



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 10:34 AM UTC

PDB ID : 3SQ9 / pdb_00003sq9
Title : Crystal Structures of the Ligand Binding Domain of a Pentameric Alpha7 Nicotinic Receptor Chimera
Authors : Li, S.-X.; Huang, S.; Bren, N.; Noridomi, K.; Dellisanti, C.; Sine, S.; Chen, L.
Deposited on : 2011-07-05
Resolution : 3.10 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

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

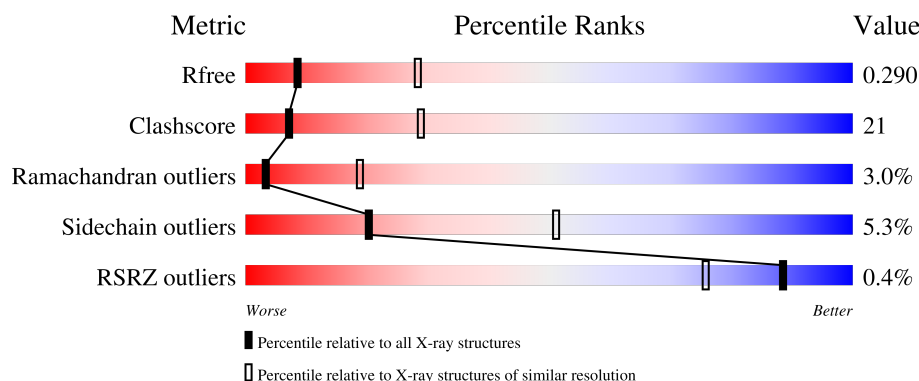
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 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	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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	204	 65% 27% . .
1	B	204	 61% 30% 6% . .
1	C	204	 60% 34% . .
1	D	204	 62% 30% 5% . .
1	E	204	 63% 30% . .

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Mol	Chain	Length	Quality of chain
1	F	204	64%29%
1	G	204	61%31%5%
1	H	204	% 63%31%
1	I	204	62%30%5%
1	J	204	61%32%
2	K	2	100%
2	L	2	100%
2	M	2	50%50%

2 Entry composition [i](#)

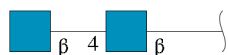
There are 3 unique types of molecules in this entry. The entry contains 16750 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	B	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	C	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	D	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	E	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	F	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	G	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	H	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	I	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			
1	J	202	Total	C	N	O	S	0	0	0
			1647	1055	276	309	7			

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



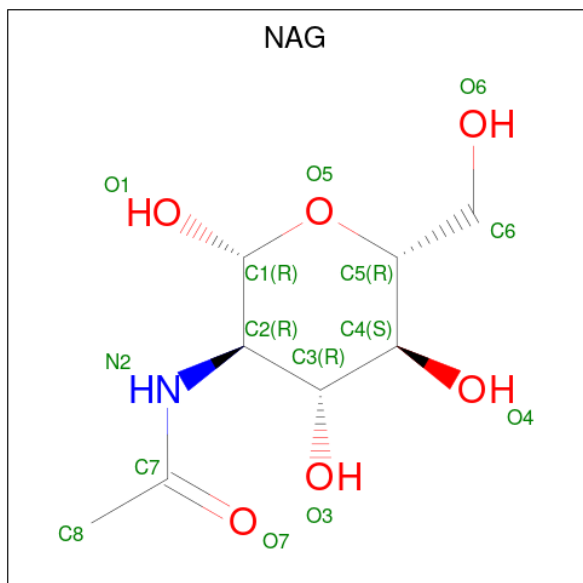
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	K	2	Total	C	N	O	0	0	0
			28	16	2	10			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	L	2	Total	C	N	O	0	0	0
			28	16	2	10			
2	M	2	Total	C	N	O	0	0	0
			28	16	2	10			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	C	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		

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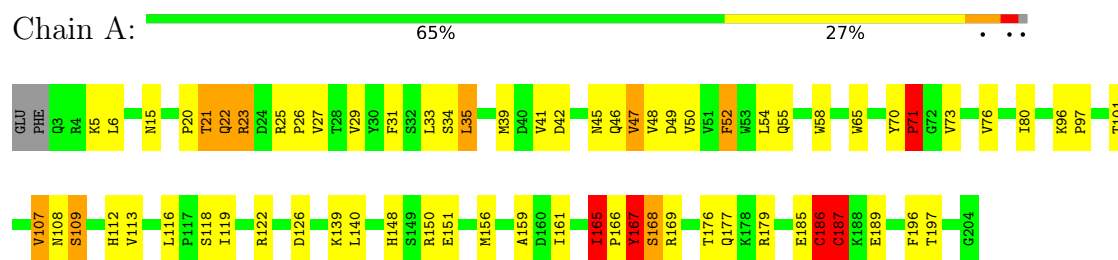
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	H	1	Total	C	N	O	0	0
			14	8	1	5		
3	I	1	Total	C	N	O	0	0
			14	8	1	5		
3	J	1	Total	C	N	O	0	0
			14	8	1	5		
3	J	1	Total	C	N	O	0	0
			14	8	1	5		

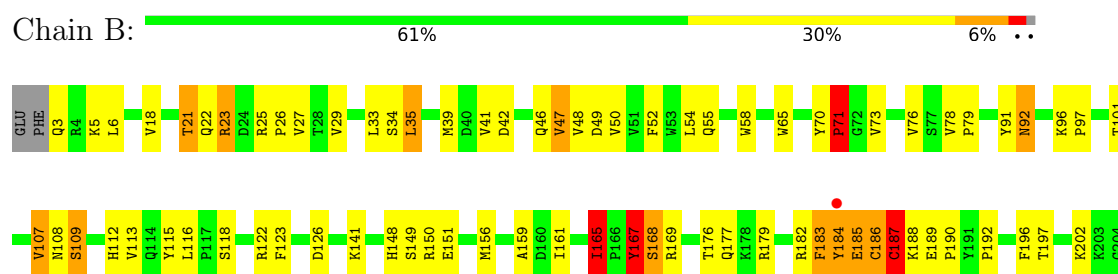
3 Residue-property plots

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.

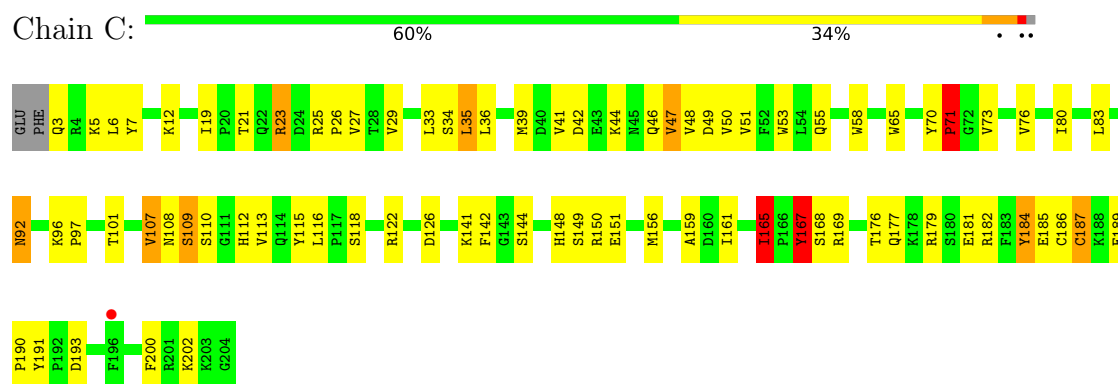
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein



- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein

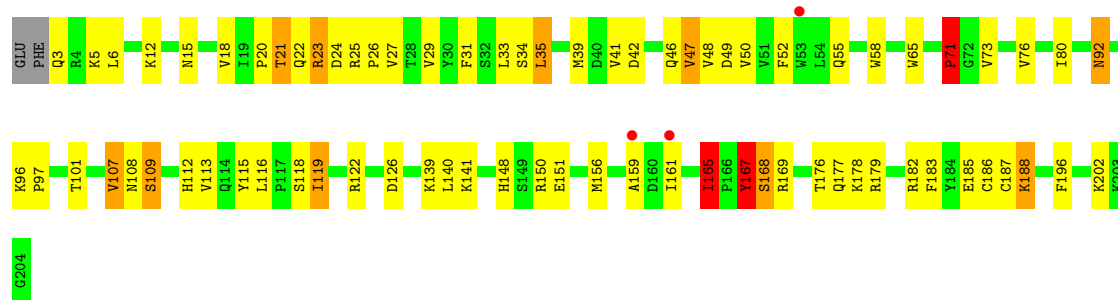


- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein



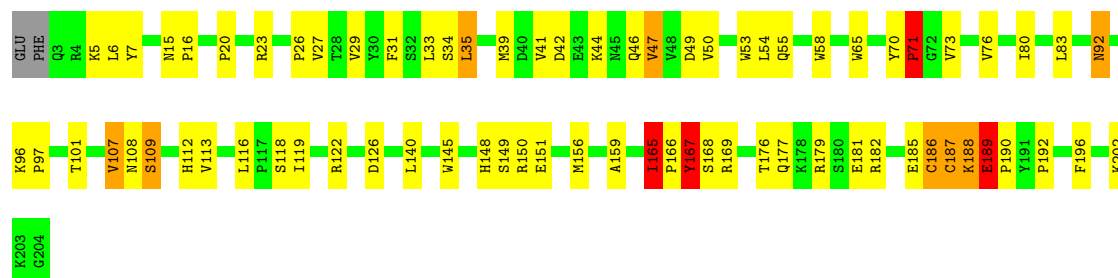
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein





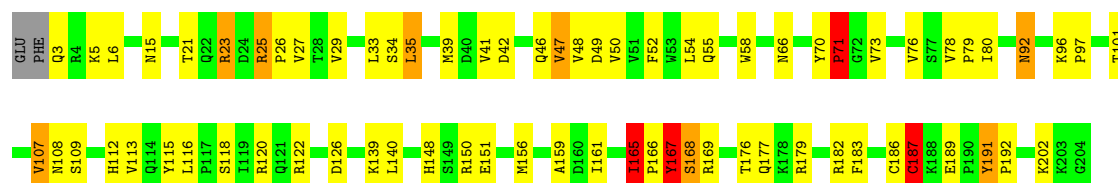
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein

Chain E: 63% 30% . . .



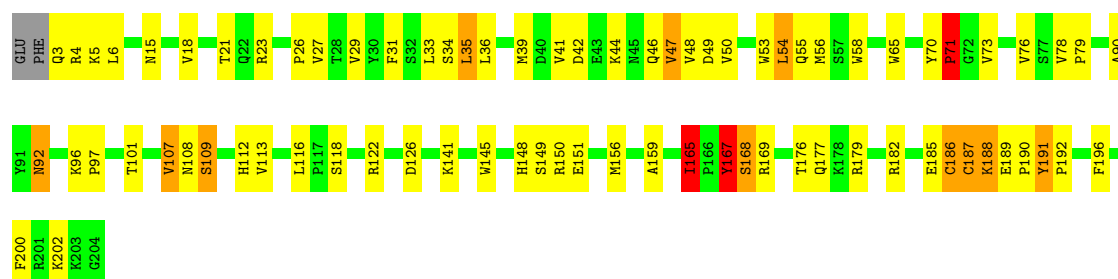
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein

Chain F: 64% 29% . . .



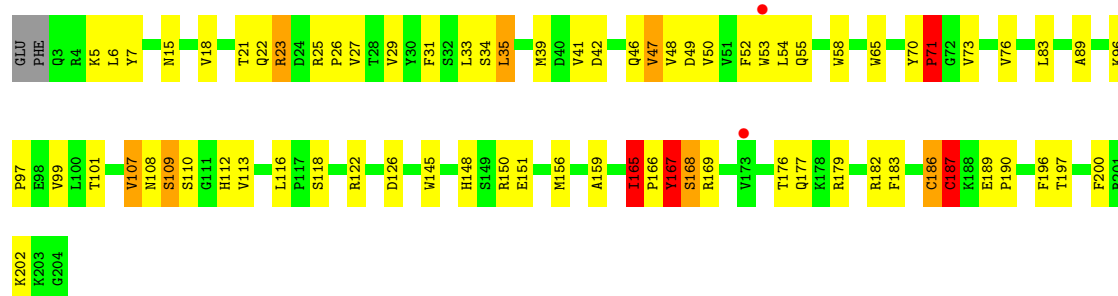
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein

Chain G: 61% 31% 5% . .



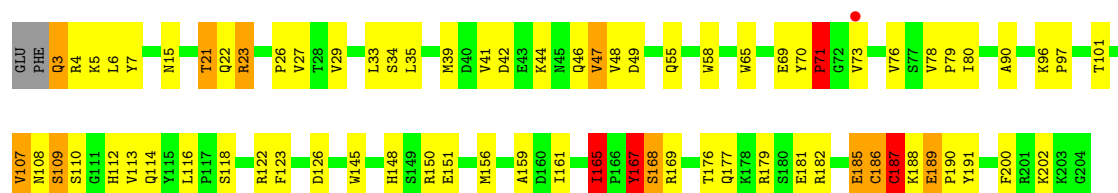
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein

Chain H: 63% 31% . . .



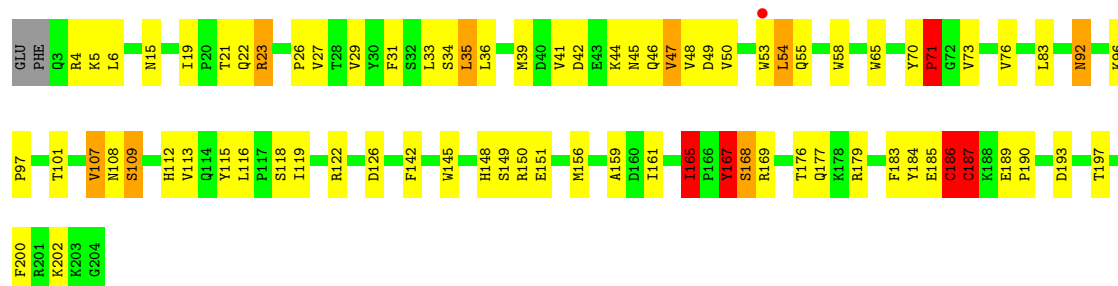
- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein

Chain I: 62% 30% 5% ..



- Molecule 1: Neuronal acetylcholine receptor subunit alpha-7, Acetylcholine-binding protein

Chain J: 61% 32% . ..



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

Chain K: 100%



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

Chain L: 100%



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

Chain M:  50% 50%

MAG1
MAG2

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	79.12Å 144.56Å 131.12Å 90.00° 102.46° 90.00°	Depositor
Resolution (Å)	46.45 – 3.10 46.45 – 3.10	Depositor EDS
% Data completeness (in resolution range)	90.4 (46.45-3.10) 90.4 (46.45-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 3.12Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.262 , 0.289 0.264 , 0.290	Depositor DCC
R_{free} test set	4737 reflections (10.05%)	wwPDB-VP
Wilson B-factor (Å ²)	79.5	Xtriage
Anisotropy	0.656	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	16750	wwPDB-VP
Average B, all atoms (Å ²)	110.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.17% 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	0.53	0/1691	0.99	6/2300 (0.3%)
1	B	0.51	0/1691	0.97	6/2300 (0.3%)
1	C	0.54	0/1691	1.00	9/2300 (0.4%)
1	D	0.54	0/1691	0.96	9/2300 (0.4%)
1	E	0.53	0/1691	1.02	9/2300 (0.4%)
1	F	0.54	0/1691	0.99	12/2300 (0.5%)
1	G	0.54	0/1691	1.01	10/2300 (0.4%)
1	H	0.54	0/1691	0.99	9/2300 (0.4%)
1	I	0.52	0/1691	0.99	9/2300 (0.4%)
1	J	0.51	0/1691	1.03	11/2300 (0.5%)
All	All	0.53	0/16910	1.00	90/23000 (0.4%)

There are no bond length outliers.

All (90) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	5	LYS	N-CA-C	-11.37	99.06	113.16
1	I	5	LYS	N-CA-C	-11.28	99.17	113.16
1	E	5	LYS	N-CA-C	-11.22	99.25	113.16
1	H	5	LYS	N-CA-C	-11.10	99.40	113.16
1	D	5	LYS	N-CA-C	-11.07	99.43	113.16
1	B	5	LYS	N-CA-C	-11.03	99.48	113.16
1	A	5	LYS	N-CA-C	-10.91	99.63	113.16
1	J	5	LYS	N-CA-C	-10.87	99.69	113.16
1	F	5	LYS	N-CA-C	-10.83	99.74	113.16
1	C	5	LYS	N-CA-C	-10.72	99.87	113.16
1	E	188	LYS	N-CA-C	-9.60	101.15	112.86
1	J	187	CYS	N-CA-C	8.02	122.14	110.28
1	E	189	GLU	CA-C-N	7.48	127.64	120.31
1	E	189	GLU	C-N-CA	7.48	127.64	120.31

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	191	TYR	CA-C-N	7.38	127.96	120.14
1	G	191	TYR	C-N-CA	7.38	127.96	120.14
1	G	188	LYS	N-CA-C	-7.22	104.05	112.86
1	J	15	ASN	CA-C-N	6.98	126.66	119.82
1	J	15	ASN	C-N-CA	6.98	126.66	119.82
1	I	189	GLU	CA-C-N	6.93	128.51	119.84
1	I	189	GLU	C-N-CA	6.93	128.51	119.84
1	J	165	ILE	CA-C-N	6.93	126.44	119.24
1	J	165	ILE	C-N-CA	6.93	126.44	119.24
1	H	165	ILE	CA-C-N	6.80	126.31	119.24
1	H	165	ILE	C-N-CA	6.80	126.31	119.24
1	G	165	ILE	CA-C-N	6.69	126.19	119.24
1	G	165	ILE	C-N-CA	6.69	126.19	119.24
1	I	165	ILE	CA-C-N	6.67	126.17	119.24
1	I	165	ILE	C-N-CA	6.67	126.17	119.24
1	H	15	ASN	CA-C-N	6.66	126.35	119.82
1	H	15	ASN	C-N-CA	6.66	126.35	119.82
1	J	186	CYS	CA-C-N	6.64	130.31	120.87
1	J	186	CYS	C-N-CA	6.64	130.31	120.87
1	D	109	SER	N-CA-C	-6.58	104.74	112.89
1	B	165	ILE	CA-C-N	6.51	126.01	119.24
1	B	165	ILE	C-N-CA	6.51	126.01	119.24
1	J	109	SER	N-CA-C	-6.46	104.88	112.89
1	D	15	ASN	CA-C-N	6.40	126.09	119.82
1	D	15	ASN	C-N-CA	6.40	126.09	119.82
1	F	165	ILE	CA-C-N	6.34	125.83	119.24
1	F	165	ILE	C-N-CA	6.34	125.83	119.24
1	C	109	SER	N-CA-C	-6.31	105.06	112.89
1	I	109	SER	N-CA-C	-6.29	105.09	112.89
1	I	15	ASN	CA-C-N	6.28	125.98	119.82
1	I	15	ASN	C-N-CA	6.28	125.98	119.82
1	C	165	ILE	CA-C-N	6.27	125.76	119.24
1	C	165	ILE	C-N-CA	6.27	125.76	119.24
1	G	109	SER	N-CA-C	-6.27	105.12	112.89
1	B	109	SER	N-CA-C	-6.25	105.14	112.89
1	A	165	ILE	CA-C-N	6.22	125.71	119.24
1	A	165	ILE	C-N-CA	6.22	125.71	119.24
1	C	25	ARG	CA-C-N	6.22	126.58	119.93
1	C	25	ARG	C-N-CA	6.22	126.58	119.93
1	F	191	TYR	CA-C-N	6.18	127.11	120.13
1	F	191	TYR	C-N-CA	6.18	127.11	120.13
1	F	109	SER	N-CA-C	-6.13	105.29	112.89

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	109	SER	N-CA-C	-6.11	105.31	112.89
1	A	109	SER	N-CA-C	-6.10	105.33	112.89
1	E	165	ILE	CA-C-N	6.09	125.58	119.24
1	E	165	ILE	C-N-CA	6.09	125.58	119.24
1	F	25	ARG	CA-C-N	6.09	126.30	119.90
1	F	25	ARG	C-N-CA	6.09	126.30	119.90
1	H	109	SER	N-CA-C	-6.07	105.36	112.89
1	J	184	TYR	N-CA-C	6.05	119.41	110.46
1	A	15	ASN	CA-C-N	6.04	125.74	119.82
1	A	15	ASN	C-N-CA	6.04	125.74	119.82
1	H	52	PHE	N-CA-C	6.03	118.58	109.23
1	G	15	ASN	CA-C-N	5.99	125.69	119.82
1	G	15	ASN	C-N-CA	5.99	125.69	119.82
1	H	187	CYS	N-CA-C	5.95	119.80	111.30
1	J	4	ARG	N-CA-C	5.88	116.83	108.54
1	D	165	ILE	CA-C-N	5.86	125.33	119.24
1	D	165	ILE	C-N-CA	5.86	125.33	119.24
1	C	184	TYR	N-CA-C	5.59	118.19	109.52
1	H	182	ARG	N-CA-C	5.54	118.59	109.72
1	D	24	ASP	N-CA-C	-5.38	107.09	113.38
1	I	4	ARG	N-CA-C	5.32	116.05	108.54
1	C	144	SER	N-CA-C	-5.32	103.12	110.35
1	D	182	ARG	N-CA-C	5.32	117.89	107.57
1	D	119	ILE	N-CA-C	5.31	117.03	108.85
1	B	183	PHE	N-CA-C	5.31	119.01	112.54
1	G	90	ALA	N-CA-C	-5.30	101.34	109.76
1	F	15	ASN	CA-C-N	5.14	126.26	119.84
1	F	15	ASN	C-N-CA	5.14	126.26	119.84
1	E	181	GLU	CA-C-N	-5.11	114.30	122.53
1	E	181	GLU	C-N-CA	-5.11	114.30	122.53
1	C	187	CYS	N-CA-CB	-5.10	103.51	111.56
1	B	91	TYR	N-CA-C	5.08	116.90	111.36
1	F	120	ARG	N-CA-C	-5.05	100.17	108.41
1	F	182	ARG	N-CA-C	5.03	117.16	109.07

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	1647	0	1610	66	0
1	B	1647	0	1610	70	0
1	C	1647	0	1609	77	0
1	D	1647	0	1611	76	0
1	E	1647	0	1610	74	0
1	F	1647	0	1610	70	0
1	G	1647	0	1611	74	0
1	H	1647	0	1609	69	0
1	I	1647	0	1610	72	0
1	J	1647	0	1609	74	0
2	K	28	0	25	0	0
2	L	28	0	25	3	0
2	M	28	0	25	1	0
3	A	28	0	26	2	0
3	B	14	0	13	2	0
3	C	28	0	26	4	0
3	D	14	0	13	0	0
3	E	28	0	26	0	0
3	F	14	0	13	2	0
3	G	14	0	13	0	0
3	H	14	0	13	0	0
3	I	14	0	13	0	0
3	J	28	0	26	0	0
All	All	16750	0	16356	687	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 21.

All (687) 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:J:92:ASN:H	1:J:92:ASN:HD22	1.07	0.99
1:D:92:ASN:HD22	1:D:92:ASN:H	1.03	0.97
1:F:92:ASN:H	1:F:92:ASN:HD22	0.96	0.93
1:C:92:ASN:H	1:C:92:ASN:HD22	1.17	0.89
1:A:185:GLU:O	1:A:186:CYS:HB3	1.71	0.88
1:J:21:THR:HG22	1:J:27:VAL:HG23	1.56	0.88
1:J:31:PHE:CE1	1:J:54:LEU:HD21	2.08	0.88
1:F:92:ASN:HD22	1:F:92:ASN:N	1.71	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:21:THR:HG23	1:D:27:VAL:HG23	1.57	0.86
1:A:186:CYS:SG	1:A:187:CYS:N	2.49	0.85
1:E:185:GLU:O	1:E:186:CYS:HB3	1.76	0.85
1:D:92:ASN:H	1:D:92:ASN:ND2	1.75	0.83
1:D:101:THR:HG23	1:D:118:SER:HB2	1.60	0.83
1:D:55:GLN:HG3	1:D:116:LEU:HD11	1.59	0.82
1:J:31:PHE:HE1	1:J:54:LEU:HD21	1.43	0.82
1:F:92:ASN:H	1:F:92:ASN:ND2	1.75	0.81
1:G:31:PHE:HE1	1:G:54:LEU:HD11	1.43	0.81
1:J:101:THR:HG23	1:J:118:SER:HB2	1.62	0.81
1:G:101:THR:HG23	1:G:118:SER:HB2	1.61	0.81
1:H:159:ALA:HB2	1:H:177:GLN:OE1	1.82	0.80
1:F:55:GLN:HG3	1:F:116:LEU:HD11	1.62	0.80
1:F:101:THR:HG23	1:F:118:SER:HB2	1.64	0.80
1:B:101:THR:HG23	1:B:118:SER:HB2	1.63	0.80
1:A:101:THR:HG23	1:A:118:SER:HB2	1.62	0.80
1:E:108:ASN:HD22	1:E:112:HIS:HB3	1.46	0.80
1:E:101:THR:HG23	1:E:118:SER:HB2	1.62	0.79
1:E:149:SER:HB2	1:E:190:PRO:HG2	1.61	0.79
1:C:26:PRO:HB3	1:C:150:ARG:O	1.84	0.78
1:D:26:PRO:HB3	1:D:150:ARG:O	1.83	0.78
1:I:159:ALA:HB2	1:I:177:GLN:OE1	1.84	0.78
1:I:101:THR:HG23	1:I:118:SER:HB2	1.65	0.78
1:G:55:GLN:HG3	1:G:116:LEU:HD11	1.66	0.78
1:H:101:THR:HG23	1:H:118:SER:HB2	1.65	0.77
1:B:159:ALA:HB2	1:B:177:GLN:OE1	1.85	0.77
1:F:108:ASN:HD22	1:F:112:HIS:HB3	1.49	0.77
1:G:149:SER:HB2	1:G:190:PRO:HG2	1.66	0.77
1:E:55:GLN:HG3	1:E:116:LEU:HD11	1.65	0.76
1:F:66:ASN:HB2	3:F:801:NAG:H2	1.67	0.76
1:A:55:GLN:HG3	1:A:116:LEU:HD11	1.68	0.76
1:I:181:GLU:HG2	1:I:190:PRO:HB2	1.67	0.76
1:D:92:ASN:HD22	1:D:92:ASN:N	1.84	0.76
1:H:148:HIS:CG	1:H:189:GLU:HG2	2.21	0.75
1:C:55:GLN:HG3	1:C:116:LEU:HD11	1.67	0.75
1:G:31:PHE:CE1	1:G:54:LEU:HD11	2.21	0.75
1:H:21:THR:HG22	1:H:27:VAL:HG23	1.68	0.75
1:A:159:ALA:HB2	1:A:177:GLN:OE1	1.84	0.75
1:F:159:ALA:HB2	1:F:177:GLN:OE1	1.87	0.75
1:J:92:ASN:HD22	1:J:92:ASN:N	1.84	0.74
1:H:55:GLN:HG3	1:H:116:LEU:HD11	1.69	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:55:GLN:HG3	1:J:116:LEU:HD11	1.69	0.74
1:C:101:THR:HG23	1:C:118:SER:HB2	1.68	0.74
1:D:108:ASN:HD22	1:D:112:HIS:HB3	1.51	0.74
1:J:26:PRO:HB3	1:J:150:ARG:O	1.87	0.73
1:D:55:GLN:HA	1:D:116:LEU:HD13	1.71	0.73
1:G:185:GLU:O	1:G:186:CYS:HB3	1.88	0.73
1:A:26:PRO:HB3	1:A:150:ARG:O	1.89	0.73
1:C:159:ALA:HB2	1:C:177:GLN:OE1	1.88	0.73
1:B:26:PRO:HB3	1:B:150:ARG:O	1.87	0.72
1:I:21:THR:HG23	1:I:27:VAL:HG23	1.71	0.72
1:B:108:ASN:HD22	1:B:112:HIS:HB3	1.55	0.72
1:D:159:ALA:HB2	1:D:177:GLN:OE1	1.89	0.72
1:F:26:PRO:HB3	1:F:150:ARG:O	1.90	0.72
1:B:55:GLN:HG3	1:B:116:LEU:HD11	1.72	0.72
1:I:26:PRO:HB3	1:I:150:ARG:O	1.91	0.71
1:H:26:PRO:HB3	1:H:150:ARG:O	1.88	0.71
1:G:159:ALA:HB2	1:G:177:GLN:OE1	1.91	0.70
1:E:26:PRO:HB3	1:E:150:ARG:O	1.91	0.70
1:J:159:ALA:HB2	1:J:177:GLN:OE1	1.91	0.70
1:F:186:CYS:C	1:F:187:CYS:SG	2.74	0.70
1:I:55:GLN:HG3	1:I:116:LEU:HD11	1.73	0.69
1:F:39:MET:HB2	1:F:49:ASP:HB3	1.74	0.69
1:E:176:THR:O	1:E:177:GLN:HB3	1.92	0.69
1:C:176:THR:O	1:C:177:GLN:HB3	1.92	0.69
1:E:159:ALA:HB2	1:E:177:GLN:OE1	1.92	0.69
2:L:1:NAG:H4	2:L:2:NAG:N2	2.07	0.69
1:F:55:GLN:HA	1:F:116:LEU:HD13	1.74	0.69
1:I:182:ARG:HG3	1:I:182:ARG:HH11	1.58	0.69
1:E:54:LEU:HD12	1:E:119:ILE:HD12	1.75	0.68
1:I:185:GLU:O	1:I:186:CYS:HB3	1.94	0.68
1:A:55:GLN:HA	1:A:116:LEU:HD13	1.74	0.68
1:G:182:ARG:O	1:G:190:PRO:HA	1.93	0.68
1:I:55:GLN:HA	1:I:116:LEU:HD13	1.76	0.67
1:B:183:PHE:O	1:B:184:TYR:O	2.13	0.67
1:E:39:MET:HB2	1:E:49:ASP:HB3	1.76	0.67
1:J:39:MET:HB2	1:J:49:ASP:HB3	1.76	0.67
1:E:55:GLN:HA	1:E:116:LEU:HD13	1.76	0.67
1:J:176:THR:O	1:J:177:GLN:HB3	1.92	0.67
1:G:26:PRO:HB3	1:G:150:ARG:O	1.93	0.67
1:J:186:CYS:SG	1:J:187:CYS:N	2.64	0.67
1:B:39:MET:HB2	1:B:49:ASP:HB3	1.75	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:165:ILE:HG22	1:F:168:SER:HB2	1.77	0.67
1:H:21:THR:CG2	1:H:27:VAL:HG23	2.25	0.67
1:G:39:MET:HB2	1:G:49:ASP:HB3	1.76	0.66
1:I:176:THR:O	1:I:177:GLN:HB3	1.93	0.66
1:A:39:MET:HB2	1:A:49:ASP:HB3	1.77	0.66
1:J:21:THR:CG2	1:J:27:VAL:HG23	2.23	0.66
1:A:176:THR:O	1:A:177:GLN:HB3	1.96	0.66
1:I:39:MET:HB2	1:I:49:ASP:HB3	1.76	0.66
1:D:39:MET:HB2	1:D:49:ASP:HB3	1.77	0.65
1:H:176:THR:O	1:H:177:GLN:HB3	1.95	0.65
1:D:23:ARG:O	1:D:25:ARG:HG2	1.96	0.65
1:I:156:MET:HE3	1:I:177:GLN:HE21	1.62	0.65
1:D:176:THR:O	1:D:177:GLN:HB3	1.96	0.65
1:I:150:ARG:NH1	1:I:190:PRO:HD2	2.11	0.65
1:C:39:MET:HB2	1:C:49:ASP:HB3	1.79	0.65
1:D:165:ILE:HG22	1:D:168:SER:HB2	1.79	0.65
1:A:108:ASN:HD22	1:A:112:HIS:HB3	1.62	0.65
1:A:116:LEU:O	1:E:145:TRP:HZ2	1.80	0.65
1:A:165:ILE:HG22	1:A:168:SER:HB2	1.79	0.64
1:G:55:GLN:HA	1:G:116:LEU:HD13	1.79	0.64
1:I:181:GLU:OE2	1:I:190:PRO:HG3	1.97	0.64
1:A:46:GLN:OE1	1:A:126:ASP:HA	1.97	0.64
1:B:176:THR:O	1:B:177:GLN:HB3	1.97	0.64
1:F:42:ASP:HB3	1:F:47:VAL:HG22	1.79	0.64
1:F:176:THR:O	1:F:177:GLN:HB3	1.95	0.64
1:G:176:THR:O	1:G:177:GLN:HB3	1.96	0.64
1:A:21:THR:HG23	1:A:27:VAL:HG23	1.80	0.64
1:I:3:GLN:N	1:I:3:GLN:OE1	2.31	0.64
1:I:169:ARG:HD2	1:I:202:LYS:HE3	1.80	0.64
1:B:55:GLN:HA	1:B:116:LEU:HD13	1.79	0.63
1:B:141:LYS:NZ	1:B:182:ARG:HH22	1.96	0.63
1:G:156:MET:HE3	1:G:177:GLN:HE21	1.63	0.63
1:H:165:ILE:HG22	1:H:168:SER:HB2	1.80	0.63
1:J:42:ASP:HB3	1:J:47:VAL:HG22	1.80	0.63
1:C:55:GLN:HA	1:C:116:LEU:HD13	1.79	0.63
1:D:21:THR:CG2	1:D:27:VAL:HG23	2.27	0.63
1:D:29:VAL:HG12	1:D:58:TRP:HB3	1.81	0.63
1:C:3:GLN:HE21	1:C:3:GLN:HA	1.62	0.63
1:A:156:MET:HE3	1:A:177:GLN:HE21	1.62	0.63
1:J:92:ASN:H	1:J:92:ASN:ND2	1.88	0.63
1:D:42:ASP:HB3	1:D:47:VAL:HG22	1.80	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:108:ASN:HB2	1:B:112:HIS:H	1.64	0.62
1:E:108:ASN:HB2	1:E:112:HIS:H	1.64	0.62
1:H:55:GLN:HA	1:H:116:LEU:HD13	1.80	0.62
1:C:46:GLN:OE1	1:C:126:ASP:HA	1.99	0.62
1:E:156:MET:HE3	1:E:177:GLN:HE21	1.65	0.62
1:B:42:ASP:HB3	1:B:47:VAL:HG22	1.79	0.62
1:I:46:GLN:OE1	1:I:126:ASP:HA	1.98	0.62
1:C:169:ARG:HD2	1:C:202:LYS:HE3	1.81	0.62
1:E:165:ILE:HG22	1:E:168:SER:HB2	1.81	0.62
1:H:46:GLN:OE1	1:H:126:ASP:HA	1.98	0.62
1:J:6:LEU:HD23	1:J:73:VAL:HG11	1.81	0.62
1:D:96:LYS:HG3	1:D:97:PRO:HD2	1.82	0.61
1:G:42:ASP:HB3	1:G:47:VAL:HG22	1.81	0.61
1:B:141:LYS:HZ2	1:B:182:ARG:HH22	1.48	0.61
1:F:23:ARG:HH11	1:G:71:PRO:HB3	1.65	0.61
1:H:156:MET:HE3	1:H:177:GLN:HE21	1.65	0.61
1:I:181:GLU:OE2	1:I:190:PRO:CG	2.48	0.61
1:G:149:SER:OG	1:G:192:PRO:HD3	2.01	0.61
1:A:42:ASP:HB3	1:A:47:VAL:HG22	1.81	0.61
1:F:29:VAL:HG12	1:F:58:TRP:HB3	1.83	0.61
1:F:39:MET:HE1	1:J:47:VAL:HG11	1.83	0.61
1:J:108:ASN:HB2	1:J:112:HIS:H	1.66	0.61
1:A:107:VAL:HG12	1:A:113:VAL:HG22	1.83	0.61
1:B:156:MET:HE3	1:B:177:GLN:HE21	1.65	0.61
1:B:6:LEU:HD23	1:B:73:VAL:HG11	1.83	0.61
1:B:22:GLN:HG3	1:B:25:ARG:HB2	1.83	0.61
1:C:108:ASN:HB2	1:C:112:HIS:H	1.65	0.61
1:J:55:GLN:HA	1:J:116:LEU:HD13	1.80	0.61
1:H:42:ASP:HB3	1:H:47:VAL:HG22	1.83	0.61
1:J:165:ILE:HG22	1:J:168:SER:HB2	1.82	0.60
1:D:156:MET:HE3	1:D:177:GLN:HE21	1.66	0.60
1:D:108:ASN:HB2	1:D:112:HIS:H	1.65	0.60
1:E:107:VAL:HG12	1:E:113:VAL:HG22	1.82	0.60
1:A:108:ASN:HB2	1:A:112:HIS:H	1.66	0.60
1:B:46:GLN:OE1	1:B:126:ASP:HA	2.01	0.60
1:H:39:MET:HB2	1:H:49:ASP:HB3	1.84	0.60
1:B:185:GLU:O	1:B:186:CYS:HB3	2.01	0.60
1:G:165:ILE:HG22	1:G:168:SER:HB2	1.81	0.60
1:H:6:LEU:HD23	1:H:73:VAL:HG11	1.83	0.60
1:C:156:MET:HE3	1:C:177:GLN:HE21	1.67	0.60
1:D:6:LEU:HD23	1:D:73:VAL:HG11	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:6:LEU:HD23	1:F:73:VAL:HG11	1.83	0.60
1:G:107:VAL:HG12	1:G:113:VAL:HG22	1.84	0.59
1:H:108:ASN:HB2	1:H:112:HIS:H	1.65	0.59
1:A:6:LEU:HD23	1:A:73:VAL:HG11	1.84	0.59
1:G:108:ASN:HB2	1:G:112:HIS:H	1.67	0.59
1:C:108:ASN:CG	3:C:901:NAG:HN2	2.10	0.59
1:C:165:ILE:HG22	1:C:168:SER:HB2	1.84	0.59
1:D:107:VAL:HG12	1:D:113:VAL:HG22	1.84	0.59
1:G:6:LEU:HD23	1:G:73:VAL:HG11	1.84	0.59
1:I:42:ASP:HB3	1:I:47:VAL:HG22	1.84	0.59
1:I:165:ILE:HG22	1:I:168:SER:HB2	1.83	0.59
1:A:148:HIS:HB2	1:A:189:GLU:HB3	1.83	0.59
1:F:169:ARG:HD2	1:F:202:LYS:HE3	1.85	0.59
3:B:901:NAG:H3	3:B:901:NAG:C8	2.33	0.59
1:D:46:GLN:OE1	1:D:126:ASP:HA	2.02	0.59
1:C:6:LEU:HD23	1:C:73:VAL:HG11	1.83	0.59
1:F:108:ASN:HB2	1:F:112:HIS:H	1.65	0.59
1:F:150:ARG:HD3	1:F:189:GLU:CG	2.33	0.59
1:J:96:LYS:HG3	1:J:97:PRO:HD2	1.85	0.59
1:F:46:GLN:OE1	1:F:126:ASP:HA	2.02	0.59
1:G:156:MET:HE3	1:G:177:GLN:NE2	2.18	0.59
1:F:96:LYS:HG3	1:F:97:PRO:HD2	1.85	0.58
1:H:22:GLN:O	1:H:23:ARG:O	2.21	0.58
1:C:21:THR:HG22	1:C:27:VAL:HG23	1.84	0.58
1:D:185:GLU:O	1:D:186:CYS:SG	2.62	0.58
1:E:187:CYS:HB3	1:E:189:GLU:HB2	1.85	0.58
1:J:46:GLN:OE1	1:J:126:ASP:HA	2.03	0.58
1:J:156:MET:HE3	1:J:177:GLN:HE21	1.67	0.58
1:G:92:ASN:HD21	1:G:141:LYS:H	1.50	0.58
1:I:108:ASN:HB2	1:I:112:HIS:H	1.67	0.58
1:B:165:ILE:HG22	1:B:168:SER:HB2	1.85	0.58
1:E:46:GLN:OE1	1:E:126:ASP:HA	2.03	0.58
1:F:107:VAL:HG12	1:F:113:VAL:HG22	1.85	0.58
1:F:156:MET:HE3	1:F:177:GLN:HE21	1.67	0.58
1:B:169:ARG:HD2	1:B:202:LYS:HE3	1.85	0.58
1:G:3:GLN:O	1:G:4:ARG:HD3	2.03	0.58
1:E:42:ASP:HB3	1:E:47:VAL:HG22	1.86	0.58
1:I:6:LEU:HD23	1:I:73:VAL:HG11	1.85	0.58
1:E:156:MET:HE3	1:E:177:GLN:NE2	2.19	0.57
1:B:29:VAL:HG12	1:B:58:TRP:HB3	1.86	0.57
1:C:3:GLN:HA	1:C:3:GLN:NE2	2.19	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:156:MET:HE3	1:A:177:GLN:NE2	2.19	0.57
1:I:156:MET:HE3	1:I:177:GLN:NE2	2.19	0.57
1:C:23:ARG:HH11	1:D:71:PRO:HB3	1.69	0.57
1:I:29:VAL:HG12	1:I:58:TRP:HB3	1.86	0.57
1:G:29:VAL:HG12	1:G:58:TRP:HB3	1.87	0.57
1:I:96:LYS:HG3	1:I:97:PRO:HD2	1.86	0.57
1:C:42:ASP:HB3	1:C:47:VAL:HG22	1.87	0.57
1:G:92:ASN:N	1:G:92:ASN:HD22	2.02	0.56
1:E:148:HIS:CD2	1:E:189:GLU:HG2	2.39	0.56
1:E:6:LEU:HD23	1:E:73:VAL:HG11	1.87	0.56
1:G:46:GLN:OE1	1:G:126:ASP:HA	2.04	0.56
1:G:108:ASN:HD22	1:G:112:HIS:HB3	1.70	0.56
1:J:22:GLN:O	1:J:23:ARG:O	2.23	0.56
1:E:187:CYS:C	1:E:189:GLU:N	2.58	0.56
1:H:23:ARG:O	1:H:25:ARG:HG2	2.06	0.56
1:B:156:MET:HE3	1:B:177:GLN:NE2	2.21	0.56
1:J:149:SER:HB2	1:J:190:PRO:HG2	1.88	0.56
1:A:29:VAL:HG12	1:A:58:TRP:HB3	1.87	0.56
1:C:107:VAL:HG12	1:C:113:VAL:HG22	1.87	0.56
1:D:55:GLN:HG3	1:D:116:LEU:CD1	2.34	0.56
1:E:29:VAL:HG12	1:E:58:TRP:HB3	1.88	0.56
1:I:150:ARG:NH1	1:I:189:GLU:HA	2.21	0.55
1:B:96:LYS:HG3	1:B:97:PRO:HD2	1.87	0.55
1:E:149:SER:CB	1:E:190:PRO:HG2	2.34	0.55
1:G:185:GLU:O	1:G:186:CYS:CB	2.55	0.55
1:H:29:VAL:HG12	1:H:58:TRP:HB3	1.88	0.55
1:B:107:VAL:HG12	1:B:113:VAL:HG22	1.87	0.55
1:F:107:VAL:CG1	1:F:113:VAL:HG22	2.37	0.55
1:G:169:ARG:HD2	1:G:202:LYS:HE3	1.89	0.55
1:G:189:GLU:HA	1:G:189:GLU:OE2	2.05	0.55
1:H:49:ASP:HA	1:H:122:ARG:HG3	1.89	0.55
1:A:101:THR:HG21	1:E:145:TRP:CH2	2.42	0.55
1:D:169:ARG:HD2	1:D:202:LYS:HE3	1.88	0.55
1:H:108:ASN:ND2	1:H:112:HIS:HB3	2.21	0.55
1:I:107:VAL:HG12	1:I:113:VAL:HG22	1.89	0.55
1:J:167:TYR:O	1:J:168:SER:HB3	2.07	0.55
1:F:3:GLN:HE22	1:J:23:ARG:NE	2.05	0.54
1:G:96:LYS:HG3	1:G:97:PRO:HD2	1.89	0.54
1:J:169:ARG:HD2	1:J:202:LYS:HE3	1.89	0.54
1:C:96:LYS:HG3	1:C:97:PRO:HD2	1.88	0.54
1:D:107:VAL:CG1	1:D:113:VAL:HG22	2.37	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:96:LYS:HG3	1:E:97:PRO:HD2	1.88	0.54
1:A:52:PHE:CD1	1:A:52:PHE:N	2.75	0.54
1:E:92:ASN:HD21	1:E:140:LEU:HA	1.72	0.54
1:H:189:GLU:HG3	1:H:190:PRO:HD2	1.89	0.54
1:J:29:VAL:HG12	1:J:58:TRP:HB3	1.87	0.54
1:B:148:HIS:CE1	1:B:189:GLU:HG2	2.43	0.54
1:C:110:SER:HB3	3:C:901:NAG:C8	2.38	0.54
1:H:107:VAL:HG12	1:H:113:VAL:HG22	1.89	0.54
1:D:76:VAL:HG23	1:D:107:VAL:HG23	1.88	0.54
1:I:108:ASN:HD22	1:I:112:HIS:HB3	1.72	0.54
1:J:107:VAL:HG12	1:J:113:VAL:HG22	1.89	0.54
1:A:156:MET:CE	1:A:177:GLN:HE21	2.21	0.54
1:A:167:TYR:O	1:A:168:SER:HB3	2.08	0.54
1:G:187:CYS:O	1:G:189:GLU:HG2	2.07	0.54
1:I:49:ASP:HA	1:I:122:ARG:HG3	1.90	0.54
1:A:39:MET:HE1	1:E:47:VAL:HG11	1.89	0.53
1:A:96:LYS:HG3	1:A:97:PRO:HD2	1.88	0.53
1:H:156:MET:HE3	1:H:177:GLN:NE2	2.22	0.53
1:A:107:VAL:CG1	1:A:113:VAL:HG22	2.38	0.53
1:G:34:SER:HB3	1:G:53:TRP:HB3	1.90	0.53
1:C:49:ASP:HA	1:C:122:ARG:HG3	1.91	0.53
1:D:156:MET:HE3	1:D:177:GLN:NE2	2.23	0.53
1:F:156:MET:HE3	1:F:177:GLN:NE2	2.24	0.53
1:B:76:VAL:HG23	1:B:107:VAL:HG23	1.91	0.53
1:C:189:GLU:OE2	1:C:190:PRO:HD2	2.08	0.53
1:E:107:VAL:CG1	1:E:113:VAL:HG22	2.39	0.53
1:J:156:MET:HE3	1:J:177:GLN:NE2	2.22	0.53
1:G:107:VAL:CG1	1:G:113:VAL:HG22	2.37	0.53
1:H:186:CYS:SG	1:H:187:CYS:N	2.81	0.53
1:D:167:TYR:O	1:D:168:SER:HB3	2.08	0.52
1:F:21:THR:HG22	1:F:27:VAL:HG23	1.90	0.52
1:H:22:GLN:HG3	1:H:25:ARG:HG3	1.91	0.52
1:G:47:VAL:HG11	1:H:39:MET:HE1	1.90	0.52
1:H:110:SER:HB3	2:L:1:NAG:H82	1.91	0.52
1:A:49:ASP:HA	1:A:122:ARG:HG3	1.92	0.52
1:B:107:VAL:CG1	1:B:113:VAL:HG22	2.39	0.52
1:C:156:MET:HE3	1:C:177:GLN:NE2	2.24	0.52
1:G:150:ARG:NH1	1:G:190:PRO:HG3	2.24	0.52
1:H:33:LEU:C	1:H:33:LEU:HD23	2.34	0.52
1:I:110:SER:OG	2:M:1:NAG:H61	2.08	0.52
1:F:76:VAL:HG23	1:F:107:VAL:HG23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:55:GLN:HG3	1:G:116:LEU:CD1	2.38	0.52
1:G:149:SER:HG	1:G:192:PRO:HD3	1.74	0.52
1:H:96:LYS:HG3	1:H:97:PRO:HD2	1.91	0.52
1:B:141:LYS:NZ	1:B:182:ARG:NH2	2.57	0.52
1:C:21:THR:CG2	1:C:27:VAL:HG23	2.40	0.52
1:H:33:LEU:HD23	1:H:34:SER:N	2.25	0.52
1:E:167:TYR:O	1:E:168:SER:HB3	2.10	0.52
1:G:21:THR:HG22	1:G:27:VAL:HG23	1.90	0.52
1:I:156:MET:CE	1:I:177:GLN:HE21	2.22	0.52
1:D:3:GLN:HA	1:D:3:GLN:NE2	2.24	0.52
1:D:188:LYS:O	1:D:188:LYS:HD3	2.10	0.52
1:C:167:TYR:O	1:C:168:SER:HB3	2.07	0.52
1:G:156:MET:CE	1:G:177:GLN:HE21	2.23	0.52
1:I:185:GLU:O	1:I:186:CYS:CB	2.58	0.52
1:I:167:TYR:O	1:I:168:SER:HB3	2.10	0.51
1:A:45:ASN:ND2	1:B:39:MET:O	2.43	0.51
1:C:29:VAL:HG12	1:C:58:TRP:HB3	1.93	0.51
1:G:186:CYS:O	1:G:188:LYS:N	2.39	0.51
1:H:165:ILE:CG2	1:H:168:SER:HB2	2.41	0.51
1:B:49:ASP:HA	1:B:122:ARG:HG3	1.91	0.51
1:E:49:ASP:HA	1:E:122:ARG:HG3	1.92	0.51
1:F:167:TYR:O	1:F:168:SER:HB3	2.10	0.51
1:D:165:ILE:CG2	1:D:168:SER:HB2	2.40	0.51
1:E:186:CYS:O	1:E:188:LYS:N	2.40	0.51
1:F:66:ASN:HB2	3:F:801:NAG:C2	2.40	0.51
1:I:107:VAL:CG1	1:I:113:VAL:HG22	2.41	0.51
1:J:34:SER:HB3	1:J:53:TRP:HB3	1.92	0.51
1:C:47:VAL:HG11	1:D:39:MET:HE1	1.92	0.51
1:D:3:GLN:HA	1:D:3:GLN:HE21	1.76	0.51
1:A:76:VAL:HG23	1:A:107:VAL:HG23	1.92	0.51
1:B:101:THR:CG2	1:B:118:SER:HB2	2.39	0.51
1:I:186:CYS:SG	1:I:187:CYS:N	2.83	0.51
1:E:101:THR:CG2	1:E:118:SER:HB2	2.39	0.51
1:E:187:CYS:C	1:E:189:GLU:H	2.19	0.51
1:I:145:TRP:CH2	1:J:101:THR:HG21	2.45	0.51
1:A:20:PRO:O	1:A:27:VAL:HG22	2.11	0.51
1:B:3:GLN:HA	1:B:3:GLN:NE2	2.25	0.51
1:H:18:VAL:HG22	1:I:7:TYR:CD2	2.46	0.51
1:H:34:SER:HB3	1:H:53:TRP:HB3	1.91	0.51
1:H:49:ASP:OD2	1:H:122:ARG:HD2	2.11	0.50
1:J:107:VAL:CG1	1:J:113:VAL:HG22	2.41	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:33:LEU:C	1:E:33:LEU:HD23	2.37	0.50
1:F:165:ILE:CG2	1:F:168:SER:HB2	2.40	0.50
1:C:148:HIS:HA	1:C:191:TYR:CD2	2.46	0.50
1:J:49:ASP:HA	1:J:122:ARG:HG3	1.93	0.50
1:A:165:ILE:CG2	1:A:168:SER:HB2	2.42	0.50
1:D:33:LEU:HD23	1:D:33:LEU:C	2.37	0.50
1:F:186:CYS:SG	1:F:187:CYS:N	2.84	0.50
1:C:107:VAL:CG1	1:C:113:VAL:HG22	2.41	0.50
1:C:161:ILE:HG12	1:C:161:ILE:O	2.11	0.50
1:D:49:ASP:HA	1:D:122:ARG:HG3	1.93	0.50
1:A:33:LEU:C	1:A:33:LEU:HD23	2.36	0.50
1:B:22:GLN:HG3	1:B:25:ARG:CB	2.42	0.50
1:G:54:LEU:HD12	1:G:56:MET:HE2	1.94	0.50
1:A:169:ARG:NH1	1:E:44:LYS:HD3	2.27	0.49
1:E:54:LEU:HD12	1:E:119:ILE:CD1	2.42	0.49
1:F:55:GLN:HG3	1:F:116:LEU:CD1	2.36	0.49
1:F:169:ARG:HH12	1:J:44:LYS:HD3	1.77	0.49
1:D:23:ARG:NH1	1:E:71:PRO:HB3	2.27	0.49
1:E:31:PHE:CE1	1:E:54:LEU:HD22	2.47	0.49
1:F:23:ARG:NH1	1:G:71:PRO:HB3	2.26	0.49
1:J:165:ILE:CG2	1:J:168:SER:HB2	2.42	0.49
1:B:92:ASN:HD22	1:B:123:PHE:HD2	1.60	0.49
1:B:150:ARG:NH1	1:B:190:PRO:CD	2.75	0.49
1:F:21:THR:CG2	1:F:27:VAL:HG23	2.42	0.49
1:G:101:THR:CG2	1:G:118:SER:HB2	2.38	0.49
1:G:145:TRP:CH2	1:H:101:THR:HG21	2.47	0.49
1:B:167:TYR:O	1:B:168:SER:HB3	2.13	0.49
1:B:186:CYS:SG	1:B:187:CYS:N	2.84	0.49
1:C:44:LYS:HD3	1:D:169:ARG:HH12	1.77	0.49
1:F:52:PHE:CE2	1:F:54:LEU:HG	2.48	0.49
1:H:55:GLN:HG3	1:H:116:LEU:CD1	2.42	0.49
1:J:142:PHE:O	1:J:193:ASP:HB2	2.13	0.49
1:C:76:VAL:HG23	1:C:107:VAL:HG23	1.94	0.49
1:E:165:ILE:CG2	1:E:168:SER:HB2	2.42	0.49
1:H:148:HIS:CE1	1:H:151:GLU:HG3	2.47	0.49
1:A:31:PHE:HE2	1:A:196:PHE:HE1	1.59	0.49
1:B:150:ARG:HD3	1:B:189:GLU:HG3	1.95	0.49
1:D:92:ASN:HD21	1:D:141:LYS:H	1.58	0.49
1:H:107:VAL:CG1	1:H:113:VAL:HG22	2.43	0.49
1:E:187:CYS:O	1:E:188:LYS:HB2	2.13	0.49
1:G:165:ILE:CG2	1:G:168:SER:HB2	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:76:VAL:HG23	1:H:107:VAL:HG23	1.94	0.49
1:B:156:MET:CE	1:B:177:GLN:HE21	2.26	0.49
1:G:49:ASP:HA	1:G:122:ARG:HG3	1.94	0.49
1:D:12:LYS:HE3	1:I:69:GLU:OE1	2.13	0.49
1:I:41:VAL:HG12	1:I:48:VAL:HG12	1.94	0.49
1:D:156:MET:CE	1:D:177:GLN:HE21	2.26	0.48
1:E:150:ARG:NH1	1:E:190:PRO:CD	2.76	0.48
1:I:21:THR:CG2	1:I:27:VAL:HG23	2.41	0.48
1:C:141:LYS:NZ	1:C:184:TYR:OH	2.44	0.48
1:F:3:GLN:HE21	1:F:3:GLN:HA	1.78	0.48
1:F:49:ASP:HA	1:F:122:ARG:HG3	1.95	0.48
1:G:76:VAL:HG23	1:G:107:VAL:HG23	1.94	0.48
1:C:92:ASN:H	1:C:92:ASN:ND2	1.97	0.48
1:E:55:GLN:HG3	1:E:116:LEU:CD1	2.38	0.48
1:E:156:MET:CE	1:E:177:GLN:HE21	2.24	0.48
1:D:65:TRP:CE2	1:D:109:SER:HA	2.49	0.48
1:E:167:TYR:CD1	1:E:167:TYR:N	2.81	0.48
1:F:92:ASN:N	1:F:92:ASN:ND2	2.39	0.48
1:J:76:VAL:HG23	1:J:107:VAL:HG23	1.94	0.48
1:J:149:SER:CB	1:J:190:PRO:HG2	2.44	0.48
1:C:92:ASN:HD22	1:C:92:ASN:N	1.95	0.48
1:D:20:PRO:O	1:D:27:VAL:HG22	2.14	0.48
1:F:33:LEU:HD23	1:F:34:SER:N	2.28	0.48
1:F:161:ILE:HG12	1:F:161:ILE:O	2.13	0.48
1:D:76:VAL:CG2	1:D:107:VAL:HG23	2.44	0.48
1:F:101:THR:HG21	1:J:145:TRP:CZ2	2.48	0.48
1:I:148:HIS:HB3	1:I:191:TYR:CZ	2.47	0.48
1:A:49:ASP:OD2	1:A:122:ARG:HD2	2.13	0.48
1:E:169:ARG:HD2	1:E:202:LYS:HE3	1.95	0.48
1:I:165:ILE:CG2	1:I:168:SER:HB2	2.43	0.48
1:G:33:LEU:HD23	1:G:34:SER:N	2.29	0.48
1:J:36:LEU:CD2	1:J:53:TRP:HB2	2.43	0.48
1:J:41:VAL:HG12	1:J:48:VAL:HG12	1.96	0.48
1:A:33:LEU:HD23	1:A:34:SER:N	2.29	0.47
1:A:169:ARG:HH12	1:E:44:LYS:HD3	1.79	0.47
1:J:161:ILE:HG12	1:J:161:ILE:O	2.13	0.47
1:C:165:ILE:CG2	1:C:168:SER:HB2	2.43	0.47
1:H:35:LEU:HD21	1:H:50:VAL:HG13	1.95	0.47
1:I:76:VAL:HG23	1:I:107:VAL:HG23	1.96	0.47
1:I:189:GLU:CD	1:I:189:GLU:H	2.22	0.47
1:B:149:SER:OG	1:B:192:PRO:HD3	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:150:ARG:NH1	1:B:190:PRO:HD3	2.28	0.47
1:C:110:SER:HB3	3:C:901:NAG:H82	1.95	0.47
1:E:33:LEU:HD23	1:E:34:SER:N	2.28	0.47
1:A:101:THR:CG2	1:A:118:SER:HB2	2.39	0.47
1:G:148:HIS:CE1	1:G:151:GLU:HG3	2.49	0.47
1:J:76:VAL:CG2	1:J:107:VAL:HG23	2.45	0.47
1:A:112:HIS:HD2	3:A:901:NAG:C8	2.28	0.47
1:B:148:HIS:CE1	1:B:151:GLU:HG3	2.50	0.47
1:F:33:LEU:HD23	1:F:33:LEU:C	2.40	0.47
1:J:36:LEU:HD21	1:J:53:TRP:HB2	1.96	0.47
1:G:92:ASN:N	1:G:92:ASN:ND2	2.61	0.47
1:F:101:THR:HG21	1:J:145:TRP:CH2	2.49	0.46
1:H:101:THR:CG2	1:H:118:SER:HB2	2.41	0.46
1:D:101:THR:CG2	1:D:118:SER:HB2	2.38	0.46
1:F:156:MET:CE	1:F:177:GLN:HE21	2.26	0.46
1:F:169:ARG:NH1	1:J:44:LYS:HD3	2.30	0.46
1:G:44:LYS:HD3	1:H:169:ARG:NH1	2.30	0.46
1:H:23:ARG:NH1	1:I:71:PRO:HB3	2.30	0.46
1:B:49:ASP:OD2	1:B:122:ARG:HD2	2.15	0.46
1:E:148:HIS:CE1	1:E:151:GLU:HG3	2.50	0.46
1:F:148:HIS:CE1	1:F:151:GLU:HG3	2.51	0.46
1:G:33:LEU:HD23	1:G:33:LEU:C	2.39	0.46
1:I:33:LEU:HD23	1:I:33:LEU:C	2.41	0.46
1:J:33:LEU:HD23	1:J:34:SER:N	2.30	0.46
1:A:55:GLN:HG3	1:A:116:LEU:CD1	2.43	0.46
1:A:76:VAL:CG2	1:A:107:VAL:HG23	2.45	0.46
1:B:21:THR:HG23	1:B:27:VAL:HG23	1.98	0.46
1:B:33:LEU:HD23	1:B:33:LEU:C	2.41	0.46
1:C:76:VAL:CG2	1:C:107:VAL:HG23	2.45	0.46
1:D:33:LEU:HD23	1:D:34:SER:N	2.30	0.46
1:E:20:PRO:HG2	1:E:27:VAL:HG21	1.96	0.46
1:H:156:MET:CE	1:H:177:GLN:HE21	2.28	0.46
1:D:41:VAL:HG12	1:D:48:VAL:HG12	1.96	0.46
1:D:76:VAL:HG23	1:D:107:VAL:CG2	2.45	0.46
1:E:76:VAL:HG23	1:E:107:VAL:HG23	1.97	0.46
1:E:150:ARG:HH11	1:E:190:PRO:CD	2.29	0.46
1:G:18:VAL:HG22	1:H:7:TYR:CD2	2.51	0.46
1:G:44:LYS:HD3	1:H:169:ARG:HH12	1.79	0.46
1:B:189:GLU:HA	1:B:190:PRO:HD3	1.83	0.46
1:C:41:VAL:HG12	1:C:48:VAL:HG12	1.98	0.46
1:C:148:HIS:CE1	1:C:151:GLU:HG3	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:VAL:CG2	1:B:107:VAL:HG23	2.46	0.46
1:G:167:TYR:O	1:G:168:SER:HB3	2.15	0.46
1:J:187:CYS:HB3	1:J:189:GLU:OE1	2.16	0.46
1:D:23:ARG:C	1:D:25:ARG:N	2.73	0.46
1:G:187:CYS:C	1:G:189:GLU:N	2.72	0.46
1:I:47:VAL:HG11	1:J:39:MET:HE1	1.98	0.46
1:I:148:HIS:CE1	1:I:151:GLU:HG3	2.51	0.46
1:A:54:LEU:HD12	1:A:119:ILE:HD12	1.98	0.46
1:D:148:HIS:CE1	1:D:151:GLU:HG3	2.50	0.45
1:H:76:VAL:CG2	1:H:107:VAL:HG23	2.46	0.45
1:I:33:LEU:HD23	1:I:34:SER:N	2.31	0.45
1:A:35:LEU:HD21	1:A:50:VAL:HG13	1.97	0.45
1:C:23:ARG:NH1	1:D:71:PRO:HB3	2.30	0.45
1:D:185:GLU:OE1	1:D:185:GLU:HA	2.16	0.45
1:H:167:TYR:O	1:H:168:SER:HB3	2.17	0.45
1:H:169:ARG:HD2	1:H:202:LYS:HE3	1.99	0.45
1:B:167:TYR:CD1	1:B:167:TYR:N	2.84	0.45
1:G:33:LEU:HB3	1:G:196:PHE:CE2	2.51	0.45
1:H:54:LEU:HA	1:H:54:LEU:HD23	1.67	0.45
1:J:156:MET:CE	1:J:177:GLN:HE21	2.29	0.45
1:E:80:ILE:C	1:E:80:ILE:HD12	2.41	0.45
1:F:76:VAL:CG2	1:F:107:VAL:HG23	2.46	0.45
1:J:54:LEU:HD12	1:J:119:ILE:HD12	1.99	0.45
1:E:34:SER:HB3	1:E:53:TRP:HB3	1.97	0.45
1:I:167:TYR:CD1	1:I:167:TYR:N	2.83	0.45
1:J:41:VAL:O	1:J:41:VAL:HG23	2.17	0.45
1:A:41:VAL:HG12	1:A:48:VAL:HG12	1.99	0.45
1:B:52:PHE:CE2	1:B:54:LEU:HG	2.51	0.45
1:F:76:VAL:HG23	1:F:107:VAL:CG2	2.47	0.45
1:G:76:VAL:CG2	1:G:107:VAL:HG23	2.47	0.45
1:A:22:GLN:HG3	1:A:25:ARG:HB2	1.98	0.45
1:A:148:HIS:CE1	1:A:151:GLU:HG3	2.51	0.45
1:C:29:VAL:HG23	1:C:29:VAL:O	2.17	0.45
1:E:76:VAL:CG2	1:E:107:VAL:HG23	2.46	0.45
1:F:150:ARG:HD3	1:F:189:GLU:HG2	1.97	0.45
1:H:167:TYR:CD1	1:H:167:TYR:N	2.82	0.45
1:B:165:ILE:CG2	1:B:168:SER:HB2	2.46	0.45
1:D:161:ILE:O	1:D:161:ILE:HG12	2.17	0.45
1:F:41:VAL:HG12	1:F:48:VAL:HG12	1.99	0.45
1:G:149:SER:CB	1:G:190:PRO:HG2	2.43	0.45
1:I:101:THR:CG2	1:I:118:SER:HB2	2.43	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:J:183:PHE:CE1	1:J:190:PRO:HB3	2.52	0.45
1:A:21:THR:CG2	1:A:27:VAL:HG23	2.47	0.45
1:H:22:GLN:HG3	1:H:25:ARG:CB	2.47	0.45
1:A:176:THR:OG1	1:A:197:THR:HB	2.17	0.44
1:D:35:LEU:HD21	1:D:50:VAL:HG13	1.99	0.44
1:F:23:ARG:C	1:F:25:ARG:H	2.23	0.44
1:B:76:VAL:HG23	1:B:107:VAL:CG2	2.47	0.44
1:C:34:SER:HB3	1:C:53:TRP:HB3	1.99	0.44
1:J:33:LEU:HD23	1:J:33:LEU:C	2.42	0.44
1:E:41:VAL:O	1:E:41:VAL:HG23	2.17	0.44
1:G:41:VAL:HG23	1:G:41:VAL:O	2.16	0.44
1:J:148:HIS:CE1	1:J:151:GLU:HG3	2.52	0.44
1:A:23:ARG:C	1:A:25:ARG:N	2.75	0.44
1:B:33:LEU:HD23	1:B:34:SER:N	2.33	0.44
3:B:901:NAG:H3	3:B:901:NAG:H83	1.99	0.44
1:G:191:TYR:HA	1:G:192:PRO:HD2	1.84	0.44
1:I:22:GLN:O	1:I:23:ARG:O	2.35	0.44
1:I:80:ILE:HD12	1:I:80:ILE:C	2.42	0.44
1:I:182:ARG:HG3	1:I:182:ARG:NH1	2.30	0.44
1:J:115:TYR:C	1:J:116:LEU:HD22	2.43	0.44
1:C:55:GLN:HG3	1:C:116:LEU:CD1	2.41	0.44
1:D:178:LYS:O	1:D:178:LYS:HG3	2.18	0.44
1:E:187:CYS:CB	1:E:189:GLU:HB2	2.46	0.44
1:H:165:ILE:HA	1:H:166:PRO:HD3	1.89	0.44
1:J:65:TRP:CE2	1:J:109:SER:HA	2.53	0.44
1:A:76:VAL:HG23	1:A:107:VAL:CG2	2.48	0.44
1:B:35:LEU:HD21	1:B:50:VAL:HG13	1.99	0.44
1:B:52:PHE:CD1	1:B:52:PHE:N	2.86	0.44
1:C:148:HIS:CG	1:C:189:GLU:HG2	2.53	0.44
1:I:90:ALA:HB1	1:I:123:PHE:HE2	1.83	0.44
1:J:49:ASP:OD2	1:J:122:ARG:HD2	2.18	0.44
1:C:115:TYR:C	1:C:116:LEU:HD22	2.43	0.43
1:D:42:ASP:CB	1:D:47:VAL:HG22	2.48	0.43
1:D:167:TYR:CD1	1:D:167:TYR:N	2.84	0.43
1:F:108:ASN:ND2	1:F:112:HIS:HB3	2.26	0.43
1:B:41:VAL:HG12	1:B:48:VAL:HG12	2.01	0.43
1:F:42:ASP:CB	1:F:47:VAL:HG22	2.47	0.43
1:F:78:VAL:HA	1:F:79:PRO:HD3	1.83	0.43
1:C:19:ILE:HG12	1:C:21:THR:HG23	1.99	0.43
1:C:41:VAL:HG23	1:C:41:VAL:O	2.17	0.43
1:D:52:PHE:CE2	1:D:119:ILE:HB	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:29:VAL:O	1:E:29:VAL:HG23	2.19	0.43
1:B:22:GLN:HG3	1:B:25:ARG:HG3	2.00	0.43
1:C:83:LEU:HD23	1:C:83:LEU:H	1.83	0.43
1:D:115:TYR:C	1:D:116:LEU:HD22	2.43	0.43
1:E:83:LEU:HD23	1:E:83:LEU:H	1.83	0.43
1:D:188:LYS:HD3	1:D:188:LYS:C	2.43	0.43
1:E:108:ASN:ND2	1:E:112:HIS:HB3	2.24	0.43
1:H:89:ALA:HB3	1:H:145:TRP:CE3	2.53	0.43
1:C:49:ASP:OD2	1:C:122:ARG:HD2	2.18	0.43
1:C:108:ASN:ND2	1:C:112:HIS:HB3	2.34	0.43
2:L:1:NAG:C4	2:L:2:NAG:N2	2.81	0.43
1:B:176:THR:OG1	1:B:197:THR:HB	2.19	0.43
1:C:167:TYR:CD1	1:C:167:TYR:N	2.82	0.43
1:C:181:GLU:HG2	1:C:190:PRO:HB2	2.01	0.43
1:H:41:VAL:HG23	1:H:41:VAL:O	2.19	0.43
1:J:70:TYR:O	1:J:71:PRO:C	2.62	0.43
1:G:36:LEU:CD2	1:G:53:TRP:HB2	2.49	0.43
1:G:76:VAL:HG23	1:G:107:VAL:CG2	2.49	0.43
1:J:167:TYR:CD1	1:J:167:TYR:N	2.82	0.43
1:E:165:ILE:HA	1:E:166:PRO:HD3	1.93	0.43
1:F:191:TYR:HA	1:F:192:PRO:HD2	1.90	0.43
1:I:200:PHE:CD1	1:I:200:PHE:N	2.87	0.43
1:B:65:TRP:CE2	1:B:109:SER:HA	2.54	0.42
1:D:41:VAL:HG23	1:D:41:VAL:O	2.19	0.42
1:D:55:GLN:CG	1:D:116:LEU:HD11	2.41	0.42
1:C:36:LEU:HB2	1:C:51:VAL:O	2.19	0.42
1:D:22:GLN:O	1:D:23:ARG:O	2.37	0.42
1:J:19:ILE:HG12	1:J:21:THR:HG23	2.01	0.42
1:B:55:GLN:HG3	1:B:116:LEU:CD1	2.46	0.42
1:C:156:MET:CE	1:C:177:GLN:HE21	2.31	0.42
1:C:200:PHE:CD1	1:C:200:PHE:N	2.88	0.42
1:D:23:ARG:C	1:D:25:ARG:H	2.27	0.42
1:H:29:VAL:HG23	1:H:29:VAL:O	2.18	0.42
1:I:49:ASP:OD2	1:I:122:ARG:HD2	2.19	0.42
1:J:83:LEU:HD23	1:J:83:LEU:H	1.84	0.42
1:A:31:PHE:CE2	1:A:196:PHE:HE1	2.37	0.42
1:B:33:LEU:HB3	1:B:196:PHE:CE2	2.53	0.42
1:C:70:TYR:O	1:C:71:PRO:C	2.62	0.42
1:F:70:TYR:O	1:F:71:PRO:C	2.62	0.42
1:I:65:TRP:CE2	1:I:109:SER:HA	2.54	0.42
1:A:70:TYR:O	1:A:71:PRO:C	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:33:LEU:HB3	1:E:196:PHE:CE2	2.55	0.42
1:G:65:TRP:CE2	1:G:109:SER:HA	2.53	0.42
1:J:35:LEU:HD21	1:J:50:VAL:HG13	2.01	0.42
1:J:55:GLN:HG3	1:J:116:LEU:CD1	2.43	0.42
1:B:141:LYS:HZ1	1:B:182:ARG:NH2	2.16	0.42
1:C:33:LEU:HD23	1:C:34:SER:N	2.34	0.42
1:C:108:ASN:CG	3:C:901:NAG:N2	2.73	0.42
1:E:83:LEU:HD23	1:E:83:LEU:N	2.35	0.42
1:D:49:ASP:OD2	1:D:122:ARG:HD2	2.20	0.42
1:E:149:SER:OG	1:E:192:PRO:HD3	2.20	0.42
1:J:42:ASP:CB	1:J:47:VAL:HG22	2.49	0.42
1:A:165:ILE:HA	1:A:166:PRO:HD3	1.91	0.42
1:B:22:GLN:CG	1:B:25:ARG:HB2	2.50	0.42
1:E:35:LEU:HD21	1:E:50:VAL:HG13	2.02	0.42
1:F:139:LYS:C	1:F:140:LEU:HD12	2.45	0.42
1:F:186:CYS:O	1:F:187:CYS:O	2.38	0.42
1:I:44:LYS:HD3	1:J:169:ARG:HH12	1.85	0.42
1:J:76:VAL:HG23	1:J:107:VAL:CG2	2.50	0.42
1:E:65:TRP:CE2	1:E:109:SER:HA	2.55	0.42
1:E:185:GLU:O	1:E:186:CYS:CB	2.54	0.42
1:G:35:LEU:HD21	1:G:50:VAL:HG13	2.02	0.42
1:G:49:ASP:OD2	1:G:122:ARG:HD2	2.19	0.42
1:I:161:ILE:O	1:I:161:ILE:HG12	2.20	0.42
1:A:42:ASP:CB	1:A:47:VAL:HG22	2.50	0.42
1:D:31:PHE:HE2	1:D:196:PHE:HE1	1.68	0.42
1:D:80:ILE:C	1:D:80:ILE:HD12	2.45	0.42
1:D:108:ASN:ND2	1:D:112:HIS:HB3	2.27	0.42
1:A:148:HIS:CB	1:A:189:GLU:HB3	2.48	0.41
1:B:115:TYR:C	1:B:116:LEU:HD22	2.45	0.41
1:H:76:VAL:HG23	1:H:107:VAL:CG2	2.49	0.41
1:I:3:GLN:HB3	1:I:71:PRO:HG2	2.00	0.41
1:J:176:THR:OG1	1:J:197:THR:HB	2.20	0.41
1:A:23:ARG:C	1:A:25:ARG:H	2.28	0.41
1:C:33:LEU:HD23	1:C:33:LEU:C	2.45	0.41
1:C:141:LYS:HZ2	1:C:182:ARG:NH1	2.18	0.41
1:C:148:HIS:HA	1:C:191:TYR:HD2	1.84	0.41
1:G:200:PHE:CD1	1:G:200:PHE:N	2.89	0.41
1:A:161:ILE:HG12	1:A:161:ILE:O	2.19	0.41
1:B:29:VAL:O	1:B:29:VAL:HG23	2.19	0.41
1:C:44:LYS:CG	1:D:169:ARG:NH1	2.84	0.41
1:E:70:TYR:O	1:E:71:PRO:C	2.63	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:29:VAL:HG23	1:I:29:VAL:O	2.19	0.41
1:I:76:VAL:CG2	1:I:107:VAL:HG23	2.50	0.41
1:C:44:LYS:HD3	1:D:169:ARG:NH1	2.34	0.41
1:H:31:PHE:HE2	1:H:196:PHE:HE1	1.68	0.41
1:H:70:TYR:O	1:H:71:PRO:C	2.63	0.41
1:I:78:VAL:HA	1:I:79:PRO:HD3	1.82	0.41
1:A:65:TRP:CE2	1:A:109:SER:HA	2.56	0.41
1:C:65:TRP:CE2	1:C:109:SER:HA	2.55	0.41
1:D:18:VAL:HG22	1:E:7:TYR:CE2	2.56	0.41
1:H:176:THR:OG1	1:H:197:THR:HB	2.20	0.41
1:C:76:VAL:HG23	1:C:107:VAL:CG2	2.50	0.41
1:E:49:ASP:OD2	1:E:122:ARG:HD2	2.20	0.41
1:I:70:TYR:O	1:I:71:PRO:C	2.63	0.41
1:E:76:VAL:HG23	1:E:107:VAL:CG2	2.51	0.41
1:G:41:VAL:HG12	1:G:48:VAL:HG12	2.02	0.41
1:I:182:ARG:HH11	1:I:182:ARG:CG	2.29	0.41
1:J:200:PHE:CD1	1:J:200:PHE:N	2.88	0.41
1:A:80:ILE:C	1:A:80:ILE:HD12	2.45	0.41
1:A:139:LYS:C	1:A:140:LEU:HD12	2.46	0.41
1:B:78:VAL:HA	1:B:79:PRO:HD3	1.82	0.41
1:B:150:ARG:HH12	1:B:190:PRO:HD3	1.86	0.41
1:C:83:LEU:HD23	1:C:83:LEU:N	2.36	0.41
1:C:142:PHE:O	1:C:193:ASP:HB2	2.20	0.41
1:F:35:LEU:HD21	1:F:50:VAL:HG13	2.02	0.41
1:A:112:HIS:HD2	3:A:901:NAG:H82	1.85	0.41
1:B:70:TYR:O	1:B:71:PRO:C	2.63	0.41
1:C:80:ILE:C	1:C:80:ILE:HD12	2.46	0.41
1:C:149:SER:HB2	1:C:190:PRO:HG2	2.02	0.41
1:F:70:TYR:O	1:F:71:PRO:O	2.39	0.41
1:F:80:ILE:HD12	1:F:80:ILE:C	2.46	0.41
1:F:115:TYR:C	1:F:116:LEU:HD22	2.45	0.41
1:G:70:TYR:O	1:G:71:PRO:C	2.63	0.41
1:G:145:TRP:HZ2	1:H:116:LEU:O	2.04	0.41
1:H:41:VAL:HG12	1:H:48:VAL:HG12	2.02	0.41
1:J:108:ASN:ND2	1:J:112:HIS:HB3	2.35	0.41
1:A:41:VAL:O	1:A:41:VAL:HG23	2.21	0.41
1:B:23:ARG:C	1:B:25:ARG:H	2.28	0.41
1:H:65:TRP:CE2	1:H:109:SER:HA	2.56	0.41
1:H:145:TRP:CZ2	1:I:101:THR:HG21	2.55	0.41
1:C:12:LYS:HE2	1:H:23:ARG:HH22	1.85	0.40
1:C:35:LEU:HD21	1:C:50:VAL:HG13	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:TYR:O	1:C:71:PRO:O	2.39	0.40
1:D:3:GLN:HE21	1:D:3:GLN:CA	2.32	0.40
1:E:15:ASN:HA	1:E:16:PRO:HD2	1.93	0.40
1:G:96:LYS:HD3	1:H:99:VAL:O	2.20	0.40
1:J:83:LEU:HD23	1:J:83:LEU:N	2.35	0.40
1:B:161:ILE:O	1:B:161:ILE:HG12	2.21	0.40
1:H:83:LEU:HD23	1:H:83:LEU:H	1.87	0.40
1:H:200:PHE:CD1	1:H:200:PHE:N	2.89	0.40
1:I:150:ARG:HH11	1:I:190:PRO:HD2	1.85	0.40
1:B:18:VAL:HG22	1:C:7:TYR:CD2	2.56	0.40
1:D:65:TRP:CZ2	1:D:109:SER:HA	2.56	0.40
1:E:182:ARG:O	1:E:190:PRO:HA	2.21	0.40
1:D:139:LYS:C	1:D:140:LEU:HD12	2.46	0.40
1:G:29:VAL:O	1:G:29:VAL:HG23	2.21	0.40
1:I:70:TYR:O	1:I:71:PRO:O	2.40	0.40
1:I:113:VAL:HG12	1:I:114:GLN:N	2.36	0.40
1:A:29:VAL:O	1:A:29:VAL:HG23	2.21	0.40
1:F:39:MET:O	1:J:45:ASN:ND2	2.55	0.40
1:F:140:LEU:HD12	1:F:140:LEU:N	2.37	0.40
1:F:165:ILE:HA	1:F:166:PRO:HD3	1.89	0.40
1:G:78:VAL:HA	1:G:79:PRO:HD3	1.85	0.40
1:I:41:VAL:HG23	1:I:41:VAL:O	2.21	0.40
1:I:186:CYS:O	1:I:187:CYS:C	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	200/204 (98%)	173 (86%)	21 (10%)	6 (3%)	3	19
1	B	200/204 (98%)	168 (84%)	23 (12%)	9 (4%)	2	12

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	200/204 (98%)	171 (86%)	26 (13%)	3 (2%)	8	32
1	D	200/204 (98%)	171 (86%)	22 (11%)	7 (4%)	3	16
1	E	200/204 (98%)	171 (86%)	24 (12%)	5 (2%)	4	21
1	F	200/204 (98%)	173 (86%)	22 (11%)	5 (2%)	4	21
1	G	200/204 (98%)	169 (84%)	25 (12%)	6 (3%)	3	19
1	H	200/204 (98%)	173 (86%)	22 (11%)	5 (2%)	4	21
1	I	200/204 (98%)	168 (84%)	24 (12%)	8 (4%)	2	13
1	J	200/204 (98%)	172 (86%)	22 (11%)	6 (3%)	3	19
All	All	2000/2040 (98%)	1709 (86%)	231 (12%)	60 (3%)	3	19

All (60) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	71	PRO
1	A	187	CYS
1	B	23	ARG
1	B	71	PRO
1	B	184	TYR
1	B	186	CYS
1	B	187	CYS
1	C	23	ARG
1	C	71	PRO
1	D	23	ARG
1	D	71	PRO
1	D	187	CYS
1	E	23	ARG
1	E	71	PRO
1	E	186	CYS
1	E	187	CYS
1	F	23	ARG
1	F	71	PRO
1	F	187	CYS
1	G	23	ARG
1	G	71	PRO
1	G	186	CYS
1	H	23	ARG
1	H	71	PRO
1	I	23	ARG
1	I	71	PRO

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Mol	Chain	Res	Type
1	I	186	CYS
1	J	23	ARG
1	J	71	PRO
1	J	186	CYS
1	A	23	ARG
1	A	167	TYR
1	B	185	GLU
1	C	167	TYR
1	D	167	TYR
1	D	188	LYS
1	E	167	TYR
1	F	167	TYR
1	G	167	TYR
1	H	167	TYR
1	I	185	GLU
1	I	187	CYS
1	J	167	TYR
1	A	186	CYS
1	B	167	TYR
1	B	188	LYS
1	D	183	PHE
1	G	187	CYS
1	I	167	TYR
1	J	185	GLU
1	I	188	LYS
1	B	168	SER
1	H	168	SER
1	H	186	CYS
1	I	168	SER
1	A	168	SER
1	D	168	SER
1	F	168	SER
1	G	168	SER
1	J	168	SER

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	188/190 (99%)	176 (94%)	12 (6%)	16	44
1	B	188/190 (99%)	178 (95%)	10 (5%)	20	51
1	C	188/190 (99%)	177 (94%)	11 (6%)	18	48
1	D	188/190 (99%)	179 (95%)	9 (5%)	23	54
1	E	188/190 (99%)	179 (95%)	9 (5%)	23	54
1	F	188/190 (99%)	178 (95%)	10 (5%)	20	51
1	G	188/190 (99%)	179 (95%)	9 (5%)	23	54
1	H	188/190 (99%)	179 (95%)	9 (5%)	23	54
1	I	188/190 (99%)	178 (95%)	10 (5%)	20	51
1	J	188/190 (99%)	178 (95%)	10 (5%)	20	51
All	All	1880/1900 (99%)	1781 (95%)	99 (5%)	20	51

All (99) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	THR
1	A	22	GLN
1	A	35	LEU
1	A	47	VAL
1	A	52	PHE
1	A	71	PRO
1	A	107	VAL
1	A	165	ILE
1	A	167	TYR
1	A	179	ARG
1	A	186	CYS
1	A	187	CYS
1	B	21	THR
1	B	35	LEU
1	B	47	VAL
1	B	71	PRO
1	B	92	ASN
1	B	107	VAL
1	B	165	ILE
1	B	167	TYR
1	B	179	ARG
1	B	187	CYS
1	C	35	LEU
1	C	47	VAL

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Mol	Chain	Res	Type
1	C	71	PRO
1	C	92	ASN
1	C	107	VAL
1	C	165	ILE
1	C	167	TYR
1	C	179	ARG
1	C	185	GLU
1	C	186	CYS
1	C	187	CYS
1	D	21	THR
1	D	35	LEU
1	D	47	VAL
1	D	71	PRO
1	D	92	ASN
1	D	107	VAL
1	D	165	ILE
1	D	167	TYR
1	D	179	ARG
1	E	35	LEU
1	E	47	VAL
1	E	71	PRO
1	E	92	ASN
1	E	107	VAL
1	E	165	ILE
1	E	167	TYR
1	E	179	ARG
1	E	189	GLU
1	F	35	LEU
1	F	47	VAL
1	F	71	PRO
1	F	92	ASN
1	F	107	VAL
1	F	165	ILE
1	F	167	TYR
1	F	179	ARG
1	F	183	PHE
1	F	187	CYS
1	G	35	LEU
1	G	47	VAL
1	G	54	LEU
1	G	71	PRO
1	G	92	ASN

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Mol	Chain	Res	Type
1	G	107	VAL
1	G	165	ILE
1	G	167	TYR
1	G	179	ARG
1	H	35	LEU
1	H	47	VAL
1	H	71	PRO
1	H	107	VAL
1	H	165	ILE
1	H	167	TYR
1	H	179	ARG
1	H	183	PHE
1	H	187	CYS
1	I	3	GLN
1	I	21	THR
1	I	35	LEU
1	I	47	VAL
1	I	71	PRO
1	I	107	VAL
1	I	165	ILE
1	I	167	TYR
1	I	179	ARG
1	I	187	CYS
1	J	35	LEU
1	J	47	VAL
1	J	54	LEU
1	J	71	PRO
1	J	92	ASN
1	J	107	VAL
1	J	165	ILE
1	J	167	TYR
1	J	179	ARG
1	J	187	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (80) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	GLN
1	A	112	HIS
1	A	157	GLN
1	A	177	GLN
1	B	3	GLN

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Mol	Chain	Res	Type
1	B	37	GLN
1	B	45	ASN
1	B	103	GLN
1	B	148	HIS
1	B	157	GLN
1	B	177	GLN
1	C	3	GLN
1	C	37	GLN
1	C	45	ASN
1	C	92	ASN
1	C	103	GLN
1	C	112	HIS
1	C	148	HIS
1	C	157	GLN
1	C	177	GLN
1	D	3	GLN
1	D	22	GLN
1	D	37	GLN
1	D	92	ASN
1	D	103	GLN
1	D	108	ASN
1	D	112	HIS
1	D	148	HIS
1	D	157	GLN
1	D	177	GLN
1	E	3	GLN
1	E	37	GLN
1	E	45	ASN
1	E	103	GLN
1	E	112	HIS
1	E	157	GLN
1	E	177	GLN
1	F	3	GLN
1	F	37	GLN
1	F	92	ASN
1	F	103	GLN
1	F	108	ASN
1	F	148	HIS
1	F	157	GLN
1	F	177	GLN
1	G	3	GLN
1	G	22	GLN

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Mol	Chain	Res	Type
1	G	37	GLN
1	G	45	ASN
1	G	92	ASN
1	G	103	GLN
1	G	148	HIS
1	G	157	GLN
1	G	177	GLN
1	H	37	GLN
1	H	45	ASN
1	H	92	ASN
1	H	103	GLN
1	H	112	HIS
1	H	121	GLN
1	H	157	GLN
1	I	3	GLN
1	I	22	GLN
1	I	37	GLN
1	I	45	ASN
1	I	103	GLN
1	I	112	HIS
1	I	148	HIS
1	I	157	GLN
1	I	177	GLN
1	J	3	GLN
1	J	37	GLN
1	J	45	ASN
1	J	92	ASN
1	J	103	GLN
1	J	112	HIS
1	J	114	GLN
1	J	147	HIS
1	J	157	GLN
1	J	177	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

6 monosaccharides are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	NAG	K	1	2,1	14,14,15	0.93	1 (7%)	17,19,21	0.62	0
2	NAG	K	2	2	14,14,15	0.98	1 (7%)	17,19,21	0.88	1 (5%)
2	NAG	L	1	2,1	14,14,15	1.15	2 (14%)	17,19,21	1.24	3 (17%)
2	NAG	L	2	2	14,14,15	1.14	2 (14%)	17,19,21	0.72	0
2	NAG	M	1	2,1	14,14,15	0.88	0	17,19,21	0.88	1 (5%)
2	NAG	M	2	2	14,14,15	0.99	1 (7%)	17,19,21	0.96	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	NAG	K	1	2,1	-	2/6/23/26	0/1/1/1
2	NAG	K	2	2	-	4/6/23/26	0/1/1/1
2	NAG	L	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	L	2	2	-	2/6/23/26	0/1/1/1
2	NAG	M	1	2,1	-	4/6/23/26	0/1/1/1
2	NAG	M	2	2	-	5/6/23/26	0/1/1/1

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	2	NAG	C1-C2	3.11	1.56	1.52
2	K	2	NAG	C1-C2	2.53	1.55	1.52
2	L	1	NAG	C3-C2	2.52	1.57	1.52
2	L	1	NAG	C4-C3	2.38	1.58	1.52
2	M	2	NAG	C1-C2	2.34	1.55	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	K	1	NAG	C1-C2	2.22	1.55	1.52
2	L	2	NAG	C3-C2	2.12	1.57	1.52

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	L	1	NAG	C4-C3-C2	3.11	115.57	111.02
2	M	2	NAG	C1-O5-C5	3.08	116.31	112.19
2	M	1	NAG	C1-O5-C5	2.73	115.84	112.19
2	L	1	NAG	O5-C1-C2	-2.30	107.73	111.29
2	K	2	NAG	C1-O5-C5	2.29	115.26	112.19
2	L	1	NAG	C3-C4-C5	2.09	114.02	110.23

There are no chirality outliers.

All (21) torsion outliers are listed below:

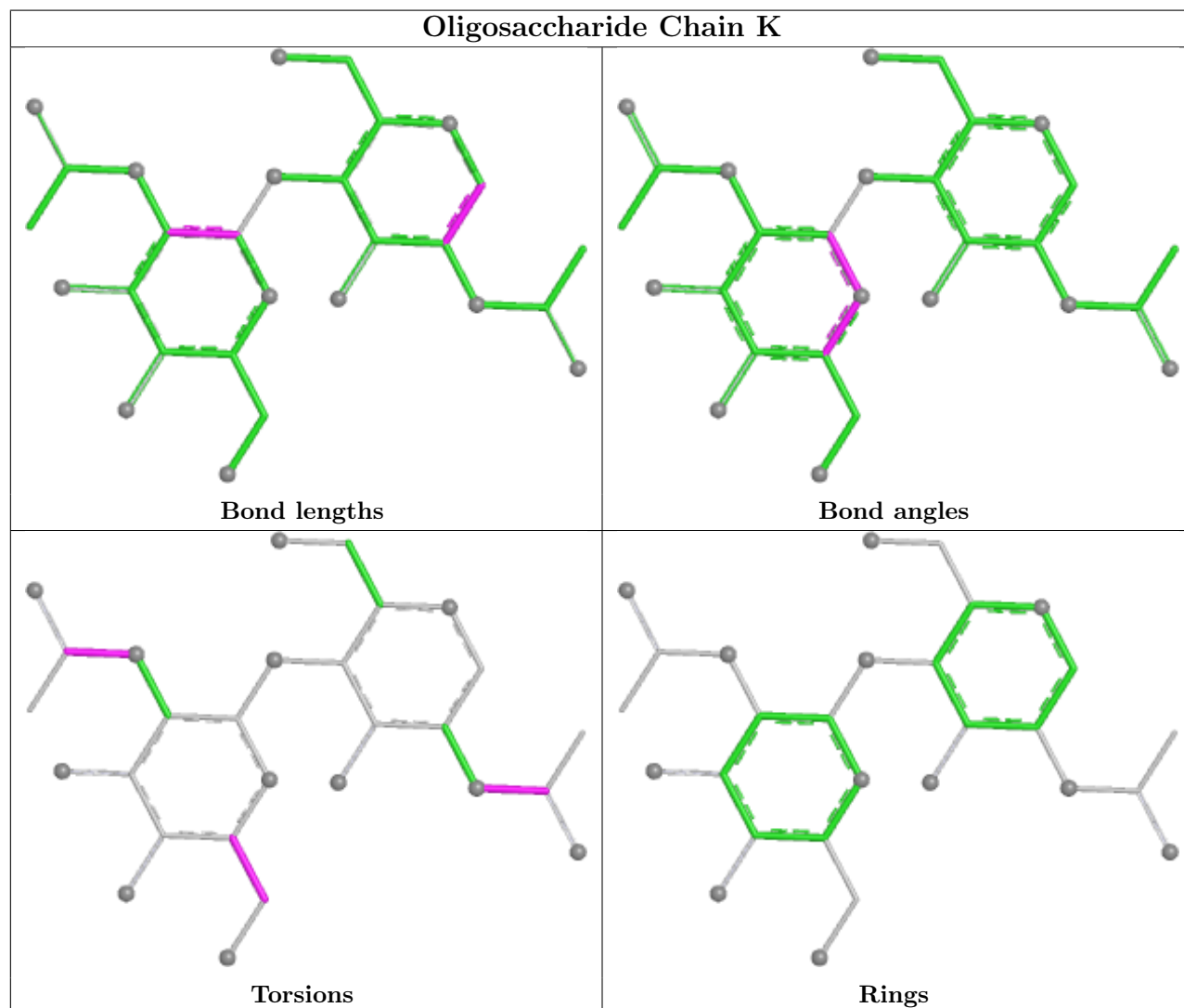
Mol	Chain	Res	Type	Atoms
2	M	2	NAG	C1-C2-N2-C7
2	M	2	NAG	C8-C7-N2-C2
2	M	2	NAG	O7-C7-N2-C2
2	M	1	NAG	C8-C7-N2-C2
2	M	1	NAG	O7-C7-N2-C2
2	M	2	NAG	C4-C5-C6-O6
2	L	2	NAG	O5-C5-C6-O6
2	M	1	NAG	O5-C5-C6-O6
2	M	1	NAG	C4-C5-C6-O6
2	L	1	NAG	C8-C7-N2-C2
2	L	2	NAG	C4-C5-C6-O6
2	K	1	NAG	C8-C7-N2-C2
2	K	1	NAG	O7-C7-N2-C2
2	L	1	NAG	O7-C7-N2-C2
2	K	2	NAG	C8-C7-N2-C2
2	M	2	NAG	O5-C5-C6-O6
2	K	2	NAG	O5-C5-C6-O6
2	L	1	NAG	C4-C5-C6-O6
2	K	2	NAG	O7-C7-N2-C2
2	K	2	NAG	C4-C5-C6-O6
2	L	1	NAG	O5-C5-C6-O6

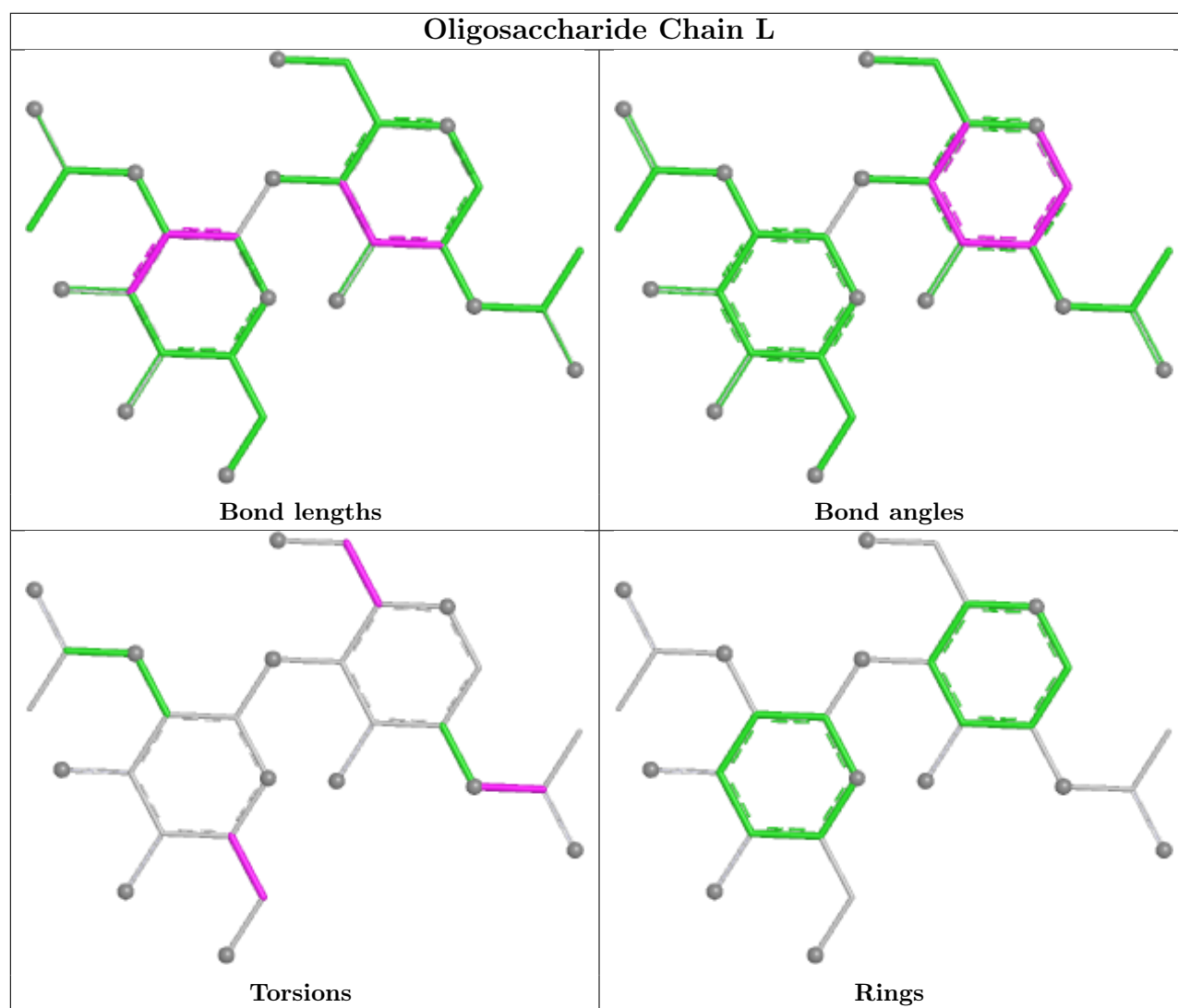
There are no ring outliers.

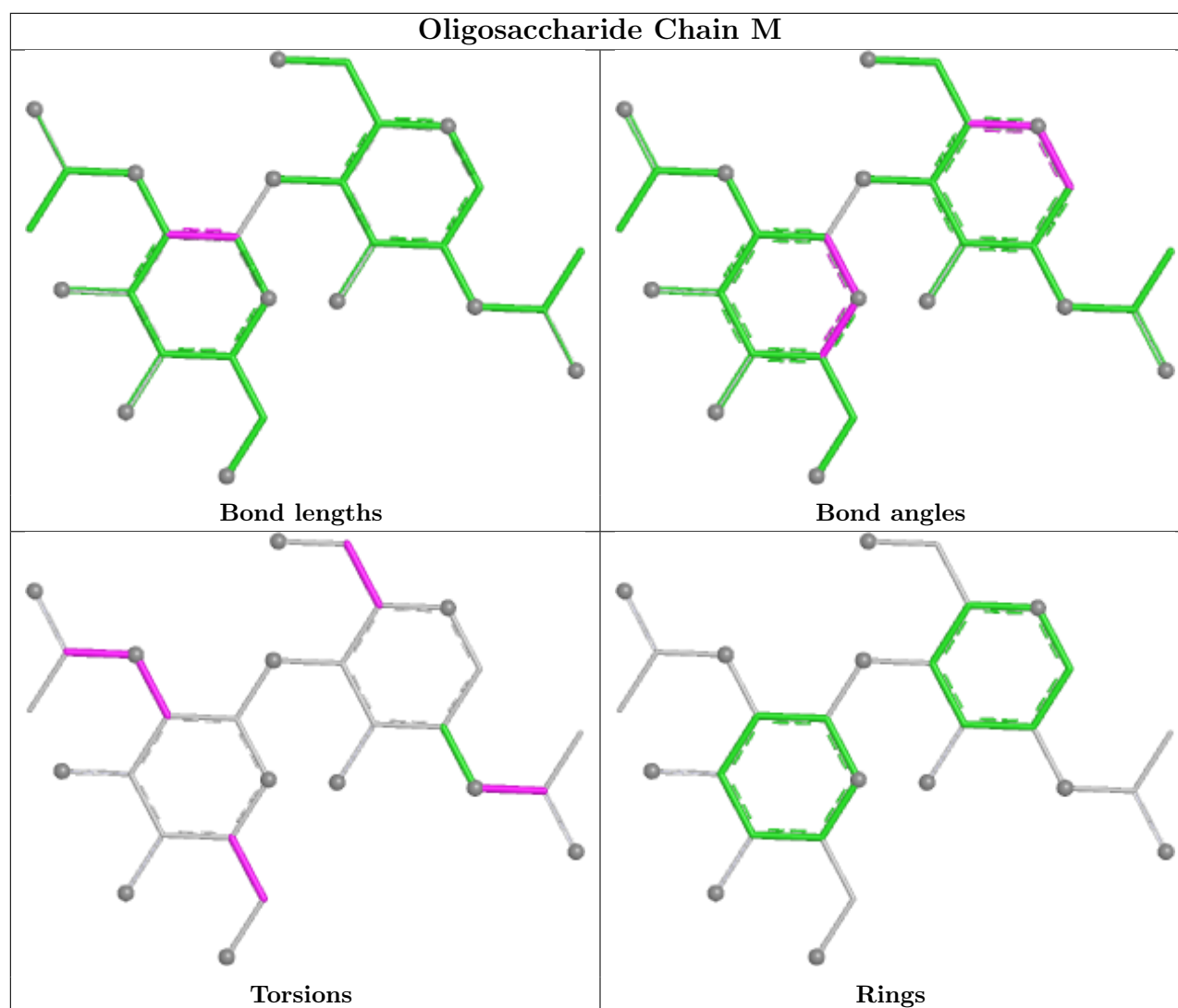
3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	M	1	NAG	1	0
2	L	1	NAG	3	0
2	L	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](#)

14 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	$\# Z > 2$
3	NAG	J	801	1	14,14,15	0.95	1 (7%)	17,19,21	0.58	0
3	NAG	A	901	1	14,14,15	0.77	0	17,19,21	0.67	0
3	NAG	C	901	1	14,14,15	0.98	1 (7%)	17,19,21	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	G	901	1	14,14,15	0.95	2 (14%)	17,19,21	1.35	2 (11%)
3	NAG	E	901	1	14,14,15	0.79	0	17,19,21	0.93	0
3	NAG	J	901	1	14,14,15	0.81	0	17,19,21	0.77	0
3	NAG	F	801	1	14,14,15	0.98	1 (7%)	17,19,21	0.99	1 (5%)
3	NAG	B	901	1	14,14,15	1.10	2 (14%)	17,19,21	1.21	1 (5%)
3	NAG	C	801	1	14,14,15	0.87	1 (7%)	17,19,21	0.84	1 (5%)
3	NAG	H	801	1	14,14,15	0.82	0	17,19,21	0.75	0
3	NAG	A	801	1	14,14,15	1.05	2 (14%)	17,19,21	1.16	1 (5%)
3	NAG	I	801	1	14,14,15	0.82	0	17,19,21	0.76	0
3	NAG	E	801	1	14,14,15	0.66	0	17,19,21	0.56	0
3	NAG	D	801	1	14,14,15	0.74	0	17,19,21	0.90	1 (5%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	J	801	1	-	3/6/23/26	0/1/1/1
3	NAG	A	901	1	-	4/6/23/26	0/1/1/1
3	NAG	C	901	1	-	4/6/23/26	0/1/1/1
3	NAG	G	901	1	-	5/6/23/26	0/1/1/1
3	NAG	E	901	1	-	3/6/23/26	0/1/1/1
3	NAG	J	901	1	-	3/6/23/26	0/1/1/1
3	NAG	F	801	1	-	2/6/23/26	0/1/1/1
3	NAG	B	901	1	-	5/6/23/26	0/1/1/1
3	NAG	C	801	1	-	4/6/23/26	0/1/1/1
3	NAG	H	801	1	-	2/6/23/26	0/1/1/1
3	NAG	A	801	1	-	1/6/23/26	0/1/1/1
3	NAG	I	801	1	-	0/6/23/26	0/1/1/1
3	NAG	E	801	1	-	4/6/23/26	0/1/1/1
3	NAG	D	801	1	-	2/6/23/26	0/1/1/1

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	901	NAG	C1-C2	2.88	1.56	1.52
3	C	901	NAG	C1-C2	2.66	1.56	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	J	801	NAG	C1-C2	2.62	1.55	1.52
3	F	801	NAG	C1-C2	2.56	1.55	1.52
3	A	801	NAG	C3-C2	2.51	1.57	1.52
3	C	801	NAG	C1-C2	2.47	1.55	1.52
3	A	801	NAG	C1-C2	2.42	1.55	1.52
3	B	901	NAG	C3-C2	2.13	1.57	1.52
3	G	901	NAG	C3-C2	2.03	1.56	1.52
3	G	901	NAG	C1-C2	2.01	1.55	1.52

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	G	901	NAG	C4-C3-C2	4.58	117.73	111.02
3	A	801	NAG	C4-C3-C2	3.85	116.66	111.02
3	B	901	NAG	C4-C3-C2	3.54	116.21	111.02
3	F	801	NAG	C1-O5-C5	2.87	116.03	112.19
3	C	801	NAG	C4-C3-C2	2.27	114.35	111.02
3	D	801	NAG	C2-N2-C7	-2.13	120.04	122.90
3	G	901	NAG	C3-C4-C5	2.04	113.94	110.23

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	901	NAG	C1-C2-N2-C7
3	A	901	NAG	C8-C7-N2-C2
3	A	901	NAG	O7-C7-N2-C2
3	B	901	NAG	C1-C2-N2-C7
3	B	901	NAG	C8-C7-N2-C2
3	B	901	NAG	O7-C7-N2-C2
3	C	901	NAG	C3-C2-N2-C7
3	C	901	NAG	C8-C7-N2-C2
3	C	901	NAG	O7-C7-N2-C2
3	E	901	NAG	C8-C7-N2-C2
3	E	901	NAG	O7-C7-N2-C2
3	G	901	NAG	C1-C2-N2-C7
3	G	901	NAG	C8-C7-N2-C2
3	G	901	NAG	O7-C7-N2-C2
3	J	901	NAG	C8-C7-N2-C2
3	J	901	NAG	O7-C7-N2-C2
3	J	801	NAG	C8-C7-N2-C2
3	J	801	NAG	O7-C7-N2-C2

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Mol	Chain	Res	Type	Atoms
3	C	801	NAG	O5-C5-C6-O6
3	E	801	NAG	O5-C5-C6-O6
3	E	801	NAG	C4-C5-C6-O6
3	F	801	NAG	C4-C5-C6-O6
3	F	801	NAG	O5-C5-C6-O6
3	B	901	NAG	C4-C5-C6-O6
3	E	801	NAG	C8-C7-N2-C2
3	B	901	NAG	O5-C5-C6-O6
3	G	901	NAG	C4-C5-C6-O6
3	C	801	NAG	C4-C5-C6-O6
3	C	801	NAG	C8-C7-N2-C2
3	E	801	NAG	O7-C7-N2-C2
3	E	901	NAG	O5-C5-C6-O6
3	G	901	NAG	O5-C5-C6-O6
3	C	801	NAG	O7-C7-N2-C2
3	C	901	NAG	O5-C5-C6-O6
3	A	801	NAG	O5-C5-C6-O6
3	J	801	NAG	O5-C5-C6-O6
3	A	901	NAG	O5-C5-C6-O6
3	J	901	NAG	O5-C5-C6-O6
3	D	801	NAG	C4-C5-C6-O6
3	D	801	NAG	O5-C5-C6-O6
3	H	801	NAG	C4-C5-C6-O6
3	H	801	NAG	O5-C5-C6-O6

There are no ring outliers.

4 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	901	NAG	2	0
3	C	901	NAG	4	0
3	F	801	NAG	2	0
3	B	901	NAG	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	202/204 (99%)	0.06	0 100 100	72, 108, 139, 150	0
1	B	202/204 (99%)	0.08	1 (0%) 87 73	74, 110, 137, 154	0
1	C	202/204 (99%)	0.15	1 (0%) 87 73	70, 108, 138, 158	0
1	D	202/204 (99%)	0.16	3 (1%) 72 52	74, 108, 138, 156	0
1	E	202/204 (99%)	0.12	0 100 100	75, 107, 136, 148	0
1	F	202/204 (99%)	0.15	0 100 100	73, 112, 140, 161	0
1	G	202/204 (99%)	0.05	0 100 100	75, 108, 138, 148	0
1	H	202/204 (99%)	0.14	2 (0%) 79 61	70, 107, 140, 153	0
1	I	202/204 (99%)	0.12	1 (0%) 87 73	76, 108, 138, 154	0
1	J	202/204 (99%)	0.11	1 (0%) 87 73	76, 110, 140, 155	0
All	All	2020/2040 (99%)	0.12	9 (0%) 88 76	70, 109, 140, 161	0

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	173	VAL	3.4
1	J	53	TRP	2.7
1	C	196	PHE	2.6
1	I	73	VAL	2.4
1	D	159	ALA	2.3
1	D	53	TRP	2.2
1	B	184	TYR	2.1
1	H	53	TRP	2.1
1	D	161	ILE	2.0

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

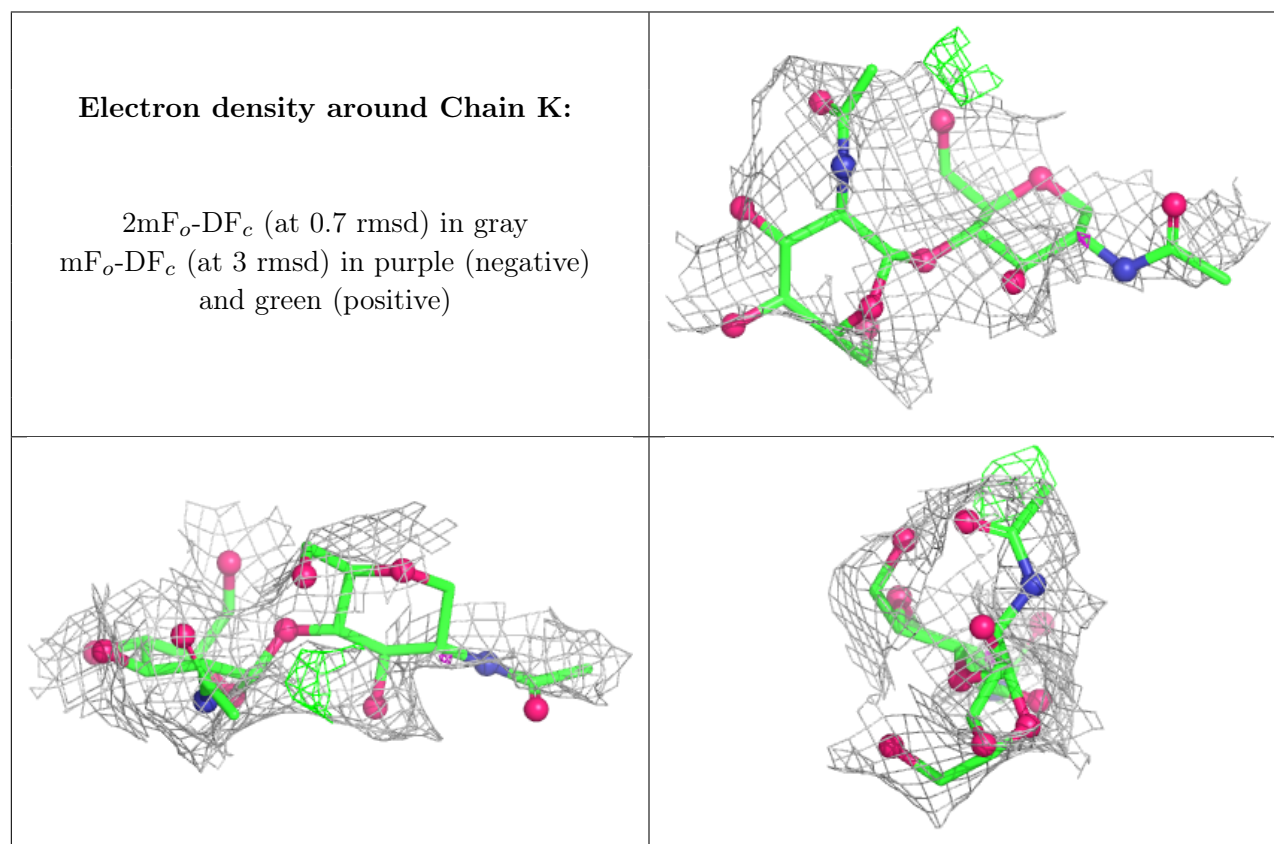
There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

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

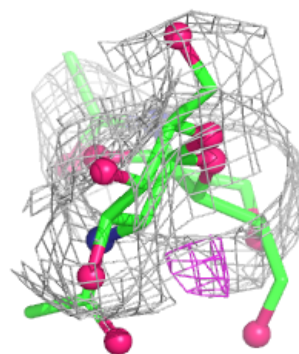
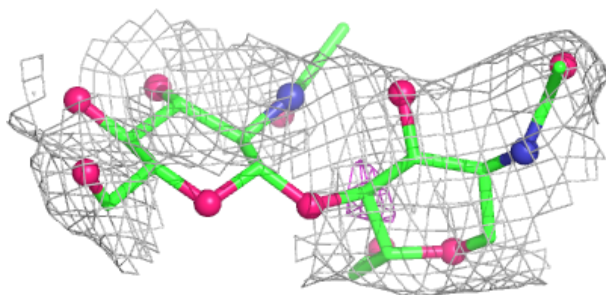
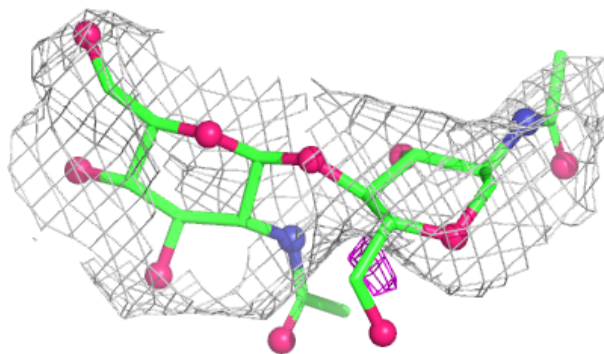
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	NAG	K	2	14/15	0.32	0.12	145,164,167,167	0
2	NAG	K	1	14/15	0.47	0.12	148,161,165,167	0
2	NAG	L	2	14/15	0.48	0.12	129,145,155,157	0
2	NAG	L	1	14/15	0.52	0.12	138,149,154,156	0
2	NAG	M	1	14/15	0.54	0.15	145,151,153,154	0
2	NAG	M	2	14/15	0.58	0.12	133,149,156,157	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.

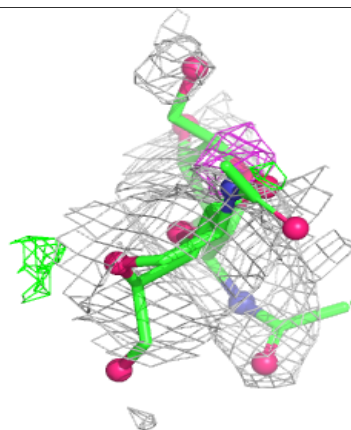
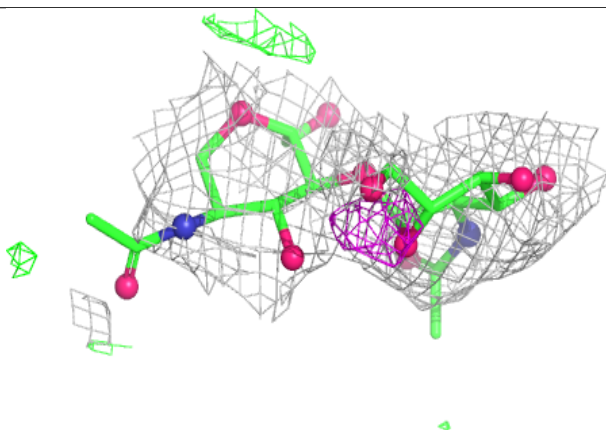
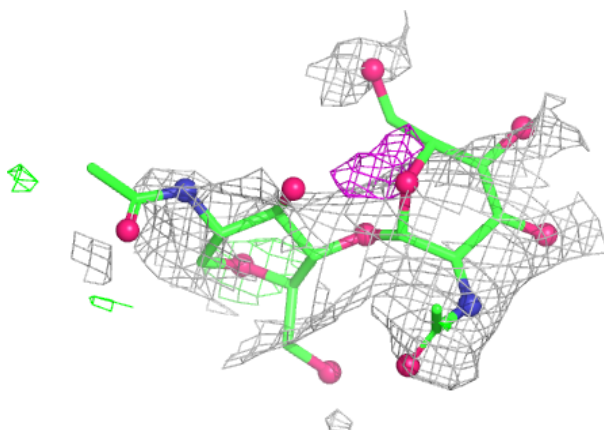


Electron density around Chain L:

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

**Electron density around Chain M:**

$2mF_o-DF_c$ (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
3	NAG	E	901	14/15	0.22	0.14	127,145,147,150	0
3	NAG	J	901	14/15	0.28	0.14	142,153,156,158	0
3	NAG	D	801	14/15	0.37	0.14	160,170,171,173	0
3	NAG	J	801	14/15	0.39	0.13	147,161,162,163	0
3	NAG	I	801	14/15	0.40	0.13	133,151,155,156	0
3	NAG	F	801	14/15	0.45	0.11	141,162,164,165	0
3	NAG	A	901	14/15	0.45	0.14	140,149,153,156	0
3	NAG	E	801	14/15	0.48	0.11	138,161,166,167	0
3	NAG	G	901	14/15	0.50	0.13	145,151,155,155	0
3	NAG	C	801	14/15	0.52	0.12	139,159,162,163	0
3	NAG	H	801	14/15	0.54	0.11	129,149,155,155	0
3	NAG	B	901	14/15	0.57	0.13	115,127,130,132	0
3	NAG	A	801	14/15	0.58	0.11	132,144,152,152	0
3	NAG	C	901	14/15	0.67	0.11	125,139,146,148	0

6.5 Other polymers [i](#)

There are no such residues in this entry.