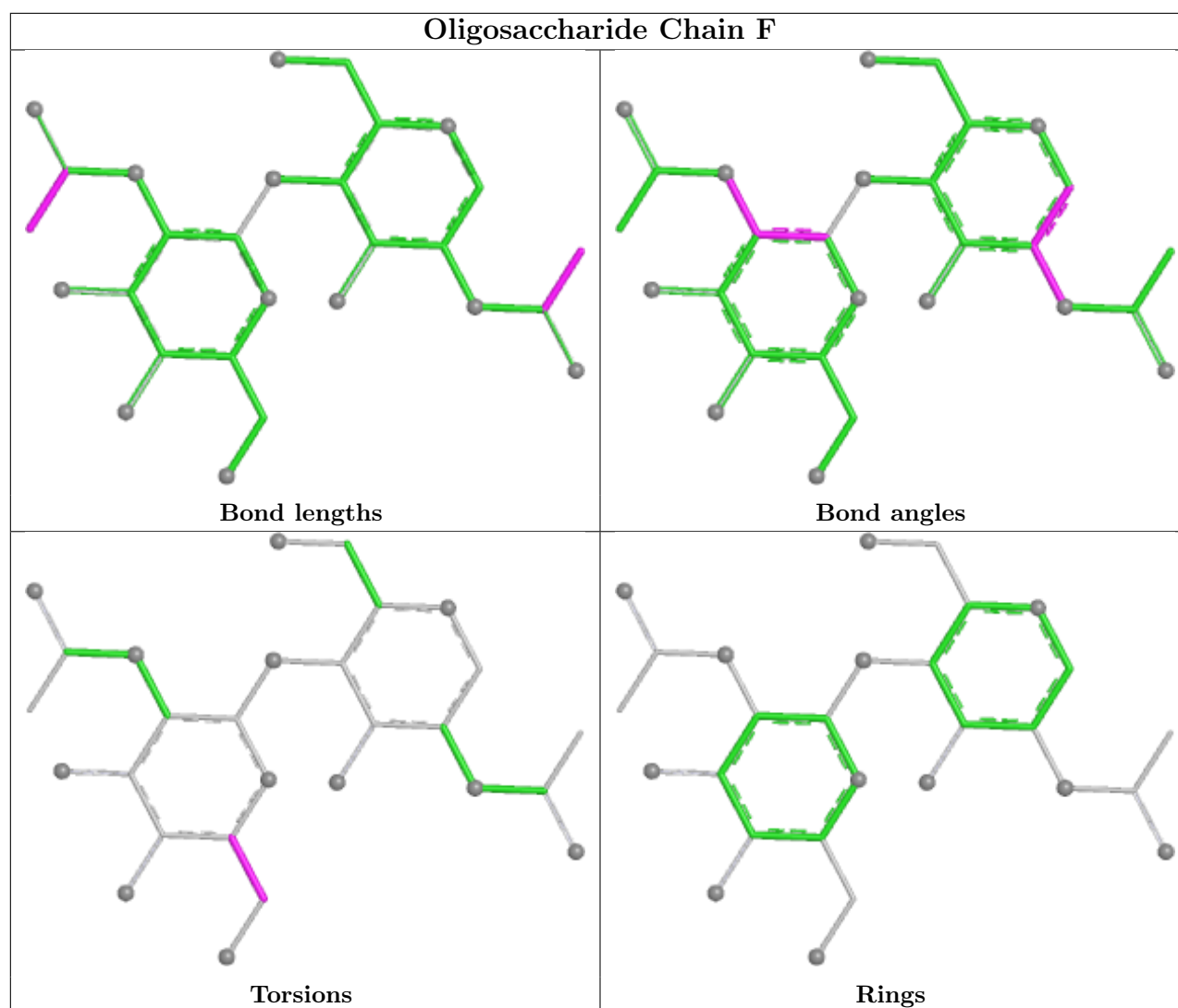
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	1	NAG	1	0
2	C	1	NAG	1	1
2	F	1	NAG	2	0
2	E	1	NAG	3	0
2	D	2	NAG	6	0

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](#)

Of 14 ligands modelled in this entry, 4 are monoatomic - leaving 10 for Mogul analysis.

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	$\# Z > 2$
6	GOL	B	706	-	5,5,5	3.49	3 (60%)	5,5,5	2.22	4 (80%)
6	GOL	A	707	-	5,5,5	3.46	3 (60%)	5,5,5	2.04	4 (80%)
3	NAG	A	702	1	14,14,15	0.58	0	17,19,21	0.75	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	ACT	B	705	-	3,3,3	1.36	1 (33%)	3,3,3	1.60	1 (33%)
3	NAG	A	703	1	14,14,15	0.57	0	17,19,21	0.75	1 (5%)
3	NAG	B	703	1	14,14,15	0.57	0	17,19,21	0.75	1 (5%)
3	NAG	B	702	1	14,14,15	3.79	2 (14%)	17,19,21	1.63	3 (17%)
5	ACT	A	705	-	3,3,3	1.24	0	3,3,3	1.50	1 (33%)
3	NAG	A	701	1	14,14,15	3.99	3 (21%)	17,19,21	1.93	4 (23%)
5	ACT	A	706	-	3,3,3	1.36	0	3,3,3	1.57	1 (33%)

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
6	GOL	B	706	-	-	1/4/4/4	-
6	GOL	A	707	-	-	1/4/4/4	-
3	NAG	A	702	1	-	0/6/23/26	0/1/1/1
3	NAG	A	703	1	-	0/6/23/26	0/1/1/1
3	NAG	B	703	1	-	0/6/23/26	0/1/1/1
3	NAG	B	702	1	-	5/6/23/26	0/1/1/1
3	NAG	A	701	1	-	5/6/23/26	0/1/1/1

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	701	NAG	C8-C7	-13.57	1.22	1.50
3	B	702	NAG	C8-C7	-13.53	1.22	1.50
6	B	706	GOL	C3-C2	-6.59	1.26	1.51
6	A	707	GOL	C3-C2	-6.53	1.26	1.51
3	A	701	NAG	O5-C5	4.55	1.52	1.43
3	A	701	NAG	O7-C7	4.13	1.32	1.23
3	B	702	NAG	O7-C7	4.10	1.32	1.23
6	B	706	GOL	O2-C2	2.81	1.51	1.43
6	A	707	GOL	C1-C2	2.69	1.62	1.51
6	A	707	GOL	O2-C2	2.64	1.51	1.43
6	B	706	GOL	C1-C2	2.45	1.61	1.51
5	B	705	ACT	O-C	2.06	1.31	1.22

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	701	NAG	C1-O5-C5	-4.45	106.22	112.19
3	A	701	NAG	C8-C7-N2	4.14	122.98	116.12
3	B	702	NAG	C8-C7-N2	4.12	122.96	116.12
3	B	702	NAG	C1-C2-N2	-3.45	104.99	110.43
3	A	701	NAG	C1-C2-N2	-3.43	105.03	110.43
3	A	701	NAG	O7-C7-N2	-3.11	116.49	121.98
3	B	702	NAG	O7-C7-N2	-3.07	116.55	121.98
6	B	706	GOL	O2-C2-C3	-2.76	97.75	109.18
6	A	707	GOL	O2-C2-C3	-2.52	98.73	109.18
6	B	706	GOL	O2-C2-C1	-2.51	98.79	109.18
6	B	706	GOL	O3-C3-C2	-2.41	99.51	110.38
6	A	707	GOL	O2-C2-C1	-2.35	99.46	109.18
6	B	706	GOL	C3-C2-C1	2.21	119.90	111.80
5	B	705	ACT	O-C-CH3	-2.19	113.56	122.53
6	A	707	GOL	O3-C3-C2	-2.18	100.57	110.38
5	A	706	ACT	O-C-CH3	-2.16	113.68	122.53
3	A	703	NAG	C2-N2-C7	-2.15	120.02	122.90
3	B	703	NAG	C2-N2-C7	-2.14	120.03	122.90
3	A	702	NAG	C2-N2-C7	-2.13	120.05	122.90
5	A	705	ACT	O-C-CH3	-2.07	114.05	122.53
6	A	707	GOL	C3-C2-C1	2.04	119.26	111.80

There are no chirality outliers.

All (12) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	702	NAG	C3-C2-N2-C7
3	A	701	NAG	O5-C5-C6-O6
3	A	701	NAG	C4-C5-C6-O6
3	B	702	NAG	O5-C5-C6-O6
3	A	701	NAG	C8-C7-N2-C2
3	A	701	NAG	O7-C7-N2-C2
3	B	702	NAG	C8-C7-N2-C2
3	B	702	NAG	O7-C7-N2-C2
3	B	702	NAG	C4-C5-C6-O6
6	A	707	GOL	O2-C2-C3-O3
6	B	706	GOL	O2-C2-C3-O3
3	A	701	NAG	C3-C2-N2-C7

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	706	GOL	1	0
6	A	707	GOL	1	0
3	A	703	NAG	1	0
3	B	702	NAG	2	0
3	A	701	NAG	3	0

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	612/612 (100%)	0.16	21 (3%)	48 27	12, 35, 58, 82	3 (0%)
1	B	611/612 (99%)	0.28	32 (5%)	33 16	12, 38, 67, 84	6 (0%)
All	All	1223/1224 (99%)	0.22	53 (4%)	40 21	12, 37, 64, 84	9 (0%)

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	8	GLY	6.0
1	A	612	ASP	5.6
1	B	77	GLU	4.8
1	B	12	ALA	4.4
1	B	278	ASP	4.3
1	A	412	ASP	4.2
1	B	410	LEU	4.0
1	B	13	ASP	4.0
1	A	135	THR	3.9
1	A	414	VAL	3.9
1	A	413	ARG	3.9
1	B	133	THR	3.7
1	B	11	SER	3.7
1	A	134	ALA	3.7
1	B	134	ALA	3.7
1	A	140	ASP	3.6
1	B	611	ILE	3.5
1	A	609	GLU	3.4
1	B	610	GLY	3.4
1	B	412	ASP	3.3
1	A	278	ASP	3.2
1	B	76	TYR	3.1
1	B	135	THR	3.1
1	B	2	ASP	3.0

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Mol	Chain	Res	Type	RSRZ
1	B	612	ASP	3.0
1	B	276	ASN	3.0
1	A	415	THR	3.0
1	B	609	GLU	2.9
1	A	350	ARG	2.9
1	B	9	ASN	2.8
1	A	132	LYS	2.8
1	A	131	ASN	2.8
1	B	374	ASP	2.8
1	A	136	CYS	2.7
1	B	277	LEU	2.6
1	B	79	ILE	2.6
1	B	15	ALA	2.6
1	B	608	PRO	2.4
1	B	274	LYS	2.4
1	A	142	ASP	2.4
1	B	5	LEU	2.4
1	B	23	SER	2.3
1	B	70	GLN	2.3
1	A	611	ILE	2.3
1	B	414	VAL	2.3
1	B	44	THR	2.2
1	A	292	HIS	2.2
1	A	330	CYS	2.2
1	B	406	HIS	2.1
1	A	133	THR	2.1
1	A	289	ASN	2.1
1	B	19	LEU	2.1
1	B	352	THR	2.0

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

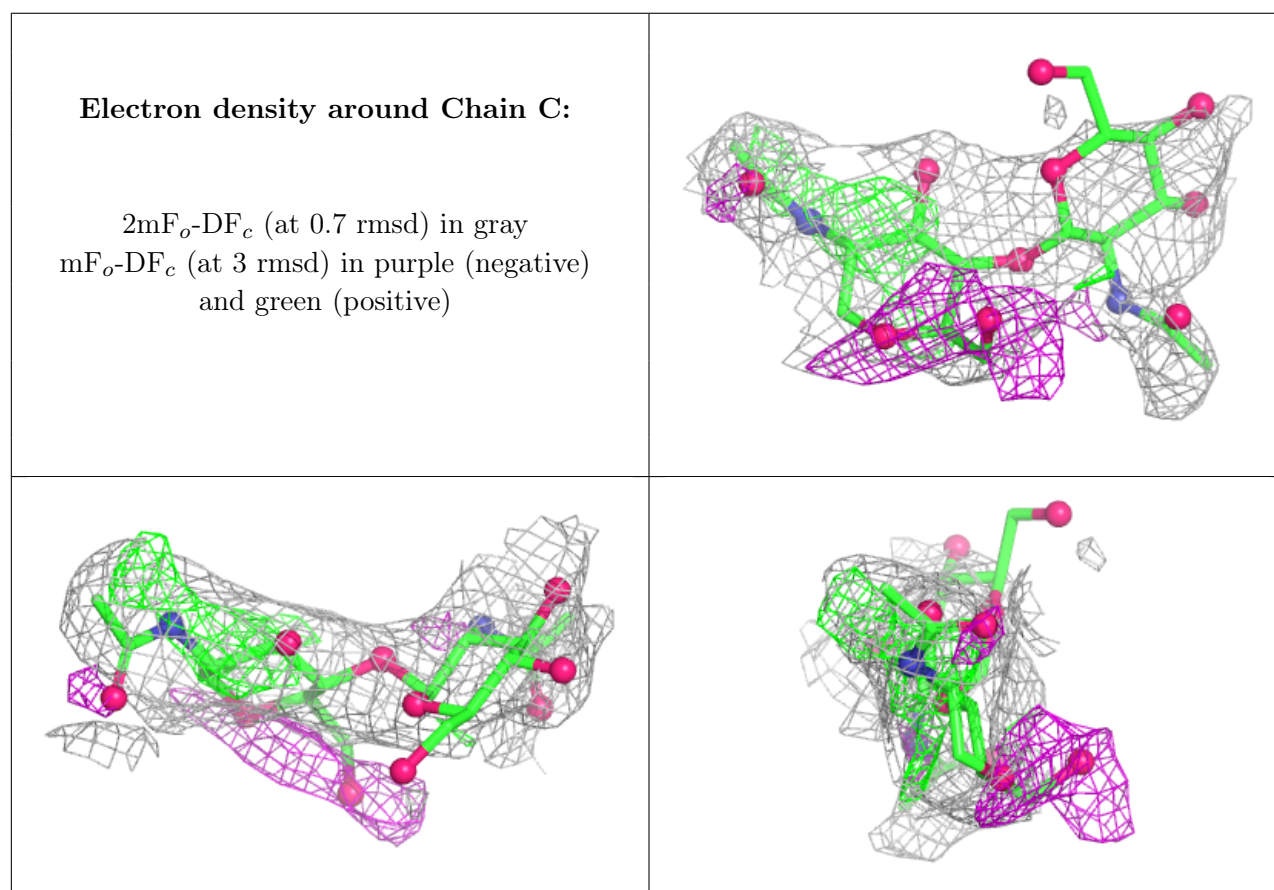
There are no non-standard protein/DNA/RNA residues in this entry.

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.

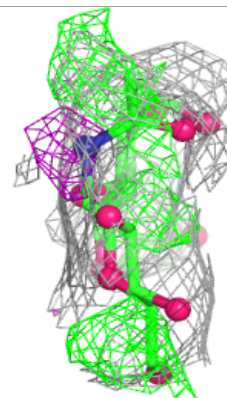
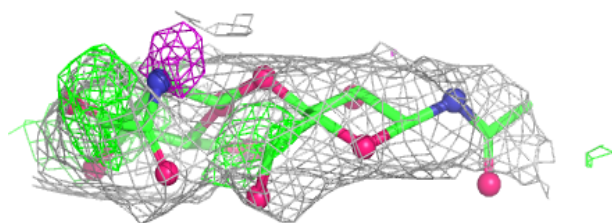
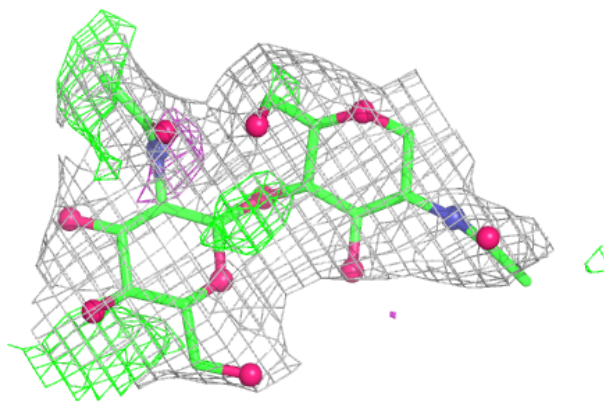
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	E	2	14/15	0.45	0.23	86,90,91,91	0
2	NAG	D	2	14/15	0.47	0.29	69,72,73,75	0
2	NAG	C	2	14/15	0.54	0.21	95,96,97,97	0
2	NAG	F	2	14/15	0.60	0.24	79,80,81,81	0
2	NAG	C	1	14/15	0.64	0.29	56,60,68,72	0
2	NAG	E	1	14/15	0.69	0.26	55,57,62,69	0
2	NAG	F	1	14/15	0.70	0.22	61,65,67,68	0
2	NAG	D	1	14/15	0.82	0.19	56,58,60,60	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.



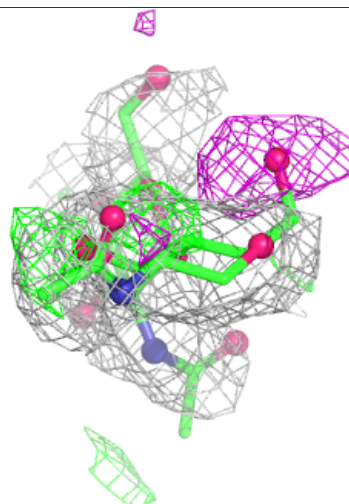
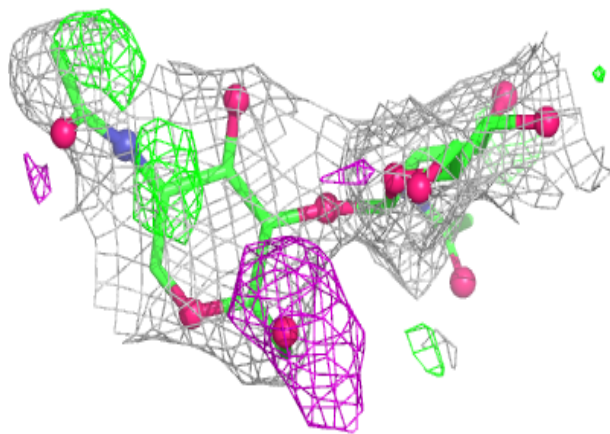
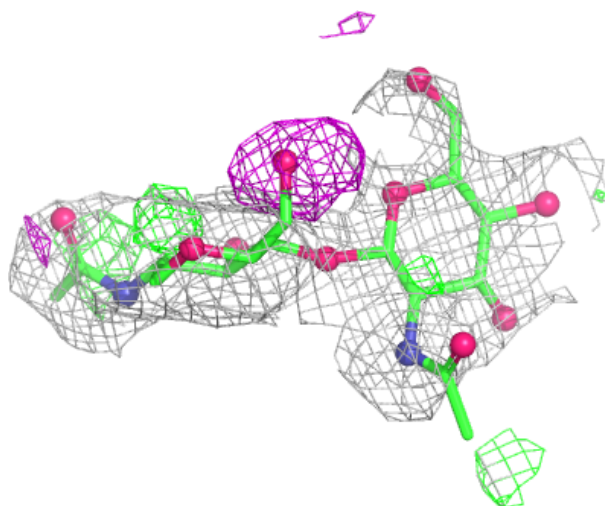
Electron density around Chain D:

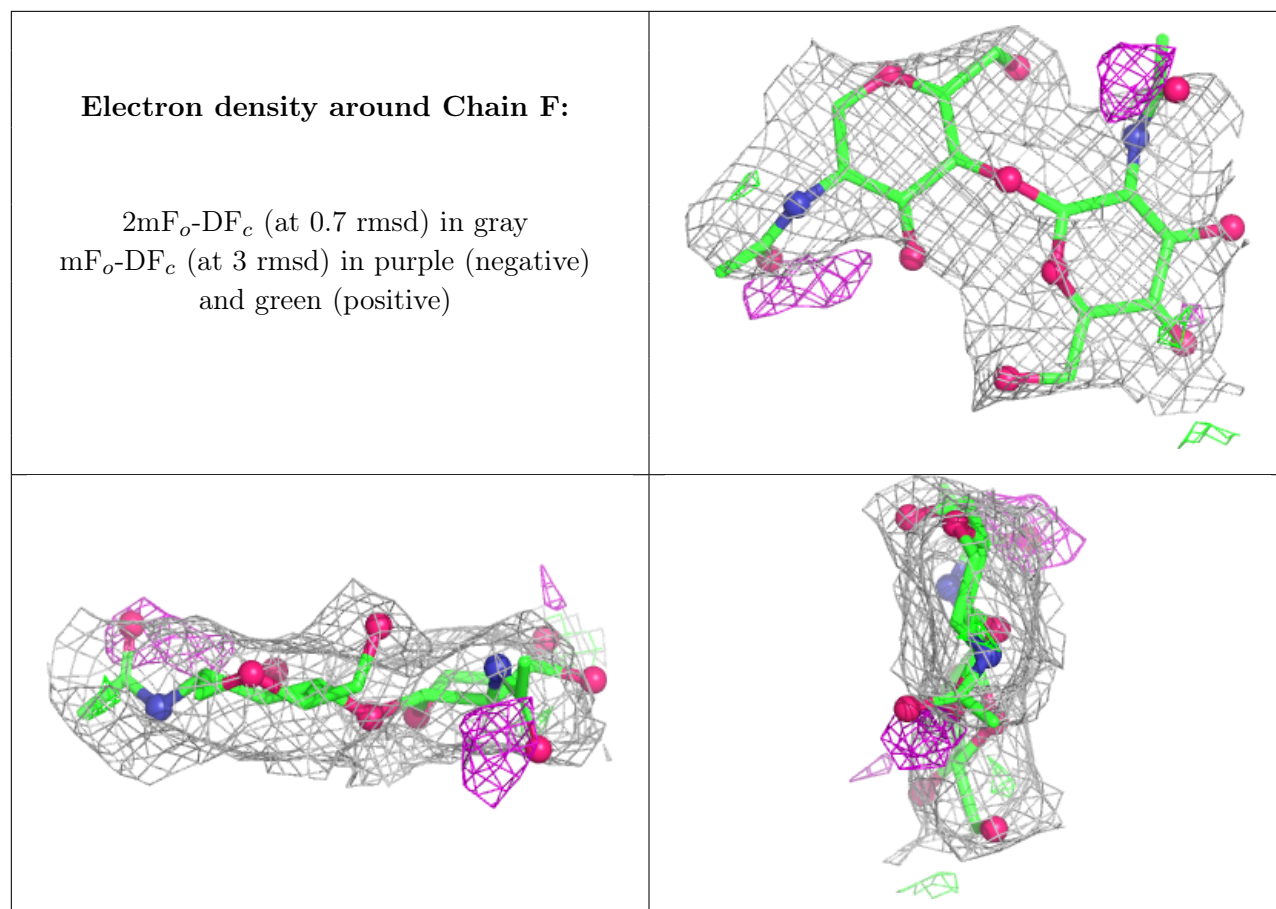
$2mF_o - DF_c$ (at 0.7 rmsd) in gray
 $mF_o - DF_c$ (at 3 rmsd) in purple (negative)
and green (positive)



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)





6.4 Ligands ⓘ

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
3	NAG	A	703	14/15	0.40	0.29	116,117,118,119	0
3	NAG	B	702	14/15	0.52	0.27	83,86,86,87	0
3	NAG	A	701	14/15	0.55	0.20	66,68,70,71	0
5	ACT	A	705	4/4	0.64	0.30	49,51,52,53	0
3	NAG	B	703	14/15	0.77	0.21	76,78,80,83	0
3	NAG	A	702	14/15	0.77	0.20	75,76,79,79	0
5	ACT	A	706	4/4	0.81	0.20	32,33,34,35	0
5	ACT	B	705	4/4	0.84	0.23	52,52,53,53	0
6	GOL	A	707	6/6	0.91	0.16	29,31,32,33	0
6	GOL	B	706	6/6	0.91	0.18	31,34,35,39	0
7	CL	A	708	1/1	0.91	0.07	29,29,29,29	0
7	CL	B	701	1/1	0.95	0.06	37,37,37,37	0
4	ZN	B	704	1/1	0.99	0.05	29,29,29,29	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	ZN	A	704	1/1	1.00	0.02	31,31,31,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.