



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 08:13 AM UTC

PDB ID : 2WW1 / pdb_00002ww1
Title : Structure of the Family GH92 Inverting Mannosidase BT3990 from Bacteroides
thetaiotaomicron VPI-5482 in complex with Thiomannobioside
Authors : Suits, M.D.L.; Zhu, Y.; Thompson, A.; Gilbert, H.J.; Davies, G.J.
Deposited on : 2009-10-21
Resolution : 2.25 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

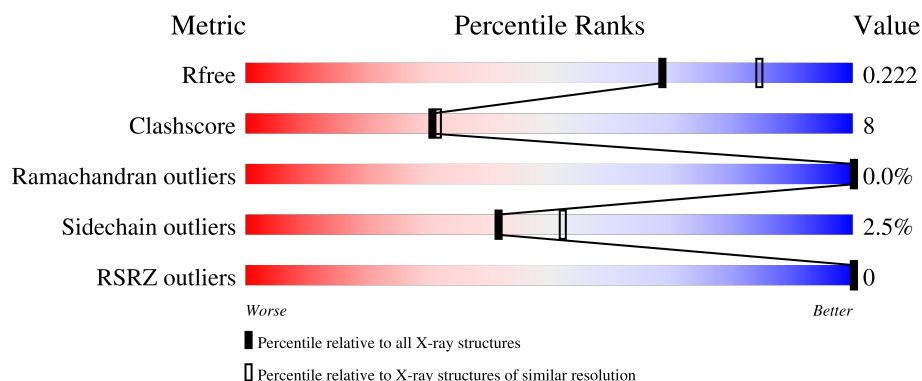
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	744	
1	B	744	
1	C	744	
1	D	744	
2	E	2	

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Mol	Chain	Length	Quality of chain
2	F	2	 100%
2	G	2	 50%  50%
2	H	2	 100%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 25408 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 PUTATIVE ALPHA-1,2-MANNOSIDASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	737	Total	C	N	O	S	0	6	0
			6002	3851	986	1130	35			
1	B	737	Total	C	N	O	S	0	4	0
			5988	3844	984	1125	35			
1	C	737	Total	C	N	O	S	0	2	0
			5972	3831	984	1122	35			
1	D	737	Total	C	N	O	S	0	4	0
			5990	3842	985	1129	34			

- Molecule 2 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-methyl 2-thio-alpha-D-mannopyranoside.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	E	2	Total	C	O	S	0	0	0
			24	13	10	1			
2	F	2	Total	C	O	S	0	0	0
			24	13	10	1			
2	G	2	Total	C	O	S	0	0	0
			24	13	10	1			
2	H	2	Total	C	O	S	0	0	0
			24	13	10	1			

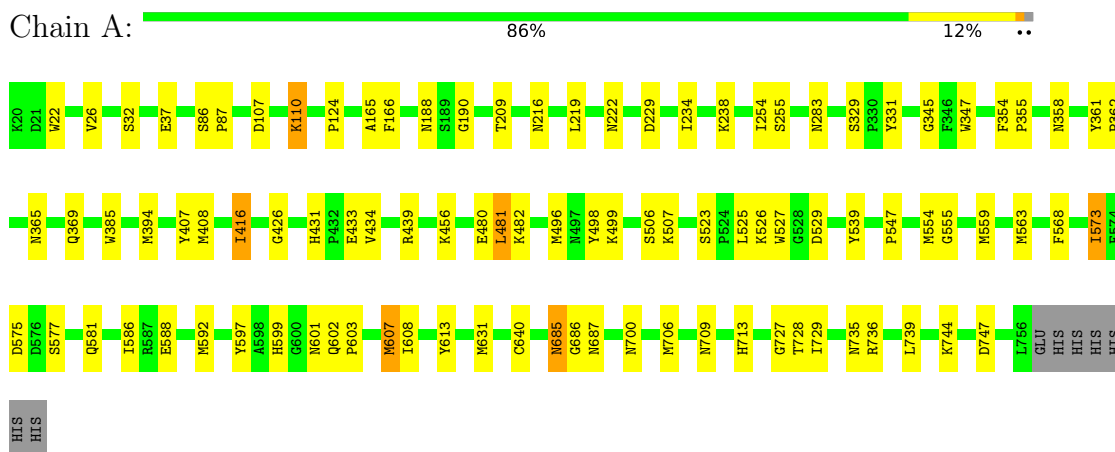
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	346	Total	O	0	0
			346	346		
3	B	364	Total	O	0	0
			364	364		
3	C	342	Total	O	0	0
			342	342		
3	D	308	Total	O	0	0
			308	308		

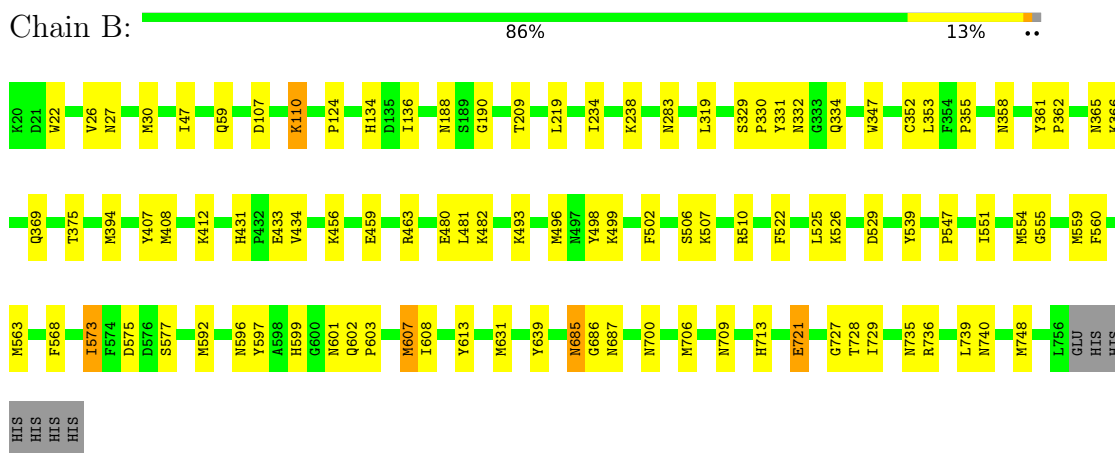
3 Residue-property plots

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.

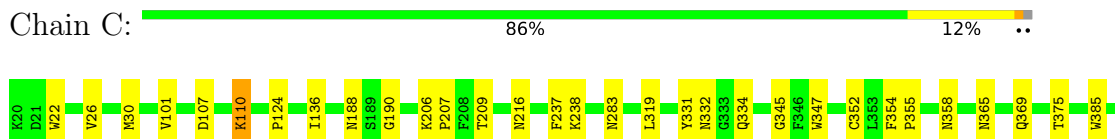
• Molecule 1: PUTATIVE ALPHA-1,2-MANNOSIDASE



• Molecule 1: PUTATIVE ALPHA-1,2-MANNOSIDASE



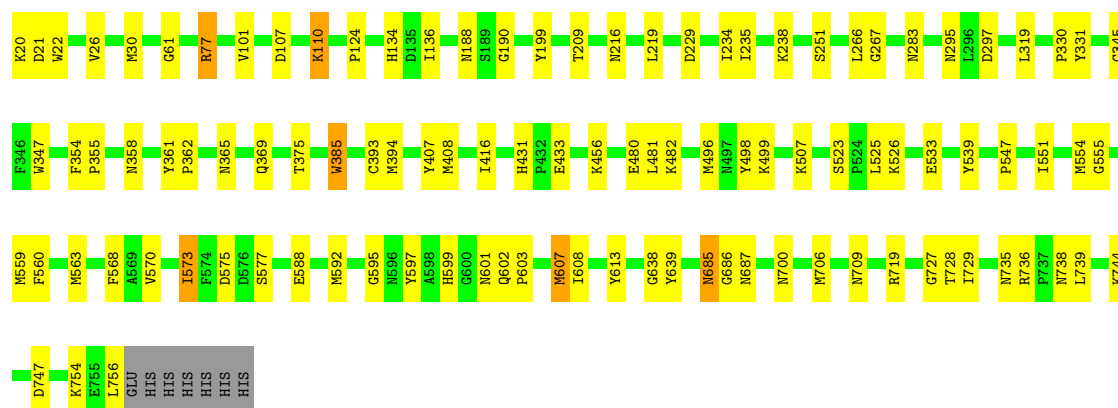
• Molecule 1: PUTATIVE ALPHA-1,2-MANNOSIDASE





- Molecule 1: PUTATIVE ALPHA-1,2-MANNOSIDASE

Chain D: 85% 13% ..



- Molecule 2: alpha-D-mannopyranose-(1-2)-methyl 2-thio-alpha-D-mannopyranoside

Chain E: 100%



- Molecule 2: alpha-D-mannopyranose-(1-2)-methyl 2-thio-alpha-D-mannopyranoside

Chain F: 100%



- Molecule 2: alpha-D-mannopyranose-(1-2)-methyl 2-thio-alpha-D-mannopyranoside

Chain G: 50% 50%



- Molecule 2: alpha-D-mannopyranose-(1-2)-methyl 2-thio-alpha-D-mannopyranoside

Chain H: 100%

Z5L1
MAN2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	66.24Å 153.11Å 185.21Å 90.00° 91.37° 90.00°	Depositor
Resolution (Å)	185.16 – 2.25 185.16 – 2.25	Depositor EDS
% Data completeness (in resolution range)	100.0 (185.16-2.25) 99.9 (185.16-2.25)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.66 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.4.0077	Depositor
R, R_{free}	0.188 , 0.218 0.191 , 0.222	Depositor DCC
R_{free} test set	8730 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	21.4	Xtriage
Anisotropy	0.013	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 8.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	0.156 for h,-k,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	25408	wwPDB-VP
Average B, all atoms (Å ²)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.46% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: Z5L, MAN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.64	0/6204	0.80	2/8416 (0.0%)
1	B	0.66	0/6181	0.81	2/8383 (0.0%)
1	C	0.66	0/6159	0.81	1/8355 (0.0%)
1	D	0.64	0/6183	0.81	4/8388 (0.0%)
All	All	0.65	0/24727	0.81	9/33542 (0.0%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	638	GLY	N-CA-C	6.55	122.14	113.37
1	B	721	GLU	CA-CB-CG	5.89	125.88	114.10
1	A	523	SER	CA-C-N	5.58	125.42	119.28
1	A	523	SER	C-N-CA	5.58	125.42	119.28
1	D	523	SER	CA-C-N	5.46	125.12	119.56
1	D	523	SER	C-N-CA	5.46	125.12	119.56
1	C	721	GLU	CA-CB-CG	5.39	124.89	114.10
1	B	330	PRO	N-CA-C	-5.10	109.38	114.68
1	D	330	PRO	N-CA-C	-5.09	109.39	114.68

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	6002	0	5681	95	1
1	B	5988	0	5672	102	1
1	C	5972	0	5644	103	0
1	D	5990	0	5664	95	0
2	E	24	0	10	0	0
2	F	24	0	10	0	0
2	G	24	0	10	1	0
2	H	24	0	10	0	0
3	A	346	0	0	6	0
3	B	364	0	0	8	0
3	C	342	0	0	8	0
3	D	308	0	0	4	0
All	All	25408	0	22701	386	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 8.

All (386) 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:332:ASN:HB2	1:C:334:GLN:NE2	1.14	1.46
1:B:332:ASN:CB	1:C:334:GLN:HE22	1.30	1.44
1:B:332:ASN:CB	1:C:334:GLN:NE2	1.92	1.27
3:C:2095:HOH:O	1:D:416:ILE:HB	1.34	1.25
1:C:581:GLN:HA	3:C:2265:HOH:O	1.50	1.07
1:D:496:MET:CE	1:D:499:LYS:HE3	1.86	1.06
1:D:555:GLY:HA3	1:D:559[B]:MET:CE	1.89	1.02
1:A:555:GLY:HA3	1:A:559[B]:MET:CE	1.90	1.01
1:C:496:MET:CE	1:C:499:LYS:HE3	1.95	0.97
1:B:496:MET:CE	1:B:499:LYS:HE3	1.95	0.97
1:B:555:GLY:O	1:B:559[B]:MET:HE2	1.64	0.95
1:C:555:GLY:O	1:C:559[B]:MET:HE2	1.65	0.95
1:D:407:TYR:HD1	1:D:408:MET:CE	1.80	0.94
1:C:496:MET:HE2	1:C:499:LYS:NZ	1.83	0.94
1:B:407:TYR:HD1	1:B:408:MET:HE2	1.30	0.94
1:A:555:GLY:HA3	1:A:559[B]:MET:HE1	1.50	0.93
1:A:496:MET:CE	1:A:499:LYS:HE3	1.99	0.92
1:C:407:TYR:HD1	1:C:408:MET:CE	1.82	0.92
1:D:555:GLY:HA3	1:D:559[B]:MET:HE1	1.49	0.92
1:A:407:TYR:HD1	1:A:408:MET:HE2	1.34	0.92
1:C:555:GLY:HA3	1:C:559[B]:MET:CE	1.98	0.92
1:D:555:GLY:O	1:D:559[B]:MET:HE2	1.71	0.90

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:736:ARG:HG2	1:D:736:ARG:HH11	1.36	0.90
1:B:555:GLY:HA3	1:B:559[B]:MET:CE	2.01	0.89
1:B:407:TYR:HD1	1:B:408:MET:CE	1.85	0.89
1:A:188:ASN:HD22	1:A:190:GLY:H	1.16	0.89
1:A:407:TYR:HD1	1:A:408:MET:CE	1.84	0.89
1:B:573:ILE:HD11	3:B:2275:HOH:O	1.73	0.89
1:A:416[A]:ILE:HD11	1:A:481:LEU:HD21	1.55	0.88
1:C:407:TYR:HD1	1:C:408:MET:HE2	1.36	0.87
1:C:555:GLY:HA3	1:C:559[B]:MET:HE1	1.54	0.87
1:C:188:ASN:HD22	1:C:190:GLY:H	1.22	0.86
1:B:555:GLY:HA3	1:B:559[B]:MET:HE1	1.55	0.86
1:B:358:ASN:HD21	1:B:365:ASN:HD22	1.23	0.85
1:D:407:TYR:HD1	1:D:408:MET:HE2	1.40	0.84
1:B:496:MET:HE2	1:B:499:LYS:NZ	1.93	0.84
1:A:581:GLN:HG2	3:A:2255:HOH:O	1.78	0.83
1:C:358:ASN:HD21	1:C:365:ASN:HD22	1.23	0.82
1:C:407:TYR:CD1	1:C:408:MET:CE	2.63	0.81
1:A:358:ASN:HD21	1:A:365:ASN:HD22	1.29	0.81
1:D:188:ASN:HD22	1:D:190:GLY:H	1.28	0.80
1:D:407:TYR:CD1	1:D:408:MET:CE	2.64	0.80
1:A:736:ARG:HH11	1:A:736:ARG:HG2	1.46	0.80
1:A:555:GLY:O	1:A:559[B]:MET:HE2	1.80	0.80
1:B:188:ASN:HD22	1:B:190:GLY:H	1.27	0.80
1:B:573:ILE:CD1	3:B:2275:HOH:O	2.30	0.79
1:B:407:TYR:CD1	1:B:408:MET:HE2	2.18	0.78
1:A:713:HIS:HD2	3:A:2322:HOH:O	1.66	0.78
1:D:496:MET:HE2	1:D:499:LYS:NZ	1.99	0.78
1:B:407:TYR:CD1	1:B:408:MET:CE	2.67	0.77
1:D:496:MET:CE	1:D:499:LYS:CE	2.61	0.77
1:D:555:GLY:CA	1:D:559[B]:MET:CE	2.61	0.76
1:C:496:MET:CE	1:C:499:LYS:CE	2.62	0.76
1:A:407:TYR:CD1	1:A:408:MET:CE	2.67	0.76
1:B:334:GLN:HE21	1:C:332:ASN:HD22	1.33	0.76
1:B:736:ARG:HG2	1:B:736:ARG:HH11	1.49	0.76
1:A:555:GLY:CA	1:A:559[B]:MET:CE	2.63	0.75
1:C:736:ARG:HH11	1:C:736:ARG:HG2	1.50	0.75
1:A:496:MET:HE2	1:A:499:LYS:NZ	2.02	0.74
1:B:555:GLY:O	1:B:559[B]:MET:CE	2.35	0.74
1:C:555:GLY:O	1:C:559[B]:MET:CE	2.35	0.74
1:B:496:MET:CE	1:B:499:LYS:CE	2.66	0.73
1:B:358:ASN:ND2	1:B:365:ASN:HD22	1.84	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:555:GLY:HA3	1:A:559[B]:MET:HE3	1.70	0.73
1:A:358:ASN:ND2	1:A:365:ASN:HD22	1.86	0.73
1:C:407:TYR:CD1	1:C:408:MET:HE2	2.23	0.72
1:A:407:TYR:CD1	1:A:408:MET:HE2	2.22	0.72
1:A:22:TRP:H	1:A:283:ASN:ND2	1.88	0.72
1:C:555:GLY:CA	1:C:559[B]:MET:CE	2.67	0.72
1:A:22:TRP:H	1:A:283:ASN:HD21	1.38	0.71
1:C:216:ASN:ND2	3:C:2113:HOH:O	2.23	0.71
1:D:22:TRP:H	1:D:283:ASN:ND2	1.88	0.71
1:C:358:ASN:ND2	1:C:365:ASN:HD22	1.87	0.71
1:B:555:GLY:CA	1:B:559[B]:MET:CE	2.68	0.70
1:D:407:TYR:CD1	1:D:408:MET:HE2	2.27	0.69
1:D:555:GLY:HA3	1:D:559[B]:MET:HE3	1.72	0.69
1:D:685:ASN:ND2	1:D:687:ASN:H	1.89	0.69
1:B:332:ASN:HB2	1:C:334:GLN:HE21	1.46	0.69
1:A:496:MET:CE	1:A:499:LYS:CE	2.71	0.69
1:A:588:GLU:O	1:A:592[B]:MET:HG3	1.93	0.69
1:B:334:GLN:NE2	1:C:332:ASN:HD22	1.91	0.69
1:D:496:MET:HE2	1:D:499:LYS:CE	2.23	0.68
1:D:736:ARG:HG2	1:D:736:ARG:NH1	2.09	0.68
1:D:555:GLY:C	1:D:559[B]:MET:CE	2.67	0.68
1:D:685:ASN:C	1:D:685:ASN:HD22	2.02	0.67
1:C:555:GLY:C	1:C:559[B]:MET:CE	2.67	0.67
1:B:555:GLY:C	1:B:559[B]:MET:CE	2.67	0.67
1:A:685:ASN:ND2	1:A:687:ASN:H	1.94	0.66
1:D:358:ASN:HD21	1:D:365:ASN:HD22	1.41	0.66
1:D:22:TRP:H	1:D:283:ASN:HD21	1.39	0.66
1:B:408:MET:HE1	1:B:480:GLU:HG2	1.78	0.66
1:C:685:ASN:ND2	1:C:687:ASN:H	1.94	0.66
1:C:408:MET:HE1	1:C:480:GLU:HG2	1.77	0.66
1:B:525:LEU:HD13	1:B:573:ILE:HG12	1.78	0.65
1:A:431:HIS:HE1	3:A:2164:HOH:O	1.79	0.65
1:D:358:ASN:ND2	1:D:365:ASN:HD22	1.94	0.65
1:B:559[A]:MET:HE2	1:B:563:MET:HG2	1.77	0.65
1:C:496:MET:HE2	1:C:499:LYS:CE	2.24	0.65
1:C:554:MET:O	1:C:559[B]:MET:HE3	1.97	0.65
1:D:555:GLY:O	1:D:559[B]:MET:CE	2.44	0.65
1:B:22:TRP:H	1:B:283:ASN:ND2	1.95	0.65
1:B:559[A]:MET:CE	1:B:563:MET:HG2	2.27	0.64
1:B:332:ASN:CA	1:C:334:GLN:NE2	2.60	0.64
1:B:685:ASN:ND2	1:B:687:ASN:H	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:347:TRP:CZ3	1:D:394:MET:HE2	2.32	0.64
1:D:496:MET:HE1	1:D:499:LYS:HE3	1.74	0.64
1:C:507:LYS:HE3	1:C:559[A]:MET:SD	2.38	0.64
1:D:554:MET:O	1:D:559[B]:MET:HE3	1.97	0.64
1:A:555:GLY:C	1:A:559[B]:MET:CE	2.70	0.64
1:B:554:MET:O	1:B:559[B]:MET:HE3	1.98	0.64
1:A:555:GLY:CA	1:A:559[B]:MET:HE3	2.27	0.64
1:A:496:MET:HE2	1:A:499:LYS:CE	2.28	0.63
1:D:555:GLY:CA	1:D:559[B]:MET:HE3	2.28	0.63
1:A:685:ASN:HD22	1:A:685:ASN:C	2.06	0.63
1:A:216:ASN:HD21	1:A:229[B]:ASP:HB2	1.62	0.63
1:A:736:ARG:HG2	1:A:736:ARG:NH1	2.14	0.62
1:B:592[B]:MET:HE1	1:B:599:HIS:H	1.63	0.62
1:A:592[B]:MET:HE1	1:A:599:HIS:H	1.63	0.62
1:B:496:MET:HE2	1:B:499:LYS:CE	2.27	0.62
1:C:496:MET:HE2	1:C:499:LYS:HZ2	1.65	0.62
1:C:525:LEU:HD13	1:C:573:ILE:HG12	1.81	0.62
1:D:394:MET:HE1	1:D:533:GLU:HG3	1.80	0.62
1:C:22:TRP:H	1:C:283:ASN:ND2	1.98	0.62
1:B:507:LYS:HE3	1:B:559[A]:MET:SD	2.40	0.61
1:C:555:GLY:CA	1:C:559[B]:MET:HE3	2.30	0.61
1:C:559[A]:MET:HE2	1:C:563:MET:HG2	1.81	0.61
1:C:736:ARG:HG2	1:C:736:ARG:NH1	2.16	0.61
1:A:408:MET:HE1	1:A:480:GLU:HG2	1.82	0.61
1:B:597:TYR:OH	1:B:599:HIS:HD2	1.84	0.60
1:B:22:TRP:H	1:B:283:ASN:HD21	1.47	0.60
1:B:332:ASN:HB3	1:C:334:GLN:HE22	1.54	0.60
1:B:685:ASN:C	1:B:685:ASN:HD22	2.09	0.60
1:C:431:HIS:HD2	1:C:433:GLU:H	1.49	0.60
1:D:597:TYR:OH	1:D:599:HIS:HD2	1.84	0.60
1:B:332:ASN:HB2	1:C:334:GLN:HE22	0.56	0.60
1:C:685:ASN:C	1:C:685:ASN:HD22	2.09	0.60
1:C:709:ASN:HD21	1:C:727:GLY:HA3	1.65	0.60
1:A:188:ASN:ND2	1:A:190:GLY:H	1.93	0.60
1:A:559[A]:MET:CE	1:A:563:MET:HG2	2.32	0.60
1:A:554:MET:O	1:A:559[B]:MET:HE3	2.01	0.59
1:C:496:MET:HE1	1:C:499:LYS:HE3	1.82	0.59
1:D:408:MET:HE1	1:D:480:GLU:HG2	1.83	0.59
1:D:709:ASN:HD21	1:D:727:GLY:HA3	1.67	0.59
1:A:555:GLY:O	1:A:559[B]:MET:CE	2.49	0.59
1:A:597:TYR:OH	1:A:599:HIS:HD2	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:736:ARG:HG2	1:B:736:ARG:NH1	2.15	0.59
1:A:601[B]:ASN:OD1	1:A:603:PRO:HD2	2.03	0.59
1:C:559[A]:MET:CE	1:C:563:MET:HG2	2.33	0.58
1:B:709:ASN:ND2	1:B:728:THR:H	2.00	0.58
1:C:188:ASN:ND2	1:C:190:GLY:H	1.95	0.58
1:B:709:ASN:HD21	1:B:727:GLY:HA3	1.67	0.58
1:C:347:TRP:CZ3	1:C:394:MET:HE2	2.38	0.58
1:D:431:HIS:HD2	1:D:433:GLU:H	1.51	0.58
1:D:507:LYS:HE3	1:D:559[A]:MET:SD	2.44	0.58
1:D:496:MET:HE3	1:D:499:LYS:HE3	1.81	0.58
1:C:22:TRP:H	1:C:283:ASN:HD21	1.50	0.58
1:C:555:GLY:HA3	1:C:559[B]:MET:HE3	1.80	0.58
1:D:20:LYS:N	3:D:2001:HOH:O	2.37	0.57
1:D:77:ARG:NH2	3:D:2041:HOH:O	2.37	0.57
1:B:526:LYS:HA	1:B:575:ASP:HB3	1.85	0.57
1:C:496:MET:HE2	1:C:499:LYS:HZ1	1.65	0.57
1:C:755:GLU:O	1:C:756:LEU:C	2.47	0.57
1:B:601:ASN:HB2	3:B:2288:HOH:O	2.05	0.57
1:D:525:LEU:HD13	1:D:573:ILE:HG12	1.87	0.57
1:B:496:MET:HE3	1:B:499:LYS:HE3	1.84	0.56
1:C:685:ASN:HD22	1:C:687:ASN:H	1.53	0.56
1:A:601[B]:ASN:CG	1:A:603:PRO:HD2	2.31	0.56
1:B:555:GLY:CA	1:B:559[B]:MET:HE3	2.35	0.56
1:C:431:HIS:CD2	1:C:433:GLU:H	2.24	0.56
1:D:709:ASN:ND2	1:D:728:THR:H	2.04	0.56
1:A:709:ASN:ND2	1:A:728:THR:H	2.04	0.56
1:C:547:PRO:HG2	1:C:613:TYR:CE2	2.41	0.56
1:A:700:ASN:ND2	1:A:735:ASN:HB3	2.22	0.55
1:D:685:ASN:HD22	1:D:687:ASN:H	1.55	0.55
1:B:496:MET:HE1	1:B:499:LYS:HE3	1.85	0.55
1:A:496:MET:HE1	1:A:499:LYS:HE3	1.88	0.55
1:A:431:HIS:HD2	1:A:433:GLU:H	1.55	0.55
1:A:507:LYS:HE3	1:A:559[A]:MET:SD	2.47	0.55
1:B:431:HIS:HD2	1:B:433:GLU:H	1.54	0.55
1:C:709:ASN:ND2	1:C:728:THR:H	2.04	0.55
1:A:709:ASN:HD21	1:A:727:GLY:HA3	1.72	0.55
1:D:607:MET:HE3	1:D:608:ILE:HA	1.89	0.55
1:C:592[B]:MET:HE1	1:C:599:HIS:H	1.71	0.54
2:G:1:Z5L:C5	2:G:1:Z5L:C1M	2.86	0.54
1:C:496:MET:HE3	1:C:499:LYS:HE3	1.85	0.54
1:C:597:TYR:OH	1:C:599:HIS:HD2	1.91	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:525:LEU:HD13	1:B:573:ILE:CG1	2.37	0.54
1:C:347:TRP:CE3	1:C:394:MET:HE2	2.42	0.54
1:B:555:GLY:HA3	1:B:559[B]:MET:HE3	1.86	0.54
1:D:559[A]:MET:CE	1:D:563:MET:HG2	2.37	0.53
1:D:431:HIS:CD2	1:D:433:GLU:H	2.26	0.53
1:A:607:MET:HE2	1:A:608:ILE:HG13	1.91	0.53
1:B:547:PRO:HG2	1:B:613:TYR:CE2	2.44	0.53
1:D:216:ASN:HD21	1:D:229[B]:ASP:HB2	1.72	0.53
1:D:347:TRP:CE3	1:D:394:MET:HE2	2.43	0.53
1:B:592[B]:MET:HE3	1:B:631:MET:SD	2.49	0.53
1:C:352:CYS:C	1:C:355:PRO:HD2	2.34	0.52
1:D:394:MET:HE1	1:D:533:GLU:CG	2.40	0.52
1:A:496:MET:HE1	1:A:498:TYR:OH	2.09	0.52
1:A:525:LEU:HD13	1:A:573:ILE:HG12	1.91	0.52
1:B:431:HIS:CD2	1:B:433:GLU:H	2.27	0.52
1:D:607:MET:HE2	1:D:608:ILE:HG13	1.92	0.52
1:B:496:MET:HE2	1:B:499:LYS:HZ2	1.75	0.52
1:D:700:ASN:ND2	1:D:735:ASN:HB3	2.25	0.52
1:B:607:MET:HE2	1:B:608:ILE:HG13	1.92	0.51
1:B:685:ASN:HD22	1:B:687:ASN:H	1.59	0.51
1:A:496:MET:HE2	1:A:499:LYS:HZ1	1.73	0.51
1:A:607:MET:HE3	1:A:608:ILE:HA	1.91	0.51
1:C:607:MET:HE2	1:C:608:ILE:HG13	1.92	0.51
1:C:525:LEU:HD13	1:C:573:ILE:CG1	2.40	0.51
1:A:365:ASN:O	1:A:369:GLN:HG2	2.10	0.51
1:A:431:HIS:CD2	1:A:433:GLU:H	2.28	0.51
1:D:188:ASN:ND2	1:D:190:GLY:H	2.04	0.51
1:D:559[A]:MET:HE2	1:D:563:MET:HG2	1.93	0.50
1:A:559[A]:MET:HE2	1:A:563:MET:HG2	1.93	0.50
1:B:568:PHE:CZ	1:B:607:MET:HE1	2.45	0.50
1:D:496:MET:HE1	1:D:498:TYR:OH	2.12	0.50
1:A:700:ASN:HD22	1:A:735:ASN:HB3	1.76	0.50
1:D:547:PRO:HG2	1:D:613:TYR:CE2	2.46	0.50
1:C:555:GLY:C	1:C:559[B]:MET:HE3	2.36	0.50
1:B:713:HIS:HD2	3:B:2344:HOH:O	1.93	0.50
1:B:602:GLN:N	1:B:603:PRO:CD	2.75	0.49
1:D:361:TYR:N	1:D:362:PRO:CD	2.75	0.49
1:D:525:LEU:HD13	1:D:573:ILE:CG1	2.42	0.49
1:A:685:ASN:HD22	1:A:686:GLY:N	2.10	0.49
1:C:700:ASN:ND2	1:C:735:ASN:HB3	2.27	0.49
1:B:709:ASN:HD21	1:B:728:THR:H	1.61	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:506:SER:O	1:B:507:LYS:HB2	2.12	0.49
1:C:744:LYS:HE2	1:C:747:ASP:OD2	2.13	0.49
1:C:744:LYS:HE3	3:C:2334:HOH:O	2.13	0.48
1:A:426:GLY:O	1:A:439:ARG:HG3	2.14	0.48
1:A:568:PHE:CZ	1:A:607:MET:HE1	2.49	0.48
1:C:700:ASN:HD22	1:C:735:ASN:HB3	1.78	0.48
1:C:756:LEU:HA	3:C:2341:HOH:O	2.14	0.48
1:C:568:PHE:CZ	1:C:607:MET:HE1	2.48	0.47
1:A:709:ASN:HD21	1:A:728:THR:H	1.62	0.47
1:B:601:ASN:HB2	3:B:2287:HOH:O	2.14	0.47
1:C:751:SER:HB2	1:C:754:LYS:HD2	1.96	0.47
1:A:592[A]:MET:HE3	1:A:640:CYS:HB2	1.96	0.47
1:C:630:ARG:HD3	3:C:2283:HOH:O	2.14	0.47
1:D:26:VAL:HG11	1:D:124:PRO:HG3	1.97	0.47
1:A:254:ILE:O	1:A:255:SER:HB3	2.15	0.47
1:B:26:VAL:HG11	1:B:124:PRO:HG3	1.96	0.47
1:B:188:ASN:ND2	1:B:190:GLY:H	2.06	0.47
1:B:496:MET:HE2	1:B:499:LYS:HZ1	1.75	0.47
1:B:700:ASN:ND2	1:B:735:ASN:HB3	2.29	0.47
1:D:555:GLY:C	1:D:559[B]:MET:HE2	2.31	0.47
1:B:607:MET:HE3	1:B:608:ILE:HA	1.96	0.47
1:D:685:ASN:HD22	1:D:686:GLY:N	2.13	0.47
1:A:592[B]:MET:HE3	1:A:631:MET:SD	2.54	0.47
1:C:331:TYR:CZ	1:C:375:THR:HG23	2.49	0.47
1:B:347:TRP:CZ3	1:B:394:MET:HE2	2.50	0.46
1:D:709:ASN:HD21	1:D:728:THR:H	1.64	0.46
1:A:32:SER:HB3	3:A:2022:HOH:O	2.15	0.46
1:A:559[A]:MET:HE3	1:A:563:MET:HG2	1.96	0.46
1:A:347:TRP:CE3	1:A:394:MET:HE2	2.51	0.46
1:B:529:ASP:HB3	3:B:2265:HOH:O	2.16	0.46
1:D:588:GLU:O	1:D:592:MET:HG2	2.15	0.46
1:B:568:PHE:HZ	1:B:607:MET:HE1	1.81	0.46
1:D:719:ARG:HD2	3:D:2291:HOH:O	2.16	0.46
1:D:754:LYS:O	1:D:756:LEU:HG	2.16	0.46
1:A:107:ASP:HB3	1:A:110:LYS:HB2	1.98	0.46
1:C:573:ILE:HD13	1:C:573:ILE:H	1.81	0.46
1:A:358:ASN:HD21	1:A:365:ASN:ND2	2.07	0.45
1:A:685:ASN:ND2	1:A:685:ASN:C	2.73	0.45
1:D:107:ASP:HB3	1:D:110:LYS:HB2	1.98	0.45
1:D:607:MET:CE	1:D:608:ILE:HG13	2.46	0.45
1:A:219:LEU:HD11	1:A:234:ILE:HD12	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:59:GLN:NE2	3:B:2028:HOH:O	2.49	0.45
1:A:744:LYS:HE2	1:A:747:ASP:OD2	2.16	0.45
1:C:602:GLN:N	1:C:603:PRO:CD	2.80	0.45
1:D:134:HIS:CB	1:D:136:ILE:HD12	2.47	0.45
1:A:573:ILE:HD13	1:A:573:ILE:H	1.82	0.45
1:A:685:ASN:HD22	1:A:687:ASN:H	1.62	0.45
1:C:496:MET:HE1	1:C:498:TYR:OH	2.17	0.45
1:D:568:PHE:CZ	1:D:607:MET:HE1	2.51	0.45
1:C:706:MET:HE3	1:C:729:ILE:HG21	1.97	0.45
1:D:61:GLY:O	1:D:77:ARG:HD3	2.16	0.45
1:D:736:ARG:NH1	1:D:736:ARG:CG	2.74	0.45
1:A:37:GLU:HG2	3:A:2015:HOH:O	2.17	0.45
1:A:525:LEU:HD13	1:A:573:ILE:CG1	2.46	0.45
1:C:745:GLU:OE2	1:C:754:LYS:NZ	2.50	0.45
1:A:347:TRP:CZ3	1:A:394:MET:HE2	2.52	0.45
1:B:107:ASP:HB3	1:B:110:LYS:HB2	1.99	0.45
1:D:551:ILE:HG12	1:D:560:PHE:CG	2.52	0.44
1:C:416:ILE:H	1:C:416:ILE:HG12	1.57	0.44
1:B:219:LEU:HD11	1:B:234:ILE:HD12	1.99	0.44
1:A:602:GLN:N	1:A:603:PRO:CD	2.80	0.44
1:B:30:MET:HE2	1:B:639:TYR:CZ	2.53	0.44
1:B:365:ASN:O	1:B:369:GLN:HG2	2.16	0.44
1:D:744:LYS:HE2	1:D:747:ASP:OD2	2.17	0.44
1:C:528:GLY:HA2	1:C:531:PHE:O	2.18	0.44
1:D:365:ASN:O	1:D:369:GLN:HG2	2.18	0.44
1:A:555:GLY:C	1:A:559[B]:MET:HE3	2.42	0.44
1:B:431:HIS:HD2	1:B:434:VAL:H	1.66	0.44
1:B:459:GLU:O	1:B:463:ARG:HG3	2.18	0.44
1:B:700:ASN:HD22	1:B:735:ASN:HB3	1.82	0.44
1:D:354:PHE:N	1:D:355:PRO:CD	2.81	0.44
1:B:559[A]:MET:HE2	1:B:563:MET:CG	2.47	0.43
1:C:107:ASP:HB3	1:C:110:LYS:HB2	2.00	0.43
1:A:547:PRO:HG2	1:A:613:TYR:CE2	2.53	0.43
1:B:134:HIS:CB	1:B:136:ILE:HD12	2.48	0.43
1:B:347:TRP:CE3	1:B:394:MET:HE2	2.52	0.43
1:B:748:MET:HE2	1:B:748:MET:HA	2.00	0.43
1:C:206:LYS:HA	1:C:207:PRO:HD3	1.87	0.43
1:C:588:GLU:OE1	3:C:2267:HOH:O	2.21	0.43
1:A:361:TYR:N	1:A:362:PRO:CD	2.81	0.43
1:D:408:MET:HE2	1:D:408:MET:HA	2.01	0.43
1:C:496:MET:CE	1:C:499:LYS:NZ	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:26:VAL:HG11	1:C:124:PRO:HG3	1.99	0.43
1:A:607:MET:CE	1:A:608:ILE:HG13	2.49	0.43
1:C:459:GLU:O	1:C:463:ARG:HG3	2.19	0.43
1:A:222:ASN:HA	3:A:2094:HOH:O	2.19	0.43
1:C:408:MET:HE2	1:C:408:MET:HA	2.00	0.43
1:D:700:ASN:HD22	1:D:735:ASN:HB3	1.83	0.43
1:B:706:MET:HE3	1:B:729:ILE:HG21	2.00	0.43
1:C:526:LYS:HA	1:C:575:ASP:HB3	2.00	0.43
1:C:709:ASN:HD21	1:C:728:THR:H	1.67	0.43
1:D:235:ILE:C	1:D:235:ILE:HD12	2.44	0.43
1:B:366:LYS:HE2	1:B:412:LYS:O	2.19	0.43
1:B:573:ILE:HD13	1:B:573:ILE:H	1.84	0.43
1:C:354:PHE:N	1:C:355:PRO:CD	2.82	0.43
1:D:526:LYS:HA	1:D:575:ASP:HB3	2.01	0.43
1:A:507:LYS:NZ	1:A:559[A]:MET:HG2	2.34	0.42
1:B:555:GLY:C	1:B:559[B]:MET:HE3	2.41	0.42
1:B:353:LEU:C	1:B:353:LEU:HD13	2.44	0.42
1:C:345:GLY:HA3	1:C:385:TRP:CE2	2.53	0.42
1:C:568:PHE:HZ	1:C:607:MET:HE1	1.84	0.42
1:D:295:ASN:OD1	1:D:297:ASP:HB2	2.20	0.42
1:B:47:ILE:N	1:B:47:ILE:HD13	2.33	0.42
1:D:22:TRP:NE1	3:D:2003:HOH:O	2.16	0.42
1:D:602:GLN:N	1:D:603:PRO:CD	2.83	0.42
1:B:352:CYS:C	1:B:355:PRO:HD2	2.44	0.42
1:C:506:SER:O	1:C:507:LYS:HB2	2.20	0.42
1:B:510:ARG:NE	1:B:522:PHE:HB2	2.35	0.42
1:C:417:LYS:HD3	3:C:2217:HOH:O	2.19	0.42
1:D:331:TYR:CZ	1:D:375:THR:HG23	2.54	0.42
1:D:30:MET:HE2	1:D:639:TYR:CZ	2.55	0.42
1:D:219:LEU:HD11	1:D:234:ILE:HD12	2.01	0.42
1:A:431:HIS:HD2	1:A:434:VAL:H	1.66	0.42
1:D:738:ASN:OD1	1:D:738:ASN:C	2.63	0.42
1:A:165:ALA:O	1:A:166:PHE:HB2	2.20	0.42
1:A:345:GLY:HA3	1:A:385:TRP:CE2	2.55	0.42
1:B:496:MET:HE1	1:B:498:TYR:OH	2.20	0.41
1:C:685:ASN:ND2	1:C:685:ASN:C	2.77	0.41
1:A:527:TRP:CD2	1:A:586:ILE:HG12	2.54	0.41
1:A:706:MET:HE3	1:A:729:ILE:HG21	2.02	0.41
1:A:26:VAL:HG11	1:A:124:PRO:HG3	2.01	0.41
1:C:736:ARG:NH1	1:C:736:ARG:CG	2.82	0.41
1:B:27:ASN:HB2	3:B:2008:HOH:O	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:199:TYR:O	1:D:251:SER:HA	2.21	0.41
1:A:86:SER:HB2	1:A:87:PRO:HD2	2.03	0.41
1:A:329:SER:C	1:A:331:TYR:H	2.28	0.41
1:C:365:ASN:O	1:C:369:GLN:HG2	2.20	0.41
1:C:30:MET:HE2	1:C:639:TYR:CZ	2.56	0.41
1:C:685:ASN:HD22	1:C:686:GLY:N	2.18	0.41
1:D:101:VAL:HG13	1:D:136:ILE:HD11	2.02	0.41
1:D:555:GLY:C	1:D:559[B]:MET:HE3	2.43	0.41
1:B:332:ASN:HB3	1:C:334:GLN:NE2	2.15	0.41
1:B:361:TYR:N	1:B:362:PRO:CD	2.84	0.41
1:C:607:MET:HE3	1:C:608:ILE:HA	2.03	0.41
1:D:570:VAL:O	1:D:595:GLY:HA2	2.21	0.41
1:D:685:ASN:ND2	1:D:685:ASN:C	2.70	0.41
1:A:354:PHE:N	1:A:355:PRO:CD	2.84	0.41
1:A:496:MET:HE3	1:A:499:LYS:HE3	1.95	0.41
1:A:506:SER:O	1:A:507:LYS:HB2	2.21	0.41
1:A:526:LYS:HA	1:A:575:ASP:HB3	2.03	0.41
1:B:331:TYR:CZ	1:B:375:THR:HG23	2.56	0.41
1:B:502:PHE:CZ	1:B:507:LYS:HA	2.55	0.41
1:C:551:ILE:HG12	1:C:560:PHE:CG	2.56	0.41
1:C:101:VAL:HG13	1:C:136:ILE:HD11	2.02	0.40
1:B:329:SER:C	1:B:331:TYR:H	2.28	0.40
1:C:597:TYR:CE2	1:C:599:HIS:HB2	2.56	0.40
1:D:345:GLY:HA3	1:D:385:TRP:CE2	2.56	0.40
1:D:385:TRP:CD1	1:D:393:CYS:HB3	2.56	0.40
1:D:559[A]:MET:HE3	1:D:563:MET:HG2	2.04	0.40
1:D:266:LEU:O	1:D:267:GLY:C	2.63	0.40
1:D:706:MET:HE3	1:D:729:ILE:HG21	2.03	0.40
1:B:685:ASN:HD22	1:B:686:GLY:N	2.20	0.40
1:B:551:ILE:HG12	1:B:560:PHE:CG	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:229[A]:ASP:OD2	1:B:740:ASN:ND2[1_455]	1.90	0.30

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	742/744 (100%)	714 (96%)	27 (4%)	1 (0%)	48	56
1	B	739/744 (99%)	713 (96%)	26 (4%)	0	100	100
1	C	737/744 (99%)	712 (97%)	25 (3%)	0	100	100
1	D	739/744 (99%)	713 (96%)	26 (4%)	0	100	100
All	All	2957/2976 (99%)	2852 (96%)	104 (4%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	529	ASP

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	640/643 (100%)	626 (98%)	14 (2%)	45	56
1	B	637/643 (99%)	621 (98%)	16 (2%)	42	52
1	C	634/643 (99%)	616 (97%)	18 (3%)	38	48
1	D	638/643 (99%)	622 (98%)	16 (2%)	42	52
All	All	2549/2572 (99%)	2485 (98%)	64 (2%)	42	52

All (64) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	110	LYS
1	A	209	THR
1	A	238	LYS
1	A	416[A]	ILE
1	A	416[B]	ILE
1	A	456	LYS
1	A	481	LEU
1	A	482	LYS
1	A	539	TYR
1	A	573	ILE
1	A	577	SER
1	A	607	MET
1	A	685	ASN
1	A	739	LEU
1	B	110	LYS
1	B	209	THR
1	B	238	LYS
1	B	319	LEU
1	B	456	LYS
1	B	481	LEU
1	B	482	LYS
1	B	493	LYS
1	B	539	TYR
1	B	573	ILE
1	B	577	SER
1	B	596	ASN
1	B	607	MET
1	B	685	ASN
1	B	721	GLU
1	B	739	LEU
1	C	110	LYS
1	C	209	THR
1	C	237	PHE
1	C	238	LYS
1	C	319	LEU
1	C	416	ILE
1	C	456	LYS
1	C	459	GLU
1	C	481	LEU
1	C	482	LYS
1	C	493	LYS
1	C	539	TYR
1	C	573	ILE

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Mol	Chain	Res	Type
1	C	577	SER
1	C	607	MET
1	C	685	ASN
1	C	721	GLU
1	C	739	LEU
1	D	21	ASP
1	D	77	ARG
1	D	110	LYS
1	D	209	THR
1	D	238	LYS
1	D	319	LEU
1	D	385	TRP
1	D	456	LYS
1	D	481	LEU
1	D	482	LYS
1	D	539	TYR
1	D	573	ILE
1	D	577	SER
1	D	607	MET
1	D	685	ASN
1	D	739	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (93) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	84	GLN
1	A	103	GLN
1	A	188	ASN
1	A	216	ASN
1	A	220	GLN
1	A	226	GLN
1	A	244	GLN
1	A	246	ASN
1	A	264	ASN
1	A	283	ASN
1	A	334	GLN
1	A	358	ASN
1	A	369	GLN
1	A	374	ASN
1	A	425	HIS
1	A	431	HIS

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Mol	Chain	Res	Type
1	A	446	ASN
1	A	548	GLN
1	A	593	ASN
1	A	599	HIS
1	A	685	ASN
1	A	700	ASN
1	A	709	ASN
1	A	713	HIS
1	A	716	ASN
1	B	59	GLN
1	B	74	ASN
1	B	84	GLN
1	B	94	GLN
1	B	103	GLN
1	B	188	ASN
1	B	220	GLN
1	B	226	GLN
1	B	244	GLN
1	B	246	ASN
1	B	258	GLN
1	B	264	ASN
1	B	283	ASN
1	B	334	GLN
1	B	358	ASN
1	B	374	ASN
1	B	431	HIS
1	B	446	ASN
1	B	593	ASN
1	B	596	ASN
1	B	599	HIS
1	B	685	ASN
1	B	700	ASN
1	B	709	ASN
1	B	713	HIS
1	C	59	GLN
1	C	84	GLN
1	C	103	GLN
1	C	158	HIS
1	C	188	ASN
1	C	220	GLN
1	C	226	GLN
1	C	244	GLN

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Mol	Chain	Res	Type
1	C	264	ASN
1	C	283	ASN
1	C	334	GLN
1	C	358	ASN
1	C	374	ASN
1	C	398	ASN
1	C	431	HIS
1	C	593	ASN
1	C	599	HIS
1	C	685	ASN
1	C	700	ASN
1	C	709	ASN
1	D	59	GLN
1	D	81	GLN
1	D	84	GLN
1	D	103	GLN
1	D	158	HIS
1	D	188	ASN
1	D	216	ASN
1	D	220	GLN
1	D	226	GLN
1	D	246	ASN
1	D	264	ASN
1	D	283	ASN
1	D	358	ASN
1	D	374	ASN
1	D	431	HIS
1	D	446	ASN
1	D	593	ASN
1	D	599	HIS
1	D	685	ASN
1	D	700	ASN
1	D	709	ASN
1	D	716	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

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	Z5L	E	1	2	12,13,13	1.00	1 (8%)	14,18,18	3.21	7 (50%)
2	MAN	E	2	2	11,11,12	0.87	1 (9%)	15,15,17	1.48	2 (13%)
2	Z5L	F	1	2	12,13,13	0.84	0	14,18,18	1.88	6 (42%)
2	MAN	F	2	2	11,11,12	0.87	1 (9%)	15,15,17	1.29	2 (13%)
2	Z5L	G	1	2	12,13,13	1.15	1 (8%)	14,18,18	1.41	2 (14%)
2	MAN	G	2	2	11,11,12	0.75	0	15,15,17	0.69	0
2	Z5L	H	1	2	12,13,13	0.72	0	14,18,18	2.07	3 (21%)
2	MAN	H	2	2	11,11,12	0.89	0	15,15,17	1.09	1 (6%)

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	Z5L	E	1	2	-	0/4/24/24	0/1/1/1
2	MAN	E	2	2	-	0/2/19/22	0/1/1/1
2	Z5L	F	1	2	-	0/4/24/24	0/1/1/1
2	MAN	F	2	2	-	1/2/19/22	0/1/1/1
2	Z5L	G	1	2	-	2/4/24/24	0/1/1/1
2	MAN	G	2	2	-	0/2/19/22	0/1/1/1
2	Z5L	H	1	2	-	0/4/24/24	0/1/1/1
2	MAN	H	2	2	-	0/2/19/22	0/1/1/1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	G	1	Z5L	O1-C1	3.37	1.45	1.40
2	E	2	MAN	O5-C1	-2.27	1.39	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	1	Z5L	O5-C5	-2.24	1.38	1.44
2	F	2	MAN	O5-C1	2.01	1.47	1.43

All (23) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	1	Z5L	C1M-O1-C1	-6.03	104.10	113.26
2	E	1	Z5L	O1-C1-C2	-5.33	100.01	108.10
2	H	1	Z5L	O1-C1-C2	-4.83	100.76	108.10
2	E	1	Z5L	O5-C1-C2	4.68	119.42	110.62
2	E	1	Z5L	O5-C1-O1	-4.55	100.34	110.94
2	F	1	Z5L	C1-O5-C5	-3.96	105.98	113.72
2	E	1	Z5L	C6-C5-C4	-3.62	104.13	113.02
2	E	1	Z5L	C1-O5-C5	-3.50	106.89	113.72
2	E	2	MAN	C1-O5-C5	3.46	116.83	112.19
2	H	1	Z5L	C6-C5-C4	-3.37	104.73	113.02
2	H	1	Z5L	C1M-O1-C1	-3.21	108.38	113.26
2	G	1	Z5L	O5-C5-C6	-2.97	99.09	106.44
2	F	2	MAN	O5-C5-C6	2.64	112.81	107.66
2	E	2	MAN	C1-C2-C3	2.63	113.48	109.64
2	F	1	Z5L	O5-C5-C6	2.41	112.42	106.44
2	F	2	MAN	C1-O5-C5	2.37	115.36	112.19
2	H	2	MAN	C1-C2-C3	2.32	113.02	109.64
2	F	1	Z5L	O5-C5-C4	-2.23	105.68	109.70
2	F	1	Z5L	O1-C1-C2	-2.22	104.72	108.10
2	E	1	Z5L	O5-C5-C6	-2.20	100.98	106.44
2	F	1	Z5L	C3-C4-C5	-2.18	106.28	110.23
2	G	1	Z5L	O1-C1-C2	2.07	111.25	108.10
2	F	1	Z5L	O5-C1-O1	-2.04	106.19	110.94

There are no chirality outliers.

All (3) torsion outliers are listed below:

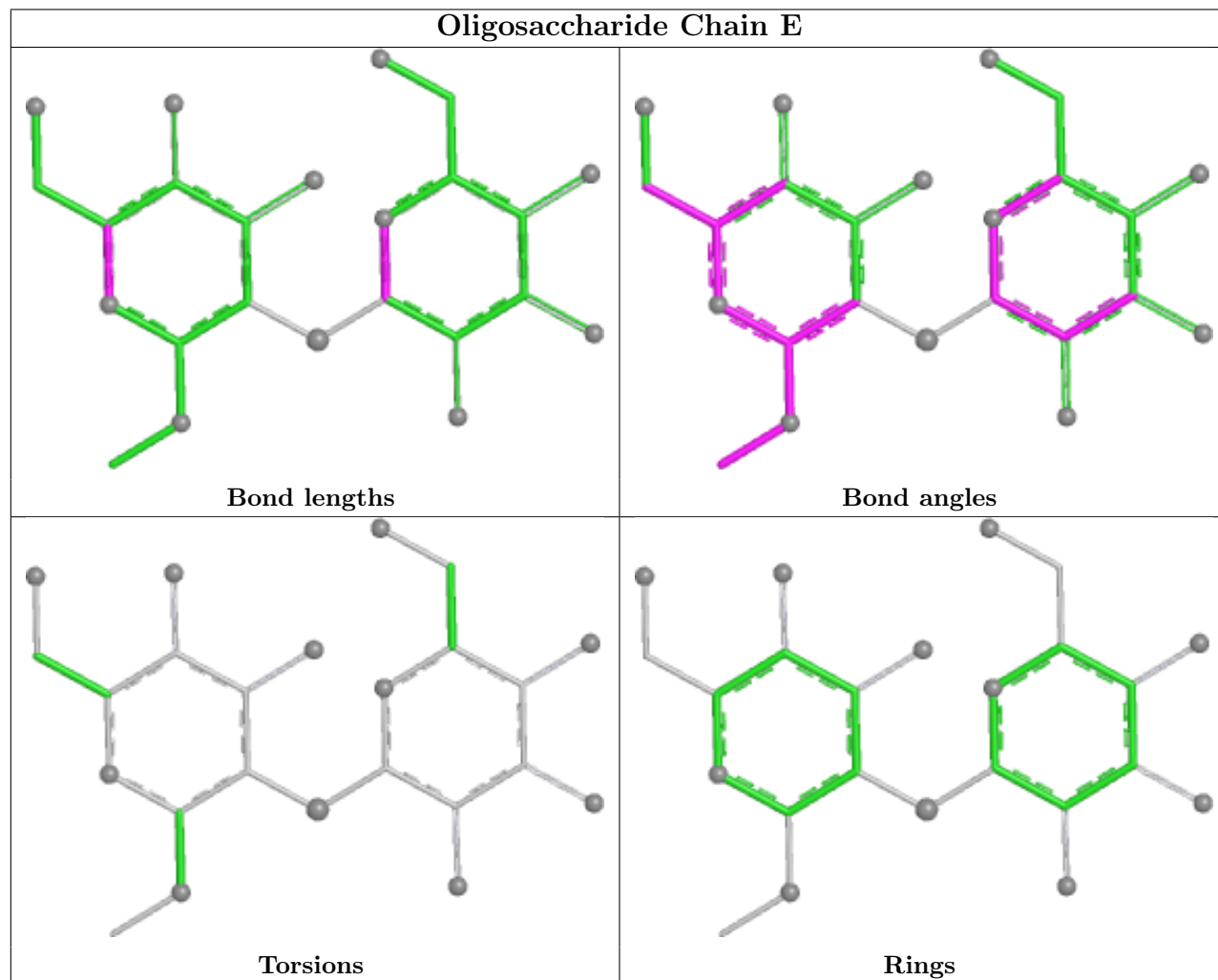
Mol	Chain	Res	Type	Atoms
2	G	1	Z5L	O5-C1-O1-C1M
2	G	1	Z5L	O5-C5-C6-O6
2	F	2	MAN	C4-C5-C6-O6

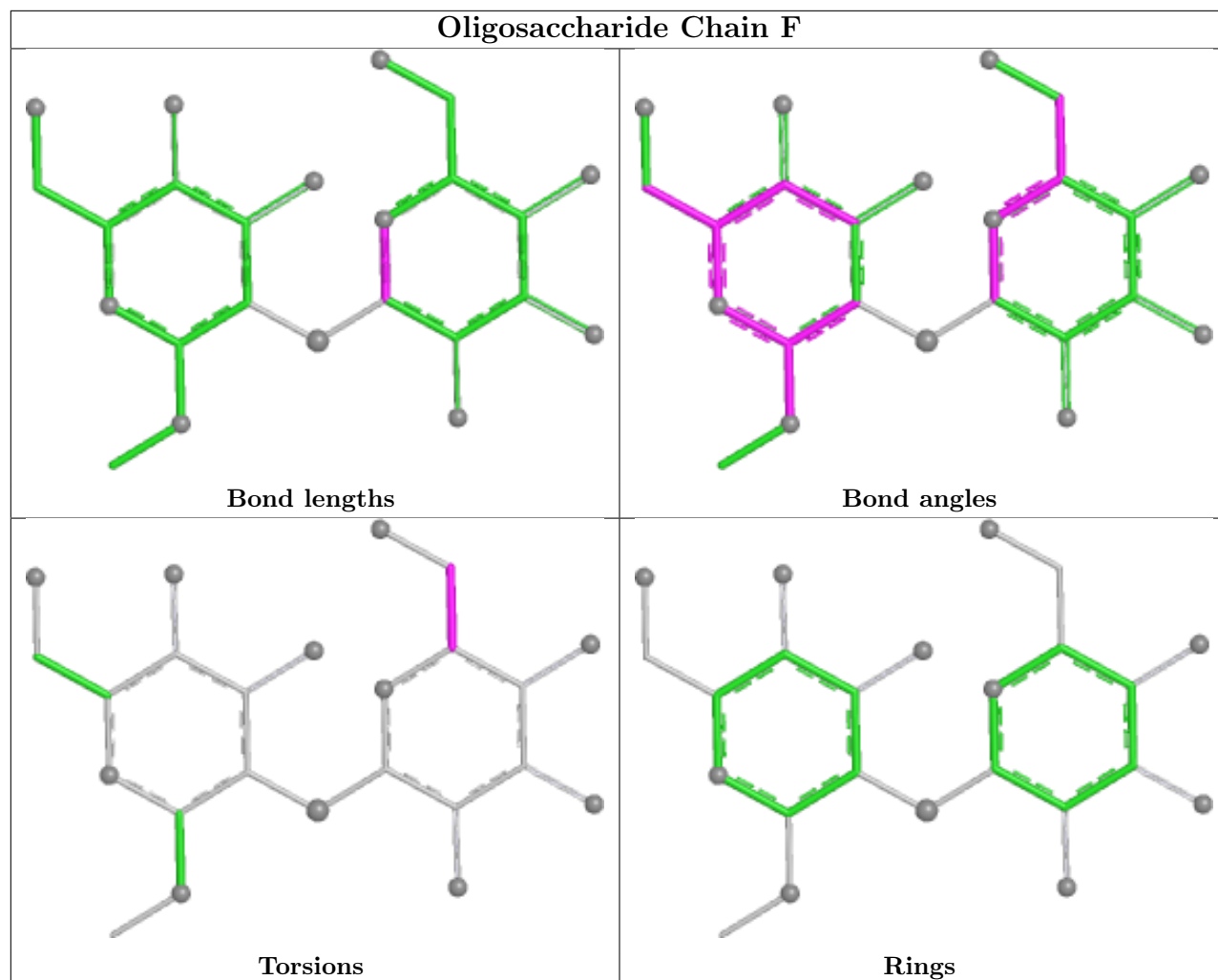
There are no ring outliers.

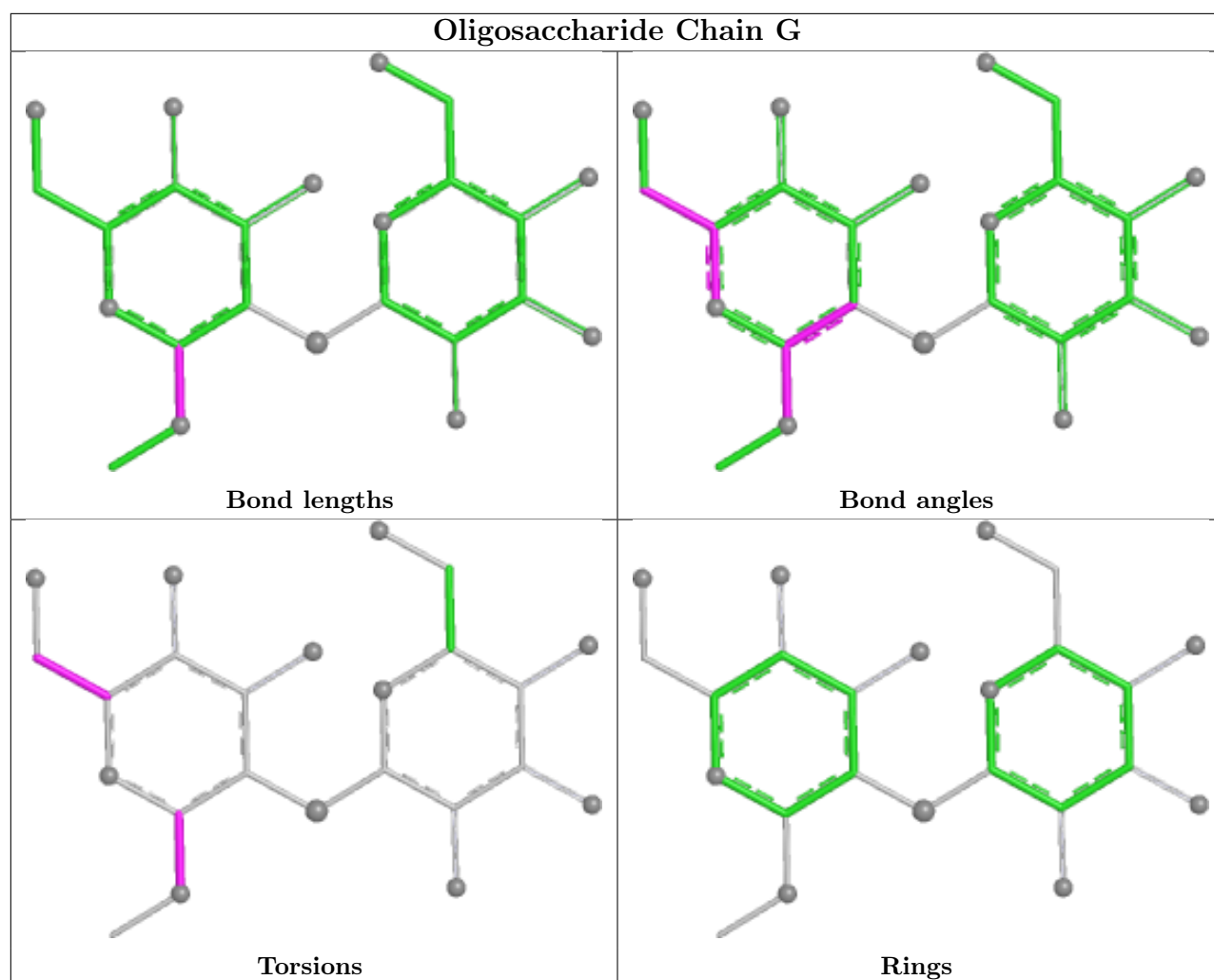
1 monomer is involved in 1 short contact:

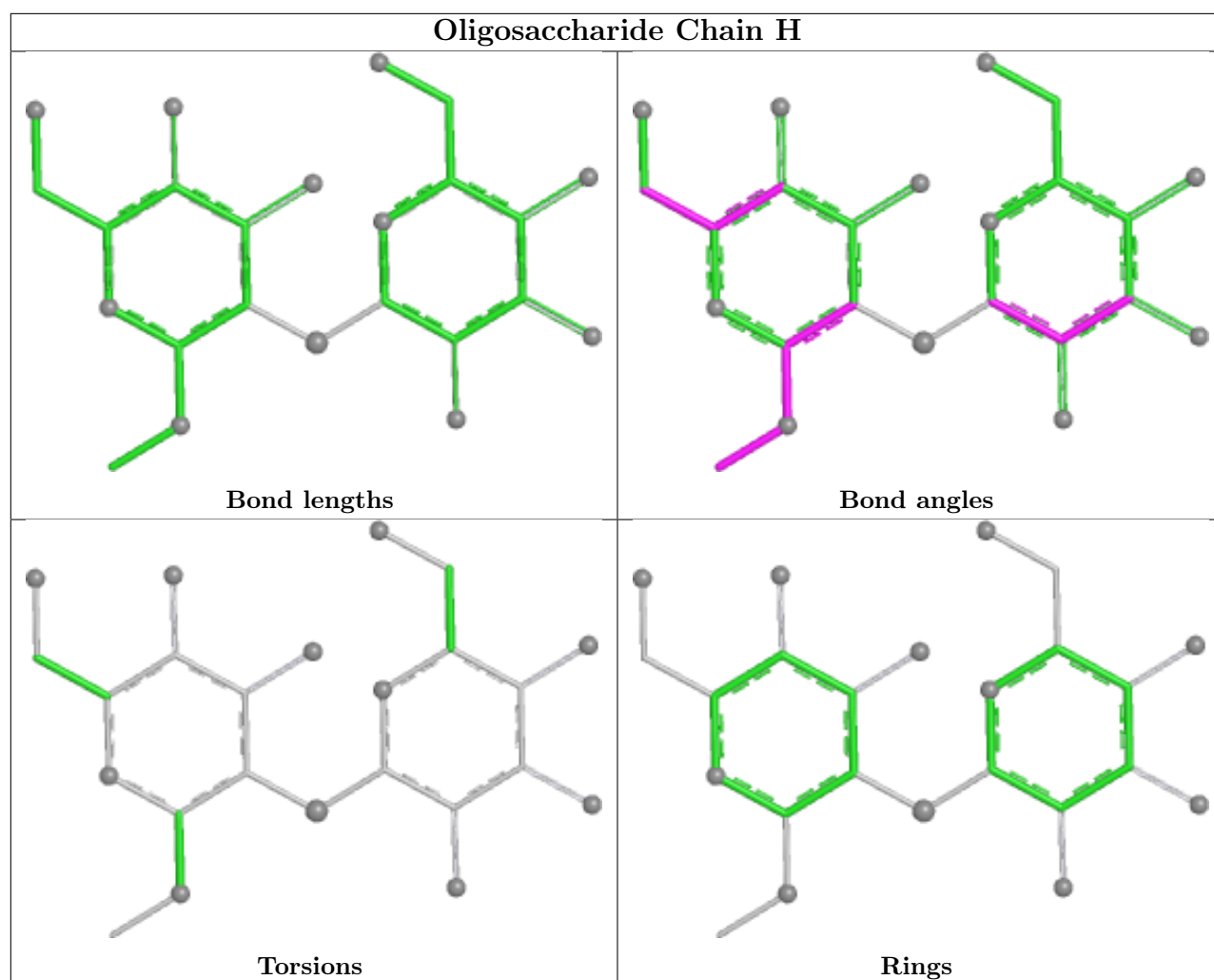
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	1	Z5L	1	0

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









5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	737/744 (99%)	-1.67	0 100 100	6, 12, 24, 33	6 (0%)
1	B	737/744 (99%)	-1.68	0 100 100	6, 12, 24, 35	4 (0%)
1	C	737/744 (99%)	-1.66	0 100 100	6, 12, 24, 35	2 (0%)
1	D	737/744 (99%)	-1.66	0 100 100	6, 12, 24, 35	4 (0%)
All	All	2948/2976 (99%)	-1.67	0 100 100	6, 12, 24, 35	16 (0%)

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

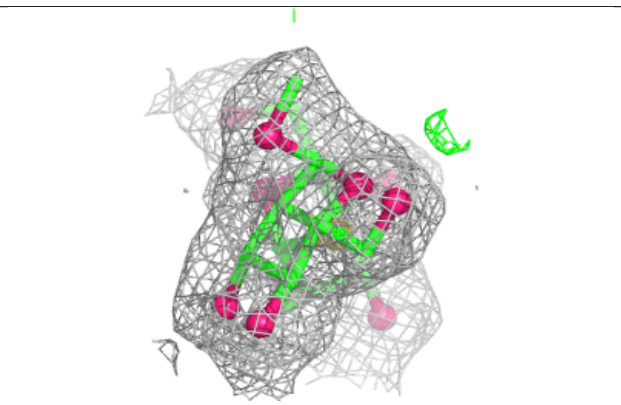
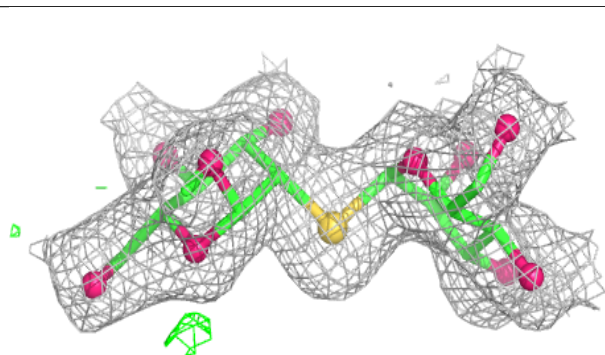
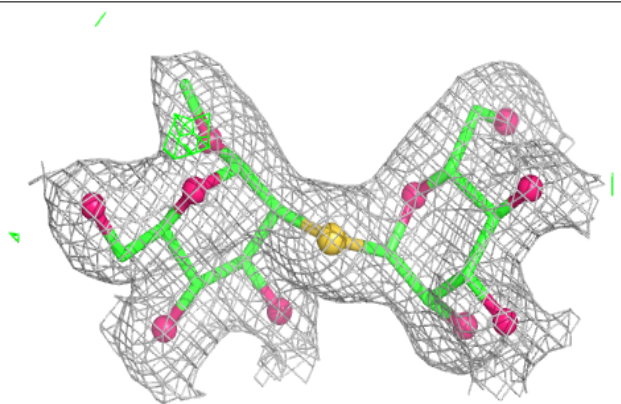
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	MAN	F	2	11/12	0.99	0.02	6,9,13,13	0
2	Z5L	G	1	13/13	0.99	0.03	16,20,23,24	0
2	MAN	G	2	11/12	0.99	0.02	9,14,16,16	0
2	Z5L	E	1	13/13	1.00	0.02	7,13,15,16	0
2	MAN	E	2	11/12	1.00	0.01	9,10,12,12	0
2	Z5L	F	1	13/13	1.00	0.03	18,23,26,27	0
2	Z5L	H	1	13/13	1.00	0.02	10,14,19,19	0
2	MAN	H	2	11/12	1.00	0.02	8,10,11,12	0

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

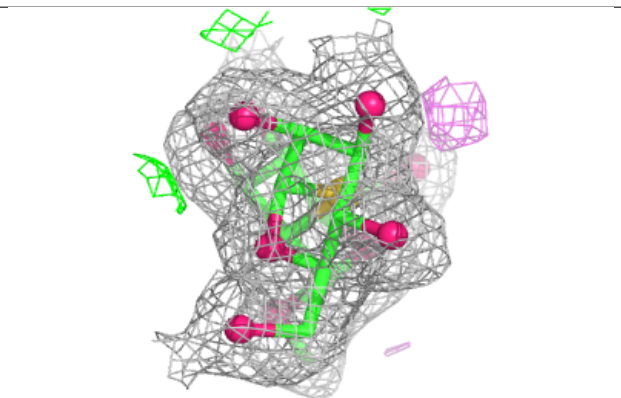
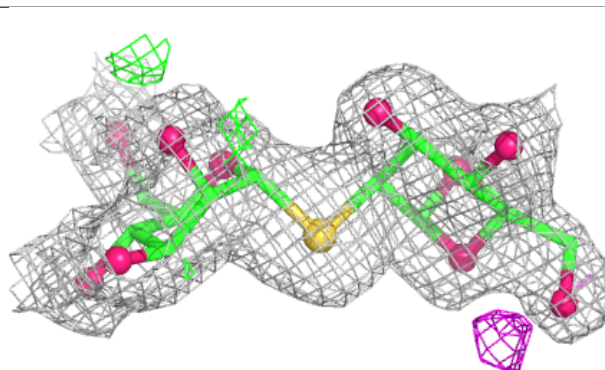
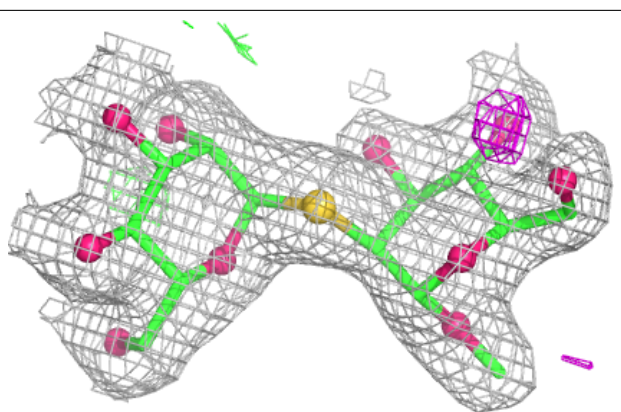
Electron density around Chain E:

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



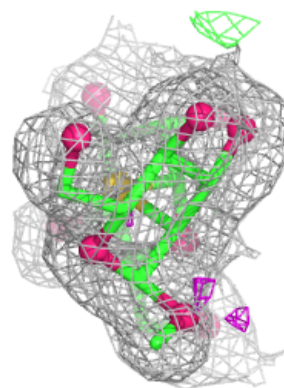
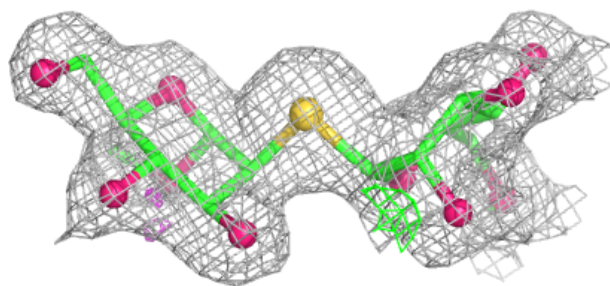
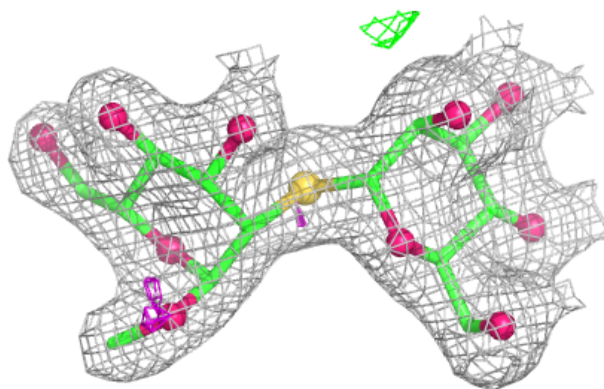
Electron density around Chain F:

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

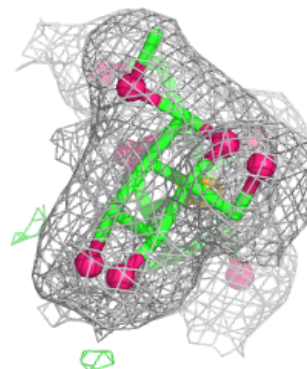
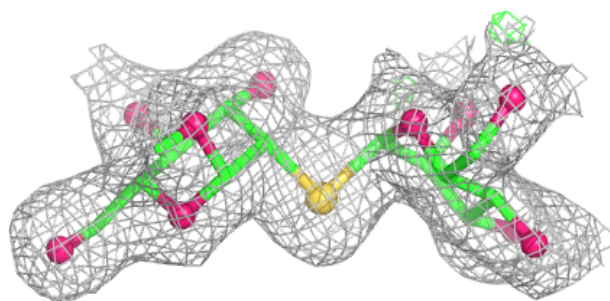
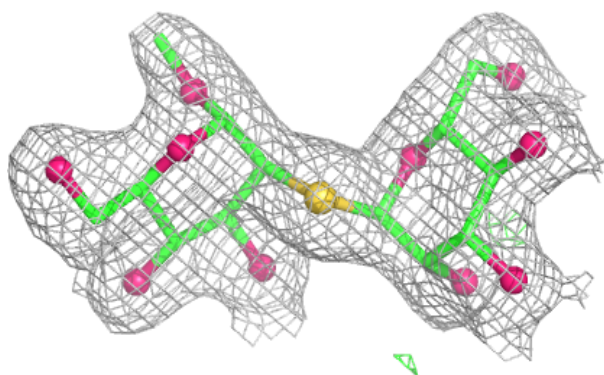


Electron density around Chain G:

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

**Electron density around Chain H:**

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



6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.