



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 13, 2026 – 07:21 PM UTC

PDB ID : 1HJV / pdb_00001hvj
Title : Crystal structure of hcgp-39 in complex with chitin tetramer
Authors : Houston, D.R.; Recklies, A.D.; Krupa, J.C.; Van Aalten, D.M.F.
Deposited on : 2003-02-28
Resolution : 2.75 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

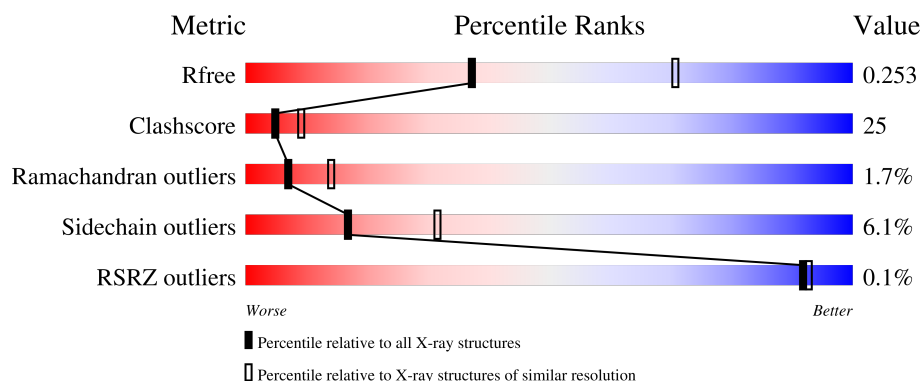
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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.



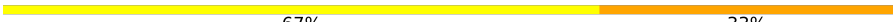
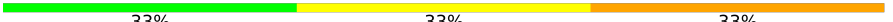
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1009 (2.76-2.76)
Clashscore	190562	1044 (2.76-2.76)
Ramachandran outliers	187476	1024 (2.76-2.76)
Sidechain outliers	187428	1024 (2.76-2.76)
RSRZ outliers	180081	1009 (2.76-2.76)

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	362	58% 37% • •
1	B	362	61% 35% •
1	C	362	42% 51% 6% •
1	D	362	53% 41% 5%
2	E	3	33% 33% 33%

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Mol	Chain	Length	Quality of chain
2	G	3	 100%
2	I	3	 67% 33%
2	K	3	 33% 33% 33%
3	F	4	 50% 50%
3	H	4	 25% 25% 50%
3	J	4	 25% 75%
3	L	4	 25% 25% 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
3	NAG	J	3	-	-	X	-
3	NAG	L	3	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 12296 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.

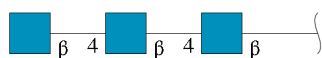
- Molecule 1 is a protein called CHITINASE-3 LIKE PROTEIN 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	362	Total	C	N	O	S	0	2	0
			2871	1833	495	532	11			
1	B	362	Total	C	N	O	S	0	1	0
			2868	1831	496	530	11			
1	C	362	Total	C	N	O	S	0	2	0
			2873	1834	496	532	11			
1	D	362	Total	C	N	O	S	0	1	0
			2866	1830	494	531	11			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	311	ILE	THR	conflict	UNP P36222
B	311	ILE	THR	conflict	UNP P36222
A	311	ILE	THR	conflict	UNP P36222
B	311	ILE	THR	conflict	UNP P36222

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	3	Total	C	N	O	0	0	0
			43	24	3	16			
2	G	3	Total	C	N	O	0	0	0
			43	24	3	16			
2	I	3	Total	C	N	O	0	0	0
			43	24	3	16			

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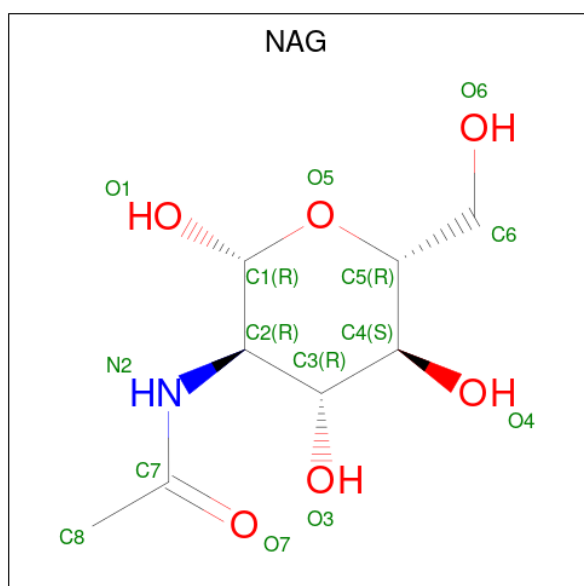
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	K	3	Total	C	N	O	0	0	0
			43	24	3	16			

- Molecule 3 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



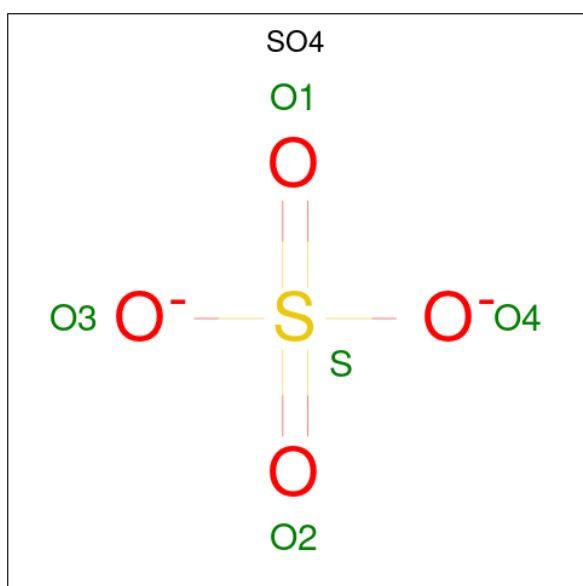
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
3	F	4	Total	C	N	O	0	0	0
			57	32	4	21			
3	H	4	Total	C	N	O	0	0	0
			57	32	4	21			
3	J	4	Total	C	N	O	0	0	0
			57	32	4	21			
3	L	4	Total	C	N	O	0	0	0
			57	32	4	21			

- Molecule 4 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



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		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	D	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 5 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	C	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		
5	D	1	Total	O	S	0	0
			5	4	1		

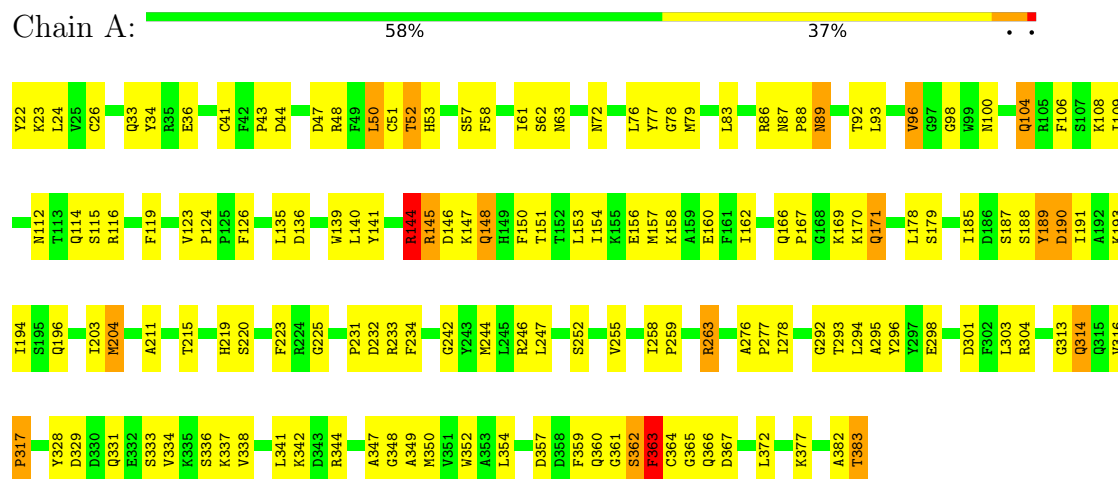
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	101	Total	O	0	0
			101	101		
6	B	80	Total	O	0	0
			80	80		
6	C	53	Total	O	0	0
			53	53		
6	D	68	Total	O	0	0
			68	68		

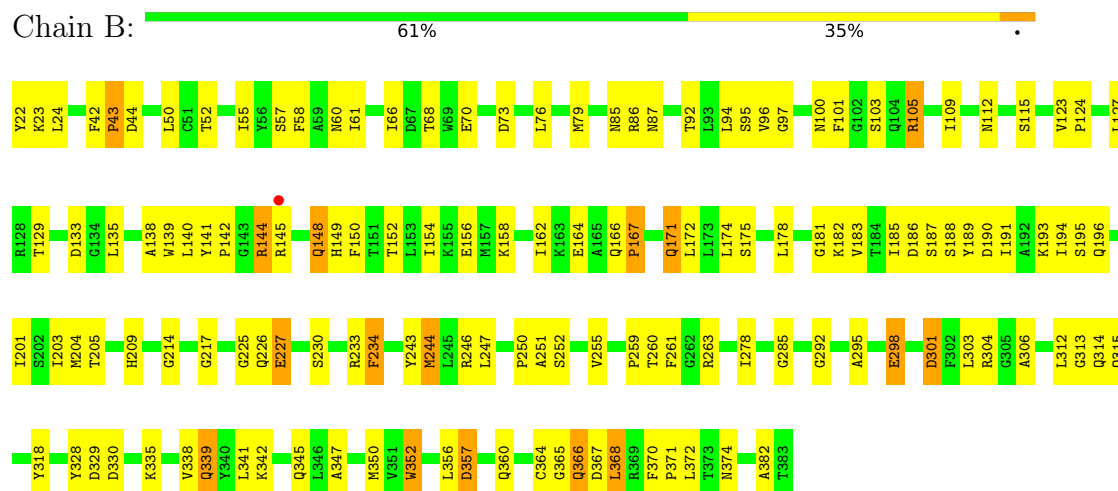
3 Residue-property plots

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: CHITINASE-3 LIKE PROTEIN 1

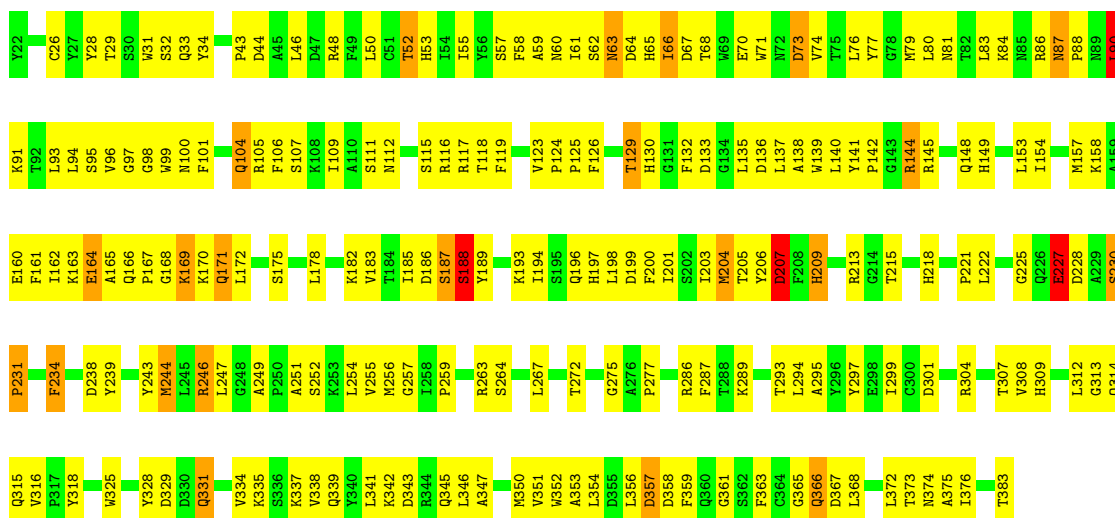


• Molecule 1: CHITINASE-3 LIKE PROTEIN 1



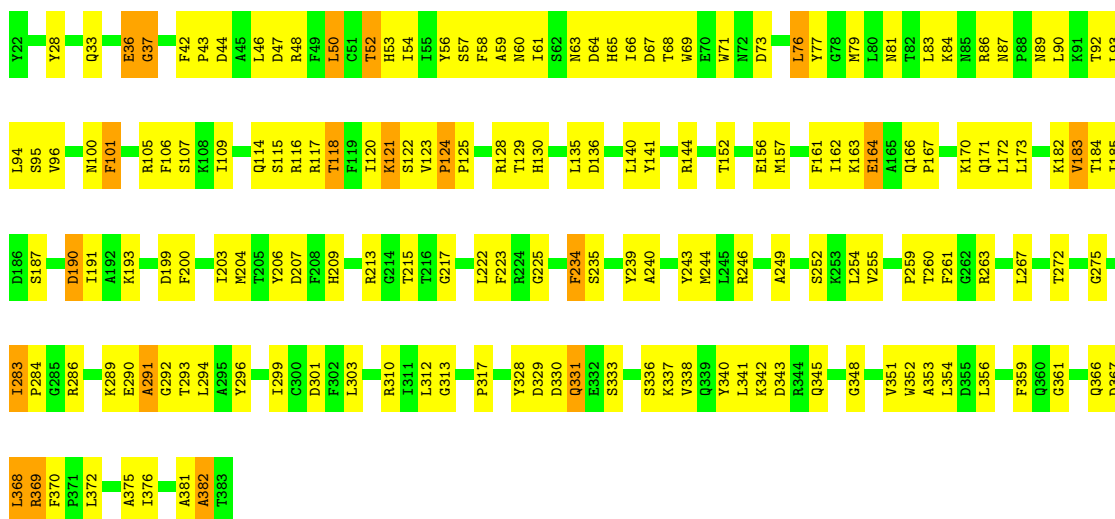
• Molecule 1: CHITINASE-3 LIKE PROTEIN 1





• Molecule 1: CHITINASE-3 LIKE PROTEIN 1

Chain D: 53% 41% 5%



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

Chain E: 33% 33% 33%



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

Chain G: 100%

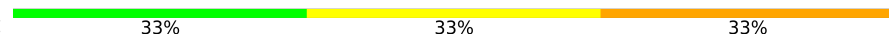


- Molecule 2: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  67% 33%



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

Chain K:  33% 33% 33%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  50% 50%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain H:  25% 25% 50%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  25% 75%



- Molecule 3: 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain L:  25% 25% 50%



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	108.64Å 123.32Å 135.49Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.79 – 2.75 24.79 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.0 (24.79-2.75) 98.8 (24.79-2.75)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.94 (at 2.72Å)	Xtrriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.212 , 0.253 0.227 , 0.253	Depositor DCC
R_{free} test set	983 reflections (2.05%)	wwPDB-VP
Wilson B-factor (Å ²)	55.6	Xtrriage
Anisotropy	0.500	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	12296	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 65.69 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 6.8110e-06. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: SO4, NAG

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	0.57	1/2955 (0.0%)	1.06	20/4002 (0.5%)
1	B	0.58	0/2948	1.06	13/3992 (0.3%)
1	C	0.49	0/2957	1.03	18/4004 (0.4%)
1	D	0.53	0/2946	1.04	12/3990 (0.3%)
All	All	0.54	1/11806 (0.0%)	1.05	63/15988 (0.4%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	204	MET	SD-CE	-5.81	1.65	1.79

All (63) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	357	ASP	N-CA-C	-8.78	99.21	110.53
1	D	234	PHE	N-CA-C	7.90	122.69	112.26
1	A	50	LEU	N-CA-C	7.58	119.18	111.07
1	C	312	LEU	N-CA-C	7.57	120.25	111.02
1	B	357	ASP	N-CA-C	-7.50	100.01	110.35
1	C	227	GLU	N-CA-C	7.26	120.11	111.33
1	C	357	ASP	N-CA-C	-7.20	100.83	110.55
1	B	312	LEU	N-CA-C	7.15	119.75	111.02
1	A	360	GLN	N-CA-C	-7.15	102.90	111.69
1	D	190	ASP	N-CA-C	-6.92	99.28	109.11
1	C	234	PHE	N-CA-C	6.82	118.88	109.54
1	C	287	PHE	N-CA-C	6.81	120.06	111.69
1	D	260	THR	N-CA-C	-6.72	104.40	112.59
1	D	283	ILE	CA-C-N	6.69	126.45	119.76
1	D	283	ILE	C-N-CA	6.69	126.45	119.76
1	C	52	THR	N-CA-C	-6.66	105.19	113.38

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	190	ASP	N-CA-C	-6.49	99.90	109.11
1	B	144	ARG	N-CA-C	-6.37	104.44	111.82
1	A	77	TYR	N-CA-C	-6.35	103.62	111.75
1	A	144	ARG	N-CA-C	-6.35	104.45	111.82
1	D	361	GLY	N-CA-C	-6.26	105.88	114.64
1	B	234	PHE	N-CA-C	6.22	121.03	113.38
1	B	171	GLN	N-CA-C	6.22	119.50	110.42
1	A	190	ASP	N-CA-C	-6.15	99.87	108.54
1	B	360	GLN	N-CA-C	-6.09	104.56	111.14
1	A	189	TYR	N-CA-C	6.08	119.45	109.85
1	C	90	LEU	N-CA-C	-6.00	100.02	109.50
1	A	295	ALA	N-CA-C	-5.94	102.53	110.55
1	A	314	GLN	N-CA-C	5.90	120.64	113.38
1	A	258	ILE	CA-C-N	5.88	126.34	119.99
1	A	258	ILE	C-N-CA	5.88	126.34	119.99
1	A	188	SER	N-CA-C	5.80	120.64	113.50
1	C	199	ASP	N-CA-C	-5.64	105.66	112.54
1	C	375	ALA	N-CA-C	-5.58	105.28	111.36
1	D	52	THR	N-CA-C	-5.57	106.64	113.50
1	C	230	SER	CA-C-N	5.57	126.80	119.84
1	C	230	SER	C-N-CA	5.57	126.80	119.84
1	C	188	SER	N-CA-C	5.51	122.54	110.80
1	C	171	GLN	N-CA-C	5.51	117.40	110.24
1	B	158	LYS	N-CA-C	-5.51	105.29	112.23
1	B	50	LEU	N-CA-C	5.49	116.95	110.97
1	B	227	GLU	N-CA-C	5.42	117.64	111.02
1	A	363	PHE	N-CA-C	5.40	122.31	110.80
1	C	205	THR	N-CA-C	5.39	117.96	109.39
1	C	46	LEU	N-CA-C	5.38	118.32	109.76
1	A	87	ASN	CA-C-N	5.35	125.43	119.87
1	A	87	ASN	C-N-CA	5.35	125.43	119.87
1	B	352	TRP	N-CA-C	-5.32	98.71	108.02
1	A	171	GLN	N-CA-C	5.31	117.73	110.24
1	A	52	THR	N-CA-C	-5.29	106.96	113.41
1	A	234	PHE	N-CA-C	5.29	119.33	111.87
1	B	298	GLU	N-CA-C	-5.29	105.57	112.23
1	A	361	GLY	N-CA-C	-5.24	107.44	114.25
1	D	301	ASP	N-CA-C	-5.20	105.26	111.03
1	D	375	ALA	N-CA-C	-5.18	105.54	111.14
1	D	50	LEU	N-CA-C	5.11	116.54	111.07
1	A	298	GLU	N-CA-C	-5.11	105.41	111.69
1	D	336	SER	N-CA-C	-5.11	105.61	111.07

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	361	GLY	N-CA-C	-5.07	107.69	115.00
1	D	124	PRO	N-CA-C	5.06	116.87	110.70
1	C	87	ASN	CA-C-N	5.05	124.77	119.82
1	C	87	ASN	C-N-CA	5.05	124.77	119.82
1	B	295	ALA	N-CA-C	-5.03	103.89	110.53

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2871	0	2789	133	0
1	B	2868	0	2790	113	0
1	C	2873	0	2791	194	0
1	D	2866	0	2785	146	0
2	E	43	0	39	3	0
2	G	43	0	39	0	0
2	I	43	0	39	4	0
2	K	43	0	39	1	0
3	F	57	0	51	11	0
3	H	57	0	51	6	0
3	J	57	0	51	15	0
3	L	57	0	51	11	0
4	A	14	0	13	0	0
4	B	14	0	13	0	0
4	C	14	0	13	1	0
4	D	14	0	13	0	0
5	A	15	0	0	0	0
5	B	15	0	0	0	0
5	C	15	0	0	0	0
5	D	15	0	0	0	0
6	A	101	0	0	25	0
6	B	80	0	0	12	0
6	C	53	0	0	2	0
6	D	68	0	0	9	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	12296	0	11567	596	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 25.

All (596) 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:D:162:ILE:HG12	1:D:171:GLN:HE22	1.31	0.95
1:A:145:ARG:HG3	6:A:2041:HOH:O	1.64	0.94
1:C:193:LYS:O	1:C:196:GLN:HG2	1.69	0.93
1:C:144:ARG:HB2	1:C:144:ARG:HH11	1.34	0.92
1:D:352:TRP:HB3	6:D:2061:HOH:O	1.69	0.91
1:A:112:ASN:HB3	1:A:115:SER:HB2	1.52	0.91
1:D:166:GLN:HG3	1:D:167:PRO:HD3	1.52	0.91
1:D:47:ASP:HB3	1:D:50:LEU:HB2	1.54	0.90
1:B:94:LEU:O	1:B:135:LEU:HD12	1.72	0.90
1:A:162:ILE:HG12	1:A:171:GLN:OE1	1.76	0.85
1:B:204:MET:HE1	3:H:3:NAG:C1	2.06	0.85
1:C:154:ILE:HD13	1:C:198:LEU:HD21	1.58	0.85
1:D:48:ARG:HE	1:D:86:ARG:HB2	1.43	0.84
1:B:261:PHE:HD1	6:B:2072:HOH:O	1.60	0.84
1:A:57:SER:HB2	1:A:58:PHE:CD1	2.12	0.83
1:C:209:HIS:CE1	1:C:213:ARG:HH11	1.98	0.81
1:A:61:ILE:HG21	1:A:109:ILE:HD13	1.61	0.81
3:J:3:NAG:H61	3:J:4:NAG:H82	1.60	0.81
1:B:261:PHE:CD1	6:B:2072:HOH:O	2.33	0.80
1:D:163:LYS:HA	1:D:166:GLN:HE21	1.48	0.79
1:A:48:ARG:HB3	1:A:83:LEU:HD22	1.63	0.79
1:A:204:MET:HE2	3:F:2:NAG:H62	1.65	0.79
1:C:62:SER:O	1:C:63:ASN:HB2	1.82	0.79
1:D:123:VAL:HB	1:D:124:PRO:HD3	1.65	0.79
1:C:352:TRP:CZ3	3:J:3:NAG:H83	2.18	0.78
1:D:204:MET:HE1	3:L:3:NAG:C1	2.14	0.77
1:C:93:LEU:HD23	1:C:135:LEU:O	1.85	0.77
1:A:334:VAL:O	1:A:338:VAL:HG23	1.84	0.77
1:C:203:ILE:HD12	1:C:244:MET:SD	2.25	0.77
1:C:144:ARG:HB2	1:C:144:ARG:NH1	1.99	0.76
1:A:89:ASN:N	1:A:89:ASN:HD22	1.84	0.76
1:B:352:TRP:O	6:B:2072:HOH:O	2.04	0.76
1:D:54:ILE:HG13	1:D:90:LEU:HD11	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:LEU:HD12	1:C:201:ILE:HG12	1.69	0.74
1:D:164:GLU:O	1:D:167:PRO:HD2	1.87	0.74
1:C:104:GLN:HE21	1:C:104:GLN:HA	1.51	0.74
1:C:185:ILE:HA	1:C:189:TYR:HD2	1.53	0.73
1:A:193:LYS:O	1:A:196:GLN:HG2	1.87	0.73
1:A:144:ARG:HD3	1:A:144:ARG:N	2.03	0.73
1:D:162:ILE:HG12	1:D:171:GLN:NE2	2.05	0.72
1:D:331:GLN:H	1:D:331:GLN:HE21	1.38	0.72
1:D:166:GLN:CG	1:D:167:PRO:HD3	2.20	0.71
1:A:204:MET:CE	3:F:2:NAG:H62	2.20	0.71
1:B:328:TYR:HA	6:B:2068:HOH:O	1.91	0.71
1:A:204:MET:HE1	3:F:3:NAG:C1	2.21	0.71
1:C:74:VAL:HG13	1:C:130:HIS:CE1	2.26	0.70
1:C:198:LEU:HD12	1:C:201:ILE:CG1	2.20	0.70
1:C:33:GLN:HG3	1:C:34:TYR:CD1	2.26	0.70
1:C:111:SER:HA	1:C:149:HIS:HD2	1.56	0.70
1:C:209:HIS:CE1	1:C:213:ARG:NH1	2.59	0.70
1:C:158:LYS:HD2	1:C:197:HIS:O	1.92	0.70
1:C:185:ILE:HA	1:C:189:TYR:CD2	2.27	0.69
1:C:185:ILE:HD13	1:C:244:MET:HE3	1.72	0.69
1:B:162:ILE:HD11	6:B:2032:HOH:O	1.91	0.69
1:A:277:PRO:HB2	1:C:144:ARG:NH2	2.08	0.69
1:B:44:ASP:H	1:B:79:MET:HE2	1.58	0.69
1:C:87:ASN:ND2	1:C:90:LEU:HB3	2.07	0.69
1:C:123:VAL:HB	1:C:124:PRO:HD3	1.75	0.69
1:C:70:GLU:O	1:C:73:ASP:HB2	1.93	0.68
1:C:206:TYR:O	1:C:207:ASP:HB2	1.93	0.68
1:B:366:GLN:HB2	1:B:368:LEU:HD12	1.74	0.68
1:A:169:LYS:NZ	6:A:2050:HOH:O	2.25	0.68
1:A:150:PHE:O	6:A:2047:HOH:O	2.12	0.67
1:C:295:ALA:O	1:C:299:ILE:HG13	1.94	0.67
1:C:341:LEU:HD12	1:C:342:LYS:N	2.09	0.67
1:C:161:PHE:CD2	1:C:172:LEU:HB2	2.30	0.66
1:B:127:LEU:HD12	1:B:172:LEU:HD13	1.76	0.66
1:C:64:ASP:OD1	1:C:115:SER:HB3	1.95	0.66
1:C:141:TYR:HH	3:J:2:NAG:HO6	1.39	0.66
1:C:198:LEU:CD1	1:C:201:ILE:HG12	2.25	0.65
1:C:112:ASN:HB3	1:C:115:SER:HB2	1.77	0.65
1:D:368:LEU:HD12	1:D:368:LEU:C	2.22	0.65
1:C:366:GLN:HB3	1:C:368:LEU:HG	1.78	0.65
1:A:365:GLY:C	1:A:367:ASP:H	2.03	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:225:GLY:HA2	1:A:313:GLY:HA3	1.77	0.65
1:C:163:LYS:O	1:C:166:GLN:HG2	1.97	0.65
1:B:105:ARG:HG2	1:B:105:ARG:HH21	1.61	0.64
1:B:185:ILE:HD13	1:B:244:MET:HE3	1.80	0.64
1:D:105:ARG:HH21	1:D:105:ARG:HG2	1.62	0.64
1:B:44:ASP:HB3	1:B:79:MET:HE3	1.80	0.64
1:C:334:VAL:O	1:C:338:VAL:HG23	1.98	0.64
1:A:158:LYS:O	1:A:162:ILE:HG13	1.98	0.64
1:A:61:ILE:HG21	1:A:109:ILE:CD1	2.26	0.63
1:D:66:ILE:HG23	1:D:66:ILE:O	1.97	0.63
1:B:203:ILE:HD12	1:B:244:MET:SD	2.39	0.63
1:C:29:THR:OG1	1:C:32:SER:HB3	1.98	0.63
1:C:160:GLU:OE1	1:C:163:LYS:HD3	1.98	0.63
3:J:3:NAG:C6	3:J:4:NAG:H82	2.29	0.63
1:A:76:LEU:HD13	6:A:2022:HOH:O	1.98	0.63
1:D:65:HIS:CE1	1:D:118:THR:HG22	2.33	0.63
1:B:100:ASN:CG	3:H:4:NAG:H62	2.24	0.62
1:B:246:ARG:HD2	6:B:2047:HOH:O	1.99	0.62
1:D:48:ARG:NE	1:D:86:ARG:HB2	2.12	0.62
1:C:352:TRP:CZ2	3:J:3:NAG:H3	2.35	0.62
1:A:43:PRO:HB2	1:A:79:MET:HB2	1.81	0.62
1:A:144:ARG:CD	1:A:144:ARG:H	2.13	0.62
1:C:50:LEU:HD13	1:C:374:ASN:OD1	2.00	0.62
1:C:204:MET:HE1	3:J:3:NAG:C7	2.30	0.62
1:A:41:CYS:HB2	1:A:359:PHE:CZ	2.36	0.61
1:C:55:ILE:CG2	1:C:95:SER:HB2	2.30	0.61
1:B:135:LEU:HD23	6:B:2025:HOH:O	2.00	0.61
1:C:87:ASN:HD22	1:C:90:LEU:HB3	1.64	0.61
1:C:209:HIS:HE1	1:C:213:ARG:HH11	1.49	0.61
1:A:112:ASN:HB3	1:A:115:SER:CB	2.29	0.61
1:B:186:ASP:HA	6:B:2035:HOH:O	2.01	0.61
1:C:100:ASN:ND2	3:J:4:NAG:H62	2.15	0.61
1:A:301:ASP:HA	1:A:304:ARG:NH1	2.16	0.61
1:B:368:LEU:O	1:B:368:LEU:HD13	2.01	0.61
1:B:338:VAL:HA	1:B:341:LEU:CD2	2.31	0.61
1:B:182:LYS:NZ	1:B:186:ASP:OD2	2.34	0.60
1:C:162:ILE:HG12	1:C:171:GLN:HE22	1.66	0.60
1:D:144:ARG:HA	1:D:187:SER:O	2.01	0.60
1:A:252:SER:O	1:A:347:ALA:HB2	2.01	0.60
2:I:2:NAG:H62	2:I:3:NAG:HN2	1.65	0.60
1:B:96:VAL:HG13	1:B:96:VAL:O	2.01	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:42:PHE:HB3	1:B:79:MET:HE1	1.84	0.59
1:B:204:MET:HE3	3:H:3:NAG:H82	1.84	0.59
1:C:60:ASN:CG	1:C:61:ILE:H	2.11	0.59
2:E:2:NAG:H61	2:E:3:NAG:H82	1.83	0.59
1:A:92:THR:C	1:A:93:LEU:HD12	2.27	0.59
1:B:61:ILE:HG21	1:B:109:ILE:HD13	1.84	0.59
1:D:261:PHE:HB3	1:D:356:LEU:HD13	1.84	0.59
1:D:331:GLN:H	1:D:331:GLN:NE2	2.00	0.59
1:A:147:LYS:HE3	1:A:189:TYR:O	2.02	0.59
1:B:260:THR:N	6:B:2072:HOH:O	2.36	0.59
1:C:90:LEU:HD12	1:C:90:LEU:C	2.27	0.59
2:E:2:NAG:H61	2:E:3:NAG:C8	2.32	0.59
1:B:112:ASN:HB3	1:B:115:SER:HB2	1.85	0.59
1:B:100:ASN:ND2	3:H:4:NAG:H62	2.17	0.58
1:C:365:GLY:C	1:C:367:ASP:H	2.11	0.58
1:D:73:ASP:HB2	6:D:2014:HOH:O	2.03	0.58
6:D:2003:HOH:O	3:L:2:NAG:H83	2.03	0.58
1:A:72:ASN:ND2	1:A:76:LEU:HD13	2.19	0.58
1:B:162:ILE:O	1:B:166:GLN:HG2	2.04	0.58
1:C:162:ILE:HG12	1:C:171:GLN:NE2	2.18	0.58
1:C:331:GLN:H	1:C:331:GLN:NE2	2.01	0.58
1:D:144:ARG:NH1	1:D:187:SER:CB	2.67	0.58
1:D:234:PHE:HB3	1:D:239:TYR:CD1	2.38	0.58
1:C:84:LYS:HZ3	1:C:133:ASP:CG	2.12	0.58
1:D:343:ASP:C	1:D:345:GLN:H	2.12	0.58
1:D:381:ALA:O	1:D:382:ALA:HB2	2.03	0.58
1:A:162:ILE:O	1:A:166:GLN:HG2	2.04	0.58
2:I:2:NAG:H62	2:I:3:NAG:N2	2.19	0.58
1:A:296:TYR:CG	1:A:372:LEU:HD11	2.38	0.57
1:D:372:LEU:O	1:D:376:ILE:HG13	2.04	0.57
1:C:104:GLN:HA	1:C:104:GLN:NE2	2.18	0.57
1:D:290:GLU:O	1:D:291:ALA:C	2.46	0.57
1:B:86[A]:ARG:HG2	1:B:86[A]:ARG:HH11	1.70	0.57
1:D:352:TRP:O	6:D:2061:HOH:O	2.17	0.57
1:D:60:ASN:HB2	1:D:69:TRP:CE3	2.40	0.57
1:C:58:PHE:CZ	1:C:98:GLY:HA2	2.40	0.57
1:C:175:SER:HB2	1:C:200:PHE:CE1	2.39	0.57
1:C:43:PRO:HB2	1:C:79:MET:HB2	1.85	0.57
1:C:178:LEU:O	1:C:204:MET:HG3	2.04	0.57
1:B:101:PHE:CD1	1:B:101:PHE:C	2.83	0.57
1:D:61:ILE:HA	1:D:65:HIS:O	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:164:GLU:O	1:C:169:LYS:HD2	2.05	0.56
1:D:56:TYR:HB3	1:D:94:LEU:HD12	1.87	0.56
1:A:86:ARG:HB2	1:A:86:ARG:CZ	2.35	0.56
1:C:28:TYR:OH	1:C:43:PRO:HD3	2.06	0.56
1:C:194:ILE:O	1:C:198:LEU:HG	2.06	0.56
1:D:191:ILE:CG2	1:D:249:ALA:HB2	2.36	0.56
1:C:329:ASP:OD2	1:C:337:LYS:HE3	2.06	0.55
1:C:368:LEU:C	1:C:368:LEU:HD12	2.30	0.55
1:B:43:PRO:HB2	1:B:79:MET:HB3	1.88	0.55
1:D:141:TYR:HH	3:L:2:NAG:HO6	1.49	0.55
1:B:142:PRO:HB2	1:B:188:SER:HB3	1.89	0.55
1:A:144:ARG:HG2	6:A:2041:HOH:O	2.06	0.55
1:A:148:GLN:NE2	6:A:2044:HOH:O	2.40	0.55
1:C:142:PRO:O	1:C:188:SER:HB3	2.06	0.55
1:D:191:ILE:HD11	1:D:244:MET:HE2	1.88	0.55
1:A:36:GLU:HB3	6:A:2012:HOH:O	2.05	0.55
1:A:154:ILE:N	6:A:2047:HOH:O	2.11	0.55
1:C:221:PRO:HB2	1:C:314:GLN:HB3	1.89	0.55
1:B:203:ILE:HG23	1:B:205:THR:HG23	1.88	0.55
1:A:61:ILE:HD11	1:A:119:PHE:CZ	2.42	0.55
1:B:55:ILE:CG2	1:B:95:SER:HB2	2.37	0.55
1:A:22:TYR:CE2	1:A:342:LYS:HE3	2.42	0.54
1:A:89:ASN:N	1:A:89:ASN:ND2	2.51	0.54
1:A:144:ARG:N	6:A:2041:HOH:O	2.41	0.54
1:D:100:ASN:O	1:D:101:PHE:C	2.50	0.54
1:C:153:LEU:O	1:C:157:MET:HB2	2.07	0.54
1:C:277:PRO:HG2	6:C:2039:HOH:O	2.07	0.54
3:F:2:NAG:H62	3:F:3:NAG:C1	2.37	0.54
1:C:301:ASP:O	1:C:304:ARG:HG3	2.08	0.54
1:C:341:LEU:HB2	1:C:346:LEU:HD12	1.90	0.54
1:A:57:SER:N	6:A:2010:HOH:O	2.41	0.54
1:C:126:PHE:O	1:C:129:THR:HG22	2.07	0.54
1:D:44:ASP:HB3	1:D:79:MET:HE2	1.90	0.54
1:B:230:SER:HB2	6:B:2042:HOH:O	2.06	0.54
1:C:29:THR:HB	1:C:31:TRP:CE2	2.42	0.54
1:C:105:ARG:HH21	1:C:105:ARG:HG2	1.73	0.54
1:D:343:ASP:O	1:D:345:GLN:HG3	2.07	0.54
1:A:344:ARG:NH2	6:A:2090:HOH:O	2.40	0.54
1:C:239:TYR:CD1	1:C:239:TYR:C	2.86	0.54
1:D:59:ALA:O	1:D:96:VAL:HG23	2.08	0.54
1:C:97:GLY:HA3	1:C:138:ALA:O	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:48:ARG:HH21	1:D:86:ARG:H	1.55	0.53
1:C:74:VAL:HG13	1:C:130:HIS:HE1	1.72	0.53
1:C:341:LEU:CB	1:C:346:LEU:HD12	2.39	0.53
1:A:352:TRP:CZ3	3:F:3:NAG:H83	2.44	0.53
1:A:341:LEU:HD21	1:A:349:ALA:HB2	1.89	0.53
1:B:124:PRO:HB3	1:B:164:GLU:HG3	1.90	0.53
1:B:225:GLY:HA2	1:B:313:GLY:HA3	1.90	0.53
1:C:165:ALA:HA	1:C:169:LYS:HG3	1.91	0.53
1:D:352:TRP:CE2	3:L:3:NAG:H5	2.43	0.53
1:A:116:ARG:NE	1:A:156:GLU:OE1	2.41	0.53
1:C:119:PHE:O	1:C:123:VAL:HG23	2.08	0.53
1:C:185:ILE:HD13	1:C:244:MET:CE	2.37	0.53
1:D:87:ASN:ND2	1:D:90:LEU:HB2	2.24	0.53
1:A:242:GLY:O	1:A:246:ARG:HB2	2.09	0.53
1:B:148:GLN:HB2	6:B:2030:HOH:O	2.07	0.53
1:B:203:ILE:HG23	1:B:203:ILE:O	2.07	0.53
1:D:259:PRO:HA	6:D:2061:HOH:O	2.08	0.53
1:A:96:VAL:CG1	1:A:135:LEU:HD11	2.39	0.53
1:A:96:VAL:HG12	1:A:135:LEU:HD11	1.91	0.53
1:A:204:MET:CE	3:F:3:NAG:C1	2.87	0.53
1:D:352:TRP:CZ2	3:L:3:NAG:H3	2.45	0.52
1:A:48:ARG:HB3	1:A:83:LEU:CD2	2.38	0.52
1:D:215:THR:HB	1:D:275:GLY:HA2	1.91	0.52
1:B:44:ASP:H	1:B:79:MET:CE	2.23	0.52
1:C:33:GLN:HG3	1:C:34:TYR:CE1	2.44	0.52
1:C:96:VAL:HG13	1:C:96:VAL:O	2.09	0.52
1:A:24:LEU:O	1:A:51:CYS:HB3	2.08	0.52
1:D:116:ARG:O	1:D:120:ILE:HG13	2.09	0.52
1:D:296:TYR:HA	1:D:299:ILE:HD12	1.92	0.52
1:D:92:THR:C	1:D:93:LEU:HD12	2.34	0.52
1:B:301:ASP:O	1:B:304:ARG:HG3	2.09	0.52
1:C:111:SER:HA	1:C:149:HIS:CD2	2.41	0.52
1:C:97:GLY:O	1:C:101:PHE:HB2	2.09	0.52
1:D:95:SER:HA	1:D:136:ASP:HB3	1.92	0.52
1:B:338:VAL:HA	1:B:341:LEU:HD23	1.92	0.51
1:B:365:GLY:C	1:B:367:ASP:H	2.16	0.51
1:D:117:ARG:O	1:D:121:LYS:HB2	2.09	0.51
1:C:251:ALA:HB1	1:C:345:GLN:O	2.09	0.51
1:C:55:ILE:HG21	1:C:95:SER:HB2	1.91	0.51
1:D:353:ALA:O	1:D:354:LEU:C	2.54	0.51
1:A:328:TYR:CD1	1:A:328:TYR:N	2.79	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:99:TRP:HB2	3:J:3:NAG:O3	2.10	0.51
1:D:352:TRP:CZ3	3:L:3:NAG:H83	2.46	0.51
1:C:60:ASN:CG	1:C:61:ILE:N	2.69	0.51
1:C:80:LEU:HD12	1:C:83:LEU:HD12	1.93	0.51
1:C:90:LEU:HD12	1:C:91:LYS:N	2.25	0.51
1:A:263:ARG:NH1	1:A:293:THR:OG1	2.43	0.51
1:B:191:ILE:O	1:B:195:SER:HB2	2.11	0.51
1:C:359:PHE:HD1	1:C:359:PHE:H	1.59	0.51
1:C:59:ALA:O	1:C:96:VAL:HG23	2.11	0.51
1:D:338:VAL:O	1:D:341:LEU:HG	2.11	0.51
1:D:225:GLY:HA2	1:D:313:GLY:HA3	1.93	0.50
1:D:341:LEU:HD12	1:D:342:LYS:N	2.25	0.50
1:A:255:VAL:CG1	1:A:350:MET:HB2	2.41	0.50
1:B:368:LEU:HD12	1:B:368:LEU:H	1.76	0.50
1:C:357:ASP:OD1	1:C:358:ASP:N	2.45	0.50
1:B:204:MET:HE2	3:H:2:NAG:H61	1.93	0.50
1:A:76:LEU:C	1:A:78:GLY:N	2.66	0.50
1:C:200:PHE:CD1	1:C:200:PHE:C	2.90	0.50
1:C:297:TYR:HB2	1:C:363:PHE:CG	2.45	0.50
1:D:144:ARG:NH1	1:D:187:SER:HB2	2.26	0.50
1:D:96:VAL:HG13	1:D:96:VAL:O	2.12	0.50
1:D:144:ARG:HH11	1:D:187:SER:HB2	1.75	0.50
1:A:76:LEU:HD22	6:A:2022:HOH:O	2.12	0.50
1:B:233:ARG:HG2	1:B:233:ARG:HH11	1.77	0.50
1:B:338:VAL:O	1:B:341:LEU:HG	2.11	0.50
1:C:77:TYR:O	1:C:81:ASN:ND2	2.45	0.50
1:D:267:LEU:HD13	1:D:272:THR:HG22	1.92	0.50
3:J:2:NAG:H62	3:J:3:NAG:C1	2.41	0.50
1:A:341:LEU:HD12	1:A:342:LYS:N	2.27	0.50
1:D:293:THR:CG2	1:D:294:LEU:N	2.75	0.50
1:B:370:PHE:O	1:B:374:ASN:HB2	2.12	0.49
1:B:85:ASN:C	1:B:87:ASN:H	2.20	0.49
1:B:150:PHE:CZ	1:B:154:ILE:HD11	2.47	0.49
1:C:166:GLN:N	1:C:167:PRO:CD	2.74	0.49
1:C:124:PRO:HB2	1:C:125:PRO:CD	2.42	0.49
1:C:246[A]:ARG:HH11	1:C:246[A]:ARG:HG3	1.76	0.49
1:D:100:ASN:ND2	3:L:4:NAG:H62	2.27	0.49
1:B:145:ARG:HH21	1:B:149:HIS:CE1	2.30	0.49
1:B:178:LEU:HG	1:B:201:ILE:HG23	1.93	0.49
1:B:204:MET:CE	3:H:3:NAG:C1	2.85	0.49
1:B:174:LEU:HD13	6:B:2025:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:185:ILE:HD13	1:A:244:MET:HE3	1.94	0.49
1:B:182:LYS:HD2	1:B:234:PHE:CD2	2.48	0.49
1:B:185:ILE:HA	1:B:189:TYR:CD2	2.48	0.49
1:C:162:ILE:HA	1:C:171:GLN:HE21	1.78	0.49
1:D:213:ARG:HD3	6:D:2027:HOH:O	2.12	0.49
1:A:341:LEU:HD12	1:A:341:LEU:C	2.38	0.49
1:D:343:ASP:C	1:D:345:GLN:N	2.69	0.49
1:C:215:THR:HB	1:C:275:GLY:HA2	1.95	0.48
1:A:43:PRO:HB2	1:A:79:MET:CB	2.44	0.48
1:B:303:LEU:HA	1:B:306:ALA:HB3	1.94	0.48
1:C:100:ASN:ND2	3:J:4:NAG:C6	2.77	0.48
1:D:234:PHE:HB3	1:D:239:TYR:CE1	2.48	0.48
1:C:86:ARG:C	1:C:88:PRO:HD3	2.38	0.48
1:D:368:LEU:HD12	1:D:369:ARG:N	2.28	0.48
1:B:144:ARG:HG2	1:B:187:SER:HB2	1.94	0.48
1:D:234:PHE:CB	1:D:239:TYR:CD1	2.97	0.48
1:C:71:TRP:H	1:C:71:TRP:CD1	2.30	0.48
1:B:22:TYR:CE2	1:B:342:LYS:HE3	2.48	0.48
1:A:57:SER:HB2	1:A:58:PHE:CG	2.48	0.48
1:A:100:ASN:ND2	3:F:4:NAG:H62	2.28	0.48
1:A:135:LEU:HG	1:A:136:ASP:N	2.29	0.48
1:A:47:ASP:OD1	1:A:50:LEU:N	2.46	0.48
1:B:261:PHE:HB3	1:B:356:LEU:HD13	1.96	0.48
1:C:341:LEU:HD12	1:C:342:LYS:HG3	1.94	0.48
1:C:352:TRP:CH2	3:J:3:NAG:H3	2.49	0.48
1:A:220:SER:O	1:A:316:VAL:HG11	2.14	0.48
1:A:263:ARG:HG3	6:A:2057:HOH:O	2.13	0.47
1:A:352:TRP:CE2	3:F:3:NAG:H5	2.49	0.47
1:D:48:ARG:NH1	1:D:83:LEU:HD23	2.29	0.47
1:A:211:ALA:O	1:A:278:ILE:HD11	2.15	0.47
1:B:185:ILE:HA	1:B:189:TYR:HD2	1.78	0.47
1:C:141:TYR:OH	3:J:2:NAG:O6	2.20	0.47
1:C:148:GLN:O	1:C:149:HIS:C	2.56	0.47
1:D:84:LYS:HE2	1:D:92:THR:HG23	1.96	0.47
1:A:139:TRP:C	1:A:140:LEU:HD23	2.39	0.47
1:D:286:ARG:HD3	6:D:2045:HOH:O	2.14	0.47
1:C:225:GLY:HA2	1:C:313:GLY:HA3	1.95	0.47
1:C:256:MET:CE	1:C:337:LYS:HB3	2.43	0.47
1:A:104:GLN:HG2	6:A:2032:HOH:O	2.13	0.47
1:B:92:THR:O	1:B:133:ASP:HB2	2.14	0.47
1:C:351:VAL:HG12	1:C:352:TRP:N	2.30	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:190:ASP:OD1	1:D:193:LYS:HB2	2.15	0.47
1:D:352:TRP:CH2	3:L:3:NAG:H3	2.49	0.47
1:D:52:THR:C	1:D:53:HIS:ND1	2.72	0.47
1:A:139:TRP:CE3	1:A:150:PHE:HD1	2.32	0.47
1:A:153:LEU:N	6:A:2047:HOH:O	2.47	0.47
1:A:178:LEU:HD22	1:A:189:TYR:CE1	2.50	0.47
1:A:204:MET:HE2	3:F:2:NAG:C6	2.40	0.47
1:B:285:GLY:HA3	1:B:298:GLU:OE2	2.15	0.47
1:D:166:GLN:N	1:D:167:PRO:CD	2.78	0.47
1:D:182:LYS:O	1:D:183:VAL:C	2.58	0.47
1:D:359:PHE:HB2	1:D:370:PHE:CZ	2.50	0.47
1:A:76:LEU:HB2	6:A:2022:HOH:O	2.14	0.47
1:A:215:THR:HA	1:A:276:ALA:O	2.15	0.47
1:B:174:LEU:HD12	1:B:175:SER:N	2.30	0.47
1:B:330:ASP:HA	1:B:372:LEU:HD21	1.97	0.47
1:C:48:ARG:CD	1:C:86:ARG:HB3	2.44	0.47
1:C:55:ILE:HG12	1:C:93:LEU:HB2	1.97	0.47
1:C:66:ILE:HG23	1:C:66:ILE:O	2.14	0.47
1:C:70:GLU:HB2	1:C:73:ASP:OD1	2.15	0.47
1:C:372:LEU:O	1:C:376:ILE:HG13	2.15	0.47
1:D:36:GLU:O	1:D:37:GLY:C	2.58	0.47
1:D:47:ASP:HB3	1:D:50:LEU:CB	2.37	0.47
1:C:64:ASP:OD1	1:C:115:SER:CB	2.63	0.47
1:C:293:THR:O	1:C:294:LEU:HD23	2.15	0.47
1:D:135:LEU:N	1:D:173:LEU:O	2.44	0.47
1:D:203:ILE:HD11	1:D:240:ALA:HB1	1.97	0.47
1:A:141:TYR:CE1	1:A:179:SER:HB2	2.49	0.46
1:A:48:ARG:CD	1:A:86:ARG:HB3	2.46	0.46
1:A:140:LEU:HA	1:A:141:TYR:HA	1.64	0.46
1:B:68:THR:HB	1:B:73:ASP:HB2	1.97	0.46
1:C:57:SER:HB2	1:C:58:PHE:CD1	2.50	0.46
1:C:62:SER:O	1:C:65:HIS:CE1	2.67	0.46
1:C:104:GLN:HE21	1:C:104:GLN:CA	2.16	0.46
1:D:206:TYR:O	1:D:207:ASP:HB2	2.14	0.46
1:D:366:GLN:O	1:D:367:ASP:HB2	2.15	0.46
1:B:209:HIS:CD2	1:B:217:GLY:HA3	2.51	0.46
1:C:247:LEU:HD23	1:C:247:LEU:HA	1.74	0.46
1:D:299:ILE:O	1:D:303:LEU:HG	2.15	0.46
1:B:60:ASN:CG	1:B:61:ILE:H	2.22	0.46
1:C:139:TRP:O	1:C:141:TYR:HA	2.14	0.46
1:C:209:HIS:NE2	1:C:213:ARG:NH1	2.64	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203:ILE:HG23	1:A:203:ILE:O	2.15	0.46
1:C:126:PHE:CD1	1:C:126:PHE:C	2.94	0.46
1:C:186:ASP:OD1	1:C:243:TYR:OH	2.30	0.46
1:C:135:LEU:HD21	1:C:137:LEU:HD21	1.98	0.46
1:D:261:PHE:HE1	6:D:2061:HOH:O	1.99	0.46
1:B:251:ALA:HB1	1:B:345:GLN:O	2.16	0.46
1:D:162:ILE:HA	1:D:171:GLN:HE21	1.79	0.46
1:D:252:SER:HA	6:D:2037:HOH:O	2.15	0.46
1:D:310:ARG:NH2	1:D:330:ASP:OD2	2.45	0.46
1:A:44:ASP:CG	1:A:79:MET:HE2	2.41	0.46
1:A:72:ASN:ND2	6:A:2022:HOH:O	2.49	0.46
1:C:116:ARG:O	1:C:117:ARG:C	2.59	0.46
1:A:153:LEU:HB3	6:A:2047:HOH:O	2.15	0.46
1:A:293:THR:HG22	1:A:294:LEU:N	2.30	0.46
1:A:365:GLY:O	1:A:366:GLN:HB2	2.16	0.46
1:C:55:ILE:HA	1:C:93:LEU:O	2.15	0.46
1:D:89:ASN:N	1:D:89:ASN:HD22	2.12	0.46
1:B:55:ILE:HG22	1:B:95:SER:HB2	1.97	0.46
1:C:95:SER:HA	1:C:136:ASP:HB3	1.98	0.46
1:D:57:SER:HB2	1:D:58:PHE:CD1	2.51	0.46
1:D:184:THR:O	1:D:185:ILE:C	2.58	0.46
1:D:340:TYR:O	1:D:343:ASP:HB2	2.16	0.46
1:D:351:VAL:HG12	1:D:352:TRP:N	2.31	0.46
6:A:2008:HOH:O	3:F:1:NAG:H2	2.16	0.45
1:B:66:ILE:HG23	1:B:66:ILE:O	2.17	0.45
1:B:263:ARG:HA	1:B:292:GLY:O	2.16	0.45
1:D:57:SER:HA	1:D:58:PHE:HA	1.52	0.45
1:D:123:VAL:CB	1:D:124:PRO:HD3	2.41	0.45
1:D:283:ILE:HA	1:D:284:PRO:HD3	1.76	0.45
1:A:263:ARG:HD2	1:A:292:GLY:O	2.16	0.45
1:C:328:TYR:CD1	1:C:328:TYR:N	2.84	0.45
1:A:72:ASN:HD22	1:A:76:LEU:HD13	1.81	0.45
1:A:377:LYS:HD3	6:A:2096:HOH:O	2.15	0.45
1:B:318:TYR:CD1	1:B:318:TYR:C	2.95	0.45
1:D:28:TYR:OH	1:D:43:PRO:HD3	2.17	0.45
1:A:365:GLY:C	1:A:367:ASP:N	2.71	0.45
6:A:2001:HOH:O	3:F:4:NAG:H3	2.17	0.45
1:D:243:TYR:O	1:D:246:ARG:HB3	2.17	0.45
1:A:146:ASP:O	1:A:147:LYS:C	2.60	0.45
1:B:181:GLY:O	1:B:185:ILE:HG13	2.17	0.45
1:B:350:MET:SD	1:B:350:MET:C	3.00	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:227:GLU:HG3	1:C:228:ASP:N	2.27	0.45
1:A:23:LYS:O	1:A:348:GLY:HA3	2.16	0.45
1:A:150:PHE:C	6:A:2047:HOH:O	2.55	0.45
1:A:190:ASP:O	1:A:191:ILE:C	2.59	0.45
1:B:252:SER:O	1:B:347:ALA:HB2	2.15	0.45
1:C:62:SER:HA	4:C:1384:NAG:H82	1.98	0.45
1:A:203:ILE:HD12	1:A:244:MET:SD	2.56	0.45
1:C:124:PRO:HD2	1:C:125:PRO:HD2	1.97	0.45
1:D:296:TYR:CD1	1:D:372:LEU:HD11	2.52	0.45
2:E:2:NAG:C6	2:E:3:NAG:H82	2.46	0.45
1:A:193:LYS:O	1:A:194:ILE:C	2.56	0.45
1:A:362:SER:O	1:A:364:CYS:N	2.50	0.45
1:C:57:SER:HA	1:C:58:PHE:HA	1.65	0.45
1:A:98:GLY:HA3	6:A:2030:HOH:O	2.16	0.44
1:A:160:GLU:OE1	1:A:160:GLU:HA	2.17	0.44
1:A:185:ILE:CD1	1:A:244:MET:HE3	2.47	0.44
1:D:71:TRP:CE2	2:K:2:NAG:H2	2.52	0.44
1:A:247:LEU:HD23	1:A:247:LEU:HA	1.68	0.44
1:C:249:ALA:HB1	1:C:254:LEU:HD13	2.00	0.44
1:A:26:CYS:HB3	1:A:354:LEU:HG	1.99	0.44
1:A:52:THR:OG1	1:A:53:HIS:ND1	2.45	0.44
1:A:88:PRO:C	1:A:89:ASN:HD22	2.25	0.44
1:A:151:THR:C	6:A:2047:HOH:O	2.60	0.44
1:B:315:GLN:HE22	1:C:315:GLN:NE2	2.16	0.44
1:C:112:ASN:HB3	1:C:115:SER:CB	2.46	0.44
1:C:308:VAL:HG12	1:C:309:HIS:N	2.32	0.44
1:C:339:GLN:O	1:C:343:ASP:OD1	2.35	0.44
1:C:352:TRP:CE2	3:J:3:NAG:H5	2.53	0.44
1:D:263:ARG:HA	1:D:292:GLY:O	2.18	0.44
1:C:178:LEU:HD22	1:C:189:TYR:CE1	2.52	0.44
1:A:144:ARG:HA	1:A:187:SER:O	2.17	0.44
1:B:57:SER:HB2	1:B:58:PHE:CD1	2.53	0.44
1:C:142:PRO:O	1:C:188:SER:CB	2.66	0.44
1:C:204:MET:CE	3:J:3:NAG:C7	2.95	0.44
1:C:309:HIS:HB2	1:C:318:TYR:CE2	2.53	0.44
1:C:368:LEU:HD12	1:C:368:LEU:O	2.18	0.44
1:A:139:TRP:C	1:A:139:TRP:CD1	2.96	0.44
1:B:203:ILE:O	1:B:203:ILE:CG2	2.66	0.44
1:B:44:ASP:OD1	1:B:79:MET:CE	2.65	0.44
1:B:57:SER:HA	1:B:58:PHE:HA	1.61	0.44
1:C:109:ILE:HG12	6:C:2010:HOH:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:203:ILE:HG23	1:C:203:ILE:O	2.18	0.44
1:C:234:PHE:HB3	1:C:239:TYR:CD1	2.52	0.44
1:C:255:VAL:HG12	1:C:350:MET:HB2	1.99	0.44
1:D:68:THR:HB	1:D:73:ASP:HB2	1.98	0.44
1:D:77:TYR:O	1:D:81:ASN:ND2	2.51	0.44
1:A:225:GLY:HA2	1:A:313:GLY:CA	2.45	0.43
1:C:48:ARG:HD3	1:C:86:ARG:HB3	2.00	0.43
1:C:267:LEU:CD1	1:C:272:THR:HG22	2.48	0.43
1:A:123:VAL:O	1:A:124:PRO:C	2.62	0.43
1:A:223:PHE:O	1:A:314:GLN:HG2	2.18	0.43
1:A:341:LEU:CD2	1:A:349:ALA:HB2	2.48	0.43
1:B:43:PRO:HB2	1:B:79:MET:CB	2.47	0.43
1:C:213:ARG:NH1	1:C:231:PRO:HD3	2.33	0.43
1:C:257:GLY:HA2	1:C:350:MET:HB3	2.00	0.43
1:D:369:ARG:H	1:D:369:ARG:HG3	1.68	0.43
1:B:70:GLU:N	1:B:73:ASP:OD2	2.51	0.43
1:C:70:GLU:H	1:C:73:ASP:CG	2.25	0.43
1:D:161:PHE:HD2	1:D:172:LEU:O	2.01	0.43
1:B:342:LYS:NZ	1:B:382:ALA:O	2.51	0.43
1:C:140:LEU:HA	1:C:141:TYR:HA	1.77	0.43
1:D:123:VAL:HG21	1:D:157:MET:HE1	1.99	0.43
1:B:166:GLN:HB2	1:B:167:PRO:HD3	2.00	0.43
1:C:107:SER:OG	1:C:145:ARG:NH2	2.51	0.43
1:C:140:LEU:HA	1:C:141:TYR:CD1	2.53	0.43
1:C:153:LEU:O	1:C:157:MET:CB	2.66	0.43
1:B:24:LEU:H	1:B:52:THR:HG1	1.64	0.43
1:B:123:VAL:HG11	1:B:135:LEU:HD22	2.01	0.43
1:C:44:ASP:HB3	1:C:79:MET:HE2	2.00	0.43
1:C:116:ARG:O	1:C:119:PHE:N	2.52	0.43
1:D:43:PRO:O	1:D:83:LEU:HD11	2.19	0.43
1:D:66:ILE:O	1:D:66:ILE:CG2	2.67	0.43
1:D:122:SER:O	1:D:125:PRO:HG2	2.18	0.43
1:A:26:CYS:CB	1:A:354:LEU:HD11	2.49	0.43
1:B:357:ASP:OD1	1:B:371:PRO:HD2	2.19	0.43
1:B:366:GLN:HB2	1:B:368:LEU:CD1	2.46	0.43
1:C:64:ASP:HB3	1:C:118:THR:HB	2.00	0.43
1:A:126:PHE:CD1	1:A:126:PHE:C	2.96	0.43
1:A:333:SER:O	1:A:337:LYS:HG3	2.19	0.43
1:B:338:VAL:O	1:B:342:LYS:HG3	2.19	0.43
1:D:56:TYR:OH	1:D:76:LEU:HD23	2.19	0.43
1:D:183:VAL:HG12	1:D:184:THR:N	2.33	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:219:HIS:CD2	1:A:263:ARG:HB2	2.54	0.43
1:B:57:SER:HB3	1:B:95:SER:HB3	1.99	0.43
1:B:185:ILE:CD1	1:B:244:MET:HE3	2.48	0.43
1:C:365:GLY:C	1:C:367:ASP:N	2.77	0.43
1:C:366:GLN:O	1:C:367:ASP:HB2	2.18	0.43
1:D:60:ASN:CG	1:D:61:ILE:N	2.77	0.43
1:C:187:SER:O	1:C:188:SER:HB3	2.19	0.42
1:C:252:SER:O	1:C:347:ALA:HB2	2.19	0.42
1:A:166:GLN:OE1	1:A:166:GLN:HA	2.19	0.42
1:C:234:PHE:HB3	1:C:239:TYR:CG	2.54	0.42
1:D:200:PHE:CD1	1:D:200:PHE:C	2.97	0.42
1:A:255:VAL:HG11	1:A:350:MET:HB2	2.02	0.42
1:C:352:TRP:HA	1:C:353:ALA:HA	1.63	0.42
1:C:356:LEU:HD23	1:C:356:LEU:HA	1.71	0.42
1:D:28:TYR:OH	1:D:42:PHE:HA	2.19	0.42
1:D:222:LEU:HG	1:D:223:PHE:CE2	2.54	0.42
1:A:57:SER:HB2	1:A:58:PHE:CE1	2.53	0.42
1:A:255:VAL:HG12	1:A:350:MET:HB2	2.00	0.42
1:B:97:GLY:HA3	1:B:138:ALA:O	2.19	0.42
1:D:116:ARG:NE	1:D:156:GLU:OE1	2.53	0.42
1:A:57:SER:HA	1:A:58:PHE:HA	1.64	0.42
1:B:60:ASN:CG	1:B:61:ILE:N	2.77	0.42
1:A:303:LEU:HD23	1:A:303:LEU:HA	1.86	0.42
1:C:94:LEU:HD22	1:C:132:PHE:CD2	2.54	0.42
1:C:164:GLU:OE2	1:C:172:LEU:HD11	2.20	0.42
1:C:264:SER:HA	1:C:325:TRP:O	2.18	0.42
1:D:54:ILE:CG1	1:D:90:LEU:HD11	2.46	0.42
1:D:57:SER:HB2	1:D:58:PHE:CG	2.54	0.42
1:D:105:ARG:O	1:D:109:ILE:HG13	2.19	0.42
1:C:76:LEU:O	1:C:77:TYR:C	2.61	0.42
1:D:140:LEU:HA	1:D:141:TYR:CD1	2.54	0.42
1:A:86:ARG:NH2	6:A:2025:HOH:O	2.52	0.42
1:A:135:LEU:CD2	1:A:157:MET:HE3	2.50	0.42
1:C:26:CYS:HB2	1:C:354:LEU:HD11	2.02	0.42
1:D:191:ILE:HG21	1:D:249:ALA:HB2	2.01	0.42
1:B:365:GLY:C	1:B:367:ASP:N	2.78	0.42
1:C:43:PRO:HB2	1:C:79:MET:CB	2.49	0.42
1:C:66:ILE:HG21	1:C:123:VAL:HA	2.02	0.42
1:C:230:SER:HA	1:C:231:PRO:HD3	1.88	0.42
1:D:124:PRO:O	1:D:128:ARG:HG2	2.19	0.42
1:D:254:LEU:HA	1:D:254:LEU:HD12	1.80	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:GLN:HG2	1:A:104:GLN:H	1.49	0.41
1:B:194:ILE:HG13	1:B:195:SER:N	2.34	0.41
1:C:135:LEU:HG	1:C:136:ASP:N	2.35	0.41
1:C:182:LYS:O	1:C:183:VAL:C	2.63	0.41
1:D:163:LYS:HA	1:D:166:GLN:NE2	2.26	0.41
1:D:381:ALA:O	1:D:382:ALA:CB	2.68	0.41
1:D:52:THR:C	1:D:53:HIS:HD1	2.28	0.41
1:D:60:ASN:CG	1:D:61:ILE:H	2.27	0.41
1:D:352:TRP:HZ3	3:L:3:NAG:H83	1.86	0.41
1:A:362:SER:OG	1:A:363:PHE:CD1	2.72	0.41
1:B:366:GLN:CB	1:B:368:LEU:HD12	2.48	0.41
1:C:100:ASN:CG	3:J:4:NAG:H62	2.45	0.41
1:C:293:THR:HG22	1:C:294:LEU:N	2.35	0.41
1:D:71:TRP:CD1	1:D:71:TRP:N	2.88	0.41
1:D:121:LYS:O	1:D:124:PRO:HD2	2.20	0.41
1:D:222:LEU:O	1:D:222:LEU:HD12	2.21	0.41
1:A:263:ARG:HD2	1:A:263:ARG:HA	1.96	0.41
1:B:314:GLN:O	1:B:315:GLN:HB2	2.20	0.41
1:D:204:MET:CE	3:L:3:NAG:C1	2.93	0.41
1:B:243:TYR:CE2	1:B:247:LEU:HD11	2.56	0.41
1:A:33:GLN:HG3	1:A:34:TYR:CD1	2.54	0.41
1:A:232:ASP:HB2	1:A:233:ARG:H	1.69	0.41
1:B:243:TYR:CZ	1:B:247:LEU:HD11	2.56	0.41
1:C:52:THR:C	1:C:53:HIS:ND1	2.78	0.41
1:C:218:HIS:NE2	1:C:316:VAL:HB	2.36	0.41
1:C:256:MET:HE1	1:C:337:LYS:HB3	2.03	0.41
1:C:267:LEU:HD13	1:C:272:THR:HG22	2.03	0.41
1:D:328:TYR:HD2	1:D:329:ASP:O	2.03	0.41
1:D:356:LEU:HD23	1:D:356:LEU:HA	1.87	0.41
1:B:193:LYS:HA	1:B:196:GLN:OE1	2.20	0.41
1:B:96:VAL:O	1:B:96:VAL:CG1	2.66	0.41
1:C:58:PHE:CD2	1:C:98:GLY:HA3	2.56	0.41
1:C:71:TRP:CZ2	2:I:2:NAG:H2	2.55	0.41
1:C:175:SER:CB	1:C:200:PHE:CE1	3.02	0.41
1:A:150:PHE:CZ	1:A:154:ILE:HD11	2.56	0.41
1:A:372:LEU:HD23	1:A:372:LEU:HA	1.82	0.41
1:B:139:TRP:C	1:B:140:LEU:HD23	2.46	0.41
1:B:214:GLY:C	1:B:278:ILE:HG12	2.46	0.41
1:B:335:LYS:HG2	1:B:339:GLN:NE2	2.35	0.41
1:C:94:LEU:HD22	1:C:132:PHE:CE2	2.56	0.41
1:C:254:LEU:HD23	1:C:346:LEU:HD22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:36:GLU:O	1:D:37:GLY:O	2.39	0.41
1:D:128:ARG:C	1:D:130:HIS:H	2.27	0.41
1:D:199:ASP:O	1:D:200:PHE:HB3	2.20	0.41
1:D:209:HIS:CD2	1:D:217:GLY:HA3	2.56	0.41
1:D:333:SER:O	1:D:337:LYS:HG3	2.21	0.41
1:D:33:GLN:NE2	1:D:76:LEU:HD21	2.35	0.41
1:D:140:LEU:HA	1:D:141:TYR:HA	1.80	0.41
1:B:140:LEU:HA	1:B:141:TYR:HA	1.78	0.40
1:C:286:ARG:O	1:C:289:LYS:HG3	2.21	0.40
1:B:255:VAL:HG12	1:B:350:MET:HB2	2.03	0.40
1:D:64:ASP:OD1	1:D:115:SER:HB3	2.21	0.40
1:A:123:VAL:HB	1:A:124:PRO:HD3	2.03	0.40
1:C:66:ILE:HB	1:C:119:PHE:HE1	1.86	0.40
1:D:255:VAL:HA	1:D:348:GLY:O	2.22	0.40
1:D:296:TYR:CG	1:D:372:LEU:HD11	2.57	0.40
3:L:3:NAG:H61	3:L:4:NAG:C7	2.50	0.40
1:A:62:SER:O	1:A:63:ASN:CB	2.70	0.40
1:A:382:ALA:O	1:A:383:THR:HG23	2.21	0.40
1:B:226:GLN:HG3	1:C:238:ASP:OD1	2.20	0.40
1:C:293:THR:CG2	1:C:294:LEU:N	2.85	0.40
1:C:297:TYR:HB2	1:C:363:PHE:CD1	2.56	0.40
1:D:293:THR:HG22	1:D:294:LEU:N	2.37	0.40
1:B:152:THR:O	1:B:156:GLU:HG3	2.21	0.40
1:B:250:PRO:O	1:B:251:ALA:C	2.65	0.40
1:C:34:TYR:CZ	2:I:1:NAG:H3	2.57	0.40
1:C:234:PHE:HB3	1:C:239:TYR:CE1	2.57	0.40
1:D:46:LEU:HD23	1:D:46:LEU:N	2.36	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	362/362 (100%)	329 (91%)	29 (8%)	4 (1%)	11	22
1	B	361/362 (100%)	336 (93%)	24 (7%)	1 (0%)	36	56
1	C	362/362 (100%)	309 (85%)	44 (12%)	9 (2%)	4	8
1	D	361/362 (100%)	304 (84%)	47 (13%)	10 (3%)	4	6
All	All	1446/1448 (100%)	1278 (88%)	144 (10%)	24 (2%)	7	14

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	362	SER
1	A	363	PHE
1	C	66	ILE
1	C	68	THR
1	C	188	SER
1	D	36	GLU
1	D	37	GLY
1	A	231	PRO
1	C	63	ASN
1	D	101	PHE
1	D	114	GLN
1	D	291	ALA
1	D	312	LEU
1	D	382	ALA
1	B	366	GLN
1	C	204	MET
1	D	63	ASN
1	D	289	LYS
1	A	317	PRO
1	C	168	GLY
1	C	207	ASP
1	C	187	SER
1	C	231	PRO
1	D	183	VAL

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	304/302 (101%)	285 (94%)	19 (6%)	16	30
1	B	303/302 (100%)	285 (94%)	18 (6%)	18	33
1	C	304/302 (101%)	279 (92%)	25 (8%)	10	20
1	D	303/302 (100%)	289 (95%)	14 (5%)	24	45
All	All	1214/1208 (100%)	1138 (94%)	76 (6%)	17	30

All (76) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	89	ASN
1	A	96	VAL
1	A	104	GLN
1	A	106	PHE
1	A	108	LYS
1	A	114[A]	GLN
1	A	114[B]	GLN
1	A	144	ARG
1	A	145	ARG
1	A	148	GLN
1	A	167	PRO
1	A	170	LYS
1	A	259	PRO
1	A	263	ARG
1	A	317	PRO
1	A	329	ASP
1	A	331	GLN
1	A	336	SER
1	A	383	THR
1	B	23	LYS
1	B	43	PRO
1	B	76	LEU
1	B	103	SER
1	B	105	ARG
1	B	129	THR
1	B	148	GLN
1	B	167	PRO
1	B	171	GLN
1	B	183	VAL
1	B	227	GLU
1	B	244	MET
1	B	259	PRO
1	B	301	ASP

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Mol	Chain	Res	Type
1	B	329	ASP
1	B	339	GLN
1	B	364	CYS
1	B	368	LEU
1	C	67	ASP
1	C	73	ASP
1	C	90	LEU
1	C	104	GLN
1	C	106	PHE
1	C	129	THR
1	C	144	ARG
1	C	164	GLU
1	C	169	LYS
1	C	170	LYS
1	C	207	ASP
1	C	209	HIS
1	C	222	LEU
1	C	227	GLU
1	C	244	MET
1	C	246[A]	ARG
1	C	246[B]	ARG
1	C	259	PRO
1	C	263	ARG
1	C	307	THR
1	C	331	GLN
1	C	335	LYS
1	C	366	GLN
1	C	373	THR
1	C	383	THR
1	D	67	ASP
1	D	76	LEU
1	D	106	PHE
1	D	118	THR
1	D	121	LYS
1	D	129	THR
1	D	152	THR
1	D	164	GLU
1	D	170	LYS
1	D	235	SER
1	D	317	PRO
1	D	331	GLN
1	D	368	LEU

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Mol	Chain	Res	Type
1	D	369	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (35) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	85	ASN
1	A	87	ASN
1	A	89	ASN
1	A	100	ASN
1	A	315	GLN
1	A	331	GLN
1	B	85	ASN
1	B	89	ASN
1	B	148	GLN
1	B	149	HIS
1	B	339	GLN
1	C	65	HIS
1	C	85	ASN
1	C	87	ASN
1	C	89	ASN
1	C	100	ASN
1	C	104	GLN
1	C	130	HIS
1	C	171	GLN
1	C	196	GLN
1	C	197	HIS
1	C	209	HIS
1	C	315	GLN
1	C	331	GLN
1	C	366	GLN
1	D	33	GLN
1	D	81	ASN
1	D	85	ASN
1	D	89	ASN
1	D	166	GLN
1	D	171	GLN
1	D	196	GLN
1	D	197	HIS
1	D	315	GLN
1	D	331	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 ⓘ

28 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
2	NAG	E	1	2	15,15,15	0.52	0	21,21,21	0.87	0
2	NAG	E	2	2	14,14,15	0.58	0	17,19,21	0.75	1 (5%)
2	NAG	E	3	2	14,14,15	0.61	0	17,19,21	0.63	0
3	NAG	F	1	3	15,15,15	0.64	0	21,21,21	0.92	1 (4%)
3	NAG	F	2	3	14,14,15	0.68	0	17,19,21	0.86	1 (5%)
3	NAG	F	3	3	14,14,15	0.79	0	17,19,21	0.81	0
3	NAG	F	4	3	14,14,15	0.63	0	17,19,21	0.72	0
2	NAG	G	1	2	15,15,15	0.51	0	21,21,21	0.68	0
2	NAG	G	2	2	14,14,15	0.74	0	17,19,21	0.79	0
2	NAG	G	3	2	14,14,15	0.69	0	17,19,21	0.68	0
3	NAG	H	1	3	15,15,15	0.48	0	21,21,21	0.62	0
3	NAG	H	2	3	14,14,15	0.58	0	17,19,21	0.87	1 (5%)
3	NAG	H	3	3	14,14,15	0.91	0	17,19,21	0.98	2 (11%)
3	NAG	H	4	3	14,14,15	0.55	0	17,19,21	0.67	0
2	NAG	I	1	2	15,15,15	0.58	0	21,21,21	0.60	0
2	NAG	I	2	2	14,14,15	0.71	0	17,19,21	0.81	1 (5%)
2	NAG	I	3	2	14,14,15	0.63	0	17,19,21	0.66	0
3	NAG	J	1	3	15,15,15	0.57	0	21,21,21	0.94	1 (4%)
3	NAG	J	2	3	14,14,15	0.76	0	17,19,21	0.90	1 (5%)
3	NAG	J	3	3	14,14,15	0.77	0	17,19,21	1.08	1 (5%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	J	4	3	14,14,15	0.66	0	17,19,21	0.88	1 (5%)
2	NAG	K	1	2	15,15,15	0.68	0	21,21,21	0.95	1 (4%)
2	NAG	K	2	2	14,14,15	0.66	0	17,19,21	0.72	1 (5%)
2	NAG	K	3	2	14,14,15	0.64	0	17,19,21	0.69	0
3	NAG	L	1	3	15,15,15	0.58	0	21,21,21	0.73	0
3	NAG	L	2	3	14,14,15	0.58	0	17,19,21	0.78	1 (5%)
3	NAG	L	3	3	14,14,15	0.85	0	17,19,21	1.06	2 (11%)
3	NAG	L	4	3	14,14,15	0.67	0	17,19,21	0.60	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	E	1	2	-	4/6/26/26	0/1/1/1
2	NAG	E	2	2	-	2/6/23/26	0/1/1/1
2	NAG	E	3	2	-	3/6/23/26	0/1/1/1
3	NAG	F	1	3	-	2/6/26/26	0/1/1/1
3	NAG	F	2	3	-	2/6/23/26	0/1/1/1
3	NAG	F	3	3	-	2/6/23/26	0/1/1/1
3	NAG	F	4	3	-	0/6/23/26	0/1/1/1
2	NAG	G	1	2	-	5/6/26/26	0/1/1/1
2	NAG	G	2	2	-	2/6/23/26	0/1/1/1
2	NAG	G	3	2	-	3/6/23/26	0/1/1/1
3	NAG	H	1	3	-	1/6/26/26	0/1/1/1
3	NAG	H	2	3	-	2/6/23/26	0/1/1/1
3	NAG	H	3	3	-	2/6/23/26	0/1/1/1
3	NAG	H	4	3	-	2/6/23/26	0/1/1/1
2	NAG	I	1	2	-	6/6/26/26	0/1/1/1
2	NAG	I	2	2	-	2/6/23/26	0/1/1/1
2	NAG	I	3	2	-	3/6/23/26	0/1/1/1
3	NAG	J	1	3	-	4/6/26/26	0/1/1/1
3	NAG	J	2	3	-	2/6/23/26	0/1/1/1
3	NAG	J	3	3	-	2/6/23/26	0/1/1/1
3	NAG	J	4	3	-	4/6/23/26	0/1/1/1
2	NAG	K	1	2	-	4/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	K	2	2	-	4/6/23/26	0/1/1/1
2	NAG	K	3	2	-	3/6/23/26	0/1/1/1
3	NAG	L	1	3	-	4/6/26/26	0/1/1/1
3	NAG	L	2	3	-	4/6/23/26	0/1/1/1
3	NAG	L	3	3	-	0/6/23/26	0/1/1/1
3	NAG	L	4	3	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	J	3	NAG	C1-O5-C5	2.76	115.89	112.19
3	F	2	NAG	C2-N2-C7	-2.73	119.25	122.90
3	H	3	NAG	C2-N2-C7	-2.60	119.42	122.90
3	J	4	NAG	C2-N2-C7	-2.59	119.42	122.90
2	K	1	NAG	C1-C2-C3	-2.57	107.04	110.54
3	L	3	NAG	C2-N2-C7	-2.57	119.46	122.90
3	J	2	NAG	C2-N2-C7	-2.51	119.53	122.90
2	E	2	NAG	C2-N2-C7	-2.48	119.58	122.90
3	H	2	NAG	C2-N2-C7	-2.45	119.62	122.90
2	I	2	NAG	C2-N2-C7	-2.34	119.76	122.90
3	L	3	NAG	C1-O5-C5	2.27	115.22	112.19
3	J	1	NAG	C4-C3-C2	2.24	113.66	110.40
3	H	3	NAG	C1-O5-C5	2.23	115.17	112.19
3	L	2	NAG	C2-N2-C7	-2.18	119.97	122.90
3	F	1	NAG	C3-C4-C5	2.18	114.18	110.23
2	K	2	NAG	C2-N2-C7	-2.10	120.09	122.90

There are no chirality outliers.

All (74) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O7-C7-N2-C2
2	E	2	NAG	C8-C7-N2-C2
2	E	2	NAG	O7-C7-N2-C2
2	G	1	NAG	C8-C7-N2-C2
2	G	1	NAG	O7-C7-N2-C2
2	G	2	NAG	C8-C7-N2-C2
2	G	2	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
2	G	3	NAG	C8-C7-N2-C2
2	G	3	NAG	O7-C7-N2-C2
2	I	1	NAG	C1-C2-N2-C7
2	I	1	NAG	C8-C7-N2-C2
2	I	1	NAG	O7-C7-N2-C2
2	I	3	NAG	C8-C7-N2-C2
2	I	3	NAG	O7-C7-N2-C2
2	K	1	NAG	C8-C7-N2-C2
2	K	1	NAG	O7-C7-N2-C2
2	K	2	NAG	C8-C7-N2-C2
2	K	2	NAG	O7-C7-N2-C2
2	K	3	NAG	C8-C7-N2-C2
2	K	3	NAG	O7-C7-N2-C2
3	F	2	NAG	C8-C7-N2-C2
3	F	2	NAG	O7-C7-N2-C2
3	J	2	NAG	C8-C7-N2-C2
3	J	2	NAG	O7-C7-N2-C2
3	L	2	NAG	C8-C7-N2-C2
3	L	2	NAG	O7-C7-N2-C2
3	J	4	NAG	O5-C5-C6-O6
2	I	1	NAG	C4-C5-C6-O6
3	H	2	NAG	O5-C5-C6-O6
2	E	1	NAG	C4-C5-C6-O6
3	L	1	NAG	C4-C5-C6-O6
3	F	1	NAG	O5-C5-C6-O6
3	J	4	NAG	C4-C5-C6-O6
3	F	1	NAG	C4-C5-C6-O6
3	J	1	NAG	C8-C7-N2-C2
2	E	1	NAG	O5-C5-C6-O6
2	I	1	NAG	O5-C5-C6-O6
3	L	1	NAG	O5-C5-C6-O6
3	H	3	NAG	C8-C7-N2-C2
3	J	1	NAG	O7-C7-N2-C2
3	J	4	NAG	C8-C7-N2-C2
3	J	4	NAG	O7-C7-N2-C2
3	H	2	NAG	C4-C5-C6-O6
2	E	3	NAG	C8-C7-N2-C2
2	I	2	NAG	C8-C7-N2-C2
3	L	1	NAG	C8-C7-N2-C2
3	H	4	NAG	O5-C5-C6-O6
3	L	2	NAG	O5-C5-C6-O6
2	E	3	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	H	3	NAG	O7-C7-N2-C2
3	L	1	NAG	O7-C7-N2-C2
3	J	3	NAG	C4-C5-C6-O6
2	I	3	NAG	O5-C5-C6-O6
3	F	3	NAG	C4-C5-C6-O6
2	I	2	NAG	O7-C7-N2-C2
2	G	1	NAG	C1-C2-N2-C7
2	K	3	NAG	O5-C5-C6-O6
3	F	3	NAG	O5-C5-C6-O6
2	G	3	NAG	O5-C5-C6-O6
2	E	3	NAG	O5-C5-C6-O6
3	J	1	NAG	C4-C5-C6-O6
3	J	3	NAG	O5-C5-C6-O6
2	K	1	NAG	O5-C5-C6-O6
3	H	4	NAG	C4-C5-C6-O6
2	K	2	NAG	C1-C2-N2-C7
3	L	2	NAG	C4-C5-C6-O6
2	K	2	NAG	C3-C2-N2-C7
3	H	1	NAG	O5-C5-C6-O6
2	K	1	NAG	C1-C2-N2-C7
3	J	1	NAG	O5-C5-C6-O6
2	G	1	NAG	C3-C2-N2-C7
2	I	1	NAG	C3-C2-N2-C7
2	G	1	NAG	O5-C5-C6-O6

There are no ring outliers.

19 monomers are involved in 51 short contacts:

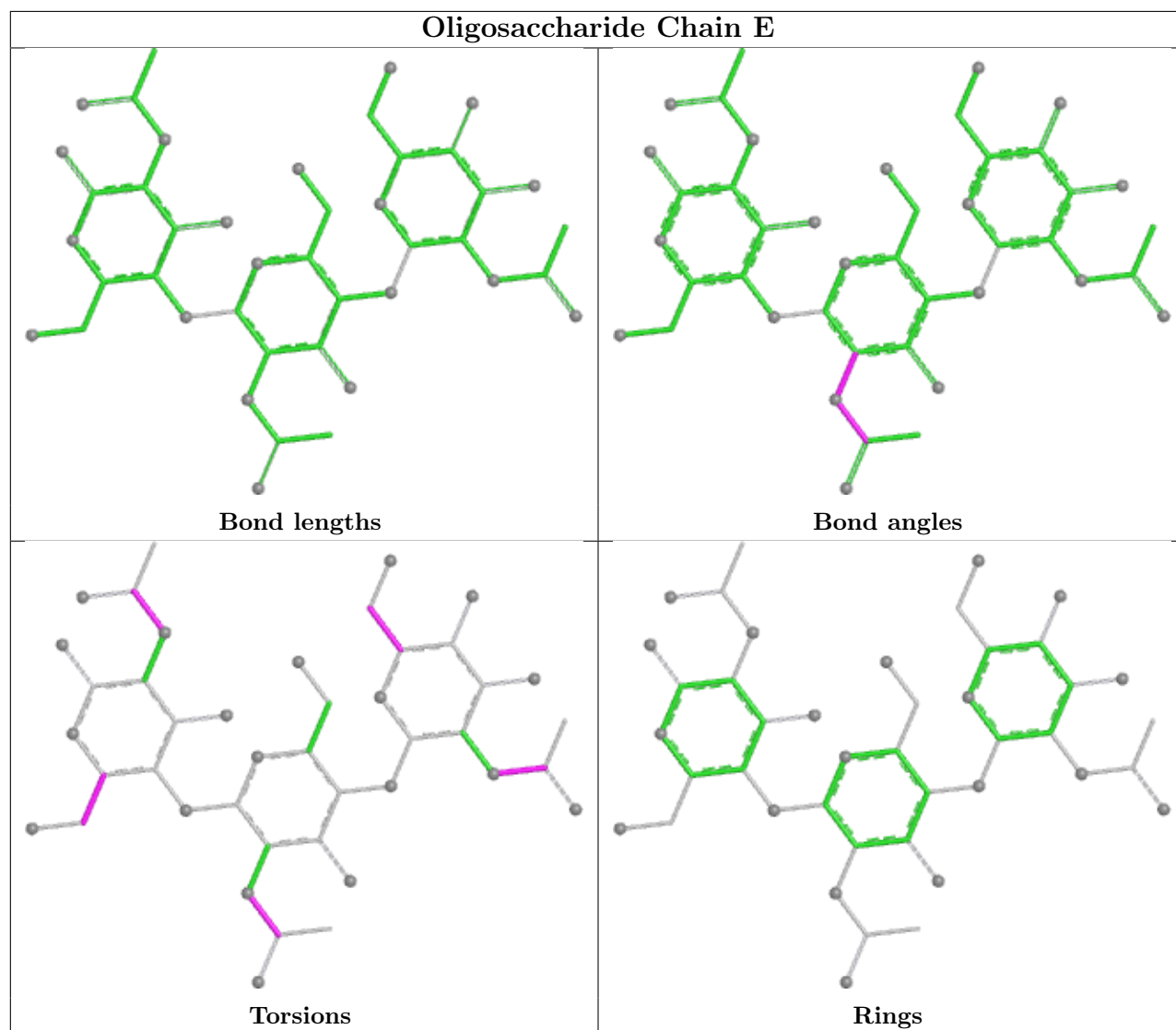
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	J	4	NAG	5	0
3	F	1	NAG	1	0
3	F	2	NAG	4	0
2	K	2	NAG	1	0
3	F	4	NAG	2	0
3	F	3	NAG	5	0
3	J	2	NAG	3	0
3	L	2	NAG	2	0
2	I	1	NAG	1	0
3	H	2	NAG	1	0
3	L	4	NAG	2	0
3	J	3	NAG	10	0
2	E	2	NAG	3	0

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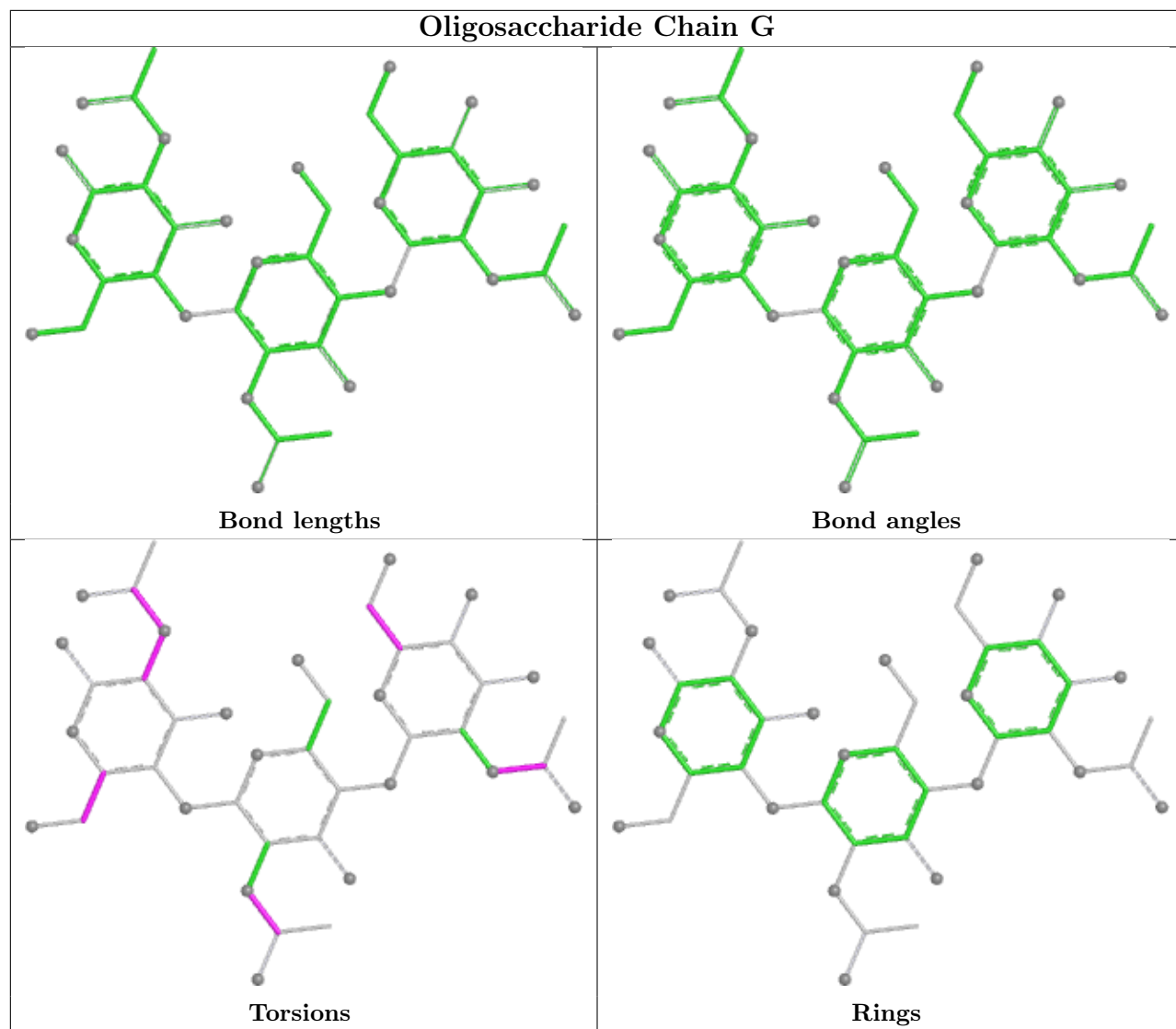
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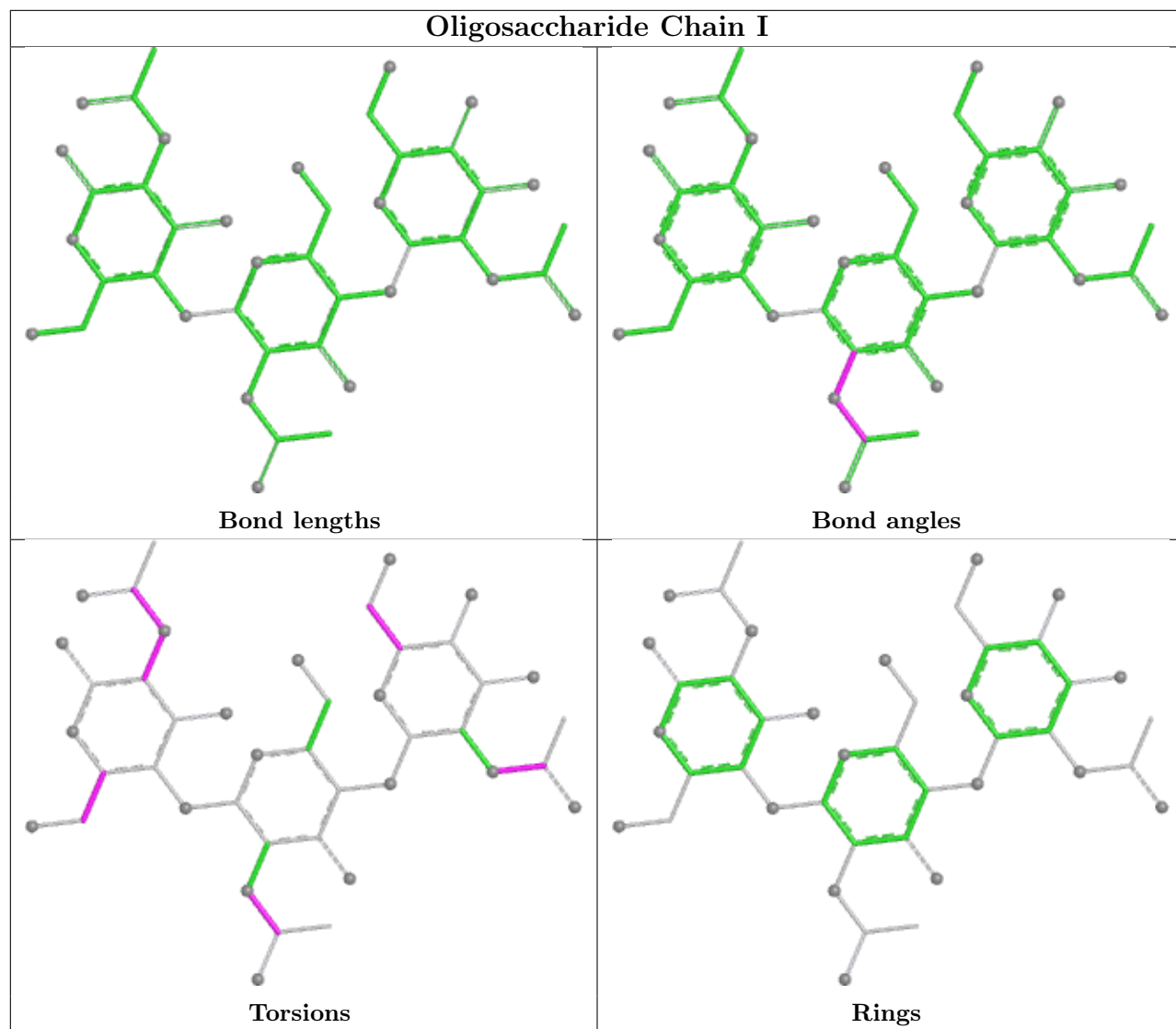
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	H	4	NAG	2	0
2	I	3	NAG	2	0
3	L	3	NAG	8	0
2	E	3	NAG	3	0
2	I	2	NAG	3	0
3	H	3	NAG	3	0

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

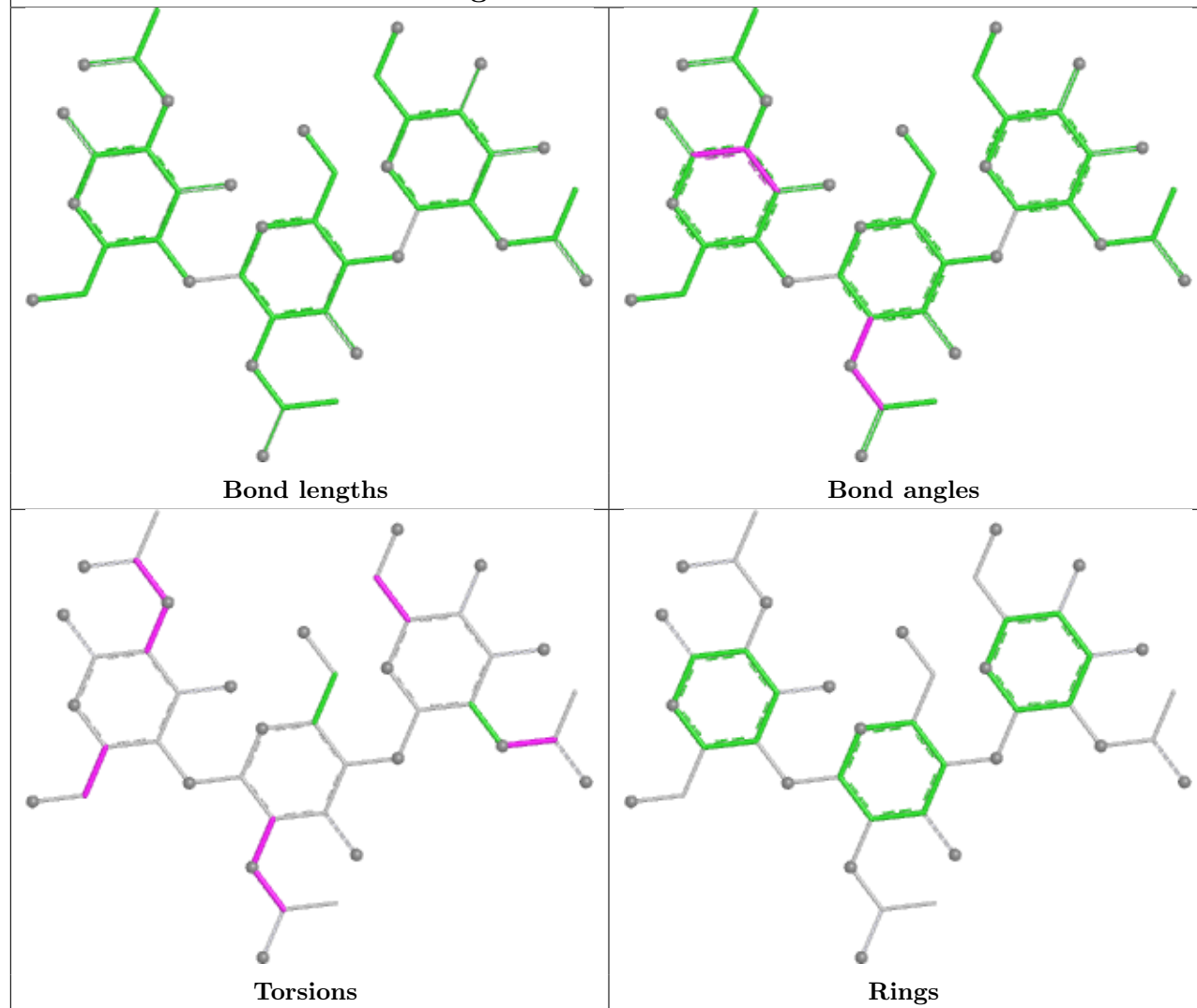


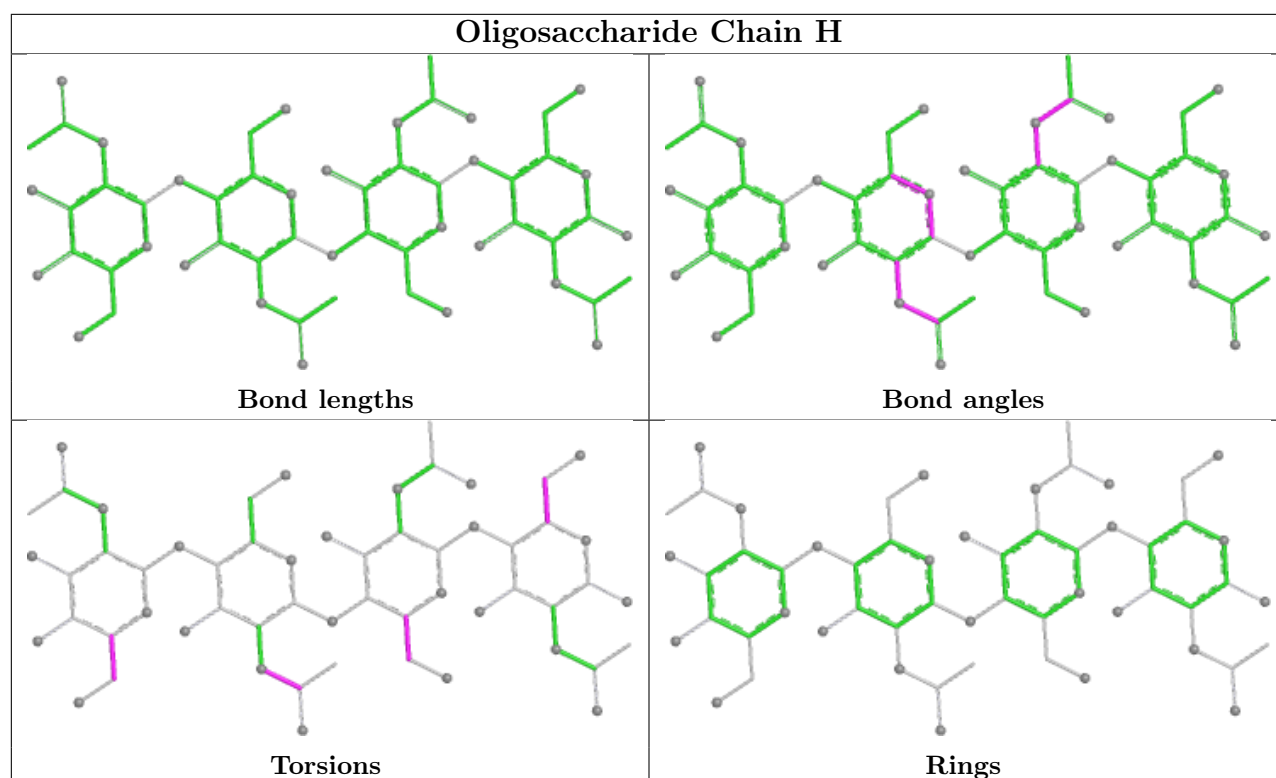
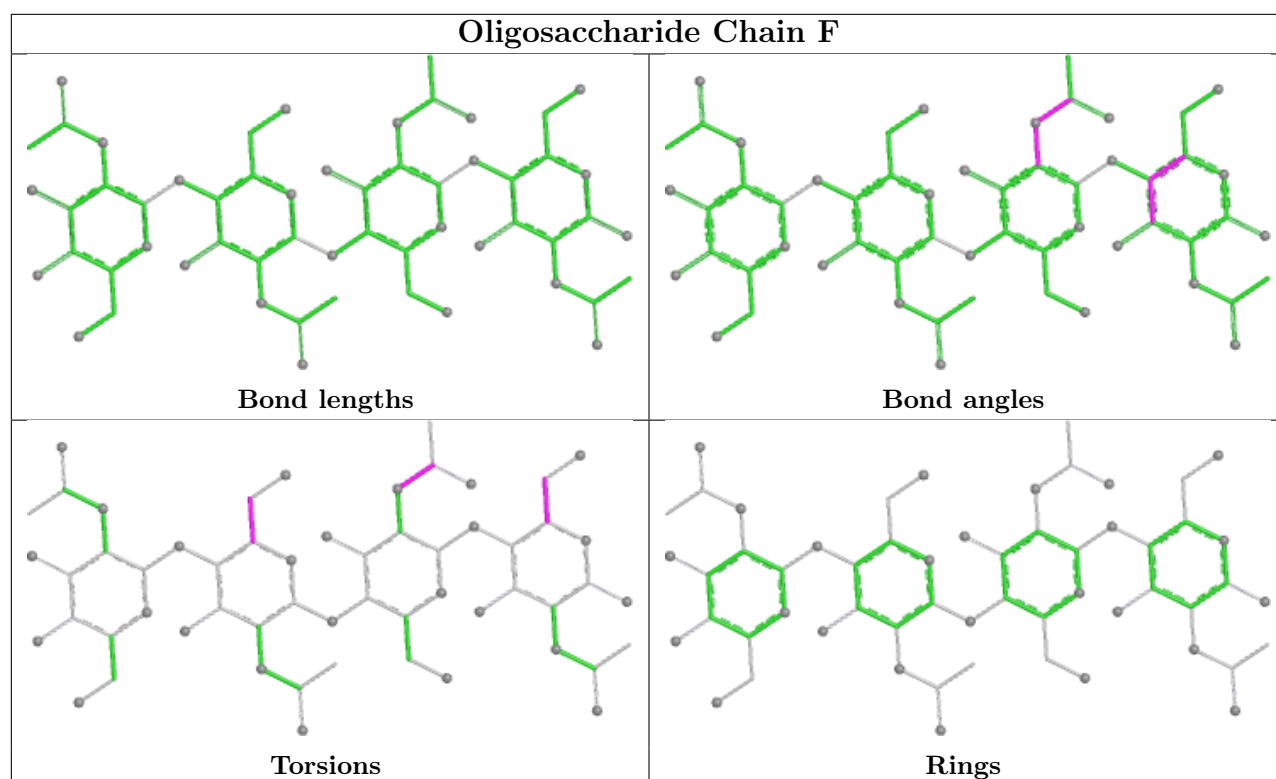
Oligosaccharide Chain G

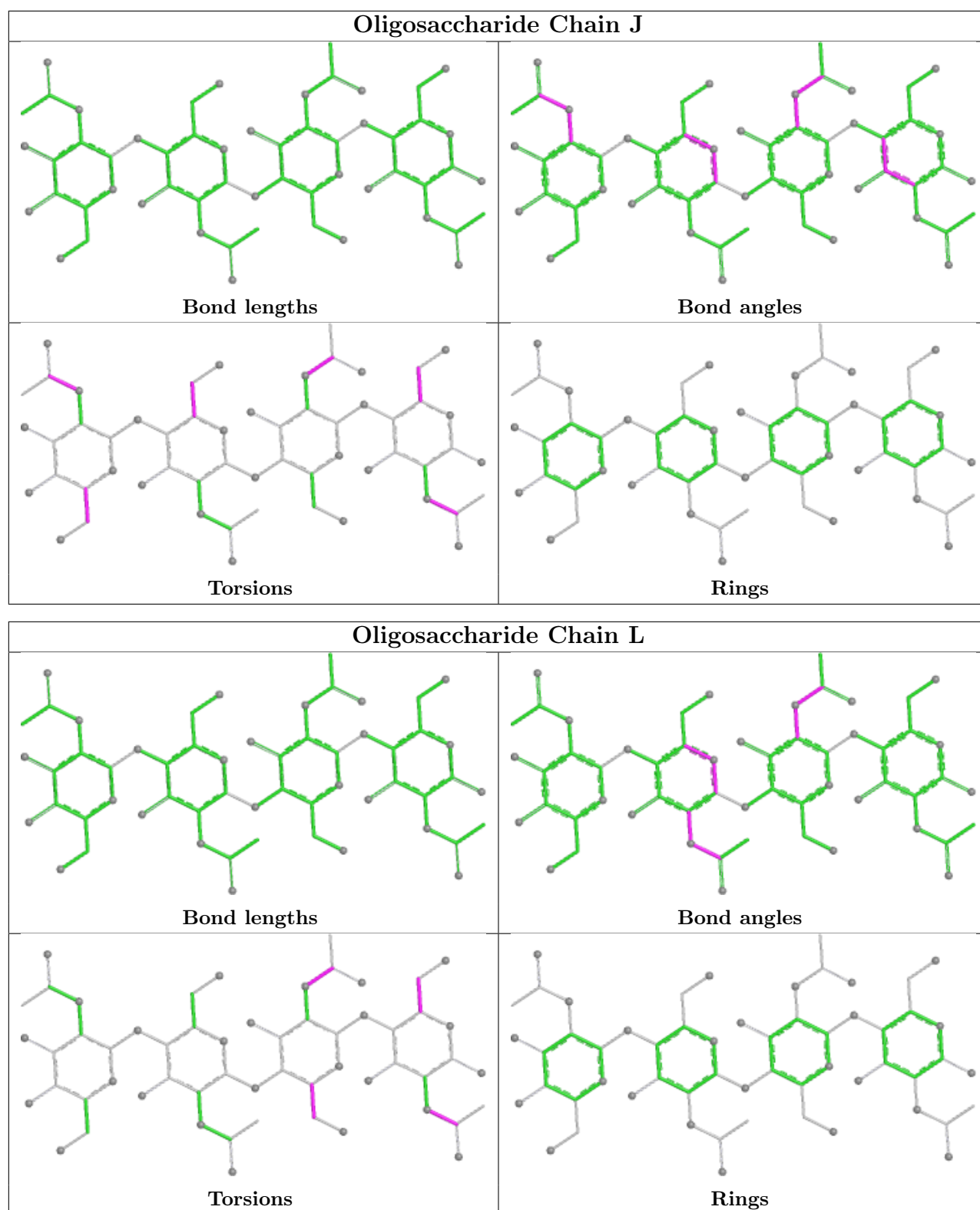




Oligosaccharide Chain K







5.6 Ligand geometry [i](#)

16 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	$\# Z > 2$
5	SO4	A	1387	-	4,4,4	0.41	0	6,6,6	0.13	0
4	NAG	C	1384	1	14,14,15	0.85	1 (7%)	17,19,21	0.69	0
4	NAG	B	1384	1	14,14,15	0.79	0	17,19,21	0.81	1 (5%)
5	SO4	A	1385	-	4,4,4	0.36	0	6,6,6	0.17	0
4	NAG	D	1384	1	14,14,15	0.67	0	17,19,21	0.66	0
5	SO4	C	1387	-	4,4,4	0.39	0	6,6,6	0.10	0
5	SO4	A	1386	-	4,4,4	0.40	0	6,6,6	0.23	0
5	SO4	B	1386	-	4,4,4	0.37	0	6,6,6	0.19	0
4	NAG	A	1384	1	14,14,15	0.74	0	17,19,21	1.11	1 (5%)
5	SO4	C	1385	-	4,4,4	0.39	0	6,6,6	0.04	0
5	SO4	D	1385	-	4,4,4	0.37	0	6,6,6	0.21	0
5	SO4	D	1386	-	4,4,4	0.36	0	6,6,6	0.13	0
5	SO4	B	1385	-	4,4,4	0.39	0	6,6,6	0.05	0
5	SO4	B	1387	-	4,4,4	0.42	0	6,6,6	0.11	0
5	SO4	D	1387	-	4,4,4	0.37	0	6,6,6	0.20	0
5	SO4	C	1386	-	4,4,4	0.35	0	6,6,6	0.09	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	B	1384	1	-	4/6/23/26	0/1/1/1
4	NAG	C	1384	1	-	5/6/23/26	0/1/1/1
4	NAG	A	1384	1	-	3/6/23/26	0/1/1/1
4	NAG	D	1384	1	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	1384	NAG	C1-C2	2.44	1.55	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1384	NAG	C2-N2-C7	-2.73	119.25	122.90
4	B	1384	NAG	C2-N2-C7	-2.13	120.05	122.90

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	C	1384	NAG	C3-C2-N2-C7
4	C	1384	NAG	C8-C7-N2-C2
4	C	1384	NAG	O7-C7-N2-C2
4	D	1384	NAG	C8-C7-N2-C2
4	D	1384	NAG	O7-C7-N2-C2
4	C	1384	NAG	C4-C5-C6-O6
4	C	1384	NAG	O5-C5-C6-O6
4	B	1384	NAG	C8-C7-N2-C2
4	B	1384	NAG	O7-C7-N2-C2
4	A	1384	NAG	C8-C7-N2-C2
4	B	1384	NAG	C4-C5-C6-O6
4	B	1384	NAG	O5-C5-C6-O6
4	A	1384	NAG	O7-C7-N2-C2
4	A	1384	NAG	C4-C5-C6-O6

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	1384	NAG	1	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 [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	362/362 (100%)	-0.56	0 100 100	27, 49, 75, 102	2 (0%)
1	B	362/362 (100%)	-0.54	1 (0%) 90 91	32, 48, 73, 92	1 (0%)
1	C	362/362 (100%)	-0.13	0 100 100	26, 71, 106, 117	2 (0%)
1	D	362/362 (100%)	-0.28	0 100 100	31, 61, 91, 108	1 (0%)
All	All	1448/1448 (100%)	-0.38	1 (0%) 92 93	26, 54, 95, 117	6 (0%)

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	145	ARG	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	G	3	14/15	0.64	0.10	98,100,102,104	0
2	NAG	K	1	15/15	0.67	0.12	93,95,96,99	0
2	NAG	I	3	14/15	0.71	0.10	129,130,133,133	0
2	NAG	I	1	15/15	0.73	0.13	115,117,119,120	0
2	NAG	G	1	15/15	0.74	0.12	89,92,93,93	0
2	NAG	E	3	14/15	0.75	0.09	109,111,111,112	0

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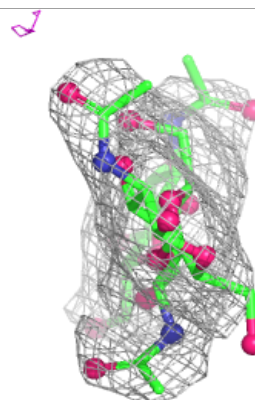
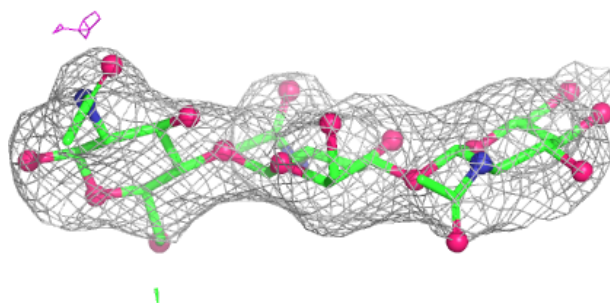
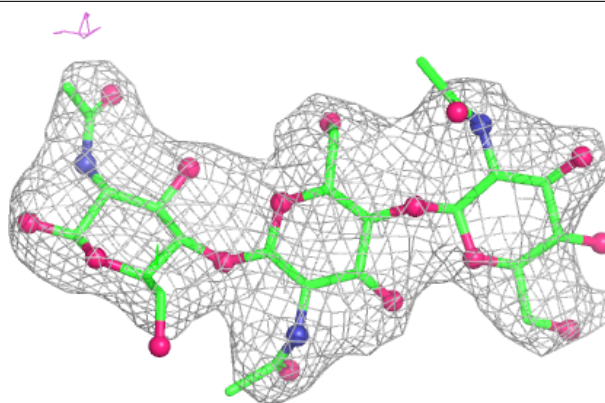
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	NAG	I	2	14/15	0.75	0.11	118,122,123,127	0
2	NAG	K	3	14/15	0.75	0.10	98,99,101,101	0
3	NAG	J	2	14/15	0.78	0.11	94,99,99,99	0
2	NAG	G	2	14/15	0.79	0.10	90,94,95,98	0
2	NAG	E	1	15/15	0.84	0.09	93,98,100,100	0
2	NAG	K	2	14/15	0.84	0.08	93,95,96,97	0
3	NAG	J	4	14/15	0.87	0.11	93,100,102,103	0
3	NAG	F	2	14/15	0.89	0.08	66,68,70,71	0
3	NAG	H	2	14/15	0.89	0.09	65,68,70,72	0
3	NAG	L	2	14/15	0.89	0.09	69,71,75,77	0
3	NAG	J	1	15/15	0.90	0.12	98,100,101,101	0
2	NAG	E	2	14/15	0.90	0.08	97,101,103,106	0
3	NAG	J	3	14/15	0.90	0.11	98,99,100,100	0
3	NAG	H	1	15/15	0.90	0.07	64,66,68,69	0
3	NAG	F	1	15/15	0.90	0.08	63,66,69,69	0
3	NAG	L	4	14/15	0.90	0.09	58,66,67,68	0
3	NAG	F	4	14/15	0.91	0.09	60,65,66,68	0
3	NAG	L	1	15/15	0.91	0.08	69,74,78,80	0
3	NAG	F	3	14/15	0.92	0.10	57,62,64,64	0
3	NAG	L	3	14/15	0.93	0.09	61,67,68,69	0
3	NAG	H	3	14/15	0.93	0.08	64,67,68,68	0
3	NAG	H	4	14/15	0.95	0.07	55,65,68,69	0

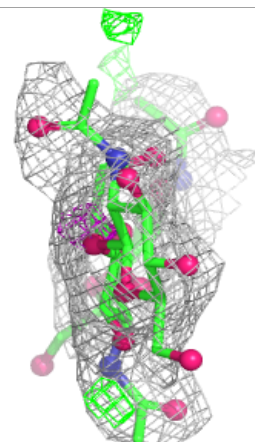
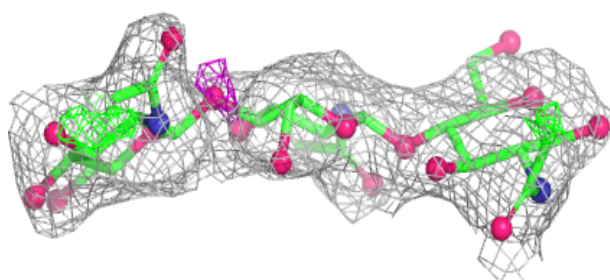
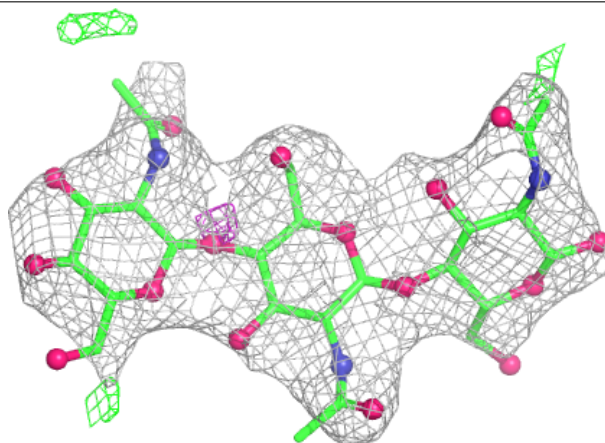
The following is a graphical depiction of the model fit to experimental electron density for oligosaccharide. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around Chain E:

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

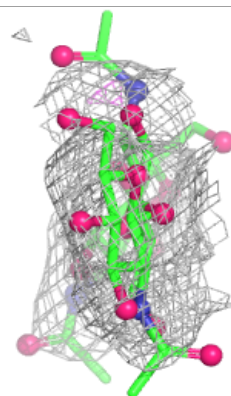
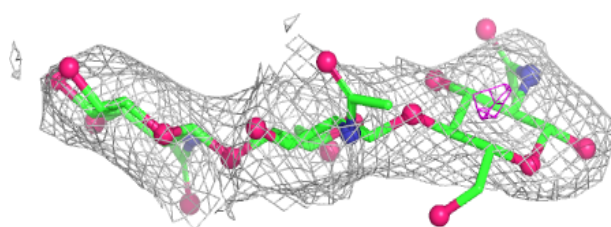
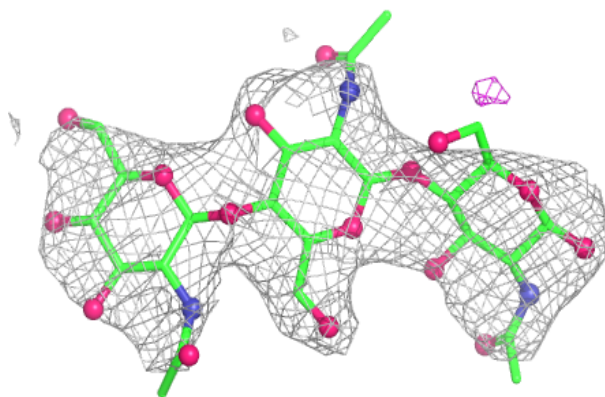
**Electron density around Chain G:**

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

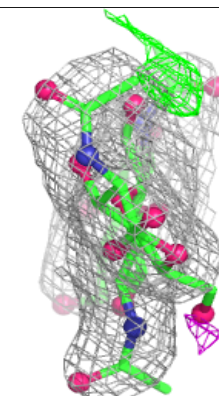
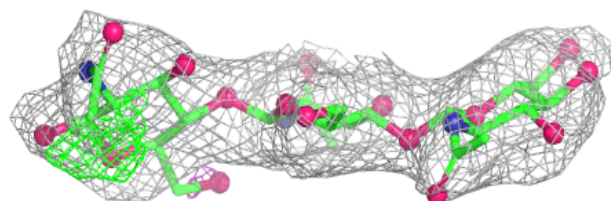
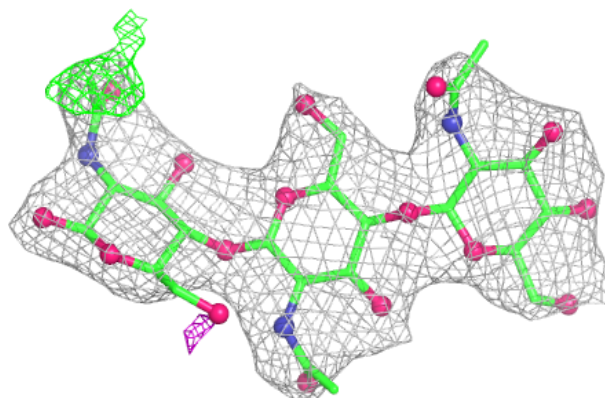


Electron density around Chain I:

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

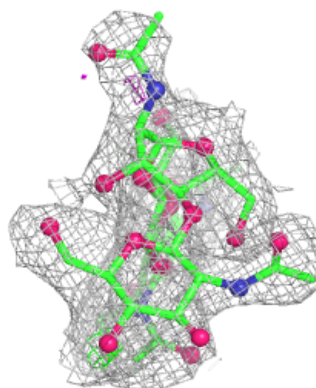
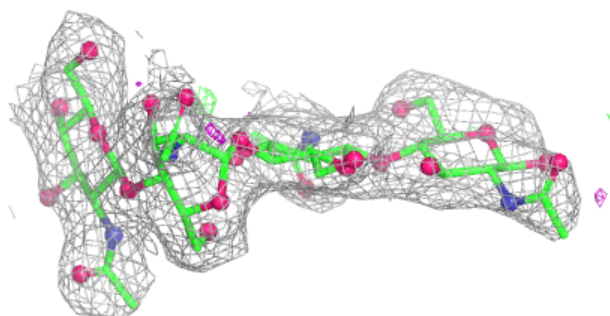
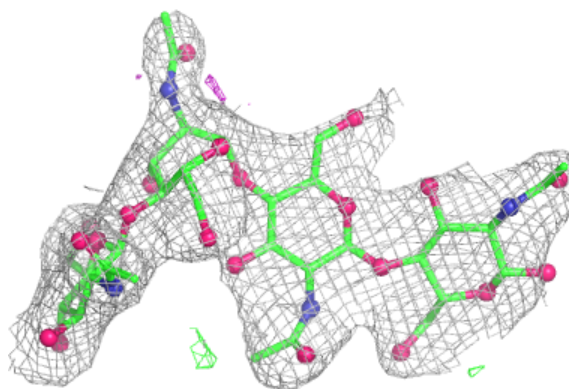
**Electron density around Chain K:**

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

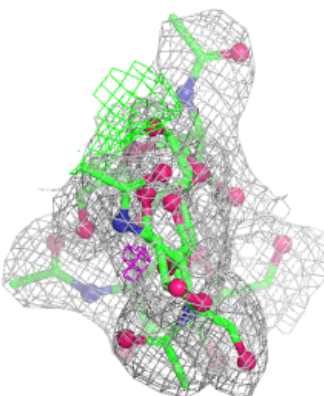
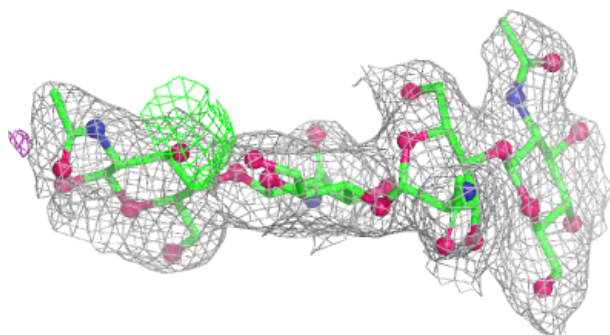
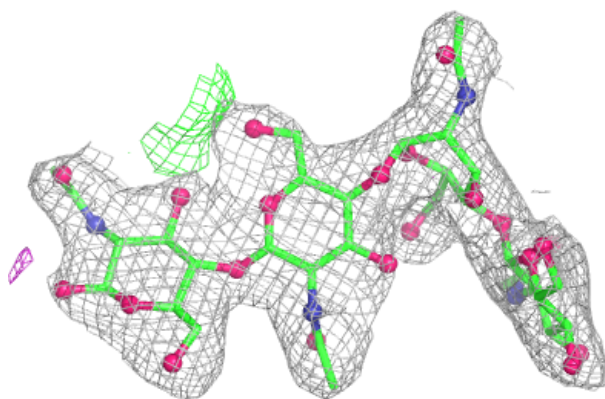


Electron density around Chain F:

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and green (positive)

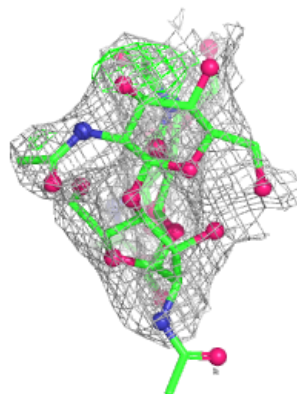
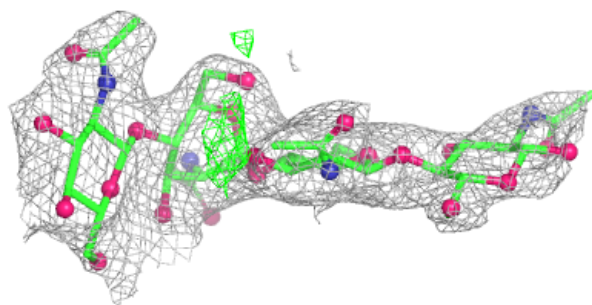
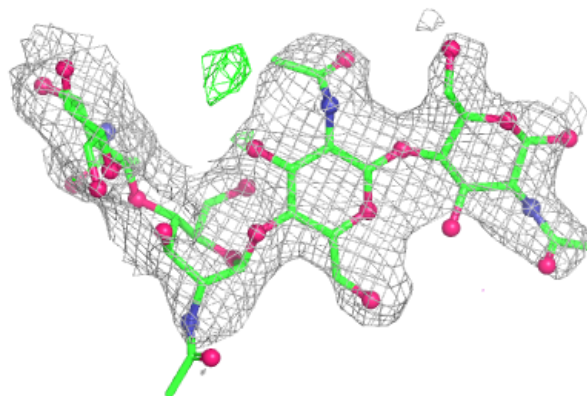
**Electron density around Chain H:**

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

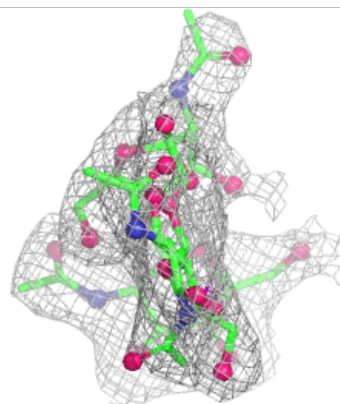
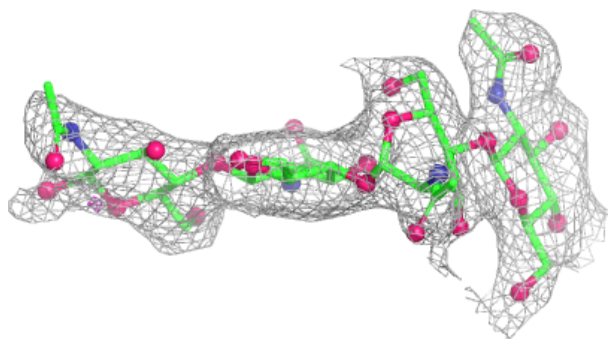
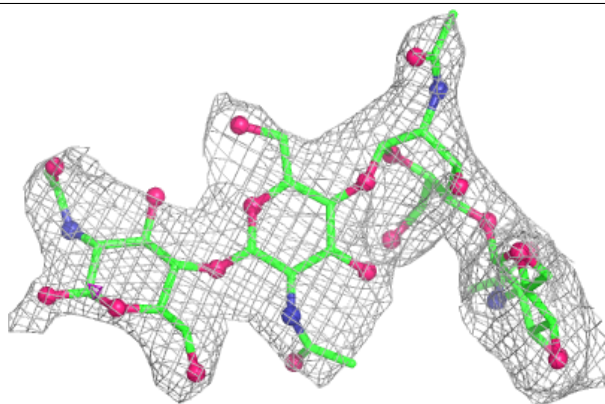


Electron density around Chain J:

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

**Electron density around Chain L:**

$2mF_o-DF_c$ (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($\text{\AA}^2$)	Q<0.9
4	NAG	C	1384	14/15	0.71	0.10	102,104,105,107	0
5	SO4	C	1387	5/5	0.71	0.07	139,140,140,140	0
5	SO4	C	1385	5/5	0.81	0.12	137,137,138,138	0
5	SO4	B	1385	5/5	0.81	0.10	138,138,138,138	0
4	NAG	A	1384	14/15	0.82	0.09	62,64,67,67	0
5	SO4	A	1386	5/5	0.83	0.08	103,104,105,105	0
5	SO4	C	1386	5/5	0.83	0.10	142,142,142,142	0
5	SO4	B	1386	5/5	0.83	0.08	115,115,116,116	0
4	NAG	B	1384	14/15	0.84	0.10	64,66,69,71	0
5	SO4	D	1386	5/5	0.84	0.13	117,118,119,119	0
5	SO4	D	1387	5/5	0.84	0.08	102,102,103,103	0
5	SO4	A	1385	5/5	0.86	0.13	102,102,103,103	0
5	SO4	A	1387	5/5	0.87	0.13	101,101,102,102	0
4	NAG	D	1384	14/15	0.88	0.08	86,87,88,89	0
5	SO4	B	1387	5/5	0.93	0.08	87,87,88,88	0
5	SO4	D	1385	5/5	0.95	0.08	86,87,88,89	0

6.5 Other polymers

There are no such residues in this entry.