



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 04:49 PM UTC

PDB ID : 5GS0 / pdb_00005gs0
Title : Crystal structure of the complex of TLR3 and bi-specific diabody
Authors : Kim, J.H.; Song, D.H.; Youn, S.J.; Kim, J.W.; Cho, G.; Lee, H.; Lee, J.O.
Deposited on : 2016-08-13
Resolution : 3.27 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

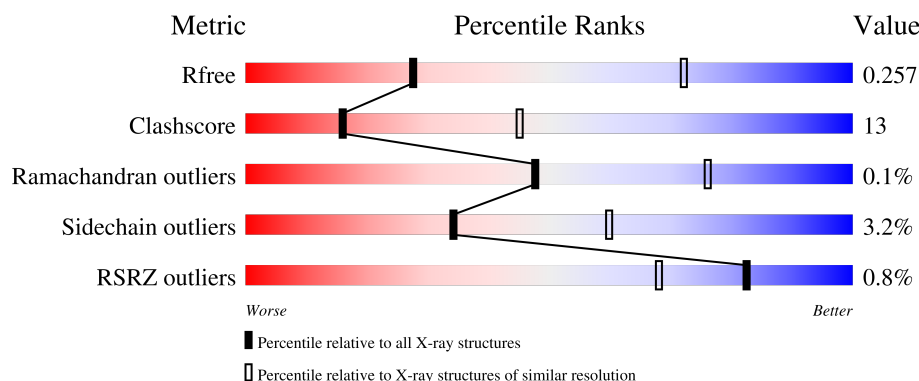
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.27 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	1303 (3.30-3.26)
Clashscore	190562	1354 (3.30-3.26)
Ramachandran outliers	187476	1334 (3.30-3.26)
Sidechain outliers	187428	1333 (3.30-3.26)
RSRZ outliers	180081	1303 (3.30-3.26)

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

Mol	Chain	Length	Quality of chain
1	A	671	69% 28% ..
1	B	671	73% 24% ..
2	C	107	68% 29% .
2	E	107	73% 21% 5% .
3	D	127	66% 29% ..

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Mol	Chain	Length	Quality of chain
3	F	127	% 65% 28%
4	H	121	3% 67% 28%
4	X	121	60% 32% 6%
5	L	107	6% 64% 32%
5	Y	107	2% 63% 34%

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 18219 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 Toll-like receptor 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	663	Total	C	N	O	S	0	0	0
			5323	3407	903	996	17			
1	B	663	Total	C	N	O	S	0	0	0
			5323	3407	903	996	17			

- Molecule 2 is a protein called light chain (anti-TLR3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	107	Total	C	N	O	S	0	0	0
			809	504	133	170	2			
2	E	107	Total	C	N	O	S	0	0	0
			809	504	133	170	2			

- Molecule 3 is a protein called heavy chain (anti-TLR3).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	122	Total	C	N	O	S	0	0	0
			956	609	161	184	2			
3	F	122	Total	C	N	O	S	0	0	0
			956	609	161	184	2			

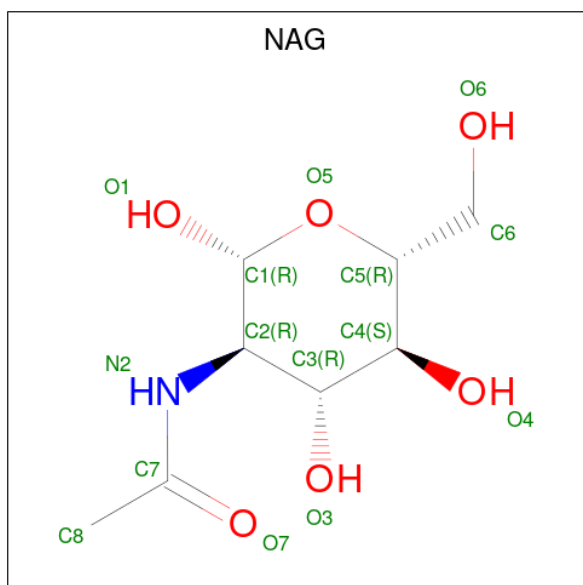
- Molecule 4 is a protein called heavy chain (anti-Lid).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	X	114	Total	C	N	O	S	0	0	0
			909	582	151	173	3			
4	H	116	Total	C	N	O	S	0	0	0
			921	588	153	177	3			

- Molecule 5 is a protein called light chain (anti-Lid).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
5	Y	107	Total	C	N	O	S	0	0	0
			824	514	141	167	2			
5	L	105	Total	C	N	O	S	0	0	0
			807	502	138	165	2			

- Molecule 6 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		

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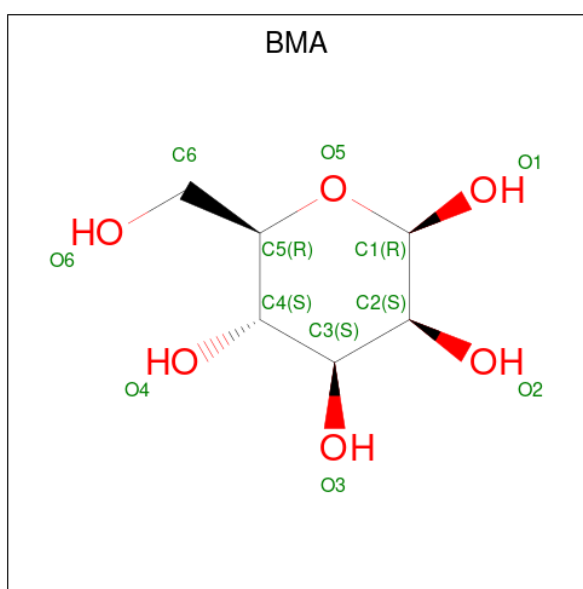
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	A	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		

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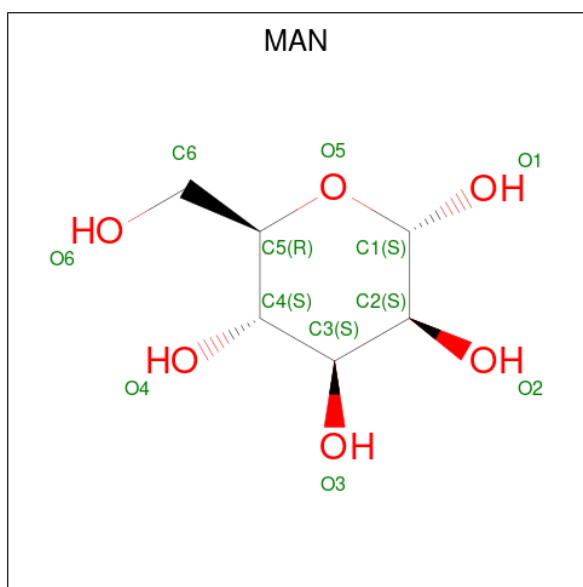
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		
6	B	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 7 is beta-D-mannopyranose (CCD ID: BMA) (formula: $C_6H_{12}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
7	A	1	Total	C	O	0	0
			12	6	6		
7	A	1	Total	C	O	0	0
			12	6	6		
7	B	1	Total	C	O	0	0
			12	6	6		
7	B	1	Total	C	O	0	0
			12	6	6		

- Molecule 8 is alpha-D-mannopyranose (CCD ID: MAN) (formula: $C_6H_{12}O_6$).

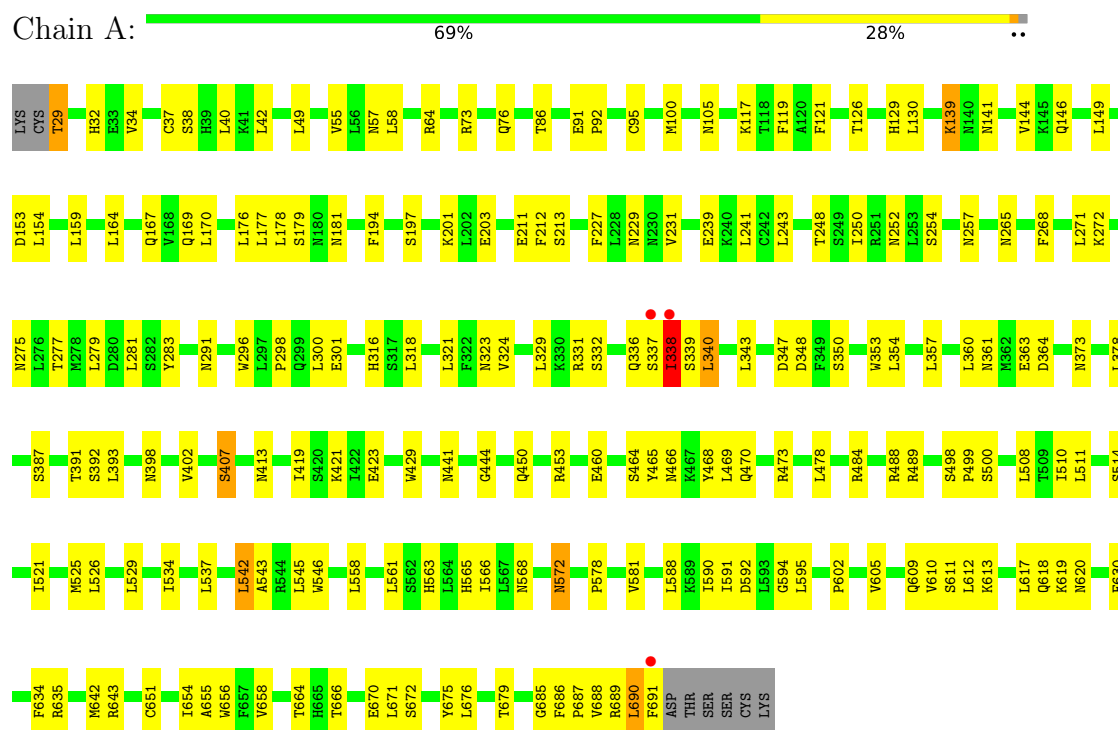


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
8	A	1	Total	C	O	0	0
			12	6	6		
8	B	1	Total	C	O	0	0
			12	6	6		

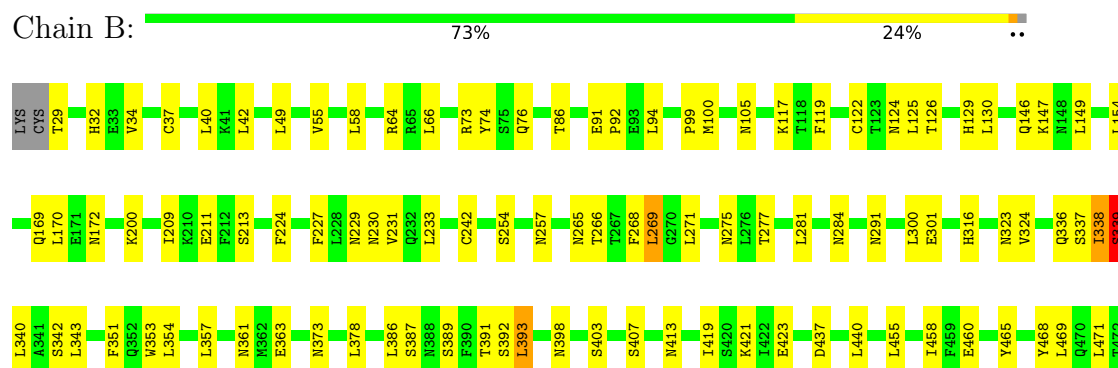
3 Residue-property plots

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: Toll-like receptor 3



• Molecule 1: Toll-like receptor 3





- Molecule 2: light chain (anti-TLR3)

Chain C: 68% 29% .



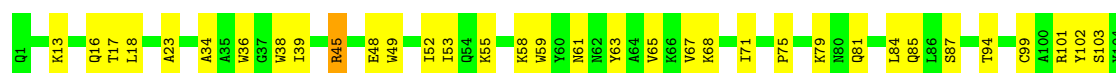
- Molecule 2: light chain (anti-TLR3)

Chain E: 73% 21% 5% .



- Molecule 3: heavy chain (anti-TLR3)

Chain D: 66% 29% . .



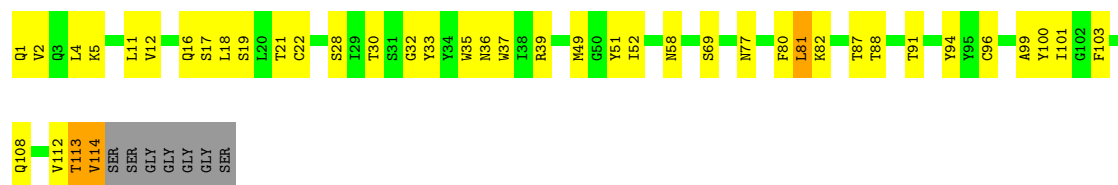
- Molecule 3: heavy chain (anti-TLR3)

Chain F: 65% 28% . .

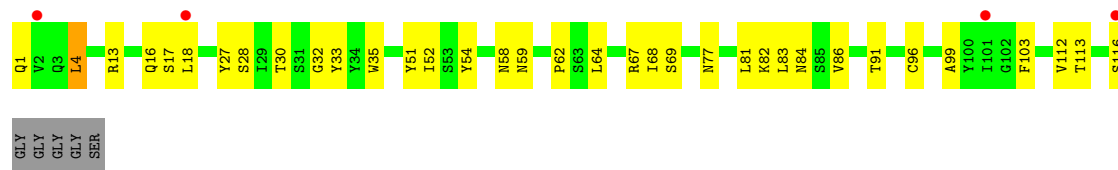


- Molecule 4: heavy chain (anti-Lid)

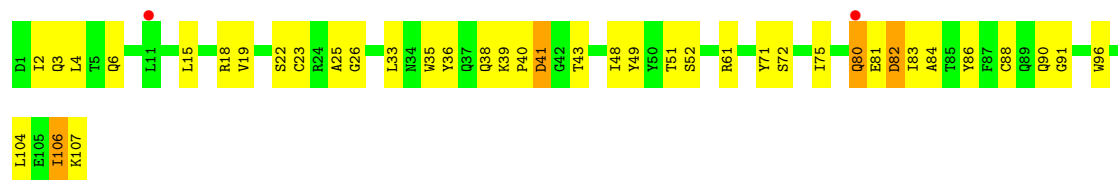
Chain X: 60% 32% . 6%



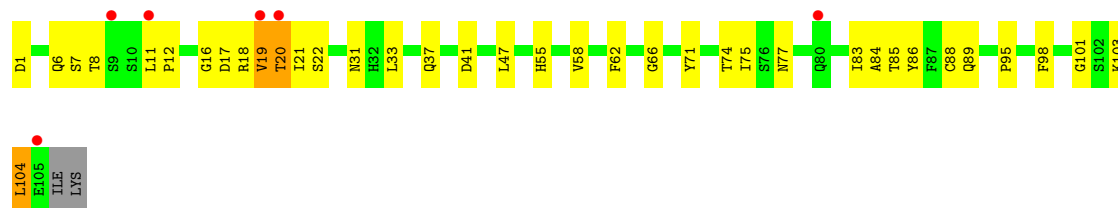
- Molecule 4: heavy chain (anti-Lid)



- Molecule 5: light chain (anti-Lid)



- Molecule 5: light chain (anti-Lid)



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	98.52Å 141.33Å 150.83Å 90.00° 106.88° 90.00°	Depositor
Resolution (Å)	35.07 – 3.27 35.07 – 3.27	Depositor EDS
% Data completeness (in resolution range)	95.7 (35.07-3.27) 92.2 (35.07-3.27)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.91 (at 3.25Å)	Xtriage
Refinement program	PHENIX 1.10.1_2155	Depositor
R, R_{free}	0.192 , 0.248 0.209 , 0.257	Depositor DCC
R_{free} test set	2795 reflections (4.59%)	wwPDB-VP
Wilson B-factor (Å ²)	87.3	Xtriage
Anisotropy	0.375	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 54.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.018 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	18219	wwPDB-VP
Average B, all atoms (Å ²)	102.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.80% 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: BMA, NAG, MAN

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.50	1/5436 (0.0%)	0.73	7/7380 (0.1%)
1	B	0.48	0/5436	0.74	3/7380 (0.0%)
2	C	0.49	0/829	0.76	2/1132 (0.2%)
2	E	0.65	2/829 (0.2%)	1.56	17/1132 (1.5%)
3	D	0.47	0/981	0.73	1/1337 (0.1%)
3	F	0.40	0/981	0.73	2/1337 (0.1%)
4	H	0.38	0/946	0.73	0/1287
4	X	0.42	0/934	0.69	1/1271 (0.1%)
5	L	0.37	0/824	0.88	3/1119 (0.3%)
5	Y	0.48	0/841	0.70	1/1141 (0.1%)
All	All	0.48	3/18037 (0.0%)	0.80	37/24516 (0.2%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	93	PRO	N-CD	9.80	1.61	1.47
2	E	54	PRO	N-CD	6.47	1.56	1.47
1	A	407	SER	C-N	5.89	1.47	1.33

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	91	ASP	CB-CA-C	19.86	151.35	111.22
2	E	91	ASP	N-CA-C	-19.35	81.82	109.69
2	E	54	PRO	N-CA-CB	-16.05	86.40	103.25
5	L	19	VAL	CB-CA-C	14.48	130.53	110.84
2	E	54	PRO	CB-CA-C	-12.16	91.50	111.56
5	L	19	VAL	N-CA-C	-12.12	91.03	108.36
2	E	53	ARG	CB-CA-C	11.34	130.48	109.51
2	E	91	ASP	CA-C-N	10.58	131.57	119.83

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	E	91	ASP	C-N-CA	10.58	131.57	119.83
2	E	54	PRO	N-CD-CG	-10.22	87.86	103.20
5	L	20	THR	N-CA-C	9.42	124.24	111.24
2	E	54	PRO	N-CA-C	-9.30	93.32	112.47
1	A	340	LEU	N-CA-C	-8.46	93.85	108.56
1	A	339	SER	CA-C-N	-8.17	109.90	122.16
1	A	339	SER	C-N-CA	-8.17	109.90	122.16
1	B	339	SER	N-CA-C	-7.82	97.52	109.95
2	E	55	SER	N-CA-CB	-7.76	98.35	110.06
3	F	65	VAL	CB-CA-C	7.52	123.63	111.29
1	B	680	PRO	N-CA-C	7.42	119.75	110.70
3	D	65	VAL	N-CA-C	-7.29	99.15	110.09
2	E	54	PRO	CA-N-CD	-7.23	101.87	112.00
3	F	65	VAL	N-CA-C	-7.02	94.74	109.34
2	E	49	GLU	CA-C-N	6.97	137.29	125.02
2	E	49	GLU	C-N-CA	6.97	137.29	125.02
1	A	338	ILE	N-CA-CB	6.87	122.56	111.23
2	E	93	PRO	N-CA-C	-6.74	98.59	112.47
2	E	53	ARG	C-N-CD	-6.60	97.93	125.00
1	A	339	SER	CB-CA-C	-6.43	99.29	111.06
2	C	49	GLU	CA-C-N	6.36	136.21	125.02
2	C	49	GLU	C-N-CA	6.36	136.21	125.02
2	E	92	ASP	N-CA-CB	-6.25	104.45	110.39
1	A	690	LEU	CA-C-N	-5.57	111.67	121.70
1	A	690	LEU	C-N-CA	-5.57	111.67	121.70
1	B	339	SER	N-CA-CB	5.51	119.66	110.85
5	Y	52	SER	N-CA-C	5.50	119.25	112.54
2	E	55	SER	N-CA-C	5.21	118.44	110.20
4	X	32	GLY	N-CA-C	5.15	119.98	111.59

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	5323	0	5345	133	0
1	B	5323	0	5345	118	0
2	C	809	0	755	20	0
2	E	809	0	755	17	0
3	D	956	0	938	25	0
3	F	956	0	938	30	0
4	H	921	0	885	23	0
4	X	909	0	875	30	0
5	L	807	0	772	32	0
5	Y	824	0	796	39	0
6	A	255	0	255	18	0
6	B	255	0	255	15	0
7	A	24	0	24	2	0
7	B	24	0	24	2	0
8	A	12	0	12	0	0
8	B	12	0	12	0	0
All	All	18219	0	17986	476	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 13.

All (476) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:338:ILE:HG12	1:B:338:ILE:O	1.50	1.06
4:X:12:VAL:O	4:X:114:VAL:HG22	1.60	1.01
2:E:53:ARG:NH1	2:E:61:PHE:O	2.03	0.92
4:X:69:SER:HB3	4:X:82:LYS:HB3	1.52	0.91
4:X:28:SER:HA	4:X:77:ASN:HD21	1.36	0.90
1:A:687:PRO:HG2	1:A:690:LEU:HD13	1.54	0.89
1:B:473:ARG:HD3	1:B:499:PRO:HD2	1.55	0.88
1:B:100:MET:HG2	6:B:704:NAG:H4	1.54	0.88
1:B:338:ILE:O	1:B:338:ILE:CG1	2.20	0.87
1:B:230:ASN:HD22	1:B:257:ASN:HD22	1.20	0.86
5:L:86:TYR:HE2	5:L:104:LEU:HD11	1.39	0.86
5:Y:19:VAL:HG11	5:Y:104:LEU:HD11	1.57	0.86
6:A:814:NAG:O4	6:A:815:NAG:H1	1.75	0.86
5:L:18:ARG:HD2	5:L:75:ILE:H	1.37	0.85
5:Y:81:GLU:O	5:Y:83:ILE:N	2.08	0.85
1:A:658:VAL:HG21	1:A:691:PHE:HE2	1.44	0.82
5:L:18:ARG:NH1	5:L:21:ILE:H	1.80	0.80
5:Y:81:GLU:C	5:Y:83:ILE:H	1.90	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:19:ARG:HG2	2:C:73:THR:HG23	1.65	0.79
1:A:100:MET:HG2	6:A:803:NAG:H4	1.64	0.78
1:B:339:SER:HB2	1:B:342:SER:HB3	1.64	0.78
5:Y:80:GLN:HE22	5:Y:83:ILE:HG13	1.47	0.78
1:B:291:ASN:HD22	1:B:316:HIS:HB2	1.47	0.78
4:H:28:SER:HA	4:H:77:ASN:HD21	1.48	0.76
5:Y:80:GLN:OE1	5:Y:106:ILE:HD11	1.84	0.76
5:Y:39:LYS:HG3	5:Y:41:ASP:H	1.51	0.75
6:B:715:NAG:O4	6:B:716:NAG:H1	1.87	0.75
5:L:11:LEU:HD12	5:L:12:PRO:HD2	1.66	0.75
1:B:119:PHE:O	1:B:146:GLN:NE2	2.21	0.73
5:L:85:THR:CG2	5:L:103:LYS:HG3	2.19	0.73
1:A:594:GLY:HA3	1:A:618:GLN:HE21	1.53	0.73
4:X:12:VAL:O	4:X:114:VAL:CG2	2.36	0.73
1:B:227:PHE:HD1	1:B:254:SER:HB3	1.55	0.72
6:B:713:NAG:H83	6:B:713:NAG:H3	1.72	0.71
3:F:17:THR:HG22	3:F:87:SER:HA	1.73	0.71
1:A:473:ARG:HH21	1:A:498:SER:H	1.38	0.71
2:E:48:TYR:CD1	2:E:54:PRO:HD3	2.26	0.70
5:L:85:THR:HG22	5:L:103:LYS:HG3	1.73	0.70
1:B:489:ARG:O	1:B:489:ARG:HD3	1.91	0.70
5:L:66:GLY:HA3	5:L:71:TYR:HA	1.72	0.70
1:A:291:ASN:HD22	1:A:316:HIS:HB2	1.57	0.69
4:H:69:SER:HB3	4:H:82:LYS:HB3	1.73	0.69
5:Y:48:ILE:HG22	5:Y:51:THR:O	1.93	0.69
5:Y:80:GLN:NE2	5:Y:83:ILE:HG13	2.07	0.68
6:A:812:NAG:H83	6:A:812:NAG:H3	1.73	0.68
1:B:29:THR:N	1:B:37:CYS:HG	1.90	0.68
3:F:34:ALA:HA	3:F:103:SER:HA	1.76	0.68
5:Y:39:LYS:HD2	5:Y:40:PRO:HD2	1.76	0.68
1:B:265:ASN:ND2	6:B:708:NAG:H61	2.09	0.68
2:E:1:ASP:HB3	2:E:91:ASP:OD2	1.94	0.67
1:A:40:LEU:HB2	1:A:42:LEU:HG	1.76	0.66
3:F:64:ALA:C	3:F:65:VAL:O	2.31	0.66
1:B:211:GLU:OE2	1:B:213:SER:OG	2.12	0.66
6:A:811:NAG:O7	6:A:811:NAG:O1	2.14	0.66
1:B:351:PHE:HB3	1:B:378:LEU:HD21	1.79	0.65
1:A:169:GLN:HG3	1:A:170:LEU:HG	1.78	0.65
1:A:689:ARG:HG3	1:A:690:LEU:HD12	1.78	0.65
1:A:268:PHE:HB3	1:A:271:LEU:HD12	1.78	0.64
1:A:211:GLU:OE2	1:A:213:SER:OG	2.16	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:L:18:ARG:HD2	5:L:75:ILE:N	2.12	0.64
1:B:37:CYS:HB2	1:B:58:LEU:HD23	1.79	0.63
5:Y:80:GLN:NE2	5:Y:80:GLN:HA	2.13	0.63
4:H:16:GLN:HG2	4:H:17:SER:H	1.63	0.62
1:A:473:ARG:HH21	1:A:498:SER:HB3	1.64	0.62
1:A:489:ARG:O	1:A:489:ARG:HD3	2.00	0.62
5:Y:81:GLU:C	5:Y:83:ILE:N	2.56	0.62
3:D:17:THR:HG22	3:D:87:SER:HA	1.82	0.62
1:A:521:ILE:HD12	1:A:542:LEU:HD13	1.82	0.62
1:A:658:VAL:HG21	1:A:691:PHE:CE2	2.32	0.62
1:A:688:VAL:HG12	1:A:688:VAL:O	2.00	0.62
2:E:91:ASP:O	2:E:94:ASN:O	2.18	0.61
6:B:716:NAG:O3	7:B:717:BMA:H1	2.00	0.61
3:F:53:ILE:HG22	3:F:61:ASN:HB3	1.80	0.61
1:A:393:LEU:O	1:A:419:ILE:HA	2.01	0.61
1:A:450:GLN:OE1	1:A:453:ARG:NH2	2.34	0.60
1:B:169:GLN:HG3	1:B:170:LEU:HG	1.82	0.60
1:B:508:LEU:HD21	1:B:511:LEU:HD13	1.82	0.60
1:B:566:ILE:HD12	1:B:590:ILE:HD12	1.83	0.60
1:A:591:ILE:HG22	1:A:612:LEU:HD11	1.83	0.60
1:A:227:PHE:HD1	1:A:254:SER:HB3	1.66	0.60
1:A:29:THR:N	1:A:37:CYS:HG	1.98	0.60
5:L:18:ARG:HE	5:L:21:ILE:HD12	1.67	0.60
1:B:275:ASN:OD1	6:B:710:NAG:O1	2.19	0.59
5:Y:80:GLN:OE1	5:Y:106:ILE:CD1	2.51	0.59
3:D:36:TRP:CH2	3:D:101:ARG:HD2	2.37	0.59
4:H:52:ILE:HD12	4:H:58:ASN:HB3	1.85	0.59
1:A:37:CYS:HB2	1:A:58:LEU:HD23	1.85	0.58
1:A:510:ILE:HG13	1:A:534:ILE:HB	1.84	0.58
4:X:35:TRP:HB2	4:X:52:ILE:HG23	1.85	0.58
1:A:252:ASN:ND2	6:A:806:NAG:O7	2.35	0.58
4:H:67:ARG:HB2	4:H:84:ASN:HB2	1.86	0.58
1:B:32:HIS:O	1:B:32:HIS:ND1	2.33	0.58
4:H:67:ARG:HA	4:H:84:ASN:HD22	1.68	0.58
2:C:23:SER:OG	2:C:69:THR:HG22	2.03	0.58
5:Y:61:ARG:HH11	5:Y:82:ASP:CG	2.11	0.58
1:A:500:SER:HB2	1:A:525:MET:HA	1.86	0.58
1:A:378:LEU:O	1:A:407:SER:HB3	2.03	0.58
3:D:38:TRP:CE2	3:D:84:LEU:HB2	2.38	0.58
3:F:101:ARG:NH1	3:F:110:ASP:HB2	2.19	0.57
4:X:114:VAL:HG12	4:X:114:VAL:O	2.02	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:94:ASN:HA	3:F:49:TRP:CZ3	2.38	0.57
4:H:35:TRP:HB2	4:H:52:ILE:HG23	1.85	0.57
6:A:810:NAG:O3	6:A:811:NAG:O5	2.23	0.57
1:A:49:LEU:O	1:A:73:ARG:NH2	2.37	0.57
5:Y:80:GLN:O	5:Y:83:ILE:HG12	2.04	0.57
1:A:651:CYS:HA	1:A:655:ALA:HB2	1.86	0.57
1:B:40:LEU:HB2	1:B:42:LEU:HG	1.85	0.57
1:B:598:LEU:H	1:B:620:ASN:HD22	1.51	0.57
1:A:588:LEU:HD21	1:A:591:ILE:HB	1.86	0.57
4:X:91:THR:HG23	4:X:113:THR:HA	1.86	0.57
4:X:18:LEU:HD22	4:X:112:VAL:HG11	1.86	0.56
1:A:336:GLN:HB2	1:A:343:LEU:HD23	1.88	0.56
3:D:13:LYS:O	3:D:16:GLN:HB2	2.05	0.56
5:Y:3:GLN:OE1	5:Y:26:GLY:HA3	2.05	0.56
5:L:37:GLN:HB2	5:L:47:LEU:HD11	1.87	0.56
3:D:79:LYS:HB3	3:D:81:GLN:HG2	1.88	0.56
1:A:347:ASP:O	1:A:350:SER:OG	2.23	0.56
1:B:683:TYR:HE2	1:B:691:PHE:CE2	2.24	0.56
1:B:473:ARG:NH2	1:B:498:SER:HB3	2.20	0.56
5:L:31:ASN:HA	5:L:71:TYR:CE2	2.41	0.56
1:B:688:VAL:O	1:B:690:LEU:N	2.39	0.55
3:F:28:SER:HB3	3:F:31:SER:HB2	1.88	0.55
2:C:107:LYS:HG3	2:C:107:LYS:O	2.05	0.55
3:D:58:LYS:CD	3:D:59:TRP:H	2.19	0.55
5:L:18:ARG:HH11	5:L:21:ILE:H	1.52	0.55
1:A:546:TRP:HB3	1:A:578:PRO:HD3	1.89	0.55
3:F:13:LYS:O	3:F:16:GLN:HB2	2.07	0.55
2:C:100:GLY:O	3:D:45:ARG:NH1	2.38	0.55
1:A:473:ARG:HE	1:A:499:PRO:HD2	1.71	0.55
6:B:712:NAG:O7	6:B:712:NAG:O1	2.20	0.55
1:A:563:HIS:O	1:A:565:HIS:ND1	2.38	0.55
1:B:683:TYR:CE2	1:B:691:PHE:CE2	2.94	0.55
3:F:101:ARG:HH12	3:F:110:ASP:HB2	1.72	0.55
4:X:33:TYR:HD1	4:X:100:TYR:HB2	1.70	0.55
1:B:32:HIS:HD1	1:B:32:HIS:C	2.15	0.54
4:H:28:SER:HA	4:H:77:ASN:ND2	2.19	0.54
6:A:807:NAG:H3	6:A:808:NAG:H1	1.89	0.54
3:F:94:THR:HG23	3:F:119:THR:HA	1.90	0.54
4:H:4:LEU:HD13	4:H:96:CYS:SG	2.47	0.54
1:A:159:LEU:HB2	1:A:181:ASN:HD22	1.72	0.54
1:B:579:VAL:HG13	1:B:602:PRO:HB2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:X:51:TYR:O	4:X:58:ASN:HB2	2.06	0.54
1:A:348:ASP:HB3	1:A:373:ASN:HB2	1.89	0.54
1:A:578:PRO:HB2	1:A:581:VAL:HG13	1.89	0.54
6:A:802:NAG:O7	6:A:802:NAG:O1	2.23	0.54
1:B:76:GLN:HA	1:B:100:MET:HE3	1.90	0.54
3:F:40:ARG:NH1	3:F:93:ASP:OD1	2.41	0.54
5:Y:80:GLN:NE2	5:Y:80:GLN:O	2.40	0.54
3:D:23:ALA:HA	3:D:81:GLN:HB3	1.89	0.54
1:A:167:GLN:NE2	3:D:105:PRO:O	2.42	0.53
4:X:87:THR:O	4:X:114:VAL:HG12	2.09	0.53
4:H:51:TYR:O	4:H:58:ASN:HB2	2.08	0.53
4:X:99:ALA:HB2	4:X:103:PHE:HA	1.91	0.53
3:D:59:TRP:CE2	3:D:75:PRO:HG3	2.43	0.53
5:Y:80:GLN:NE2	5:Y:80:GLN:CA	2.72	0.53
1:A:643:ARG:NH1	1:A:670:GLU:OE1	2.36	0.53
1:A:654:ILE:O	1:A:658:VAL:HG23	2.08	0.53
1:B:473:ARG:HH21	1:B:499:PRO:HD2	1.73	0.53
5:Y:83:ILE:CD1	5:Y:106:ILE:HG23	2.40	0.52
5:Y:91:GLY:HA2	5:Y:96:TRP:CD1	2.44	0.52
1:A:664:THR:OG1	1:A:666:THR:HG23	2.09	0.52
5:L:1:ASP:H1	5:L:95:PRO:HD2	1.74	0.52
1:A:543:ALA:HB2	1:A:572:ASN:O	2.09	0.52
1:B:49:LEU:O	1:B:73:ARG:NH2	2.41	0.52
1:A:594:GLY:HA2	1:A:620:ASN:HD21	1.74	0.52
4:X:21:THR:HG1	4:X:80:PHE:HD1	1.56	0.52
1:A:154:LEU:HB2	1:A:178:LEU:HD23	1.92	0.52
4:H:91:THR:HG23	4:H:113:THR:HA	1.91	0.52
1:A:265:ASN:ND2	6:A:807:NAG:H61	2.25	0.52
2:C:94:ASN:HA	3:D:49:TRP:CZ3	2.45	0.52
1:B:465:TYR:HA	1:B:489:ARG:HD2	1.91	0.52
3:D:94:THR:HG23	3:D:119:THR:HA	1.92	0.52
1:B:521:ILE:HD11	1:B:542:LEU:HD22	1.92	0.52
4:X:37:TRP:CZ3	4:X:96:CYS:HB2	2.45	0.52
1:B:266:THR:O	1:B:269:LEU:HB2	2.10	0.52
3:D:36:TRP:CZ3	3:D:101:ARG:HB2	2.45	0.51
2:E:6:GLN:HB2	2:E:7:PRO:HD2	1.92	0.51
1:A:119:PHE:O	1:A:146:GLN:NE2	2.42	0.51
1:A:357:LEU:HD21	1:A:360:LEU:HB2	1.92	0.51
1:A:611:SER:O	1:A:613:LYS:HG3	2.10	0.51
1:B:99:PRO:O	1:B:124:ASN:ND2	2.44	0.51
1:B:623:THR:HG22	1:B:646:PRO:O	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
5:L:85:THR:HG22	5:L:103:LYS:CG	2.40	0.51
5:L:8:THR:HG21	5:L:20:THR:O	2.10	0.51
5:L:86:TYR:CE2	5:L:104:LEU:HD11	2.31	0.51
1:B:337:SER:C	1:B:338:ILE:HG23	2.36	0.51
5:L:84:ALA:H	5:L:104:LEU:HD12	1.74	0.51
1:A:595:LEU:HD12	1:A:619:LYS:HB2	1.93	0.51
1:A:473:ARG:NH2	1:A:498:SER:HB3	2.24	0.51
1:B:291:ASN:ND2	6:B:711:NAG:H61	2.26	0.51
1:B:229:ASN:C	1:B:231:VAL:H	2.19	0.51
5:Y:3:GLN:O	5:Y:25:ALA:HA	2.10	0.51
1:A:76:GLN:HA	1:A:100:MET:HE3	1.92	0.51
1:A:398:ASN:CG	6:A:812:NAG:H61	2.35	0.51
1:A:464:SER:HB3	1:A:488:ARG:HB3	1.93	0.51
3:D:58:LYS:HD3	3:D:59:TRP:H	1.75	0.51
6:B:719:NAG:O6	7:B:720:BMA:H3	2.11	0.51
5:Y:83:ILE:HD12	5:Y:106:ILE:HG23	1.92	0.51
4:X:39:ARG:HB3	4:X:94:TYR:CE1	2.46	0.51
5:L:85:THR:HG23	5:L:103:LYS:HG3	1.90	0.51
1:A:568:ASN:HA	1:A:592:ASP:HB3	1.93	0.50
1:B:469:LEU:HD12	1:B:490:VAL:HG11	1.93	0.50
4:X:16:GLN:HG2	4:X:17:SER:H	1.76	0.50
1:A:167:GLN:HE22	3:D:105:PRO:HA	1.76	0.50
1:B:664:THR:OG1	1:B:666:THR:HG23	2.11	0.50
4:X:49:MET:SD	4:X:94:TYR:HE1	2.34	0.50
1:B:526:LEU:HD22	1:B:529:LEU:HD12	1.92	0.50
1:B:602:PRO:HG2	1:B:605:VAL:HB	1.93	0.50
3:F:30:SER:O	3:F:55:LYS:HE3	2.12	0.50
1:A:361:ASN:OD1	1:A:363:GLU:HG2	2.11	0.50
2:C:58:PRO:HG2	2:C:61:PHE:CE1	2.47	0.50
1:B:336:GLN:HB2	1:B:343:LEU:HD23	1.93	0.50
4:H:99:ALA:HB2	4:H:103:PHE:HA	1.93	0.50
1:B:361:ASN:OD1	1:B:363:GLU:HG2	2.11	0.49
1:A:272:LYS:HD3	1:A:296:TRP:HE3	1.77	0.49
1:B:630:PHE:CE1	1:B:642:MET:HE1	2.47	0.49
1:B:591:ILE:HG22	1:B:612:LEU:HD11	1.92	0.49
1:B:678:ASN:OD1	1:B:684:HIS:HE1	1.96	0.49
2:E:107:LYS:N	2:E:107:LYS:CD	2.75	0.49
3:D:53:ILE:HG22	3:D:61:ASN:HB3	1.95	0.49
1:B:339:SER:HB2	1:B:342:SER:CB	2.36	0.49
1:A:465:TYR:HD1	1:A:489:ARG:HD2	1.77	0.49
1:B:354:LEU:HB3	1:B:357:LEU:HB2	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:X:5:LYS:NZ	4:X:108:GLN:HB2	2.28	0.49
1:B:680:PRO:HG2	1:B:683:TYR:HB2	1.95	0.49
1:A:617:LEU:HD13	1:A:642:MET:HE3	1.95	0.49
2:E:38:LYS:HB3	2:E:39:PRO:HD2	1.94	0.49
1:A:618:GLN:O	1:A:643:ARG:O	2.31	0.49
1:B:130:LEU:HB2	1:B:154:LEU:HD23	1.94	0.49
3:D:52:ILE:HG22	3:D:102:TYR:CZ	2.48	0.48
1:B:473:ARG:HH22	1:B:498:SER:HB3	1.78	0.48
5:Y:18:ARG:HA	5:Y:75:ILE:O	2.13	0.48
1:A:64:ARG:HG2	1:A:86:THR:HB	1.93	0.48
1:A:688:VAL:O	1:A:688:VAL:CG1	2.61	0.48
2:C:91:ASP:OD2	2:C:92:ASP:N	2.46	0.48
6:B:703:NAG:O7	6:B:703:NAG:O1	2.24	0.48
4:X:88:THR:HA	4:X:114:VAL:O	2.13	0.48
1:A:676:LEU:HD23	1:A:685:GLY:HA2	1.95	0.48
2:C:38:LYS:HG2	2:C:83:ALA:HB2	1.95	0.48
1:B:393:LEU:O	1:B:419:ILE:HA	2.14	0.48
4:H:67:ARG:HD2	4:H:84:ASN:O	2.13	0.48
1:A:179:SER:HB3	1:A:203:GLU:HG3	1.96	0.48
1:A:275:ASN:OD1	6:A:809:NAG:O1	2.30	0.48
4:H:13:ARG:NH1	4:H:116:SER:O	2.46	0.48
1:B:421:LYS:HE3	1:B:423:GLU:HG2	1.96	0.48
4:H:18:LEU:HD22	4:H:112:VAL:HG11	1.95	0.48
1:A:316:HIS:HA	1:A:353:TRP:CH2	2.48	0.48
1:B:473:ARG:NH2	1:B:498:SER:H	2.11	0.48
2:E:4:LEU:HD12	2:E:22:CYS:SG	2.54	0.48
1:B:122:CYS:HB3	1:B:125:LEU:HG	1.94	0.48
4:H:64:LEU:HD13	4:H:68:ILE:HD12	1.95	0.48
1:B:460:GLU:OE1	1:B:484:ARG:HD2	2.13	0.48
2:E:32:VAL:H	2:E:50:ASP:HB3	1.79	0.48
4:H:62:PRO:HD2	5:L:95:PRO:HG3	1.96	0.48
1:A:391:THR:OG1	1:A:392:SER:N	2.47	0.48
1:A:130:LEU:HB2	1:A:154:LEU:HD23	1.96	0.47
1:A:413:ASN:CG	6:A:814:NAG:O1	2.57	0.47
2:C:10:VAL:HG13	2:C:104:LEU:HD13	1.95	0.47
1:B:543:ALA:HB2	1:B:572:ASN:O	2.14	0.47
4:X:22:CYS:HB2	4:X:37:TRP:CH2	2.49	0.47
4:X:28:SER:OG	4:X:30:THR:HG22	2.15	0.47
1:B:413:ASN:HD22	1:B:437:ASP:HB3	1.78	0.47
2:E:92:ASP:HB3	2:E:93:PRO:HD3	1.96	0.47
1:B:268:PHE:HB3	1:B:271:LEU:HD12	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:715:NAG:O4	6:B:716:NAG:N2	2.48	0.47
3:F:67:VAL:O	3:F:71:ILE:HG22	2.13	0.47
1:B:578:PRO:HB2	1:B:581:VAL:HG13	1.96	0.47
1:B:580:GLU:O	1:B:583:LYS:HB2	2.15	0.47
1:A:34:VAL:HG13	1:A:55:VAL:HB	1.97	0.47
1:A:146:GLN:HB3	1:A:149:LEU:HB2	1.97	0.47
1:A:387:SER:HB2	1:A:413:ASN:OD1	2.13	0.47
1:A:444:GLY:HA2	1:A:468:TYR:O	2.15	0.47
6:A:815:NAG:O3	7:A:816:BMA:H1	2.14	0.47
5:Y:36:TYR:O	5:Y:86:TYR:HA	2.14	0.47
5:L:55:HIS:HB3	5:L:58:VAL:HG23	1.96	0.47
1:B:64:ARG:HG2	1:B:86:THR:HB	1.95	0.47
1:A:610:VAL:HA	1:A:635:ARG:HH21	1.80	0.46
3:D:67:VAL:O	3:D:71:ILE:HG22	2.15	0.46
1:B:373:ASN:HA	1:B:403:SER:HB3	1.96	0.46
2:E:90:TYR:HB2	2:E:95:PHE:HA	1.96	0.46
5:L:83:ILE:CG2	5:L:104:LEU:O	2.63	0.46
1:A:239:GLU:O	1:A:243:LEU:HG	2.14	0.46
1:A:279:LEU:HD11	1:A:281:LEU:HD21	1.98	0.46
2:C:107:LYS:HB2	2:C:107:LYS:HE3	1.40	0.46
1:B:497:SER:OG	1:B:499:PRO:O	2.30	0.46
1:B:680:PRO:O	1:B:681:PRO:C	2.58	0.46
5:Y:86:TYR:HE2	5:Y:104:LEU:HD22	1.79	0.46
1:A:126:THR:HA	1:A:149:LEU:HA	1.98	0.46
1:A:398:ASN:ND2	6:A:812:NAG:H61	2.30	0.46
1:B:398:ASN:ND2	6:B:713:NAG:H61	2.31	0.46
1:B:546:TRP:HB3	1:B:578:PRO:HD3	1.97	0.46
4:X:36:ASN:OD1	4:X:51:TYR:HB3	2.16	0.46
5:Y:80:GLN:O	5:Y:83:ILE:CG1	2.63	0.46
1:A:105:ASN:HA	1:A:129:HIS:HB2	1.97	0.46
2:E:88:SER:HA	2:E:97:VAL:O	2.16	0.46
1:A:301:GLU:HA	1:A:324:VAL:HA	1.97	0.46
2:E:107:LYS:N	2:E:107:LYS:HD3	2.31	0.46
1:A:521:ILE:H	1:A:545:LEU:HD11	1.81	0.46
3:F:13:LYS:HB2	3:F:16:GLN:HG3	1.98	0.46
4:X:52:ILE:HA	4:X:58:ASN:HB3	1.98	0.46
1:A:421:LYS:HE3	1:A:423:GLU:HG2	1.96	0.46
1:A:602:PRO:HG2	1:A:605:VAL:HB	1.97	0.46
1:B:337:SER:O	1:B:338:ILE:HG23	2.16	0.46
2:C:58:PRO:HG2	2:C:61:PHE:CD1	2.51	0.46
3:F:18:LEU:O	3:F:85:GLN:HA	2.16	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:617:LEU:HD13	1:A:642:MET:CE	2.46	0.46
3:F:49:TRP:HZ2	3:F:52:ILE:HG23	1.81	0.46
3:D:18:LEU:O	3:D:85:GLN:HA	2.15	0.45
3:D:49:TRP:HZ2	3:D:52:ILE:HG23	1.80	0.45
1:B:593:LEU:HD12	1:B:617:LEU:HG	1.98	0.45
1:A:609:GLN:HB3	1:A:612:LEU:HB2	1.98	0.45
4:X:91:THR:OG1	4:X:114:VAL:O	2.33	0.45
1:A:176:LEU:C	1:A:177:LEU:HD12	2.42	0.45
1:B:32:HIS:ND1	1:B:32:HIS:C	2.75	0.45
1:B:386:LEU:HA	1:B:389:SER:OG	2.16	0.45
3:F:58:LYS:HD2	3:F:58:LYS:HA	1.58	0.45
4:H:32:GLY:O	4:H:54:TYR:HB3	2.16	0.45
1:A:316:HIS:CD2	1:A:353:TRP:CZ2	3.05	0.45
1:B:34:VAL:HG13	1:B:55:VAL:HB	1.98	0.45
1:B:465:TYR:HA	1:B:489:ARG:CD	2.46	0.45
3:D:34:ALA:HA	3:D:103:SER:HA	1.97	0.45
3:F:4:LEU:HB3	3:F:22:CYS:SG	2.56	0.45
3:F:72:THR:HB	3:F:85:GLN:HB2	1.98	0.45
5:L:16:GLY:HA2	5:L:77:ASN:HA	1.99	0.45
1:B:594:GLY:HA3	1:B:618:GLN:HE21	1.82	0.45
3:F:49:TRP:CE3	3:F:64:ALA:HB2	2.51	0.45
4:X:101:ILE:HG12	5:Y:49:TYR:CD2	2.52	0.45
5:L:6:GLN:C	5:L:8:THR:H	2.23	0.45
1:A:318:LEU:HB3	1:A:321:LEU:HD12	1.98	0.45
1:A:469:LEU:HD23	1:A:469:LEU:HA	1.70	0.45
1:B:105:ASN:HA	1:B:129:HIS:HB2	1.99	0.45
3:D:38:TRP:CZ3	3:D:99:CYS:HB3	2.52	0.45
5:Y:35:TRP:CZ3	5:Y:88:CYS:HB3	2.51	0.45
4:X:87:THR:O	4:X:114:VAL:CG1	2.64	0.45
4:H:51:TYR:CE2	4:H:59:ASN:HB3	2.52	0.45
5:L:31:ASN:HA	5:L:71:TYR:HE2	1.82	0.45
5:L:83:ILE:HG23	5:L:104:LEU:HB2	1.99	0.45
5:L:89:GLN:HB2	5:L:98:PHE:CD2	2.52	0.44
2:E:18:ALA:HB3	2:E:74:ILE:HB	2.00	0.44
1:A:508:LEU:HD21	1:A:511:LEU:HD13	1.99	0.44
1:B:92:PRO:HG3	1:B:117:LYS:HB3	2.00	0.44
3:F:38:TRP:CE2	3:F:84:LEU:HB2	2.52	0.44
1:A:298:PRO:O	1:A:323:ASN:HB2	2.17	0.44
2:C:22:CYS:N	2:C:70:ALA:O	2.50	0.44
1:B:589:LYS:O	1:B:612:LEU:HD12	2.18	0.44
4:H:83:LEU:HD23	4:H:86:VAL:HG12	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:277:THR:HA	1:A:300:LEU:HA	1.99	0.44
1:A:465:TYR:HA	1:A:489:ARG:HD2	2.00	0.44
1:B:610:VAL:HA	1:B:635:ARG:HH21	1.82	0.44
3:F:64:ALA:O	3:F:65:VAL:C	2.57	0.44
1:A:153:ASP:HA	1:A:177:LEU:HB2	2.00	0.44
2:C:82:GLU:HG3	2:C:104:LEU:O	2.18	0.44
3:F:36:TRP:CH2	3:F:101:ARG:HD2	2.53	0.44
2:C:38:LYS:HB3	2:C:39:PRO:HD2	1.98	0.44
3:D:63:TYR:HB2	3:D:68:LYS:HD2	1.99	0.44
1:B:301:GLU:HA	1:B:324:VAL:HA	1.99	0.44
5:Y:48:ILE:CG2	5:Y:51:THR:O	2.63	0.44
4:H:28:SER:OG	4:H:30:THR:HG22	2.18	0.44
5:L:62:PHE:HB2	5:L:74:THR:O	2.18	0.44
6:A:818:NAG:O6	7:A:819:BMA:H3	2.18	0.43
2:C:6:GLN:HB2	2:C:7:PRO:HD2	2.00	0.43
3:F:100:ALA:HB1	3:F:109:ILE:CG2	2.48	0.43
6:A:802:NAG:HO1	6:A:802:NAG:C7	2.29	0.43
4:H:27:TYR:HE2	4:H:33:TYR:HD2	1.66	0.43
1:A:177:LEU:HD12	1:A:177:LEU:N	2.33	0.43
1:A:498:SER:OG	1:A:499:PRO:HD3	2.18	0.43
1:A:671:LEU:HA	1:A:675:TYR:CD1	2.53	0.43
1:B:316:HIS:HD2	1:B:353:TRP:CZ2	2.36	0.43
1:B:378:LEU:O	1:B:407:SER:HB3	2.17	0.43
4:X:101:ILE:HG12	5:Y:49:TYR:HD2	1.83	0.43
5:Y:33:LEU:HD22	5:Y:71:TYR:CB	2.48	0.43
1:A:40:LEU:HD12	1:A:42:LEU:HD11	1.99	0.43
1:A:257:ASN:HA	1:A:283:TYR:O	2.18	0.43
5:Y:15:LEU:HA	5:Y:15:LEU:HD12	1.82	0.43
5:L:7:SER:HB2	5:L:22:SER:OG	2.18	0.43
1:A:316:HIS:HD2	1:A:353:TRP:CZ2	2.36	0.43
1:A:478:LEU:HA	1:A:478:LEU:HD23	1.63	0.43
1:B:455:LEU:HB3	1:B:458:ILE:HB	2.01	0.43
1:B:516:ASN:HB2	1:B:518:ILE:HG13	2.01	0.43
1:A:654:ILE:HD12	1:A:654:ILE:HA	1.79	0.43
1:A:672:SER:HB3	1:A:689:ARG:NH1	2.33	0.43
1:B:469:LEU:HD23	1:B:469:LEU:HA	1.71	0.43
2:C:43:PRO:HD2	3:D:112:TRP:CE3	2.54	0.43
1:B:391:THR:OG1	1:B:392:SER:N	2.51	0.43
1:B:510:ILE:HG13	1:B:534:ILE:HB	2.00	0.43
3:F:49:TRP:CZ2	3:F:52:ILE:HG23	2.53	0.43
4:X:19:SER:HA	4:X:81:LEU:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:440:LEU:HD23	1:B:465:TYR:CB	2.49	0.43
2:E:7:PRO:O	2:E:102:THR:OG1	2.30	0.43
1:A:526:LEU:HD22	1:A:529:LEU:HD12	2.01	0.43
3:F:12:VAL:O	3:F:120:VAL:HA	2.18	0.43
1:A:630:PHE:HB3	1:A:634:PHE:CE1	2.54	0.42
1:A:141:ASN:HB3	1:A:144:VAL:HB	2.00	0.42
1:B:66:LEU:HB2	1:B:94:LEU:HD22	1.99	0.42
1:B:147:LYS:O	1:B:172:ASN:ND2	2.49	0.42
1:B:398:ASN:CG	6:B:713:NAG:H61	2.44	0.42
3:F:103:SER:OG	3:F:110:ASP:OD2	2.17	0.42
5:Y:6:GLN:HE21	5:Y:6:GLN:HB3	1.60	0.42
1:A:248:THR:C	1:A:250:ILE:H	2.27	0.42
1:A:337:SER:C	1:A:338:ILE:HG13	2.42	0.42
1:B:277:THR:HA	1:B:300:LEU:HA	2.01	0.42
1:B:49:LEU:HB2	1:B:74:TYR:CE1	2.54	0.42
1:B:229:ASN:C	1:B:231:VAL:N	2.75	0.42
1:B:233:LEU:HD23	1:B:233:LEU:HA	1.82	0.42
5:L:18:ARG:HH12	5:L:21:ILE:H	1.63	0.42
1:A:558:LEU:HA	1:A:561:LEU:HD12	2.01	0.42
2:C:32:VAL:H	2:C:50:ASP:HB3	1.84	0.42
1:B:675:TYR:C	1:B:676:LEU:HD12	2.45	0.42
1:A:393:LEU:HD12	1:A:393:LEU:HA	1.69	0.42
1:B:588:LEU:HD21	1:B:591:ILE:HB	2.02	0.42
5:Y:2:ILE:HB	5:Y:90:GLN:HE21	1.84	0.42
5:L:33:LEU:HD11	5:L:88:CYS:HB2	2.02	0.42
1:A:117:LYS:HD2	2:C:28:GLY:O	2.19	0.42
1:A:331:ARG:HG3	1:A:364:ASP:HB3	2.00	0.42
1:B:617:LEU:N	1:B:617:LEU:HD12	2.35	0.42
3:F:38:TRP:CG	3:F:84:LEU:HD22	2.54	0.42
1:A:354:LEU:HB3	1:A:357:LEU:HB2	2.02	0.42
1:B:126:THR:HA	1:B:149:LEU:HA	2.01	0.42
1:B:146:GLN:HB3	1:B:149:LEU:HB2	2.02	0.42
5:Y:4:LEU:HD22	5:Y:23:CYS:SG	2.60	0.42
1:B:680:PRO:HA	1:B:681:PRO:HD3	2.00	0.42
1:A:402:VAL:HG23	1:A:429:TRP:CZ2	2.55	0.41
1:A:465:TYR:HA	1:A:489:ARG:CD	2.50	0.41
1:B:508:LEU:HD12	1:B:508:LEU:HA	1.90	0.41
6:B:714:NAG:O7	6:B:714:NAG:O1	2.35	0.41
1:A:139:LYS:HA	1:A:139:LYS:HD2	1.88	0.41
1:A:212:PHE:CE2	1:A:241:LEU:HB2	2.55	0.41
5:Y:80:GLN:HE21	5:Y:80:GLN:C	2.28	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:A:817:NAG:H4	6:A:818:NAG:HN2	1.85	0.41
2:C:2:ILE:HD13	2:C:30:TYR:CE1	2.55	0.41
1:A:38:SER:HB3	1:A:57:ASN:OD1	2.20	0.41
1:A:336:GLN:HB2	1:A:343:LEU:CD2	2.50	0.41
1:B:550:ASN:HA	1:B:551:PRO:HD3	1.92	0.41
5:Y:39:LYS:HA	5:Y:84:ALA:HB1	2.03	0.41
1:A:441:ASN:O	1:A:466:ASN:HA	2.20	0.41
5:Y:86:TYR:CE2	5:Y:104:LEU:HD22	2.54	0.41
1:A:201:LYS:HG3	1:A:227:PHE:CE2	2.56	0.41
1:A:331:ARG:CG	1:A:364:ASP:HB3	2.51	0.41
5:Y:22:SER:HA	5:Y:72:SER:HA	2.03	0.41
1:A:229:ASN:C	1:A:231:VAL:H	2.27	0.41
6:A:807:NAG:C3	6:A:808:NAG:H1	2.49	0.41
4:H:32:GLY:C	4:H:54:TYR:HB3	2.46	0.41
1:A:164:LEU:HD12	1:A:194:PHE:HZ	1.85	0.41
1:A:329:LEU:HA	1:A:332:SER:OG	2.21	0.41
1:B:680:PRO:HG2	1:B:683:TYR:CB	2.50	0.41
4:X:11:LEU:HD23	4:X:113:THR:OG1	2.21	0.41
5:L:62:PHE:HB3	5:L:75:ILE:HA	2.02	0.41
1:A:92:PRO:HG3	1:A:117:LYS:HB3	2.03	0.41
1:A:566:ILE:HD12	1:A:590:ILE:HD12	2.03	0.41
2:C:88:SER:HA	2:C:97:VAL:O	2.21	0.41
1:B:200:LYS:HA	1:B:224:PHE:HB2	2.03	0.41
1:B:281:LEU:C	1:B:284:ASN:HD22	2.29	0.41
1:B:513:LEU:HB2	1:B:537:LEU:HD23	2.02	0.41
3:F:101:ARG:NH1	3:F:111:TYR:HD2	2.19	0.41
5:L:86:TYR:O	5:L:101:GLY:HA2	2.20	0.41
1:A:468:TYR:CE2	1:A:470:GLN:HB2	2.56	0.41
3:D:39:ILE:HG23	3:D:48:GLU:O	2.20	0.41
1:B:387:SER:C	1:B:389:SER:N	2.79	0.41
1:B:473:ARG:HD3	1:B:499:PRO:CD	2.38	0.41
5:Y:38:GLN:HA	5:Y:43:THR:O	2.20	0.41
1:A:159:LEU:HB2	1:A:181:ASN:ND2	2.36	0.40
1:A:488:ARG:HG3	1:A:514:SER:OG	2.21	0.40
1:B:227:PHE:CD1	1:B:254:SER:HB3	2.46	0.40
1:B:595:LEU:HD12	1:B:619:LYS:HB2	2.03	0.40
3:F:38:TRP:CZ3	3:F:99:CYS:HB3	2.56	0.40
1:A:95:CYS:HB2	1:A:121:PHE:HB2	2.03	0.40
1:A:537:LEU:HD23	1:A:537:LEU:HA	1.85	0.40
1:B:471:LEU:HD23	1:B:471:LEU:HA	1.92	0.40
1:B:660:TRP:CZ2	1:B:664:THR:HG21	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:107:LYS:HA	2:E:107:LYS:HD2	1.90	0.40
4:X:33:TYR:CD1	4:X:100:TYR:HB2	2.53	0.40
1:A:460:GLU:OE1	1:A:484:ARG:HD2	2.20	0.40
1:B:242:CYS:HB3	1:B:271:LEU:HG	2.03	0.40
1:B:413:ASN:CG	6:B:715:NAG:O1	2.64	0.40
1:B:468:TYR:HA	1:B:490:VAL:O	2.21	0.40
1:B:654:ILE:HD12	1:B:654:ILE:HA	1.89	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	661/671 (98%)	611 (92%)	50 (8%)	0	100	100
1	B	661/671 (98%)	613 (93%)	48 (7%)	0	100	100
2	C	105/107 (98%)	98 (93%)	7 (7%)	0	100	100
2	E	105/107 (98%)	99 (94%)	6 (6%)	0	100	100
3	D	120/127 (94%)	110 (92%)	10 (8%)	0	100	100
3	F	120/127 (94%)	109 (91%)	11 (9%)	0	100	100
4	H	114/121 (94%)	103 (90%)	11 (10%)	0	100	100
4	X	112/121 (93%)	102 (91%)	10 (9%)	0	100	100
5	L	103/107 (96%)	88 (85%)	15 (15%)	0	100	100
5	Y	105/107 (98%)	93 (89%)	10 (10%)	2 (2%)	6	29
All	All	2206/2266 (97%)	2026 (92%)	178 (8%)	2 (0%)	48	76

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
5	Y	82	ASP
5	Y	106	ILE

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	618/626 (99%)	606 (98%)	12 (2%)	50	69
1	B	618/626 (99%)	601 (97%)	17 (3%)	38	62
2	C	90/90 (100%)	83 (92%)	7 (8%)	11	37
2	E	90/90 (100%)	84 (93%)	6 (7%)	15	42
3	D	107/108 (99%)	104 (97%)	3 (3%)	38	62
3	F	107/108 (99%)	103 (96%)	4 (4%)	30	57
4	H	101/102 (99%)	98 (97%)	3 (3%)	36	61
4	X	99/102 (97%)	93 (94%)	6 (6%)	17	45
5	L	91/93 (98%)	87 (96%)	4 (4%)	25	54
5	Y	93/93 (100%)	90 (97%)	3 (3%)	34	60
All	All	2014/2038 (99%)	1949 (97%)	65 (3%)	34	60

All (65) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	29	THR
1	A	32	HIS
1	A	91	GLU
1	A	139	LYS
1	A	197	SER
1	A	338	ILE
1	A	340	LEU
1	A	542	LEU
1	A	572	ASN
1	A	656	TRP
1	A	679	THR
1	A	686	PHE

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Mol	Chain	Res	Type
2	C	2	ILE
2	C	4	LEU
2	C	5	THR
2	C	52	GLU
2	C	53	ARG
2	C	94	ASN
2	C	107	LYS
3	D	45	ARG
3	D	55	LYS
3	D	121	SER
1	B	91	GLU
1	B	209	ILE
1	B	269	LEU
1	B	323	ASN
1	B	338	ILE
1	B	339	SER
1	B	340	LEU
1	B	393	LEU
1	B	492	LEU
1	B	536	ASP
1	B	542	LEU
1	B	618	GLN
1	B	651	CYS
1	B	656	TRP
1	B	678	ASN
1	B	683	TYR
1	B	689	ARG
2	E	5	THR
2	E	52	GLU
2	E	53	ARG
2	E	54	PRO
2	E	94	ASN
2	E	107	LYS
3	F	45	ARG
3	F	55	LYS
3	F	58	LYS
3	F	61	ASN
4	X	1	GLN
4	X	2	VAL
4	X	4	LEU
4	X	81	LEU
4	X	113	THR

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Mol	Chain	Res	Type
4	X	114	VAL
5	Y	41	ASP
5	Y	80	GLN
5	Y	107	LYS
4	H	1	GLN
4	H	4	LEU
4	H	81	LEU
5	L	17	ASP
5	L	19	VAL
5	L	41	ASP
5	L	104	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (40) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	GLN
1	A	60	HIS
1	A	230	ASN
1	A	232	GLN
1	A	257	ASN
1	A	265	ASN
1	A	275	ASN
1	A	398	ASN
1	A	457	ASN
1	A	515	ASN
1	A	520	ASN
1	A	541	ASN
1	A	550	ASN
1	A	597	ASN
1	A	616	ASN
1	A	662	ASN
1	A	684	HIS
3	D	3	GLN
1	B	156	HIS
1	B	169	GLN
1	B	174	GLN
1	B	257	ASN
1	B	265	ASN
1	B	287	ASN
1	B	388	ASN
1	B	398	ASN
1	B	413	ASN

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Mol	Chain	Res	Type
1	B	417	ASN
1	B	515	ASN
1	B	539	HIS
1	B	597	ASN
1	B	684	HIS
2	E	78	GLN
3	F	3	GLN
3	F	54	GLN
4	X	1	GLN
4	X	77	ASN
5	Y	92	ASN
4	H	77	ASN
5	L	77	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

40 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	$\# Z > 2$
6	NAG	A	814	-	15,15,15	0.39	0	21,21,21	0.85	1 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	A	808	-	15,15,15	0.85	1 (6%)	21,21,21	0.89	2 (9%)
6	NAG	B	713	-	15,15,15	1.14	2 (13%)	21,21,21	1.23	2 (9%)
6	NAG	A	802	-	15,15,15	0.32	0	21,21,21	0.62	0
6	NAG	B	719	-	15,15,15	0.71	1 (6%)	21,21,21	1.11	3 (14%)
6	NAG	A	817	-	15,15,15	0.54	0	21,21,21	0.65	0
6	NAG	B	709	-	15,15,15	1.01	1 (6%)	21,21,21	0.96	1 (4%)
6	NAG	B	703	-	15,15,15	0.54	0	21,21,21	0.66	0
6	NAG	A	805	-	15,15,15	0.28	0	21,21,21	0.40	0
6	NAG	A	813	-	15,15,15	0.88	1 (6%)	21,21,21	0.56	0
6	NAG	A	818	-	15,15,15	0.71	1 (6%)	21,21,21	1.10	2 (9%)
6	NAG	B	704	-	15,15,15	0.37	0	21,21,21	0.52	0
6	NAG	B	708	-	15,15,15	0.74	1 (6%)	21,21,21	0.57	0
8	MAN	B	701	-	12,12,12	1.10	0	17,17,17	1.16	1 (5%)
8	MAN	A	820	-	12,12,12	1.10	1 (8%)	17,17,17	1.16	1 (5%)
6	NAG	A	809	-	15,15,15	0.25	0	21,21,21	0.37	0
6	NAG	B	711	-	15,15,15	0.23	0	21,21,21	0.68	0
7	BMA	B	717	-	12,12,12	1.13	1 (8%)	17,17,17	0.78	0
6	NAG	A	811	-	15,15,15	0.39	0	21,21,21	0.89	1 (4%)
6	NAG	A	804	-	15,15,15	0.59	0	21,21,21	0.92	1 (4%)
6	NAG	A	803	-	15,15,15	0.41	0	21,21,21	0.52	0
7	BMA	B	720	-	12,12,12	0.96	0	17,17,17	0.77	0
6	NAG	A	806	-	15,15,15	1.44	2 (13%)	21,21,21	1.32	3 (14%)
6	NAG	B	706	-	15,15,15	0.43	0	21,21,21	0.48	0
6	NAG	B	705	-	15,15,15	0.58	0	21,21,21	0.91	1 (4%)
6	NAG	A	815	-	15,15,15	0.48	0	21,21,21	0.73	0
6	NAG	B	702	-	15,15,15	0.74	1 (6%)	21,21,21	0.40	0
6	NAG	B	707	-	15,15,15	1.26	2 (13%)	21,21,21	1.18	2 (9%)
6	NAG	B	718	-	15,15,15	0.52	0	21,21,21	0.58	0
6	NAG	B	712	-	15,15,15	0.49	0	21,21,21	1.20	2 (9%)
6	NAG	B	716	-	15,15,15	0.50	0	21,21,21	0.84	0
6	NAG	A	801	-	15,15,15	0.78	1 (6%)	21,21,21	0.37	0
7	BMA	A	819	-	12,12,12	1.17	1 (8%)	17,17,17	0.69	0
6	NAG	A	812	-	15,15,15	0.93	1 (6%)	21,21,21	1.17	2 (9%)
7	BMA	A	816	-	12,12,12	1.20	0	17,17,17	0.72	0
6	NAG	A	810	-	15,15,15	0.51	0	21,21,21	0.45	0
6	NAG	A	807	-	15,15,15	0.90	1 (6%)	21,21,21	0.71	0
6	NAG	B	715	-	15,15,15	0.46	0	21,21,21	0.85	1 (4%)
6	NAG	B	714	-	15,15,15	0.71	1 (6%)	21,21,21	0.70	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	NAG	B	710	-	15,15,15	0.30	0	21,21,21	0.34	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	A	814	-	-	0/6/26/26	0/1/1/1
6	NAG	A	808	-	-	4/6/26/26	0/1/1/1
6	NAG	B	713	-	-	4/6/26/26	0/1/1/1
6	NAG	A	802	-	-	4/6/26/26	0/1/1/1
6	NAG	B	719	-	-	4/6/26/26	0/1/1/1
6	NAG	A	817	-	-	1/6/26/26	0/1/1/1
6	NAG	B	709	-	-	4/6/26/26	0/1/1/1
6	NAG	B	703	-	-	2/6/26/26	0/1/1/1
6	NAG	A	805	-	-	2/6/26/26	0/1/1/1
6	NAG	A	813	-	-	2/6/26/26	0/1/1/1
6	NAG	A	818	-	-	4/6/26/26	0/1/1/1
6	NAG	B	704	-	-	3/6/26/26	0/1/1/1
6	NAG	B	708	-	-	1/6/26/26	0/1/1/1
8	MAN	B	701	-	-	2/2/22/22	0/1/1/1
8	MAN	A	820	-	-	2/2/22/22	0/1/1/1
6	NAG	A	809	-	-	2/6/26/26	0/1/1/1
6	NAG	B	711	-	-	1/6/26/26	0/1/1/1
7	BMA	B	717	-	-	2/2/22/22	0/1/1/1
6	NAG	A	811	-	-	3/6/26/26	0/1/1/1
6	NAG	A	804	-	-	2/6/26/26	0/1/1/1
6	NAG	A	803	-	-	2/6/26/26	0/1/1/1
7	BMA	B	720	-	-	1/2/22/22	0/1/1/1
6	NAG	A	806	-	-	3/6/26/26	0/1/1/1
6	NAG	B	706	-	-	2/6/26/26	0/1/1/1
6	NAG	B	705	-	-	2/6/26/26	0/1/1/1
6	NAG	A	815	-	-	2/6/26/26	0/1/1/1
6	NAG	B	702	-	-	1/6/26/26	0/1/1/1
6	NAG	B	707	-	-	2/6/26/26	0/1/1/1
6	NAG	B	718	-	-	2/6/26/26	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	NAG	B	712	-	-	3/6/26/26	0/1/1/1
6	NAG	B	716	-	-	2/6/26/26	0/1/1/1
6	NAG	A	801	-	-	1/6/26/26	0/1/1/1
7	BMA	A	819	-	-	1/2/22/22	0/1/1/1
6	NAG	A	812	-	-	4/6/26/26	0/1/1/1
7	BMA	A	816	-	-	2/2/22/22	0/1/1/1
6	NAG	A	810	-	-	1/6/26/26	0/1/1/1
6	NAG	A	807	-	-	1/6/26/26	0/1/1/1
6	NAG	B	715	-	-	0/6/26/26	0/1/1/1
6	NAG	B	714	-	-	2/6/26/26	0/1/1/1
6	NAG	B	710	-	-	2/6/26/26	0/1/1/1

All (20) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	806	NAG	C1-C2	4.69	1.58	1.52
6	B	707	NAG	C1-C2	3.86	1.57	1.52
6	B	709	NAG	C1-C2	3.44	1.57	1.52
6	B	713	NAG	C1-C2	3.29	1.56	1.52
6	A	807	NAG	C1-C2	3.07	1.56	1.52
6	A	812	NAG	C1-C2	3.07	1.56	1.52
6	A	813	NAG	C1-C2	2.90	1.56	1.52
6	A	808	NAG	C1-C2	2.78	1.56	1.52
6	A	801	NAG	C1-C2	2.65	1.56	1.52
6	B	713	NAG	O5-C1	2.60	1.49	1.42
6	A	818	NAG	C1-C2	2.43	1.55	1.52
6	B	714	NAG	C1-C2	2.33	1.55	1.52
6	B	707	NAG	C3-C2	2.32	1.57	1.53
6	B	702	NAG	C1-C2	2.29	1.55	1.52
7	A	819	BMA	C4-C5	2.24	1.57	1.53
6	B	719	NAG	C1-C2	2.20	1.55	1.52
6	A	806	NAG	C3-C2	2.18	1.57	1.53
8	A	820	MAN	C4-C3	2.12	1.57	1.52
7	B	717	BMA	C1-C2	2.10	1.57	1.52
6	B	708	NAG	C1-C2	2.04	1.55	1.52

All (26) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	713	NAG	C1-C2-N2	4.32	115.73	110.73
6	B	712	NAG	C1-C2-N2	3.78	115.11	110.73

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	B	707	NAG	C4-C3-C2	3.39	115.33	110.40
6	A	806	NAG	C4-C3-C2	3.22	115.09	110.40
6	A	814	NAG	O1-C1-O5	-3.07	101.29	110.41
6	A	812	NAG	C1-C2-N2	3.03	114.23	110.73
6	B	715	NAG	O1-C1-C2	-2.88	103.23	109.22
6	A	812	NAG	C2-N2-C7	2.70	129.43	123.11
6	A	806	NAG	C2-N2-C7	2.60	129.19	123.11
6	B	713	NAG	C2-N2-C7	2.60	129.19	123.11
6	B	719	NAG	C1-O5-C5	2.55	118.58	113.65
6	A	818	NAG	C1-O5-C5	2.45	118.40	113.65
6	B	705	NAG	C1-O5-C5	2.43	118.35	113.65
6	A	804	NAG	C1-O5-C5	2.42	118.33	113.65
6	B	712	NAG	C2-N2-C7	2.36	128.63	123.11
6	A	806	NAG	C1-C2-C3	2.28	113.65	110.54
6	A	818	NAG	C3-C4-C5	2.26	114.33	110.23
8	B	701	MAN	O2-C2-C3	-2.23	105.11	110.38
6	B	719	NAG	C3-C4-C5	2.17	114.17	110.23
6	A	808	NAG	O5-C5-C4	-2.15	105.83	109.70
6	B	707	NAG	C1-C2-C3	2.14	113.47	110.54
6	A	808	NAG	C4-C3-C2	2.12	113.48	110.40
6	B	709	NAG	C4-C3-C2	2.04	113.37	110.40
6	B	719	NAG	O5-C5-C4	2.03	113.35	109.70
6	A	811	NAG	C1-C2-N2	2.01	113.05	110.73
8	A	820	MAN	C1-O5-C5	2.00	117.53	113.65

There are no chirality outliers.

All (85) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	A	806	NAG	C1-C2-N2-C7
6	A	811	NAG	C1-C2-N2-C7
6	B	707	NAG	C1-C2-N2-C7
6	B	712	NAG	C1-C2-N2-C7
6	A	818	NAG	C4-C5-C6-O6
6	B	706	NAG	C4-C5-C6-O6
6	B	705	NAG	O5-C5-C6-O6
6	A	815	NAG	O5-C5-C6-O6
6	B	718	NAG	C4-C5-C6-O6
6	B	719	NAG	C4-C5-C6-O6
6	A	803	NAG	O5-C5-C6-O6
6	A	805	NAG	C4-C5-C6-O6
6	B	710	NAG	O5-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	B	712	NAG	O5-C5-C6-O6
8	A	820	MAN	O5-C5-C6-O6
8	B	701	MAN	O5-C5-C6-O6
6	A	809	NAG	O5-C5-C6-O6
6	A	804	NAG	C4-C5-C6-O6
6	A	813	NAG	O5-C5-C6-O6
6	A	802	NAG	O5-C5-C6-O6
6	A	804	NAG	O5-C5-C6-O6
6	B	718	NAG	O5-C5-C6-O6
6	B	710	NAG	C4-C5-C6-O6
6	A	806	NAG	O5-C5-C6-O6
6	A	806	NAG	C4-C5-C6-O6
6	A	805	NAG	O5-C5-C6-O6
6	B	719	NAG	O5-C5-C6-O6
6	A	815	NAG	C4-C5-C6-O6
6	A	818	NAG	O5-C5-C6-O6
6	B	706	NAG	O5-C5-C6-O6
6	A	809	NAG	C4-C5-C6-O6
6	B	705	NAG	C4-C5-C6-O6
6	A	802	NAG	C4-C5-C6-O6
6	A	813	NAG	C4-C5-C6-O6
6	B	712	NAG	C4-C5-C6-O6
7	A	816	BMA	C4-C5-C6-O6
7	B	717	BMA	C4-C5-C6-O6
6	A	802	NAG	C3-C2-N2-C7
6	A	808	NAG	C3-C2-N2-C7
6	A	803	NAG	C4-C5-C6-O6
6	A	812	NAG	C8-C7-N2-C2
6	A	812	NAG	O7-C7-N2-C2
6	A	818	NAG	C8-C7-N2-C2
6	A	818	NAG	O7-C7-N2-C2
6	B	713	NAG	C8-C7-N2-C2
6	B	713	NAG	O7-C7-N2-C2
6	B	719	NAG	C8-C7-N2-C2
6	B	719	NAG	O7-C7-N2-C2
7	A	816	BMA	O5-C5-C6-O6
7	B	717	BMA	O5-C5-C6-O6
6	B	703	NAG	C3-C2-N2-C7
6	B	709	NAG	C3-C2-N2-C7
6	B	714	NAG	C3-C2-N2-C7
8	B	701	MAN	C4-C5-C6-O6
8	A	820	MAN	C4-C5-C6-O6

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Mol	Chain	Res	Type	Atoms
6	A	811	NAG	C4-C5-C6-O6
6	A	817	NAG	O5-C5-C6-O6
6	A	808	NAG	O5-C5-C6-O6
6	A	812	NAG	O5-C5-C6-O6
6	A	808	NAG	C4-C5-C6-O6
6	A	810	NAG	O5-C5-C6-O6
6	B	704	NAG	C4-C5-C6-O6
6	A	812	NAG	C3-C2-N2-C7
6	B	713	NAG	C3-C2-N2-C7
7	B	720	BMA	O5-C5-C6-O6
6	B	704	NAG	O5-C5-C6-O6
6	B	713	NAG	O5-C5-C6-O6
6	A	801	NAG	O5-C5-C6-O6
6	B	702	NAG	O5-C5-C6-O6
6	B	708	NAG	O5-C5-C6-O6
7	A	819	BMA	O5-C5-C6-O6
6	A	807	NAG	O5-C5-C6-O6
6	B	711	NAG	O5-C5-C6-O6
6	B	709	NAG	C4-C5-C6-O6
6	B	703	NAG	C1-C2-N2-C7
6	A	811	NAG	O5-C5-C6-O6
6	B	709	NAG	O5-C5-C6-O6
6	A	802	NAG	C1-C2-N2-C7
6	A	808	NAG	C1-C2-N2-C7
6	B	709	NAG	C1-C2-N2-C7
6	B	714	NAG	C1-C2-N2-C7
6	B	716	NAG	C4-C5-C6-O6
6	B	707	NAG	C3-C2-N2-C7
6	B	704	NAG	C1-C2-N2-C7
6	B	716	NAG	O5-C5-C6-O6

There are no ring outliers.

28 monomers are involved in 33 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	814	NAG	2	0
6	A	808	NAG	2	0
6	B	713	NAG	3	0
6	A	802	NAG	2	0
6	B	719	NAG	1	0
6	A	817	NAG	1	0
6	B	703	NAG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	818	NAG	2	0
6	B	704	NAG	1	0
6	B	708	NAG	1	0
6	A	809	NAG	1	0
6	B	711	NAG	1	0
7	B	717	BMA	1	0
6	A	811	NAG	2	0
6	A	803	NAG	1	0
7	B	720	BMA	1	0
6	A	806	NAG	1	0
6	A	815	NAG	2	0
6	B	712	NAG	1	0
6	B	716	NAG	3	0
7	A	819	BMA	1	0
6	A	812	NAG	3	0
7	A	816	BMA	1	0
6	A	810	NAG	1	0
6	A	807	NAG	3	0
6	B	715	NAG	3	0
6	B	714	NAG	1	0
6	B	710	NAG	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

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	663/671 (98%)	-0.43	3 (0%) 87 75	58, 85, 118, 174	0
1	B	663/671 (98%)	-0.43	2 (0%) 90 82	53, 88, 119, 184	0
2	C	107/107 (100%)	-0.45	0 100 100	61, 83, 109, 136	0
2	E	107/107 (100%)	-0.43	0 100 100	70, 88, 113, 145	0
3	D	122/127 (96%)	-0.38	0 100 100	58, 90, 124, 145	0
3	F	122/127 (96%)	-0.20	1 (0%) 82 68	81, 116, 135, 157	0
4	H	116/121 (95%)	0.05	4 (3%) 48 32	106, 152, 186, 196	0
4	X	114/121 (94%)	0.23	0 100 100	90, 141, 186, 204	0
5	L	105/107 (98%)	0.38	6 (5%) 29 20	96, 176, 209, 243	0
5	Y	107/107 (100%)	-0.23	2 (1%) 66 48	65, 109, 136, 157	0
All	All	2226/2266 (98%)	-0.31	18 (0%) 82 68	53, 94, 168, 243	0

All (18) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	691	PHE	4.2
5	L	105	GLU	3.6
1	B	679	THR	3.6
4	H	2	VAL	3.2
5	L	80	GLN	3.1
1	B	691	PHE	2.8
4	H	18	LEU	2.7
4	H	101	ILE	2.7
4	H	116	SER	2.6
5	Y	11	LEU	2.6
5	L	11	LEU	2.6
5	L	19	VAL	2.6
5	Y	80	GLN	2.5

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Mol	Chain	Res	Type	RSRZ
5	L	20	THR	2.2
5	L	9	SER	2.2
1	A	337	SER	2.1
3	F	122	SER	2.1
1	A	338	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](#)

There are no oligosaccharides in this entry.

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
6	NAG	A	804	15/15	0.11	0.16	178,194,206,206	0
6	NAG	B	705	15/15	0.24	0.12	180,202,208,209	0
6	NAG	B	706	15/15	0.30	0.14	150,171,187,187	0
7	BMA	A	819	12/12	0.31	0.13	169,179,186,188	0
6	NAG	A	813	15/15	0.32	0.13	161,184,192,194	0
6	NAG	A	808	15/15	0.34	0.13	135,159,166,174	0
8	MAN	A	820	12/12	0.35	0.10	194,200,204,204	0
6	NAG	B	714	15/15	0.47	0.18	133,159,168,176	0
7	BMA	B	717	12/12	0.52	0.08	174,186,193,193	0
6	NAG	B	712	15/15	0.58	0.11	152,173,179,179	0
6	NAG	A	805	15/15	0.60	0.10	140,164,172,174	0
8	MAN	B	701	12/12	0.60	0.08	192,213,216,217	0
6	NAG	A	807	15/15	0.62	0.12	111,128,145,145	0
6	NAG	B	709	15/15	0.63	0.10	140,160,166,167	0
7	BMA	B	720	12/12	0.64	0.09	157,176,180,180	0
7	BMA	A	816	12/12	0.67	0.08	165,176,181,184	0
6	NAG	B	708	15/15	0.68	0.11	116,145,157,162	0
6	NAG	B	702	15/15	0.70	0.10	104,136,158,159	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
6	NAG	B	719	15/15	0.70	0.13	151,158,173,176	0
6	NAG	A	812	15/15	0.72	0.12	117,140,153,154	0
6	NAG	B	710	15/15	0.74	0.09	93,126,154,156	0
6	NAG	A	818	15/15	0.75	0.12	143,158,163,165	0
6	NAG	A	811	15/15	0.76	0.11	148,169,183,183	0
6	NAG	B	704	15/15	0.77	0.12	129,144,159,162	0
6	NAG	A	801	15/15	0.79	0.08	104,132,144,145	0
6	NAG	B	707	15/15	0.80	0.12	109,133,145,156	0
6	NAG	B	713	15/15	0.80	0.09	91,108,123,131	0
6	NAG	A	809	15/15	0.83	0.08	96,124,135,138	0
6	NAG	B	716	15/15	0.84	0.09	110,121,139,154	0
6	NAG	A	815	15/15	0.84	0.09	117,131,140,143	0
6	NAG	A	802	15/15	0.86	0.10	97,118,126,129	0
6	NAG	A	810	15/15	0.86	0.08	94,109,143,145	0
6	NAG	A	806	15/15	0.87	0.11	106,129,147,147	0
6	NAG	B	718	15/15	0.87	0.10	90,96,99,100	0
6	NAG	B	703	15/15	0.87	0.12	117,133,145,147	0
6	NAG	B	711	15/15	0.88	0.10	97,107,148,154	0
6	NAG	A	803	15/15	0.90	0.09	104,135,142,142	0
6	NAG	A	817	15/15	0.91	0.07	80,92,102,104	0
6	NAG	A	814	15/15	0.96	0.05	49,66,82,82	0
6	NAG	B	715	15/15	0.97	0.06	58,64,84,87	0

6.5 Other polymers

There are no such residues in this entry.