



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 17, 2026 – 10:24 PM UTC

PDB ID : 5V7J / pdb_00005v7j
Title : Crystal Structure at 3.7 Å Resolution of Glycosylated HIV-1 Clade A BG505 SOSIP.664 Prefusion Env Trimer with Four Glycans (N197, N276, N362, and N462) removed in Complex with Neutralizing Antibodies 3H+109L and 35O22.
Authors : Stewart-Jones, G.B.E.; Zhou, T.; Kwong, P.D.
Deposited on : 2017-03-20
Resolution : 2.91 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

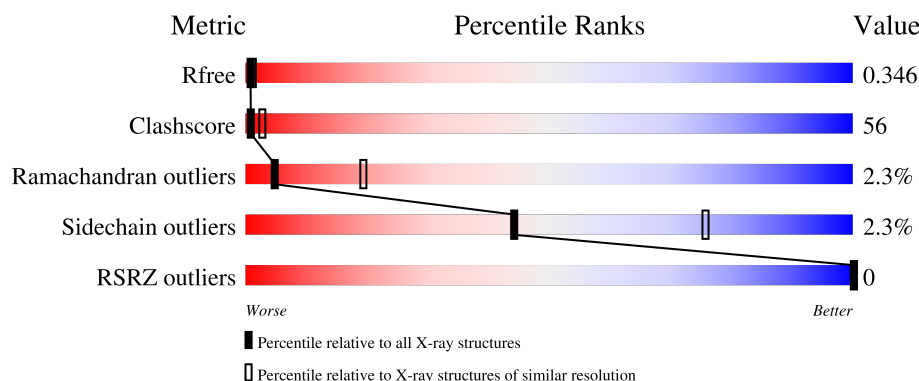
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.91 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.



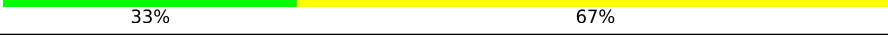
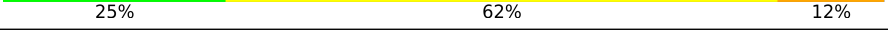
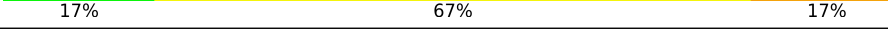
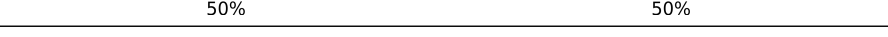
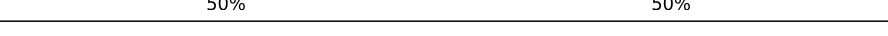
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	2481 (2.90-2.90)
Clashscore	190562	2690 (2.90-2.90)
Ramachandran outliers	187476	2623 (2.90-2.90)
Sidechain outliers	187428	2625 (2.90-2.90)
RSRZ outliers	180081	2481 (2.90-2.90)

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	G	480	35% 55% 6%
2	B	153	50% 43% ...
3	L	218	34% 59% ...
4	H	236	31% 62% 5%
5	D	240	38% 52% 8%

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Mol	Chain	Length	Quality of chain
6	E	216	
7	A	6	
8	C	7	
9	F	7	
9	K	7	
10	I	3	
10	S	3	
11	J	8	
11	M	8	
12	N	6	
13	O	10	
14	P	7	
15	Q	2	
15	R	2	

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
7	NAG	A	1	-	-	X	-

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 12436 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 Envelope glycoprotein gp160.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	452	Total	C	N	O	S	0	0	0
			3538	2222	625	664	27			

There are 11 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	199	ALA	SER	engineered mutation	UNP Q2N0S6
G	278	ALA	THR	engineered mutation	UNP Q2N0S6
G	332	ASN	THR	engineered mutation	UNP Q2N0S6
G	365	ALA	SER	engineered mutation	UNP Q2N0S6
G	464	ALA	THR	engineered mutation	UNP Q2N0S6
G	501	CYS	ALA	engineered mutation	UNP Q2N0S6
G	509	ARG	-	expression tag	UNP Q2N0S6
G	510	ARG	-	expression tag	UNP Q2N0S6
G	511	ARG	-	expression tag	UNP Q2N0S6
G	512	ARG	-	expression tag	UNP Q2N0S6
G	513	ARG	-	expression tag	UNP Q2N0S6

- Molecule 2 is a protein called Envelope glycoprotein gp160.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	148	Total	C	N	O	S	0	0	0
			1167	739	203	219	6			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	559	PRO	ILE	engineered mutation	UNP Q2N0S6
B	605	CYS	THR	engineered mutation	UNP Q2N0S6

- Molecule 3 is a protein called Antibody 3H+109L Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	211	Total	C	N	O	S	0	0	0
			1603	1007	276	315	5			

- Molecule 4 is a protein called Antibody 3H+109L Fab heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	231	Total	C	N	O	S	0	0	0
			1744	1108	283	347	6			

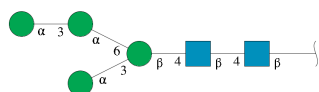
- Molecule 5 is a protein called Antibody 35O22 Fab light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	D	240	Total	C	N	O	S	0	0	0
			1813	1150	303	352	8			

- Molecule 6 is a protein called Antibody 35O22 Fab heavy chain.

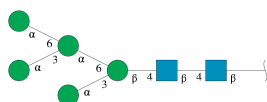
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
6	E	213	Total	C	N	O	S	0	0	0
			1615	1012	267	328	8			

- Molecule 7 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



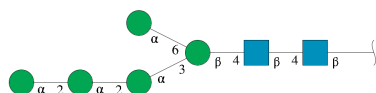
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
7	A	6	Total	C	N	O		0	0	0
			72	40	2	30				

- Molecule 8 is an oligosaccharide called alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



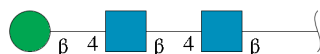
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
8	C	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 9 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



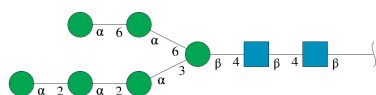
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
9	F	7	Total	C	N	O	0	0	0
			83	46	2	35			
9	K	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 10 is an oligosaccharide called beta-D-mannopyranose-(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
10	I	3	Total	C	N	O	0	0	0
			39	22	2	15			
10	S	3	Total	C	N	O	0	0	0
			39	22	2	15			

- Molecule 11 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



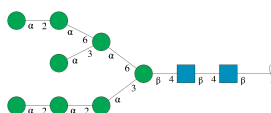
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
11	J	8	Total	C	N	O	0	0	0
			94	52	2	40			
11	M	8	Total	C	N	O	0	0	0
			94	52	2	40			

- Molecule 12 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



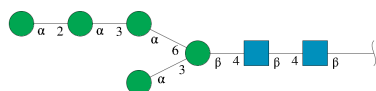
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
12	N	6	Total	C	N	O	0	0	0
			72	40	2	30			

- Molecule 13 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(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
13	O	10	Total	C	N	O	0	0	0
			116	64	2	50			

- Molecule 14 is an oligosaccharide called alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



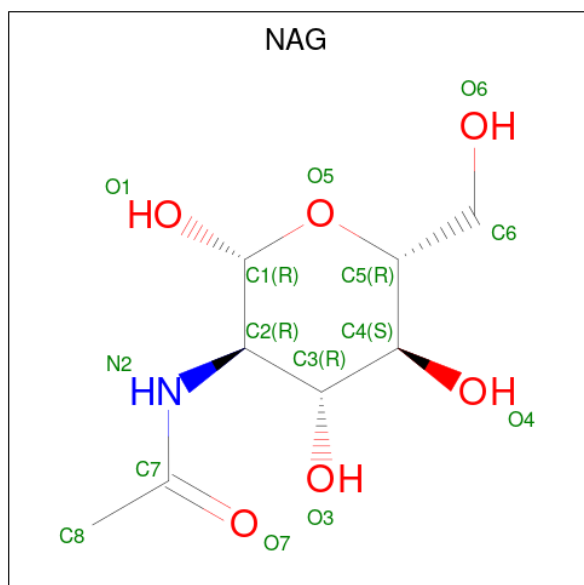
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
14	P	7	Total	C	N	O	0	0	0
			83	46	2	35			

- Molecule 15 is an oligosaccharide called 2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose.



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

- Molecule 16 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: C₈H₁₅NO₆).

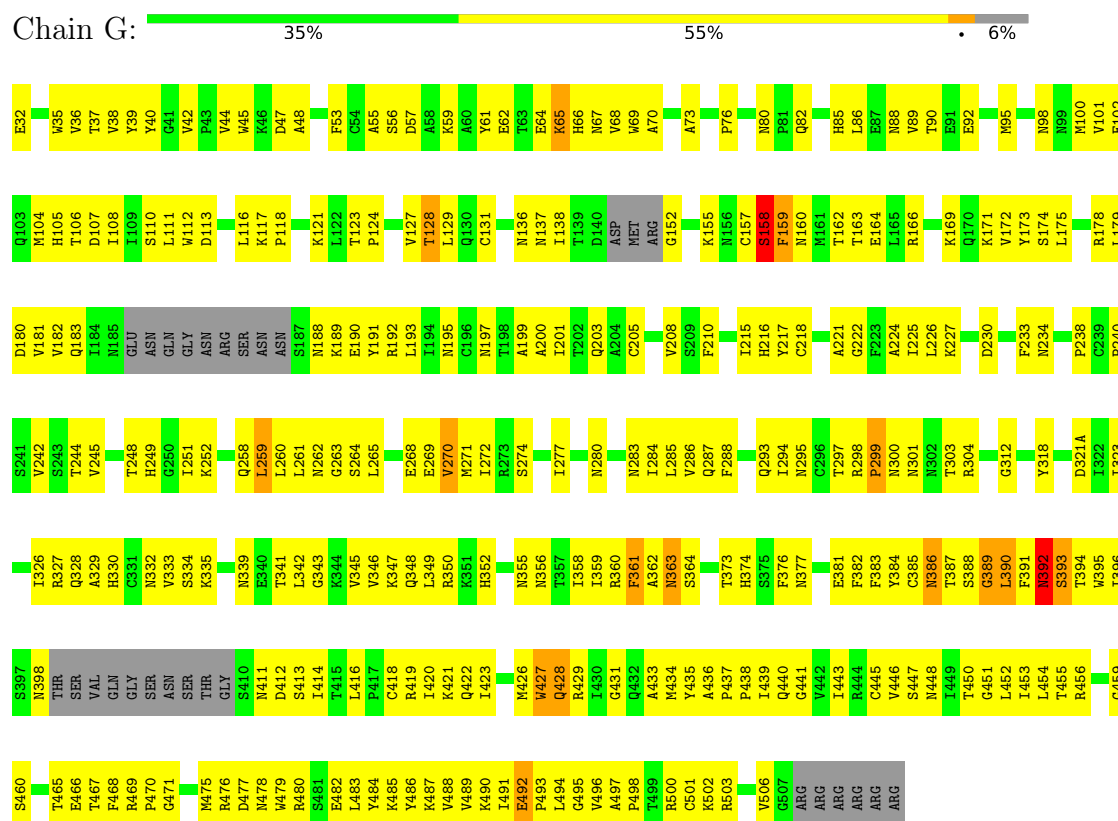


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
16	G	1	Total	C	N	O	0	0
			14	8	1	5		
16	B	1	Total	C	N	O	0	0
			14	8	1	5		
16	B	1	Total	C	N	O	0	0
			14	8	1	5		

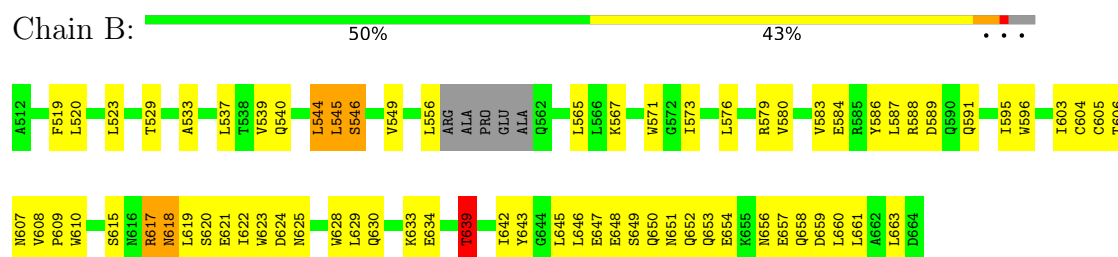
3 Residue-property plots

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.

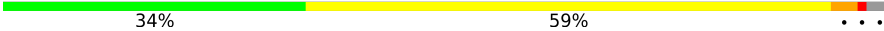
• Molecule 1: Envelope glycoprotein gp160

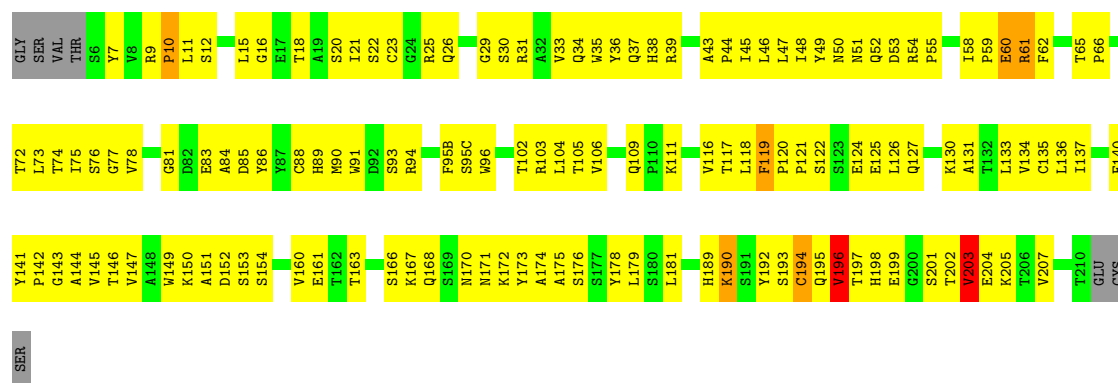


• Molecule 2: Envelope glycoprotein gp160



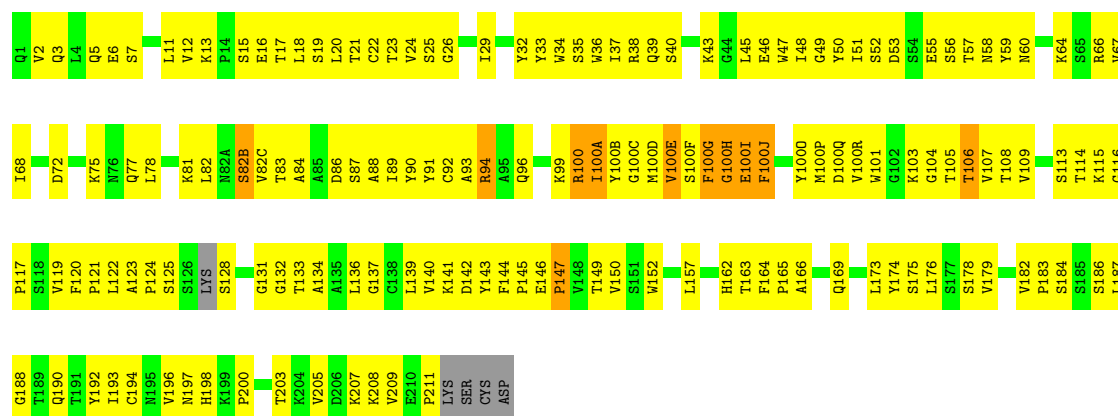
• Molecule 3: Antibody 3H+109L Fab light chain

Chain L:  34% 59% . . .



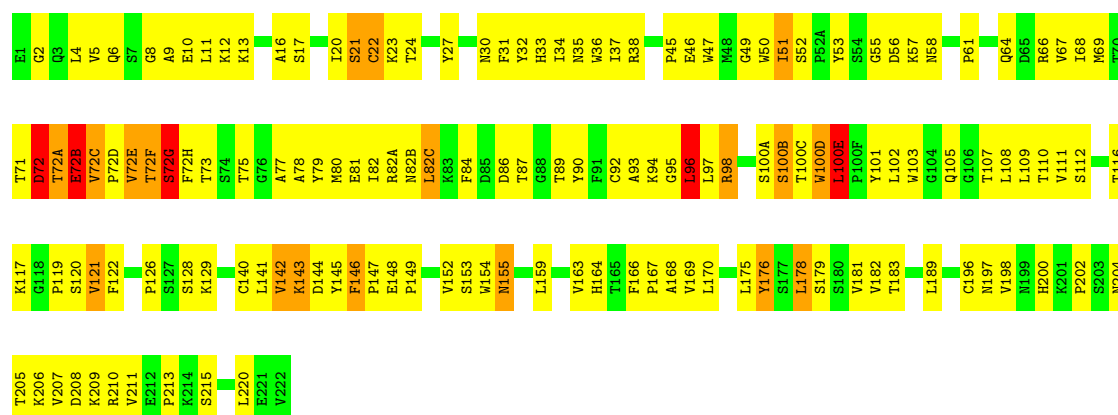
• Molecule 4: Antibody 3H+109L Fab heavy chain

Chain H:  31% 62% 5% .

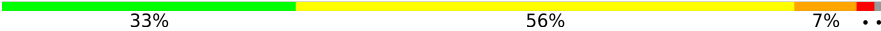


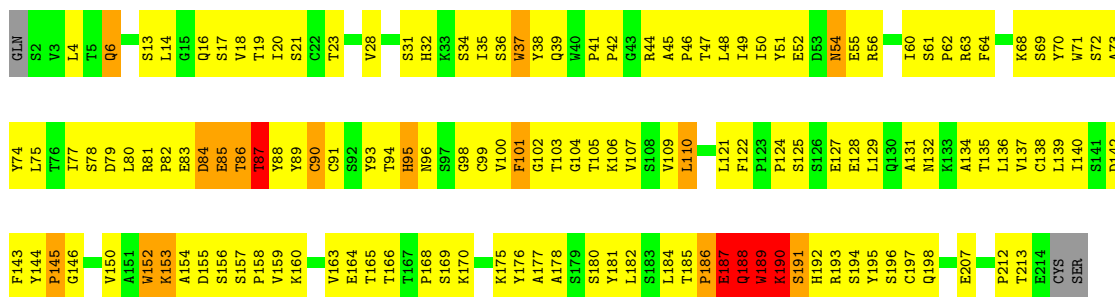
• Molecule 5: Antibody 35O22 Fab light chain

Chain D:  38% 52% 8% .

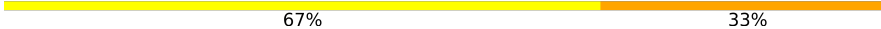


• Molecule 6: Antibody 35O22 Fab heavy chain

Chain E:  33% 56% 7% ..



- Molecule 7: alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain A:  67% 33%



- Molecule 8: alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain C:  43% 57%



- Molecule 9: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain F:  43% 57%

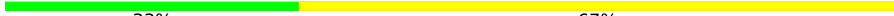


- Molecule 9: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain K:  43% 57%



- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain I:  33% 67%

MAN1
MAN2
MAN3

- Molecule 10: beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain S:  67% 33%

MAN1
MAN2
MAN3

- Molecule 11: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain J:  38% 50% 12%

MAN1
MAN2
MAN3
MAN4
MAN5
MAN6
MAN7
MAN8

- Molecule 11: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-6)-alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain M:  25% 62% 12%

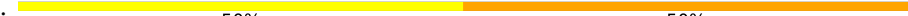
MAN1
MAN2
MAN3
MAN4
MAN5
MAN6
MAN7
MAN8

- Molecule 12: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-alpha-D-mannopyranose-(1-6)-beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain N:  17% 67% 17%

MAN1
MAN2
MAN3
MAN4
MAN5
MAN6

- Molecule 13: alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-3)-[alpha-D-mannopyranose-(1-2)-alpha-D-mannopyranose-(1-6)-[alpha-D-mannopyranose-(1-3)]alpha-D-mannopyranose-(1-6)]beta-D-mannopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose-(1-4)-2-acetamido-2-deoxy-beta-D-glucopyranose

Chain O:  50% 50%

MAN1
MAN2
MAN3
MAN4
MAN5
MAN6
MAN7
MAN8
MAN9
MAN10

- Molecule 14: α -D-mannopyranose-(1-2)- α -D-mannopyranose-(1-3)- α -D-mannopyranose-(1-6)-[α -D-mannopyranose-(1-3)] β -D-mannopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain P:  86% 14%

MAG1
MAG2
EMA3
MAN4
MAN5
MAN6
MAN7

- Molecule 15: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain Q:  50% 50%

MAG1
MAG2

- Molecule 15: 2-acetamido-2-deoxy- β -D-glucopyranose-(1-4)-2-acetamido-2-deoxy- β -D-glucopyranose

Chain R:  50% 50%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 63	Depositor
Cell constants a, b, c, α , β , γ	131.16Å 131.16Å 315.24Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	41.42 – 2.91 41.42 – 2.91	Depositor EDS
% Data completeness (in resolution range)	49.9 (41.42-2.91) 49.9 (41.42-2.91)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.45 (at 2.90Å)	Xtriage
Refinement program	PHENIX 1.11.1_2575	Depositor
R, R_{free}	0.316 , 0.339 0.320 , 0.346	Depositor DCC
R_{free} test set	1972 reflections (5.48%)	wwPDB-VP
Wilson B-factor (Å ²)	65.3	Xtriage
Anisotropy	0.129	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.15 , 69.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.340 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	12436	wwPDB-VP
Average B, all atoms (Å ²)	76.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.34% 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: MAN, NAG, BMA

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	G	0.44	1/3611 (0.0%)	0.79	10/4903 (0.2%)
2	B	0.45	0/1186	0.81	5/1608 (0.3%)
3	L	0.62	3/1646 (0.2%)	0.90	6/2247 (0.3%)
4	H	0.40	0/1787	0.81	2/2436 (0.1%)
5	D	0.51	2/1860 (0.1%)	0.94	10/2533 (0.4%)
6	E	0.88	9/1659 (0.5%)	1.23	20/2269 (0.9%)
All	All	0.55	15/11749 (0.1%)	0.91	53/15996 (0.3%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	G	0	3
3	L	0	3
4	H	0	2
5	D	0	4
6	E	0	7
All	All	0	19

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	E	85	GLU	CA-C	-15.07	1.36	1.53
6	E	187	GLU	CA-C	-12.47	1.35	1.52
3	L	61	ARG	CA-C	-11.94	1.40	1.52
3	L	196	VAL	CA-C	-9.61	1.41	1.53
6	E	85	GLU	CA-CB	9.34	1.64	1.52
1	G	390	LEU	N-CA	-8.47	1.36	1.46
6	E	188	GLN	CA-C	7.64	1.63	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	D	176	TYR	CA-C	-7.30	1.42	1.53
3	L	22	SER	CA-C	6.09	1.61	1.53
6	E	86	THR	CA-CB	6.00	1.62	1.53
6	E	189	TRP	CA-C	-5.92	1.45	1.52
6	E	190	LYS	CA-CB	-5.61	1.43	1.53
6	E	85	GLU	N-CA	-5.57	1.39	1.46
5	D	22	CYS	CB-SG	-5.57	1.62	1.81
6	E	191	SER	CA-C	5.13	1.59	1.52

All (53) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	E	85	GLU	N-CA-C	15.83	133.36	109.62
1	G	390	LEU	N-CA-C	12.89	125.06	111.14
3	L	196	VAL	N-CA-C	12.88	124.77	110.21
6	E	87	THR	N-CA-C	11.26	126.87	110.10
6	E	54	ASN	N-CA-C	11.17	123.54	111.36
6	E	188	GLN	CA-C-N	10.56	138.34	120.71
6	E	188	GLN	C-N-CA	10.56	138.34	120.71
2	B	544	LEU	N-CA-C	9.66	121.57	111.14
6	E	86	THR	OG1-CB-CG2	-8.88	91.54	109.30
1	G	393	SER	N-CA-C	8.52	122.10	109.24
6	E	140	ILE	N-CA-C	8.31	119.81	108.17
6	E	191	SER	N-CA-C	-7.98	102.56	111.82
6	E	187	GLU	N-CA-C	7.72	121.96	112.54
6	E	86	THR	CA-C-N	7.63	131.59	120.82
6	E	86	THR	C-N-CA	7.63	131.59	120.82
1	G	389	GLY	CA-C-N	-7.61	110.09	120.44
1	G	389	GLY	C-N-CA	-7.61	110.09	120.44
2	B	618	ASN	N-CA-C	7.30	120.77	109.60
5	D	176	TYR	N-CA-C	7.27	121.14	110.17
5	D	178	LEU	CD1-CG-CD2	-7.02	95.35	110.80
2	B	639	THR	N-CA-C	6.74	118.28	111.07
5	D	98	ARG	N-CA-C	6.59	126.47	113.29
3	L	61	ARG	N-CA-C	6.58	121.69	112.93
1	G	363	ASN	N-CA-C	6.50	120.14	110.52
3	L	21	ILE	N-CA-CB	6.39	121.78	111.23
6	E	190	LYS	N-CA-CB	6.39	120.78	110.42
5	D	96	LEU	N-CA-C	6.37	116.98	108.38
3	L	60	GLU	CB-CA-C	-6.26	103.45	111.22
5	D	82(C)	LEU	CB-CG-CD1	-6.19	92.14	110.70
1	G	158	SER	CA-C-N	-6.13	109.47	121.91

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	158	SER	C-N-CA	-6.13	109.47	121.91
6	E	188	GLN	CB-CG-CD	-6.12	102.20	112.60
1	G	159	PHE	N-CA-C	6.05	117.74	108.42
3	L	20	SER	CB-CA-C	-6.04	97.83	111.30
5	D	100(E)	LEU	CA-CB-CG	5.87	136.84	116.30
2	B	639	THR	CB-CA-C	-5.70	101.93	110.88
4	H	194	CYS	CA-CB-SG	5.70	127.51	114.40
5	D	22	CYS	CB-CA-C	-5.60	103.01	110.79
5	D	176	TYR	CA-C-O	-5.59	114.61	121.88
6	E	86	THR	CB-CA-C	5.46	122.58	111.78
6	E	188	GLN	CA-CB-CG	5.41	124.92	114.10
3	L	20	SER	N-CA-C	5.35	116.82	110.19
2	B	617	ARG	N-CA-C	5.34	117.42	110.53
4	H	106	THR	N-CA-C	5.31	118.28	109.46
5	D	143	LYS	N-CA-C	5.25	117.60	110.35
1	G	392	ASN	N-CA-CB	-5.22	101.62	111.43
5	D	22	CYS	CA-CB-SG	5.22	126.40	114.40
6	E	85	GLU	N-CA-CB	-5.17	104.22	110.53
6	E	85	GLU	CA-CB-CG	5.15	124.39	114.10
1	G	386	ASN	OD1-CG-ND2	5.08	127.69	122.60
6	E	84	ASP	CA-C-N	-5.04	114.60	122.05
6	E	84	ASP	C-N-CA	-5.04	114.60	122.05
6	E	189	TRP	N-CA-C	5.03	117.54	111.40

There are no chirality outliers.

All (19) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
5	D	100(B)	SER	Peptide
5	D	146	PHE	Peptide
5	D	21	SER	Peptide
5	D	96	LEU	Mainchain
6	E	101	PHE	Peptide
6	E	187	GLU	Mainchain
6	E	188	GLN	Mainchain
6	E	28	VAL	Peptide
6	E	34	SER	Peptide
6	E	37	TRP	Peptide
6	E	90	CYS	Peptide
1	G	128	THR	Peptide
1	G	158	SER	Peptide
1	G	492	GLU	Peptide

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Mol	Chain	Res	Type	Group
4	H	100(H)	GLY	Peptide
4	H	100(J)	PHE	Peptide
3	L	119	PHE	Peptide
3	L	202	THR	Peptide
3	L	203	VAL	Peptide

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	G	3538	0	3469	386	0
2	B	1167	0	1158	118	0
3	L	1603	0	1546	199	0
4	H	1744	0	1710	241	0
5	D	1813	0	1784	244	0
6	E	1615	0	1544	225	0
7	A	72	0	61	9	0
8	C	83	0	70	9	0
9	F	83	0	70	6	0
9	K	83	0	70	8	0
10	I	39	0	34	2	0
10	S	39	0	34	0	0
11	J	94	0	79	1	0
11	M	94	0	79	2	0
12	N	72	0	61	1	0
13	O	116	0	97	11	0
14	P	83	0	70	3	0
15	Q	28	0	25	2	0
15	R	28	0	25	0	0
16	B	28	0	26	4	0
16	G	14	0	13	1	0
All	All	12436	0	12025	1360	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 56.

All (1360) 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:G:234:ASN:HD21	9:K:1:NAG:C1	1.12	1.60
5:D:22:CYS:O	5:D:23:LYS:HE2	1.42	1.15
5:D:22:CYS:C	5:D:23:LYS:HE2	1.72	1.13
4:H:113:SER:O	4:H:144:PHE:CE2	2.01	1.13
1:G:95:MET:HE2	1:G:484:TYR:HB2	1.30	1.12
1:G:226:LEU:HD11	1:G:244:THR:HA	1.28	1.08
2:B:618:ASN:HB2	2:B:621:GLU:CD	1.79	1.07
1:G:456:ARG:HA	1:G:468:PHE:HB2	1.30	1.06
3:L:37:GLN:HA	3:L:86:TYR:HD1	1.20	1.05
1:G:234:ASN:CG	9:K:1:NAG:C1	2.30	1.05
5:D:69:MET:HG2	5:D:80:MET:HA	1.38	1.05
1:G:452:LEU:HD12	1:G:454:LEU:HD11	1.36	1.04
5:D:34:ILE:O	5:D:35:ASN:ND2	1.88	1.03
1:G:364:SER:HA	1:G:469:ARG:HD3	1.39	1.03
1:G:271:MET:C	1:G:272:ILE:HD12	1.84	1.03
1:G:385:CYS:SG	1:G:418:CYS:SG	1.18	1.00
5:D:166:PHE:O	5:D:178:LEU:HD22	1.59	1.00
3:L:150:LYS:HB3	3:L:193:SER:HB2	1.43	0.99
6:E:4:LEU:HD11	6:E:102:GLY:HA2	1.45	0.99
6:E:49:ILE:HD11	6:E:64:PHE:HB3	1.43	0.99
2:B:639:THR:O	2:B:643:TYR:CD1	2.17	0.98
4:H:2:VAL:HG23	4:H:26:GLY:HA3	1.45	0.98
6:E:153:LYS:HE2	6:E:159:VAL:H	1.27	0.97
1:G:160:ASN:HB3	7:A:1:NAG:C7	1.96	0.96
4:H:149:THR:HG22	4:H:197:ASN:OD1	1.66	0.96
1:G:100:MET:HB2	1:G:483:LEU:HD21	1.47	0.96
5:D:98:ARG:HH12	5:D:100(C):THR:HA	1.30	0.95
1:G:259:LEU:HD11	1:G:374:HIS:NE2	1.80	0.95
4:H:33:TYR:HE1	4:H:53:ASP:H	1.16	0.93
4:H:119:VAL:HG11	4:H:205:VAL:HG11	1.48	0.93
3:L:54:ARG:HH11	3:L:58:ILE:HD13	1.29	0.93
4:H:33:TYR:HB3	4:H:50:TYR:HE2	1.33	0.93
1:G:385:CYS:SG	1:G:418:CYS:CB	2.58	0.92
6:E:68:LYS:HE2	6:E:72:SER:HB3	1.52	0.92
1:G:38:VAL:HG12	1:G:496:VAL:HG12	1.51	0.91
1:G:261:LEU:HD11	1:G:447:SER:HB3	1.53	0.91
6:E:49:ILE:HD11	6:E:64:PHE:CB	2.01	0.91
1:G:129:LEU:HG	1:G:159:PHE:HB2	1.50	0.91
3:L:37:GLN:HA	3:L:86:TYR:CD1	2.05	0.90
3:L:54:ARG:NH2	3:L:62:PHE:O	2.05	0.89
1:G:272:ILE:HG12	1:G:286:VAL:HA	1.54	0.89
5:D:122:PHE:O	5:D:141:LEU:HB2	1.71	0.89

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:259:LEU:HD11	1:G:374:HIS:CD2	2.08	0.89
4:H:88:ALA:HB3	4:H:90:TYR:HE1	1.34	0.89
5:D:98:ARG:NH1	5:D:100(C):THR:OG1	2.06	0.88
5:D:144:ASP:HB2	5:D:175:LEU:HD13	1.55	0.88
5:D:66:ARG:HH22	5:D:82(C):LEU:HD11	1.37	0.88
1:G:506:VAL:HG21	2:B:658:GLN:HE21	1.39	0.87
5:D:23:LYS:HD2	5:D:77:ALA:HA	1.55	0.87
1:G:361:PHE:CE1	1:G:470:PRO:HD3	2.09	0.87
1:G:234:ASN:OD1	9:K:1:NAG:C1	2.23	0.86
1:G:391:PHE:HE2	1:G:470:PRO:HG3	1.38	0.86
2:B:648:GLU:HA	2:B:652:GLN:HB2	1.57	0.86
5:D:152:VAL:HG12	5:D:198:VAL:HA	1.57	0.86
2:B:618:ASN:CB	2:B:621:GLU:CD	2.49	0.85
6:E:14:LEU:HG	6:E:81:ARG:HH12	1.42	0.85
1:G:252:LYS:NZ	1:G:262:ASN:O	2.10	0.85
1:G:385:CYS:CB	1:G:418:CYS:SG	2.64	0.85
3:L:10:PRO:HA	3:L:103:ARG:HB3	1.59	0.85
3:L:54:ARG:NH1	3:L:59:PRO:O	2.10	0.85
5:D:66:ARG:NH2	5:D:82(C):LEU:HD11	1.92	0.85
6:E:137:VAL:HG12	6:E:181:TYR:CB	2.06	0.85
5:D:96:LEU:HD11	5:D:100(E):LEU:N	1.92	0.85
5:D:10:GLU:HG2	5:D:202:PRO:HD3	1.57	0.84
5:D:23:LYS:HZ1	5:D:78:ALA:HB3	1.42	0.84
6:E:80:LEU:HD21	6:E:84:ASP:HB2	1.58	0.84
4:H:137:GLY:HA3	4:H:179:VAL:HG22	1.59	0.84
3:L:137:ILE:HD11	3:L:175:ALA:HB3	1.58	0.84
1:G:335:LYS:O	1:G:339:ASN:ND2	2.09	0.84
6:E:143:PHE:CE2	6:E:146:GLY:HA2	2.12	0.84
1:G:391:PHE:CE2	1:G:470:PRO:HG3	2.13	0.84
3:L:54:ARG:NH1	3:L:58:ILE:HD13	1.92	0.83
1:G:160:ASN:HB3	7:A:1:NAG:N2	1.92	0.83
5:D:35:ASN:HB2	5:D:94:LYS:HE3	1.61	0.83
6:E:121:LEU:HG	6:E:138:CYS:HB2	1.60	0.82
3:L:195:GLN:HB3	3:L:204:GLU:HB3	1.61	0.82
2:B:607:ASN:HD21	2:B:650:GLN:HA	1.44	0.82
5:D:93:ALA:HB3	5:D:103:TRP:HA	1.60	0.81
1:G:304:ARG:HH11	1:G:440:GLN:HB3	1.45	0.81
6:E:13:SER:OG	6:E:110:LEU:O	1.96	0.81
1:G:201:ILE:HG13	1:G:433:ALA:HB3	1.63	0.81
5:D:121:VAL:HG23	5:D:141:LEU:HD12	1.63	0.81
5:D:141:LEU:HD21	5:D:143:LYS:HB3	1.63	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:35:ASN:HD21	5:D:50:TRP:HB3	1.44	0.81
1:G:203:GLN:HE21	1:G:437:PRO:HG3	1.45	0.80
4:H:115:LYS:HB3	4:H:144:PHE:HE1	1.44	0.80
5:D:12:LYS:NZ	5:D:13:LYS:O	2.14	0.80
1:G:361:PHE:HE1	1:G:470:PRO:HD3	1.47	0.80
1:G:454:LEU:HA	1:G:470:PRO:HA	1.63	0.80
2:B:618:ASN:OD1	16:B:702:NAG:N2	2.15	0.80
2:B:639:THR:O	2:B:643:TYR:HD1	1.60	0.80
1:G:95:MET:CE	1:G:484:TYR:HB2	2.10	0.80
1:G:469:ARG:HG3	1:G:470:PRO:HD2	1.64	0.79
1:G:333:VAL:O	1:G:414:ILE:N	2.16	0.79
2:B:596:TRP:HE1	2:B:647:GLU:CD	1.90	0.79
3:L:143:GLY:HA3	3:L:173:TYR:HD2	1.46	0.79
1:G:386:ASN:HD21	14:P:1:NAG:C7	1.96	0.79
2:B:639:THR:O	2:B:643:TYR:CE1	2.34	0.79
5:D:166:PHE:CE2	6:E:177:ALA:HB3	2.17	0.79
6:E:79:ASP:OD1	6:E:81:ARG:NH2	2.14	0.79
7:A:1:NAG:H83	7:A:1:NAG:H3	1.65	0.79
1:G:476:ARG:HG3	1:G:480:ARG:HH12	1.47	0.78
6:E:176:TYR:HD1	6:E:177:ALA:H	1.31	0.78
3:L:96:TRP:HA	4:H:47:TRP:CZ3	2.18	0.78
4:H:114:THR:HG21	4:H:200:PRO:C	2.09	0.78
3:L:167:LYS:HB2	3:L:173:TYR:HD1	1.48	0.77
5:D:35:ASN:ND2	5:D:50:TRP:HB3	1.99	0.77
6:E:68:LYS:HD3	6:E:74:TYR:HB2	1.64	0.77
1:G:38:VAL:HG22	2:B:604:CYS:HB2	1.66	0.77
2:B:639:THR:HG22	2:B:643:TYR:CE1	2.20	0.77
1:G:381:GLU:OE1	1:G:438:PRO:HA	1.82	0.76
3:L:84:ALA:H	3:L:104:LEU:HD11	1.50	0.76
5:D:20:ILE:HG12	5:D:107:THR:HG21	1.66	0.76
4:H:115:LYS:HB3	4:H:144:PHE:CE1	2.21	0.76
6:E:14:LEU:HG	6:E:81:ARG:NH1	2.01	0.76
4:H:113:SER:O	4:H:144:PHE:HE2	1.66	0.76
5:D:121:VAL:HA	5:D:141:LEU:HD12	1.68	0.76
5:D:4:LEU:HD21	5:D:22:CYS:HB3	1.66	0.75
1:G:131:CYS:HB3	1:G:157:CYS:HA	1.67	0.75
1:G:364:SER:CA	1:G:469:ARG:HD3	2.17	0.75
3:L:134:VAL:HG11	3:L:178:TYR:CD2	2.22	0.75
5:D:23:LYS:NZ	5:D:78:ALA:HB3	2.00	0.75
5:D:142:VAL:HG21	5:D:178:LEU:HG	1.67	0.75
1:G:188:ASN:N	1:G:189:LYS:HA	2.01	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:271:MET:O	1:G:272:ILE:HD12	1.86	0.75
3:L:142:PRO:O	3:L:198:HIS:NE2	2.19	0.75
5:D:22:CYS:O	5:D:23:LYS:CE	2.30	0.75
6:E:49:ILE:HG13	6:E:60:ILE:HD13	1.67	0.75
1:G:475:MET:O	1:G:478:ASN:N	2.16	0.75
4:H:134:ALA:N	4:H:182:VAL:O	2.17	0.74
3:L:195:GLN:HE22	3:L:205:LYS:C	1.94	0.74
2:B:608:VAL:H	2:B:650:GLN:NE2	1.85	0.74
3:L:30:SER:HB3	4:H:100:ARG:HH21	1.52	0.74
4:H:11:LEU:HD12	4:H:108:THR:O	1.86	0.74
5:D:166:PHE:CD2	5:D:167:PRO:HD2	2.23	0.74
5:D:96:LEU:HD11	5:D:100(E):LEU:HB3	1.70	0.74
6:E:155:ASP:HB2	6:E:192:HIS:HB2	1.68	0.74
1:G:452:LEU:HD12	1:G:454:LEU:CD1	2.16	0.74
3:L:35:TRP:HB3	3:L:47:LEU:HD21	1.69	0.74
6:E:86:THR:HG22	6:E:87:THR:HG23	1.69	0.74
1:G:164:GLU:HA	1:G:312:GLY:HA2	1.70	0.73
6:E:42:PRO:HD3	6:E:86:THR:HG21	1.70	0.73
4:H:136:LEU:HD11	4:H:209:VAL:HG11	1.70	0.73
6:E:42:PRO:CD	6:E:86:THR:HG21	2.17	0.73
5:D:72(E):VAL:HB	5:D:73:THR:HG23	1.70	0.73
1:G:195:ASN:OD1	1:G:201:ILE:HD11	1.87	0.73
5:D:119:PRO:HA	5:D:145:TYR:CE1	2.24	0.73
3:L:204:GLU:C	3:L:205:LYS:HD3	2.14	0.73
4:H:132:GLY:O	4:H:184:SER:N	2.17	0.73
2:B:596:TRP:HD1	2:B:647:GLU:HB3	1.54	0.73
2:B:606:THR:HG23	2:B:650:GLN:HE22	1.54	0.72
1:G:358:ILE:HG12	1:G:465:THR:HA	1.71	0.72
2:B:606:THR:HG23	2:B:650:GLN:NE2	2.04	0.72
5:D:51:ILE:HG13	5:D:52:SER:H	1.52	0.72
6:E:23:THR:HG22	6:E:72:SER:CB	2.18	0.72
3:L:31:ARG:NH2	3:L:65:THR:O	2.22	0.72
3:L:48:ILE:HG22	3:L:54:ARG:HG2	1.71	0.72
2:B:618:ASN:HB3	2:B:621:GLU:OE1	1.89	0.72
6:E:47:THR:C	6:E:48:LEU:HD23	2.14	0.72
1:G:439:ILE:HD12	1:G:443:ILE:HG21	1.71	0.72
3:L:167:LYS:HB2	3:L:173:TYR:CD1	2.24	0.72
6:E:13:SER:CB	6:E:110:LEU:O	2.37	0.72
3:L:38:HIS:O	3:L:84:ALA:HB1	1.90	0.72
6:E:44:ARG:NH1	6:E:45:ALA:O	2.23	0.72
5:D:69:MET:HG2	5:D:80:MET:CA	2.19	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:134:VAL:HG12	3:L:178:TYR:HA	1.72	0.71
2:B:639:THR:HG22	2:B:643:TYR:HE1	1.55	0.71
3:L:150:LYS:NZ	3:L:154:SER:H	1.88	0.71
3:L:54:ARG:HH11	3:L:58:ILE:CD1	2.03	0.71
6:E:189:TRP:O	6:E:192:HIS:ND1	2.23	0.71
2:B:596:TRP:CD1	2:B:647:GLU:HB3	2.25	0.71
3:L:54:ARG:HH12	3:L:59:PRO:C	1.97	0.71
6:E:49:ILE:HD11	6:E:64:PHE:CG	2.25	0.71
6:E:137:VAL:HG12	6:E:181:TYR:HB2	1.72	0.71
1:G:121:LYS:HZ1	1:G:200:ALA:HB1	1.55	0.71
1:G:171:LYS:HG3	7:A:1:NAG:O7	1.91	0.71
6:E:184:LEU:HA	6:E:189:TRP:CZ2	2.26	0.71
5:D:98:ARG:HH12	5:D:100(C):THR:CA	2.04	0.71
6:E:16:GLN:NE2	6:E:18:VAL:HB	2.06	0.71
5:D:169:VAL:HG22	5:D:170:LEU:H	1.56	0.70
1:G:476:ARG:HG3	1:G:480:ARG:NH1	2.06	0.70
4:H:152:TRP:HB3	4:H:157:LEU:HD21	1.73	0.70
5:D:154:TRP:CE2	5:D:196:CYS:HB3	2.26	0.70
1:G:360:ARG:HG3	1:G:467:THR:HA	1.73	0.70
3:L:189:HIS:HD1	3:L:190:LYS:H	1.38	0.70
5:D:72(C):VAL:N	5:D:75:THR:O	2.21	0.70
4:H:16:GLU:O	4:H:82(B):SER:N	2.23	0.70
4:H:88:ALA:HB3	4:H:90:TYR:CE1	2.23	0.70
2:B:647:GLU:O	2:B:651:ASN:N	2.17	0.70
3:L:124:GLU:O	3:L:127:GLN:NE2	2.21	0.70
4:H:20:LEU:HD12	4:H:36:TRP:HH2	1.55	0.70
4:H:48:ILE:HA	4:H:60:ASN:HB2	1.73	0.70
5:D:51:ILE:HD13	5:D:57:LYS:HE2	1.73	0.70
3:L:121:PRO:HD3	3:L:133:LEU:HB2	1.72	0.70
1:G:269:GLU:HB2	1:G:271:MET:HE1	1.72	0.70
4:H:149:THR:HG22	4:H:197:ASN:CG	2.16	0.70
1:G:428:GLN:O	1:G:428:GLN:NE2	2.23	0.70
5:D:153:SER:C	5:D:154:TRP:HD1	1.99	0.70
4:H:35:SER:N	4:H:93:ALA:O	2.22	0.70
5:D:142:VAL:HG21	5:D:178:LEU:CG	2.22	0.70
9:K:1:NAG:H4	9:K:2:NAG:N2	2.06	0.69
1:G:39:TYR:HD2	2:B:603:ILE:HG22	1.57	0.69
2:B:654:GLU:O	2:B:658:GLN:NE2	2.26	0.69
4:H:149:THR:CG2	4:H:197:ASN:OD1	2.39	0.69
5:D:121:VAL:HA	5:D:141:LEU:CD1	2.22	0.69
5:D:69:MET:HG2	5:D:81:GLU:H	1.56	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:69:TRP:HA	1:G:111:LEU:HD11	1.74	0.69
1:G:226:LEU:HD11	1:G:244:THR:CA	2.15	0.69
1:G:271:MET:C	1:G:272:ILE:CD1	2.64	0.69
4:H:3:GLN:OE1	4:H:3:GLN:N	2.26	0.69
3:L:195:GLN:HB3	3:L:204:GLU:CB	2.23	0.69
4:H:99:LYS:HD3	13:O:9:MAN:H62	1.74	0.69
6:E:184:LEU:HA	6:E:189:TRP:CH2	2.27	0.69
1:G:288:PHE:N	1:G:450:THR:O	2.24	0.69
3:L:43:ALA:HA	4:H:91:TYR:CE2	2.29	0.68
4:H:35:SER:HB2	4:H:100(P):MET:HE1	1.73	0.68
4:H:113:SER:O	4:H:144:PHE:CD2	2.46	0.68
3:L:149:TRP:CE3	3:L:194:CYS:HB2	2.28	0.68
9:F:3:BMA:H4	9:F:4:MAN:H2	1.75	0.68
1:G:183:GLN:HB3	1:G:191:TYR:CD1	2.29	0.68
3:L:61:ARG:NH2	3:L:74:THR:O	2.26	0.68
5:D:196:CYS:SG	5:D:209:LYS:HD3	2.34	0.68
1:G:129:LEU:HG	1:G:159:PHE:CB	2.22	0.68
2:B:576:LEU:HA	2:B:579:ARG:HE	1.58	0.68
6:E:96:ASN:HB2	8:C:6:MAN:H2	1.74	0.68
3:L:18:THR:OG1	3:L:76:SER:HA	1.93	0.68
5:D:84:PHE:HE2	5:D:112:SER:HA	1.58	0.68
4:H:59:TYR:HB2	4:H:64:LYS:HE2	1.76	0.68
4:H:100(A):ILE:HG22	4:H:100(J):PHE:HB3	1.75	0.68
1:G:138:ILE:HG23	10:I:1:NAG:H83	1.76	0.67
1:G:303:THR:OG1	1:G:321(A):ASP:O	2.10	0.67
1:G:350:ARG:NH1	1:G:398:ASN:OD1	2.28	0.67
4:H:94:ARG:H	4:H:101:TRP:CD1	2.11	0.67
1:G:210:PHE:HE2	1:G:434:MET:HE1	1.60	0.67
3:L:163:THR:HG21	3:L:175:ALA:HA	1.75	0.67
4:H:20:LEU:HD12	4:H:36:TRP:CH2	2.29	0.67
4:H:100(D):MET:HB3	13:O:2:NAG:H2	1.77	0.67
1:G:152:GLY:HA3	1:G:178:ARG:HE	1.60	0.67
2:B:586:TYR:CD1	2:B:587:LEU:HD23	2.29	0.67
3:L:37:GLN:HB3	3:L:45:ILE:HG13	1.75	0.67
1:G:270:VAL:C	1:G:272:ILE:HD13	2.20	0.67
3:L:35:TRP:HE1	3:L:73:LEU:HD11	1.59	0.67
3:L:95(C):SER:O	4:H:47:TRP:HZ3	1.77	0.67
2:B:545:LEU:HD23	2:B:586:TYR:CG	2.29	0.67
4:H:22:CYS:HB3	4:H:78:LEU:HD21	1.77	0.67
4:H:100(D):MET:C	4:H:100(I):GLU:HG3	2.20	0.67
6:E:110:LEU:N	6:E:110:LEU:HD12	2.10	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:544:LEU:HD11	2:B:586:TYR:HB2	1.77	0.67
1:G:261:LEU:HG	1:G:448:ASN:C	2.20	0.67
4:H:33:TYR:HE1	4:H:53:ASP:N	1.92	0.67
4:H:207:LYS:NZ	4:H:208:LYS:O	2.28	0.67
1:G:502:LYS:HA	2:B:605:CYS:HB3	1.76	0.66
6:E:49:ILE:CD1	6:E:64:PHE:CD2	2.79	0.66
1:G:106:THR:O	1:G:110:SER:OG	2.12	0.66
1:G:453:ILE:C	1:G:454:LEU:HD12	2.21	0.66
5:D:2:GLY:HA3	5:D:27:TYR:CD1	2.30	0.66
6:E:152:TRP:CZ2	6:E:160:LYS:HA	2.30	0.66
1:G:37:THR:HB	2:B:604:CYS:O	1.96	0.66
1:G:225:ILE:HA	1:G:488:VAL:HA	1.77	0.66
3:L:35:TRP:O	3:L:46:LEU:HD11	1.96	0.66
5:D:93:ALA:CB	5:D:103:TRP:HA	2.24	0.66
5:D:169:VAL:HG23	6:E:166:THR:HG22	1.78	0.66
9:K:1:NAG:H4	9:K:2:NAG:C7	2.26	0.66
5:D:11:LEU:HB2	5:D:147:PRO:HG3	1.78	0.66
5:D:122:PHE:H	5:D:141:LEU:HG	1.61	0.66
7:A:1:NAG:H3	7:A:1:NAG:C8	2.25	0.66
1:G:323:ILE:HD11	12:N:1:NAG:H61	1.76	0.65
2:B:618:ASN:HB2	2:B:621:GLU:OE2	1.95	0.65
6:E:68:LYS:HG2	6:E:73:ALA:HA	1.77	0.65
3:L:44:PRO:HD3	4:H:91:TYR:HE2	1.61	0.65
4:H:117:PRO:HD2	4:H:203:THR:HG21	1.77	0.65
1:G:350:ARG:HG3	1:G:355:ASN:HA	1.77	0.65
1:G:422:GLN:O	1:G:436:ALA:N	2.17	0.65
4:H:100:ARG:NH2	13:O:4:MAN:O6	2.29	0.65
5:D:71:THR:HB	5:D:78:ALA:HA	1.78	0.65
1:G:304:ARG:HH12	1:G:439:ILE:HA	1.59	0.65
5:D:108:LEU:HD22	5:D:149:PRO:HD3	1.78	0.65
3:L:126:LEU:HD23	3:L:127:GLN:H	1.60	0.65
9:K:3:BMA:H62	9:K:7:MAN:H5	1.77	0.65
6:E:50:ILE:HD11	6:E:54:ASN:C	2.21	0.65
1:G:37:THR:HG22	2:B:606:THR:HB	1.79	0.65
4:H:125:SER:HG	4:H:128:SER:N	1.94	0.65
1:G:386:ASN:ND2	14:P:1:NAG:C7	2.59	0.65
4:H:84:ALA:HA	4:H:109:VAL:HG13	1.78	0.65
5:D:126:PRO:HB2	5:D:215:SER:HB2	1.79	0.65
1:G:112:TRP:CG	1:G:427:TRP:HH2	2.14	0.65
1:G:362:ALA:C	1:G:469:ARG:HH21	2.05	0.65
4:H:141:LYS:NZ	4:H:142:ASP:OD1	2.30	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:304:ARG:HH12	1:G:439:ILE:CA	2.10	0.64
1:G:476:ARG:HA	1:G:479:TRP:CD1	2.32	0.64
2:B:618:ASN:CB	2:B:621:GLU:OE1	2.45	0.64
6:E:189:TRP:CD1	6:E:189:TRP:N	2.65	0.64
4:H:59:TYR:HB2	4:H:64:LYS:HB3	1.78	0.64
3:L:104:LEU:HD12	3:L:104:LEU:O	1.98	0.64
1:G:127:VAL:C	1:G:129:LEU:HD13	2.23	0.64
1:G:271:MET:SD	1:G:271:MET:N	2.70	0.64
2:B:544:LEU:HD12	2:B:586:TYR:HD2	1.62	0.64
3:L:170:ASN:O	3:L:171:ASN:HB2	1.96	0.64
4:H:123:ALA:HB1	4:H:211:PRO:HA	1.79	0.64
6:E:190:LYS:HZ1	6:E:212:PRO:C	2.05	0.64
1:G:327:ARG:HH12	4:H:100(C):GLY:HA3	1.63	0.64
6:E:129:LEU:O	6:E:132:ASN:ND2	2.28	0.64
1:G:421:LYS:NZ	1:G:423:ILE:O	2.23	0.64
3:L:35:TRP:HD1	3:L:48:ILE:HG12	1.61	0.64
3:L:195:GLN:NE2	3:L:205:LYS:O	2.30	0.64
5:D:4:LEU:O	5:D:105:GLN:NE2	2.31	0.64
5:D:100(D):TRP:CZ2	6:E:93:TYR:HB2	2.33	0.64
1:G:359:ILE:N	1:G:396:ILE:HD11	2.12	0.64
3:L:25:ARG:NH2	3:L:88:CYS:SG	2.71	0.63
3:L:149:TRP:CZ3	3:L:179:LEU:HG	2.33	0.63
5:D:93:ALA:C	5:D:94:LYS:HD2	2.23	0.63
1:G:131:CYS:HB3	1:G:157:CYS:CA	2.28	0.63
4:H:18:LEU:O	4:H:81:LYS:NZ	2.27	0.63
1:G:286:VAL:HB	1:G:452:LEU:HG	1.79	0.63
5:D:168:ALA:HA	5:D:178:LEU:HB3	1.80	0.63
2:B:653:GLN:HG2	2:B:656:ASN:ND2	2.13	0.63
5:D:72(B):GLU:CD	5:D:72(B):GLU:H	2.06	0.63
6:E:42:PRO:HD3	6:E:86:THR:CG2	2.27	0.63
9:K:3:BMA:O4	9:K:7:MAN:O6	2.15	0.63
5:D:5:VAL:HA	5:D:105:GLN:NE2	2.14	0.63
1:G:48:ALA:HB2	1:G:490:LYS:HB2	1.78	0.63
6:E:136:LEU:HD21	6:E:195:TYR:CE2	2.33	0.63
4:H:186:SER:O	4:H:190:GLN:NE2	2.26	0.63
5:D:100(D):TRP:HZ2	6:E:93:TYR:HB2	1.64	0.63
1:G:53:PHE:O	1:G:218:CYS:N	2.32	0.62
4:H:100(D):MET:SD	13:O:1:NAG:H5	2.40	0.62
1:G:65:LYS:HB3	1:G:66:HIS:HD2	1.65	0.62
1:G:270:VAL:O	1:G:272:ILE:HD13	1.99	0.62
4:H:72:ASP:OD1	4:H:75:LYS:N	2.31	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:90:THR:HG22	1:G:240:PRO:HA	1.81	0.62
2:B:544:LEU:HD13	2:B:544:LEU:C	2.24	0.62
3:L:171:ASN:C	3:L:172:LYS:HD3	2.25	0.62
5:D:30:ASN:HD21	5:D:53:TYR:HB2	1.64	0.62
5:D:81:GLU:OE1	5:D:82(A):ARG:NH1	2.31	0.62
5:D:95:GLY:O	5:D:96:LEU:HG	1.99	0.62
8:C:2:NAG:H83	8:C:2:NAG:H3	1.81	0.62
1:G:259:LEU:O	1:G:260:LEU:HG	1.99	0.62
5:D:96:LEU:HD11	5:D:100(E):LEU:CB	2.29	0.62
1:G:57:ASP:O	1:G:59:LYS:NZ	2.33	0.62
1:G:261:LEU:CD1	1:G:447:SER:HB3	2.28	0.62
3:L:58:ILE:HB	3:L:62:PHE:HZ	1.63	0.62
6:E:137:VAL:HG12	6:E:181:TYR:HB3	1.77	0.62
1:G:393:SER:HB3	1:G:395:TRP:NE1	2.15	0.62
3:L:44:PRO:HD3	4:H:91:TYR:CE2	2.35	0.62
4:H:100(D):MET:N	4:H:100(I):GLU:OE2	2.28	0.62
4:H:169:GLN:HG2	4:H:173:LEU:H	1.63	0.62
5:D:168:ALA:HB1	5:D:176:TYR:HB3	1.80	0.62
4:H:133:THR:HA	4:H:183:PRO:HA	1.82	0.62
6:E:6:GLN:HE21	6:E:102:GLY:HA3	1.65	0.62
6:E:20:ILE:HG21	6:E:88:TYR:HB2	1.80	0.62
1:G:361:PHE:CZ	1:G:470:PRO:HD3	2.35	0.62
3:L:61:ARG:NH2	3:L:76:SER:H	1.98	0.62
4:H:34:TRP:N	4:H:50:TYR:OH	2.26	0.62
6:E:36:SER:C	6:E:37:TRP:CD1	2.78	0.62
1:G:284:ILE:N	1:G:454:LEU:O	2.32	0.62
1:G:298:ARG:HG2	1:G:420:ILE:HD11	1.80	0.62
4:H:100(A):ILE:HD11	13:O:3:BMA:H62	1.81	0.62
1:G:39:TYR:CD2	2:B:603:ILE:HG22	2.35	0.61
1:G:113:ASP:HA	1:G:116:LEU:HG	1.81	0.61
3:L:39:ARG:NH2	3:L:81:GLY:O	2.33	0.61
4:H:100(P):MET:HB3	4:H:101:TRP:HE1	1.65	0.61
1:G:421:LYS:HE2	1:G:422:GLN:H	1.64	0.61
3:L:25:ARG:NE	3:L:88:CYS:O	2.34	0.61
4:H:143:TYR:OH	4:H:176:LEU:N	2.32	0.61
4:H:140:VAL:HG12	4:H:143:TYR:CD2	2.35	0.61
5:D:120:SER:O	5:D:145:TYR:OH	2.17	0.61
5:D:154:TRP:O	5:D:155:ASN:ND2	2.33	0.61
5:D:164:HIS:H	5:D:182:VAL:HG22	1.65	0.61
5:D:10:GLU:HG2	5:D:202:PRO:CD	2.29	0.61
6:E:37:TRP:HA	6:E:38:TYR:CE2	2.35	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:335:LYS:HB3	1:G:414:ILE:HD11	1.81	0.61
6:E:31:SER:OG	6:E:32:HIS:N	2.34	0.61
5:D:96:LEU:CD1	5:D:100(E):LEU:HB3	2.30	0.61
6:E:41:PRO:HB2	6:E:44:ARG:HB3	1.80	0.61
1:G:101:VAL:HA	1:G:104:MET:HG2	1.81	0.61
3:L:195:GLN:CD	3:L:204:GLU:HB3	2.26	0.61
6:E:37:TRP:CE2	6:E:90:CYS:HB2	2.36	0.61
1:G:225:ILE:HG21	1:G:486:TYR:HD2	1.65	0.61
4:H:91:TYR:CE1	4:H:104:GLY:HA3	2.35	0.61
6:E:35:ILE:HG23	6:E:91:CYS:O	2.01	0.61
1:G:334:SER:HA	1:G:413:SER:HA	1.83	0.61
3:L:84:ALA:N	3:L:104:LEU:HD11	2.16	0.61
5:D:143:LYS:HG3	5:D:145:TYR:CZ	2.36	0.61
6:E:23:THR:HG22	6:E:72:SER:OG	2.00	0.61
6:E:84:ASP:OD1	6:E:88:TYR:OH	2.15	0.61
3:L:84:ALA:HB3	3:L:86:TYR:CZ	2.36	0.60
3:L:86:TYR:HE2	3:L:104:LEU:HD21	1.64	0.60
3:L:143:GLY:HA3	3:L:173:TYR:CD2	2.33	0.60
1:G:248:THR:HA	1:G:486:TYR:CE2	2.36	0.60
1:G:393:SER:CB	1:G:395:TRP:HE1	2.14	0.60
6:E:49:ILE:HD11	6:E:64:PHE:CD2	2.35	0.60
4:H:113:SER:C	4:H:144:PHE:CE2	2.80	0.60
3:L:37:GLN:HG3	3:L:86:TYR:HE1	1.65	0.60
3:L:61:ARG:CZ	3:L:75:ILE:HA	2.30	0.60
4:H:33:TYR:HB3	4:H:50:TYR:CE2	2.25	0.60
4:H:34:TRP:H	4:H:50:TYR:HH	1.45	0.60
1:G:259:LEU:HD21	1:G:374:HIS:CG	2.37	0.60
6:E:50:ILE:HD12	6:E:56:ARG:HB3	1.83	0.60
1:G:359:ILE:H	1:G:396:ILE:HD11	1.67	0.60
5:D:121:VAL:HA	5:D:141:LEU:HG	1.84	0.60
1:G:304:ARG:NH2	1:G:438:PRO:O	2.29	0.60
3:L:61:ARG:NH1	3:L:75:ILE:HA	2.17	0.60
5:D:68:ILE:O	5:D:81:GLU:HG2	2.02	0.60
5:D:96:LEU:HD11	5:D:100(E):LEU:CA	2.32	0.60
6:E:6:GLN:NE2	6:E:102:GLY:HA3	2.17	0.60
3:L:59:PRO:HG2	3:L:62:PHE:HE1	1.66	0.60
6:E:80:LEU:HD21	6:E:84:ASP:CB	2.29	0.60
3:L:189:HIS:HD1	3:L:190:LYS:N	2.00	0.60
4:H:115:LYS:NZ	4:H:142:ASP:OD2	2.29	0.60
1:G:39:TYR:HH	2:B:623:TRP:HH2	1.49	0.60
2:B:544:LEU:HD12	2:B:586:TYR:CD2	2.37	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:36:SER:OG	6:E:51:TYR:HA	2.02	0.60
6:E:16:GLN:CD	6:E:18:VAL:H	2.10	0.59
5:D:53:TYR:CD2	8:C:1:NAG:H5	2.37	0.59
5:D:96:LEU:HD11	5:D:100(D):TRP:C	2.28	0.59
1:G:163:THR:HG22	1:G:164:GLU:H	1.66	0.59
5:D:34:ILE:HD13	5:D:51:ILE:HG23	1.84	0.59
5:D:119:PRO:HD2	5:D:205:THR:HB	1.85	0.59
1:G:295:ASN:OD1	1:G:332:ASN:HB2	2.02	0.59
1:G:363:ASN:ND2	1:G:388:SER:HA	2.17	0.59
4:H:39:GLN:O	4:H:88:ALA:HB1	2.02	0.59
6:E:19:THR:HG21	6:E:75:LEU:HG	1.85	0.59
1:G:95:MET:HE2	1:G:484:TYR:CB	2.20	0.59
6:E:13:SER:O	6:E:16:GLN:HB2	2.02	0.59
1:G:327:ARG:HH11	4:H:100(B):TYR:HD2	1.43	0.59
1:G:491:ILE:HG21	2:B:544:LEU:HD23	1.85	0.59
3:L:125:GLU:HG2	3:L:130:LYS:O	2.03	0.59
4:H:150:VAL:HG12	4:H:196:VAL:HG12	1.85	0.59
4:H:190:GLN:HG2	4:H:192:TYR:CE1	2.38	0.59
6:E:143:PHE:HE2	6:E:146:GLY:HA2	1.66	0.59
1:G:39:TYR:CE1	1:G:497:ALA:HB3	2.37	0.59
1:G:363:ASN:HD21	1:G:388:SER:HA	1.68	0.59
2:B:533:ALA:HB3	2:B:628:TRP:HE1	1.66	0.59
1:G:280:ASN:HA	1:G:456:ARG:HB2	1.84	0.58
3:L:33:VAL:HG22	3:L:90:MET:HA	1.84	0.58
5:D:119:PRO:HA	5:D:145:TYR:CZ	2.37	0.58
6:E:68:LYS:HZ2	6:E:74:TYR:HD2	1.51	0.58
3:L:31:ARG:NH1	3:L:51:ASN:OD1	2.37	0.58
5:D:69:MET:CG	5:D:80:MET:HA	2.24	0.58
6:E:37:TRP:HD1	6:E:91:CYS:HB3	1.69	0.58
6:E:176:TYR:HD1	6:E:177:ALA:N	1.99	0.58
5:D:96:LEU:CG	5:D:100(E):LEU:HB3	2.34	0.58
6:E:49:ILE:HD12	6:E:64:PHE:CD2	2.39	0.58
3:L:119:PHE:H	3:L:133:LEU:HD11	1.69	0.58
5:D:12:LYS:HD2	5:D:111:VAL:HG22	1.86	0.58
1:G:195:ASN:CG	1:G:201:ILE:HD11	2.28	0.58
6:E:189:TRP:O	6:E:192:HIS:CG	2.56	0.58
5:D:69:MET:HG2	5:D:81:GLU:N	2.18	0.58
5:D:89:THR:HA	5:D:108:LEU:HA	1.86	0.58
6:E:164:GLU:N	6:E:164:GLU:OE1	2.36	0.58
6:E:6:GLN:NE2	6:E:103:THR:H	2.02	0.58
6:E:168:PRO:HB2	6:E:176:TYR:CE1	2.39	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:95(B):PHE:CE1	4:H:47:TRP:HH2	2.22	0.57
3:L:168:GLN:HE22	3:L:173:TYR:C	2.12	0.57
4:H:5:GLN:HE21	4:H:5:GLN:HA	1.68	0.57
5:D:51:ILE:HD12	5:D:57:LYS:N	2.18	0.57
1:G:501:CYS:SG	1:G:502:LYS:N	2.78	0.57
3:L:61:ARG:HG2	3:L:61:ARG:HH21	1.68	0.57
4:H:149:THR:HG22	4:H:197:ASN:H	1.68	0.57
6:E:170:LYS:H	6:E:170:LYS:HD3	1.69	0.57
3:L:12:SER:OG	3:L:103:ARG:NH1	2.37	0.57
4:H:11:LEU:HD11	4:H:108:THR:H	1.69	0.57
4:H:12:VAL:HG23	4:H:16:GLU:OE1	2.03	0.57
4:H:94:ARG:O	4:H:94:ARG:HG3	2.04	0.57
5:D:189:LEU:HD22	5:D:213:PRO:HB2	1.86	0.57
5:D:90:TYR:N	5:D:107:THR:O	2.20	0.57
4:H:18:LEU:HD23	4:H:19:SER:N	2.19	0.57
5:D:9:ALA:CB	5:D:108:LEU:H	2.17	0.57
1:G:469:ARG:HG3	1:G:470:PRO:CD	2.32	0.57
2:B:586:TYR:HD1	2:B:587:LEU:HD23	1.70	0.57
3:L:46:LEU:HD12	3:L:47:LEU:H	1.70	0.57
6:E:68:LYS:HG3	6:E:69:SER:H	1.70	0.57
1:G:36:VAL:HG13	2:B:610:TRP:HB2	1.86	0.57
5:D:6:GLN:H	5:D:105:GLN:HG2	1.70	0.57
13:O:1:NAG:H3	13:O:1:NAG:H83	1.86	0.57
2:B:621:GLU:OE1	2:B:621:GLU:N	2.37	0.57
1:G:116:LEU:HD12	1:G:117:LYS:N	2.20	0.57
2:B:620:SER:OG	5:D:97:LEU:CD2	2.53	0.57
3:L:167:LYS:HD2	3:L:171:ASN:C	2.30	0.57
2:B:643:TYR:HA	2:B:646:LEU:CD1	2.35	0.56
3:L:47:LEU:O	3:L:54:ARG:HB3	2.04	0.56
5:D:142:VAL:CG2	5:D:178:LEU:HG	2.34	0.56
7:A:1:NAG:H83	7:A:1:NAG:C3	2.33	0.56
4:H:29:ILE:HA	4:H:34:TRP:CH2	2.39	0.56
4:H:37:ILE:HB	4:H:45:LEU:HD21	1.88	0.56
4:H:144:PHE:HA	4:H:173:LEU:HD11	1.88	0.56
5:D:36:TRP:HE1	5:D:80:MET:HE3	1.70	0.56
6:E:94:THR:O	6:E:96:ASN:N	2.39	0.56
1:G:127:VAL:HG23	1:G:129:LEU:CD2	2.35	0.56
1:G:159:PHE:CE1	1:G:173:TYR:HA	2.39	0.56
1:G:304:ARG:NH1	1:G:440:GLN:N	2.54	0.56
1:G:502:LYS:NZ	1:G:503:ARG:O	2.38	0.56
4:H:51:ILE:HD13	4:H:57:THR:HB	1.86	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:109:LEU:CD2	5:D:111:VAL:HG23	2.36	0.56
2:B:629:LEU:HD12	2:B:629:LEU:H	1.71	0.56
5:D:78:ALA:HB1	5:D:80:MET:HE1	1.88	0.56
5:D:120:SER:N	5:D:145:TYR:OH	2.38	0.56
1:G:363:ASN:OD1	1:G:388:SER:HA	2.05	0.56
4:H:15:SER:HA	4:H:82(B):SER:HA	1.87	0.56
5:D:34:ILE:HD13	5:D:51:ILE:CG2	2.36	0.56
5:D:100(C):THR:HG22	5:D:100(D):TRP:CZ3	2.41	0.56
1:G:271:MET:SD	1:G:348:GLN:OE1	2.63	0.56
1:G:439:ILE:CD1	1:G:443:ILE:HG21	2.35	0.56
5:D:84:PHE:CZ	5:D:111:VAL:HG12	2.41	0.56
5:D:179:SER:HB3	6:E:139:LEU:CD2	2.36	0.56
1:G:506:VAL:HG21	2:B:658:GLN:NE2	2.15	0.56
3:L:109:GLN:OE1	3:L:109:GLN:N	2.36	0.56
5:D:72(A):THR:HG23	5:D:72(B):GLU:OE1	2.06	0.56
5:D:167:PRO:HG3	6:E:169:SER:HB3	1.88	0.56
1:G:181:VAL:HG12	1:G:193:LEU:HA	1.88	0.56
2:B:618:ASN:OD1	16:B:702:NAG:C2	2.52	0.56
5:D:69:MET:CG	5:D:81:GLU:H	2.18	0.56
1:G:271:MET:HA	1:G:348:GLN:CD	2.31	0.55
1:G:362:ALA:HB3	1:G:469:ARG:NH2	2.20	0.55
2:B:523:LEU:HD12	2:B:540:GLN:NE2	2.21	0.55
3:L:86:TYR:CE2	3:L:104:LEU:HD21	2.40	0.55
9:F:3:BMA:O2	9:F:4:MAN:H3	2.07	0.55
1:G:393:SER:HB3	1:G:395:TRP:HE1	1.72	0.55
3:L:149:TRP:CD2	3:L:194:CYS:HB2	2.41	0.55
4:H:115:LYS:HG2	4:H:116:GLY:N	2.20	0.55
6:E:100:VAL:C	6:E:101:PHE:HD1	2.14	0.55
6:E:153:LYS:HD2	6:E:157:SER:N	2.22	0.55
1:G:270:VAL:HG13	1:G:287:GLN:O	2.06	0.55
2:B:619:LEU:HD22	2:B:623:TRP:HE1	1.70	0.55
4:H:140:VAL:HG12	4:H:143:TYR:CE2	2.40	0.55
5:D:32:TYR:CD1	5:D:98:ARG:HG3	2.42	0.55
5:D:51:ILE:HG13	5:D:55:GLY:HA2	1.88	0.55
5:D:153:SER:C	5:D:154:TRP:CD1	2.83	0.55
10:I:1:NAG:H62	10:I:2:NAG:N2	2.21	0.55
1:G:427:TRP:HD1	1:G:429:ARG:HB2	1.71	0.55
5:D:68:ILE:C	5:D:69:MET:HG3	2.31	0.55
1:G:188:ASN:H	1:G:189:LYS:HA	1.72	0.55
3:L:146:THR:O	3:L:197:THR:HG23	2.05	0.55
5:D:50:TRP:NE1	5:D:58:ASN:HB3	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:68:LYS:HG2	6:E:72:SER:C	2.32	0.55
1:G:128:THR:C	1:G:129:LEU:CD1	2.80	0.55
1:G:387:THR:HG23	1:G:390:LEU:HD12	1.89	0.55
5:D:72:ASP:HB2	5:D:77:ALA:HB3	1.89	0.55
6:E:61:SER:HB3	6:E:62:PRO:CD	2.36	0.55
1:G:259:LEU:CD1	1:G:374:HIS:NE2	2.64	0.55
2:B:656:ASN:HA	2:B:659:ASP:HB3	1.89	0.55
2:B:660:LEU:HA	2:B:663:LEU:HD12	1.89	0.55
3:L:205:LYS:HD3	3:L:205:LYS:N	2.22	0.55
5:D:23:LYS:CE	5:D:23:LYS:HA	2.35	0.55
5:D:38:ARG:HG2	5:D:46:GLU:HB3	1.89	0.55
6:E:134:ALA:HB3	6:E:184:LEU:HG	1.89	0.55
4:H:149:THR:HG22	4:H:197:ASN:N	2.21	0.55
2:B:588:ARG:NH1	2:B:589:ASP:OD1	2.40	0.55
2:B:647:GLU:HB2	2:B:652:GLN:OE1	2.07	0.55
4:H:67:VAL:HG12	4:H:82:LEU:HD12	1.89	0.55
6:E:68:LYS:HG3	6:E:69:SER:N	2.22	0.55
1:G:95:MET:CE	1:G:484:TYR:CB	2.83	0.55
1:G:230:ASP:OD1	1:G:233:PHE:HB2	2.07	0.55
3:L:50:ASN:HD21	13:O:5:MAN:H61	1.70	0.55
4:H:16:GLU:OE2	4:H:16:GLU:N	2.40	0.55
4:H:83:THR:O	4:H:109:VAL:HG11	2.07	0.55
1:G:85:HIS:HD1	1:G:86:LEU:N	2.05	0.54
1:G:387:THR:HG22	1:G:387:THR:O	2.06	0.54
2:B:651:ASN:HA	2:B:654:GLU:HB3	1.88	0.54
3:L:58:ILE:HB	3:L:62:PHE:CZ	2.42	0.54
3:L:145:VAL:HG23	3:L:198:HIS:N	2.23	0.54
4:H:145:PRO:HA	4:H:174:TYR:CE1	2.42	0.54
6:E:84:ASP:O	6:E:88:TYR:OH	2.24	0.54
6:E:192:HIS:HD2	6:E:194:SER:N	2.05	0.54
1:G:203:GLN:HB2	1:G:435:TYR:HD2	1.73	0.54
2:B:573:ILE:HG13	2:B:576:LEU:HD22	1.89	0.54
1:G:295:ASN:HA	1:G:445:CYS:O	2.07	0.54
3:L:86:TYR:O	3:L:102:THR:HG22	2.06	0.54
3:L:151:ALA:O	3:L:154:SER:N	2.41	0.54
4:H:100(J):PHE:N	4:H:100(J):PHE:CD1	2.75	0.54
1:G:67:ASN:OD1	1:G:68:VAL:HG12	2.08	0.54
1:G:121:LYS:NZ	1:G:201:ILE:H	2.06	0.54
3:L:58:ILE:HD12	3:L:58:ILE:O	2.08	0.54
3:L:15:LEU:HD13	3:L:106:VAL:HG11	1.89	0.54
4:H:100(D):MET:O	4:H:100(I):GLU:HG3	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:31:PHE:HA	8:C:1:NAG:O6	2.07	0.54
5:D:121:VAL:CG2	5:D:141:LEU:HD12	2.36	0.54
5:D:121:VAL:HG22	5:D:209:LYS:NZ	2.22	0.54
1:G:65:LYS:HB3	1:G:66:HIS:CD2	2.43	0.54
1:G:287:GLN:HE21	1:G:450:THR:CB	2.21	0.54
2:B:565:LEU:O	2:B:567:LYS:NZ	2.41	0.54
4:H:82:LEU:HD22	4:H:82(C):VAL:HG22	1.90	0.54
4:H:136:LEU:HD13	4:H:137:GLY:N	2.22	0.54
5:D:23:LYS:HE2	5:D:23:LYS:HA	1.90	0.54
6:E:72:SER:OG	6:E:73:ALA:N	2.40	0.54
1:G:174:SER:OG	1:G:175:LEU:N	2.41	0.54
1:G:395:TRP:O	1:G:396:ILE:HD13	2.08	0.54
3:L:83:GLU:HA	3:L:104:LEU:CD1	2.38	0.54
3:L:150:LYS:NZ	3:L:154:SER:N	2.56	0.54
3:L:167:LYS:HA	3:L:173:TYR:HA	1.89	0.54
6:E:75:LEU:O	6:E:75:LEU:HD12	2.08	0.54
1:G:159:PHE:O	1:G:171:LYS:HA	2.08	0.53
1:G:200:ALA:O	1:G:433:ALA:N	2.24	0.53
2:B:624:ASP:O	5:D:97:LEU:O	2.26	0.53
3:L:18:THR:HG23	3:L:75:ILE:C	2.32	0.53
3:L:96:TRP:HA	4:H:47:TRP:HZ3	1.71	0.53
3:L:116:VAL:HG22	3:L:137:ILE:HG22	1.90	0.53
3:L:35:TRP:HE1	3:L:73:LEU:CD1	2.20	0.53
3:L:203:VAL:O	3:L:204:GLU:HG3	2.08	0.53
4:H:66:ARG:NH1	4:H:82:LEU:HD21	2.23	0.53
6:E:37:TRP:CD1	6:E:90:CYS:HA	2.43	0.53
1:G:268:GLU:HG2	1:G:269:GLU:H	1.74	0.53
3:L:34:GLN:OE1	3:L:91:TRP:NE1	2.42	0.53
4:H:60:ASN:C	4:H:64:LYS:HE3	2.33	0.53
4:H:100(G):PHE:HB3	4:H:100(I):GLU:CG	2.38	0.53
6:E:37:TRP:CZ2	6:E:90:CYS:HB2	2.43	0.53
6:E:63:ARG:HB3	6:E:78:SER:O	2.09	0.53
1:G:183:GLN:HA	1:G:191:TYR:HA	1.91	0.53
2:B:617:ARG:HH12	2:B:621:GLU:HB3	1.74	0.53
3:L:199:GLU:CG	3:L:201:SER:HB2	2.39	0.53
4:H:149:THR:CG2	4:H:197:ASN:CG	2.80	0.53
6:E:39:GLN:HA	6:E:88:TYR:CD1	2.44	0.53
4:H:139:LEU:HD23	4:H:140:VAL:N	2.23	0.53
6:E:16:GLN:HE22	6:E:18:VAL:HB	1.72	0.53
4:H:11:LEU:HD11	4:H:108:THR:N	2.23	0.53
4:H:152:TRP:HH2	4:H:178:SER:HB2	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:150:VAL:HG23	6:E:198:GLN:O	2.09	0.53
1:G:259:LEU:HA	1:G:452:LEU:HA	1.90	0.53
3:L:50:ASN:C	3:L:52:GLN:H	2.17	0.53
3:L:195:GLN:CB	3:L:204:GLU:HB3	2.36	0.53
6:E:105:THR:O	6:E:106:LYS:HG2	2.08	0.53
6:E:129:LEU:HD21	6:E:134:ALA:HB2	1.89	0.53
1:G:389:GLY:HA2	15:Q:1:NAG:O6	2.08	0.53
4:H:121:PRO:HB2	4:H:209:VAL:HG13	1.91	0.53
5:D:9:ALA:HB1	5:D:108:LEU:H	1.74	0.53
6:E:39:GLN:HA	6:E:88:TYR:HD1	1.74	0.53
1:G:39:TYR:OH	2:B:623:TRP:HH2	1.92	0.53
3:L:25:ARG:HG3	3:L:90:MET:SD	2.49	0.53
4:H:115:LYS:CB	4:H:144:PHE:CE1	2.92	0.53
5:D:34:ILE:O	5:D:50:TRP:HB3	2.09	0.53
6:E:168:PRO:HB3	6:E:178:ALA:HB2	1.91	0.53
1:G:73:ALA:HB3	2:B:571:TRP:HD1	1.74	0.53
1:G:491:ILE:HG22	1:G:493:PRO:HD3	1.90	0.53
5:D:121:VAL:HG22	5:D:209:LYS:HD2	1.91	0.53
6:E:20:ILE:HB	6:E:105:THR:HG21	1.91	0.53
6:E:36:SER:O	6:E:38:TYR:CZ	2.61	0.53
2:B:645:LEU:HD23	2:B:648:GLU:HB3	1.90	0.52
4:H:96:GLN:N	4:H:100(O):TYR:O	2.42	0.52
5:D:34:ILE:O	5:D:35:ASN:CG	2.52	0.52
5:D:206:LYS:HE2	5:D:208:ASP:OD1	2.10	0.52
6:E:37:TRP:HA	6:E:38:TYR:CD2	2.44	0.52
1:G:335:LYS:HE2	1:G:412:ASP:HB3	1.91	0.52
2:B:607:ASN:ND2	2:B:650:GLN:HA	2.21	0.52
2:B:625:ASN:HB2	5:D:97:LEU:HG	1.91	0.52
1:G:107:ASP:OD2	1:G:217:TYR:OH	2.22	0.52
1:G:298:ARG:HB3	1:G:329:ALA:HA	1.91	0.52
1:G:342:LEU:O	1:G:346:VAL:HG23	2.08	0.52
1:G:360:ARG:HE	1:G:467:THR:C	2.17	0.52
5:D:121:VAL:HA	5:D:141:LEU:CG	2.38	0.52
5:D:155:ASN:ND2	5:D:159:LEU:HB3	2.24	0.52
1:G:39:TYR:HE1	1:G:497:ALA:HB3	1.74	0.52
1:G:427:TRP:CD1	1:G:429:ARG:HB2	2.45	0.52
4:H:6:GLU:HB3	4:H:20:LEU:HD11	1.90	0.52
5:D:140:CYS:SG	5:D:209:LYS:HD3	2.49	0.52
6:E:38:TYR:O	6:E:88:TYR:HA	2.09	0.52
2:B:544:LEU:CD1	2:B:586:TYR:HB2	2.39	0.52
2:B:545:LEU:HD23	2:B:586:TYR:CD1	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:9:ALA:H	5:D:10:GLU:CD	2.18	0.52
5:D:11:LEU:HD11	5:D:112:SER:HB2	1.90	0.52
6:E:124:PRO:HB2	6:E:128:GLU:OE1	2.09	0.52
1:G:205:CYS:O	1:G:208:VAL:HG22	2.10	0.52
1:G:287:GLN:HA	1:G:450:THR:O	2.10	0.52
5:D:23:LYS:NZ	5:D:78:ALA:CB	2.73	0.52
6:E:17:SER:O	6:E:17:SER:OG	2.23	0.52
6:E:68:LYS:HG2	6:E:73:ALA:CA	2.37	0.52
6:E:87:THR:HG22	6:E:106:LYS:HG2	1.92	0.52
6:E:110:LEU:N	6:E:110:LEU:CD1	2.73	0.52
6:E:135:THR:HA	6:E:182:LEU:O	2.10	0.52
3:L:34:GLN:HG2	4:H:100(O):TYR:HB3	1.90	0.52
3:L:86:TYR:HE2	3:L:104:LEU:CD2	2.22	0.52
3:L:119:PHE:CZ	4:H:125:SER:HB3	2.45	0.52
5:D:95:GLY:HA3	5:D:101:TYR:HB3	1.90	0.52
1:G:69:TRP:HA	1:G:111:LEU:CD1	2.40	0.52
1:G:265:LEU:HD11	1:G:288:PHE:HB2	1.92	0.52
1:G:294:ILE:HD13	1:G:333:VAL:HG22	1.90	0.52
3:L:54:ARG:NH1	3:L:60:GLU:HA	2.24	0.52
3:L:122:SER:CB	4:H:120:PHE:HB3	2.40	0.52
3:L:122:SER:HB2	4:H:120:PHE:HB3	1.91	0.52
5:D:121:VAL:CA	5:D:141:LEU:HD12	2.39	0.52
6:E:192:HIS:CE1	6:E:195:TYR:CZ	2.97	0.52
16:B:702:NAG:H83	6:E:55:GLU:CD	2.34	0.52
4:H:100(H):GLY:N	4:H:100(I):GLU:HG2	2.25	0.52
5:D:21:SER:HA	5:D:36:TRP:HH2	1.74	0.52
1:G:101:VAL:HG11	1:G:480:ARG:CD	2.40	0.52
1:G:299:PRO:HD2	1:G:327:ARG:HB3	1.90	0.52
3:L:35:TRP:O	3:L:36:TYR:CG	2.63	0.52
4:H:136:LEU:O	4:H:179:VAL:HG13	2.10	0.52
1:G:181:VAL:HG12	1:G:193:LEU:HG	1.92	0.51
3:L:84:ALA:H	3:L:104:LEU:CD1	2.20	0.51
3:L:119:PHE:C	3:L:133:LEU:HD13	2.35	0.51
5:D:9:ALA:HB1	5:D:108:LEU:HB3	1.92	0.51
6:E:99:CYS:HB3	6:E:101:PHE:HE1	1.75	0.51
6:E:153:LYS:HD2	6:E:157:SER:H	1.74	0.51
7:A:1:NAG:C8	7:A:1:NAG:C3	2.87	0.51
1:G:261:LEU:HD21	1:G:448:ASN:H	1.76	0.51
1:G:364:SER:HA	1:G:469:ARG:CD	2.26	0.51
2:B:647:GLU:CG	2:B:652:GLN:HE22	2.23	0.51
4:H:59:TYR:HD2	4:H:64:LYS:HB2	1.74	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:21:SER:OG	5:D:79:TYR:HA	2.10	0.51
6:E:134:ALA:N	6:E:184:LEU:HG	2.25	0.51
1:G:80:ASN:O	1:G:82:GLN:NE2	2.38	0.51
1:G:271:MET:HA	1:G:348:GLN:OE1	2.10	0.51
1:G:361:PHE:HE1	1:G:470:PRO:CD	2.21	0.51
3:L:37:GLN:OE1	3:L:45:ILE:HD11	2.10	0.51
4:H:100(B):TYR:CE2	4:H:100(I):GLU:HA	2.45	0.51
4:H:166:ALA:HB2	4:H:176:LEU:HB2	1.92	0.51
5:D:51:ILE:HG13	5:D:52:SER:N	2.21	0.51
1:G:327:ARG:NH1	4:H:100(C):GLY:HA3	2.24	0.51
1:G:376:PHE:CE1	1:G:383:PHE:HB2	2.46	0.51
4:H:145:PRO:HD2	4:H:198:HIS:CE1	2.45	0.51
1:G:62:GLU:OE2	1:G:62:GLU:N	2.41	0.51
1:G:102:GLU:O	1:G:106:THR:OG1	2.20	0.51
1:G:107:ASP:O	1:G:110:SER:N	2.43	0.51
2:B:647:GLU:HG3	2:B:652:GLN:HE22	1.74	0.51
4:H:66:ARG:HH11	4:H:82:LEU:HD11	1.76	0.51
5:D:11:LEU:HD11	5:D:112:SER:CB	2.40	0.51
5:D:98:ARG:CZ	5:D:98:ARG:HA	2.40	0.51
5:D:121:VAL:CG2	5:D:209:LYS:HD2	2.40	0.51
1:G:101:VAL:HG13	1:G:483:LEU:HD23	1.93	0.51
3:L:38:HIS:HB3	3:L:85:ASP:O	2.11	0.51
4:H:5:GLN:HA	4:H:5:GLN:NE2	2.25	0.51
4:H:16:GLU:HG2	4:H:17:THR:N	2.26	0.51
4:H:149:THR:HG22	4:H:197:ASN:CB	2.39	0.51
6:E:86:THR:HG22	6:E:87:THR:H	1.75	0.51
6:E:152:TRP:HD1	6:E:152:TRP:O	1.94	0.51
1:G:116:LEU:HD12	1:G:117:LYS:HG2	1.92	0.51
1:G:346:VAL:HG13	1:G:359:ILE:CD1	2.41	0.51
1:G:359:ILE:O	1:G:394:THR:HG23	2.10	0.51
1:G:363:ASN:CG	1:G:388:SER:HA	2.35	0.51
3:L:37:GLN:HE21	3:L:86:TYR:HE1	1.53	0.51
4:H:89:ILE:HD11	4:H:104:GLY:HA3	1.93	0.51
5:D:35:ASN:HD21	5:D:50:TRP:HE3	1.57	0.51
5:D:144:ASP:HA	5:D:176:TYR:O	2.10	0.51
6:E:23:THR:CG2	6:E:72:SER:CB	2.89	0.51
4:H:124:PRO:HD2	4:H:211:PRO:HA	1.93	0.51
5:D:100(A):SER:OG	8:C:7:MAN:H3	2.11	0.51
1:G:363:ASN:O	1:G:469:ARG:NE	2.44	0.51
2:B:576:LEU:HA	2:B:579:ARG:NE	2.24	0.51
2:B:657:GLU:O	2:B:661:LEU:N	2.44	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:163:VAL:HG13	5:D:182:VAL:HG21	1.92	0.51
6:E:13:SER:HB2	6:E:110:LEU:O	2.10	0.51
6:E:131:ALA:N	6:E:132:ASN:HA	2.26	0.51
1:G:152:GLY:N	1:G:178:ARG:HH21	2.08	0.51
1:G:259:LEU:C	1:G:260:LEU:HG	2.35	0.51
4:H:117:PRO:HA	4:H:142:ASP:O	2.11	0.51
4:H:164:PHE:O	4:H:176:LEU:HG	2.11	0.51
1:G:44:VAL:HG13	1:G:492:GLU:HB2	1.93	0.50
1:G:200:ALA:C	1:G:201:ILE:HD12	2.35	0.50
1:G:327:ARG:NH1	4:H:100(B):TYR:CD2	2.70	0.50
1:G:373:THR:OG1	1:G:385:CYS:O	2.27	0.50
1:G:104:MET:O	1:G:108:ILE:HG13	2.12	0.50
1:G:287:GLN:HE21	1:G:450:THR:HB	1.75	0.50
4:H:100:ARG:O	4:H:100(A):ILE:HG12	2.11	0.50
5:D:96:LEU:CD1	5:D:100(D):TRP:C	2.84	0.50
5:D:142:VAL:HG11	5:D:178:LEU:HD12	1.93	0.50
6:E:37:TRP:CD1	6:E:91:CYS:HB3	2.45	0.50
6:E:185:THR:HG22	6:E:186:PRO:HD2	1.92	0.50
1:G:203:GLN:HE22	1:G:318:TYR:HB2	1.75	0.50
2:B:576:LEU:HA	2:B:579:ARG:HH21	1.76	0.50
3:L:72:THR:C	3:L:73:LEU:HD12	2.35	0.50
6:E:38:TYR:CE2	6:E:101:PHE:CE2	2.99	0.50
3:L:36:TYR:HA	3:L:46:LEU:CD1	2.41	0.50
4:H:117:PRO:HB3	4:H:143:TYR:HB3	1.93	0.50
4:H:149:THR:CG2	4:H:197:ASN:N	2.73	0.50
5:D:103:TRP:CE3	6:E:46:PRO:HD2	2.45	0.50
6:E:49:ILE:HD13	6:E:75:LEU:HD13	1.93	0.50
1:G:40:TYR:HA	1:G:494:LEU:HD22	1.92	0.50
1:G:285:LEU:HD21	1:G:477:ASP:HB3	1.93	0.50
3:L:36:TYR:HA	3:L:46:LEU:HD13	1.91	0.50
1:G:469:ARG:CZ	1:G:469:ARG:HB2	2.41	0.50
3:L:189:HIS:HD2	3:L:192:TYR:CE2	2.30	0.50
4:H:33:TYR:O	4:H:94:ARG:NH2	2.45	0.50
4:H:51:ILE:HA	4:H:56:SER:O	2.12	0.50
1:G:264:SER:H	1:G:482:GLU:CD	2.20	0.50
2:B:523:LEU:HD12	2:B:540:GLN:HE21	1.77	0.50
2:B:624:ASP:HB3	5:D:97:LEU:HB3	1.93	0.50
3:L:37:GLN:CG	3:L:86:TYR:HE1	2.25	0.50
4:H:198:HIS:NE2	4:H:200:PRO:HG2	2.27	0.50
5:D:17:SER:OG	5:D:82(A):ARG:HD3	2.10	0.50
1:G:127:VAL:HG23	1:G:129:LEU:HD21	1.94	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:96:LEU:HD21	5:D:100(E):LEU:HB2	1.94	0.50
1:G:350:ARG:CG	1:G:355:ASN:HA	2.41	0.50
6:E:68:LYS:CG	6:E:73:ALA:HA	2.41	0.50
1:G:121:LYS:HZ1	1:G:201:ILE:H	1.60	0.49
1:G:346:VAL:HG13	1:G:359:ILE:HD11	1.94	0.49
2:B:573:ILE:HA	2:B:576:LEU:HB3	1.94	0.49
3:L:143:GLY:CA	3:L:173:TYR:HD2	2.21	0.49
5:D:31:PHE:CE2	8:C:1:NAG:H4	2.46	0.49
3:L:197:THR:O	3:L:197:THR:OG1	2.28	0.49
5:D:142:VAL:O	5:D:142:VAL:HG13	2.12	0.49
3:L:29:GLY:HA3	3:L:93:SER:OG	2.11	0.49
5:D:11:LEU:HD22	5:D:147:PRO:HD3	1.95	0.49
5:D:96:LEU:HD21	5:D:100(E):LEU:CB	2.42	0.49
6:E:134:ALA:H	6:E:184:LEU:HG	1.77	0.49
1:G:335:LYS:C	1:G:339:ASN:HD22	2.16	0.49
4:H:51:ILE:HD12	4:H:56:SER:C	2.37	0.49
4:H:91:TYR:HE1	4:H:104:GLY:HA3	1.77	0.49
4:H:100(D):MET:O	4:H:100(F):SER:N	2.45	0.49
4:H:149:THR:CG2	4:H:197:ASN:H	2.24	0.49
14:P:1:NAG:O3	14:P:2:NAG:O5	2.23	0.49
2:B:608:VAL:HG21	2:B:649:SER:HB3	1.94	0.49
3:L:171:ASN:O	3:L:172:LYS:HD3	2.11	0.49
4:H:162:HIS:C	4:H:162:HIS:CD2	2.90	0.49
5:D:71:THR:CB	5:D:78:ALA:HA	2.41	0.49
5:D:93:ALA:HB1	5:D:102:LEU:O	2.12	0.49
6:E:152:TRP:H	6:E:152:TRP:CD1	2.31	0.49
1:G:327:ARG:HA	4:H:100(G):PHE:HZ	1.78	0.49
4:H:6:GLU:HG2	4:H:22:CYS:HB2	1.94	0.49
4:H:145:PRO:HB2	4:H:147:PRO:HD2	1.95	0.49
1:G:68:VAL:HG13	1:G:69:TRP:H	1.77	0.49
2:B:591:GLN:O	2:B:595:ILE:HG12	2.13	0.49
5:D:6:GLN:CD	5:D:22:CYS:SG	2.96	0.49
6:E:80:LEU:HD22	6:E:81:ARG:N	2.26	0.49
1:G:327:ARG:NH1	4:H:100(I):GLU:OE1	2.46	0.49
3:L:149:TRP:CE3	3:L:179:LEU:HG	2.47	0.49
5:D:155:ASN:HD21	5:D:159:LEU:N	2.11	0.49
6:E:23:THR:CG2	6:E:72:SER:HB2	2.43	0.49
6:E:49:ILE:O	6:E:56:ARG:HB2	2.13	0.49
6:E:192:HIS:HE1	6:E:195:TYR:CZ	2.31	0.49
1:G:216:HIS:ND1	1:G:248:THR:O	2.42	0.49
1:G:227:LYS:HG3	1:G:485:LYS:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:274:SER:OG	1:G:277:ILE:HG13	2.13	0.49
1:G:392:ASN:OD1	15:Q:1:NAG:O5	2.21	0.49
2:B:606:THR:CG2	2:B:650:GLN:HE22	2.23	0.49
5:D:61:PRO:HA	5:D:64:GLN:CD	2.38	0.49
6:E:49:ILE:HG23	6:E:50:ILE:H	1.78	0.49
1:G:294:ILE:O	1:G:447:SER:HB2	2.13	0.49
3:L:61:ARG:NH1	3:L:62:PHE:HB3	2.28	0.49
4:H:146:GLU:HB2	4:H:174:TYR:CD2	2.48	0.49
5:D:96:LEU:HD13	5:D:98:ARG:HH22	1.77	0.49
5:D:100(E):LEU:HD22	5:D:101:TYR:HB2	1.94	0.49
5:D:197:ASN:HB2	5:D:208:ASP:OD1	2.13	0.49
2:B:642:ILE:O	2:B:646:LEU:HD12	2.12	0.48
3:L:93:SER:OG	3:L:94:ARG:HG3	2.13	0.48
4:H:46:GLU:HG2	4:H:48:ILE:CD1	2.43	0.48
4:H:183:PRO:O	4:H:186:SER:HB3	2.12	0.48
6:E:121:LEU:HD12	6:E:197:CYS:SG	2.53	0.48
1:G:226:LEU:HD11	1:G:245:VAL:H	1.78	0.48
2:B:617:ARG:NH1	2:B:621:GLU:HB3	2.27	0.48
3:L:7:TYR:H	3:L:9:ARG:NH1	2.11	0.48
3:L:7:TYR:H	3:L:9:ARG:HH12	1.61	0.48
3:L:66:PRO:HB3	13:O:5:MAN:O3	2.13	0.48
3:L:160:VAL:N	3:L:161:GLU:OE2	2.46	0.48
4:H:123:ALA:HB1	4:H:124:PRO:HD2	1.95	0.48
5:D:198:VAL:HG12	5:D:207:VAL:H	1.77	0.48
6:E:49:ILE:CD1	6:E:64:PHE:CG	2.96	0.48
1:G:107:ASP:O	1:G:111:LEU:HD23	2.12	0.48
3:L:199:GLU:HG2	3:L:201:SER:HB2	1.94	0.48
4:H:86:ASP:O	4:H:90:TYR:OH	2.27	0.48
1:G:455:THR:HG23	1:G:471:GLY:HA3	1.95	0.48
3:L:47:LEU:HB2	3:L:62:PHE:CD2	2.48	0.48
3:L:51:ASN:N	3:L:51:ASN:HD22	2.11	0.48
3:L:146:THR:O	3:L:197:THR:N	2.47	0.48
5:D:100(D):TRP:HH2	6:E:94:THR:O	1.96	0.48
6:E:129:LEU:HD13	6:E:132:ASN:HB3	1.95	0.48
1:G:446:VAL:O	11:M:1:NAG:H3	2.14	0.48
1:G:258:GLN:OE1	1:G:470:PRO:HB2	2.13	0.48
2:B:618:ASN:CB	2:B:621:GLU:OE2	2.57	0.48
6:E:49:ILE:HG23	6:E:50:ILE:N	2.27	0.48
6:E:188:GLN:HB2	6:E:189:TRP:CD1	2.48	0.48
1:G:92:GLU:OE2	1:G:487:LYS:NZ	2.41	0.48
4:H:25:SER:O	4:H:25:SER:OG	2.32	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:271:MET:HA	1:G:348:GLN:NE2	2.28	0.48
2:B:630:GLN:O	2:B:633:LYS:HE3	2.14	0.48
3:L:48:ILE:HA	3:L:54:ARG:HA	1.96	0.48
3:L:118:LEU:HD12	3:L:133:LEU:HD21	1.94	0.48
5:D:163:VAL:HG22	5:D:182:VAL:HG11	1.95	0.48
6:E:23:THR:HB	6:E:71:TRP:O	2.13	0.48
1:G:485:LYS:HB2	1:G:486:TYR:CD1	2.48	0.48
3:L:23:CYS:HB3	3:L:25:ARG:NH1	2.28	0.48
6:E:74:TYR:HD1	6:E:75:LEU:H	1.60	0.48
1:G:35:TRP:CE3	2:B:609:PRO:HA	2.49	0.48
1:G:62:GLU:HA	1:G:65:LYS:HE3	1.96	0.48
1:G:272:ILE:CG1	1:G:286:VAL:HA	2.36	0.48
3:L:144:ALA:O	3:L:198:HIS:HD2	1.97	0.48
4:H:20:LEU:HD21	4:H:105:THR:HG21	1.94	0.48
5:D:12:LYS:HZ1	5:D:16:ALA:N	2.12	0.48
5:D:38:ARG:O	5:D:45:PRO:HA	2.14	0.48
5:D:84:PHE:O	5:D:87:THR:HG22	2.14	0.48
5:D:86:ASP:O	5:D:90:TYR:OH	2.18	0.48
6:E:44:ARG:HG2	6:E:45:ALA:N	2.29	0.48
1:G:86:LEU:HD12	1:G:89:VAL:HG21	1.96	0.47
2:B:596:TRP:CH2	2:B:643:TYR:HD2	2.32	0.47
5:D:128:SER:HB2	5:D:220:LEU:HB2	1.95	0.47
4:H:100(P):MET:HB2	4:H:101:TRP:HZ2	1.79	0.47
4:H:115:LYS:CB	4:H:144:PHE:HE1	2.21	0.47
5:D:51:ILE:HD13	5:D:57:LYS:CE	2.41	0.47
5:D:152:VAL:CG1	5:D:198:VAL:HA	2.36	0.47
1:G:47:ASP:OD1	1:G:487:LYS:NZ	2.46	0.47
1:G:56:SER:O	1:G:76:PRO:HA	2.13	0.47
1:G:163:THR:HG22	1:G:164:GLU:N	2.28	0.47
1:G:225:ILE:HG22	1:G:226:LEU:N	2.28	0.47
5:D:143:LYS:HG3	5:D:145:TYR:OH	2.14	0.47
6:E:23:THR:HG22	6:E:72:SER:HA	1.96	0.47
6:E:37:TRP:NE1	6:E:91:CYS:H	2.13	0.47
6:E:153:LYS:HE2	6:E:159:VAL:N	2.11	0.47
1:G:67:ASN:HD22	1:G:70:ALA:HB3	1.80	0.47
1:G:226:LEU:CD1	1:G:245:VAL:H	2.27	0.47
3:L:117:THR:OG1	3:L:136:LEU:HD23	2.15	0.47
5:D:67:VAL:HG13	5:D:69:MET:HE1	1.96	0.47
6:E:189:TRP:N	6:E:189:TRP:HD1	2.09	0.47
1:G:347:LYS:HA	1:G:347:LYS:HD3	1.65	0.47
1:G:502:LYS:HE2	2:B:607:ASN:HB2	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:82(B):ASN:C	5:D:82(C):LEU:HD12	2.39	0.47
6:E:38:TYR:HD1	6:E:48:LEU:CD2	2.27	0.47
6:E:136:LEU:HD21	6:E:195:TYR:CD2	2.49	0.47
1:G:101:VAL:HG11	1:G:480:ARG:HD3	1.94	0.47
1:G:158:SER:OG	1:G:173:TYR:HB3	2.14	0.47
5:D:8:GLY:O	5:D:107:THR:OG1	2.25	0.47
1:G:37:THR:CG2	2:B:605:CYS:HA	2.44	0.47
1:G:128:THR:C	1:G:129:LEU:HD13	2.38	0.47
1:G:249:HIS:HD1	1:G:249:HIS:H	1.63	0.47
1:G:459:GLY:O	1:G:460:SER:OG	2.29	0.47
4:H:23:THR:HA	4:H:77:GLN:HB3	1.95	0.47
4:H:48:ILE:HG23	4:H:60:ASN:HB3	1.96	0.47
4:H:100(B):TYR:CZ	4:H:100(I):GLU:HA	2.50	0.47
4:H:100(G):PHE:HD1	4:H:100(G):PHE:O	1.97	0.47
4:H:100(P):MET:HB3	4:H:101:TRP:NE1	2.29	0.47
4:H:150:VAL:HG12	4:H:196:VAL:CG1	2.43	0.47
4:H:157:LEU:H	4:H:157:LEU:HD23	1.80	0.47
4:H:166:ALA:HB1	4:H:175:SER:H	1.80	0.47
5:D:6:GLN:NE2	5:D:22:CYS:SG	2.88	0.47
5:D:33:HIS:C	5:D:94:LYS:HZ2	2.23	0.47
5:D:142:VAL:CG1	5:D:179:SER:HA	2.45	0.47
6:E:63:ARG:O	6:E:78:SER:N	2.48	0.47
6:E:101:PHE:HD1	6:E:101:PHE:N	2.12	0.47
6:E:121:LEU:HA	6:E:138:CYS:HB2	1.96	0.47
1:G:349:LEU:O	1:G:352:HIS:ND1	2.47	0.47
1:G:502:LYS:HE2	2:B:607:ASN:CB	2.45	0.47
2:B:580:VAL:O	2:B:584:GLU:HG3	2.15	0.47
4:H:94:ARG:O	4:H:100(P):MET:HA	2.14	0.47
4:H:114:THR:HG21	4:H:200:PRO:O	2.12	0.47
2:B:533:ALA:HB3	2:B:628:TRP:NE1	2.29	0.47
4:H:89:ILE:HD12	4:H:105:THR:O	2.14	0.47
4:H:100(J):PHE:N	4:H:100(J):PHE:HD1	2.13	0.47
6:E:121:LEU:HG	6:E:138:CYS:CB	2.40	0.47
6:E:144:TYR:HA	6:E:145:PRO:C	2.40	0.47
1:G:182:VAL:C	1:G:191:TYR:HB3	2.39	0.47
1:G:304:ARG:HH12	1:G:440:GLN:N	2.13	0.47
6:E:96:ASN:HB2	8:C:6:MAN:C2	2.45	0.47
6:E:135:THR:HG21	6:E:181:TYR:CD1	2.50	0.47
1:G:493:PRO:C	1:G:494:LEU:HD23	2.40	0.46
4:H:17:THR:HA	4:H:82:LEU:O	2.15	0.46
5:D:179:SER:HB3	6:E:139:LEU:HD22	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:125:SER:O	6:E:128:GLU:HB3	2.16	0.46
1:G:287:GLN:HA	1:G:451:GLY:HA2	1.97	0.46
1:G:393:SER:HB2	1:G:395:TRP:HE1	1.80	0.46
4:H:46:GLU:HG2	4:H:48:ILE:HD11	1.96	0.46
5:D:37:ILE:HD11	5:D:103:TRP:CH2	2.50	0.46
5:D:84:PHE:CE2	5:D:111:VAL:HG12	2.50	0.46
5:D:147:PRO:O	5:D:200:HIS:NE2	2.47	0.46
6:E:4:LEU:HD11	6:E:102:GLY:CA	2.31	0.46
1:G:35:TRP:CD1	1:G:502:LYS:HB2	2.50	0.46
1:G:178:ARG:NH1	1:G:419:ARG:HH22	2.14	0.46
1:G:377:ASN:ND2	1:G:382:PHE:HE1	2.13	0.46
6:E:56:ARG:HG3	6:E:60:ILE:HD11	1.97	0.46
1:G:192:ARG:HH21	1:G:197:ASN:HB2	1.81	0.46
3:L:163:THR:HG22	3:L:176:SER:H	1.80	0.46
4:H:33:TYR:H	4:H:94:ARG:HH22	1.62	0.46
5:D:23:LYS:HE2	5:D:23:LYS:CA	2.46	0.46
6:E:101:PHE:N	6:E:101:PHE:CD1	2.84	0.46
2:B:596:TRP:NE1	2:B:647:GLU:CD	2.67	0.46
3:L:46:LEU:HD12	3:L:46:LEU:HA	1.41	0.46
4:H:94:ARG:HG2	4:H:100(R):VAL:H	1.80	0.46
5:D:36:TRP:NE1	5:D:80:MET:HG2	2.31	0.46
6:E:37:TRP:HB3	6:E:89:TYR:O	2.16	0.46
1:G:128:THR:C	1:G:129:LEU:HD12	2.40	0.46
2:B:645:LEU:HD23	2:B:648:GLU:CG	2.46	0.46
5:D:31:PHE:CE1	8:C:2:NAG:H82	2.51	0.46
4:H:45:LEU:HD23	4:H:46:GLU:N	2.31	0.46
4:H:64:LYS:HE2	4:H:64:LYS:HB3	1.55	0.46
5:D:33:HIS:HE1	5:D:98:ARG:NE	2.13	0.46
5:D:100(B):SER:HB3	6:E:52:GLU:OE1	2.14	0.46
6:E:184:LEU:HD23	6:E:189:TRP:CZ3	2.50	0.46
1:G:480:ARG:HG2	1:G:480:ARG:HH11	1.81	0.46
3:L:167:LYS:HD3	3:L:173:TYR:CE1	2.51	0.46
4:H:13:LYS:O	4:H:16:GLU:CD	2.58	0.46
5:D:35:ASN:CB	5:D:94:LYS:HE3	2.40	0.46
6:E:190:LYS:HZ1	6:E:213:THR:N	2.12	0.46
1:G:62:GLU:HA	1:G:65:LYS:CE	2.45	0.46
1:G:323:ILE:O	1:G:323:ILE:HG13	2.15	0.46
4:H:29:ILE:HA	4:H:34:TRP:HH2	1.81	0.46
5:D:72(C):VAL:O	5:D:75:THR:N	2.48	0.46
5:D:96:LEU:HG	5:D:100(E):LEU:HB3	1.97	0.46
6:E:129:LEU:CD2	6:E:134:ALA:HB2	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:85:HIS:ND1	1:G:242:VAL:O	2.48	0.46
1:G:478:ASN:O	1:G:478:ASN:ND2	2.49	0.46
2:B:545:LEU:HD13	2:B:545:LEU:O	2.16	0.46
3:L:37:GLN:NE2	3:L:86:TYR:CE1	2.76	0.46
4:H:24:VAL:HG22	4:H:26:GLY:H	1.80	0.46
4:H:125:SER:OG	4:H:128:SER:N	2.49	0.46
6:E:99:CYS:SG	6:E:100:VAL:N	2.88	0.46
1:G:68:VAL:HG13	1:G:69:TRP:N	2.31	0.45
3:L:137:ILE:HD13	3:L:140:PHE:CE1	2.51	0.45
5:D:47:TRP:CZ2	5:D:49:GLY:HA2	2.51	0.45
5:D:68:ILE:O	5:D:69:MET:HG3	2.15	0.45
5:D:167:PRO:HG3	6:E:169:SER:CB	2.46	0.45
6:E:95:HIS:CE1	6:E:96:ASN:HB3	2.51	0.45
6:E:127:GLU:H	6:E:127:GLU:CD	2.24	0.45
6:E:192:HIS:HD2	6:E:193:ARG:C	2.24	0.45
1:G:295:ASN:HD21	9:F:1:NAG:C7	2.28	0.45
1:G:301:ASN:HB3	1:G:441:GLY:HA2	1.98	0.45
3:L:16:GLY:H	3:L:78:VAL:HG23	1.81	0.45
3:L:61:ARG:NH2	3:L:61:ARG:HG2	2.31	0.45
3:L:195:GLN:OE1	3:L:204:GLU:HB3	2.16	0.45
4:H:38:ARG:HE	4:H:38:ARG:HB2	1.59	0.45
5:D:30:ASN:ND2	5:D:53:TYR:HB2	2.31	0.45
5:D:155:ASN:HD21	5:D:159:LEU:H	1.64	0.45
1:G:335:LYS:HG2	1:G:412:ASP:O	2.15	0.45
3:L:75:ILE:HD13	3:L:78:VAL:CG1	2.46	0.45
3:L:84:ALA:HB3	3:L:86:TYR:CE1	2.51	0.45
4:H:52:SER:OG	4:H:56:SER:OG	2.33	0.45
5:D:11:LEU:CD1	5:D:110:THR:HG22	2.47	0.45
5:D:72(F):THR:HG22	5:D:72(G):SER:H	1.81	0.45
1:G:42:VAL:HG22	1:G:493:PRO:O	2.16	0.45
1:G:95:MET:HA	1:G:98:ASN:HB2	1.99	0.45
4:H:82(C):VAL:HG13	4:H:86:ASP:HB2	1.97	0.45
1:G:272:ILE:HG21	1:G:286:VAL:HA	1.98	0.45
1:G:300:ASN:OD1	1:G:327:ARG:HB2	2.17	0.45
3:L:16:GLY:N	3:L:78:VAL:HG23	2.31	0.45
5:D:168:ALA:CB	5:D:176:TYR:HB3	2.46	0.45
6:E:144:TYR:CD1	6:E:145:PRO:HA	2.51	0.45
1:G:221:ALA:HB1	2:B:544:LEU:O	2.16	0.45
1:G:383:PHE:CD1	1:G:383:PHE:N	2.85	0.45
4:H:20:LEU:HD13	4:H:21:THR:N	2.31	0.45
6:E:68:LYS:CB	6:E:73:ALA:HA	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:E:152:TRP:O	6:E:152:TRP:CD1	2.69	0.45
6:E:170:LYS:HD3	6:E:170:LYS:N	2.31	0.45
6:E:196:SER:HB3	6:E:207:GLU:OE1	2.16	0.45
1:G:111:LEU:HD23	1:G:111:LEU:H	1.80	0.45
1:G:269:GLU:CB	1:G:271:MET:HE1	2.45	0.45
4:H:187:LEU:HA	4:H:188:GLY:HA2	1.67	0.45
6:E:35:ILE:HA	6:E:91:CYS:O	2.16	0.45
1:G:101:VAL:HG21	1:G:480:ARG:NE	2.31	0.45
1:G:181:VAL:CG1	1:G:193:LEU:HG	2.46	0.45
1:G:384:TYR:HE1	1:G:421:LYS:HB2	1.81	0.45
2:B:539:VAL:HG13	2:B:540:GLN:OE1	2.16	0.45
3:L:83:GLU:HA	3:L:104:LEU:HD11	1.99	0.45
5:D:140:CYS:HB2	5:D:154:TRP:CZ2	2.52	0.45
1:G:304:ARG:HD2	1:G:440:GLN:HB3	1.99	0.45
2:B:650:GLN:O	2:B:654:GLU:HB2	2.16	0.45
4:H:107:VAL:HG23	4:H:107:VAL:O	2.16	0.45
6:E:131:ALA:H	6:E:132:ASN:HA	1.81	0.45
1:G:37:THR:HG22	2:B:606:THR:CB	2.46	0.45
1:G:123:THR:HG23	1:G:124:PRO:HD3	1.98	0.45
1:G:128:THR:N	1:G:129:LEU:HD13	2.32	0.45
3:L:37:GLN:O	3:L:44:PRO:HA	2.17	0.45
5:D:116:THR:C	5:D:117:LYS:HD2	2.42	0.45
1:G:234:ASN:ND2	9:K:1:NAG:O5	2.23	0.44
2:B:519:PHE:HA	2:B:520:LEU:HA	1.52	0.44
2:B:639:THR:O	2:B:642:ILE:HG22	2.17	0.44
3:L:150:LYS:NZ	3:L:153:SER:N	2.65	0.44
5:D:24:THR:HG23	5:D:27:TYR:CZ	2.52	0.44
5:D:169:VAL:HG22	5:D:170:LEU:N	2.29	0.44
1:G:35:TRP:O	1:G:498:PRO:HA	2.17	0.44
1:G:104:MET:HE3	1:G:104:MET:HB2	1.78	0.44
1:G:128:THR:O	1:G:129:LEU:HD12	2.17	0.44
1:G:298:ARG:H	1:G:443:ILE:HD11	1.81	0.44
2:B:639:THR:CG2	2:B:643:TYR:HE1	2.24	0.44
3:L:59:PRO:HG2	3:L:62:PHE:CE1	2.49	0.44
3:L:109:GLN:O	3:L:109:GLN:HG2	2.16	0.44
1:G:360:ARG:O	1:G:361:PHE:HB2	2.17	0.44
3:L:135:CYS:O	3:L:136:LEU:HD13	2.16	0.44
4:H:87:SER:O	4:H:88:ALA:HB2	2.18	0.44
4:H:146:GLU:N	4:H:147:PRO:HD2	2.32	0.44
1:G:76:PRO:HG3	2:B:556:LEU:HB3	1.99	0.44
1:G:238:PRO:HG2	5:D:72(F):THR:HA	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:304:ARG:HH12	1:G:439:ILE:C	2.26	0.44
2:B:656:ASN:OD1	2:B:656:ASN:C	2.59	0.44
5:D:98:ARG:NE	5:D:98:ARG:CA	2.81	0.44
6:E:37:TRP:CG	6:E:90:CYS:HA	2.53	0.44
6:E:165:THR:HB	6:E:180:SER:CB	2.47	0.44
1:G:361:PHE:HD1	1:G:362:ALA:H	1.66	0.44
3:L:96:TRP:HH2	4:H:49:GLY:HA2	1.83	0.44
4:H:23:THR:HA	4:H:77:GLN:CB	2.47	0.44
4:H:51:ILE:HG13	4:H:52:SER:N	2.32	0.44
5:D:72(F):THR:HG22	5:D:72(G):SER:N	2.32	0.44
6:E:99:CYS:SG	6:E:101:PHE:HE1	2.41	0.44
6:E:192:HIS:HE1	6:E:195:TYR:CE2	2.36	0.44
1:G:105:HIS:HB3	1:G:479:TRP:NE1	2.32	0.44
1:G:261:LEU:CD2	1:G:448:ASN:H	2.31	0.44
4:H:66:ARG:NH1	4:H:82:LEU:HD11	2.31	0.44
4:H:99:LYS:O	4:H:100:ARG:HB3	2.18	0.44
4:H:149:THR:CB	4:H:197:ASN:OD1	2.66	0.44
5:D:154:TRP:CD2	5:D:196:CYS:HB3	2.52	0.44
1:G:283:ASN:CG	1:G:455:THR:HG22	2.43	0.44
1:G:304:ARG:HH22	1:G:438:PRO:C	2.22	0.44
1:G:356:ASN:OD1	16:G:1663:NAG:H2	2.18	0.44
3:L:50:ASN:OD1	3:L:51:ASN:N	2.51	0.44
4:H:5:GLN:NE2	4:H:103:LYS:HE3	2.33	0.44
5:D:121:VAL:HG23	5:D:141:LEU:CD1	2.40	0.44
6:E:23:THR:HG22	6:E:72:SER:CA	2.47	0.44
6:E:37:TRP:CE3	6:E:75:LEU:HD23	2.53	0.44
1:G:48:ALA:HB2	1:G:490:LYS:N	2.33	0.44
1:G:327:ARG:HE	4:H:100(B):TYR:HD2	1.65	0.44
3:L:10:PRO:O	3:L:11:LEU:HD13	2.17	0.44
3:L:46:LEU:HD12	3:L:47:LEU:N	2.32	0.44
5:D:119:PRO:HG2	5:D:205:THR:HB	2.00	0.44
3:L:194:CYS:O	3:L:194:CYS:SG	2.76	0.43
4:H:67:VAL:O	4:H:68:ILE:HG13	2.18	0.43
6:E:85:GLU:OE1	6:E:109:VAL:HG22	2.18	0.43
7:A:2:NAG:O7	7:A:2:NAG:O3	2.34	0.43
4:H:131:GLY:HA2	4:H:132:GLY:C	2.43	0.43
5:D:38:ARG:O	5:D:46:GLU:N	2.50	0.43
6:E:185:THR:HB	6:E:188:GLN:HG3	2.00	0.43
1:G:270:VAL:C	1:G:272:ILE:CD1	2.88	0.43
1:G:359:ILE:HG12	1:G:396:ILE:HD11	2.00	0.43
1:G:387:THR:HA	1:G:416:LEU:HD21	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:32:TYR:OH	4:H:96:GLN:HA	2.18	0.43
4:H:57:THR:OG1	4:H:58:ASN:N	2.51	0.43
5:D:11:LEU:HA	5:D:110:THR:O	2.18	0.43
6:E:20:ILE:HD12	6:E:21:SER:H	1.83	0.43
6:E:64:PHE:HD2	6:E:77:ILE:HG22	1.83	0.43
6:E:189:TRP:O	6:E:192:HIS:CE1	2.72	0.43
9:F:5:MAN:H61	9:F:6:MAN:O6	2.19	0.43
1:G:61:TYR:O	1:G:61:TYR:CG	2.72	0.43
1:G:178:ARG:HB3	1:G:179:LEU:HD22	2.00	0.43
1:G:373:THR:HG21	1:G:384:TYR:HB3	2.00	0.43
3:L:43:ALA:HA	4:H:91:TYR:HE2	1.81	0.43
3:L:150:LYS:HZ2	3:L:154:SER:H	1.62	0.43
4:H:100(G):PHE:HB3	4:H:100(I):GLU:CD	2.43	0.43
4:H:141:LYS:HG3	4:H:142:ASP:HB2	1.99	0.43
4:H:173:LEU:HD12	4:H:173:LEU:HA	1.58	0.43
6:E:134:ALA:CB	6:E:184:LEU:HG	2.48	0.43
9:F:3:BMA:C4	9:F:4:MAN:H2	2.46	0.43
5:D:68:ILE:N	5:D:69:MET:HE2	2.32	0.43
5:D:198:VAL:O	5:D:206:LYS:HA	2.19	0.43
6:E:153:LYS:HG3	6:E:156:SER:HA	1.99	0.43
1:G:39:TYR:HD1	1:G:495:GLY:O	2.01	0.43
1:G:64:GLU:OE2	11:M:4:MAN:O4	2.30	0.43
1:G:102:GLU:HA	1:G:105:HIS:CD2	2.53	0.43
1:G:201:ILE:HD12	1:G:201:ILE:N	2.34	0.43
1:G:270:VAL:HG22	1:G:288:PHE:C	2.44	0.43
1:G:360:ARG:NE	1:G:468:PHE:HD1	2.16	0.43
2:B:586:TYR:CE1	2:B:587:LEU:HD23	2.53	0.43
4:H:6:GLU:HA	4:H:22:CYS:HA	2.00	0.43
1:G:55:ALA:O	1:G:216:HIS:HB2	2.19	0.43
1:G:136:ASN:OD1	1:G:137:ASN:N	2.51	0.43
1:G:373:THR:HG1	1:G:385:CYS:C	2.23	0.43
2:B:645:LEU:HD23	2:B:648:GLU:CB	2.48	0.43
4:H:16:GLU:N	4:H:82(B):SER:HA	2.34	0.43
6:E:143:PHE:CD2	6:E:146:GLY:HA2	2.52	0.43
1:G:108:ILE:HD12	1:G:479:TRP:CH2	2.54	0.43
1:G:469:ARG:HH11	1:G:469:ARG:HD2	1.68	0.43
2:B:573:ILE:O	2:B:576:LEU:HB3	2.18	0.43
3:L:131:ALA:O	3:L:181:LEU:HB3	2.18	0.43
4:H:81:LYS:HE3	4:H:81:LYS:HB3	1.77	0.43
4:H:149:THR:CG2	4:H:197:ASN:CB	2.97	0.43
5:D:72(B):GLU:OE1	5:D:72(B):GLU:N	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:98:ARG:NH1	5:D:100(C):THR:CA	2.79	0.43
1:G:152:GLY:CA	1:G:178:ARG:HE	2.29	0.43
2:B:546:SER:HB2	2:B:549:VAL:HG12	2.00	0.43
2:B:622:ILE:HG13	2:B:623:TRP:N	2.34	0.43
3:L:96:TRP:HZ3	4:H:48:ILE:C	2.27	0.43
3:L:152:ASP:OD1	3:L:152:ASP:N	2.51	0.43
4:H:39:GLN:OE1	4:H:91:TYR:CZ	2.71	0.43
6:E:192:HIS:CE1	6:E:195:TYR:CE2	3.07	0.43
1:G:298:ARG:HG3	1:G:443:ILE:HD11	2.01	0.43
3:L:111:LYS:HB3	3:L:141:TYR:O	2.18	0.43
4:H:32:TYR:C	4:H:33:TYR:CG	2.96	0.43
4:H:50:TYR:O	4:H:58:ASN:N	2.52	0.43
5:D:33:HIS:O	5:D:94:LYS:HG3	2.19	0.43
1:G:330:HIS:HA	1:G:416:LEU:O	2.19	0.42
2:B:633:LYS:HD2	2:B:634:GLU:N	2.34	0.42
3:L:36:TYR:HE2	3:L:89:HIS:HB2	1.84	0.42
3:L:50:ASN:HB3	3:L:52:GLN:HG3	2.01	0.42
3:L:137:ILE:HD12	3:L:137:ILE:O	2.19	0.42
6:E:6:GLN:NE2	6:E:103:THR:N	2.66	0.42
6:E:84:ASP:O	6:E:107:VAL:HG11	2.19	0.42
1:G:172:VAL:HG22	1:G:173:TYR:N	2.35	0.42
3:L:120:PRO:O	4:H:122:LEU:HG	2.19	0.42
5:D:52:SER:OG	8:C:2:NAG:O7	2.32	0.42
5:D:179:SER:HB3	6:E:139:LEU:HD21	2.00	0.42
1:G:112:TRP:CZ2	1:G:426:MET:HE1	2.54	0.42
1:G:469:ARG:HB2	1:G:469:ARG:NH1	2.34	0.42
2:B:618:ASN:OD1	16:B:702:NAG:C7	2.67	0.42
2:B:643:TYR:HA	2:B:646:LEU:HD13	2.02	0.42
6:E:16:GLN:NE2	6:E:18:VAL:H	2.17	0.42
6:E:17:SER:HA	6:E:77:ILE:O	2.19	0.42
6:E:192:HIS:CD2	6:E:193:ARG:N	2.87	0.42
1:G:32:GLU:N	1:G:500:ARG:HH21	2.17	0.42
3:L:50:ASN:C	3:L:52:GLN:N	2.77	0.42
3:L:119:PHE:HA	3:L:120:PRO:HD3	1.92	0.42
4:H:29:ILE:HG23	4:H:34:TRP:CZ3	2.54	0.42
4:H:136:LEU:HA	4:H:136:LEU:HD22	1.67	0.42
6:E:142:ASP:CG	6:E:175:LYS:HD3	2.45	0.42
1:G:40:TYR:O	2:B:537:LEU:HB3	2.19	0.42
1:G:162:THR:OG1	1:G:166:ARG:NH2	2.31	0.42
1:G:190:GLU:CD	1:G:191:TYR:H	2.27	0.42
1:G:215:ILE:N	1:G:251:ILE:O	2.47	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:222:GLY:C	1:G:491:ILE:HD12	2.45	0.42
1:G:224:ALA:O	1:G:489:VAL:N	2.52	0.42
1:G:263:GLY:H	1:G:450:THR:HG21	1.84	0.42
1:G:373:THR:OG1	1:G:385:CYS:C	2.62	0.42
1:G:395:TRP:C	1:G:396:ILE:HD13	2.45	0.42
3:L:166:SER:O	3:L:174:ALA:N	2.44	0.42
4:H:193:ILE:HD13	4:H:208:LYS:HZ2	1.84	0.42
5:D:67:VAL:HG23	5:D:82:ILE:HA	2.02	0.42
6:E:68:LYS:HG2	6:E:73:ALA:N	2.34	0.42
6:E:81:ARG:HB2	6:E:83:GLU:OE2	2.20	0.42
6:E:81:ARG:O	6:E:83:GLU:N	2.52	0.42
6:E:188:GLN:O	6:E:191:SER:HB2	2.18	0.42
1:G:155:LYS:NZ	1:G:191:TYR:HE2	2.17	0.42
1:G:478:ASN:ND2	1:G:478:ASN:C	2.78	0.42
2:B:608:VAL:H	2:B:650:GLN:HE22	1.62	0.42
3:L:25:ARG:N	11:J:5:MAN:O6	2.53	0.42
3:L:150:LYS:NZ	3:L:153:SER:H	2.17	0.42
4:H:121:PRO:HD3	4:H:207:LYS:HE3	2.02	0.42
5:D:100(C):THR:HG23	5:D:100(C):THR:O	2.19	0.42
5:D:181:VAL:HG21	6:E:122:PHE:CE2	2.54	0.42
6:E:153:LYS:HD3	6:E:154:ALA:N	2.34	0.42
1:G:297:THR:O	1:G:329:ALA:HA	2.19	0.42
1:G:382:PHE:O	1:G:420:ILE:HA	2.20	0.42
3:L:172:LYS:HD3	3:L:172:LYS:N	2.33	0.42
5:D:148:GLU:HB2	5:D:149:PRO:HA	2.02	0.42
5:D:153:SER:O	5:D:154:TRP:HD1	2.03	0.42
6:E:38:TYR:OH	6:E:91:CYS:SG	2.76	0.42
6:E:70:TYR:HD1	6:E:70:TYR:H	1.67	0.42
6:E:152:TRP:HE3	6:E:163:VAL:HG23	1.84	0.42
1:G:298:ARG:NH1	1:G:300:ASN:O	2.53	0.42
1:G:327:ARG:NH1	4:H:100(B):TYR:HD2	2.12	0.42
2:B:544:LEU:HD13	2:B:544:LEU:O	2.19	0.42
3:L:25:ARG:CZ	3:L:88:CYS:SG	3.08	0.42
3:L:95(B):PHE:CE1	4:H:47:TRP:CH2	3.06	0.42
4:H:163:THR:HA	4:H:178:SER:HA	2.02	0.42
5:D:100(D):TRP:CD1	5:D:100(E):LEU:N	2.87	0.42
1:G:328:GLN:H	4:H:100(G):PHE:HZ	1.66	0.42
1:G:385:CYS:HA	1:G:418:CYS:HA	2.02	0.42
1:G:455:THR:CG2	1:G:471:GLY:HA3	2.50	0.42
4:H:33:TYR:CE1	4:H:52:SER:HA	2.54	0.42
4:H:100(D):MET:HB2	13:O:2:NAG:O7	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:193:ILE:HD13	4:H:208:LYS:NZ	2.35	0.42
5:D:52:SER:HB2	5:D:56:ASP:OD1	2.20	0.42
5:D:110:THR:HG21	5:D:147:PRO:HB3	2.00	0.42
1:G:411:ASN:OD1	1:G:412:ASP:N	2.53	0.42
2:B:583:VAL:O	2:B:587:LEU:N	2.37	0.42
3:L:104:LEU:HD12	3:L:104:LEU:C	2.44	0.42
4:H:6:GLU:HG2	4:H:92:CYS:SG	2.59	0.42
4:H:87:SER:HA	4:H:108:THR:HA	2.02	0.42
4:H:100(D):MET:CB	13:O:2:NAG:H2	2.48	0.42
5:D:10:GLU:HG2	5:D:202:PRO:CG	2.50	0.42
6:E:89:TYR:OH	6:E:104:GLY:HA2	2.19	0.42
1:G:332:ASN:HA	1:G:414:ILE:O	2.20	0.41
4:H:59:TYR:CD2	4:H:64:LYS:HB2	2.52	0.41
4:H:166:ALA:HB1	4:H:175:SER:N	2.35	0.41
5:D:210:ARG:NH1	5:D:211:VAL:O	2.53	0.41
6:E:41:PRO:HG2	6:E:44:ARG:HD3	2.01	0.41
6:E:180:SER:O	6:E:181:TYR:HB3	2.20	0.41
1:G:226:LEU:N	1:G:487:LYS:O	2.52	0.41
1:G:497:ALA:HA	1:G:498:PRO:HD3	1.89	0.41
5:D:33:HIS:CE1	5:D:98:ARG:CD	3.03	0.41
5:D:153:SER:O	5:D:154:TRP:CD1	2.72	0.41
6:E:99:CYS:SG	6:E:101:PHE:CE1	3.13	0.41
1:G:179:LEU:HD21	1:G:419:ARG:NH2	2.35	0.41
1:G:182:VAL:O	1:G:192:ARG:N	2.44	0.41
1:G:391:PHE:N	1:G:391:PHE:CD1	2.85	0.41
3:L:49:TYR:HB2	4:H:100(O):TYR:CG	2.56	0.41
4:H:59:TYR:CB	4:H:64:LYS:HB3	2.46	0.41
4:H:100(D):MET:O	4:H:100(D):MET:HG2	2.20	0.41
5:D:166:PHE:O	5:D:178:LEU:HD13	2.20	0.41
6:E:87:THR:HG22	6:E:105:THR:O	2.20	0.41
1:G:162:THR:HG1	1:G:166:ARG:HH22	1.59	0.41
1:G:502:LYS:HD2	2:B:607:ASN:HA	2.02	0.41
4:H:40:SER:OG	4:H:43:LYS:HB2	2.21	0.41
4:H:51:ILE:HD11	4:H:55:GLU:HA	2.02	0.41
5:D:72(E):VAL:O	5:D:72(F):THR:OG1	2.33	0.41
6:E:81:ARG:HA	6:E:81:ARG:HD3	1.84	0.41
1:G:166:ARG:HH21	1:G:169:LYS:HZ2	1.69	0.41
2:B:576:LEU:HA	2:B:579:ARG:NH2	2.35	0.41
2:B:595:ILE:HG21	2:B:647:GLU:OE2	2.20	0.41
2:B:619:LEU:O	2:B:622:ILE:HG12	2.20	0.41
2:B:654:GLU:C	2:B:657:GLU:OE1	2.64	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:47:LEU:H	3:L:47:LEU:HD23	1.86	0.41
3:L:52:GLN:OE1	3:L:53:ASP:HB3	2.20	0.41
3:L:96:TRP:CH2	4:H:49:GLY:HA2	2.56	0.41
3:L:106:VAL:O	3:L:109:GLN:NE2	2.52	0.41
5:D:66:ARG:HD2	5:D:66:ARG:HA	1.38	0.41
5:D:67:VAL:C	5:D:69:MET:HE2	2.45	0.41
6:E:138:CYS:N	6:E:139:LEU:HD12	2.36	0.41
6:E:187:GLU:O	6:E:189:TRP:N	2.54	0.41
1:G:104:MET:SD	1:G:479:TRP:HB3	2.60	0.41
1:G:180:ASP:OD1	1:G:180:ASP:N	2.52	0.41
3:L:35:TRP:HB3	3:L:47:LEU:CD2	2.45	0.41
4:H:146:GLU:OE1	4:H:165:PRO:HB3	2.20	0.41
4:H:190:GLN:HG2	4:H:192:TYR:CZ	2.56	0.41
5:D:82:ILE:O	5:D:82(A):ARG:HB2	2.21	0.41
5:D:121:VAL:CG2	5:D:141:LEU:HB3	2.51	0.41
6:E:42:PRO:HD2	6:E:86:THR:HG21	2.00	0.41
1:G:423:ILE:HA	1:G:434:MET:O	2.20	0.41
1:G:460:SER:HB2	1:G:466:GLU:OE1	2.20	0.41
1:G:503:ARG:HD3	1:G:503:ARG:HA	1.88	0.41
2:B:580:VAL:O	2:B:583:VAL:HG22	2.20	0.41
5:D:72(C):VAL:HB	5:D:75:THR:HB	2.01	0.41
5:D:96:LEU:HD13	5:D:98:ARG:NH2	2.34	0.41
6:E:23:THR:HG21	6:E:72:SER:HB2	2.03	0.41
6:E:192:HIS:CD2	6:E:194:SER:N	2.88	0.41
1:G:117:LYS:N	1:G:118:PRO:HD2	2.36	0.41
1:G:178:ARG:CZ	1:G:419:ARG:HH22	2.34	0.41
1:G:270:VAL:HG13	1:G:287:GLN:C	2.46	0.41
1:G:271:MET:HG3	1:G:348:GLN:OE1	2.20	0.41
3:L:50:ASN:OD1	13:O:5:MAN:O4	2.37	0.41
3:L:133:LEU:HG	3:L:134:VAL:N	2.36	0.41
3:L:147:VAL:HA	3:L:196:VAL:HA	2.03	0.41
4:H:7:SER:O	4:H:20:LEU:HD22	2.20	0.41
5:D:37:ILE:HD11	5:D:103:TRP:CZ3	2.56	0.41
5:D:197:ASN:HB2	5:D:208:ASP:CG	2.46	0.41
1:G:269:GLU:C	1:G:271:MET:SD	3.04	0.41
1:G:272:ILE:HG23	1:G:285:LEU:O	2.21	0.41
1:G:293:GLN:OE1	9:F:1:NAG:H3	2.20	0.41
1:G:363:ASN:HD22	1:G:363:ASN:N	2.19	0.41
3:L:120:PRO:HG3	3:L:207:VAL:HG21	2.02	0.41
3:L:167:LYS:HD3	3:L:173:TYR:CD1	2.56	0.41
4:H:94:ARG:C	4:H:94:ARG:HE	2.29	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:100(A):ILE:HA	4:H:100(I):GLU:O	2.20	0.41
4:H:146:GLU:CD	4:H:166:ALA:H	2.29	0.41
5:D:50:TRP:CH2	6:E:98:GLY:HA2	2.56	0.41
5:D:72:ASP:OD1	5:D:72:ASP:N	2.53	0.41
6:E:49:ILE:HD13	6:E:49:ILE:HG21	1.86	0.41
6:E:87:THR:HB	6:E:89:TYR:HE1	1.86	0.41
6:E:152:TRP:CD1	6:E:152:TRP:C	2.99	0.41
6:E:190:LYS:HA	6:E:190:LYS:HD2	1.25	0.41
1:G:171:LYS:CG	7:A:1:NAG:O7	2.65	0.41
1:G:199:ALA:HB1	1:G:431:GLY:O	2.21	0.41
1:G:298:ARG:N	1:G:443:ILE:HD11	2.36	0.41
1:G:343:GLY:O	1:G:346:VAL:HB	2.21	0.41
3:L:37:GLN:NE2	3:L:86:TYR:HE1	2.18	0.41
3:L:199:GLU:HG2	3:L:201:SER:H	1.86	0.41
4:H:34:TRP:N	4:H:50:TYR:CZ	2.89	0.41
5:D:129:LYS:NZ	5:D:183:THR:HG21	2.36	0.41
6:E:190:LYS:NZ	6:E:212:PRO:HB2	2.36	0.41
1:G:195:ASN:OD1	1:G:195:ASN:C	2.63	0.40
1:G:271:MET:CG	1:G:348:GLN:OE1	2.68	0.40
3:L:76:SER:OG	3:L:77:GLY:N	2.54	0.40
3:L:90:MET:C	3:L:91:TRP:CD1	2.99	0.40
6:E:192:HIS:CD2	6:E:192:HIS:C	2.99	0.40
1:G:45:TRP:HB2	1:G:489:VAL:HG22	2.04	0.40
1:G:421:LYS:HD2	1:G:421:LYS:HA	1.90	0.40
3:L:35:TRP:CE3	3:L:35:TRP:HA	2.56	0.40
3:L:103:ARG:CZ	3:L:105:THR:HG22	2.51	0.40
3:L:189:HIS:ND1	3:L:190:LYS:N	2.67	0.40
4:H:100(E):VAL:HG23	4:H:100(F):SER:N	2.37	0.40
5:D:100(B):SER:HA	6:E:93:TYR:CE1	2.57	0.40
1:G:166:ARG:HH21	1:G:169:LYS:NZ	2.20	0.40
2:B:529:THR:HB	2:B:623:TRP:O	2.21	0.40
4:H:13:LYS:HB2	4:H:13:LYS:HE3	1.89	0.40
4:H:38:ARG:CZ	4:H:90:TYR:CZ	3.05	0.40
4:H:50:TYR:HD1	4:H:50:TYR:HA	1.63	0.40
4:H:115:LYS:HG2	4:H:116:GLY:H	1.87	0.40
4:H:157:LEU:HD23	4:H:157:LEU:N	2.36	0.40
5:D:37:ILE:HG22	5:D:47:TRP:HA	2.04	0.40
1:G:127:VAL:HG23	1:G:129:LEU:CD1	2.52	0.40
1:G:326:ILE:O	1:G:327:ARG:HG2	2.21	0.40
1:G:360:ARG:HE	1:G:467:THR:CA	2.34	0.40
3:L:145:VAL:CG2	3:L:196:VAL:HB	2.51	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:D:94:LYS:HA	5:D:94:LYS:HE2	2.04	0.40
1:G:113:ASP:N	1:G:113:ASP:OD1	2.55	0.40
1:G:274:SER:OG	1:G:277:ILE:N	2.53	0.40
1:G:341:THR:O	1:G:345:VAL:HG23	2.21	0.40
1:G:362:ALA:CA	1:G:469:ARG:HH21	2.35	0.40
3:L:125:GLU:CD	3:L:131:ALA:HA	2.46	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	G	444/480 (92%)	378 (85%)	62 (14%)	4 (1%)	14	41
2	B	144/153 (94%)	127 (88%)	16 (11%)	1 (1%)	18	48
3	L	209/218 (96%)	171 (82%)	34 (16%)	4 (2%)	6	23
4	H	227/236 (96%)	184 (81%)	35 (15%)	8 (4%)	3	12
5	D	238/240 (99%)	190 (80%)	36 (15%)	12 (5%)	1	6
6	E	211/216 (98%)	177 (84%)	29 (14%)	5 (2%)	4	18
All	All	1473/1543 (96%)	1227 (83%)	212 (14%)	34 (2%)	5	19

All (34) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	G	88	ASN
4	H	100	ARG
4	H	100(E)	VAL
4	H	100(G)	PHE
5	D	72(B)	GLU
5	D	100(E)	LEU

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Mol	Chain	Res	Type
6	E	95	HIS
1	G	427	TRP
4	H	100(A)	ILE
5	D	51	ILE
5	D	92	CYS
5	D	100(D)	TRP
5	D	142	VAL
5	D	204	ASN
2	B	546	SER
3	L	26	GLN
4	H	100(Q)	ASP
5	D	72	ASP
5	D	72(G)	SER
4	H	100(I)	GLU
5	D	121	VAL
5	D	155	ASN
6	E	145	PRO
6	E	158	PRO
3	L	203	VAL
4	H	82(B)	SER
5	D	146	PHE
6	E	186	PRO
6	E	82	PRO
3	L	55	PRO
4	H	147	PRO
1	G	270	VAL
1	G	299	PRO
3	L	10	PRO

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	G	397/424 (94%)	392 (99%)	5 (1%)	61	86
2	B	126/129 (98%)	123 (98%)	3 (2%)	43	75
3	L	175/181 (97%)	172 (98%)	3 (2%)	53	82

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
4	H	200/205 (98%)	198 (99%)	2 (1%)	68	89
5	D	203/203 (100%)	193 (95%)	10 (5%)	22	54
6	E	186/189 (98%)	179 (96%)	7 (4%)	29	64
All	All	1287/1331 (97%)	1257 (98%)	30 (2%)	44	76

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	65	LYS
1	G	259	LEU
1	G	361	PHE
1	G	392	ASN
1	G	428	GLN
2	B	545	LEU
2	B	615	SER
2	B	639	THR
3	L	190	LYS
3	L	194	CYS
3	L	196	VAL
4	H	94	ARG
4	H	106	THR
5	D	72	ASP
5	D	72(A)	THR
5	D	72(B)	GLU
5	D	72(C)	VAL
5	D	72(D)	PRO
5	D	72(E)	VAL
5	D	72(F)	THR
5	D	72(G)	SER
5	D	72(H)	PHE
5	D	96	LEU
6	E	6	GLN
6	E	87	THR
6	E	110	LEU
6	E	152	TRP
6	E	153	LYS
6	E	189	TRP
6	E	190	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	66	HIS
1	G	203	GLN
1	G	246	GLN
1	G	287	GLN
1	G	289	ASN
1	G	339	ASN
1	G	440	GLN
1	G	478	ASN
2	B	564	HIS
2	B	575	GLN
2	B	607	ASN
2	B	653	GLN
3	L	168	GLN
3	L	195	GLN
4	H	5	GLN
4	H	60	ASN
4	H	97	GLN
4	H	162	HIS
4	H	198	HIS
5	D	6	GLN
5	D	33	HIS
5	D	35	ASN
5	D	64	GLN
5	D	105	GLN
6	E	32	HIS
6	E	192	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

76 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
7	NAG	A	1	7,1	14,14,15	2.46	2 (14%)	17,19,21	1.77	3 (17%)
7	NAG	A	2	7	14,14,15	1.46	2 (14%)	17,19,21	0.94	1 (5%)
7	BMA	A	3	7	11,11,12	1.64	3 (27%)	15,15,17	1.59	2 (13%)
7	MAN	A	4	7	11,11,12	1.04	1 (9%)	15,15,17	0.90	0
7	MAN	A	5	7	11,11,12	0.77	0	15,15,17	0.93	1 (6%)
7	MAN	A	6	7	11,11,12	0.72	0	15,15,17	0.86	1 (6%)
8	NAG	C	1	1,8	14,14,15	1.12	1 (7%)	17,19,21	1.67	2 (11%)
8	NAG	C	2	8	14,14,15	1.08	2 (14%)	17,19,21	1.37	2 (11%)
8	BMA	C	3	8	11,11,12	0.90	1 (9%)	15,15,17	0.89	0
8	MAN	C	4	8	11,11,12	0.96	1 (9%)	15,15,17	1.22	3 (20%)
8	MAN	C	5	8	11,11,12	0.57	0	15,15,17	0.98	2 (13%)
8	MAN	C	6	8	11,11,12	0.90	1 (9%)	15,15,17	1.41	3 (20%)
8	MAN	C	7	8	11,11,12	0.85	1 (9%)	15,15,17	1.15	2 (13%)
9	NAG	F	1	9,1	14,14,15	0.22	0	17,19,21	0.54	0
9	NAG	F	2	9	14,14,15	0.59	0	17,19,21	1.03	1 (5%)
9	BMA	F	3	9	11,11,12	1.84	2 (18%)	15,15,17	1.17	2 (13%)
9	MAN	F	4	9	11,11,12	2.06	3 (27%)	15,15,17	1.45	3 (20%)
9	MAN	F	5	9	11,11,12	1.16	2 (18%)	15,15,17	1.57	3 (20%)
9	MAN	F	6	9	11,11,12	0.73	0	15,15,17	0.98	2 (13%)
9	MAN	F	7	9	11,11,12	0.65	0	15,15,17	0.98	2 (13%)
10	NAG	I	1	1,10	14,14,15	0.31	0	17,19,21	0.69	0
10	NAG	I	2	10	14,14,15	0.26	0	17,19,21	0.56	0
10	BMA	I	3	10	11,11,12	0.48	0	15,15,17	0.78	0
11	NAG	J	1	11,1	14,14,15	0.20	0	17,19,21	0.53	0
11	NAG	J	2	11	14,14,15	0.23	0	17,19,21	0.46	0
11	BMA	J	3	11	11,11,12	0.51	0	15,15,17	0.82	0
11	MAN	J	4	11	11,11,12	0.66	0	15,15,17	1.23	2 (13%)
11	MAN	J	5	11	11,11,12	0.77	0	15,15,17	1.11	1 (6%)
11	MAN	J	6	11	11,11,12	0.80	0	15,15,17	1.00	2 (13%)
11	MAN	J	7	11	11,11,12	0.67	0	15,15,17	1.16	2 (13%)
11	MAN	J	8	11	11,11,12	0.70	0	15,15,17	0.85	1 (6%)
9	NAG	K	1	9,1	14,14,15	0.69	1 (7%)	17,19,21	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	NAG	K	2	9	14,14,15	0.77	1 (7%)	17,19,21	0.96	1 (5%)
9	BMA	K	3	9	11,11,12	0.34	0	15,15,17	1.63	2 (13%)
9	MAN	K	4	9	11,11,12	1.30	2 (18%)	15,15,17	1.38	3 (20%)
9	MAN	K	5	9	11,11,12	1.05	0	15,15,17	1.53	3 (20%)
9	MAN	K	6	9	11,11,12	0.96	1 (9%)	15,15,17	1.05	2 (13%)
9	MAN	K	7	9	11,11,12	0.33	0	15,15,17	0.93	1 (6%)
11	NAG	M	1	11,1	14,14,15	0.22	0	17,19,21	0.56	0
11	NAG	M	2	11	14,14,15	0.20	0	17,19,21	0.42	0
11	BMA	M	3	11	11,11,12	0.58	0	15,15,17	0.69	0
11	MAN	M	4	11	11,11,12	0.80	0	15,15,17	1.25	2 (13%)
11	MAN	M	5	11	11,11,12	0.68	0	15,15,17	1.17	2 (13%)
11	MAN	M	6	11	11,11,12	0.79	1 (9%)	15,15,17	0.91	1 (6%)
11	MAN	M	7	11	11,11,12	0.76	1 (9%)	15,15,17	1.08	2 (13%)
11	MAN	M	8	11	11,11,12	0.59	0	15,15,17	0.97	2 (13%)
12	NAG	N	1	1,12	14,14,15	1.80	1 (7%)	17,19,21	1.02	0
12	NAG	N	2	12	14,14,15	0.42	0	17,19,21	0.68	1 (5%)
12	BMA	N	3	12	11,11,12	0.75	0	15,15,17	0.79	0
12	MAN	N	4	12	11,11,12	0.79	0	15,15,17	1.09	2 (13%)
12	MAN	N	5	12	11,11,12	1.38	2 (18%)	15,15,17	1.22	2 (13%)
12	MAN	N	6	12	11,11,12	0.91	1 (9%)	15,15,17	1.51	2 (13%)
13	NAG	O	1	13,1	14,14,15	0.88	1 (7%)	17,19,21	1.49	2 (11%)
13	MAN	O	10	13	11,11,12	0.57	0	15,15,17	1.09	2 (13%)
13	NAG	O	2	13	14,14,15	0.33	0	17,19,21	0.53	0
13	BMA	O	3	13	11,11,12	0.98	0	15,15,17	0.97	1 (6%)
13	MAN	O	4	13	11,11,12	0.82	1 (9%)	15,15,17	1.10	1 (6%)
13	MAN	O	5	13	11,11,12	0.92	1 (9%)	15,15,17	1.16	1 (6%)
13	MAN	O	6	13	11,11,12	0.96	1 (9%)	15,15,17	0.80	1 (6%)
13	MAN	O	7	13	11,11,12	0.56	0	15,15,17	1.15	2 (13%)
13	MAN	O	8	13	11,11,12	0.73	0	15,15,17	1.00	1 (6%)
13	MAN	O	9	13	11,11,12	0.66	0	15,15,17	1.17	2 (13%)
14	NAG	P	1	1,14	14,14,15	2.21	4 (28%)	17,19,21	0.97	1 (5%)
14	NAG	P	2	14	14,14,15	0.35	0	17,19,21	0.61	0
14	BMA	P	3	14	11,11,12	0.59	0	15,15,17	1.57	2 (13%)
14	MAN	P	4	14	11,11,12	0.90	1 (9%)	15,15,17	1.34	1 (6%)
14	MAN	P	5	14	11,11,12	1.10	1 (9%)	15,15,17	1.30	3 (20%)
14	MAN	P	6	14	11,11,12	1.05	1 (9%)	15,15,17	0.95	1 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
14	MAN	P	7	14	11,11,12	1.04	1 (9%)	15,15,17	0.83	1 (6%)
15	NAG	Q	1	1,15	14,14,15	0.56	0	17,19,21	1.07	1 (5%)
15	NAG	Q	2	15	14,14,15	0.68	1 (7%)	17,19,21	1.00	1 (5%)
15	NAG	R	1	1,15	14,14,15	0.14	0	17,19,21	0.81	1 (5%)
15	NAG	R	2	15	14,14,15	0.34	0	17,19,21	0.53	0
10	NAG	S	1	10,2	14,14,15	0.42	0	17,19,21	0.61	0
10	NAG	S	2	10	14,14,15	0.24	0	17,19,21	0.57	0
10	BMA	S	3	10	11,11,12	0.52	0	15,15,17	1.10	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
7	NAG	A	1	7,1	-	5/6/23/26	0/1/1/1
7	NAG	A	2	7	-	3/6/23/26	0/1/1/1
7	BMA	A	3	7	-	2/2/19/22	0/1/1/1
7	MAN	A	4	7	-	0/2/19/22	0/1/1/1
7	MAN	A	5	7	-	0/2/19/22	0/1/1/1
7	MAN	A	6	7	-	0/2/19/22	0/1/1/1
8	NAG	C	1	1,8	-	2/6/23/26	0/1/1/1
8	NAG	C	2	8	-	4/6/23/26	0/1/1/1
8	BMA	C	3	8	-	2/2/19/22	0/1/1/1
8	MAN	C	4	8	-	0/2/19/22	0/1/1/1
8	MAN	C	5	8	-	0/2/19/22	0/1/1/1
8	MAN	C	6	8	-	0/2/19/22	0/1/1/1
8	MAN	C	7	8	-	0/2/19/22	0/1/1/1
9	NAG	F	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	F	2	9	-	2/6/23/26	0/1/1/1
9	BMA	F	3	9	-	2/2/19/22	0/1/1/1
9	MAN	F	4	9	-	2/2/19/22	0/1/1/1
9	MAN	F	5	9	-	0/2/19/22	0/1/1/1
9	MAN	F	6	9	-	0/2/19/22	0/1/1/1
9	MAN	F	7	9	-	0/2/19/22	0/1/1/1
10	NAG	I	1	1,10	-	0/6/23/26	0/1/1/1
10	NAG	I	2	10	-	2/6/23/26	0/1/1/1
10	BMA	I	3	10	-	0/2/19/22	0/1/1/1
11	NAG	J	1	11,1	-	2/6/23/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
11	NAG	J	2	11	-	2/6/23/26	0/1/1/1
11	BMA	J	3	11	-	2/2/19/22	0/1/1/1
11	MAN	J	4	11	-	1/2/19/22	0/1/1/1
11	MAN	J	5	11	-	0/2/19/22	0/1/1/1
11	MAN	J	6	11	-	0/2/19/22	0/1/1/1
11	MAN	J	7	11	-	2/2/19/22	0/1/1/1
11	MAN	J	8	11	-	0/2/19/22	0/1/1/1
9	NAG	K	1	9,1	-	0/6/23/26	0/1/1/1
9	NAG	K	2	9	-	2/6/23/26	0/1/1/1
9	BMA	K	3	9	-	2/2/19/22	0/1/1/1
9	MAN	K	4	9	-	0/2/19/22	0/1/1/1
9	MAN	K	5	9	-	2/2/19/22	0/1/1/1
9	MAN	K	6	9	-	0/2/19/22	0/1/1/1
9	MAN	K	7	9	-	0/2/19/22	0/1/1/1
11	NAG	M	1	11,1	-	2/6/23/26	0/1/1/1
11	NAG	M	2	11	-	0/6/23/26	0/1/1/1
11	BMA	M	3	11	-	0/2/19/22	0/1/1/1
11	MAN	M	4	11	-	2/2/19/22	0/1/1/1
11	MAN	M	5	11	-	2/2/19/22	0/1/1/1
11	MAN	M	6	11	-	0/2/19/22	0/1/1/1
11	MAN	M	7	11	-	0/2/19/22	0/1/1/1
11	MAN	M	8	11	-	0/2/19/22	0/1/1/1
12	NAG	N	1	1,12	-	2/6/23/26	0/1/1/1
12	NAG	N	2	12	-	2/6/23/26	0/1/1/1
12	BMA	N	3	12	-	0/2/19/22	0/1/1/1
12	MAN	N	4	12	-	0/2/19/22	0/1/1/1
12	MAN	N	5	12	-	1/2/19/22	0/1/1/1
12	MAN	N	6	12	-	2/2/19/22	0/1/1/1
13	NAG	O	1	13,1	-	4/6/23/26	0/1/1/1
13	MAN	O	10	13	-	0/2/19/22	0/1/1/1
13	NAG	O	2	13	-	2/6/23/26	0/1/1/1
13	BMA	O	3	13	-	0/2/19/22	0/1/1/1
13	MAN	O	4	13	-	2/2/19/22	0/1/1/1
13	MAN	O	5	13	-	0/2/19/22	0/1/1/1
13	MAN	O	6	13	-	0/2/19/22	0/1/1/1
13	MAN	O	7	13	-	2/2/19/22	0/1/1/1
13	MAN	O	8	13	-	2/2/19/22	0/1/1/1
13	MAN	O	9	13	-	1/2/19/22	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
14	NAG	P	1	1,14	-	2/6/23/26	0/1/1/1
14	NAG	P	2	14	-	2/6/23/26	0/1/1/1
14	BMA	P	3	14	-	2/2/19/22	0/1/1/1
14	MAN	P	4	14	-	0/2/19/22	0/1/1/1
14	MAN	P	5	14	-	0/2/19/22	0/1/1/1
14	MAN	P	6	14	-	0/2/19/22	0/1/1/1
14	MAN	P	7	14	-	0/2/19/22	0/1/1/1
15	NAG	Q	1	1,15	-	0/6/23/26	0/1/1/1
15	NAG	Q	2	15	-	2/6/23/26	0/1/1/1
15	NAG	R	1	1,15	-	2/6/23/26	0/1/1/1
15	NAG	R	2	15	-	2/6/23/26	0/1/1/1
10	NAG	S	1	10,2	-	0/6/23/26	0/1/1/1
10	NAG	S	2	10	-	2/6/23/26	0/1/1/1
10	BMA	S	3	10	-	0/2/19/22	0/1/1/1

All (46) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	1	NAG	O5-C1	8.69	1.58	1.43
12	N	1	NAG	O5-C1	6.03	1.53	1.43
14	P	1	NAG	C1-C2	-5.42	1.45	1.52
14	P	1	NAG	O5-C1	4.92	1.52	1.43
9	F	4	MAN	O2-C2	4.71	1.53	1.43
7	A	2	NAG	O5-C1	4.18	1.50	1.43
9	F	3	BMA	O3-C3	4.17	1.53	1.43
9	F	4	MAN	C1-C2	3.77	1.61	1.52
9	F	3	BMA	C2-C3	3.58	1.58	1.52
8	C	1	NAG	O5-C1	3.49	1.49	1.43
7	A	3	BMA	O5-C1	3.27	1.49	1.43
13	O	1	NAG	O5-C1	-3.13	1.38	1.43
8	C	2	NAG	C1-C2	3.11	1.56	1.52
7	A	2	NAG	C1-C2	3.00	1.56	1.52
12	N	5	MAN	O5-C1	-3.00	1.38	1.43
9	F	4	MAN	C2-C3	2.94	1.57	1.52
8	C	6	MAN	C1-C2	2.85	1.59	1.52
9	K	2	NAG	C1-C2	2.82	1.56	1.52
9	K	4	MAN	C1-C2	2.81	1.58	1.52
14	P	6	MAN	O5-C1	-2.79	1.39	1.43
14	P	7	MAN	O5-C1	-2.73	1.39	1.43
7	A	3	BMA	O5-C5	2.73	1.48	1.43
13	O	6	MAN	O5-C1	-2.73	1.39	1.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	A	1	NAG	C3-C2	2.58	1.57	1.52
8	C	4	MAN	C1-C2	2.56	1.58	1.52
9	K	4	MAN	C2-C3	2.48	1.56	1.52
15	Q	2	NAG	C1-C2	2.35	1.55	1.52
9	K	6	MAN	O5-C1	-2.35	1.39	1.43
14	P	4	MAN	C1-C2	2.34	1.57	1.52
8	C	2	NAG	O5-C1	2.33	1.47	1.43
11	M	7	MAN	C1-C2	2.31	1.57	1.52
8	C	7	MAN	C1-C2	2.20	1.57	1.52
14	P	1	NAG	O7-C7	2.20	1.28	1.23
9	K	1	NAG	O5-C1	-2.19	1.40	1.43
13	O	4	MAN	C1-C2	2.17	1.57	1.52
14	P	1	NAG	C8-C7	-2.16	1.46	1.50
9	F	5	MAN	O5-C1	-2.15	1.40	1.43
11	M	6	MAN	O5-C1	-2.14	1.40	1.43
7	A	3	BMA	C1-C2	2.14	1.57	1.52
8	C	3	BMA	C1-C2	2.10	1.57	1.52
14	P	5	MAN	O5-C1	-2.09	1.40	1.43
12	N	6	MAN	C1-C2	2.07	1.57	1.52
9	F	5	MAN	C1-C2	2.07	1.57	1.52
13	O	5	MAN	C1-C2	2.04	1.57	1.52
7	A	4	MAN	C2-C3	2.02	1.55	1.52
12	N	5	MAN	O2-C2	2.01	1.47	1.43

All (97) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	1	NAG	C1-O5-C5	5.40	119.43	112.19
7	A	1	NAG	C2-N2-C7	5.10	129.74	122.90
12	N	6	MAN	C1-O5-C5	4.93	118.79	112.19
13	O	1	NAG	C2-N2-C7	4.73	129.24	122.90
9	K	3	BMA	C1-O5-C5	4.72	118.52	112.19
7	A	3	BMA	O3-C3-C2	4.23	118.68	110.05
8	C	2	NAG	C2-N2-C7	4.10	128.40	122.90
14	P	3	BMA	C1-O5-C5	4.09	117.66	112.19
9	K	5	MAN	O2-C2-C3	-3.91	102.06	110.15
11	M	4	MAN	C1-O5-C5	3.90	117.41	112.19
9	F	5	MAN	O2-C2-C1	3.89	118.13	109.22
15	Q	2	NAG	C1-O5-C5	3.85	117.34	112.19
14	P	4	MAN	C1-O5-C5	3.74	117.20	112.19
9	F	5	MAN	O2-C2-C3	-3.63	102.62	110.15
11	J	7	MAN	C1-O5-C5	3.61	117.02	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
8	C	6	MAN	C1-O5-C5	3.56	116.95	112.19
12	N	5	MAN	C1-C2-C3	-3.55	104.48	109.64
13	O	7	MAN	C1-O5-C5	3.41	116.76	112.19
14	P	3	BMA	O3-C3-C2	3.39	116.97	110.05
7	A	2	NAG	C1-O5-C5	3.35	116.68	112.19
13	O	9	MAN	C1-O5-C5	3.32	116.64	112.19
13	O	5	MAN	O2-C2-C3	-3.23	103.47	110.15
13	O	10	MAN	C1-O5-C5	3.11	116.35	112.19
9	F	3	BMA	O3-C3-C4	3.10	117.68	110.38
11	M	5	MAN	C1-O5-C5	3.08	116.32	112.19
14	P	5	MAN	O2-C2-C3	-3.03	103.87	110.15
9	F	4	MAN	O2-C2-C1	3.03	116.15	109.22
9	K	3	BMA	C1-C2-C3	3.00	114.02	109.64
10	S	3	BMA	C1-O5-C5	3.00	116.21	112.19
7	A	3	BMA	C1-C2-C3	-2.98	105.30	109.64
8	C	4	MAN	C1-O5-C5	2.96	116.15	112.19
9	K	2	NAG	C1-O5-C5	2.92	116.10	112.19
9	K	5	MAN	C1-O5-C5	2.91	116.09	112.19
9	K	4	MAN	C1-O5-C5	2.78	115.91	112.19
14	P	5	MAN	C1-C2-C3	-2.78	105.60	109.64
11	M	5	MAN	O2-C2-C3	-2.65	104.67	110.15
11	M	7	MAN	C1-O5-C5	2.63	115.71	112.19
11	J	4	MAN	C1-O5-C5	2.61	115.69	112.19
15	Q	1	NAG	O5-C1-C2	-2.61	107.25	111.29
15	R	1	NAG	C1-O5-C5	2.59	115.66	112.19
8	C	6	MAN	C1-C2-C3	2.59	113.42	109.64
11	M	8	MAN	C1-O5-C5	2.59	115.65	112.19
8	C	2	NAG	C1-C2-N2	2.57	114.49	110.43
11	J	4	MAN	O2-C2-C3	-2.57	104.82	110.15
13	O	1	NAG	C1-C2-N2	2.56	114.47	110.43
9	F	4	MAN	O3-C3-C2	2.56	115.28	110.05
11	J	5	MAN	O2-C2-C3	-2.56	104.86	110.15
9	K	5	MAN	C1-C2-C3	-2.55	105.93	109.64
9	F	2	NAG	O4-C4-C3	-2.53	104.40	110.38
9	K	4	MAN	O2-C2-C1	2.53	115.03	109.22
9	K	4	MAN	C1-C2-C3	2.51	113.30	109.64
9	F	7	MAN	C1-O5-C5	2.49	115.52	112.19
9	F	6	MAN	C1-O5-C5	2.45	115.47	112.19
8	C	7	MAN	O2-C2-C3	-2.44	105.09	110.15
13	O	3	BMA	C1-C2-C3	2.43	113.19	109.64
12	N	5	MAN	O3-C3-C2	2.41	114.97	110.05
12	N	4	MAN	O2-C2-C3	-2.41	105.17	110.15

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
12	N	2	NAG	C1-O5-C5	2.40	115.40	112.19
9	F	6	MAN	O2-C2-C3	-2.38	105.23	110.15
13	O	8	MAN	C1-O5-C5	2.36	115.35	112.19
11	M	6	MAN	O2-C2-C3	-2.36	105.27	110.15
13	O	7	MAN	O2-C2-C3	-2.36	105.27	110.15
13	O	9	MAN	O2-C2-C3	-2.33	105.32	110.15
9	F	5	MAN	C1-O5-C5	2.33	115.31	112.19
11	J	6	MAN	C1-O5-C5	2.33	115.30	112.19
7	A	1	NAG	C1-C2-N2	2.32	114.08	110.43
14	P	7	MAN	O2-C2-C3	-2.27	105.45	110.15
8	C	5	MAN	O2-C2-C3	-2.27	105.45	110.15
13	O	6	MAN	O2-C2-C3	-2.27	105.45	110.15
14	P	6	MAN	O2-C2-C3	-2.25	105.50	110.15
9	K	6	MAN	C1-O5-C5	2.24	115.19	112.19
7	A	6	MAN	O2-C2-C3	-2.24	105.51	110.15
9	F	7	MAN	O2-C2-C3	-2.24	105.51	110.15
11	J	7	MAN	O2-C2-C3	-2.24	105.51	110.15
9	K	6	MAN	O2-C2-C3	-2.24	105.51	110.15
9	F	3	BMA	O3-C3-C2	2.23	114.61	110.05
13	O	10	MAN	O2-C2-C3	-2.23	105.53	110.15
11	J	6	MAN	O2-C2-C3	-2.22	105.55	110.15
11	J	8	MAN	O2-C2-C3	-2.22	105.55	110.15
7	A	5	MAN	O2-C2-C3	-2.22	105.55	110.15
11	M	7	MAN	O2-C2-C3	-2.22	105.56	110.15
8	C	4	MAN	C1-C2-C3	2.21	112.86	109.64
8	C	5	MAN	C1-O5-C5	2.20	115.13	112.19
9	F	4	MAN	O2-C2-C3	2.17	114.65	110.15
8	C	4	MAN	O2-C2-C3	-2.17	105.66	110.15
12	N	6	MAN	O2-C2-C3	-2.14	105.72	110.15
8	C	7	MAN	C1-O5-C5	2.13	115.04	112.19
8	C	6	MAN	O2-C2-C3	-2.10	105.80	110.15
7	A	1	NAG	C8-C7-N2	2.09	119.58	116.12
14	P	5	MAN	C1-O5-C5	2.08	114.98	112.19
8	C	1	NAG	O6-C6-C5	2.07	118.39	111.33
12	N	4	MAN	C3-C4-C5	2.07	113.98	110.23
13	O	4	MAN	O2-C2-C3	-2.07	105.87	110.15
11	M	8	MAN	O2-C2-C3	-2.06	105.89	110.15
9	K	7	MAN	C1-C2-C3	2.05	112.63	109.64
11	M	4	MAN	O2-C2-C3	-2.03	105.95	110.15
14	P	1	NAG	O5-C1-C2	-2.02	108.17	111.29

There are no chirality outliers.

All (81) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	1	NAG	C3-C2-N2-C7
8	C	1	NAG	O5-C5-C6-O6
11	J	1	NAG	O5-C5-C6-O6
14	P	1	NAG	C4-C5-C6-O6
12	N	1	NAG	O5-C5-C6-O6
8	C	3	BMA	C4-C5-C6-O6
14	P	1	NAG	O5-C5-C6-O6
9	K	2	NAG	O5-C5-C6-O6
11	J	7	MAN	O5-C5-C6-O6
11	J	3	BMA	C4-C5-C6-O6
9	K	2	NAG	C4-C5-C6-O6
7	A	1	NAG	O5-C5-C6-O6
8	C	3	BMA	O5-C5-C6-O6
11	J	2	NAG	O5-C5-C6-O6
13	O	4	MAN	O5-C5-C6-O6
8	C	1	NAG	C4-C5-C6-O6
7	A	3	BMA	O5-C5-C6-O6
12	N	2	NAG	C4-C5-C6-O6
11	J	1	NAG	C4-C5-C6-O6
13	O	8	MAN	C4-C5-C6-O6
12	N	1	NAG	C4-C5-C6-O6
11	J	3	BMA	O5-C5-C6-O6
12	N	2	NAG	O5-C5-C6-O6
14	P	2	NAG	O5-C5-C6-O6
13	O	4	MAN	C4-C5-C6-O6
13	O	2	NAG	O5-C5-C6-O6
14	P	2	NAG	C4-C5-C6-O6
9	K	5	MAN	O5-C5-C6-O6
7	A	3	BMA	C4-C5-C6-O6
11	J	7	MAN	C4-C5-C6-O6
9	K	3	BMA	C4-C5-C6-O6
10	S	2	NAG	C4-C5-C6-O6
7	A	1	NAG	C8-C7-N2-C2
7	A	1	NAG	O7-C7-N2-C2
8	C	2	NAG	C8-C7-N2-C2
8	C	2	NAG	O7-C7-N2-C2
10	I	2	NAG	C8-C7-N2-C2
10	I	2	NAG	O7-C7-N2-C2
13	O	1	NAG	C8-C7-N2-C2
13	O	1	NAG	O7-C7-N2-C2
15	R	1	NAG	C8-C7-N2-C2
15	R	1	NAG	O7-C7-N2-C2

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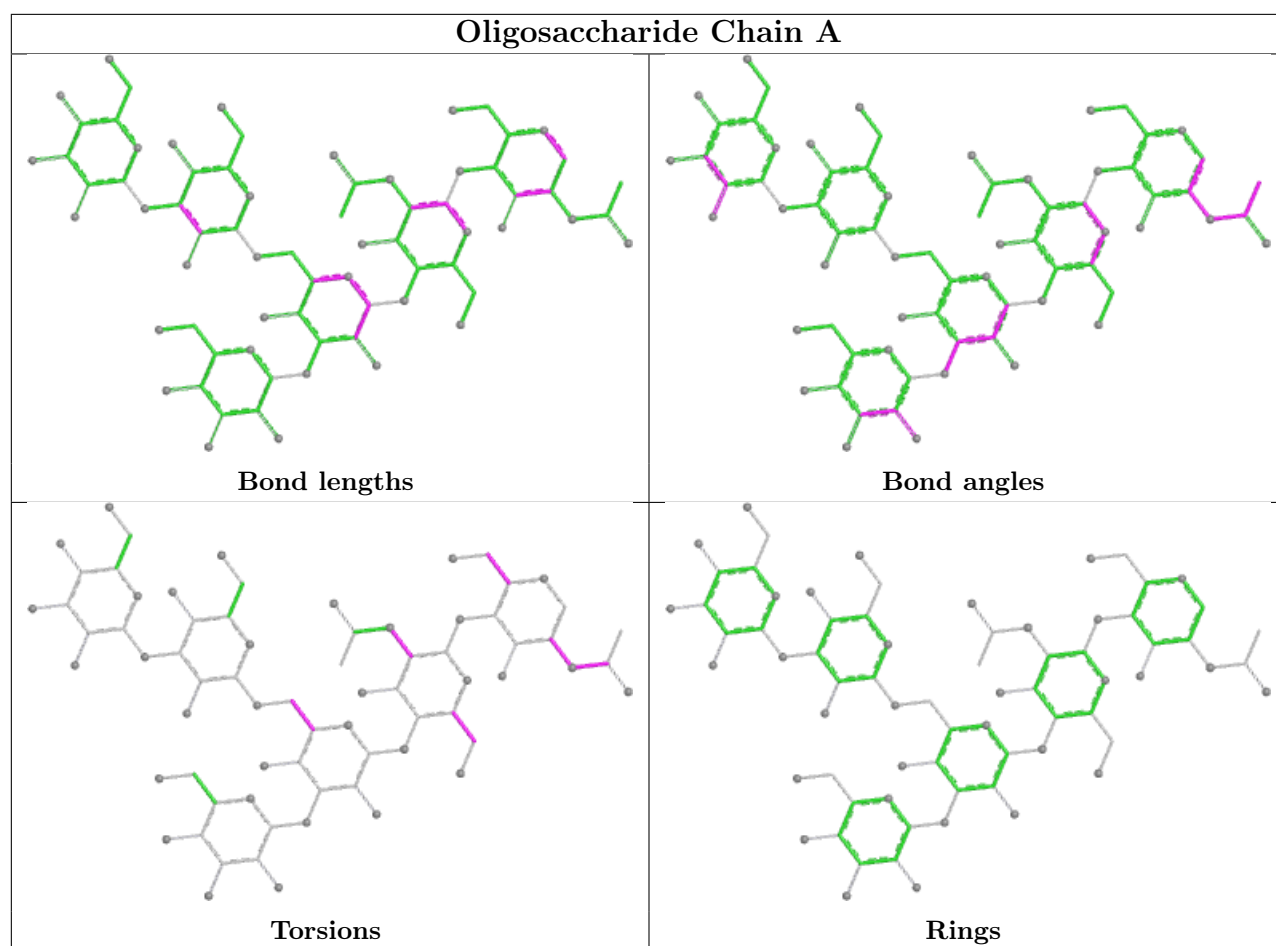
Mol	Chain	Res	Type	Atoms
9	F	4	MAN	C4-C5-C6-O6
11	M	5	MAN	O5-C5-C6-O6
9	F	2	NAG	C4-C5-C6-O6
7	A	1	NAG	C4-C5-C6-O6
11	J	2	NAG	C4-C5-C6-O6
9	K	3	BMA	O5-C5-C6-O6
15	Q	2	NAG	C4-C5-C6-O6
9	F	4	MAN	O5-C5-C6-O6
13	O	8	MAN	O5-C5-C6-O6
12	N	6	MAN	O5-C5-C6-O6
15	R	2	NAG	O5-C5-C6-O6
13	O	2	NAG	C4-C5-C6-O6
10	S	2	NAG	O5-C5-C6-O6
7	A	2	NAG	C4-C5-C6-O6
11	J	4	MAN	O5-C5-C6-O6
15	Q	2	NAG	O5-C5-C6-O6
9	K	5	MAN	C4-C5-C6-O6
7	A	2	NAG	O5-C5-C6-O6
14	P	3	BMA	C4-C5-C6-O6
9	F	3	BMA	C4-C5-C6-O6
9	F	2	NAG	O5-C5-C6-O6
11	M	1	NAG	C4-C5-C6-O6
9	F	3	BMA	O5-C5-C6-O6
13	O	7	MAN	C4-C5-C6-O6
15	R	2	NAG	C4-C5-C6-O6
8	C	2	NAG	C3-C2-N2-C7
13	O	1	NAG	C3-C2-N2-C7
14	P	3	BMA	O5-C5-C6-O6
13	O	7	MAN	O5-C5-C6-O6
11	M	1	NAG	O5-C5-C6-O6
12	N	6	MAN	C4-C5-C6-O6
8	C	2	NAG	C1-C2-N2-C7
13	O	1	NAG	C1-C2-N2-C7
11	M	4	MAN	C4-C5-C6-O6
7	A	2	NAG	C3-C2-N2-C7
13	O	9	MAN	C4-C5-C6-O6
12	N	5	MAN	C4-C5-C6-O6
11	M	5	MAN	C4-C5-C6-O6
11	M	4	MAN	O5-C5-C6-O6

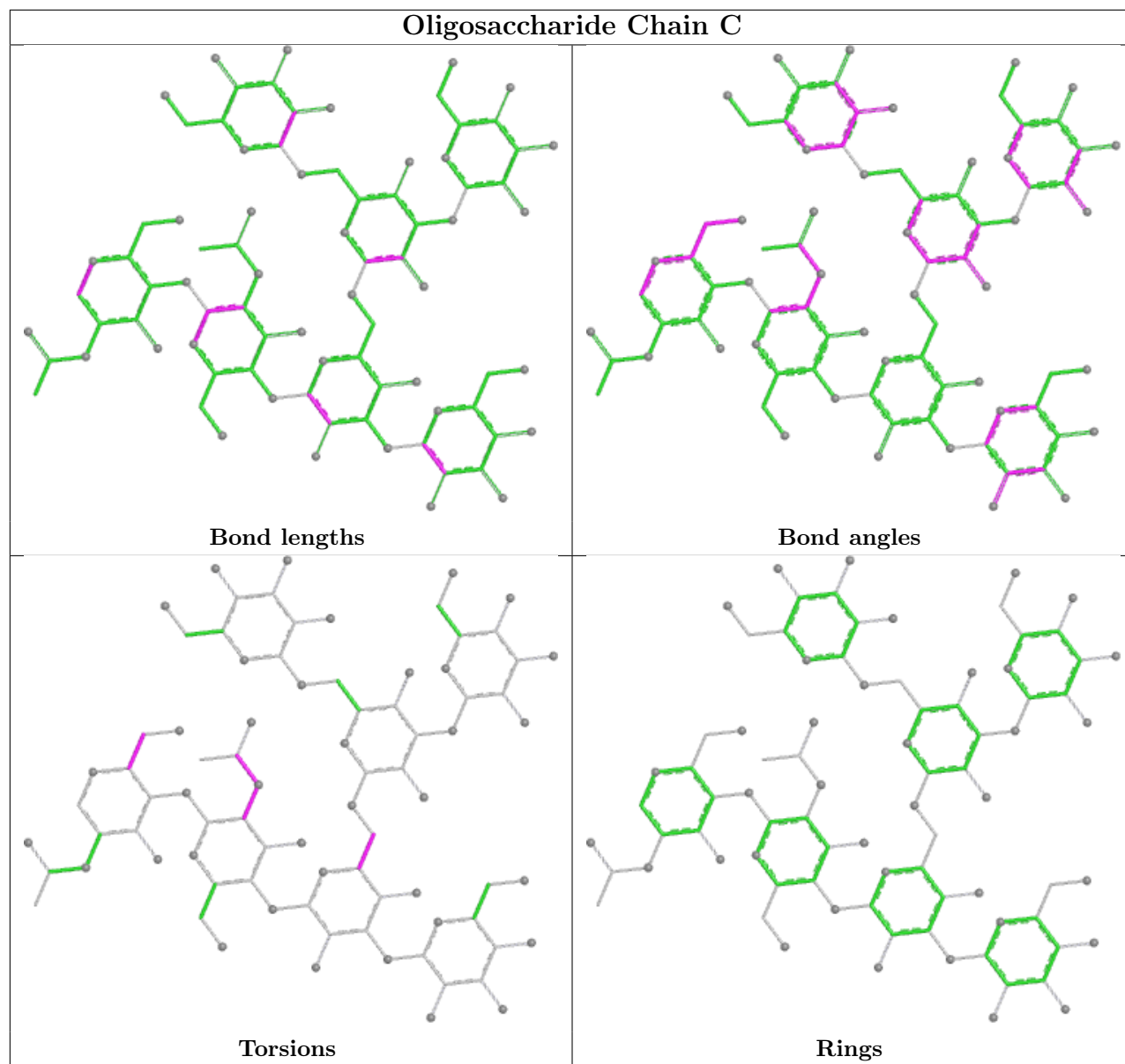
There are no ring outliers.

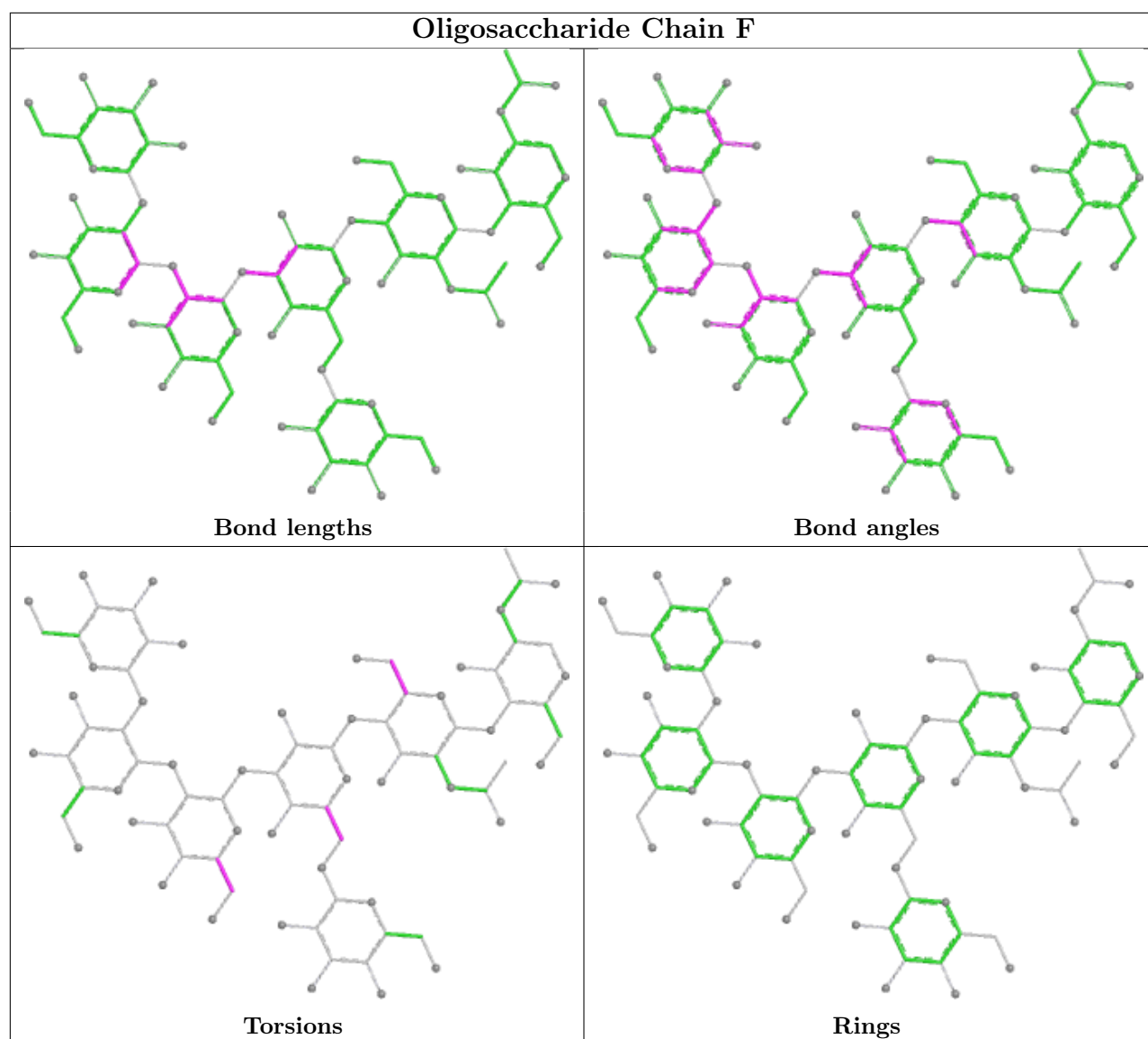
30 monomers are involved in 54 short contacts:

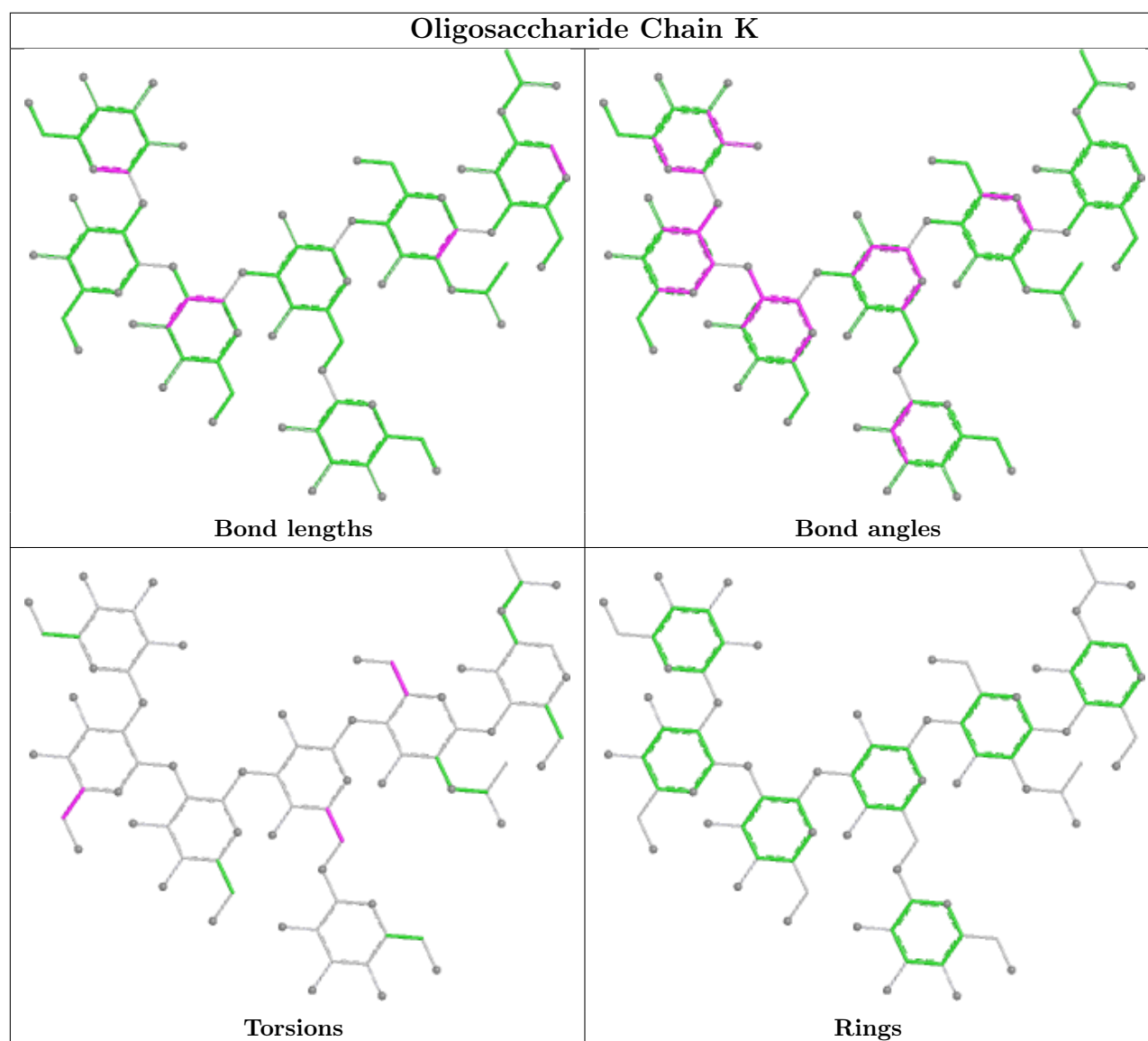
Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	F	6	MAN	1	0
13	O	2	NAG	3	0
9	K	7	MAN	2	0
9	F	3	BMA	3	0
9	F	1	NAG	2	0
11	M	4	MAN	1	0
13	O	4	MAN	1	0
14	P	1	NAG	3	0
8	C	1	NAG	3	0
12	N	1	NAG	1	0
13	O	5	MAN	3	0
11	J	5	MAN	1	0
11	M	1	NAG	1	0
8	C	7	MAN	1	0
8	C	2	NAG	3	0
13	O	9	MAN	1	0
10	I	2	NAG	1	0
13	O	1	NAG	2	0
9	K	1	NAG	6	0
15	Q	1	NAG	2	0
7	A	2	NAG	1	0
9	K	3	BMA	2	0
13	O	3	BMA	1	0
8	C	6	MAN	2	0
10	I	1	NAG	2	0
9	K	2	NAG	2	0
14	P	2	NAG	1	0
7	A	1	NAG	8	0
9	F	5	MAN	1	0
9	F	4	MAN	3	0

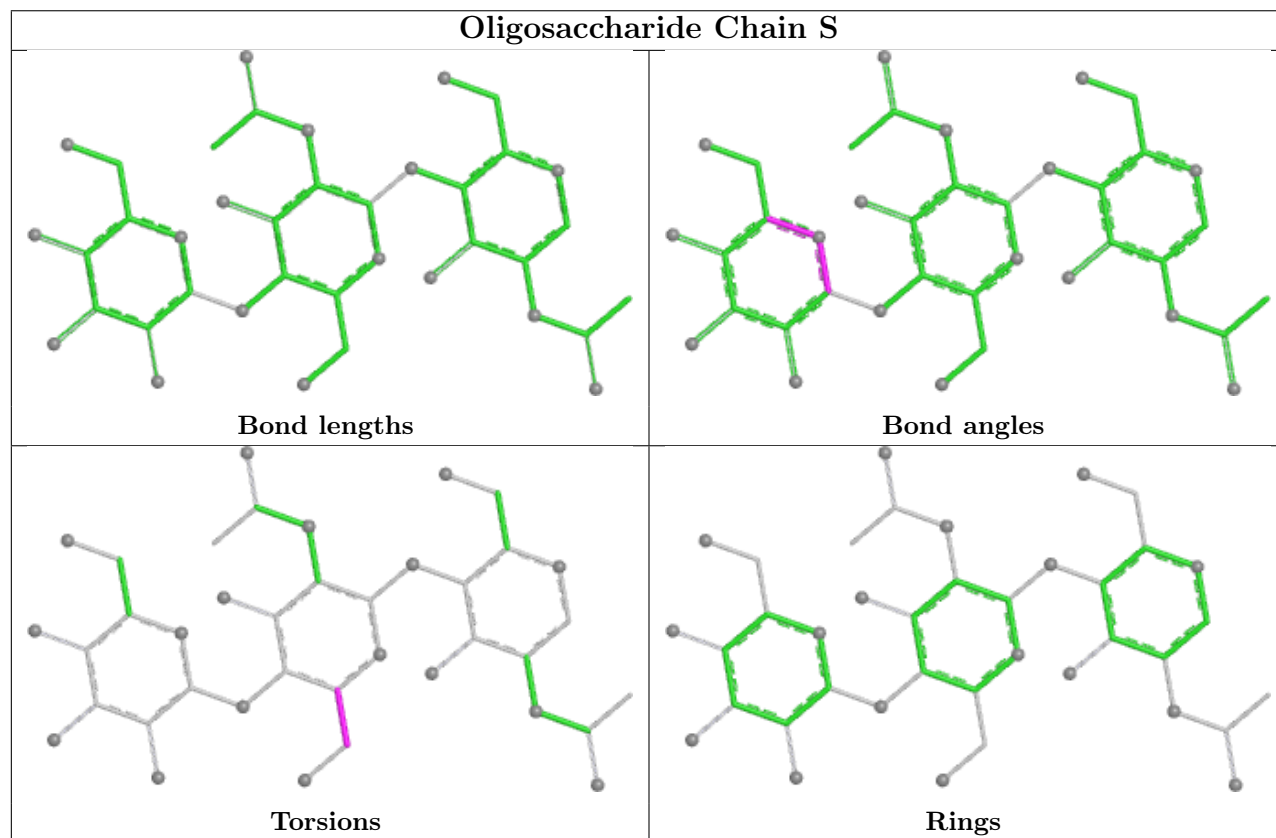
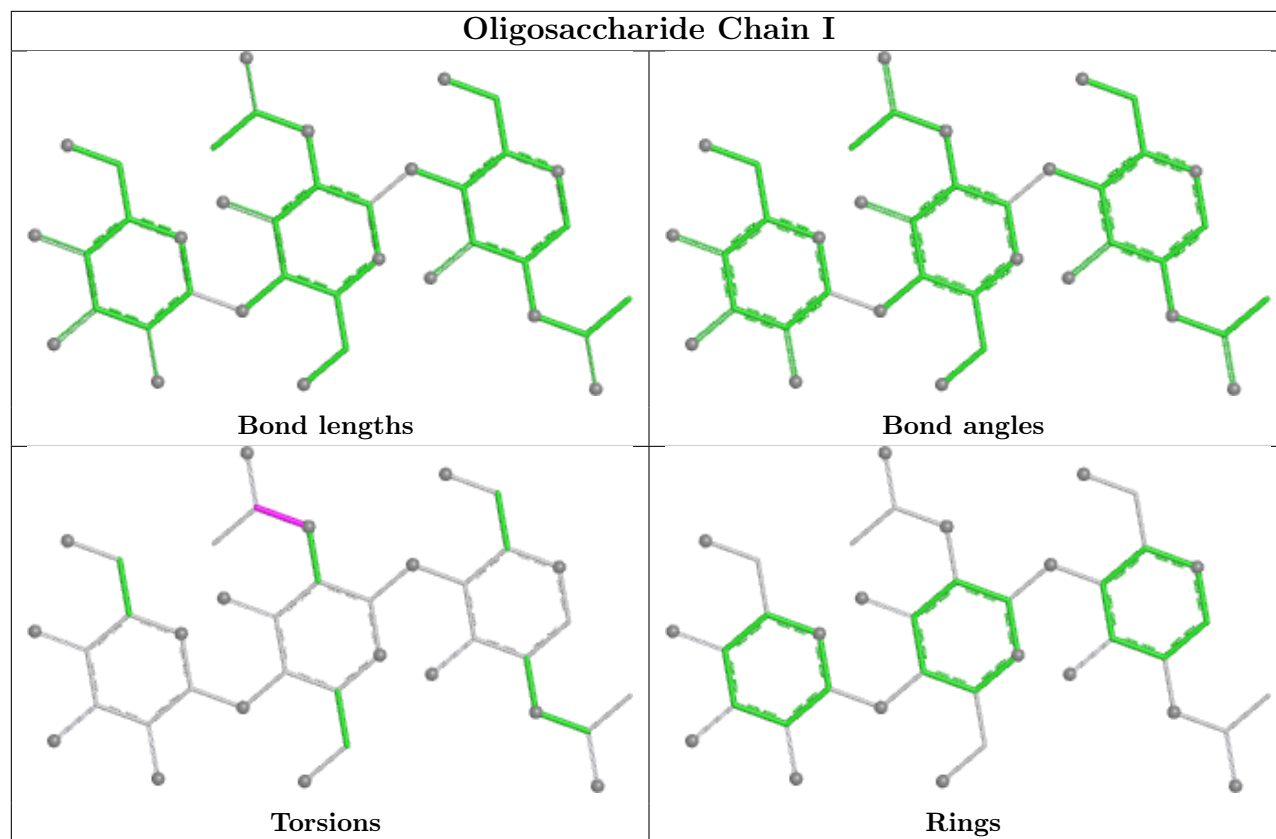
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for oligosaccharide.

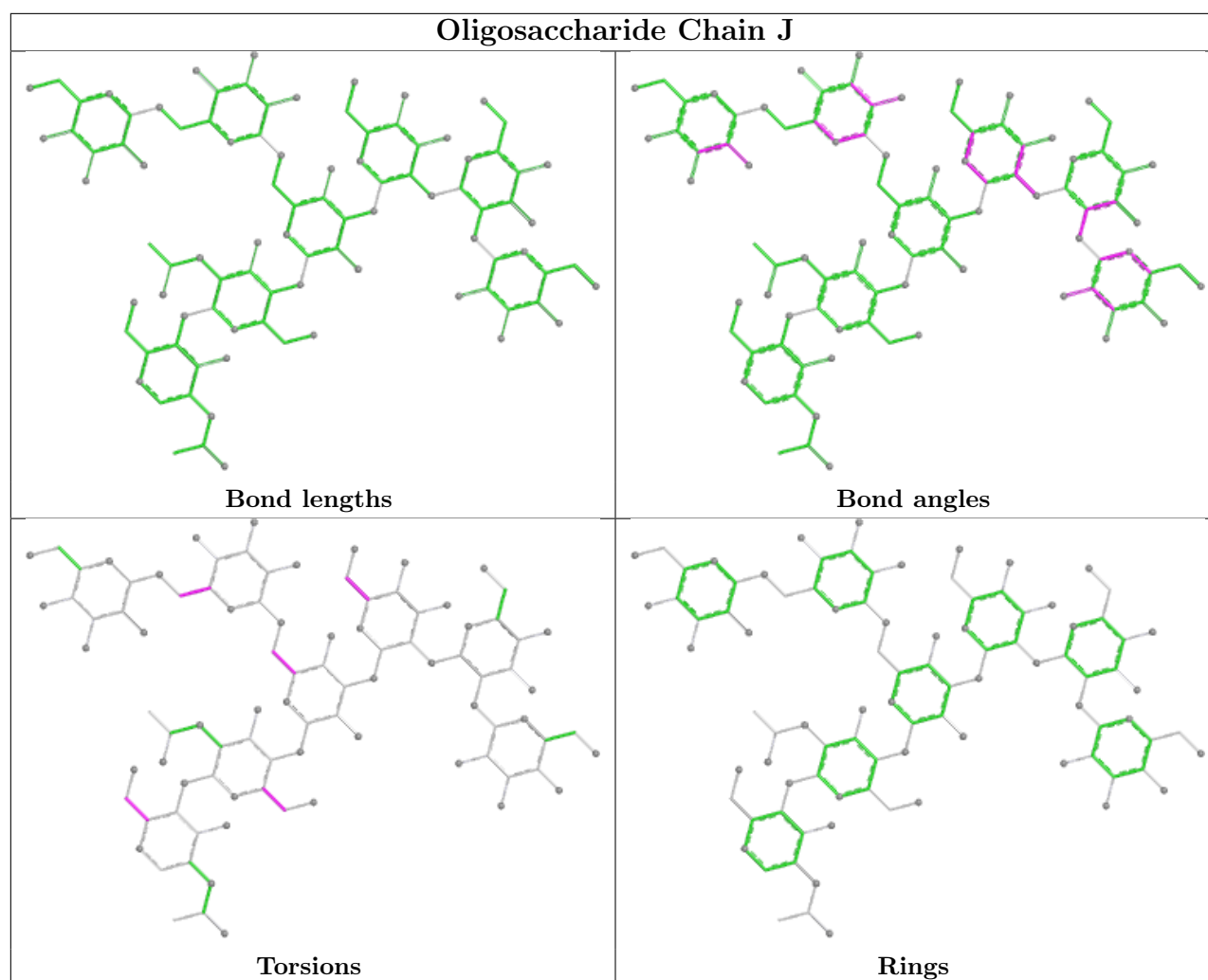




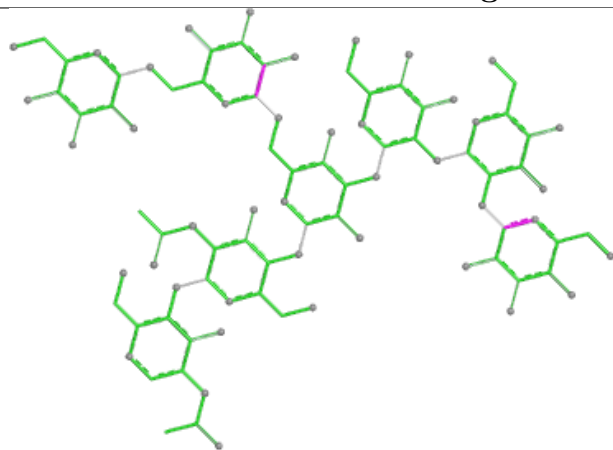




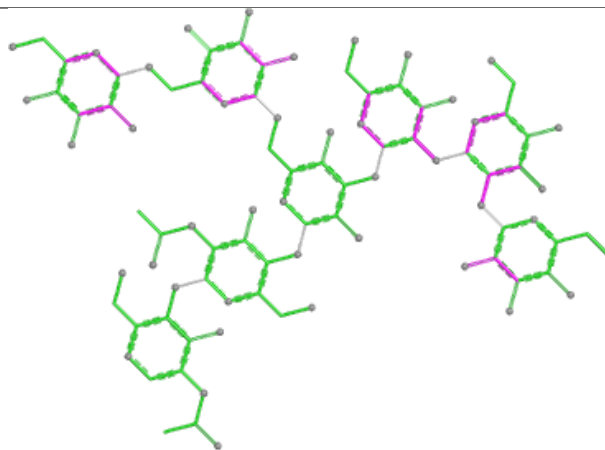




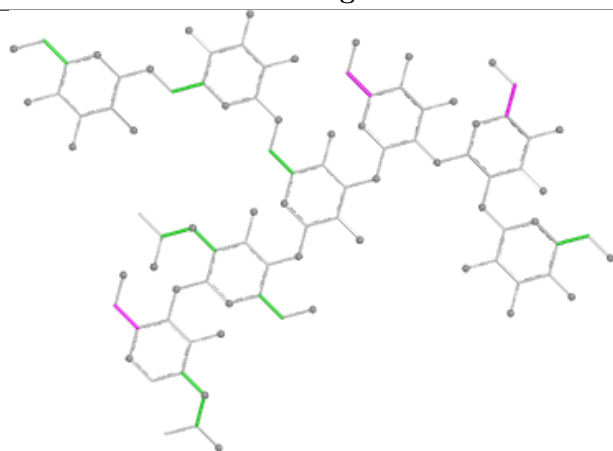
Oligosaccharide Chain M



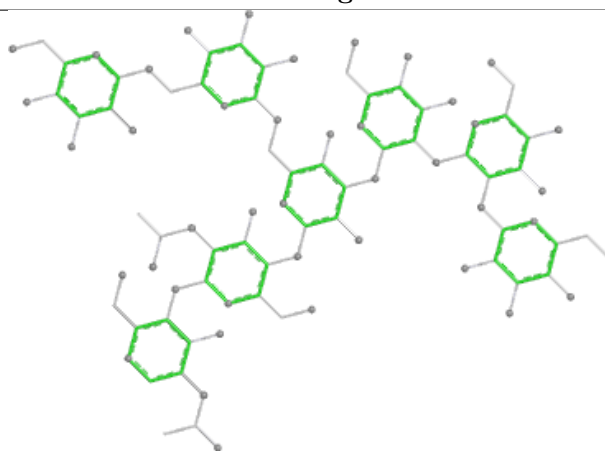
Bond lengths



Bond angles

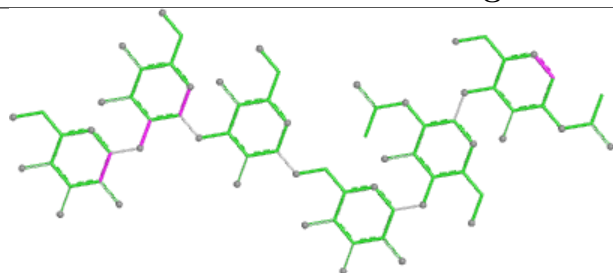


Torsions

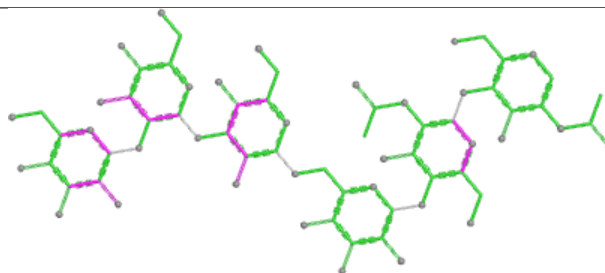


Rings

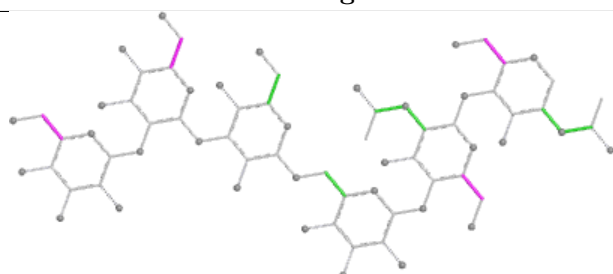
Oligosaccharide Chain N



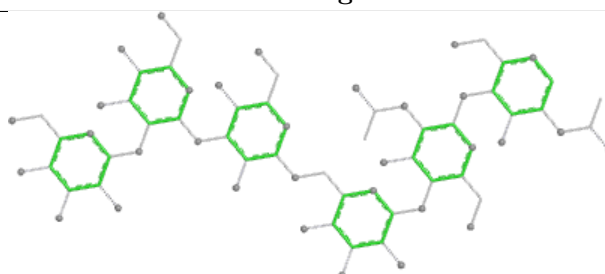
Bond lengths



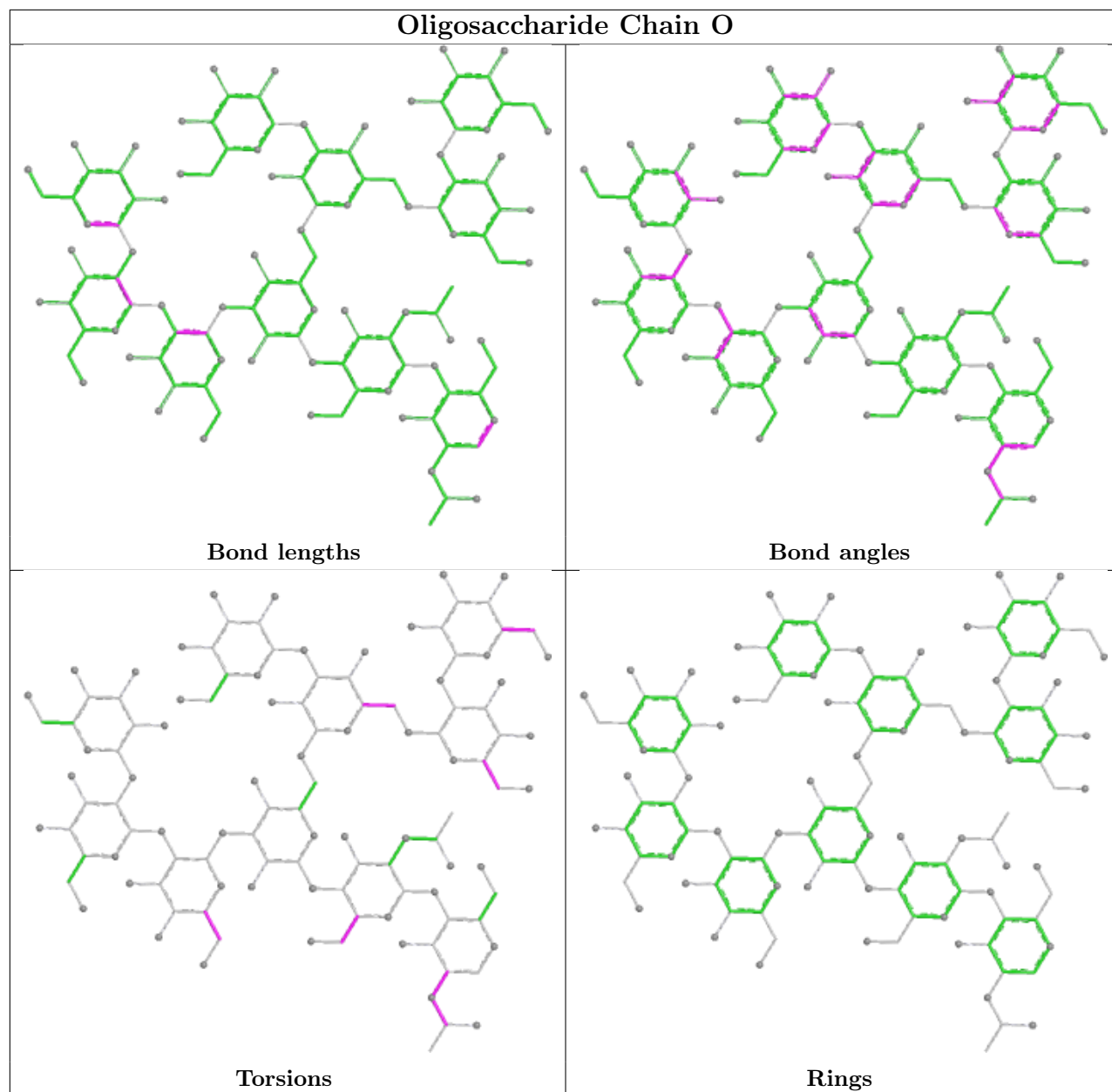
Bond angles

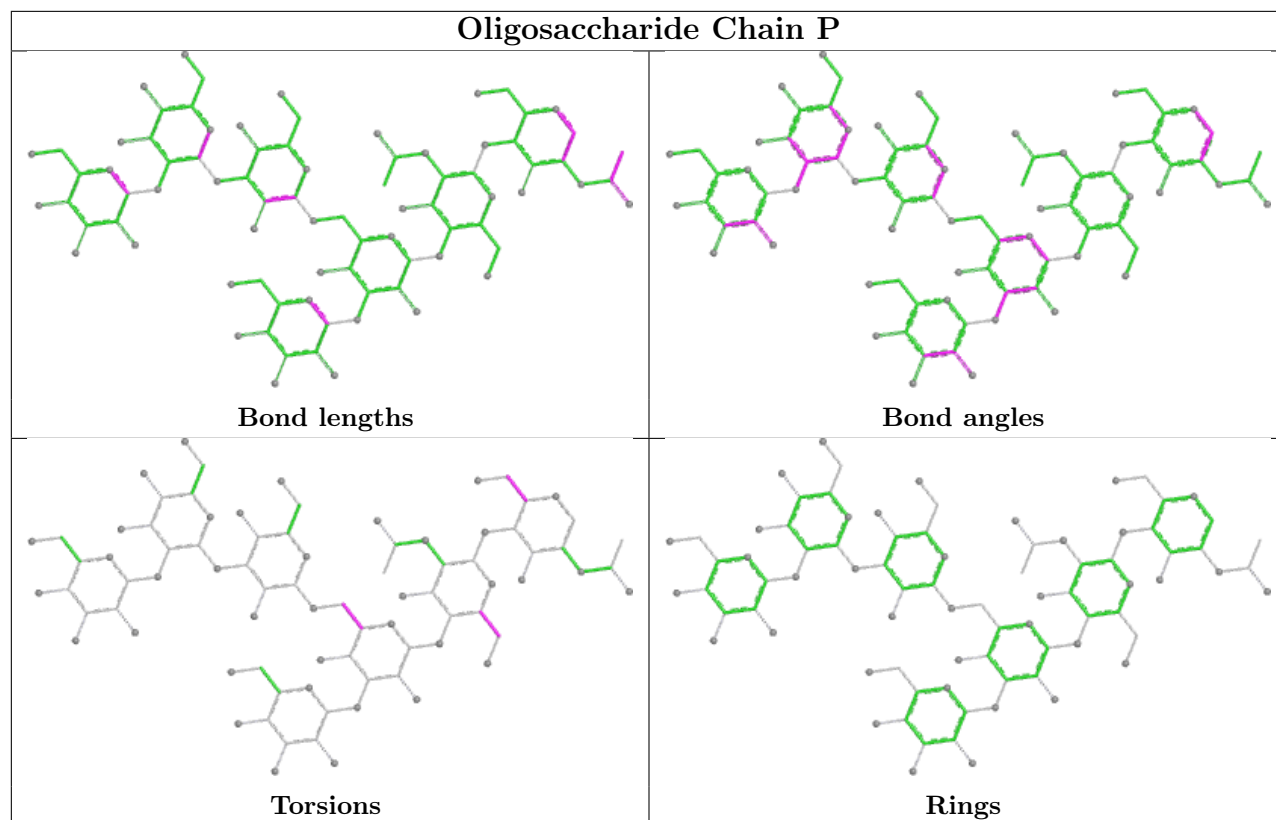


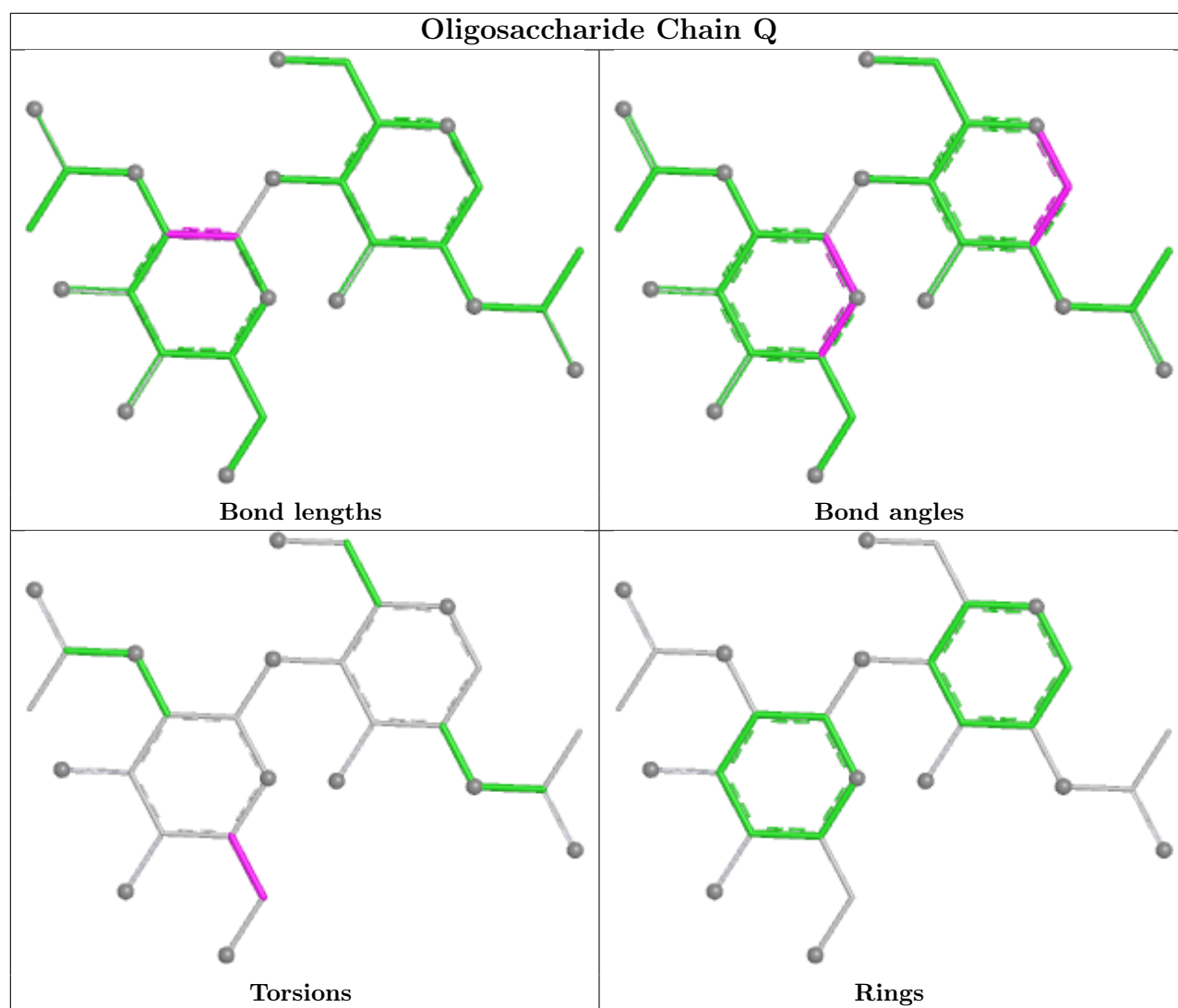
Torsions

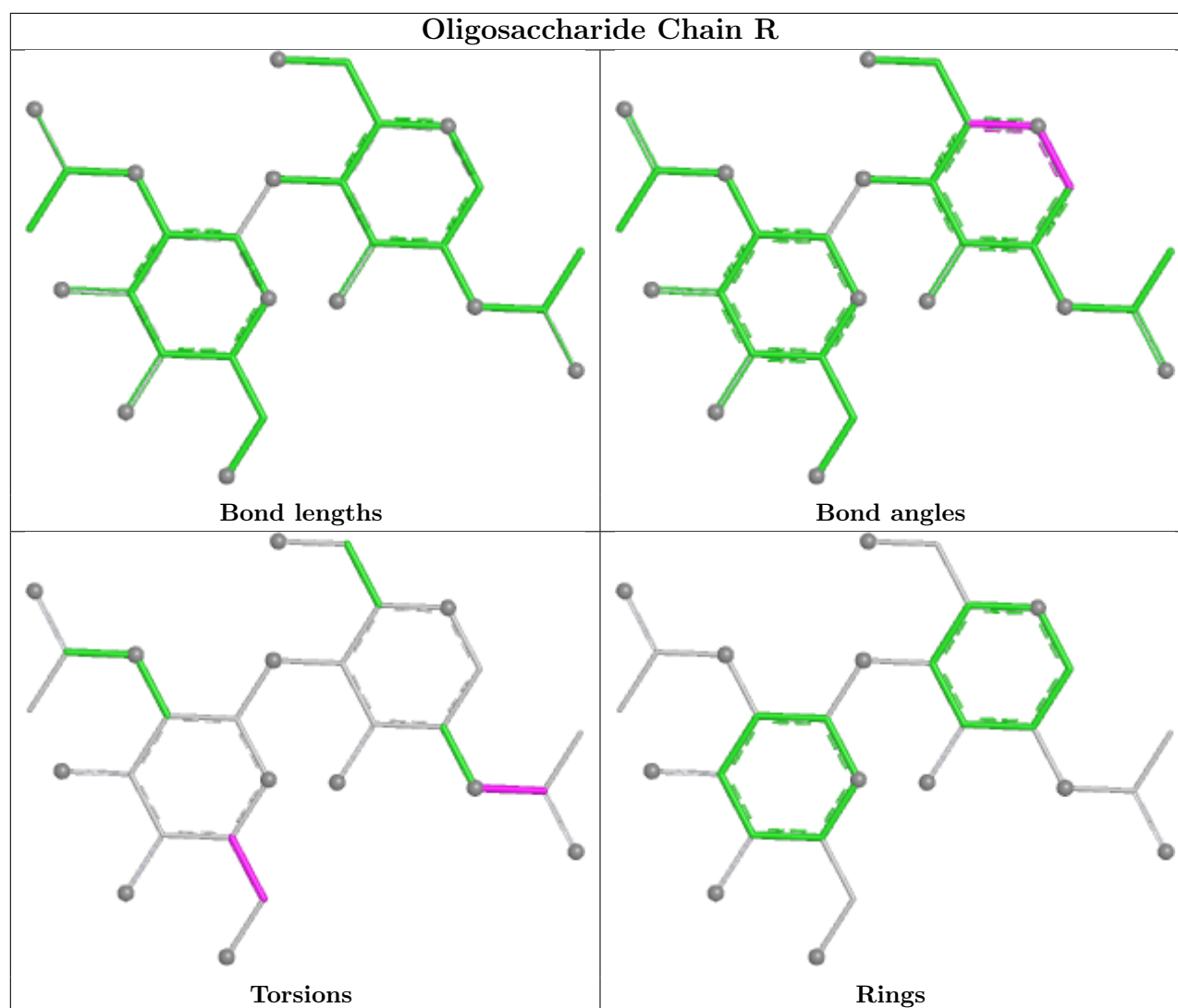


Rings









5.6 Ligand geometry [i](#)

3 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	$\# Z > 2$
16	NAG	G	1663	1	14,14,15	0.47	0	17,19,21	0.59	0
16	NAG	B	701	-	14,14,15	0.72	1 (7%)	17,19,21	0.39	0
16	NAG	B	702	2	14,14,15	2.48	2 (14%)	17,19,21	1.59	2 (11%)

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
16	NAG	G	1663	1	-	2/6/23/26	0/1/1/1
16	NAG	B	701	-	-	0/6/23/26	0/1/1/1
16	NAG	B	702	2	-	2/6/23/26	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
16	B	702	NAG	C1-C2	-7.49	1.42	1.52
16	B	702	NAG	O5-C1	5.04	1.52	1.43
16	B	701	NAG	O5-C1	-2.65	1.39	1.43

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
16	B	702	NAG	C1-O5-C5	5.17	119.11	112.19
16	B	702	NAG	O5-C5-C4	-2.10	105.72	110.83

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
16	B	702	NAG	C4-C5-C6-O6
16	B	702	NAG	O5-C5-C6-O6
16	G	1663	NAG	O5-C5-C6-O6
16	G	1663	NAG	C4-C5-C6-O6

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
16	G	1663	NAG	1	0
16	B	702	NAG	4	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	G	452/480 (94%)	-1.57	0 100 100	10, 33, 73, 120	0
2	B	148/153 (96%)	-1.54	0 100 100	9, 37, 109, 192	0
3	L	211/218 (96%)	-1.48	0 100 100	19, 44, 79, 104	0
4	H	231/236 (97%)	-1.51	0 100 100	32, 62, 91, 228	0
5	D	240/240 (100%)	-1.36	0 100 100	36, 97, 226, 263	0
6	E	213/216 (98%)	-1.36	0 100 100	55, 115, 172, 202	0
All	All	1495/1543 (96%)	-1.48	0 100 100	9, 55, 165, 263	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
7	NAG	A	1	14/15	-	-	142,152,164,166	0
7	NAG	A	2	14/15	-	-	171,175,181,186	0
7	BMA	A	3	11/12	-	-	150,162,167,168	0
7	MAN	A	4	11/12	-	-	136,139,144,148	0
7	MAN	A	5	11/12	-	-	133,137,149,150	0
7	MAN	A	6	11/12	-	-	160,167,172,173	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
8	NAG	C	1	14/15	-	-	134,135,138,145	0
8	NAG	C	2	14/15	-	-	135,141,146,148	0
8	BMA	C	3	11/12	-	-	127,140,157,159	0
8	MAN	C	4	11/12	-	-	170,172,183,187	0
8	MAN	C	5	11/12	-	-	165,172,184,185	0
8	MAN	C	6	11/12	-	-	184,191,198,199	0
8	MAN	C	7	11/12	-	-	117,120,133,134	0
9	NAG	F	1	14/15	-	-	122,130,140,142	0
9	NAG	F	2	14/15	-	-	146,153,166,167	0
9	BMA	F	3	11/12	-	-	160,165,168,172	0
9	MAN	F	4	11/12	-	-	136,155,161,171	0
9	MAN	F	5	11/12	-	-	108,116,127,131	0
9	MAN	F	6	11/12	-	-	98,102,105,106	0
9	MAN	F	7	11/12	-	-	160,171,179,190	0
11	MAN	M	6	11/12	0.96	0.04	246,261,264,268	0
12	MAN	N	6	11/12	0.96	0.05	223,225,234,237	0
14	MAN	P	5	11/12	0.96	0.04	209,221,224,226	0
11	NAG	J	2	14/15	0.97	0.04	182,193,199,201	0
10	NAG	S	2	14/15	0.97	0.06	191,196,215,219	0
14	MAN	P	6	11/12	0.97	0.07	201,208,212,213	0
14	MAN	P	7	11/12	0.97	0.06	262,267,270,273	0
15	NAG	Q	2	14/15	0.97	0.04	201,214,223,228	0
11	BMA	J	3	11/12	0.98	0.06	209,215,228,233	0
11	MAN	M	5	11/12	0.98	0.08	267,278,283,284	0
9	MAN	K	5	11/12	0.98	0.05	187,195,205,208	0
11	MAN	M	7	11/12	0.98	0.05	150,181,209,210	0
11	MAN	M	8	11/12	0.98	0.04	99,118,144,152	0
12	NAG	N	2	14/15	0.98	0.05	181,190,208,215	0
12	BMA	N	3	11/12	0.98	0.07	212,225,229,231	0
12	MAN	N	4	11/12	0.98	0.03	191,199,207,208	0
11	MAN	J	4	11/12	-	-	216,217,221,222	0
11	MAN	J	5	11/12	-	-	213,225,232,239	0
11	MAN	J	6	11/12	-	-	209,227,238,239	0
11	MAN	J	7	11/12	-	-	241,243,257,260	0
11	MAN	J	8	11/12	-	-	263,269,274,275	0
12	MAN	N	5	11/12	0.98	0.04	204,211,226,227	0
9	MAN	K	6	11/12	0.98	0.06	219,223,235,238	0
13	MAN	O	6	11/12	0.98	0.04	123,125,129,139	0
13	MAN	O	8	11/12	0.98	0.04	146,151,154,154	0
13	MAN	O	10	11/12	0.98	0.04	165,174,180,184	0
14	NAG	P	2	14/15	0.98	0.04	204,213,217,225	0
14	BMA	P	3	11/12	0.98	0.03	235,247,263,263	0

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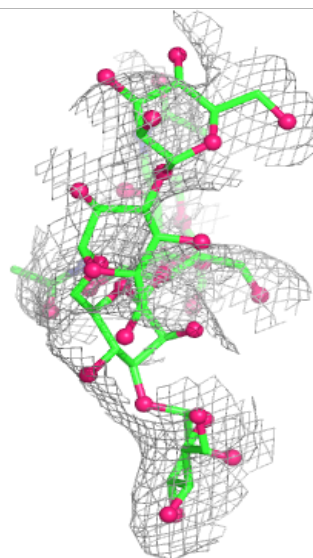
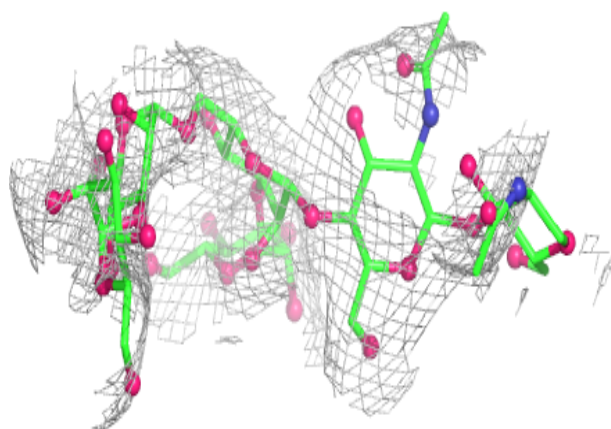
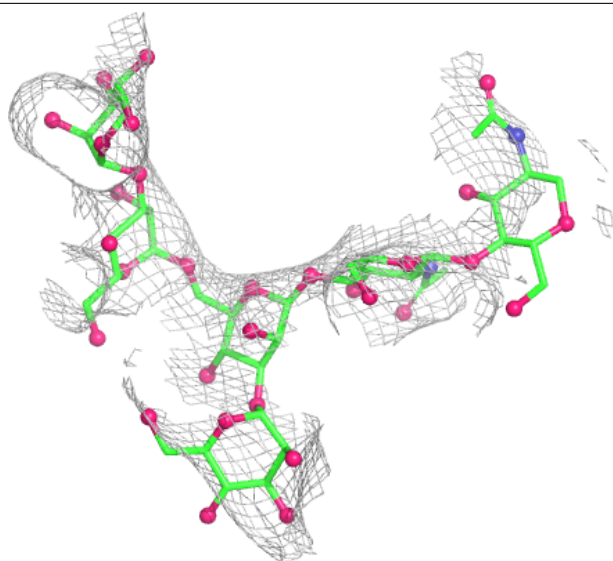
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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
14	MAN	P	4	11/12	0.98	0.03	228,240,254,255	0
9	MAN	K	7	11/12	0.98	0.04	233,243,248,257	0
9	NAG	K	2	14/15	0.98	0.07	231,240,257,259	0
11	NAG	J	1	14/15	0.98	0.10	165,173,182,184	0
9	MAN	K	4	11/12	0.98	0.03	202,211,230,239	0
11	NAG	M	1	14/15	0.99	0.06	181,187,211,221	0
11	NAG	M	2	14/15	0.99	0.05	177,196,201,205	0
13	NAG	O	1	14/15	0.99	0.04	132,137,140,141	0
13	NAG	O	2	14/15	0.99	0.08	127,131,150,172	0
13	BMA	O	3	11/12	0.99	0.04	119,126,138,145	0
13	MAN	O	4	11/12	0.99	0.03	101,102,106,109	0
13	MAN	O	5	11/12	0.99	0.03	103,107,113,114	0
11	BMA	M	3	11/12	0.99	0.04	214,223,228,232	0
13	MAN	O	7	11/12	0.99	0.04	149,154,159,163	0
11	MAN	M	4	11/12	0.99	0.05	243,244,255,265	0
13	MAN	O	9	11/12	0.99	0.04	152,157,169,174	0
10	BMA	I	3	11/12	0.99	0.03	263,271,276,280	0
14	NAG	P	1	14/15	0.99	0.04	178,191,199,199	0
10	NAG	S	1	14/15	0.99	0.04	143,170,181,191	0
9	NAG	K	1	14/15	0.99	0.04	168,198,226,229	0
10	BMA	S	3	11/12	0.99	0.05	181,186,191,191	0
12	NAG	N	1	14/15	0.99	0.06	133,150,161,169	0
9	BMA	K	3	11/12	0.99	0.02	231,234,240,242	0
10	NAG	I	1	14/15	0.99	0.04	203,222,253,261	0
15	NAG	Q	1	14/15	0.99	0.04	152,164,180,193	0
10	NAG	I	2	14/15	0.99	0.06	213,239,249,255	0
15	NAG	R	1	14/15	0.99	0.04	154,165,171,171	0
15	NAG	R	2	14/15	0.99	0.04	176,181,183,186	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.

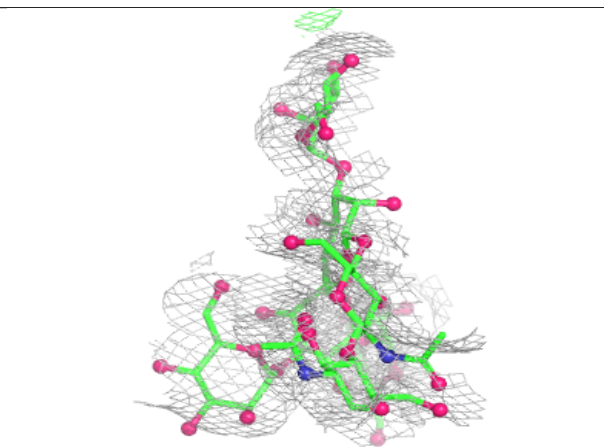
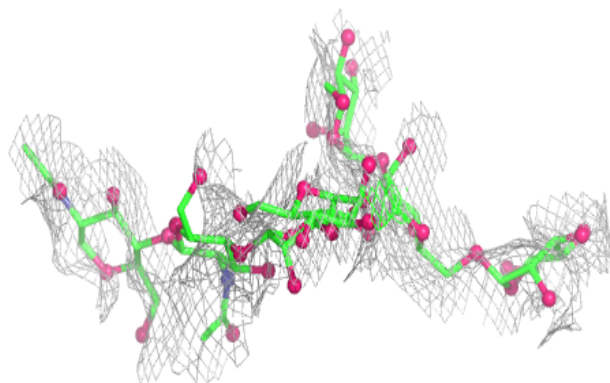
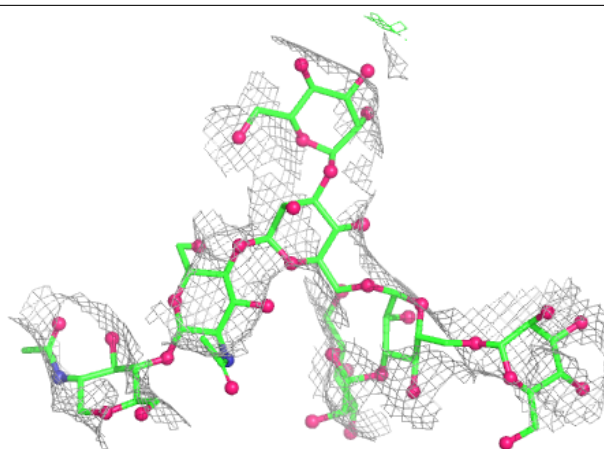
Electron density around Chain A:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

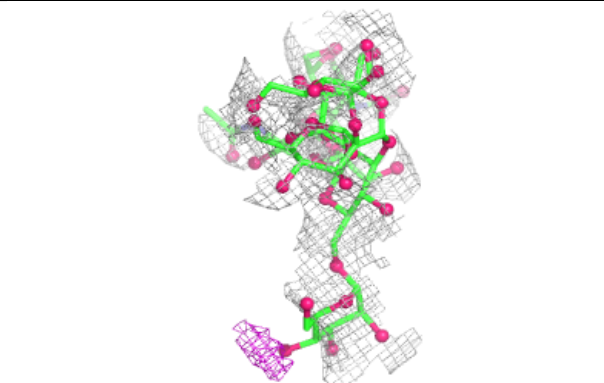
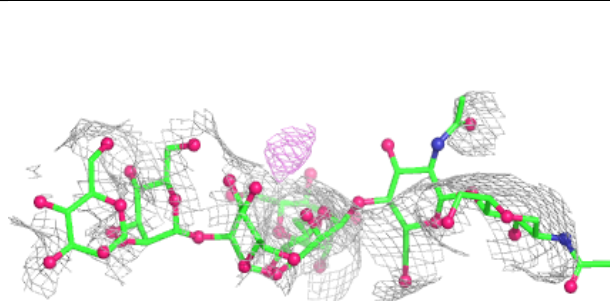
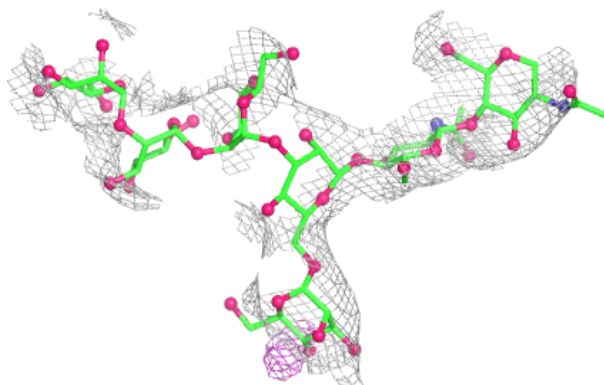


Electron density around Chain C:

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

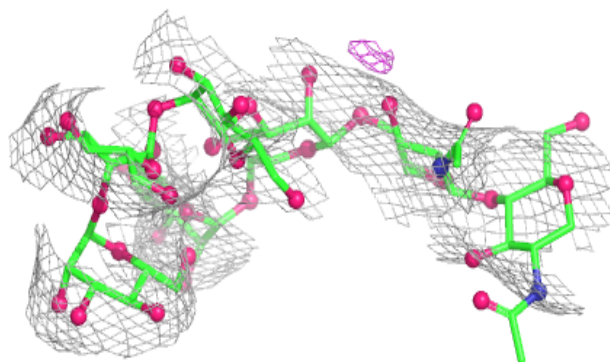
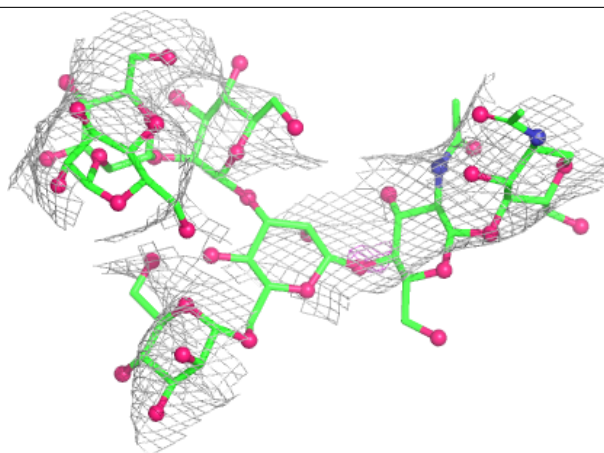
**Electron density around Chain F:**

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and green (positive)

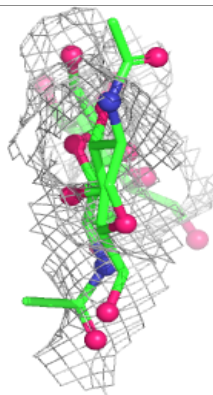
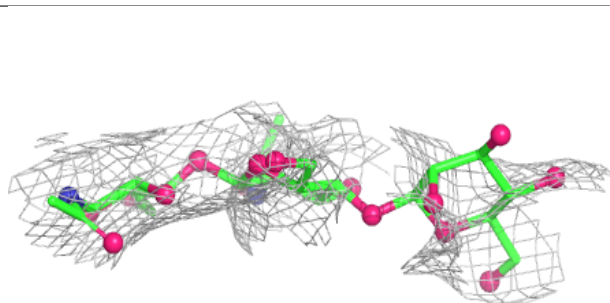
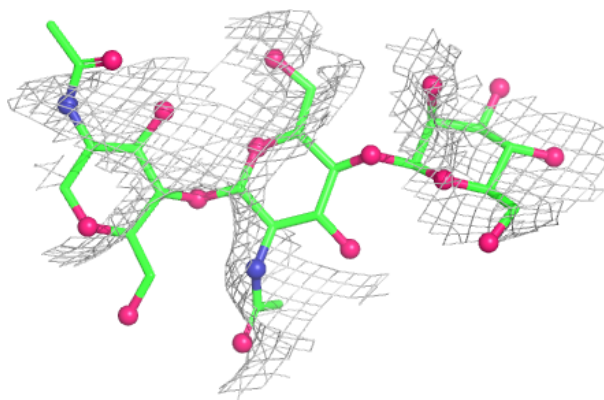


Electron density around Chain K:

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

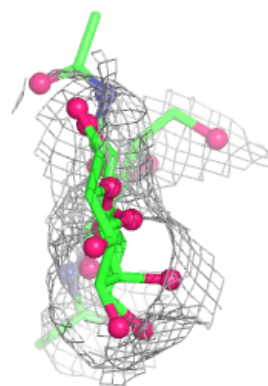
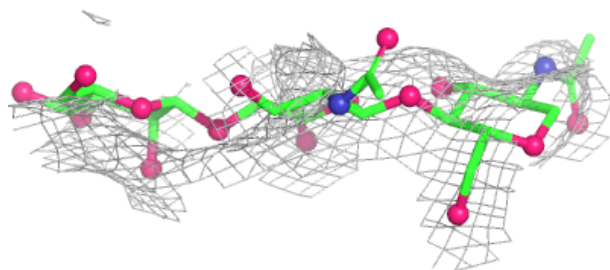
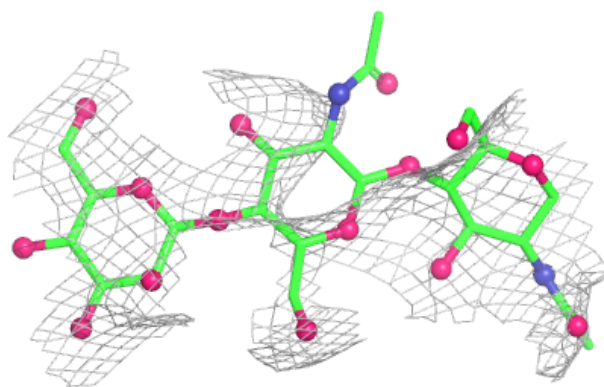
**Electron density around Chain I:**

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and green (positive)

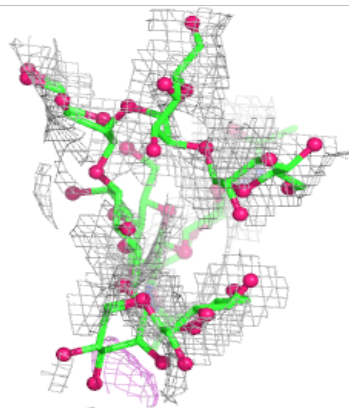
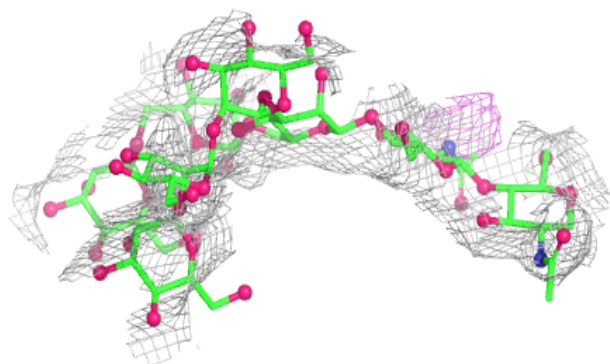
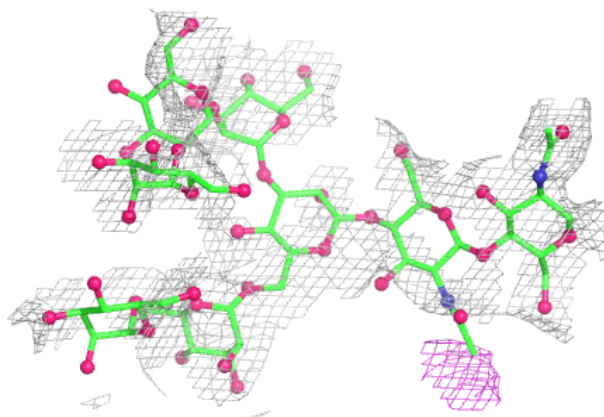


Electron density around Chain S:

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and green (positive)

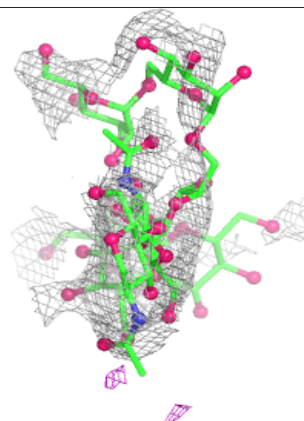
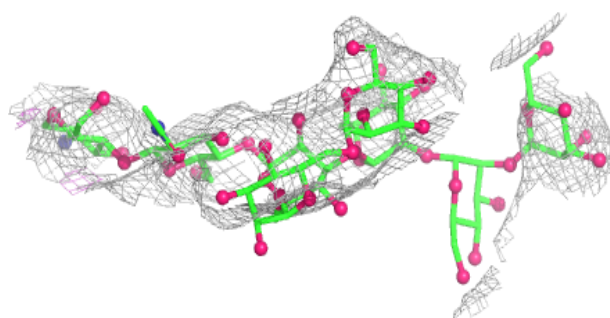
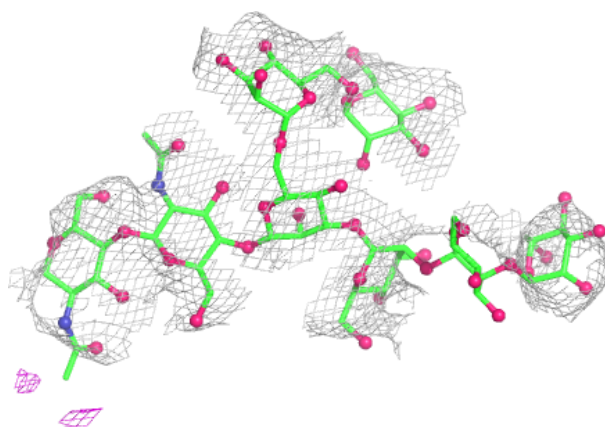
**Electron density around Chain J:**

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and green (positive)

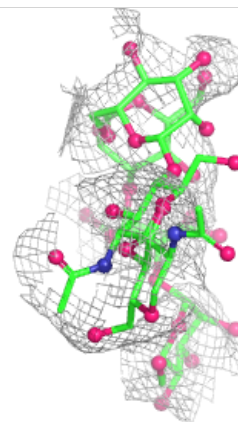
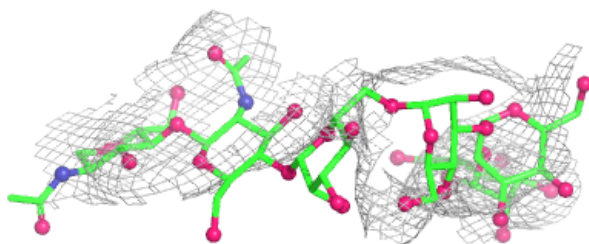
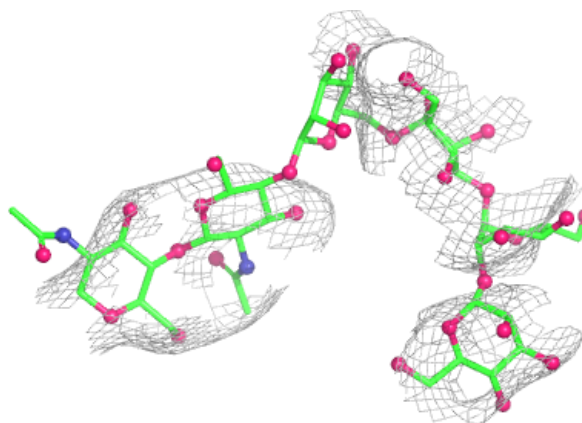


Electron density around Chain M:

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and green (positive)

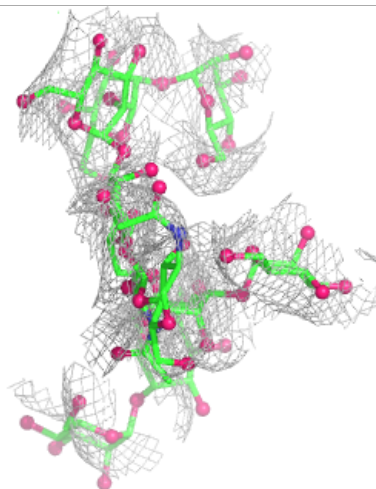
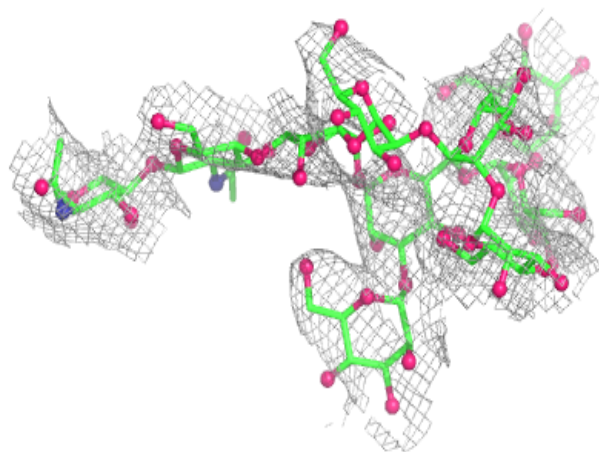
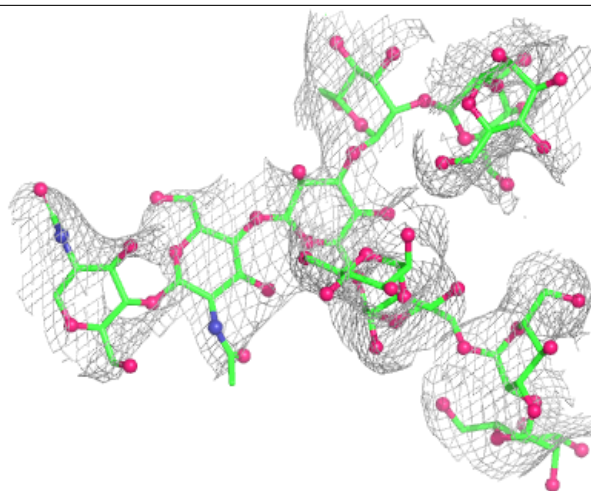
**Electron density around Chain N:**

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 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



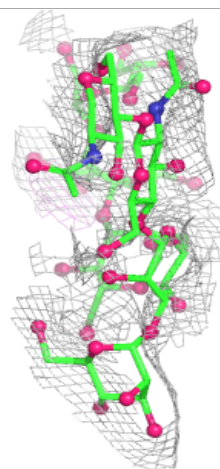
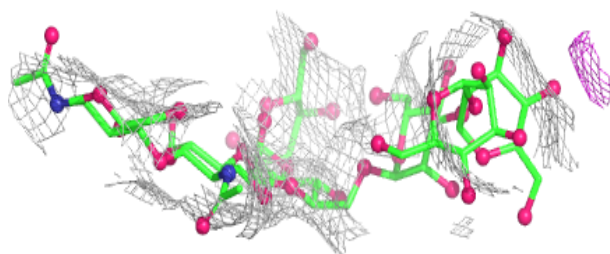
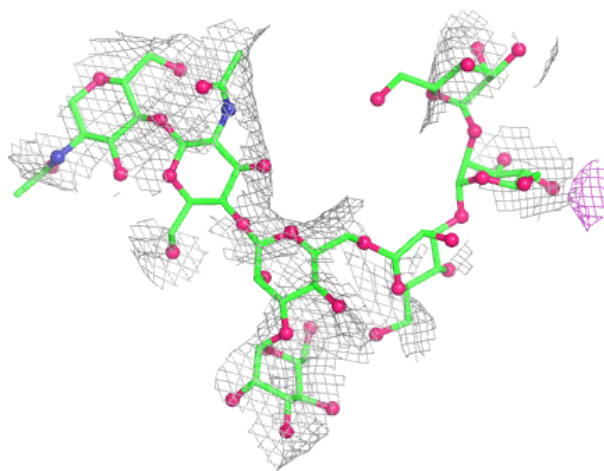
Electron density around Chain O:

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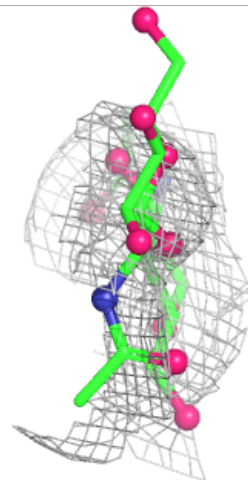
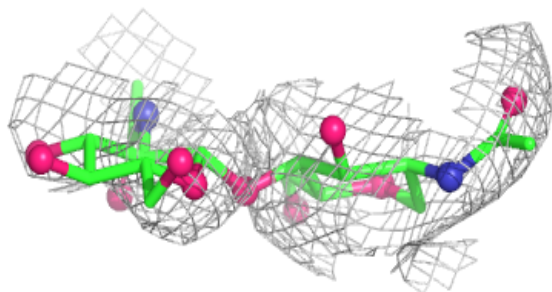
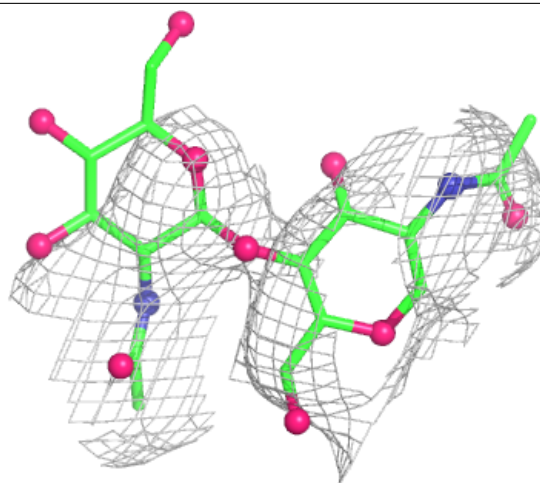
Electron density around Chain P:

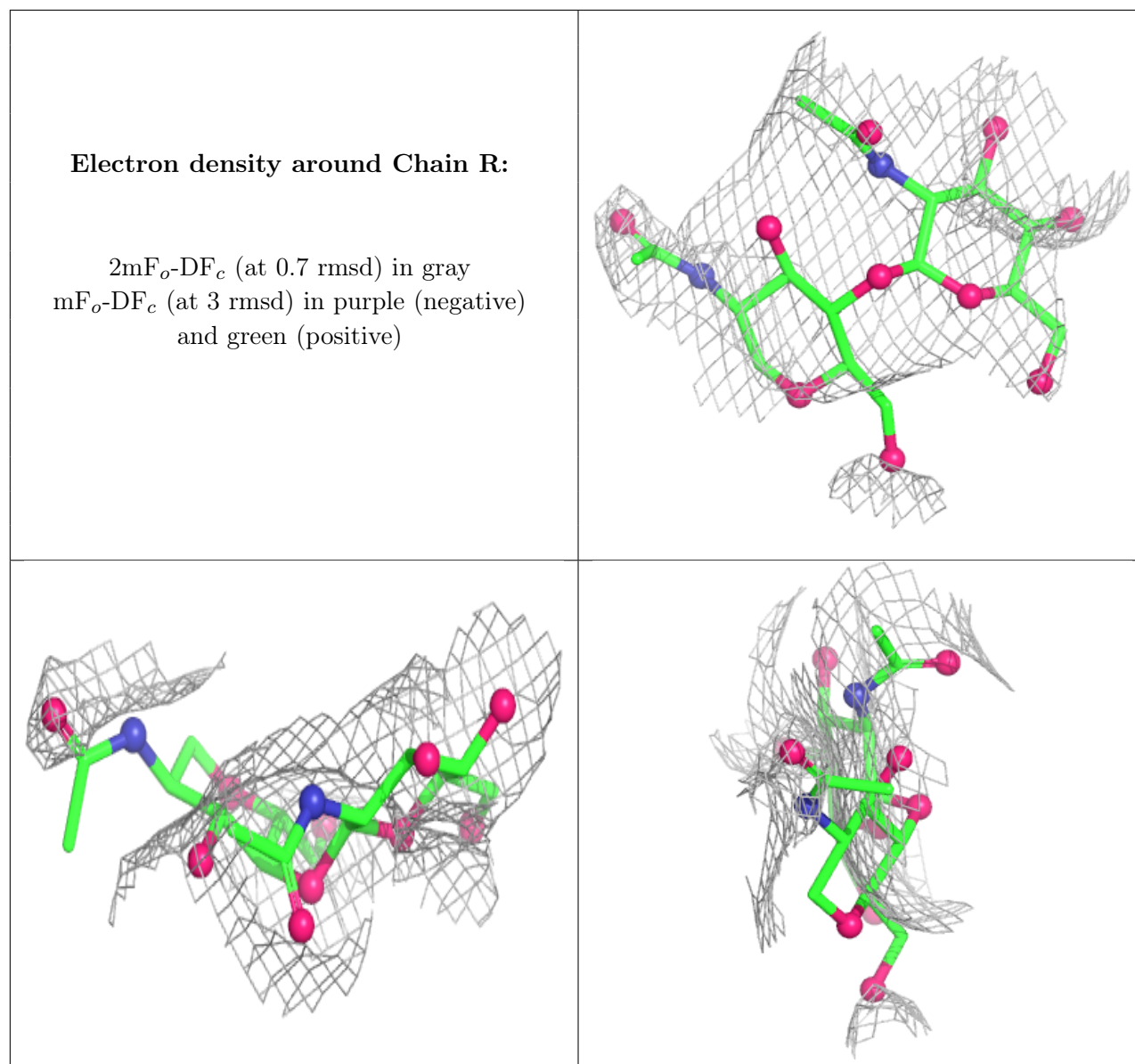
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around Chain Q:

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





6.4 Ligands [i](#)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
16	NAG	B	702	14/15	0.98	0.05	212,220,224,225	0
16	NAG	B	701	14/15	0.99	0.05	169,185,194,204	0
16	NAG	G	1663	14/15	0.99	0.04	134,144,154,161	0

6.5 Other polymers [i](#)

There are no such residues in this entry.