



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 10:23 AM UTC

PDB ID : 4JUL / pdb_00004jul
Title : Crystal structure of H5N1 influenza virus hemagglutinin, clade 2.3.4
Authors : DuBois, R.M.; Zaraket, H.; Reddivari, M.; Coop, T.; Heath, R.J.; White, S.W.; Russell, C.J.
Deposited on : 2013-03-25
Resolution : 2.79 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:

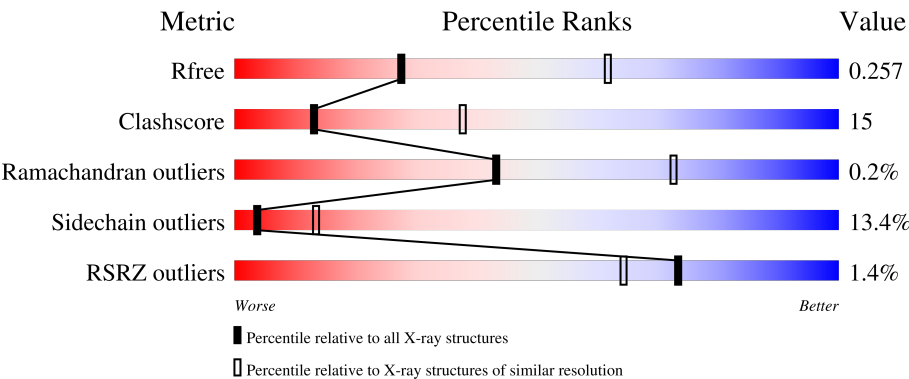
MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.79 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	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	329	67%27% . .
1	C	329	% 60%32%5% . .
1	E	329	63%30%5% .
1	H	329	% 65%27%5% .
1	J	329	% 65%29% . .

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Mol	Chain	Length	Quality of chain
1	L	329	
2	B	182	
2	D	182	
2	F	182	
2	I	182	
2	K	182	
2	M	182	
3	G	2	
3	N	2	
3	O	2	
3	P	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
4	NAG	J	401	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 23752 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 Hemagglutinin HA1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	321	Total	C	N	O	S	0	0	0
			2535	1601	436	485	13			
1	C	321	Total	C	N	O	S	0	0	0
			2535	1601	436	485	13			
1	E	322	Total	C	N	O	S	0	0	0
			2542	1606	437	486	13			
1	H	321	Total	C	N	O	S	0	0	0
			2535	1601	436	485	13			
1	J	321	Total	C	N	O	S	0	0	0
			2535	1601	436	485	13			
1	L	319	Total	C	N	O	S	0	0	0
			2523	1595	434	481	13			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	5	GLY	-	expression tag	UNP Q00G25
A	6	SER	-	expression tag	UNP Q00G25
A	7	ALA	-	expression tag	UNP Q00G25
A	8	ASP	-	expression tag	UNP Q00G25
A	9	PRO	-	expression tag	UNP Q00G25
A	10	GLY	-	expression tag	UNP Q00G25
C	5	GLY	-	expression tag	UNP Q00G25
C	6	SER	-	expression tag	UNP Q00G25
C	7	ALA	-	expression tag	UNP Q00G25
C	8	ASP	-	expression tag	UNP Q00G25
C	9	PRO	-	expression tag	UNP Q00G25
C	10	GLY	-	expression tag	UNP Q00G25
E	5	GLY	-	expression tag	UNP Q00G25
E	6	SER	-	expression tag	UNP Q00G25
E	7	ALA	-	expression tag	UNP Q00G25
E	8	ASP	-	expression tag	UNP Q00G25
E	9	PRO	-	expression tag	UNP Q00G25

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Chain	Residue	Modelled	Actual	Comment	Reference
E	10	GLY	-	expression tag	UNP Q00G25
H	5	GLY	-	expression tag	UNP Q00G25
H	6	SER	-	expression tag	UNP Q00G25
H	7	ALA	-	expression tag	UNP Q00G25
H	8	ASP	-	expression tag	UNP Q00G25
H	9	PRO	-	expression tag	UNP Q00G25
H	10	GLY	-	expression tag	UNP Q00G25
J	5	GLY	-	expression tag	UNP Q00G25
J	6	SER	-	expression tag	UNP Q00G25
J	7	ALA	-	expression tag	UNP Q00G25
J	8	ASP	-	expression tag	UNP Q00G25
J	9	PRO	-	expression tag	UNP Q00G25
J	10	GLY	-	expression tag	UNP Q00G25
L	5	GLY	-	expression tag	UNP Q00G25
L	6	SER	-	expression tag	UNP Q00G25
L	7	ALA	-	expression tag	UNP Q00G25
L	8	ASP	-	expression tag	UNP Q00G25
L	9	PRO	-	expression tag	UNP Q00G25
L	10	GLY	-	expression tag	UNP Q00G25

- Molecule 2 is a protein called Hemagglutinin HA2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	170	Total	C	N	O	S	0	0	0
			1380	859	241	272	8			
2	D	172	Total	C	N	O	S	0	0	0
			1398	869	243	278	8			
2	F	172	Total	C	N	O	S	0	0	0
			1398	869	243	278	8			
2	I	170	Total	C	N	O	S	0	0	0
			1380	859	241	272	8			
2	K	172	Total	C	N	O	S	0	0	0
			1398	869	243	278	8			
2	M	169	Total	C	N	O	S	0	0	0
			1369	853	237	271	8			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	177	ARG	-	expression tag	UNP Q00G25
B	178	SER	-	expression tag	UNP Q00G25
B	179	LEU	-	expression tag	UNP Q00G25

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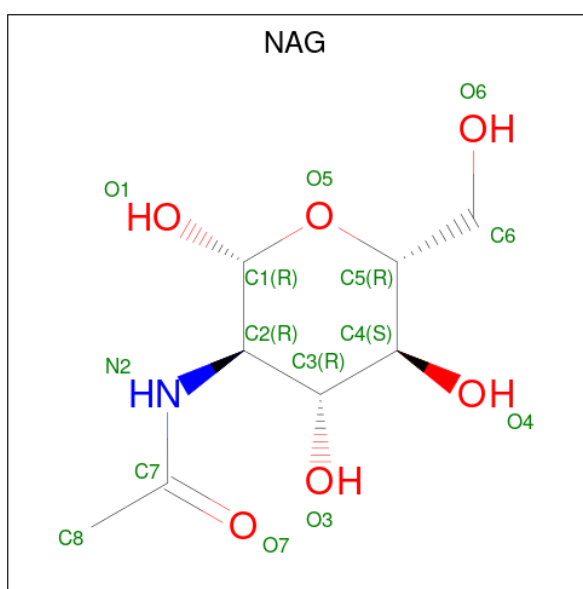
Chain	Residue	Modelled	Actual	Comment	Reference
B	180	VAL	-	expression tag	UNP Q00G25
B	181	PRO	-	expression tag	UNP Q00G25
B	182	ARG	-	expression tag	UNP Q00G25
D	177	ARG	-	expression tag	UNP Q00G25
D	178	SER	-	expression tag	UNP Q00G25
D	179	LEU	-	expression tag	UNP Q00G25
D	180	VAL	-	expression tag	UNP Q00G25
D	181	PRO	-	expression tag	UNP Q00G25
D	182	ARG	-	expression tag	UNP Q00G25
F	177	ARG	-	expression tag	UNP Q00G25
F	178	SER	-	expression tag	UNP Q00G25
F	179	LEU	-	expression tag	UNP Q00G25
F	180	VAL	-	expression tag	UNP Q00G25
F	181	PRO	-	expression tag	UNP Q00G25
F	182	ARG	-	expression tag	UNP Q00G25
I	177	ARG	-	expression tag	UNP Q00G25
I	178	SER	-	expression tag	UNP Q00G25
I	179	LEU	-	expression tag	UNP Q00G25
I	180	VAL	-	expression tag	UNP Q00G25
I	181	PRO	-	expression tag	UNP Q00G25
I	182	ARG	-	expression tag	UNP Q00G25
K	177	ARG	-	expression tag	UNP Q00G25
K	178	SER	-	expression tag	UNP Q00G25
K	179	LEU	-	expression tag	UNP Q00G25
K	180	VAL	-	expression tag	UNP Q00G25
K	181	PRO	-	expression tag	UNP Q00G25
K	182	ARG	-	expression tag	UNP Q00G25
M	177	ARG	-	expression tag	UNP Q00G25
M	178	SER	-	expression tag	UNP Q00G25
M	179	LEU	-	expression tag	UNP Q00G25
M	180	VAL	-	expression tag	UNP Q00G25
M	181	PRO	-	expression tag	UNP Q00G25
M	182	ARG	-	expression tag	UNP Q00G25

- Molecule 3 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
3	G	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	N	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	O	2	Total	C	N	O	0	0	0
			28	16	2	10			
3	P	2	Total	C	N	O	0	0	0
			28	16	2	10			

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	A	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	C	1	Total	C	N	O	0	0
			14	8	1	5		
4	E	1	Total	C	N	O	0	0
			14	8	1	5		
4	H	1	Total	C	N	O	0	0
			14	8	1	5		
4	J	1	Total	C	N	O	0	0
			14	8	1	5		
4	J	1	Total	C	N	O	0	0
			14	8	1	5		

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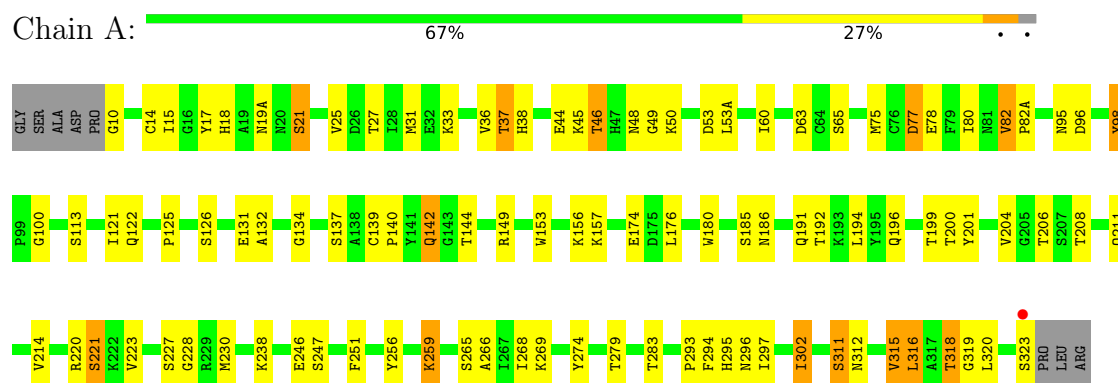
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	L	1	Total	C	N	O	0	0
			14	8	1	5		

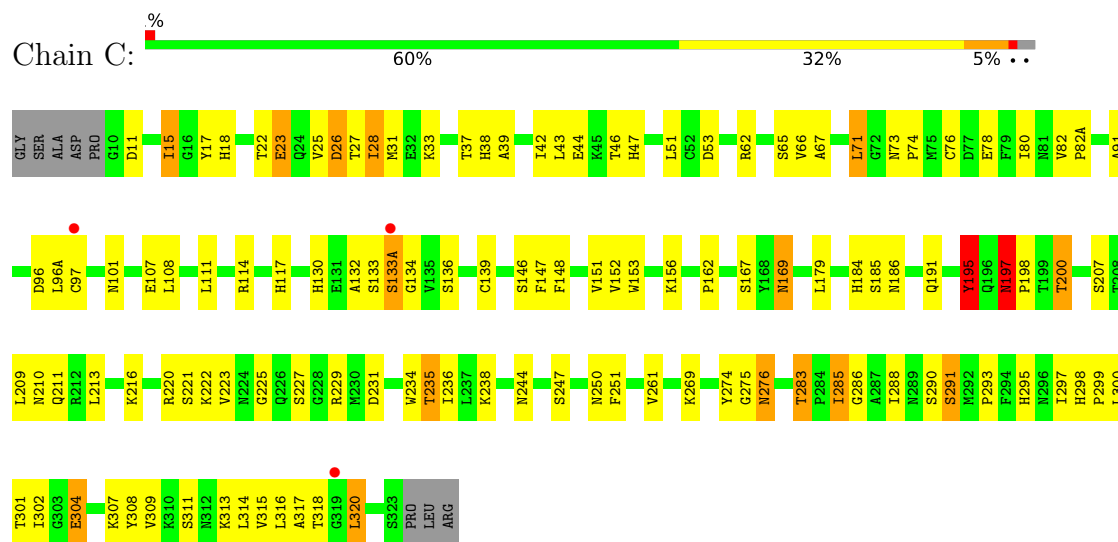
3 Residue-property plots

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.

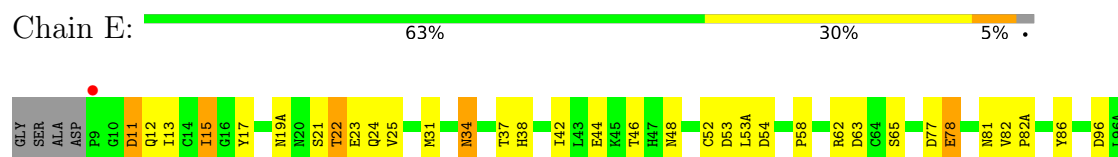
• Molecule 1: Hemagglutinin HA1

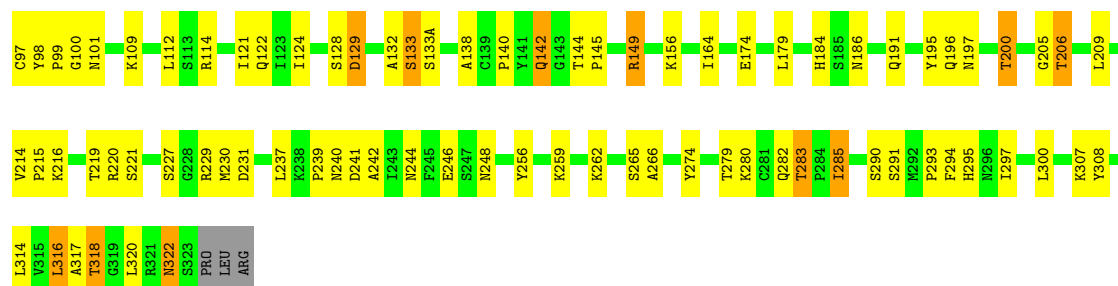


• Molecule 1: Hemagglutinin HA1

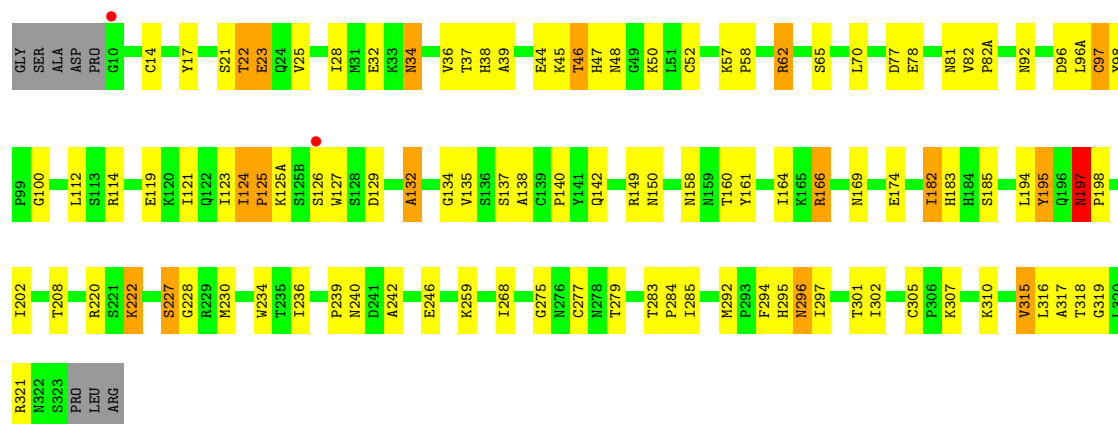


• Molecule 1: Hemagglutinin HA1

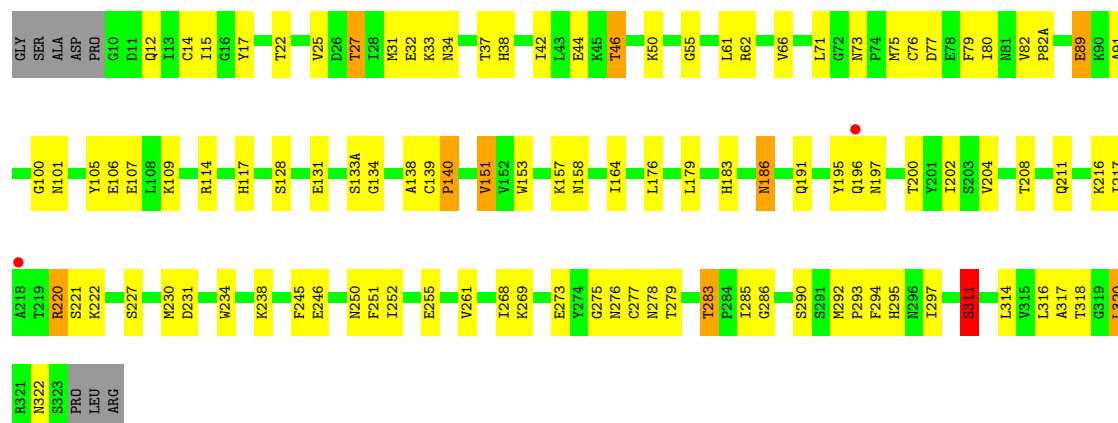




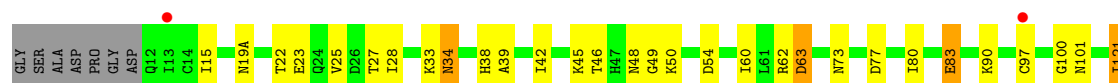
• Molecule 1: Hemagglutinin HA1

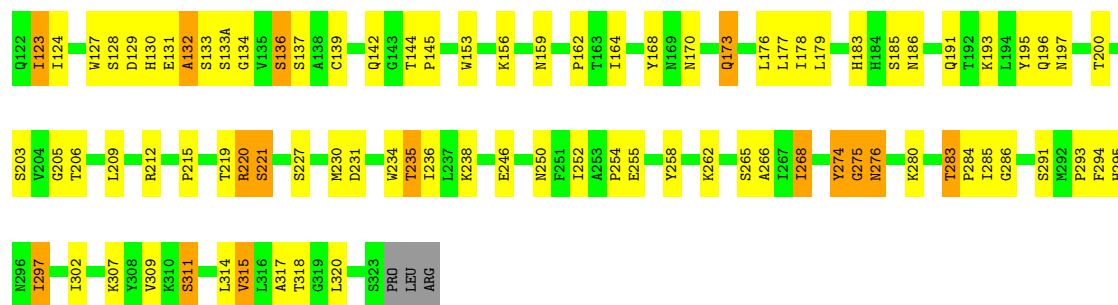


• Molecule 1: Hemagglutinin HA1

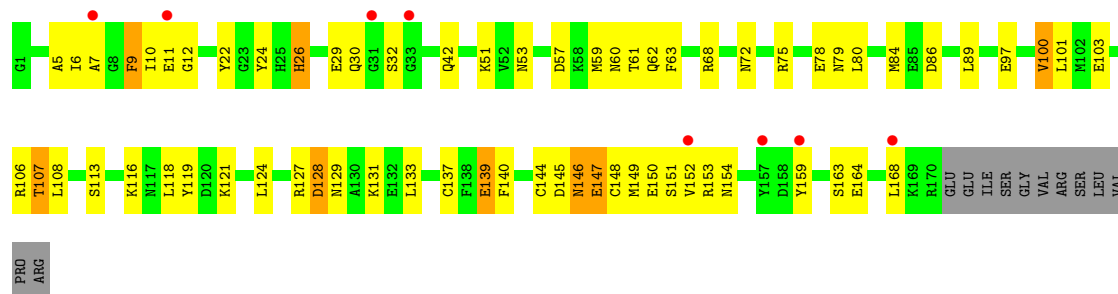


• Molecule 1: Hemagglutinin HA1

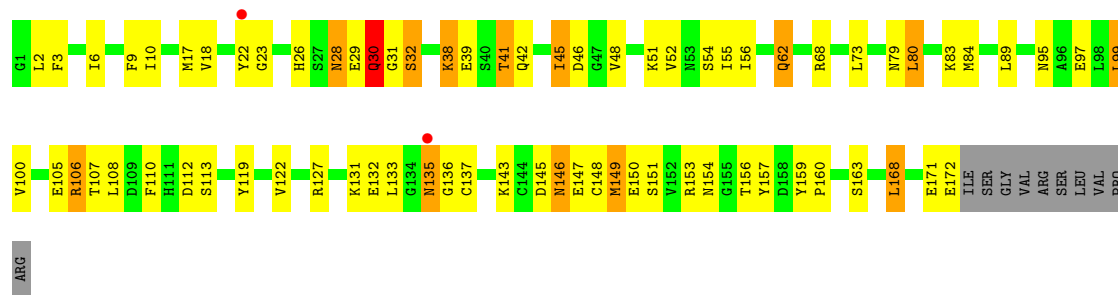




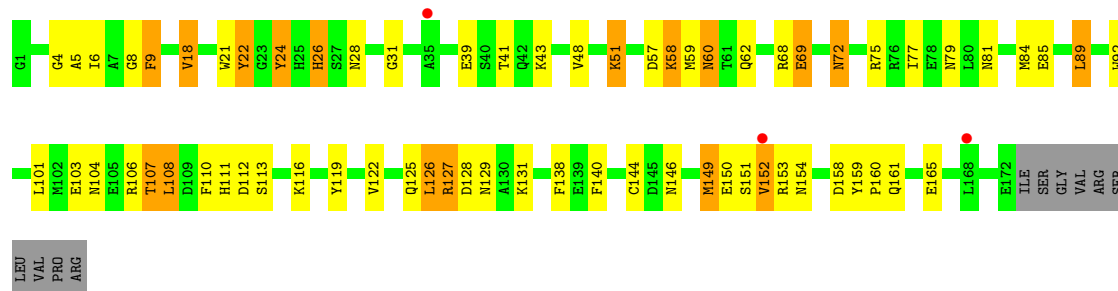
• Molecule 2: Hemagglutinin HA2



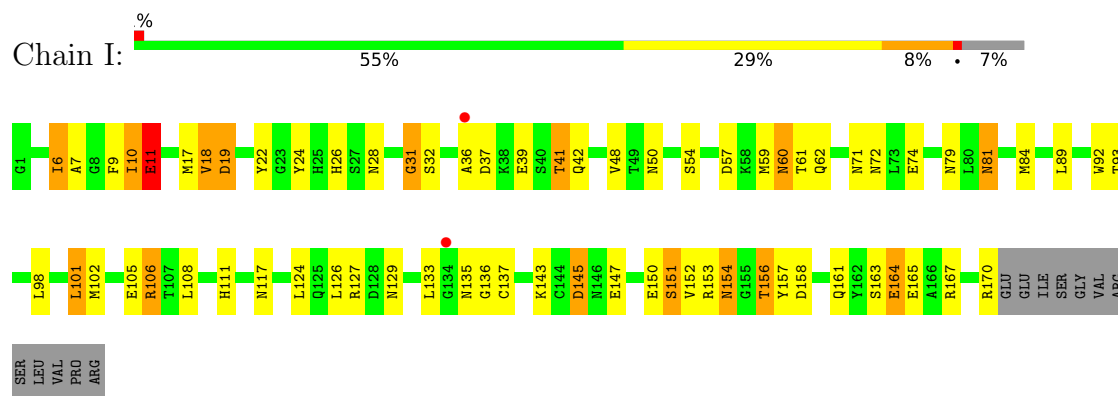
• Molecule 2: Hemagglutinin HA2



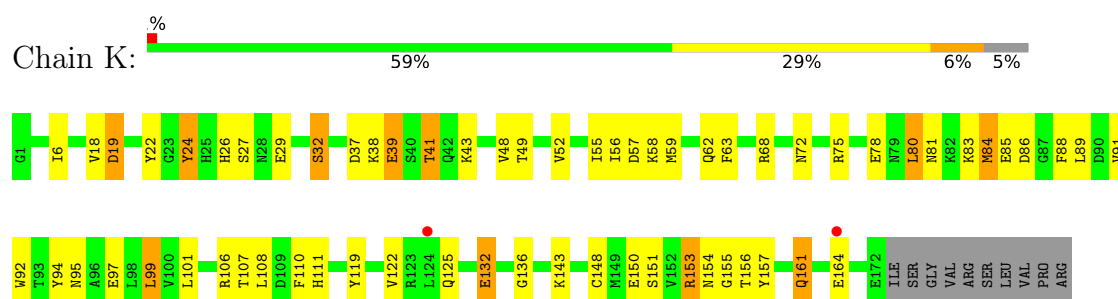
• Molecule 2: Hemagglutinin HA2



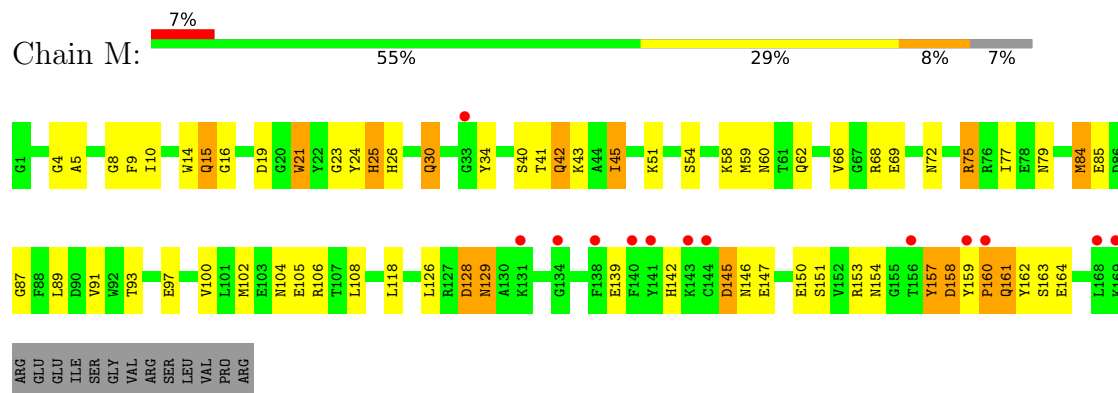
- Molecule 2: Hemagglutinin HA2



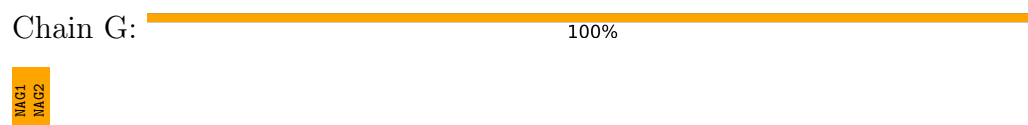
- Molecule 2: Hemagglutinin HA2



- Molecule 2: Hemagglutinin HA2



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



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



MAG1
MAG2

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

Chain O:  50% 50%MAG1
MAG2

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

Chain P:  100%MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	71.58Å 71.56Å 247.05Å 83.94° 85.46° 60.68°	Depositor
Resolution (Å)	48.87 – 2.79 48.87 – 2.79	Depositor EDS
% Data completeness (in resolution range)	81.4 (48.87-2.79) 81.4 (48.87-2.79)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.08 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.196 , 0.254 0.204 , 0.257	Depositor DCC
R_{free} test set	4355 reflections (4.13%)	wwPDB-VP
Wilson B-factor (Å ²)	46.5	Xtriage
Anisotropy	0.379	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 35.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.018 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	23752	wwPDB-VP
Average B, all atoms (Å ²)	45.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.43% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality ⓘ

5.1 Standard geometry ⓘ

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

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.66	0/2596	1.02	6/3529 (0.2%)
1	C	0.71	0/2596	1.05	13/3529 (0.4%)
1	E	0.66	0/2604	1.05	9/3540 (0.3%)
1	H	0.68	1/2596 (0.0%)	1.01	5/3529 (0.1%)
1	J	0.69	0/2596	1.03	10/3529 (0.3%)
1	L	0.65	0/2584	1.04	12/3513 (0.3%)
2	B	0.62	0/1407	1.00	3/1891 (0.2%)
2	D	0.64	0/1425	1.09	7/1915 (0.4%)
2	F	0.59	0/1425	0.97	3/1915 (0.2%)
2	I	0.66	1/1407 (0.1%)	1.08	9/1891 (0.5%)
2	K	0.63	0/1425	0.95	2/1915 (0.1%)
2	M	0.61	0/1396	1.04	5/1877 (0.3%)
All	All	0.66	2/24057 (0.0%)	1.03	84/32573 (0.3%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	18	VAL	CA-CB	-5.88	1.46	1.54
1	H	125	PRO	N-CD	-5.51	1.40	1.47

All (84) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	30	GLN	N-CA-C	-14.00	96.02	111.28
2	B	146	ASN	N-CA-C	11.30	123.60	111.28
2	I	11	GLU	N-CA-C	10.84	123.09	111.28
2	K	18	VAL	N-CA-C	9.76	122.57	113.10
2	M	5	ALA	N-CA-C	9.54	123.55	111.24
1	A	132	ALA	N-CA-C	9.15	121.34	111.36
2	F	9	PHE	N-CA-C	-8.41	102.11	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	46	THR	N-CA-C	8.32	122.97	109.40
2	I	19	ASP	N-CA-C	8.23	120.33	111.36
2	I	167	ARG	N-CA-C	-8.03	102.53	111.28
2	D	28	ASN	N-CA-C	-8.00	97.86	108.34
1	J	290	SER	N-CA-C	7.68	120.19	108.52
1	A	82	VAL	N-CA-C	7.55	116.04	109.02
2	D	146	ASN	N-CA-C	7.50	120.11	111.11
1	E	132	ALA	N-CA-C	7.50	119.53	111.36
1	C	133(A)	SER	N-CA-C	7.42	121.91	111.52
2	I	31	GLY	N-CA-C	7.40	121.22	111.52
2	I	6	ILE	CB-CA-C	-7.39	101.94	110.73
2	K	56	ILE	N-CA-C	7.30	118.09	110.72
1	H	124	ILE	CB-CA-C	-7.17	104.27	111.08
1	L	133	SER	N-CA-C	7.07	121.93	113.16
1	L	132	ALA	N-CA-C	7.00	118.99	111.36
1	L	19(A)	ASN	N-CA-C	6.80	117.21	108.24
2	I	136	GLY	N-CA-C	-6.66	106.26	114.92
1	J	322	ASN	N-CA-C	6.60	119.20	109.24
2	I	151	SER	N-CA-C	6.51	119.70	111.69
1	L	73	ASN	CA-C-N	6.44	126.66	119.32
1	L	73	ASN	C-N-CA	6.44	126.66	119.32
2	D	32	SER	N-CA-C	6.36	119.38	109.52
1	J	311	SER	N-CA-C	6.33	119.09	110.55
2	D	136	GLY	N-CA-C	-6.30	106.21	115.72
2	M	158	ASP	N-CA-C	6.25	118.60	108.41
2	B	10	ILE	N-CA-C	-6.17	98.74	107.75
1	E	12	GLN	N-CA-C	6.16	119.44	109.40
2	M	160	PRO	N-CA-C	-6.13	102.92	113.70
1	L	276	ASN	N-CA-C	6.06	120.62	113.23
1	C	197	ASN	CA-C-N	6.06	125.61	119.19
1	C	197	ASN	C-N-CA	6.06	125.61	119.19
1	J	255	GLU	N-CA-C	-6.02	105.02	112.72
1	L	34	ASN	N-CA-C	6.02	119.69	111.39
1	H	132	ALA	N-CA-C	6.01	117.92	111.36
1	E	133	SER	N-CA-C	6.00	120.30	112.92
1	E	34	ASN	N-CA-C	5.99	117.92	110.91
1	C	132	ALA	N-CA-C	5.98	118.75	111.82
1	H	197	ASN	CA-C-N	5.97	125.52	119.19
1	H	197	ASN	C-N-CA	5.97	125.52	119.19
1	E	191	GLN	N-CA-C	-5.88	104.77	111.07
2	B	147	GLU	N-CA-C	-5.85	104.98	111.71
1	E	52	CYS	N-CA-C	5.77	117.73	109.14

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	I	39	GLU	N-CA-C	5.73	117.99	111.11
2	M	40	SER	N-CA-C	-5.72	105.12	111.36
1	A	98	TYR	N-CA-C	-5.64	101.40	109.24
2	D	163	SER	N-CA-C	5.64	117.42	111.28
2	F	41	THR	N-CA-C	-5.61	105.93	112.89
1	C	195	TYR	N-CA-C	-5.60	100.84	109.52
1	L	275	GLY	N-CA-C	5.59	125.34	115.61
2	F	127	ARG	CB-CA-C	-5.48	110.27	116.63
1	C	152	VAL	N-CA-C	5.47	115.83	108.17
1	A	223	VAL	N-CA-C	-5.43	100.36	108.46
2	D	18	VAL	N-CA-C	5.43	118.42	113.20
1	L	274	TYR	N-CA-C	5.41	118.19	110.24
1	J	62	ARG	CB-CA-C	-5.39	109.90	117.23
2	M	23	GLY	N-CA-C	5.39	118.12	110.74
1	J	12	GLN	N-CA-C	5.36	117.81	108.76
1	L	297	ILE	N-CA-C	5.34	115.48	110.30
1	E	124	ILE	CA-C-N	5.33	125.64	119.93
1	E	124	ILE	C-N-CA	5.33	125.64	119.93
1	C	91	ALA	N-CA-C	-5.31	105.66	111.82
1	H	34	ASN	N-CA-C	5.30	118.56	111.24
1	C	213	LEU	CA-C-N	-5.30	118.72	123.33
1	C	213	LEU	C-N-CA	-5.30	118.72	123.33
2	I	10	ILE	N-CA-C	-5.30	98.33	109.34
1	J	140	PRO	N-CA-C	5.30	119.50	111.19
1	C	320	LEU	N-CA-C	5.24	117.45	110.53
1	J	292	MET	CA-C-N	-5.21	113.33	119.84
1	J	292	MET	C-N-CA	-5.21	113.33	119.84
1	L	63	ASP	N-CA-C	5.19	119.25	113.02
1	L	121	ILE	N-CA-C	5.17	115.71	108.17
1	A	113	SER	N-CA-C	-5.15	106.97	113.20
1	A	206	THR	N-CA-C	-5.10	102.03	108.45
1	E	262	LYS	N-CA-C	5.08	116.86	108.13
1	C	238	LYS	CA-C-N	-5.05	114.59	119.90
1	C	238	LYS	C-N-CA	-5.05	114.59	119.90
1	C	288	ILE	N-CA-C	5.05	115.24	107.77

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2535	0	2468	68	0
1	C	2535	0	2469	76	0
1	E	2542	0	2478	85	0
1	H	2535	0	2469	70	0
1	J	2535	0	2469	53	0
1	L	2523	0	2462	77	0
2	B	1380	0	1291	55	0
2	D	1398	0	1303	67	0
2	F	1398	0	1303	61	0
2	I	1380	0	1291	50	0
2	K	1398	0	1303	47	0
2	M	1369	0	1278	66	0
3	G	28	0	25	2	0
3	N	28	0	25	5	0
3	O	28	0	25	2	0
3	P	28	0	26	2	0
4	A	14	0	13	2	0
4	C	28	0	26	0	0
4	E	14	0	13	4	0
4	H	14	0	13	1	0
4	J	28	0	26	8	0
4	L	14	0	13	2	0
All	All	23752	0	22789	690	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 15.

All (690) 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:H:34:ASN:HD21	4:H:401:NAG:C1	1.04	1.59
1:E:34:ASN:HD21	4:E:401:NAG:C1	0.95	1.58
1:J:34:ASN:HD21	4:J:401:NAG:C1	1.14	1.57
1:E:53(A):LEU:O	1:E:53(A):LEU:HD23	1.44	1.15
2:M:158:ASP:O	2:M:161:GLN:HB3	1.51	1.10

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:242:ALA:HB3	3:N:1:NAG:H82	1.24	1.08
2:M:157:TYR:C	2:M:157:TYR:CD2	2.29	1.07
1:E:34:ASN:HD21	4:E:401:NAG:C2	1.72	1.02
1:A:142:GLN:HA	1:A:142:GLN:NE2	1.76	0.98
2:M:161:GLN:HG3	2:M:162:TYR:CD2	2.01	0.95
1:A:142:GLN:HA	1:A:142:GLN:HE21	1.31	0.95
1:A:49:GLY:O	1:A:50:LYS:HD3	1.69	0.93
1:J:33:LYS:HB2	4:J:401:NAG:H82	1.51	0.92
1:C:27:THR:HG22	1:C:31:MET:H	1.31	0.92
2:M:157:TYR:C	2:M:157:TYR:HD2	1.75	0.92
1:A:17:TYR:CZ	2:B:6:ILE:HG23	2.06	0.91
1:E:242:ALA:CB	3:N:1:NAG:H82	2.03	0.89
2:M:161:GLN:HG3	2:M:162:TYR:HD2	1.34	0.88
2:M:157:TYR:HD2	2:M:157:TYR:O	1.57	0.88
1:C:275:GLY:O	1:C:276:ASN:HB2	1.73	0.86
2:K:27:SER:CB	2:K:32:SER:OG	2.24	0.86
2:F:26:HIS:HD2	2:F:26:HIS:O	1.56	0.85
2:D:68:ARG:HE	2:F:79:ASN:HD22	1.22	0.85
2:F:26:HIS:O	2:F:26:HIS:CD2	2.30	0.85
2:D:9:PHE:CE1	2:D:10:ILE:HG13	2.12	0.84
2:M:26:HIS:O	2:M:26:HIS:CD2	2.30	0.84
2:B:26:HIS:O	2:B:26:HIS:CD2	2.30	0.84
1:E:206:THR:HG23	1:E:209:LEU:HB3	1.60	0.84
1:H:295:HIS:HD2	1:H:297:ILE:H	1.25	0.82
1:H:57:LYS:HG2	1:H:58:PRO:HD2	1.61	0.82
1:C:18:HIS:CD2	2:D:17:MET:O	2.33	0.82
1:A:75:MET:HB2	1:A:95:ASN:ND2	1.96	0.81
2:F:126:LEU:N	2:F:126:LEU:HD12	1.94	0.81
2:F:122:VAL:O	2:F:126:LEU:HD13	1.79	0.81
2:I:7:ALA:HB1	2:I:11:GLU:O	1.79	0.81
3:P:1:NAG:C4	3:P:2:NAG:C1	2.59	0.80
1:E:15:ILE:HD13	2:F:119:TYR:HA	1.63	0.80
1:H:242:ALA:H	3:O:1:NAG:H82	1.47	0.80
1:A:98:TYR:CE2	1:A:230:MET:HG3	2.16	0.80
1:E:53(A):LEU:O	1:E:53(A):LEU:CD2	2.30	0.79
2:D:145:ASP:O	2:D:148:CYS:HB3	1.84	0.78
2:I:7:ALA:HB1	2:I:11:GLU:HA	1.65	0.78
1:E:34:ASN:CG	4:E:401:NAG:C1	2.57	0.78
1:E:317:ALA:H	2:F:104:ASN:ND2	1.82	0.78
1:J:34:ASN:HD22	4:J:401:NAG:C1	1.98	0.77
2:D:80:LEU:C	2:D:80:LEU:HD12	2.11	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:26:HIS:O	2:B:26:HIS:HD2	1.67	0.75
2:D:127:ARG:HH12	2:F:131:LYS:HE2	1.51	0.75
1:H:98:TYR:CD2	1:H:230:MET:HG3	2.21	0.75
2:M:128:ASP:C	2:M:128:ASP:OD1	2.29	0.75
1:C:216:LYS:O	1:C:220:ARG:NH2	2.20	0.75
2:M:84:MET:HE3	2:M:85:GLU:HG3	1.69	0.75
2:F:4:GLY:O	2:F:8:GLY:HA3	1.85	0.75
1:J:25:VAL:HG21	1:J:317:ALA:HB2	1.69	0.75
2:M:26:HIS:ND1	2:M:153:ARG:NH2	2.34	0.74
2:M:159:TYR:C	2:M:161:GLN:H	1.96	0.74
2:F:28:ASN:ND2	2:F:146:ASN:OD1	2.21	0.74
1:A:295:HIS:HD2	1:A:297:ILE:H	1.35	0.74
2:I:18:VAL:HG12	2:I:18:VAL:O	1.87	0.74
1:H:25:VAL:HG21	1:H:317:ALA:HB2	1.68	0.74
1:J:275:GLY:O	1:J:276:ASN:HB2	1.87	0.74
1:H:283:THR:HG22	1:H:301:THR:HG22	1.68	0.73
1:C:169:ASN:OD1	1:C:169:ASN:C	2.30	0.73
1:C:311:SER:HB3	2:D:97:GLU:OE2	1.89	0.73
1:C:283:THR:HG22	1:C:285:ILE:H	1.52	0.73
2:F:126:LEU:N	2:F:126:LEU:CD1	2.52	0.72
1:J:33:LYS:CB	4:J:401:NAG:H82	2.20	0.72
2:F:125:GLN:C	2:F:126:LEU:HD12	2.13	0.72
2:M:142:HIS:CE1	2:M:162:TYR:CD1	2.78	0.72
1:E:11:ASP:O	2:F:140:PHE:HB2	1.90	0.72
1:L:159:ASN:O	1:L:196:GLN:NE2	2.21	0.72
1:E:295:HIS:HD2	1:E:297:ILE:H	1.37	0.72
1:E:317:ALA:H	2:F:104:ASN:HD21	1.36	0.72
2:B:147:GLU:O	2:B:151:SER:HB3	1.90	0.71
1:H:307:LYS:HE2	2:I:60:ASN:HA	1.72	0.71
1:C:51:LEU:O	1:C:275:GLY:N	2.22	0.71
1:C:191:GLN:NE2	1:C:197:ASN:O	2.23	0.71
2:D:9:PHE:CE1	2:D:10:ILE:CG1	2.74	0.70
1:C:18:HIS:HD2	2:D:17:MET:O	1.74	0.70
1:H:295:HIS:CD2	1:H:297:ILE:H	2.09	0.70
2:B:150:GLU:O	2:B:154:ASN:ND2	2.24	0.70
1:C:96:ASP:OD2	1:C:96(A):LEU:HD13	1.91	0.70
1:E:34:ASN:ND2	4:E:401:NAG:C2	2.42	0.70
1:H:294:PHE:HZ	2:I:59:MET:HG3	1.57	0.70
2:M:159:TYR:HB3	2:M:160:PRO:CD	2.21	0.70
2:B:150:GLU:OE2	2:B:153:ARG:CZ	2.41	0.69
2:M:26:HIS:O	2:M:26:HIS:HD2	1.75	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:146:ASN:O	2:F:150:GLU:HG2	1.93	0.68
2:M:151:SER:OG	2:M:157:TYR:HB2	1.93	0.68
2:I:26:HIS:HD2	2:I:153:ARG:HH12	1.41	0.68
2:I:28:ASN:O	2:I:31:GLY:N	2.25	0.68
1:J:294:PHE:HZ	2:K:59:MET:HG3	1.59	0.68
2:I:151:SER:HA	2:I:154:ASN:HB2	1.77	0.67
1:H:70:LEU:O	1:H:150:ASN:ND2	2.28	0.67
1:H:137:SER:O	1:H:140:PRO:HD3	1.94	0.67
1:A:18:HIS:HE1	1:A:19(A):ASN:HB3	1.60	0.67
1:L:203:SER:HB3	1:L:212:ARG:HD2	1.76	0.67
1:L:49:GLY:HA2	1:L:286:GLY:HA2	1.77	0.67
2:D:9:PHE:CD1	2:D:10:ILE:HG13	2.29	0.67
2:M:159:TYR:HB3	2:M:160:PRO:HD3	1.76	0.67
1:L:283:THR:HG22	1:L:285:ILE:H	1.60	0.67
2:M:157:TYR:CD2	2:M:158:ASP:N	2.63	0.67
2:B:9:PHE:CD1	2:B:9:PHE:C	2.73	0.66
1:H:77:ASP:OD1	1:H:149:ARG:HD2	1.94	0.66
1:A:37:THR:HG22	1:A:319:GLY:HA3	1.75	0.66
1:C:275:GLY:O	1:C:276:ASN:CB	2.42	0.66
1:H:125:PRO:HB2	1:H:126:SER:OG	1.95	0.66
2:B:9:PHE:HD1	2:B:9:PHE:O	1.79	0.66
1:H:100:GLY:HA3	1:H:230:MET:O	1.94	0.66
1:J:50:LYS:HD3	1:J:275:GLY:HA3	1.77	0.66
1:A:311:SER:HB3	2:B:97:GLU:OE2	1.96	0.66
2:K:27:SER:HB2	2:K:32:SER:OG	1.95	0.66
1:J:106:GLU:N	1:J:106:GLU:OE1	2.29	0.66
1:C:15:ILE:HD11	2:D:122:VAL:HG21	1.77	0.66
1:E:62:ARG:NH1	1:E:78:GLU:OE1	2.25	0.65
1:A:137:SER:O	1:A:140:PRO:HD3	1.97	0.65
1:A:316:LEU:HD13	2:B:100:VAL:HG22	1.78	0.65
2:B:30:GLN:OE1	2:B:30:GLN:N	2.30	0.65
2:B:106:ARG:NH2	2:D:106:ARG:HD2	2.12	0.65
1:J:15:ILE:HD11	2:K:122:VAL:HG21	1.79	0.65
1:J:186:ASN:OD1	1:J:186:ASN:N	2.28	0.64
1:C:27:THR:HG23	2:D:105:GLU:HB2	1.78	0.64
2:I:28:ASN:O	2:I:31:GLY:HA2	1.96	0.64
2:I:26:HIS:O	2:I:26:HIS:ND1	2.30	0.64
2:B:150:GLU:OE2	2:B:153:ARG:NE	2.29	0.64
2:D:28:ASN:OD1	2:D:28:ASN:C	2.39	0.64
1:H:197:ASN:OD1	1:H:197:ASN:N	2.27	0.64
1:C:316:LEU:HD13	2:D:100:VAL:HG22	1.80	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:17:TYR:CZ	2:I:6:ILE:HG23	2.33	0.64
1:A:15:ILE:HD12	2:B:119:TYR:HA	1.80	0.63
1:L:206:THR:OG1	1:L:209:LEU:HB3	1.98	0.63
2:M:42:GLN:HE21	2:M:42:GLN:HA	1.63	0.63
2:B:24:TYR:CE2	2:B:153:ARG:HG2	2.34	0.63
2:D:80:LEU:HD12	2:D:80:LEU:O	1.97	0.63
1:L:294:PHE:HZ	2:M:59:MET:HG3	1.64	0.63
2:K:22:TYR:OH	2:K:111:HIS:ND1	2.28	0.63
2:M:77:ILE:HD12	2:M:77:ILE:H	1.63	0.63
2:F:26:HIS:CD2	2:F:26:HIS:C	2.76	0.63
1:L:177:LEU:HB3	1:L:258:TYR:HB2	1.81	0.63
1:A:17:TYR:CE2	2:B:6:ILE:HG23	2.33	0.63
2:I:18:VAL:O	2:I:18:VAL:CG1	2.46	0.63
1:H:123:ILE:C	1:H:124:ILE:HD12	2.24	0.63
1:J:138:ALA:C	1:J:140:PRO:HD3	2.24	0.62
1:L:156:LYS:HD2	1:L:159:ASN:HA	1.81	0.62
1:H:242:ALA:N	3:O:1:NAG:H82	2.14	0.62
1:H:220:ARG:NH1	1:H:227:SER:O	2.30	0.62
2:I:106:ARG:HH21	2:K:106:ARG:HH11	1.46	0.62
1:J:179:LEU:HD23	1:J:234:TRP:HB3	1.81	0.62
1:E:24:GLN:NE2	1:E:34:ASN:HB3	2.14	0.62
1:A:31:MET:HE1	2:F:110:PHE:CZ	2.35	0.62
1:C:200:THR:HG21	1:C:250:ASN:OD1	2.00	0.61
2:D:68:ARG:HE	2:F:79:ASN:ND2	1.97	0.61
1:A:37:THR:HG23	1:A:38:HIS:ND1	2.15	0.61
1:A:75:MET:CB	1:A:95:ASN:ND2	2.62	0.61
2:I:7:ALA:CB	2:I:11:GLU:O	2.46	0.61
2:D:55:ILE:HG12	2:D:99:LEU:HD11	1.82	0.61
1:J:33:LYS:HG3	4:J:401:NAG:H82	1.82	0.61
1:L:124:ILE:HB	1:L:254:PRO:O	2.00	0.61
1:L:131:GLU:OE2	1:L:131:GLU:HA	2.00	0.61
1:A:49:GLY:C	1:A:50:LYS:HD3	2.25	0.61
2:B:140:PHE:CE2	2:B:149:MET:HE2	2.36	0.61
2:K:119:TYR:HE1	2:K:136:GLY:HA2	1.66	0.61
2:B:68:ARG:HH11	2:D:79:ASN:ND2	1.98	0.61
2:B:26:HIS:CD2	2:B:26:HIS:C	2.79	0.61
2:M:129:ASN:C	2:M:129:ASN:HD22	2.09	0.61
1:E:53(A):LEU:HD23	1:E:53(A):LEU:C	2.25	0.60
2:K:119:TYR:CE1	2:K:136:GLY:HA2	2.36	0.60
1:A:142:GLN:NE2	1:A:142:GLN:CA	2.55	0.60
2:D:29:GLU:O	2:D:30:GLN:C	2.42	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:295:HIS:HD2	1:J:297:ILE:H	1.48	0.60
2:M:129:ASN:O	2:M:129:ASN:ND2	2.30	0.60
2:B:127:ARG:HH12	2:D:131:LYS:HG2	1.66	0.60
2:K:125:GLN:HE22	2:K:155:GLY:C	2.09	0.60
1:E:205:GLY:O	1:E:206:THR:HG22	2.02	0.60
2:F:128:ASP:OD1	2:F:129:ASN:N	2.35	0.60
2:B:26:HIS:O	2:B:32:SER:HA	2.02	0.60
1:J:216:LYS:O	1:J:220:ARG:NH2	2.35	0.60
1:H:169:ASN:OD1	1:H:169:ASN:C	2.44	0.60
1:A:294:PHE:HZ	2:B:59:MET:HG3	1.65	0.59
1:C:283:THR:HB	1:C:286:GLY:O	2.01	0.59
2:K:26:HIS:HD2	2:K:153:ARG:HH12	1.50	0.59
1:A:125:PRO:HG2	1:A:126:SER:HB3	1.84	0.59
2:I:22:TYR:OH	2:I:111:HIS:HD2	1.85	0.59
1:J:134:GLY:HA3	1:J:153:TRP:HB3	1.84	0.59
2:B:68:ARG:HD3	2:D:79:ASN:HD22	1.68	0.59
1:H:97:CYS:O	1:H:138:ALA:HB1	2.03	0.59
1:J:33:LYS:CG	4:J:401:NAG:H82	2.33	0.59
1:L:200:THR:HG23	1:L:215:PRO:HG2	1.84	0.59
2:B:131:LYS:HG2	2:F:127:ARG:NH2	2.18	0.59
1:C:37:THR:HG23	1:C:320:LEU:O	2.03	0.59
1:C:295:HIS:HD2	1:C:297:ILE:H	1.49	0.59
2:F:28:ASN:HB3	2:F:149:MET:HE3	1.83	0.59
2:F:158:ASP:HB3	2:F:161:GLN:HB2	1.85	0.58
1:L:100:GLY:HA3	1:L:230:MET:O	2.03	0.58
2:M:126:LEU:O	2:M:129:ASN:HB3	2.03	0.58
2:K:68:ARG:HH21	2:M:79:ASN:ND2	2.01	0.58
1:L:311:SER:HB3	2:M:97:GLU:OE2	2.03	0.58
1:H:52:CYS:HB2	1:H:279:THR:HG22	1.84	0.58
1:J:200:THR:HG21	1:J:250:ASN:OD1	2.02	0.58
1:L:295:HIS:HD2	1:L:297:ILE:H	1.50	0.58
2:K:27:SER:HB3	2:K:32:SER:OG	2.02	0.58
2:I:79:ASN:HD22	2:M:68:ARG:HE	1.52	0.58
1:E:295:HIS:CD2	1:E:297:ILE:H	2.18	0.57
2:I:7:ALA:HB1	2:I:11:GLU:C	2.27	0.57
2:B:140:PHE:HE2	2:B:149:MET:HE2	1.67	0.57
1:L:49:GLY:HA2	1:L:286:GLY:CA	2.33	0.57
1:C:27:THR:HG22	1:C:31:MET:N	2.13	0.57
1:H:23:GLU:HG2	1:H:39:ALA:HB3	1.86	0.57
2:D:26:HIS:HB2	2:D:149:MET:HE3	1.86	0.57
1:L:314:LEU:HD22	2:M:100:VAL:HG21	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:M:159:TYR:C	2:M:161:GLN:N	2.62	0.57
1:A:201:TYR:OH	1:A:246:GLU:OE1	2.23	0.57
1:A:17:TYR:OH	2:B:6:ILE:HG23	2.04	0.56
1:J:311:SER:HB3	2:K:97:GLU:OE2	2.03	0.56
1:A:142:GLN:HE21	1:A:142:GLN:CA	2.06	0.56
2:M:21:TRP:HB2	2:M:41:THR:HG23	1.87	0.56
2:D:41:THR:HG22	2:D:45:ILE:HD11	1.87	0.56
1:E:37:THR:HG22	1:E:322:ASN:HB3	1.87	0.56
1:E:42:ILE:HD13	1:E:316:LEU:HD22	1.85	0.56
2:F:6:ILE:HD12	2:F:112:ASP:HA	1.86	0.56
2:M:161:GLN:C	2:M:161:GLN:CD	2.72	0.56
2:D:6:ILE:HG13	2:D:112:ASP:HA	1.88	0.56
1:A:131:GLU:OE2	1:A:157:LYS:HE3	2.05	0.56
1:J:77:ASP:O	1:J:80:ILE:HG12	2.05	0.56
2:B:68:ARG:HH11	2:D:79:ASN:HD21	1.53	0.56
1:E:25:VAL:HG21	1:E:317:ALA:HB2	1.88	0.56
2:D:150:GLU:OE1	2:D:154:ASN:ND2	2.39	0.56
2:I:28:ASN:O	2:I:31:GLY:CA	2.53	0.56
1:A:53(A):LEU:HD22	1:A:302:ILE:HD11	1.87	0.56
1:L:183:HIS:ND1	1:L:195:TYR:OH	2.31	0.56
1:A:185:SER:OG	1:A:191:GLN:OE1	2.22	0.55
1:C:299:PRO:HB3	2:D:89:LEU:HD11	1.88	0.55
1:H:25:VAL:HG12	1:H:315:VAL:HG22	1.88	0.55
1:L:317:ALA:H	2:M:104:ASN:HD21	1.53	0.55
1:J:100:GLY:HA3	1:J:230:MET:O	2.06	0.55
2:K:84:MET:HE3	2:K:88:PHE:HD2	1.70	0.55
1:L:179:LEU:HD23	1:L:234:TRP:HB3	1.88	0.55
1:H:45:LYS:HA	1:H:296:ASN:HD21	1.72	0.55
1:L:134:GLY:HA3	1:L:153:TRP:HB3	1.87	0.55
1:A:98:TYR:OH	1:A:228:GLY:HA3	2.07	0.55
1:E:307:LYS:HE3	2:F:62:GLN:HG2	1.88	0.55
2:F:24:TYR:CD2	2:F:153:ARG:HD3	2.42	0.55
2:F:128:ASP:OD1	2:F:128:ASP:C	2.50	0.55
2:I:7:ALA:HB1	2:I:11:GLU:CA	2.35	0.55
1:L:124:ILE:HD12	1:L:254:PRO:HG2	1.88	0.55
1:L:317:ALA:H	2:M:104:ASN:ND2	2.05	0.55
1:H:62:ARG:NH2	1:H:78:GLU:OE2	2.38	0.54
2:I:81:ASN:HB2	2:K:80:LEU:HD22	1.90	0.54
1:C:184:HIS:HB2	1:C:229:ARG:HB2	1.88	0.54
1:E:77:ASP:OD2	1:E:149:ARG:HD2	2.08	0.54
2:I:59:MET:O	2:I:62:GLN:HB2	2.08	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:27:THR:HG22	1:J:32:GLU:H	1.73	0.54
1:C:67:ALA:O	1:C:71:LEU:HB2	2.07	0.54
1:H:220:ARG:HH12	1:H:228:GLY:HA2	1.73	0.54
4:J:402:NAG:C8	1:L:221:SER:HB3	2.38	0.54
2:B:145:ASP:O	2:B:148:CYS:HB3	2.08	0.54
1:J:44:GLU:OE1	1:J:46:THR:HG22	2.08	0.53
1:J:316:LEU:HD23	2:K:52:VAL:HG22	1.89	0.53
2:B:72:ASN:OD1	2:B:75:ARG:NH1	2.42	0.53
2:B:150:GLU:OE2	2:B:153:ARG:NH2	2.39	0.53
1:C:73:ASN:ND2	1:C:97:CYS:HB3	2.23	0.53
2:F:84:MET:HE2	2:F:85:GLU:HG2	1.90	0.53
1:J:61:LEU:HA	1:J:79:PHE:CZ	2.43	0.53
1:A:33:LYS:HG2	4:A:401:NAG:H82	1.90	0.53
1:E:37:THR:OG1	1:E:38:HIS:HD2	1.91	0.53
1:E:77:ASP:OD1	1:E:149:ARG:NH1	2.41	0.53
2:F:28:ASN:OD1	2:F:31:GLY:O	2.26	0.53
2:K:26:HIS:O	2:K:32:SER:OG	2.26	0.53
2:D:133:LEU:HD12	2:D:137:CYS:CB	2.38	0.53
1:E:216:LYS:O	1:E:220:ARG:NH2	2.42	0.53
1:L:156:LYS:HE2	1:L:193:LYS:O	2.08	0.53
1:L:294:PHE:CZ	2:M:59:MET:HG3	2.43	0.53
2:M:21:TRP:CZ3	2:M:45:ILE:HG13	2.43	0.53
2:F:58:LYS:HE3	2:F:58:LYS:HA	1.90	0.53
1:H:14:CYS:O	2:I:24:TYR:HA	2.09	0.53
1:J:101:ASN:HB2	1:J:231:ASP:OD1	2.09	0.53
1:E:13:ILE:HG22	2:F:138:PHE:HB2	1.91	0.53
1:H:161:TYR:CD1	1:H:195:TYR:HB3	2.44	0.53
1:E:174:GLU:N	1:E:174:GLU:OE1	2.41	0.52
1:J:295:HIS:CD2	1:J:297:ILE:H	2.27	0.52
2:K:75:ARG:NH2	2:K:78:GLU:OE1	2.37	0.52
2:K:84:MET:HE3	2:K:88:PHE:CD2	2.44	0.52
1:A:134:GLY:HA3	1:A:153:TRP:HB3	1.89	0.52
1:L:121:ILE:HD11	1:L:176:LEU:HD11	1.92	0.52
1:A:98:TYR:CD2	1:A:230:MET:HG3	2.43	0.52
1:C:26:ASP:HB3	1:C:33:LYS:HD3	1.90	0.52
2:D:133:LEU:HD12	2:D:137:CYS:HB3	1.90	0.52
2:F:150:GLU:O	2:F:154:ASN:HB2	2.09	0.52
1:H:47:HIS:ND1	1:H:48:ASN:O	2.43	0.52
2:I:19:ASP:HB3	2:I:36:ALA:HB2	1.90	0.52
1:A:45:LYS:HG2	1:A:296:ASN:HD22	1.75	0.52
1:A:75:MET:HB2	1:A:95:ASN:HD22	1.73	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:100:GLY:HA3	1:E:230:MET:O	2.10	0.52
2:F:22:TYR:OH	2:F:111:HIS:HD2	1.93	0.52
2:B:62:GLN:HG2	2:B:63:PHE:H	1.75	0.52
2:B:103:GLU:O	2:B:107:THR:OG1	2.25	0.52
1:E:53:ASP:OD1	1:E:274:TYR:OH	2.23	0.52
1:L:27:THR:HB	2:M:105:GLU:HB2	1.92	0.52
1:A:65:SER:OG	1:A:96:ASP:OD1	2.22	0.52
1:L:309:VAL:HG13	2:M:93:THR:HA	1.91	0.52
2:I:135:ASN:OD1	2:I:137:CYS:HB2	2.09	0.52
1:A:33:LYS:CB	4:A:401:NAG:H82	2.40	0.51
1:H:37:THR:HB	1:H:319:GLY:HA3	1.92	0.51
2:B:131:LYS:HE3	2:F:127:ARG:HH22	1.75	0.51
1:H:44:GLU:OE2	1:H:46:THR:HG22	2.09	0.51
1:L:60:ILE:HD13	1:L:274:TYR:HB2	1.93	0.51
2:M:150:GLU:OE2	2:M:153:ARG:NH1	2.42	0.51
2:K:68:ARG:HE	2:M:79:ASN:HD22	1.58	0.51
1:E:98:TYR:CD2	1:E:230:MET:HG3	2.46	0.51
2:K:27:SER:CA	2:K:32:SER:OG	2.59	0.51
2:M:26:HIS:CD2	2:M:26:HIS:C	2.89	0.51
2:I:154:ASN:HB3	2:I:156:THR:HG23	1.91	0.51
2:K:132:GLU:O	2:K:132:GLU:HG2	2.09	0.51
1:L:205:GLY:O	1:L:206:THR:HG23	2.10	0.51
2:M:30:GLN:HE22	2:M:146:ASN:ND2	2.09	0.51
2:B:106:ARG:HH21	2:D:106:ARG:HD2	1.76	0.51
1:E:220:ARG:O	1:E:227:SER:HB2	2.11	0.51
2:K:58:LYS:HE3	2:M:97:GLU:HB3	1.93	0.51
2:M:159:TYR:N	2:M:160:PRO:HD2	2.26	0.51
1:H:124:ILE:O	1:H:124:ILE:HG22	2.11	0.51
1:J:38:HIS:HB2	1:J:318:THR:HG23	1.92	0.51
2:M:84:MET:HG3	2:M:85:GLU:N	2.25	0.51
1:C:23:GLU:HG2	1:C:39:ALA:HB3	1.92	0.50
2:I:9:PHE:O	2:I:135:ASN:HA	2.11	0.50
2:D:151:SER:OG	2:D:157:TYR:HA	2.11	0.50
1:H:158:ASN:O	1:H:160:THR:HG23	2.11	0.50
2:I:129:ASN:HD21	2:I:163:SER:HA	1.76	0.50
1:L:130:HIS:CE1	1:L:162:PRO:HD2	2.46	0.50
2:I:158:ASP:HB3	2:I:161:GLN:HB3	1.93	0.50
1:J:283:THR:HG22	1:J:285:ILE:H	1.76	0.50
1:A:14:CYS:HA	2:B:137:CYS:HA	1.93	0.50
1:E:48:ASN:OD1	1:E:48:ASN:O	2.29	0.50
1:L:34:ASN:ND2	4:L:402:NAG:O7	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:96:ASP:CG	1:H:96(A):LEU:HD13	2.37	0.50
2:D:119:TYR:OH	2:D:132:GLU:OE2	2.25	0.50
1:E:174:GLU:OE2	1:E:259:LYS:HE2	2.11	0.50
2:F:151:SER:HA	2:F:154:ASN:HB2	1.93	0.50
2:F:159:TYR:O	2:F:160:PRO:C	2.54	0.50
2:I:9:PHE:CD1	2:I:10:ILE:HG13	2.45	0.50
2:K:55:ILE:O	2:K:59:MET:HG2	2.12	0.50
2:B:144:CYS:SG	2:B:149:MET:HG3	2.52	0.50
2:K:125:GLN:HG2	2:K:157:TYR:HB3	1.94	0.50
1:E:101:ASN:HB2	1:E:231:ASP:OD1	2.12	0.49
1:H:183:HIS:CD2	1:H:230:MET:HE2	2.47	0.49
1:E:279:THR:OG1	1:E:280:LYS:N	2.44	0.49
1:C:25:VAL:HG11	1:C:317:ALA:HB2	1.94	0.49
1:L:307:LYS:HB3	2:M:62:GLN:NE2	2.27	0.49
2:F:62:GLN:HB2	2:F:92:TRP:CD2	2.47	0.49
1:H:119:GLU:OE1	1:H:259:LYS:NZ	2.34	0.49
2:I:150:GLU:O	2:I:154:ASN:HB2	2.12	0.49
2:I:26:HIS:ND1	2:I:26:HIS:C	2.71	0.49
2:M:30:GLN:NE2	2:M:146:ASN:ND2	2.60	0.49
1:A:48:ASN:OD1	1:A:48:ASN:O	2.30	0.49
1:H:32:GLU:CD	1:H:321:ARG:HH22	2.19	0.49
1:H:126:SER:HB3	1:H:166:ARG:NH2	2.27	0.49
1:C:220:ARG:O	1:C:227:SER:HB2	2.13	0.49
2:D:80:LEU:C	2:D:80:LEU:CD1	2.83	0.49
1:L:317:ALA:N	2:M:104:ASN:HD21	2.11	0.49
1:E:294:PHE:HZ	2:F:59:MET:HG3	1.78	0.49
1:H:294:PHE:CZ	2:I:59:MET:HG3	2.43	0.49
1:E:65:SER:OG	1:E:96:ASP:OD1	2.24	0.49
1:E:265:SER:OG	1:E:266:ALA:N	2.35	0.49
1:H:121:ILE:HG21	1:H:259:LYS:HD3	1.94	0.49
2:I:150:GLU:HA	2:I:153:ARG:HE	1.78	0.49
1:J:107:GLU:OE2	2:M:75:ARG:HB2	2.12	0.49
1:L:131:GLU:OE1	1:L:133(A):SER:OG	2.30	0.49
2:D:154:ASN:HB3	2:D:156:THR:HG23	1.95	0.49
2:M:25:HIS:HB2	2:M:34:TYR:CD1	2.48	0.49
1:C:44:GLU:HG3	1:C:290:SER:CB	2.43	0.48
1:C:207:SER:HA	1:E:229:ARG:NH2	2.27	0.48
2:D:52:VAL:O	2:D:56:ILE:HG12	2.13	0.48
1:E:128:SER:HB2	1:E:129:ASP:OD2	2.12	0.48
1:J:105:TYR:CE2	1:J:109:LYS:HE3	2.48	0.48
2:K:148:CYS:O	2:K:151:SER:HB3	2.12	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:103:GLU:OE1	1:C:28:ILE:HD12	2.12	0.48
2:F:158:ASP:OD2	2:F:161:GLN:OE1	2.30	0.48
1:A:156:LYS:HD2	1:A:196:GLN:HB2	1.95	0.48
2:B:51:LYS:NZ	2:B:107:THR:OG1	2.47	0.48
1:C:222:LYS:HD3	1:C:225:GLY:O	2.12	0.48
2:D:145:ASP:O	2:D:148:CYS:CB	2.55	0.48
2:F:5:ALA:HB2	2:F:116:LYS:HB2	1.94	0.48
1:J:128:SER:O	1:J:157:LYS:HE3	2.14	0.48
1:A:265:SER:OG	1:A:266:ALA:N	2.46	0.48
2:I:17:MET:HE1	2:I:24:TYR:N	2.27	0.48
2:I:62:GLN:HG3	2:I:92:TRP:CD2	2.47	0.48
1:J:117:HIS:O	1:J:261:VAL:HG12	2.14	0.48
2:I:145:ASP:OD2	2:I:145:ASP:N	2.40	0.48
1:A:221:SER:HB3	3:N:1:NAG:H81	1.94	0.48
1:C:134:GLY:HA2	1:C:153:TRP:HB3	1.96	0.48
1:L:77:ASP:O	1:L:80:ILE:HG12	2.13	0.48
2:B:24:TYR:CZ	2:B:153:ARG:HG2	2.48	0.48
1:C:147:PHE:O	1:C:148:PHE:C	2.55	0.48
1:H:125:PRO:O	1:H:126:SER:HB2	2.13	0.48
2:K:59:MET:HE3	2:K:62:GLN:HE21	1.79	0.48
1:L:60:ILE:CD1	1:L:274:TYR:HB2	2.44	0.48
1:C:42:ILE:O	1:C:293:PRO:HD2	2.13	0.48
1:C:53:ASP:OD2	1:C:274:TYR:OH	2.25	0.48
2:D:127:ARG:HH12	2:F:131:LYS:CE	2.24	0.48
1:A:44:GLU:OE2	1:A:46:THR:HB	2.13	0.47
1:E:239:PRO:O	1:E:240:ASN:CB	2.62	0.47
1:C:107:GLU:OE2	2:F:75:ARG:HB2	2.14	0.47
1:E:186:ASN:ND2	1:E:227:SER:OG	2.47	0.47
2:F:51:LYS:HE3	2:F:103:GLU:HB3	1.97	0.47
2:F:68:ARG:NH1	2:F:81:ASN:OD1	2.44	0.47
1:L:197:ASN:N	1:L:197:ASN:HD22	2.11	0.47
1:A:220:ARG:NH1	1:A:227:SER:O	2.46	0.47
1:C:139:CYS:HB2	1:C:146:SER:O	2.14	0.47
1:E:129:ASP:OD2	1:E:129:ASP:N	2.47	0.47
2:F:24:TYR:CE2	2:F:153:ARG:HD3	2.49	0.47
2:B:6:ILE:HA	2:B:7:ALA:HA	1.68	0.47
1:C:130:HIS:NE2	1:C:162:PRO:HD2	2.30	0.47
1:C:314:LEU:HD22	2:D:100:VAL:HG21	1.96	0.47
2:D:45:ILE:H	2:D:45:ILE:HG13	1.49	0.47
1:H:32:GLU:OE1	1:H:321:ARG:NH2	2.37	0.47
1:J:89:GLU:OE1	1:J:269:LYS:NZ	2.47	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:283:THR:HB	1:J:286:GLY:O	2.14	0.47
2:K:151:SER:OG	2:K:157:TYR:HA	2.14	0.47
1:E:200:THR:HA	1:E:248:ASN:OD1	2.14	0.47
1:E:239:PRO:O	1:E:240:ASN:HB2	2.14	0.47
2:I:126:LEU:O	2:I:129:ASN:HB2	2.15	0.47
2:K:62:GLN:HG3	2:K:92:TRP:CD2	2.50	0.47
1:L:220:ARG:O	1:L:227:SER:HB2	2.15	0.47
1:H:174:GLU:OE1	1:H:259:LYS:HE3	2.14	0.47
1:J:204:VAL:HG22	1:J:245:PHE:CD2	2.50	0.47
1:H:124:ILE:HD12	1:H:124:ILE:N	2.29	0.47
1:H:296:ASN:HD22	1:H:296:ASN:C	2.23	0.47
2:K:68:ARG:NH1	2:K:81:ASN:OD1	2.48	0.47
2:I:26:HIS:CD2	2:I:153:ARG:HH12	2.28	0.47
2:K:150:GLU:O	2:K:154:ASN:HB2	2.15	0.47
1:L:127:TRP:HB2	1:L:132:ALA:HB2	1.97	0.47
1:L:205:GLY:O	1:L:206:THR:CG2	2.63	0.47
1:C:44:GLU:HG3	1:C:290:SER:HB2	1.96	0.47
2:F:51:LYS:HE2	2:F:107:THR:OG1	2.15	0.47
1:C:38:HIS:CB	1:C:318:THR:HG23	2.46	0.46
1:E:205:GLY:C	1:E:206:THR:CG2	2.88	0.46
1:H:21:SER:OG	1:H:22:THR:N	2.47	0.46
2:M:142:HIS:ND1	2:M:162:TYR:CD1	2.83	0.46
1:A:200:THR:O	1:A:214:VAL:HG13	2.16	0.46
1:C:62:ARG:NH1	1:C:78:GLU:OE1	2.49	0.46
1:C:117:HIS:HB3	1:C:261:VAL:HG23	1.97	0.46
1:C:133(A):SER:N	1:C:134:GLY:HA3	2.30	0.46
1:E:156:LYS:HD2	1:E:196:GLN:HG2	1.97	0.46
1:H:307:LYS:HD2	2:I:62:GLN:HB3	1.97	0.46
1:L:200:THR:HG21	1:L:250:ASN:OD1	2.15	0.46
1:E:77:ASP:CG	1:E:149:ARG:HD2	2.39	0.46
2:D:41:THR:O	2:D:45:ILE:HG13	2.15	0.46
1:E:44:GLU:OE1	1:E:290:SER:OG	2.29	0.46
1:E:200:THR:O	1:E:214:VAL:HG13	2.15	0.46
1:H:38:HIS:C	1:H:318:THR:HG22	2.40	0.46
1:J:55:GLY:N	1:J:278:ASN:OD1	2.45	0.46
1:L:124:ILE:O	1:L:255:GLU:HG3	2.15	0.46
2:D:17:MET:HE2	2:D:17:MET:HB2	1.75	0.46
1:H:134:GLY:C	1:H:135:VAL:HG13	2.41	0.46
2:I:6:ILE:HA	2:I:7:ALA:HA	1.67	0.46
1:L:97:CYS:HB2	1:L:139:CYS:CA	2.45	0.46
1:L:283:THR:HG22	1:L:285:ILE:N	2.28	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:307:LYS:HA	1:L:307:LYS:HD3	1.64	0.46
3:G:1:NAG:H61	3:G:2:NAG:HN2	1.79	0.46
1:A:295:HIS:CD2	1:A:297:ILE:H	2.24	0.46
1:J:33:LYS:HB2	4:J:401:NAG:C8	2.35	0.46
1:J:139:CYS:N	1:J:140:PRO:HD3	2.30	0.46
1:L:38:HIS:C	1:L:318:THR:HG22	2.41	0.46
1:L:176:LEU:O	1:L:236:ILE:HA	2.15	0.46
2:B:146:ASN:O	2:B:150:GLU:HB2	2.16	0.46
2:D:17:MET:HE1	2:D:23:GLY:HA3	1.98	0.46
2:M:9:PHE:CD1	2:M:10:ILE:HG13	2.50	0.46
1:A:122:GLN:HB2	1:A:256:TYR:CE1	2.50	0.46
1:C:101:ASN:HB2	1:C:231:ASP:OD1	2.16	0.46
1:E:22:THR:O	1:E:24:GLN:HG3	2.15	0.46
2:I:117:ASN:HD22	2:I:117:ASN:N	2.14	0.46
1:A:15:ILE:HD13	2:B:118:LEU:HD23	1.96	0.46
1:C:186:ASN:HD21	1:C:227:SER:C	2.23	0.45
1:H:52:CYS:HB3	1:H:277:CYS:O	2.16	0.45
1:H:97:CYS:HB2	1:H:138:ALA:O	2.16	0.45
1:C:295:HIS:CD2	1:C:297:ILE:H	2.32	0.45
2:F:72:ASN:OD1	2:F:72:ASN:N	2.35	0.45
1:H:50:LYS:HD3	1:H:275:GLY:HA3	1.97	0.45
1:J:202:ILE:HD11	1:J:251:PHE:HA	1.98	0.45
1:L:275:GLY:O	1:L:276:ASN:HB2	2.16	0.45
2:F:104:ASN:HD22	2:F:104:ASN:HA	1.57	0.45
2:K:24:TYR:CE2	2:K:153:ARG:HB3	2.52	0.45
2:F:28:ASN:O	2:F:31:GLY:O	2.34	0.45
1:J:164:ILE:O	1:J:246:GLU:HA	2.16	0.45
1:L:15:ILE:HG23	2:M:118:LEU:HD23	1.97	0.45
1:H:166:ARG:HG2	1:H:166:ARG:HH11	1.80	0.45
2:I:145:ASP:OD1	2:I:147:GLU:HB2	2.16	0.45
2:K:19:ASP:OD1	2:K:19:ASP:N	2.49	0.45
1:C:133:SER:O	1:C:133(A):SER:OG	2.30	0.45
1:E:283:THR:HG22	1:E:285:ILE:H	1.82	0.45
1:J:91:ALA:HA	1:J:269:LYS:HE3	1.98	0.45
1:H:182:ILE:HB	1:H:202:ILE:HD12	1.98	0.45
1:J:151:VAL:HG13	1:J:252:ILE:HG22	1.99	0.45
1:L:173:GLN:HE21	1:L:173:GLN:HB3	1.65	0.45
1:C:108:LEU:HB2	1:C:234:TRP:CE2	2.51	0.45
1:E:97:CYS:HB2	1:E:138:ALA:O	2.16	0.45
1:E:293:PRO:HG2	1:E:294:PHE:CD1	2.52	0.45
2:B:9:PHE:C	2:B:9:PHE:HD1	2.19	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:146:ASN:C	2:D:148:CYS:N	2.71	0.45
2:D:150:GLU:HA	2:D:153:ARG:HE	1.82	0.45
2:K:27:SER:HA	2:K:32:SER:OG	2.17	0.45
1:L:136:SER:OG	1:L:137:SER:N	2.50	0.45
1:L:265:SER:OG	1:L:266:ALA:N	2.50	0.45
2:M:15:GLN:HE21	2:M:15:GLN:HB2	1.62	0.45
2:M:30:GLN:HE21	2:M:30:GLN:HB2	1.47	0.45
1:H:222:LYS:HB2	1:H:222:LYS:HE3	1.77	0.44
1:H:301:THR:HB	1:H:305:CYS:SG	2.57	0.44
1:C:301:THR:O	1:C:302:ILE:HD13	2.17	0.44
2:D:30:GLN:H	2:D:30:GLN:HG2	1.51	0.44
1:E:122:GLN:HG3	1:E:256:TYR:CE2	2.52	0.44
2:K:55:ILE:HG23	2:K:99:LEU:HD13	1.99	0.44
1:L:48:ASN:CG	1:L:48:ASN:O	2.59	0.44
2:B:29:GLU:HB3	2:B:30:GLN:OE1	2.17	0.44
1:C:307:LYS:HB3	2:D:62:GLN:CD	2.43	0.44
2:D:68:ARG:HH21	2:F:79:ASN:ND2	2.15	0.44
1:A:174:GLU:CD	1:A:259:LYS:HE3	2.42	0.44
1:A:247:SER:OG	1:A:251:PHE:HB2	2.16	0.44
1:C:65:SER:OG	1:C:96:ASP:HA	2.17	0.44
1:C:304:GLU:H	1:C:304:GLU:HG2	1.61	0.44
2:I:57:ASP:C	2:I:59:MET:H	2.26	0.44
1:J:183:HIS:HB2	1:J:230:MET:HE2	1.99	0.44
1:E:322:ASN:OD1	1:E:322:ASN:C	2.61	0.44
1:H:96:ASP:C	1:H:96(A):LEU:HD12	2.42	0.44
2:I:37:ASP:O	2:I:41:THR:HB	2.17	0.44
1:L:42:ILE:O	1:L:293:PRO:HD2	2.17	0.44
2:D:106:ARG:HH22	2:F:106:ARG:HG3	1.82	0.44
1:E:317:ALA:O	2:F:107:THR:HG21	2.18	0.44
2:K:63:PHE:CZ	2:K:84:MET:HE1	2.53	0.44
1:L:295:HIS:HD2	1:L:297:ILE:N	2.14	0.44
1:E:38:HIS:HB2	1:E:318:THR:HG22	1.99	0.44
1:L:123:ILE:HG13	1:L:168:TYR:CG	2.52	0.44
1:A:293:PRO:HB2	1:A:294:PHE:CD2	2.53	0.44
1:H:47:HIS:CE1	1:H:285:ILE:HG22	2.53	0.44
1:J:42:ILE:O	1:J:293:PRO:HD2	2.17	0.44
2:M:129:ASN:ND2	2:M:142:HIS:NE2	2.49	0.44
2:D:145:ASP:O	2:D:148:CYS:N	2.45	0.43
1:E:237:LEU:HD12	1:E:241:ASP:HB3	2.00	0.43
1:E:240:ASN:HD21	3:N:2:NAG:H61	1.83	0.43
2:F:57:ASP:O	2:F:60:ASN:ND2	2.51	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:164:GLU:H	2:I:164:GLU:HG2	1.53	0.43
2:F:149:MET:O	2:F:152:VAL:HG22	2.18	0.43
1:L:28:ILE:HD13	1:L:28:ILE:HA	1.86	0.43
1:A:75:MET:CB	1:A:95:ASN:HD21	2.30	0.43
1:A:186:ASN:HD21	1:A:227:SER:C	2.26	0.43
2:B:149:MET:C	2:B:151:SER:N	2.77	0.43
1:C:38:HIS:HB3	1:C:318:THR:HG23	2.00	0.43
1:E:144:THR:HG23	1:E:145:PRO:HD2	2.00	0.43
1:H:197:ASN:HA	1:H:198:PRO:HD3	1.72	0.43
1:A:15:ILE:HG23	2:B:118:LEU:HD23	1.99	0.43
2:B:11:GLU:CG	2:B:12:GLY:N	2.81	0.43
1:E:19(A):ASN:OD1	1:E:21:SER:HB2	2.19	0.43
1:L:23:GLU:OE2	1:L:39:ALA:HB3	2.19	0.43
1:E:17:TYR:CZ	2:F:6:ILE:HG23	2.53	0.43
1:L:164:ILE:O	1:L:246:GLU:HA	2.19	0.43
1:L:178:ILE:O	1:L:234:TRP:HA	2.19	0.43
1:L:197:ASN:N	1:L:197:ASN:ND2	2.64	0.43
1:L:295:HIS:CD2	1:L:297:ILE:HB	2.53	0.43
2:K:94:TYR:HD2	2:K:95:ASN:HD22	1.66	0.43
1:L:183:HIS:HB2	1:L:252:ILE:HD11	2.00	0.43
2:M:145:ASP:C	2:M:147:GLU:H	2.27	0.43
2:B:144:CYS:SG	2:B:149:MET:CG	3.07	0.43
1:H:239:PRO:O	1:H:240:ASN:HB2	2.18	0.43
2:D:42:GLN:HA	2:D:45:ILE:HD12	2.00	0.43
2:D:159:TYR:HB3	2:D:160:PRO:HD3	2.01	0.43
1:E:308:TYR:O	2:F:62:GLN:NE2	2.49	0.43
1:L:25:VAL:HG12	1:L:315:VAL:HG22	2.00	0.43
1:A:80:ILE:HD13	1:A:80:ILE:HA	1.90	0.43
2:D:48:VAL:O	2:D:52:VAL:HG23	2.19	0.43
3:P:1:NAG:O4	3:P:2:NAG:C2	2.57	0.43
2:B:128:ASP:OD1	2:B:159:TYR:OH	2.23	0.42
1:C:17:TYR:CZ	2:D:6:ILE:HG23	2.54	0.42
1:C:293:PRO:CG	2:D:56:ILE:HD12	2.49	0.42
1:H:220:ARG:NH1	1:H:228:GLY:HA2	2.33	0.42
1:J:27:THR:HG22	1:J:31:MET:H	1.82	0.42
2:K:95:ASN:O	2:K:99:LEU:HB2	2.19	0.42
1:A:19(A):ASN:OD1	1:A:21:SER:HB2	2.19	0.42
1:E:11:ASP:OD2	2:F:144:CYS:N	2.41	0.42
2:K:94:TYR:HD2	2:K:95:ASN:ND2	2.17	0.42
1:E:109:LYS:HD3	2:F:69:GLU:CD	2.45	0.42
1:E:121:ILE:HG21	1:E:259:LYS:HD2	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:I:9:PHE:CE1	2:I:10:ILE:HG13	2.55	0.42
2:I:71:ASN:OD1	2:I:74:GLU:HG3	2.19	0.42
1:L:209:LEU:HD22	1:L:235:THR:HG21	2.01	0.42
1:A:10:GLY:N	2:B:139:GLU:OE2	2.52	0.42
2:B:5:ALA:HB2	2:B:116:LYS:HB2	2.02	0.42
1:E:58:PRO:HB3	1:E:86:TYR:CZ	2.54	0.42
2:F:108:LEU:O	2:F:111:HIS:N	2.47	0.42
1:L:60:ILE:HG22	1:L:62:ARG:HG3	2.00	0.42
1:L:268:ILE:HD12	1:L:284:PRO:HA	2.01	0.42
2:M:16:GLY:HA3	2:M:34:TYR:CE2	2.55	0.42
2:D:146:ASN:O	2:D:149:MET:N	2.52	0.42
2:K:119:TYR:OH	2:K:132:GLU:OE1	2.29	0.42
1:A:100:GLY:HA3	1:A:230:MET:O	2.19	0.42
1:C:11:ASP:OD2	2:D:143:LYS:HD2	2.20	0.42
2:D:51:LYS:HE3	1:E:31:MET:HE2	2.01	0.42
1:E:237:LEU:CD1	1:E:241:ASP:HB3	2.49	0.42
1:E:283:THR:CG2	1:E:285:ILE:HG12	2.49	0.42
1:L:205:GLY:C	1:L:206:THR:HG23	2.45	0.42
1:A:78:GLU:O	1:A:78:GLU:HG2	2.20	0.42
2:B:79:ASN:HD22	2:F:68:ARG:HE	1.67	0.42
1:C:298:HIS:O	1:C:308:TYR:CD1	2.73	0.42
2:D:3:PHE:HB2	2:D:112:ASP:OD2	2.19	0.42
1:H:124:ILE:HG21	1:H:127:TRP:CZ2	2.55	0.42
2:M:129:ASN:C	2:M:129:ASN:ND2	2.77	0.42
2:M:129:ASN:OD1	2:M:157:TYR:OH	2.37	0.42
2:M:129:ASN:HD21	2:M:142:HIS:CD2	2.35	0.42
1:C:197:ASN:HA	1:C:198:PRO:HD3	1.74	0.42
1:C:307:LYS:HB3	2:D:62:GLN:OE1	2.19	0.42
2:D:29:GLU:C	2:D:31:GLY:N	2.64	0.42
1:L:50:LYS:HE2	1:L:50:LYS:HB3	1.73	0.42
2:M:4:GLY:O	2:M:8:GLY:HA3	2.20	0.42
1:A:25:VAL:HG12	1:A:315:VAL:HG22	2.01	0.41
2:B:53:ASN:N	2:B:53:ASN:HD22	2.17	0.41
1:H:164:ILE:O	1:H:246:GLU:HA	2.20	0.41
2:K:29:GLU:HB2	2:K:143:LYS:NZ	2.34	0.41
1:L:131:GLU:OE2	1:L:131:GLU:CA	2.64	0.41
1:A:77:ASP:OD1	1:A:149:ARG:HD2	2.20	0.41
1:H:125(A):LYS:HG2	1:H:132:ALA:HB1	2.01	0.41
2:I:98:LEU:HG	2:I:102:MET:HE3	2.02	0.41
2:I:102:MET:O	2:I:106:ARG:HG3	2.20	0.41
2:K:85:GLU:O	2:K:89:LEU:HG	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:97:CYS:HB2	1:L:139:CYS:HA	2.01	0.41
1:L:128:SER:OG	1:L:129:ASP:N	2.53	0.41
2:D:150:GLU:CA	2:D:153:ARG:HH21	2.33	0.41
1:H:268:ILE:HD12	1:H:284:PRO:HG3	2.01	0.41
2:K:107:THR:O	2:K:110:PHE:HB3	2.21	0.41
1:L:83:GLU:OE2	1:L:262:LYS:NZ	2.41	0.41
1:C:184:HIS:HB3	1:C:220:ARG:NH2	2.36	0.41
2:D:135:ASN:HD22	2:D:135:ASN:HA	1.62	0.41
1:J:37:THR:HG23	1:J:320:LEU:O	2.20	0.41
2:M:25:HIS:CD2	2:M:25:HIS:C	2.97	0.41
1:A:180:TRP:CE2	1:A:204:VAL:HG21	2.55	0.41
2:B:51:LYS:HA	1:C:28:ILE:O	2.21	0.41
1:E:282:GLN:HG3	1:E:283:THR:N	2.35	0.41
2:F:85:GLU:O	2:F:89:LEU:HD22	2.20	0.41
2:M:9:PHE:CE1	2:M:10:ILE:HG13	2.56	0.41
2:M:102:MET:O	2:M:106:ARG:HG3	2.20	0.41
1:A:38:HIS:HB2	1:A:318:THR:HG22	2.02	0.41
1:A:121:ILE:HG13	1:A:121:ILE:O	2.20	0.41
1:C:73:ASN:HA	1:C:74:PRO:HD3	1.98	0.41
1:E:240:ASN:HD21	3:N:2:NAG:C6	2.34	0.41
1:H:268:ILE:HD13	1:H:268:ILE:HA	1.92	0.41
1:A:176:LEU:HD12	1:A:259:LYS:HA	2.03	0.41
1:C:307:LYS:HD3	1:C:307:LYS:HA	1.64	0.41
1:J:131:GLU:HG2	1:J:133(A):SER:HB2	2.02	0.41
1:L:170:ASN:HD22	1:L:238:LYS:C	2.28	0.41
1:C:47:HIS:CE1	1:C:285:ILE:HG13	2.56	0.41
1:E:98:TYR:HA	1:E:99:PRO:HD3	1.89	0.41
1:E:142:GLN:HA	1:E:142:GLN:NE2	2.36	0.41
1:E:283:THR:CG2	1:E:285:ILE:H	2.33	0.41
2:M:87:GLY:O	2:M:91:VAL:HG23	2.21	0.41
1:A:122:GLN:HE21	1:A:125:PRO:HA	1.86	0.41
1:C:73:ASN:HB3	1:C:76:CYS:SG	2.60	0.41
2:D:38:LYS:HG2	2:D:39:GLU:H	1.86	0.41
1:E:42:ILE:O	1:E:293:PRO:HD2	2.19	0.41
1:E:200:THR:HG23	1:E:215:PRO:HD2	2.02	0.41
1:E:293:PRO:HG2	1:E:294:PHE:HD1	1.85	0.41
1:H:234:TRP:HZ3	1:H:236:ILE:HD13	1.86	0.41
1:J:14:CYS:O	2:K:24:TYR:HA	2.21	0.41
1:J:73:ASN:HB3	1:J:76:CYS:SG	2.61	0.41
2:K:37:ASP:O	2:K:41:THR:HB	2.20	0.41
1:L:101:ASN:HB2	1:L:231:ASP:OD1	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:210:ASN:CG	1:E:220:ARG:HE	2.28	0.41
1:L:33:LYS:CB	4:L:402:NAG:H81	2.50	0.41
1:C:42:ILE:HG12	1:C:314:LEU:O	2.21	0.40
1:C:290:SER:OG	1:C:291:SER:N	2.54	0.40
2:D:168:LEU:HD13	2:D:168:LEU:HA	1.71	0.40
2:I:101:LEU:HD12	2:I:101:LEU:HA	1.90	0.40
1:C:43:LEU:HB2	1:C:314:LEU:HB2	2.03	0.40
1:C:44:GLU:HG3	1:C:290:SER:OG	2.21	0.40
1:C:156:LYS:HB3	1:C:195:TYR:HA	2.03	0.40
1:H:65:SER:OG	1:H:96:ASP:OD1	2.34	0.40
1:A:53:ASP:OD1	1:A:274:TYR:OH	2.27	0.40
1:A:156:LYS:NZ	1:A:192:THR:O	2.52	0.40
1:A:269:LYS:HE3	1:A:269:LYS:HB3	1.92	0.40
2:B:133:LEU:HD23	2:B:133:LEU:HA	1.73	0.40
1:C:209:LEU:HD22	1:C:235:THR:HG21	2.04	0.40
2:D:107:THR:O	2:D:110:PHE:HB3	2.21	0.40
1:E:77:ASP:OD1	1:E:149:ARG:HD2	2.21	0.40
2:F:8:GLY:O	2:F:9:PHE:C	2.61	0.40
1:J:17:TYR:CZ	2:K:6:ILE:HG23	2.57	0.40
2:K:62:GLN:HG2	2:K:63:PHE:N	2.36	0.40
2:D:95:ASN:HD22	2:D:95:ASN:HA	1.57	0.40
1:E:184:HIS:HB3	1:E:220:ARG:NH2	2.36	0.40
1:E:205:GLY:O	1:E:206:THR:CG2	2.69	0.40
1:J:105:TYR:HB3	1:J:106:GLU:OE1	2.21	0.40
2:M:54:SER:O	2:M:58:LYS:HB2	2.21	0.40
1:A:122:GLN:NE2	1:A:125:PRO:HA	2.36	0.40
1:C:221:SER:HB3	3:G:1:NAG:H81	2.03	0.40
1:C:247:SER:OG	1:C:251:PHE:HB2	2.22	0.40
2:M:14:TRP:C	2:M:16:GLY:N	2.78	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	319/329 (97%)	305 (96%)	13 (4%)	1 (0%)	36	66
1	C	319/329 (97%)	305 (96%)	14 (4%)	0	100	100
1	E	320/329 (97%)	296 (92%)	22 (7%)	2 (1%)	21	51
1	H	319/329 (97%)	299 (94%)	20 (6%)	0	100	100
1	J	319/329 (97%)	298 (93%)	21 (7%)	0	100	100
1	L	317/329 (96%)	288 (91%)	29 (9%)	0	100	100
2	B	168/182 (92%)	159 (95%)	9 (5%)	0	100	100
2	D	170/182 (93%)	161 (95%)	9 (5%)	0	100	100
2	F	170/182 (93%)	154 (91%)	15 (9%)	1 (1%)	21	51
2	I	168/182 (92%)	154 (92%)	14 (8%)	0	100	100
2	K	170/182 (93%)	157 (92%)	11 (6%)	2 (1%)	10	34
2	M	167/182 (92%)	155 (93%)	12 (7%)	0	100	100
All	All	2926/3066 (95%)	2731 (93%)	189 (6%)	6 (0%)	43	72

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	K	161	GLN
1	E	140	PRO
1	A	77	ASP
1	E	78	GLU
2	K	39	GLU
2	F	18	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	286/292 (98%)	256 (90%)	30 (10%)	6	22
1	C	286/292 (98%)	249 (87%)	37 (13%)	4	14
1	E	287/292 (98%)	251 (88%)	36 (12%)	4	15

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	H	286/292 (98%)	256 (90%)	30 (10%)	6	22
1	J	286/292 (98%)	253 (88%)	33 (12%)	5	18
1	L	285/292 (98%)	257 (90%)	28 (10%)	7	25
2	B	145/156 (93%)	119 (82%)	26 (18%)	2	6
2	D	147/156 (94%)	123 (84%)	24 (16%)	2	8
2	F	147/156 (94%)	124 (84%)	23 (16%)	2	9
2	I	145/156 (93%)	115 (79%)	30 (21%)	1	4
2	K	147/156 (94%)	123 (84%)	24 (16%)	2	8
2	M	144/156 (92%)	117 (81%)	27 (19%)	1	5
All	All	2591/2688 (96%)	2243 (87%)	348 (13%)	4	13

All (348) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	SER
1	A	27	THR
1	A	36	VAL
1	A	37	THR
1	A	46	THR
1	A	60	ILE
1	A	63	ASP
1	A	82	VAL
1	A	82(A)	PRO
1	A	139	CYS
1	A	142	GLN
1	A	144	THR
1	A	194	LEU
1	A	199	THR
1	A	208	THR
1	A	211	GLN
1	A	221	SER
1	A	238	LYS
1	A	259	LYS
1	A	268	ILE
1	A	279	THR
1	A	283	THR
1	A	302	ILE
1	A	311	SER
1	A	312	ASN

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Mol	Chain	Res	Type
1	A	315	VAL
1	A	316	LEU
1	A	318	THR
1	A	320	LEU
1	A	323	SER
2	B	9	PHE
2	B	22	TYR
2	B	26	HIS
2	B	42	GLN
2	B	57	ASP
2	B	60	ASN
2	B	61	THR
2	B	78	GLU
2	B	80	LEU
2	B	84	MET
2	B	86	ASP
2	B	89	LEU
2	B	100	VAL
2	B	101	LEU
2	B	107	THR
2	B	108	LEU
2	B	113	SER
2	B	121	LYS
2	B	124	LEU
2	B	128	ASP
2	B	129	ASN
2	B	139	GLU
2	B	152	VAL
2	B	163	SER
2	B	164	GLU
2	B	168	LEU
1	C	15	ILE
1	C	22	THR
1	C	23	GLU
1	C	26	ASP
1	C	28	ILE
1	C	46	THR
1	C	66	VAL
1	C	71	LEU
1	C	80	ILE
1	C	82	VAL
1	C	82(A)	PRO

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Mol	Chain	Res	Type
1	C	111	LEU
1	C	114	ARG
1	C	136	SER
1	C	151	VAL
1	C	167	SER
1	C	169	ASN
1	C	179	LEU
1	C	185	SER
1	C	195	TYR
1	C	197	ASN
1	C	200	THR
1	C	211	GLN
1	C	223	VAL
1	C	235	THR
1	C	236	ILE
1	C	244	ASN
1	C	269	LYS
1	C	276	ASN
1	C	283	THR
1	C	285	ILE
1	C	291	SER
1	C	300	LEU
1	C	304	GLU
1	C	309	VAL
1	C	313	LYS
1	C	315	VAL
2	D	2	LEU
2	D	22	TYR
2	D	30	GLN
2	D	32	SER
2	D	38	LYS
2	D	41	THR
2	D	45	ILE
2	D	46	ASP
2	D	54	SER
2	D	62	GLN
2	D	73	LEU
2	D	80	LEU
2	D	83	LYS
2	D	84	MET
2	D	99	LEU
2	D	106	ARG

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Mol	Chain	Res	Type
2	D	108	LEU
2	D	113	SER
2	D	135	ASN
2	D	147	GLU
2	D	149	MET
2	D	168	LEU
2	D	171	GLU
2	D	172	GLU
1	E	11	ASP
1	E	15	ILE
1	E	22	THR
1	E	23	GLU
1	E	46	THR
1	E	54	ASP
1	E	63	ASP
1	E	81	ASN
1	E	82	VAL
1	E	82(A)	PRO
1	E	112	LEU
1	E	114	ARG
1	E	129	ASP
1	E	133	SER
1	E	133(A)	SER
1	E	142	GLN
1	E	149	ARG
1	E	164	ILE
1	E	179	LEU
1	E	195	TYR
1	E	197	ASN
1	E	200	THR
1	E	206	THR
1	E	219	THR
1	E	221	SER
1	E	244	ASN
1	E	246	GLU
1	E	283	THR
1	E	285	ILE
1	E	291	SER
1	E	300	LEU
1	E	314	LEU
1	E	316	LEU
1	E	318	THR

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Mol	Chain	Res	Type
1	E	320	LEU
1	E	322	ASN
2	F	18	VAL
2	F	21	TRP
2	F	22	TYR
2	F	24	TYR
2	F	26	HIS
2	F	39	GLU
2	F	43	LYS
2	F	48	VAL
2	F	51	LYS
2	F	58	LYS
2	F	60	ASN
2	F	69	GLU
2	F	72	ASN
2	F	77	ILE
2	F	89	LEU
2	F	101	LEU
2	F	107	THR
2	F	108	LEU
2	F	113	SER
2	F	126	LEU
2	F	149	MET
2	F	152	VAL
2	F	165	GLU
1	H	22	THR
1	H	23	GLU
1	H	28	ILE
1	H	36	VAL
1	H	46	THR
1	H	62	ARG
1	H	81	ASN
1	H	82	VAL
1	H	82(A)	PRO
1	H	92	ASN
1	H	97	CYS
1	H	112	LEU
1	H	114	ARG
1	H	129	ASP
1	H	142	GLN
1	H	166	ARG
1	H	182	ILE

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Mol	Chain	Res	Type
1	H	185	SER
1	H	194	LEU
1	H	195	TYR
1	H	197	ASN
1	H	208	THR
1	H	222	LYS
1	H	227	SER
1	H	292	MET
1	H	296	ASN
1	H	302	ILE
1	H	310	LYS
1	H	315	VAL
1	H	316	LEU
2	I	11	GLU
2	I	32	SER
2	I	41	THR
2	I	42	GLN
2	I	48	VAL
2	I	50	ASN
2	I	54	SER
2	I	60	ASN
2	I	61	THR
2	I	72	ASN
2	I	81	ASN
2	I	84	MET
2	I	89	LEU
2	I	93	THR
2	I	101	LEU
2	I	105	GLU
2	I	106	ARG
2	I	108	LEU
2	I	124	LEU
2	I	127	ARG
2	I	133	LEU
2	I	143	LYS
2	I	145	ASP
2	I	152	VAL
2	I	154	ASN
2	I	156	THR
2	I	157	TYR
2	I	164	GLU
2	I	165	GLU

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Mol	Chain	Res	Type
2	I	170	ARG
1	J	22	THR
1	J	27	THR
1	J	66	VAL
1	J	71	LEU
1	J	75	MET
1	J	82	VAL
1	J	82(A)	PRO
1	J	89	GLU
1	J	114	ARG
1	J	151	VAL
1	J	158	ASN
1	J	176	LEU
1	J	186	ASN
1	J	191	GLN
1	J	195	TYR
1	J	196	GLN
1	J	197	ASN
1	J	208	THR
1	J	211	GLN
1	J	217	ILE
1	J	220	ARG
1	J	221	SER
1	J	222	LYS
1	J	227	SER
1	J	238	LYS
1	J	268	ILE
1	J	273	GLU
1	J	277	CYS
1	J	279	THR
1	J	283	THR
1	J	311	SER
1	J	314	LEU
1	J	320	LEU
2	K	19	ASP
2	K	24	TYR
2	K	32	SER
2	K	38	LYS
2	K	39	GLU
2	K	41	THR
2	K	43	LYS
2	K	48	VAL

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Mol	Chain	Res	Type
2	K	49	THR
2	K	57	ASP
2	K	72	ASN
2	K	80	LEU
2	K	83	LYS
2	K	84	MET
2	K	86	ASP
2	K	91	VAL
2	K	99	LEU
2	K	101	LEU
2	K	108	LEU
2	K	132	GLU
2	K	153	ARG
2	K	156	THR
2	K	161	GLN
2	K	164	GLU
1	L	22	THR
1	L	45	LYS
1	L	46	THR
1	L	54	ASP
1	L	63	ASP
1	L	83	GLU
1	L	90	LYS
1	L	123	ILE
1	L	136	SER
1	L	142	GLN
1	L	144	THR
1	L	145	PRO
1	L	173	GLN
1	L	185	SER
1	L	186	ASN
1	L	191	GLN
1	L	219	THR
1	L	220	ARG
1	L	221	SER
1	L	235	THR
1	L	268	ILE
1	L	280	LYS
1	L	283	THR
1	L	291	SER
1	L	302	ILE
1	L	311	SER

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Mol	Chain	Res	Type
1	L	315	VAL
1	L	320	LEU
2	M	15	GLN
2	M	19	ASP
2	M	21	TRP
2	M	24	TYR
2	M	25	HIS
2	M	30	GLN
2	M	42	GLN
2	M	43	LYS
2	M	45	ILE
2	M	51	LYS
2	M	60	ASN
2	M	66	VAL
2	M	69	GLU
2	M	72	ASN
2	M	75	ARG
2	M	84	MET
2	M	89	LEU
2	M	108	LEU
2	M	128	ASP
2	M	129	ASN
2	M	139	GLU
2	M	145	ASP
2	M	154	ASN
2	M	157	TYR
2	M	161	GLN
2	M	163	SER
2	M	164	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (115) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	GLN
1	A	18	HIS
1	A	95	ASN
1	A	122	GLN
1	A	142	GLN
1	A	173	GLN
1	A	186	ASN
1	A	197	ASN
1	A	224	ASN

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Mol	Chain	Res	Type
1	A	240	ASN
1	A	244	ASN
1	A	295	HIS
1	A	322	ASN
2	B	26	HIS
2	B	42	GLN
2	B	50	ASN
2	B	53	ASN
2	B	60	ASN
2	B	79	ASN
2	B	95	ASN
2	B	111	HIS
2	B	117	ASN
2	B	125	GLN
2	B	161	GLN
1	C	18	HIS
1	C	81	ASN
1	C	92	ASN
1	C	95	ASN
1	C	186	ASN
1	C	197	ASN
1	C	211	GLN
1	C	276	ASN
1	C	322	ASN
2	D	50	ASN
2	D	79	ASN
2	D	95	ASN
2	D	117	ASN
2	D	125	GLN
2	D	129	ASN
2	D	135	ASN
1	E	24	GLN
1	E	34	ASN
1	E	38	HIS
1	E	95	ASN
1	E	117	HIS
1	E	130	HIS
1	E	142	GLN
1	E	186	ASN
1	E	197	ASN
1	E	226	GLN
1	E	240	ASN

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Mol	Chain	Res	Type
1	E	276	ASN
1	E	295	HIS
2	F	15	GLN
2	F	25	HIS
2	F	26	HIS
2	F	42	GLN
2	F	60	ASN
2	F	79	ASN
2	F	104	ASN
2	F	111	HIS
1	H	18	HIS
1	H	34	ASN
1	H	110	HIS
1	H	142	GLN
1	H	173	GLN
1	H	196	GLN
1	H	224	ASN
1	H	240	ASN
1	H	276	ASN
1	H	295	HIS
1	H	296	ASN
1	H	312	ASN
2	I	26	HIS
2	I	42	GLN
2	I	53	ASN
2	I	60	ASN
2	I	79	ASN
2	I	95	ASN
2	I	111	HIS
2	I	117	ASN
2	I	129	ASN
2	I	142	HIS
2	I	154	ASN
1	J	20	ASN
1	J	34	ASN
1	J	95	ASN
1	J	197	ASN
1	J	224	ASN
1	J	240	ASN
1	J	295	HIS
2	K	26	HIS
2	K	50	ASN

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Mol	Chain	Res	Type
2	K	62	GLN
2	K	95	ASN
2	K	117	ASN
2	K	125	GLN
1	L	20	ASN
1	L	24	GLN
1	L	38	HIS
1	L	81	ASN
1	L	130	HIS
1	L	142	GLN
1	L	172	ASN
1	L	173	GLN
1	L	186	ASN
1	L	197	ASN
1	L	295	HIS
1	L	322	ASN
2	M	25	HIS
2	M	30	GLN
2	M	42	GLN
2	M	79	ASN
2	M	104	ASN
2	M	111	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 ⓘ

8 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	G	1	3,1	14,14,15	0.86	0	17,19,21	1.58	2 (11%)
3	NAG	G	2	3	14,14,15	0.75	0	17,19,21	1.79	3 (17%)
3	NAG	N	1	3,1	14,14,15	0.75	0	17,19,21	2.85	6 (35%)
3	NAG	N	2	3	14,14,15	0.58	0	17,19,21	1.06	2 (11%)
3	NAG	O	1	3,1	14,14,15	1.01	1 (7%)	17,19,21	2.77	4 (23%)
3	NAG	O	2	3	14,14,15	0.42	0	17,19,21	2.18	7 (41%)
3	NAG	P	1	3,1	14,14,15	0.55	0	17,19,21	1.82	5 (29%)
3	NAG	P	2	3	14,14,15	0.57	0	17,19,21	1.06	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
3	NAG	G	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	G	2	3	-	1/6/23/26	0/1/1/1
3	NAG	N	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	N	2	3	-	1/6/23/26	0/1/1/1
3	NAG	O	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	O	2	3	-	4/6/23/26	0/1/1/1
3	NAG	P	1	3,1	-	2/6/23/26	0/1/1/1
3	NAG	P	2	3	-	1/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	O	1	NAG	O4-C4	2.97	1.50	1.43

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	O	1	NAG	O4-C4-C5	-9.38	86.23	109.32
3	N	1	NAG	C2-N2-C7	-8.02	112.16	122.90
3	N	1	NAG	O4-C4-C3	5.35	122.98	110.38
3	O	2	NAG	C1-O5-C5	4.36	118.02	112.19
3	G	2	NAG	C2-N2-C7	4.35	128.74	122.90
3	G	1	NAG	C2-N2-C7	-4.32	117.12	122.90
3	P	1	NAG	C1-O5-C5	3.91	117.42	112.19

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	N	1	NAG	C1-C2-N2	-3.58	104.80	110.43
3	O	1	NAG	O5-C5-C6	3.57	114.62	107.66
3	O	2	NAG	C3-C4-C5	3.57	116.70	110.23
3	N	1	NAG	O4-C4-C5	-3.41	100.93	109.32
3	O	1	NAG	C4-C3-C2	-3.30	106.18	111.02
3	O	2	NAG	O5-C1-C2	-3.15	106.42	111.29
3	O	2	NAG	C8-C7-N2	2.92	120.96	116.12
3	O	2	NAG	O5-C5-C4	2.85	117.77	110.83
3	O	2	NAG	O7-C7-C8	-2.83	117.02	122.05
3	O	1	NAG	C1-O5-C5	2.81	115.95	112.19
3	N	2	NAG	C4-C3-C2	-2.81	106.91	111.02
3	P	2	NAG	C4-C3-C2	-2.81	106.91	111.02
3	G	1	NAG	O5-C1-C2	-2.68	107.15	111.29
3	P	1	NAG	C4-C3-C2	-2.67	107.11	111.02
3	G	2	NAG	C1-O5-C5	2.65	115.73	112.19
3	O	2	NAG	C1-C2-N2	2.63	114.57	110.43
3	P	1	NAG	O3-C3-C2	-2.57	104.06	109.40
3	P	1	NAG	C1-C2-N2	2.57	114.48	110.43
3	P	2	NAG	O5-C1-C2	-2.21	107.87	111.29
3	G	2	NAG	C4-C3-C2	2.19	114.22	111.02
3	N	2	NAG	O5-C1-C2	-2.17	107.93	111.29
3	N	1	NAG	C1-O5-C5	2.13	115.04	112.19
3	N	1	NAG	O3-C3-C4	2.03	115.17	110.38
3	P	1	NAG	O7-C7-C8	-2.01	118.47	122.05

There are no chirality outliers.

All (15) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	G	1	NAG	O5-C5-C6-O6
3	N	1	NAG	O5-C5-C6-O6
3	G	1	NAG	C4-C5-C6-O6
3	O	1	NAG	O5-C5-C6-O6
3	N	1	NAG	C4-C5-C6-O6
3	P	1	NAG	O5-C5-C6-O6
3	O	2	NAG	C8-C7-N2-C2
3	O	2	NAG	O7-C7-N2-C2
3	O	1	NAG	C4-C5-C6-O6
3	O	2	NAG	C4-C5-C6-O6
3	N	2	NAG	O5-C5-C6-O6
3	P	2	NAG	O5-C5-C6-O6
3	P	1	NAG	C4-C5-C6-O6

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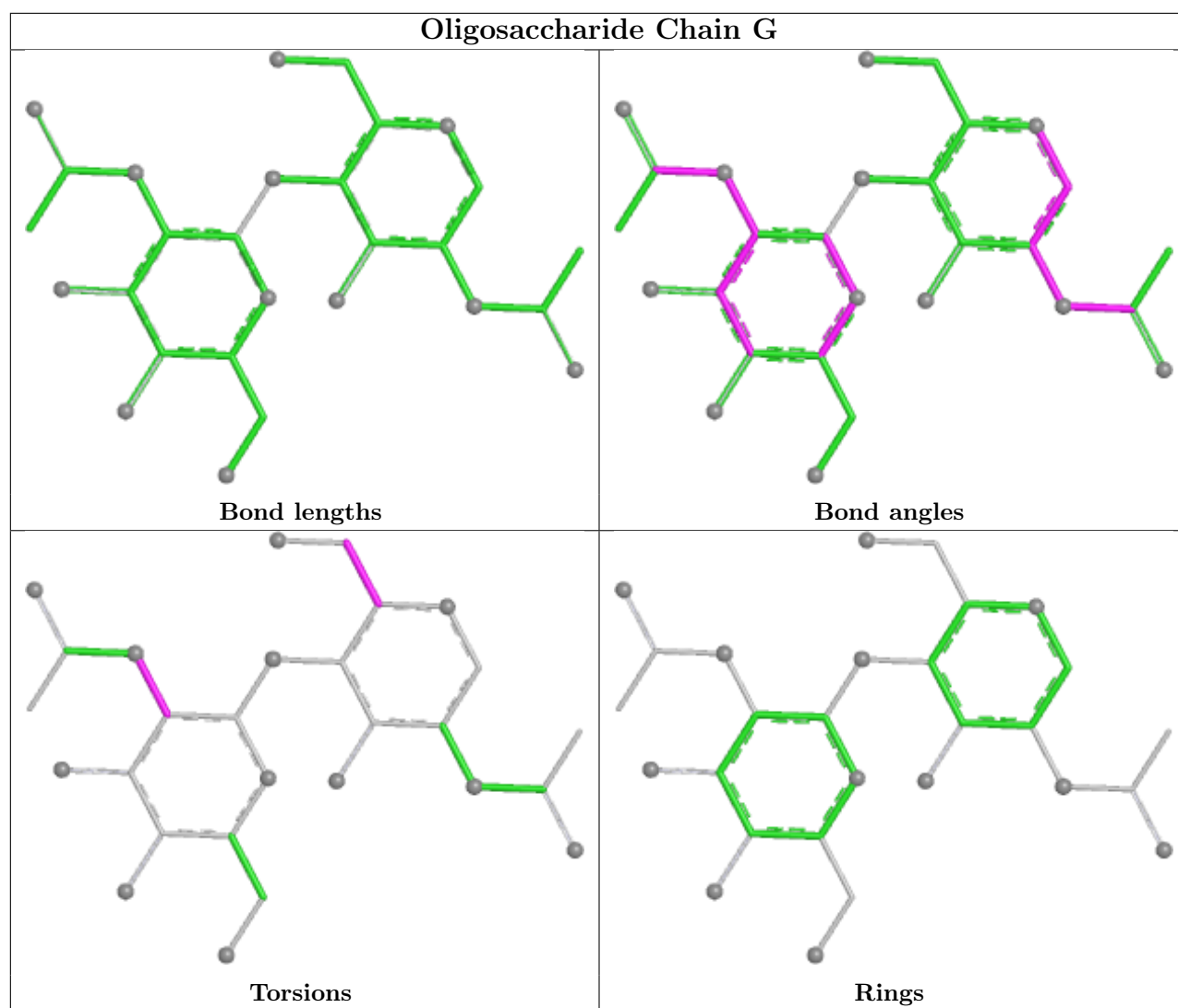
Mol	Chain	Res	Type	Atoms
3	G	2	NAG	C3-C2-N2-C7
3	O	2	NAG	O5-C5-C6-O6

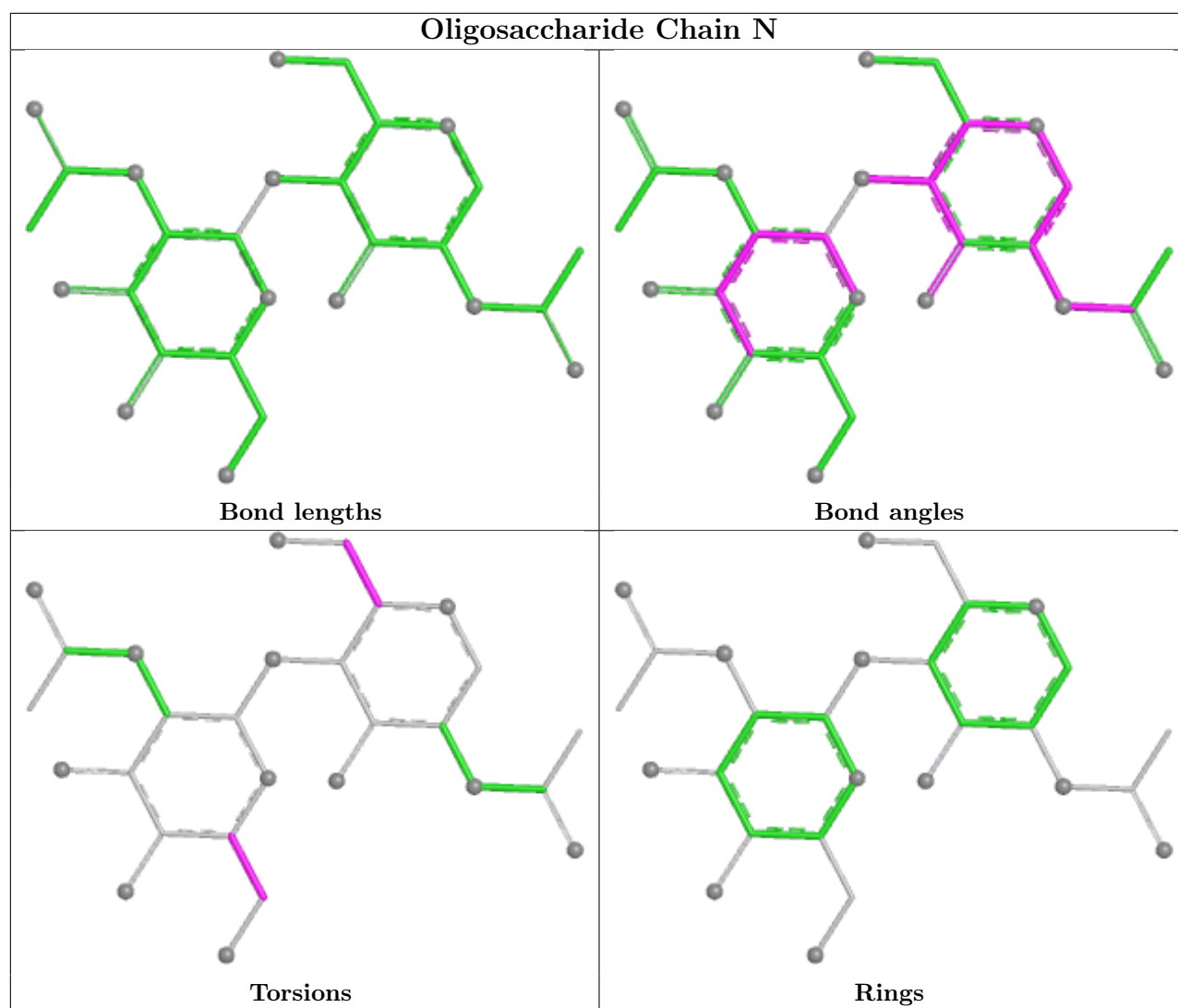
There are no ring outliers.

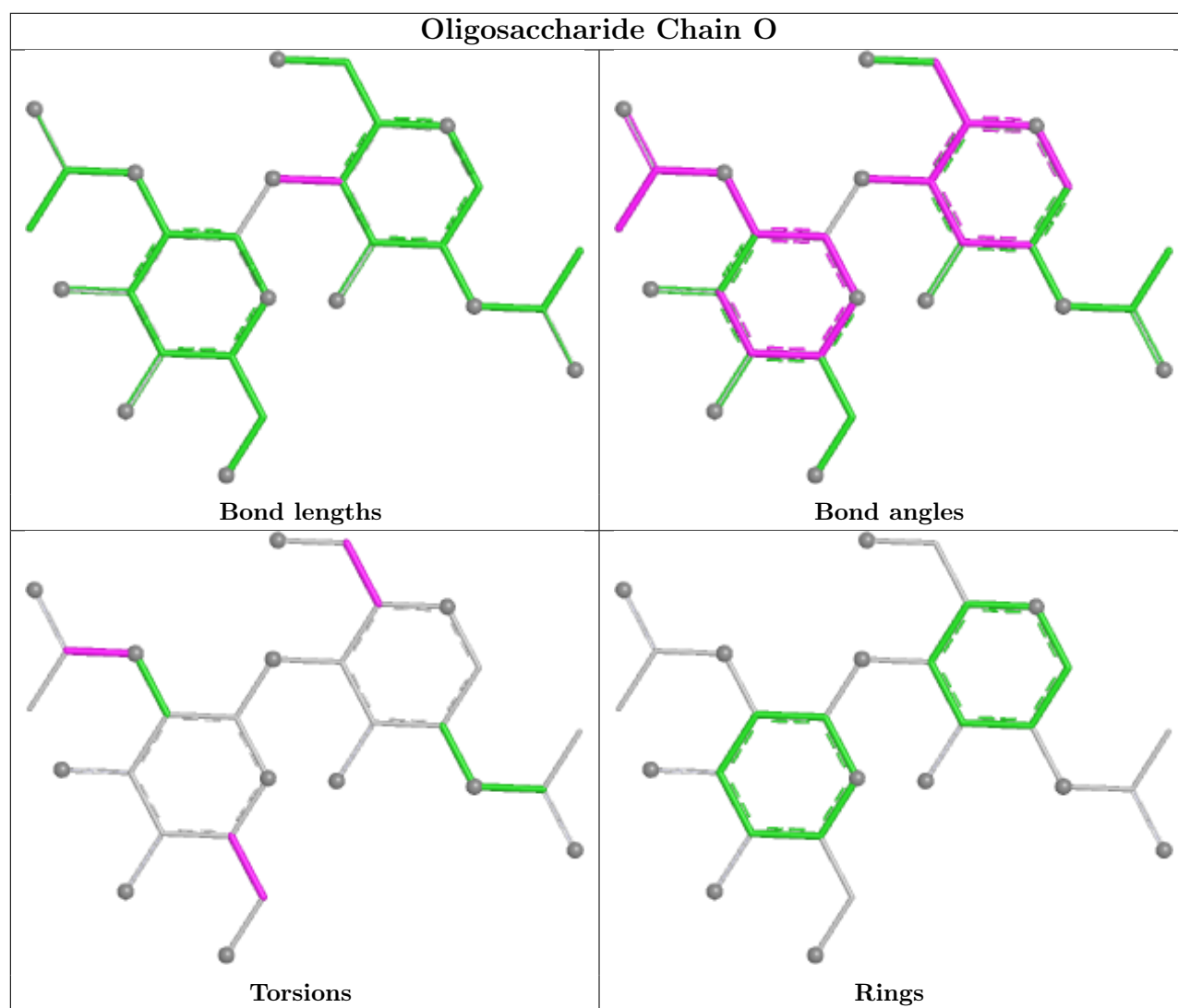
7 monomers are involved in 11 short contacts:

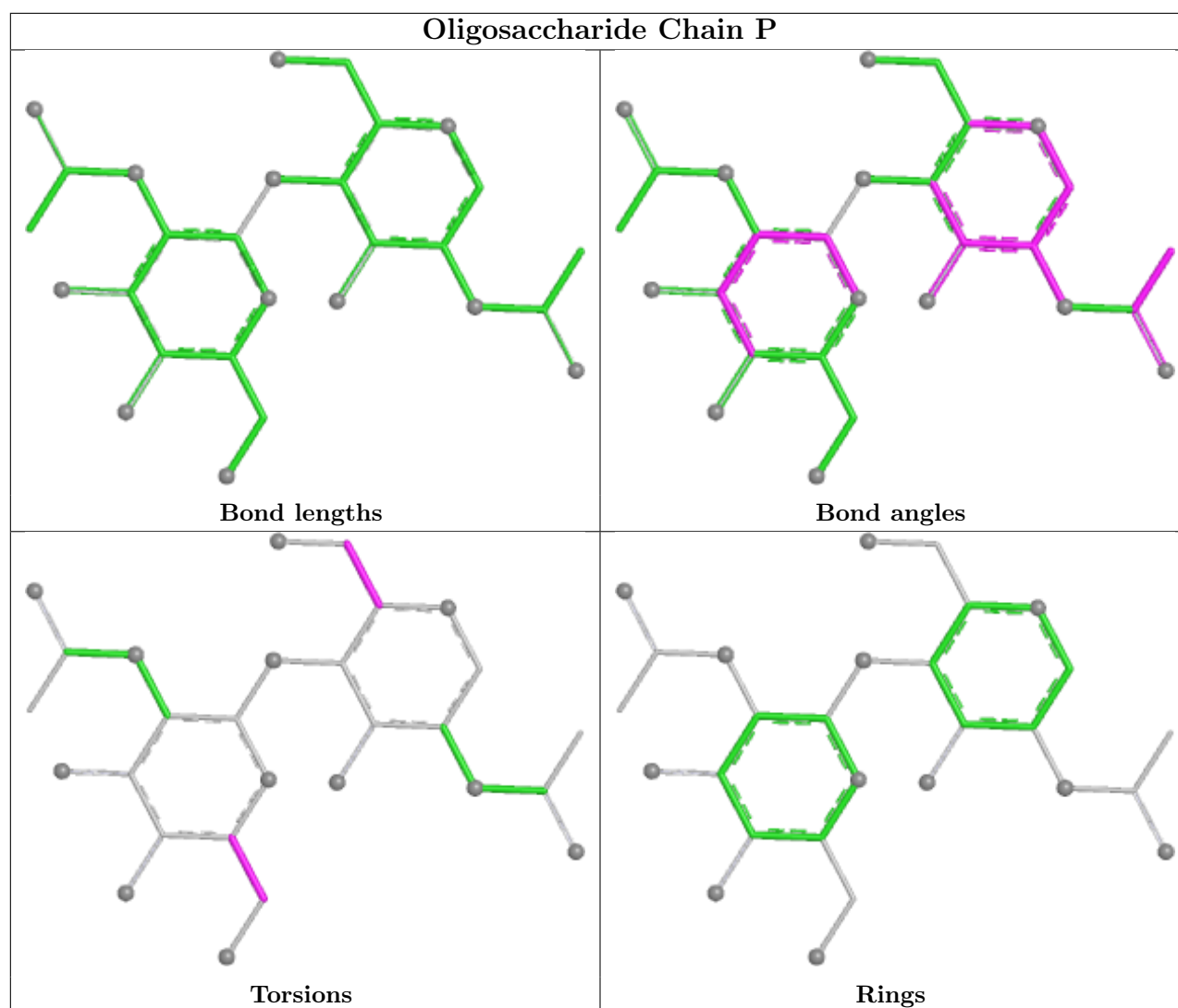
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	O	1	NAG	2	0
3	N	1	NAG	3	0
3	G	2	NAG	1	0
3	P	1	NAG	2	0
3	G	1	NAG	2	0
3	N	2	NAG	2	0
3	P	2	NAG	2	0

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









5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	NAG	J	401	1	14,14,15	0.59	0	17,19,21	1.22	1 (5%)
4	NAG	H	401	1	14,14,15	0.51	0	17,19,21	1.50	2 (11%)
4	NAG	L	402	1	14,14,15	0.66	0	17,19,21	1.52	4 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	NAG	A	401	1	14,14,15	0.50	0	17,19,21	1.51	3 (17%)
4	NAG	C	402	1	14,14,15	0.74	1 (7%)	17,19,21	2.03	8 (47%)
4	NAG	C	401	1	14,14,15	0.58	0	17,19,21	1.64	4 (23%)
4	NAG	J	402	1	14,14,15	0.56	0	17,19,21	0.74	0
4	NAG	E	401	1	14,14,15	0.60	0	17,19,21	1.64	3 (17%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	NAG	J	401	1	-	2/6/23/26	0/1/1/1
4	NAG	H	401	1	-	2/6/23/26	0/1/1/1
4	NAG	L	402	1	-	2/6/23/26	0/1/1/1
4	NAG	A	401	1	-	2/6/23/26	0/1/1/1
4	NAG	C	402	1	-	2/6/23/26	0/1/1/1
4	NAG	C	401	1	-	2/6/23/26	0/1/1/1
4	NAG	J	402	1	-	0/6/23/26	0/1/1/1
4	NAG	E	401	1	-	2/6/23/26	0/1/1/1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	402	NAG	O5-C1	-2.24	1.39	1.43

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	401	NAG	C1-O5-C5	4.26	117.89	112.19
4	H	401	NAG	C1-O5-C5	4.20	117.81	112.19
4	C	401	NAG	C2-N2-C7	-4.14	117.35	122.90
4	C	402	NAG	C3-C4-C5	3.66	116.88	110.23
4	E	401	NAG	C1-O5-C5	3.62	117.04	112.19
4	L	402	NAG	C1-O5-C5	3.51	116.89	112.19
4	E	401	NAG	O5-C5-C6	3.25	114.00	107.66
4	C	402	NAG	O5-C1-C2	-3.22	106.31	111.29
4	J	401	NAG	C1-O5-C5	3.15	116.41	112.19
4	C	402	NAG	C1-C2-N2	-2.88	105.89	110.43

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	C	402	NAG	C4-C3-C2	2.80	115.11	111.02
4	L	402	NAG	C2-N2-C7	2.61	126.40	122.90
4	C	402	NAG	O3-C3-C4	-2.61	104.23	110.38
4	A	401	NAG	C3-C4-C5	2.51	114.79	110.23
4	H	401	NAG	C3-C4-C5	2.49	114.74	110.23
4	C	401	NAG	O7-C7-C8	-2.32	117.93	122.05
4	C	401	NAG	C1-O5-C5	2.28	115.24	112.19
4	L	402	NAG	O5-C5-C6	2.24	112.02	107.66
4	E	401	NAG	O3-C3-C2	2.24	114.05	109.40
4	C	402	NAG	C1-O5-C5	2.23	115.18	112.19
4	C	402	NAG	O5-C5-C4	2.20	116.17	110.83
4	C	402	NAG	C2-N2-C7	2.19	125.83	122.90
4	L	402	NAG	C1-C2-N2	2.11	113.75	110.43
4	C	401	NAG	O4-C4-C3	2.08	115.27	110.38
4	A	401	NAG	O5-C1-C2	-2.01	108.18	111.29

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	401	NAG	O5-C5-C6-O6
4	H	401	NAG	O5-C5-C6-O6
4	A	401	NAG	C4-C5-C6-O6
4	E	401	NAG	C4-C5-C6-O6
4	H	401	NAG	C4-C5-C6-O6
4	J	401	NAG	O5-C5-C6-O6
4	C	402	NAG	O5-C5-C6-O6
4	L	402	NAG	C4-C5-C6-O6
4	L	402	NAG	O5-C5-C6-O6
4	J	401	NAG	C4-C5-C6-O6
4	C	401	NAG	C4-C5-C6-O6
4	E	401	NAG	O5-C5-C6-O6
4	C	402	NAG	C4-C5-C6-O6
4	C	401	NAG	O5-C5-C6-O6

There are no ring outliers.

6 monomers are involved in 17 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	J	401	NAG	7	0
4	H	401	NAG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L	402	NAG	2	0
4	A	401	NAG	2	0
4	J	402	NAG	1	0
4	E	401	NAG	4	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	321/329 (97%)	-0.39	1 (0%) 90 86	19, 36, 54, 81	0
1	C	321/329 (97%)	-0.31	3 (0%) 81 74	20, 39, 55, 71	0
1	E	322/329 (97%)	-0.33	1 (0%) 90 86	19, 36, 57, 93	0
1	H	321/329 (97%)	-0.25	2 (0%) 85 80	20, 40, 59, 80	0
1	J	321/329 (97%)	-0.30	2 (0%) 85 80	20, 38, 55, 65	0
1	L	319/329 (96%)	-0.18	2 (0%) 85 80	20, 40, 65, 95	0
2	B	170/182 (93%)	0.31	8 (4%) 36 29	17, 58, 92, 104	0
2	D	172/182 (94%)	0.00	2 (1%) 76 68	23, 54, 72, 84	0
2	F	172/182 (94%)	0.22	3 (1%) 69 60	18, 55, 95, 102	0
2	I	170/182 (93%)	0.13	2 (1%) 76 68	18, 56, 84, 91	0
2	K	172/182 (94%)	-0.07	2 (1%) 76 68	19, 50, 75, 99	0
2	M	169/182 (92%)	0.41	13 (7%) 19 14	17, 63, 106, 122	0
All	All	2950/3066 (96%)	-0.13	41 (1%) 73 64	17, 41, 81, 122	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	9	PRO	4.4
1	C	97	CYS	3.9
2	I	134	GLY	3.3
1	H	126	SER	3.2
1	L	13	ILE	3.2
2	M	160	PRO	3.1
2	M	141	TYR	3.0
1	J	218	ALA	3.0
2	M	140	PHE	3.0
2	F	168	LEU	2.9
2	M	168	LEU	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	168	LEU	2.7
2	M	138	PHE	2.7
2	I	36	ALA	2.7
2	M	169	LYS	2.6
1	C	133(A)	SER	2.6
2	M	33	GLY	2.6
2	B	31	GLY	2.6
2	B	157	TYR	2.6
1	H	10	GLY	2.6
1	C	319	GLY	2.5
2	M	144	CYS	2.5
2	K	124	LEU	2.5
1	J	196	GLN	2.5
2	F	152	VAL	2.4
2	M	131	LYS	2.4
2	K	164	GLU	2.4
2	D	135	ASN	2.3
1	L	97	CYS	2.3
2	D	22	TYR	2.3
2	B	159	TYR	2.2
2	M	156	THR	2.1
2	B	7	ALA	2.1
2	B	11	GLU	2.1
2	B	33	GLY	2.1
2	M	159	TYR	2.1
2	B	152	VAL	2.1
2	F	35	ALA	2.0
2	M	143	LYS	2.0
1	A	323	SER	2.0
2	M	134	GLY	2.0

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

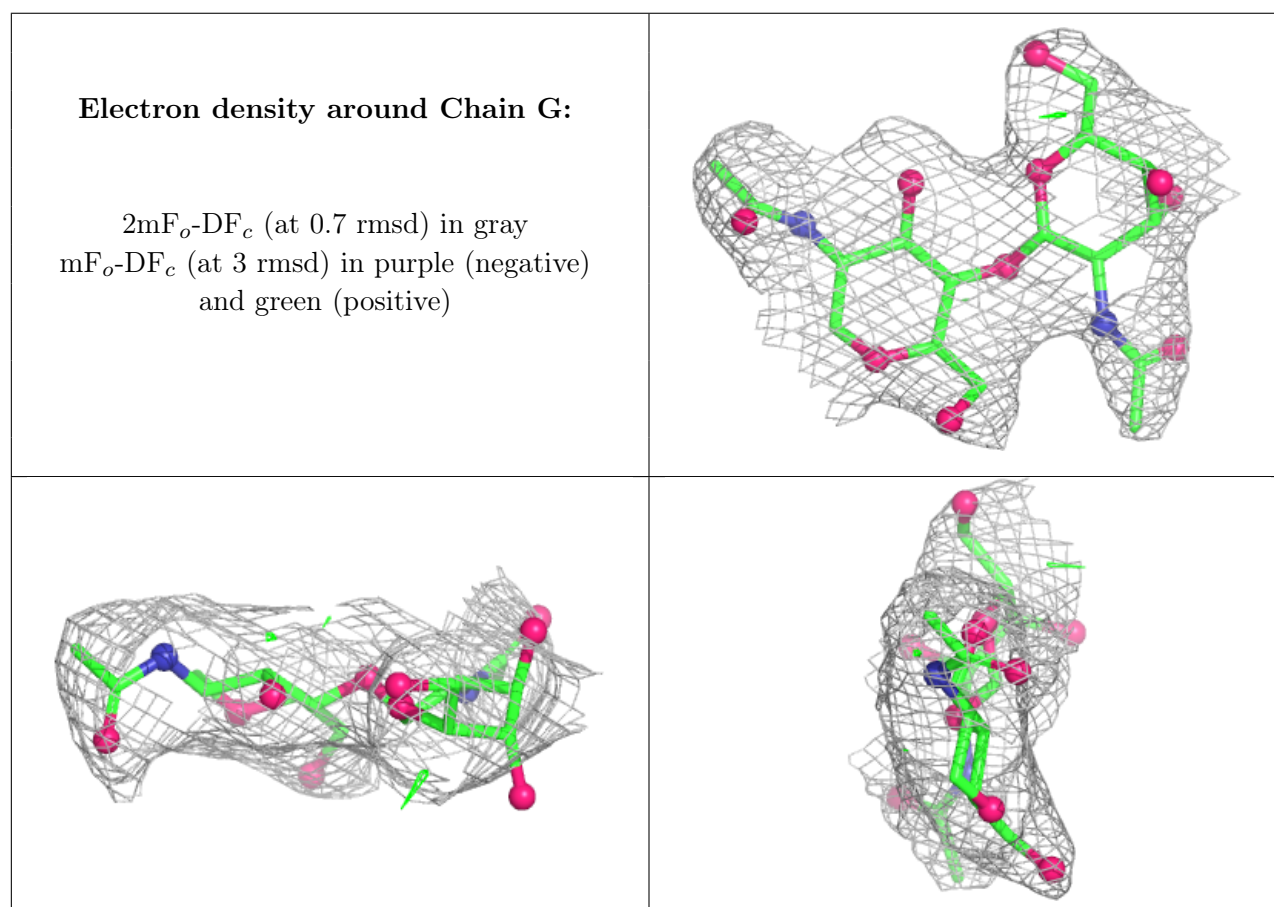
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

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

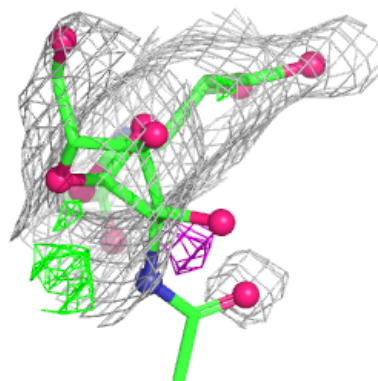
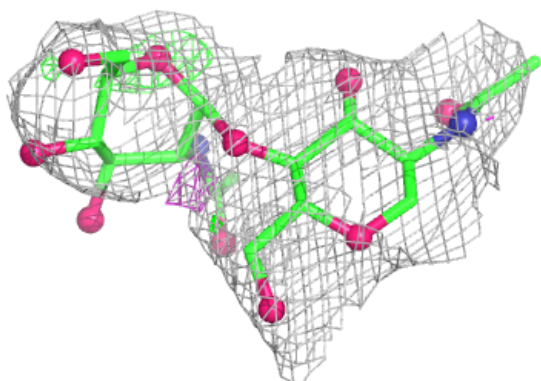
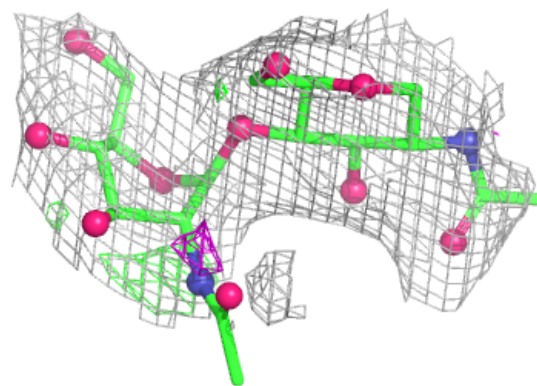
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	NAG	G	1	14/15	-	-	33,47,56,61	0
3	NAG	G	2	14/15	-	-	62,75,81,93	0
3	NAG	O	2	14/15	0.54	0.19	81,92,100,104	0
3	NAG	N	2	14/15	0.58	0.18	91,97,107,108	0
3	NAG	P	2	14/15	0.63	0.20	85,93,108,114	0
3	NAG	P	1	14/15	0.87	0.14	57,65,83,84	0
3	NAG	N	1	14/15	0.88	0.12	49,55,61,71	0
3	NAG	O	1	14/15	0.91	0.10	54,64,79,80	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.

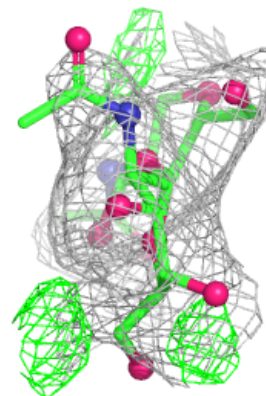
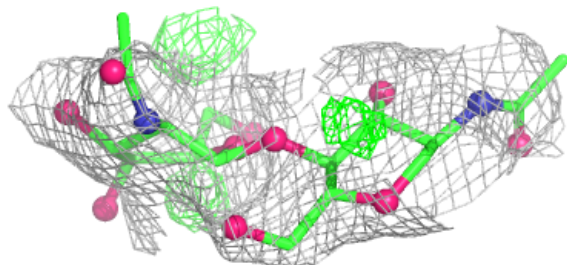
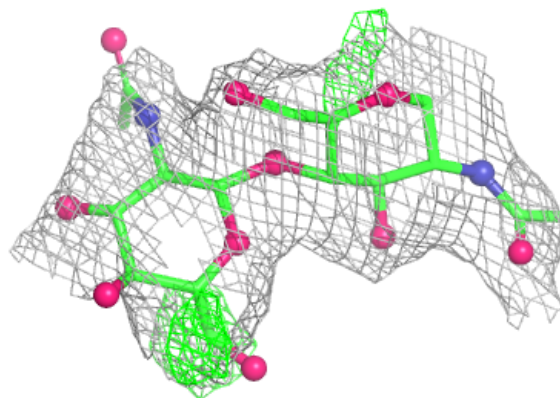


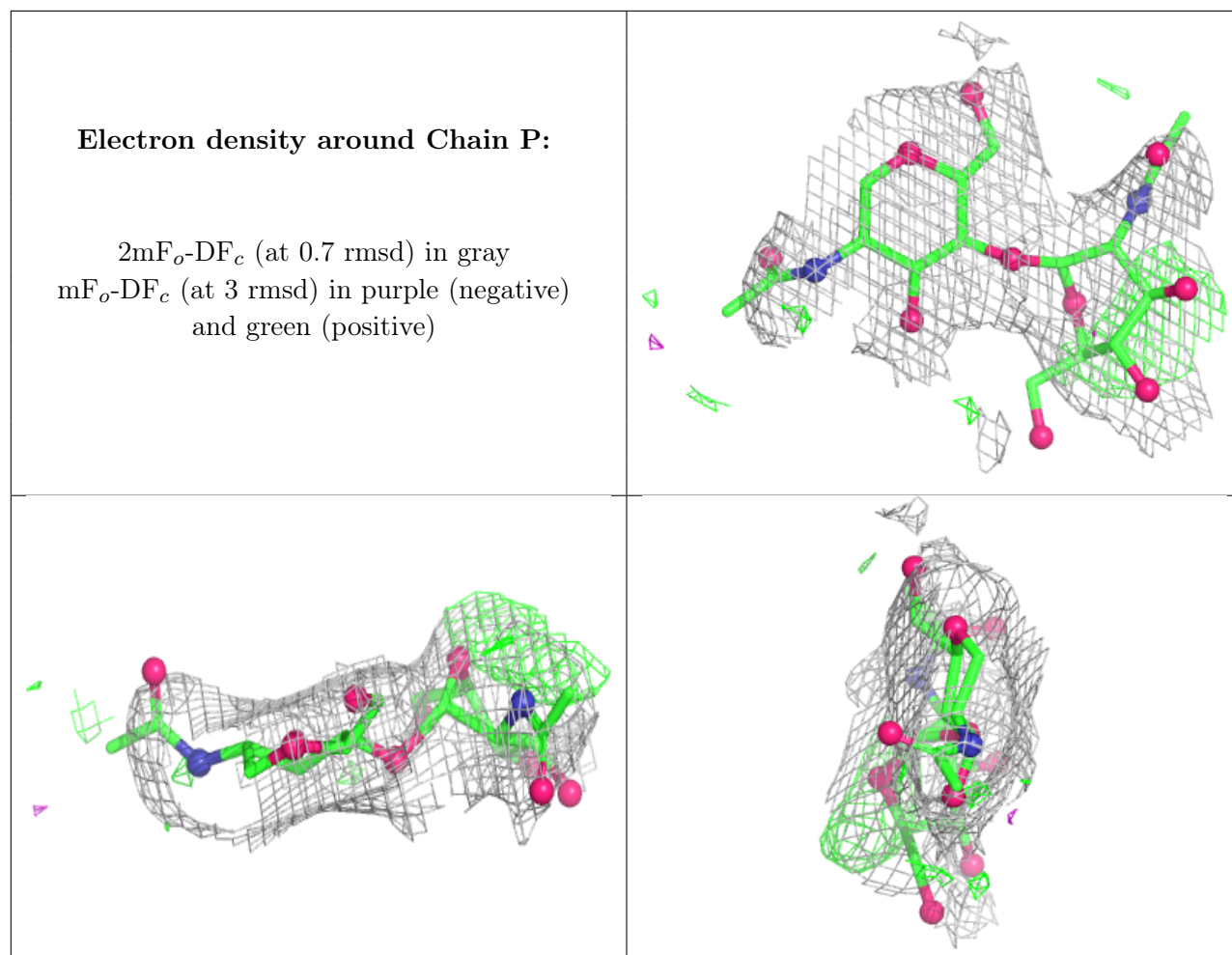
Electron density around Chain N:

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

**Electron density around Chain O:**

$2mF_o-DF_c$ (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
4	NAG	J	401	14/15	0.53	0.20	94,106,111,113	0
4	NAG	H	401	14/15	0.70	0.15	91,99,112,113	0
4	NAG	L	402	14/15	0.70	0.18	78,88,95,97	0
4	NAG	E	401	14/15	0.73	0.13	71,84,90,92	0
4	NAG	J	402	14/15	0.74	0.17	39,55,68,68	0
4	NAG	C	401	14/15	0.75	0.13	82,88,98,98	0
4	NAG	A	401	14/15	0.76	0.14	89,95,105,119	0
4	NAG	C	402	14/15	0.88	0.12	52,59,64,73	0

6.5 Other polymers [i](#)

There are no such residues in this entry.