



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 12:23 PM UTC

PDB ID : 6YFJ / pdb_00006yfq
Title : Virus-like particle of bacteriophage ESE001
Authors : Rumnieks, J.; Kalnins, G.; Sisovs, M.; Lieknina, I.; Tars, K.
Deposited on : 2020-03-26
Resolution : 3.23 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
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

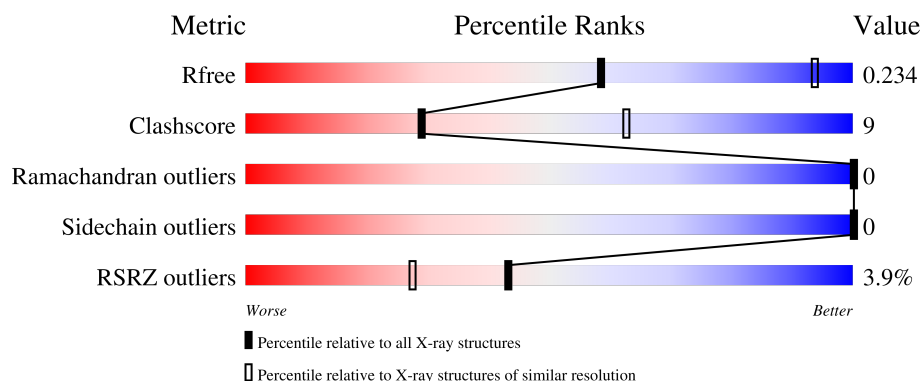
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.23 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	2153 (3.28-3.20)
Clashscore	190562	2275 (3.28-3.20)
Ramachandran outliers	187476	2233 (3.28-3.20)
Sidechain outliers	187428	2232 (3.28-3.20)
RSRZ outliers	180081	2153 (3.28-3.20)

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	AA	118	2% 73% 27%
1	AB	118	7% 80% 20%
1	AC	118	3% 81% 19%
1	AD	118	3% 81% 19%
1	AE	118	8% 81% 19%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	AF	118	
1	AG	118	
1	AH	118	
1	AI	118	
1	AJ	118	
1	AK	118	
1	AL	118	
1	AM	118	
1	AN	118	
1	AO	118	
1	AP	118	
1	AQ	118	
1	AR	118	
1	AS	118	
1	AT	118	
1	AU	118	
1	AV	118	
1	AW	118	
1	AX	118	
1	AY	118	
1	AZ	118	
1	BA	118	
1	BB	118	
1	BC	118	
1	BD	118	

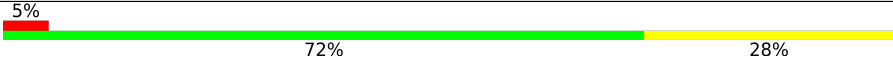
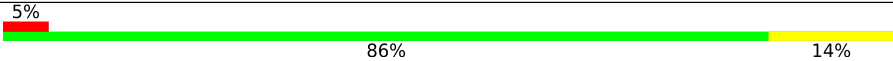
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	BE	118	4% 71% 29%
1	BF	118	4% 76% 24%
1	BG	118	3% 78% 22%
1	BH	118	% 75% 25%
1	BI	118	3% 75% 25%
1	BJ	118	3% 86% 14%
1	BK	118	7% 80% 20%
1	BL	118	5% 77% 23%
1	BM	118	2% 84% 16%
1	BN	118	5% 77% 23%
1	BO	118	8% 86% 14%
1	BP	118	4% 79% 21%
1	BQ	118	6% 80% 20%
1	BR	118	7% 79% 21%
1	BS	118	3% 85% 15%
1	BT	118	6% 79% 21%
1	BU	118	7% 78% 22%
1	BV	118	3% 85% 15%
1	BW	118	3% 78% 22%
1	BX	118	7% 84% 16%
1	BY	118	3% 83% 17%
1	BZ	118	3% 69% 31%
1	CA	118	5% 78% 22%
1	CB	118	3% 90% 10%
1	CC	118	3% 80% 20%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	CD	118	
1	CE	118	
1	CF	118	
1	CG	118	
1	CH	118	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 52380 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 coat protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AA	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AB	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AC	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AD	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AE	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AF	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AG	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AH	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AI	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AJ	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AK	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AL	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AM	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AN	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AO	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	AP	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AQ	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AR	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AS	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AT	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AU	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AV	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AW	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AX	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AY	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	AZ	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BA	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BB	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BC	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BD	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BE	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BF	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BG	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BH	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BI	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BJ	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BK	118	Total 873	C 542	N 150	O 176	S 5	0	0	0

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	BL	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BM	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BN	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BO	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BP	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BQ	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BR	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BS	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BT	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BU	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BV	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BW	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BX	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BY	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	BZ	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	CA	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	CB	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	CC	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	CD	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	CE	118	Total 873	C 542	N 150	O 176	S 5	0	0	0
1	CF	118	Total 873	C 542	N 150	O 176	S 5	0	0	0

Continued on next page...

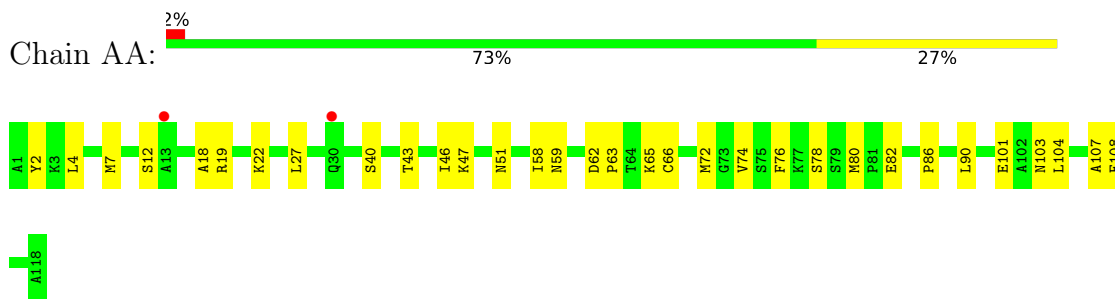
Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	CG	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			
1	CH	118	Total	C	N	O	S	0	0	0
			873	542	150	176	5			

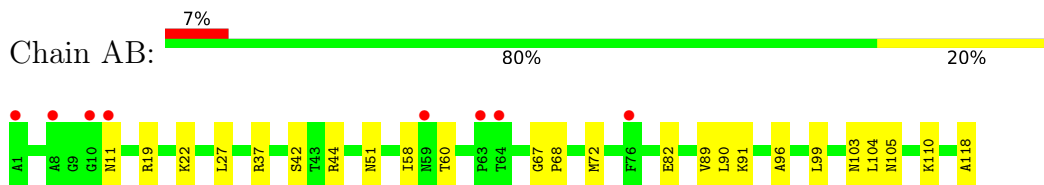
3 Residue-property plots

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.

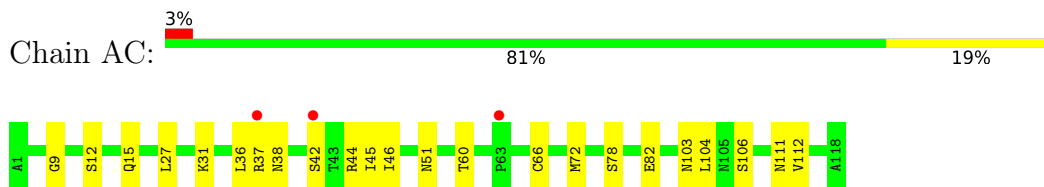
- Molecule 1: coat protein



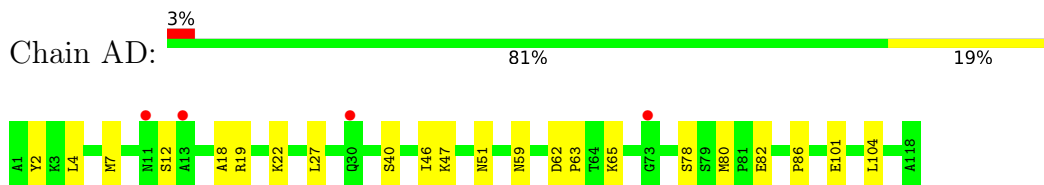
- Molecule 1: coat protein



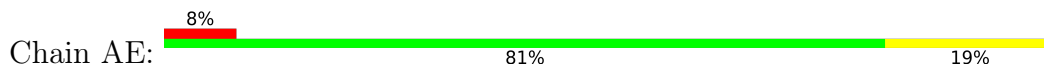
- Molecule 1: coat protein



- Molecule 1: coat protein

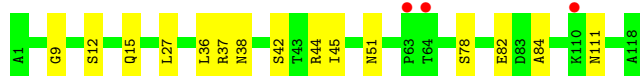
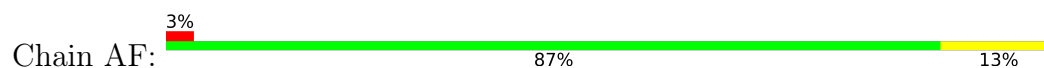


- Molecule 1: coat protein

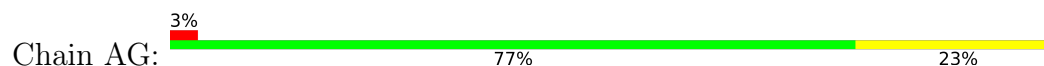




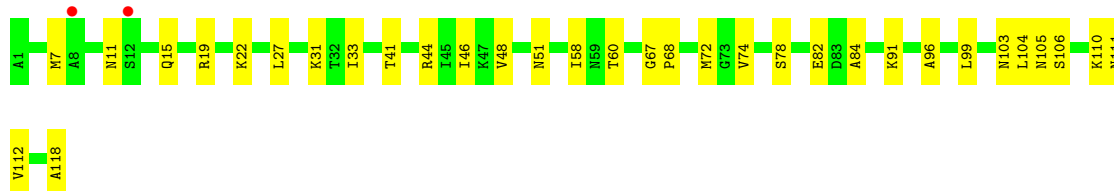
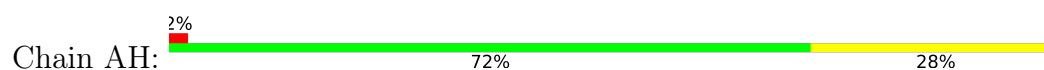
- Molecule 1: coat protein



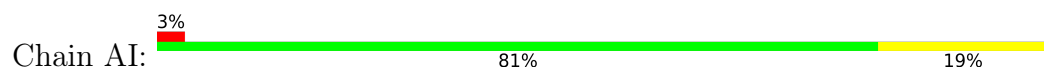
- Molecule 1: coat protein



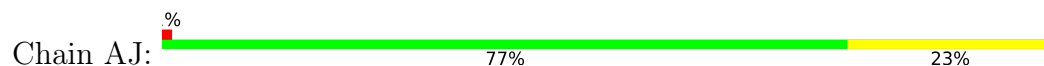
- Molecule 1: coat protein



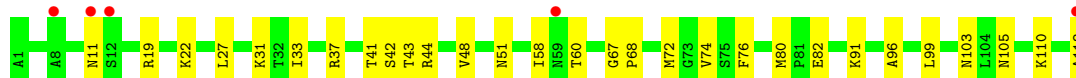
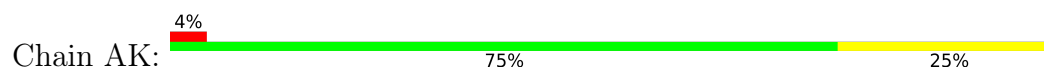
- Molecule 1: coat protein



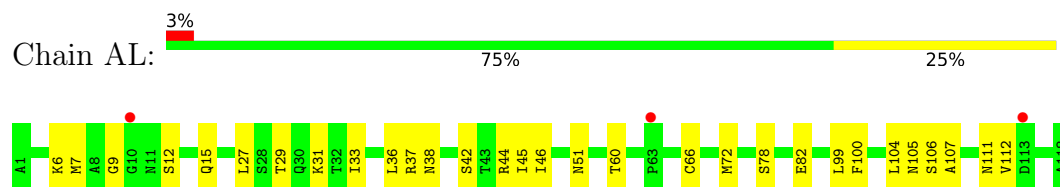
- Molecule 1: coat protein



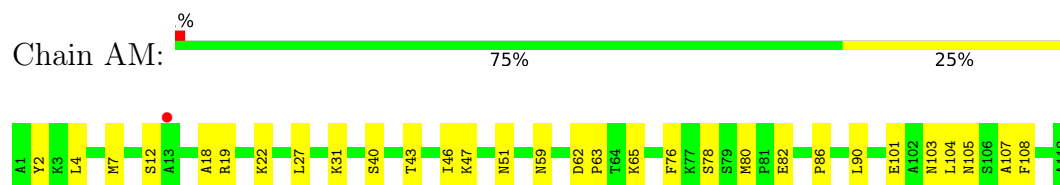
- Molecule 1: coat protein



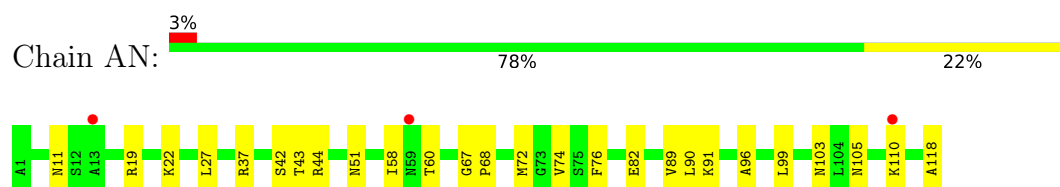
- Molecule 1: coat protein



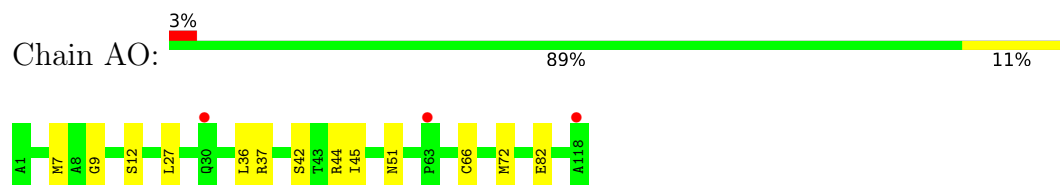
- Molecule 1: coat protein



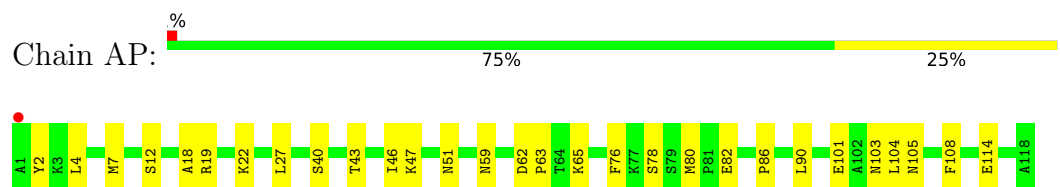
- Molecule 1: coat protein



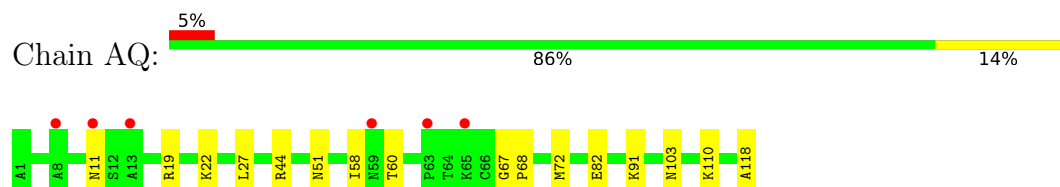
- Molecule 1: coat protein



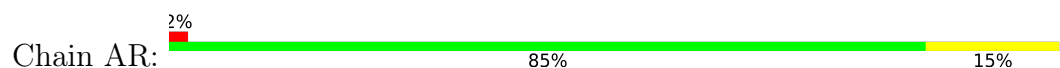
- Molecule 1: coat protein

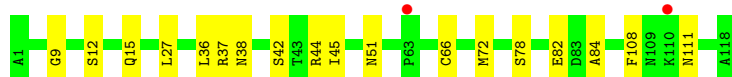


- Molecule 1: coat protein

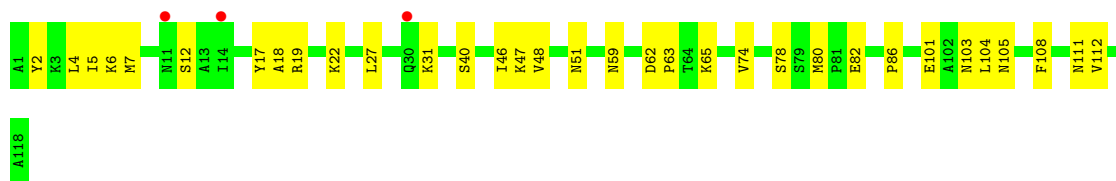
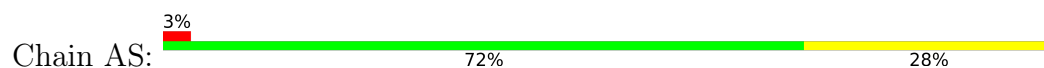


- Molecule 1: coat protein

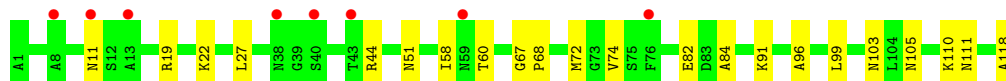
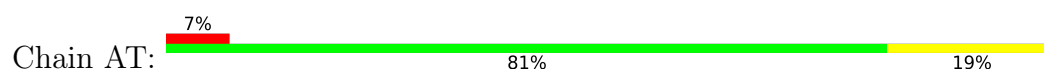




- Molecule 1: coat protein



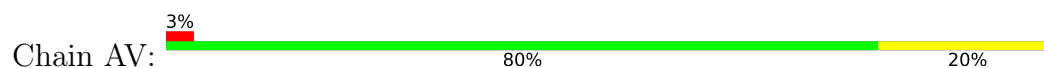
- Molecule 1: coat protein



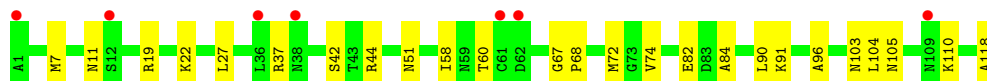
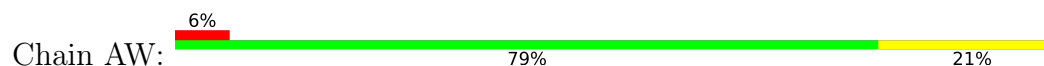
- Molecule 1: coat protein



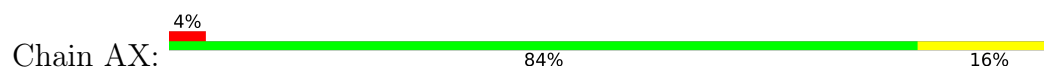
- Molecule 1: coat protein



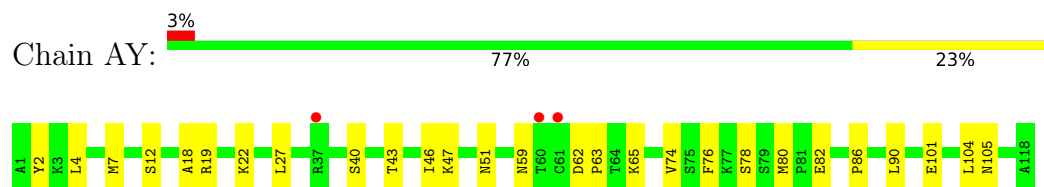
- Molecule 1: coat protein



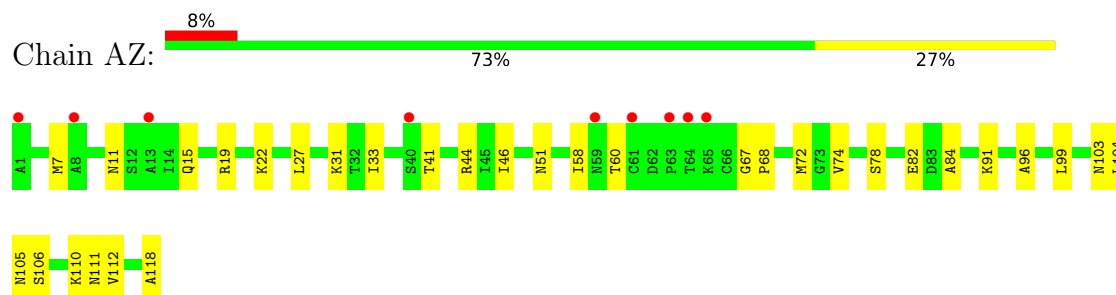
- Molecule 1: coat protein



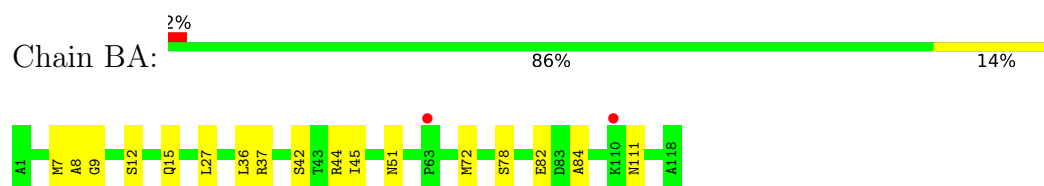
- Molecule 1: coat protein



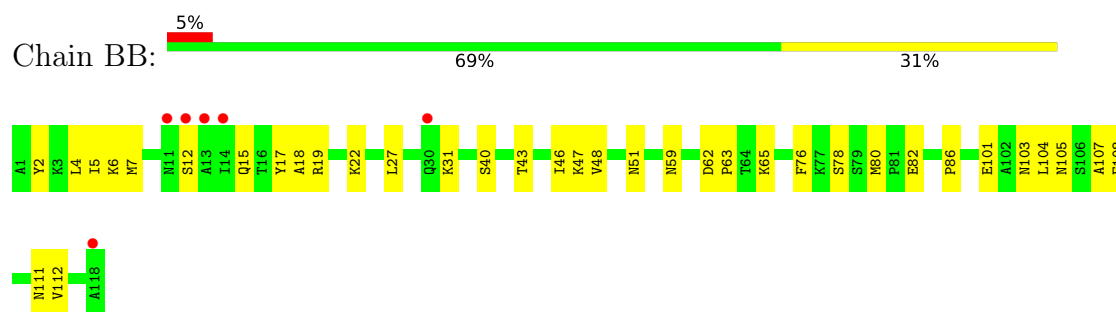
- Molecule 1: coat protein



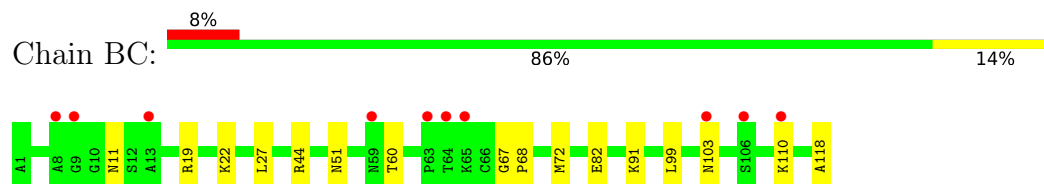
- Molecule 1: coat protein



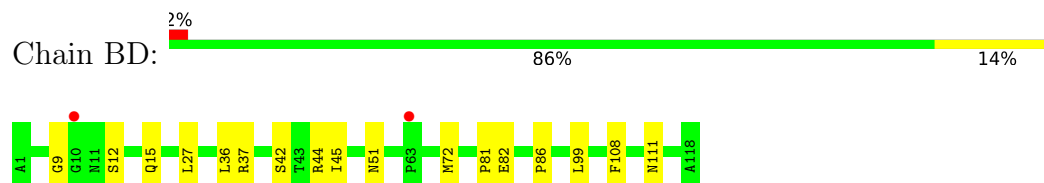
- Molecule 1: coat protein



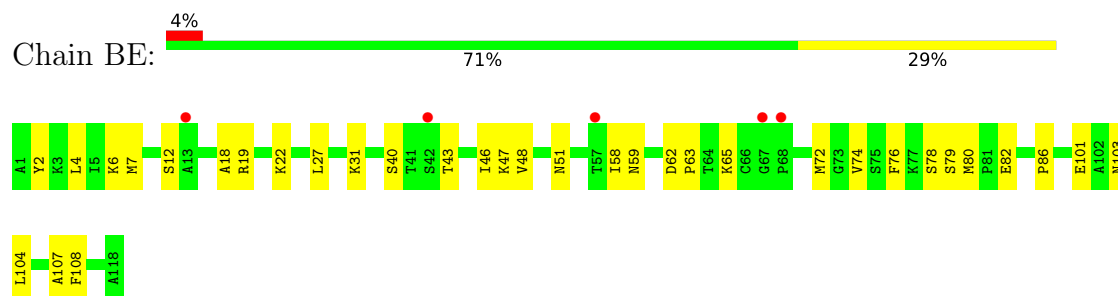
- Molecule 1: coat protein



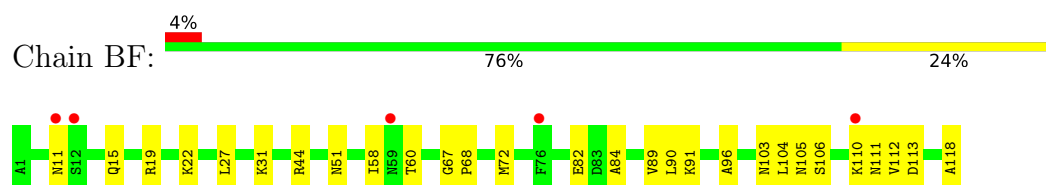
- Molecule 1: coat protein



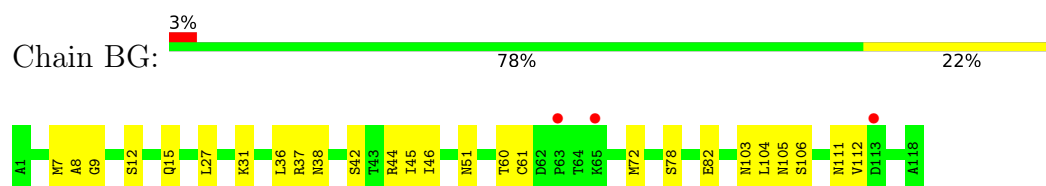
- Molecule 1: coat protein



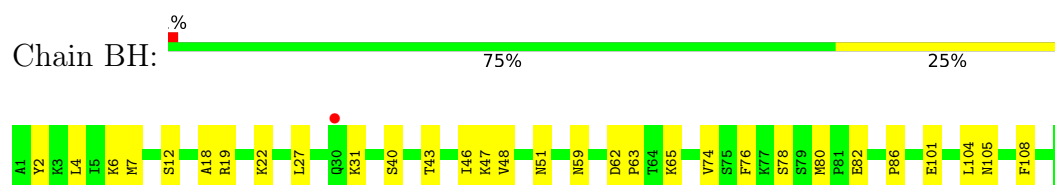
- Molecule 1: coat protein



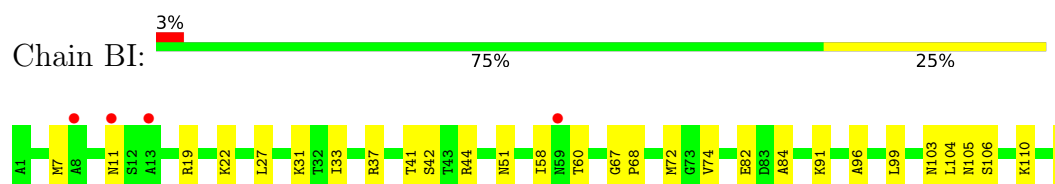
- Molecule 1: coat protein



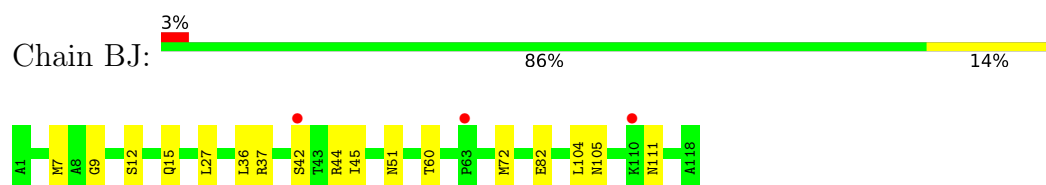
- Molecule 1: coat protein



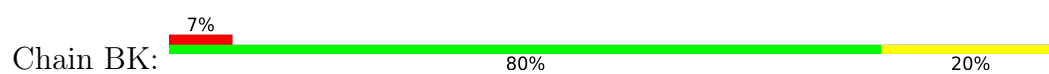
- Molecule 1: coat protein



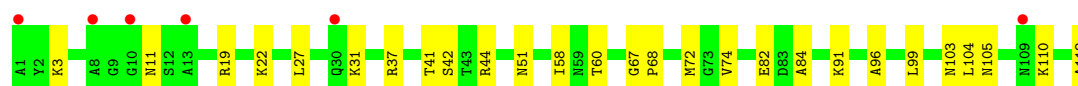
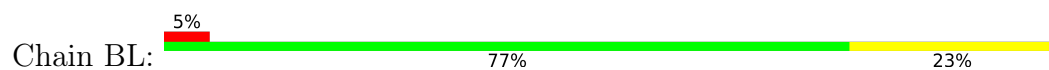
- Molecule 1: coat protein



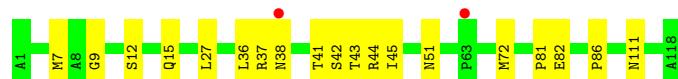
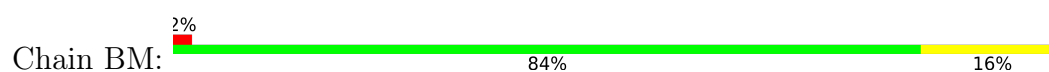
- Molecule 1: coat protein



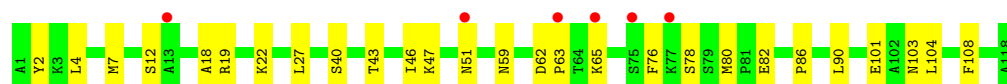
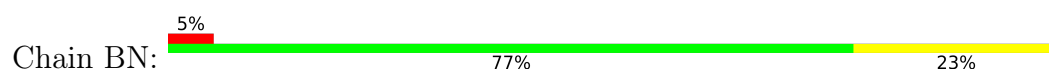
- Molecule 1: coat protein



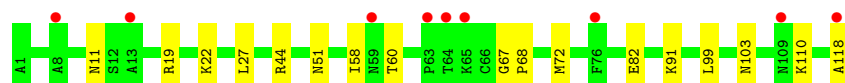
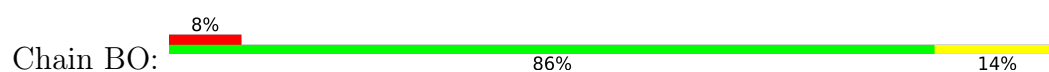
- Molecule 1: coat protein



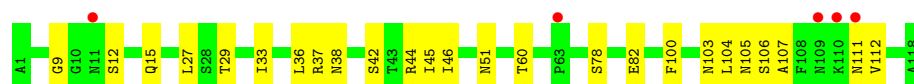
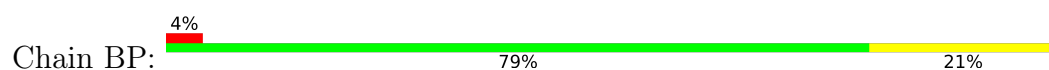
- Molecule 1: coat protein



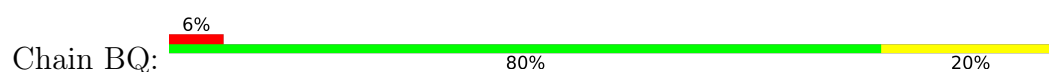
- Molecule 1: coat protein



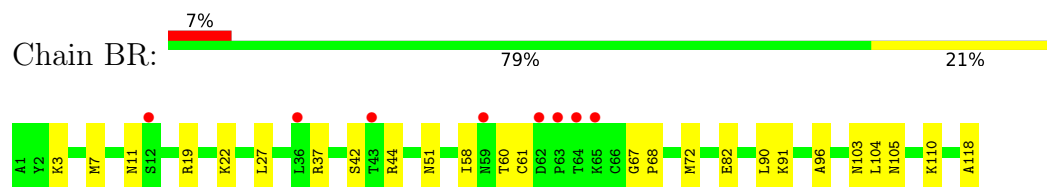
- Molecule 1: coat protein



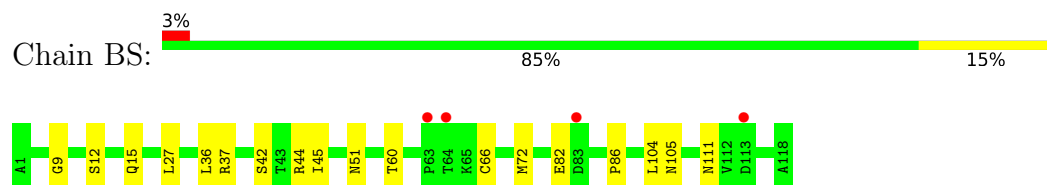
- Molecule 1: coat protein



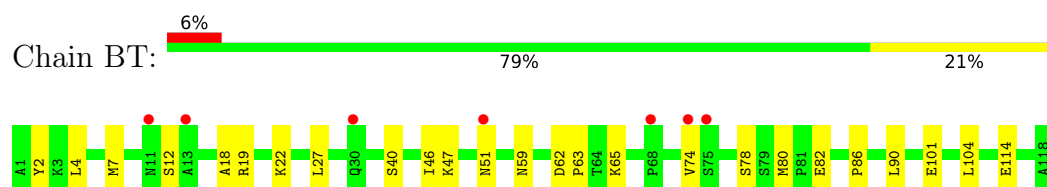
• Molecule 1: coat protein



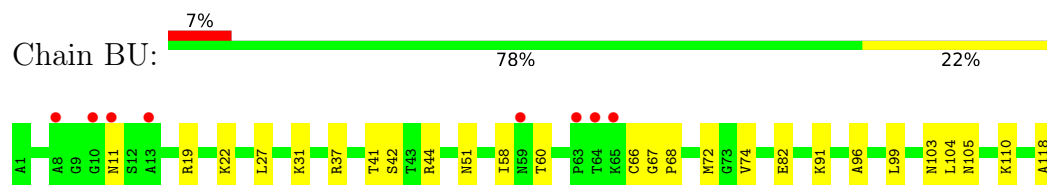
• Molecule 1: coat protein



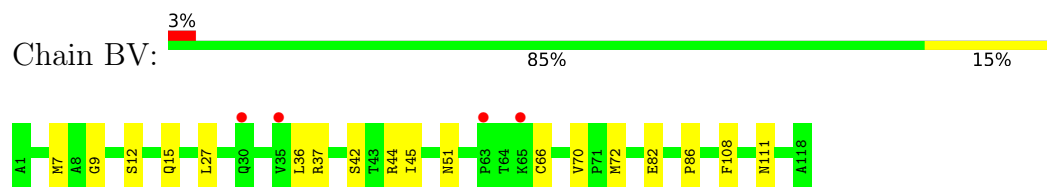
• Molecule 1: coat protein



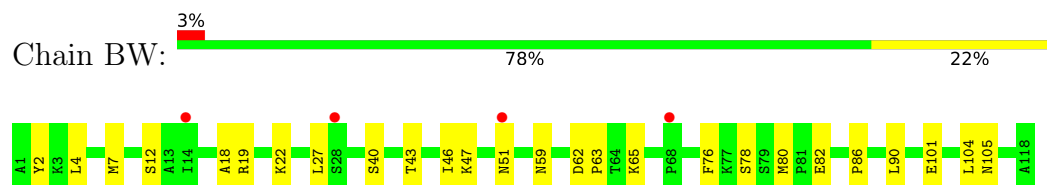
• Molecule 1: coat protein



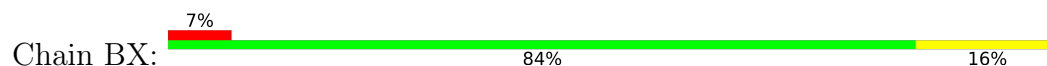
• Molecule 1: coat protein

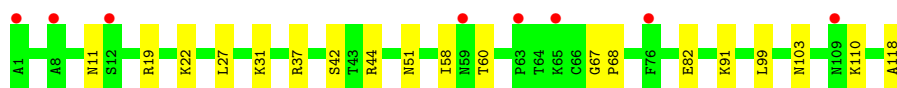


• Molecule 1: coat protein

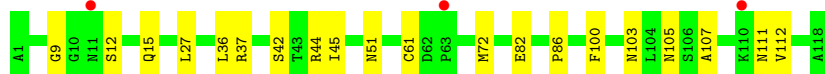
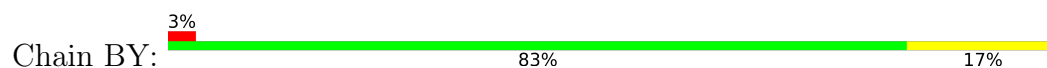


• Molecule 1: coat protein

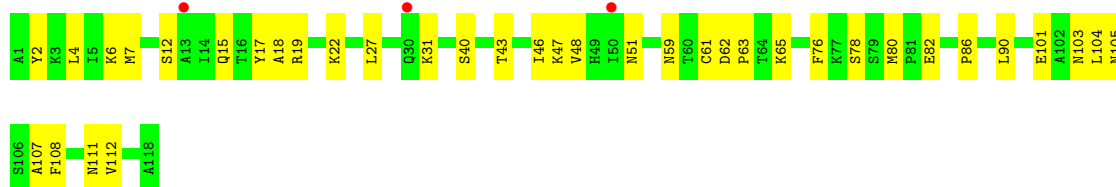




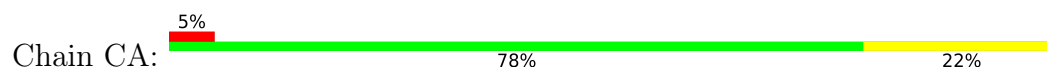
- Molecule 1: coat protein



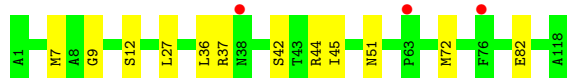
- Molecule 1: coat protein



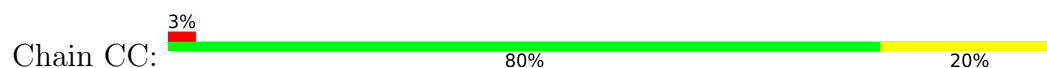
- Molecule 1: coat protein



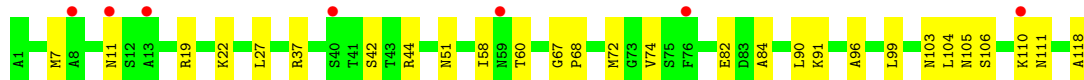
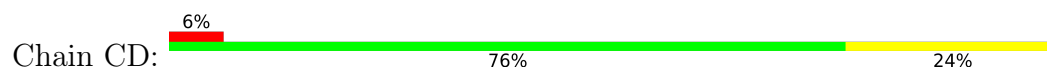
- Molecule 1: coat protein



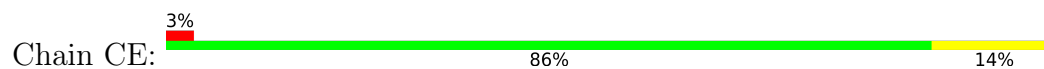
- Molecule 1: coat protein



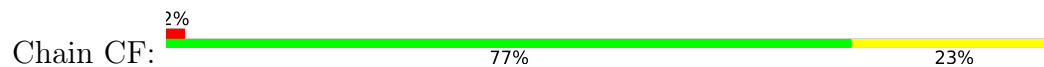
- Molecule 1: coat protein



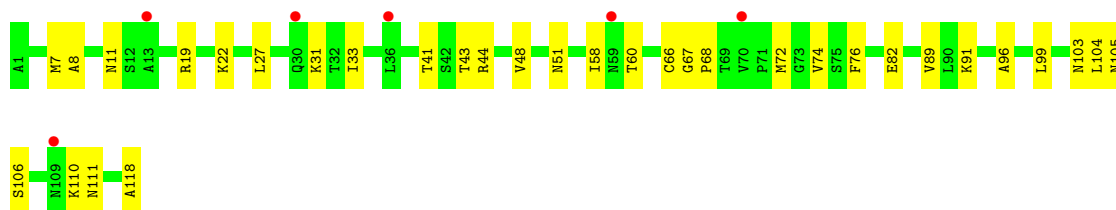
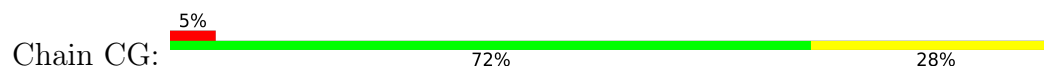
● Molecule 1: coat protein



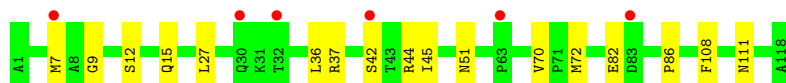
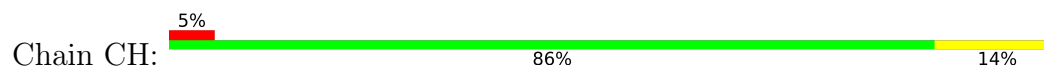
● Molecule 1: coat protein



● Molecule 1: coat protein



● Molecule 1: coat protein



4 Data and refinement statistics

Property	Value	Source
Space group	H 3	Depositor
Cell constants a, b, c, α , β , γ	283.35Å 283.35Å 666.04Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.94 – 3.23 48.94 – 3.23	Depositor EDS
% Data completeness (in resolution range)	97.5 (48.94-3.23) 97.9 (48.94-3.23)	Depositor EDS
R_{merge}	0.28	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.97 (at 3.25Å)	Xtriage
Refinement program	PHENIX 1.14_3260	Depositor
R, R_{free}	0.232 , 0.232 0.235 , 0.234	Depositor DCC
R_{free} test set	10054 reflections (3.22%)	wwPDB-VP
Wilson B-factor (Å ²)	120.1	Xtriage
Anisotropy	0.065	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 73.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	0.013 for -h,1/3*h-1/3*k-1/3*l,-4/3*h-8/3*k+1/3*l 0.009 for -1/3*h+1/3*k+1/3*l,-k,8/3*h+4/3*k+1/3*l 0.017 for -2/3*h-1/3*k-1/3*l,-1/3*h-2/3*k+1/3*l,-4/3*h+4/3*k+1/3*l 0.010 for 1/3*h+2/3*k-1/3*l,-k,-8/3*h-4/3*k-1/3*l 0.017 for -1/3*h-2/3*k+1/3*l,-2/3*h-1/3*k-1/3*l,4/3*h-4/3*k-1/3*l 0.010 for -h,2/3*h+1/3*k+1/3*l,4/3*h+8/3*k-1/3*l 0.033 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	52380	wwPDB-VP
Average B, all atoms (Å ²)	97.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.02% 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

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	AA	0.23	0/884	0.49	0/1198
1	AB	0.23	0/884	0.48	0/1198
1	AC	0.23	0/884	0.48	0/1198
1	AD	0.24	0/884	0.49	0/1198
1	AE	0.24	0/884	0.48	0/1198
1	AF	0.23	0/884	0.48	0/1198
1	AG	0.24	0/884	0.49	0/1198
1	AH	0.23	0/884	0.48	0/1198
1	AI	0.24	0/884	0.48	0/1198
1	AJ	0.23	0/884	0.49	0/1198
1	AK	0.24	0/884	0.48	0/1198
1	AL	0.24	0/884	0.48	0/1198
1	AM	0.24	0/884	0.49	0/1198
1	AN	0.24	0/884	0.48	0/1198
1	AO	0.23	0/884	0.48	0/1198
1	AP	0.24	0/884	0.49	0/1198
1	AQ	0.23	0/884	0.48	0/1198
1	AR	0.23	0/884	0.48	0/1198
1	AS	0.24	0/884	0.49	0/1198
1	AT	0.23	0/884	0.48	0/1198
1	AU	0.24	0/884	0.48	0/1198
1	AV	0.24	0/884	0.49	0/1198
1	AW	0.23	0/884	0.48	0/1198
1	AX	0.23	0/884	0.48	0/1198
1	AY	0.24	0/884	0.49	0/1198
1	AZ	0.23	0/884	0.48	0/1198
1	BA	0.23	0/884	0.48	0/1198
1	BB	0.24	0/884	0.49	0/1198
1	BC	0.23	0/884	0.48	0/1198
1	BD	0.24	0/884	0.48	0/1198
1	BE	0.23	0/884	0.49	0/1198
1	BF	0.23	0/884	0.48	0/1198
1	BG	0.23	0/884	0.48	0/1198
1	BH	0.24	0/884	0.49	0/1198

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	BI	0.23	0/884	0.48	0/1198
1	BJ	0.24	0/884	0.48	0/1198
1	BK	0.24	0/884	0.49	0/1198
1	BL	0.23	0/884	0.48	0/1198
1	BM	0.24	0/884	0.48	0/1198
1	BN	0.24	0/884	0.49	0/1198
1	BO	0.23	0/884	0.48	0/1198
1	BP	0.23	0/884	0.48	0/1198
1	BQ	0.24	0/884	0.49	0/1198
1	BR	0.23	0/884	0.48	0/1198
1	BS	0.24	0/884	0.48	0/1198
1	BT	0.24	0/884	0.49	0/1198
1	BU	0.23	0/884	0.48	0/1198
1	BV	0.24	0/884	0.48	0/1198
1	BW	0.24	0/884	0.49	0/1198
1	BX	0.23	0/884	0.48	0/1198
1	BY	0.24	0/884	0.48	0/1198
1	BZ	0.24	0/884	0.49	0/1198
1	CA	0.23	0/884	0.48	0/1198
1	CB	0.24	0/884	0.48	0/1198
1	CC	0.24	0/884	0.49	0/1198
1	CD	0.23	0/884	0.48	0/1198
1	CE	0.23	0/884	0.48	0/1198
1	CF	0.24	0/884	0.49	0/1198
1	CG	0.24	0/884	0.48	0/1198
1	CH	0.23	0/884	0.48	0/1198
All	All	0.23	0/53040	0.49	0/71880

There are no bond length outliers.

There are no bond angle outliers.

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	AA	873	0	895	32	0
1	AB	873	0	895	19	0
1	AC	873	0	895	23	0
1	AD	873	0	896	13	0
1	AE	873	0	896	20	0
1	AF	873	0	895	13	0
1	AG	873	0	895	22	0
1	AH	873	0	895	33	0
1	AI	873	0	895	25	0
1	AJ	873	0	896	25	0
1	AK	873	0	895	27	0
1	AL	873	0	895	35	0
1	AM	873	0	896	28	0
1	AN	873	0	895	23	0
1	AO	873	0	897	9	0
1	AP	873	0	895	26	0
1	AQ	873	0	896	11	0
1	AR	873	0	896	17	1
1	AS	873	0	895	32	0
1	AT	873	0	895	21	0
1	AU	873	0	895	14	0
1	AV	873	0	896	14	1
1	AW	873	0	896	22	0
1	AX	873	0	895	20	0
1	AY	873	0	895	23	0
1	AZ	873	0	895	35	0
1	BA	873	0	895	14	0
1	BB	873	0	895	38	0
1	BC	873	0	896	10	0
1	BD	873	0	896	15	0
1	BE	873	0	895	35	0
1	BF	873	0	895	26	0
1	BG	873	0	895	26	0
1	BH	873	0	895	26	0
1	BI	873	0	895	26	0
1	BJ	873	0	895	16	0
1	BK	873	0	896	16	0
1	BL	873	0	896	25	0
1	BM	873	0	895	19	0
1	BN	873	0	895	26	0
1	BO	873	0	896	12	0
1	BP	873	0	896	29	0
1	BQ	873	0	896	14	1

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	BR	873	0	896	21	0
1	BS	873	0	895	17	0
1	BT	873	0	895	21	0
1	BU	873	0	895	22	0
1	BV	873	0	895	18	0
1	BW	873	0	895	21	0
1	BX	873	0	896	11	1
1	BY	873	0	896	21	0
1	BZ	873	0	896	42	0
1	CA	873	0	895	22	0
1	CB	873	0	896	9	0
1	CC	873	0	895	21	0
1	CD	873	0	895	24	0
1	CE	873	0	895	13	0
1	CF	873	0	895	24	0
1	CG	873	0	895	34	0
1	CH	873	0	895	17	0
All	All	52380	0	53722	916	2

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 9.

All (916) 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:AI:7:MET:HG2	1:BZ:108:PHE:HE1	1.28	0.95
1:AI:7:MET:HG2	1:BZ:108:PHE:CE1	2.03	0.93
1:AL:15:GLN:HE22	1:BP:111:ASN:HA	1.38	0.89
1:AL:111:ASN:HA	1:BP:15:GLN:HE22	1.42	0.85
1:AS:27:LEU:HD21	1:BF:96:ALA:HB2	1.57	0.84
1:AH:96:ALA:HB2	1:BZ:27:LEU:HD21	1.59	0.84
1:AZ:96:ALA:HB2	1:BB:27:LEU:HD21	1.60	0.83
1:BE:27:LEU:HD21	1:CG:96:ALA:HB2	1.58	0.83
1:AX:15:GLN:HE22	1:BY:111:ASN:HA	1.46	0.80
1:AL:15:GLN:NE2	1:BP:111:ASN:HA	1.98	0.79
1:AB:96:ALA:HB2	1:CC:27:LEU:HD21	1.63	0.78
1:BH:27:LEU:HD21	1:CD:96:ALA:HB2	1.65	0.78
1:AH:31:LYS:NZ	1:BZ:103:ASN:O	2.18	0.77
1:CA:96:ALA:HB2	1:CF:27:LEU:HD21	1.67	0.77
1:AA:27:LEU:HD21	1:AK:96:ALA:HB2	1.69	0.75
1:AF:111:ASN:HA	1:BJ:15:GLN:HE22	1.50	0.75

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:7:MET:HG2	1:BB:108:PHE:HE1	1.52	0.74
1:AF:15:GLN:HE22	1:BJ:111:ASN:HA	1.52	0.74
1:AL:111:ASN:HA	1:BP:15:GLN:NE2	2.04	0.73
1:BR:105:ASN:HB3	1:BT:80:MET:HE2	1.70	0.73
1:BE:59:ASN:OD1	1:CG:41:THR:HG21	1.88	0.73
1:AR:111:ASN:HA	1:BS:15:GLN:HE22	1.53	0.72
1:AG:27:LEU:HD21	1:AT:96:ALA:HB2	1.69	0.72
1:AS:108:PHE:HE1	1:BG:7:MET:HG2	1.55	0.71
1:AS:80:MET:HE2	1:BF:105:ASN:HB3	1.71	0.71
1:BR:96:ALA:HB2	1:BT:27:LEU:HD21	1.72	0.71
1:BH:101:GLU:HA	1:BH:104:LEU:HG	1.74	0.70
1:CC:101:GLU:HA	1:CC:104:LEU:HG	1.74	0.70
1:AJ:101:GLU:HA	1:AJ:104:LEU:HG	1.74	0.70
1:BE:80:MET:HB3	1:CG:72:MET:HE3	1.72	0.70
1:BN:27:LEU:HD21	1:BU:96:ALA:HB2	1.73	0.70
1:CF:101:GLU:HA	1:CF:104:LEU:HG	1.74	0.70
1:AM:101:GLU:HA	1:AM:104:LEU:HG	1.74	0.70
1:BZ:101:GLU:HA	1:BZ:104:LEU:HG	1.74	0.70
1:AG:101:GLU:HA	1:AG:104:LEU:HG	1.74	0.70
1:BE:43:THR:HG21	1:CG:58:ILE:HG23	1.74	0.70
1:BQ:101:GLU:HA	1:BQ:104:LEU:HG	1.74	0.70
1:AE:96:ALA:HB2	1:BW:27:LEU:HD21	1.71	0.69
1:BK:101:GLU:HA	1:BK:104:LEU:HG	1.74	0.69
1:AA:101:GLU:HA	1:AA:104:LEU:HG	1.74	0.69
1:AS:101:GLU:HA	1:AS:104:LEU:HG	1.74	0.69
1:AD:101:GLU:HA	1:AD:104:LEU:HG	1.74	0.69
1:AY:101:GLU:HA	1:AY:104:LEU:HG	1.74	0.69
1:BE:101:GLU:HA	1:BE:104:LEU:HG	1.74	0.69
1:AS:108:PHE:CE1	1:BG:7:MET:HG2	2.28	0.69
1:BA:7:MET:HG2	1:BB:108:PHE:CE1	2.27	0.69
1:AM:27:LEU:HD21	1:BI:96:ALA:HB2	1.75	0.69
1:BN:101:GLU:HA	1:BN:104:LEU:HG	1.74	0.69
1:BT:101:GLU:HA	1:BT:104:LEU:HG	1.74	0.69
1:AJ:62:ASP:OD1	1:AO:66:CYS:SG	2.51	0.68
1:AP:101:GLU:HA	1:AP:104:LEU:HG	1.74	0.68
1:BW:101:GLU:HA	1:BW:104:LEU:HG	1.74	0.68
1:AC:15:GLN:HE22	1:BG:111:ASN:HA	1.57	0.68
1:AM:80:MET:HE2	1:BI:105:ASN:HB3	1.74	0.68
1:BB:101:GLU:HA	1:BB:104:LEU:HG	1.74	0.67
1:AA:108:PHE:CE1	1:AL:7:MET:HG2	2.29	0.67
1:AV:101:GLU:HA	1:AV:104:LEU:HG	1.74	0.67

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AZ:105:ASN:HB3	1:BB:80:MET:HE2	1.74	0.67
1:AH:15:GLN:HE22	1:BZ:112:VAL:H	1.42	0.67
1:AM:59:ASN:OD1	1:BI:41:THR:HG21	1.95	0.67
1:AA:103:ASN:O	1:AK:31:LYS:NZ	2.26	0.66
1:AJ:27:LEU:HD21	1:AN:96:ALA:HB2	1.77	0.66
1:AB:105:ASN:HB3	1:CC:80:MET:HE2	1.77	0.66
1:AP:27:LEU:HD21	1:BL:96:ALA:HB2	1.78	0.66
1:AX:15:GLN:NE2	1:BY:111:ASN:HA	2.10	0.65
1:BA:111:ASN:HA	1:CE:15:GLN:HE22	1.62	0.65
1:AP:80:MET:HE2	1:BL:105:ASN:HB3	1.80	0.64
1:CA:105:ASN:HB3	1:CF:80:MET:HE2	1.79	0.64
1:AD:80:MET:HE1	1:AD:86:PRO:HD3	1.80	0.64
1:CC:80:MET:HE1	1:CC:86:PRO:HD3	1.80	0.64
1:AA:59:ASN:OD1	1:AK:41:THR:HG21	1.97	0.64
1:AC:111:ASN:HA	1:BG:15:GLN:HE22	1.63	0.64
1:BD:15:GLN:HE22	1:CH:111:ASN:HA	1.63	0.64
1:BK:80:MET:HE1	1:BK:86:PRO:HD3	1.80	0.64
1:AG:80:MET:HE1	1:AG:86:PRO:HD3	1.80	0.64
1:AJ:80:MET:HE1	1:AJ:86:PRO:HD3	1.80	0.64
1:AM:80:MET:HE1	1:AM:86:PRO:HD3	1.80	0.63
1:BH:80:MET:HE1	1:BH:86:PRO:HD3	1.80	0.63
1:BE:80:MET:HE1	1:BE:86:PRO:HD3	1.80	0.63
1:BW:80:MET:HE1	1:BW:86:PRO:HD3	1.80	0.63
1:AA:80:MET:HE1	1:AA:86:PRO:HD3	1.80	0.63
1:AU:15:GLN:HE22	1:BV:111:ASN:HA	1.61	0.63
1:BQ:80:MET:HE1	1:BQ:86:PRO:HD3	1.80	0.63
1:AM:108:PHE:CE1	1:BJ:7:MET:HG2	2.33	0.63
1:AL:60:THR:HG21	1:BP:38:ASN:HD22	1.63	0.63
1:AS:80:MET:HE1	1:AS:86:PRO:HD3	1.80	0.63
1:AW:96:ALA:HB2	1:AY:27:LEU:HD21	1.81	0.63
1:BN:80:MET:HE1	1:BN:86:PRO:HD3	1.80	0.63
1:CF:80:MET:HE1	1:CF:86:PRO:HD3	1.80	0.63
1:AX:111:ASN:HA	1:BY:15:GLN:HE22	1.64	0.62
1:BA:15:GLN:HE22	1:CE:111:ASN:HA	1.64	0.62
1:AP:80:MET:HE1	1:AP:86:PRO:HD3	1.80	0.62
1:BT:80:MET:HE1	1:BT:86:PRO:HD3	1.80	0.62
1:AY:80:MET:HE1	1:AY:86:PRO:HD3	1.80	0.62
1:AV:80:MET:HE1	1:AV:86:PRO:HD3	1.80	0.62
1:BB:80:MET:HE1	1:BB:86:PRO:HD3	1.80	0.62
1:BZ:80:MET:HE1	1:BZ:86:PRO:HD3	1.80	0.62
1:BE:108:PHE:CE1	1:CH:7:MET:HG2	2.35	0.62

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BE:80:MET:HB3	1:CG:72:MET:HB2	1.82	0.62
1:AC:15:GLN:NE2	1:BG:111:ASN:HA	2.16	0.61
1:AM:108:PHE:HE1	1:BJ:7:MET:HG2	1.65	0.61
1:AG:43:THR:HG21	1:AT:58:ILE:HG23	1.82	0.61
1:AJ:74:VAL:HG11	1:AN:90:LEU:HD23	1.83	0.61
1:BH:80:MET:HB3	1:CD:72:MET:HE3	1.81	0.61
1:AC:112:VAL:H	1:BG:15:GLN:HE22	1.48	0.60
1:BH:80:MET:HE2	1:CD:105:ASN:HB3	1.82	0.60
1:AC:15:GLN:HE22	1:BG:112:VAL:H	1.47	0.60
1:AU:111:ASN:HA	1:BV:15:GLN:HE22	1.66	0.60
1:BE:76:PHE:CE2	1:CG:74:VAL:HG13	2.37	0.60
1:AA:108:PHE:HE1	1:AL:7:MET:HG2	1.65	0.60
1:AR:15:GLN:HE22	1:BS:111:ASN:HA	1.65	0.60
1:AH:105:ASN:HB3	1:BZ:80:MET:HE2	1.82	0.60
1:CA:90:LEU:HD23	1:CF:74:VAL:HG11	1.84	0.60
1:AL:38:ASN:HD22	1:BP:60:THR:HG21	1.65	0.59
1:BD:111:ASN:HA	1:CH:15:GLN:HE22	1.67	0.59
1:BE:103:ASN:O	1:CG:31:LYS:NZ	2.30	0.59
1:AA:80:MET:HB3	1:AK:72:MET:HE3	1.83	0.59
1:AG:80:MET:HB3	1:AT:72:MET:HE3	1.85	0.59
1:BN:80:MET:HE2	1:BU:105:ASN:HB3	1.85	0.59
1:AZ:111:ASN:OD1	1:BB:6:LYS:NZ	2.36	0.58
1:AH:33:ILE:HD11	1:BZ:107:ALA:HB1	1.83	0.58
1:BE:31:LYS:NZ	1:CG:106:SER:OG	2.37	0.58
1:BE:108:PHE:HE1	1:CH:7:MET:HG2	1.69	0.58
1:AM:103:ASN:O	1:BI:31:LYS:NZ	2.33	0.58
1:CB:7:MET:HG2	1:CF:108:PHE:HE1	1.69	0.58
1:AM:80:MET:HB3	1:BI:72:MET:HE3	1.85	0.58
1:AI:108:PHE:CE1	1:BK:5:ILE:HB	2.39	0.58
1:AZ:31:LYS:NZ	1:BB:103:ASN:O	2.36	0.57
1:BN:108:PHE:CE1	1:BV:7:MET:HG2	2.40	0.57
1:AZ:111:ASN:HD21	1:CH:111:ASN:ND2	2.02	0.57
1:CB:7:MET:HG2	1:CF:108:PHE:CE1	2.39	0.57
1:CA:72:MET:HE3	1:CF:80:MET:HB3	1.87	0.57
1:AI:111:ASN:HA	1:BM:15:GLN:HE22	1.69	0.56
1:AG:80:MET:HB3	1:AT:72:MET:HB2	1.88	0.56
1:AH:72:MET:HE3	1:BZ:80:MET:HB3	1.87	0.56
1:AH:111:ASN:OD1	1:BZ:6:LYS:NZ	2.39	0.56
1:AZ:72:MET:HE3	1:BB:80:MET:HB3	1.86	0.56
1:AZ:106:SER:OG	1:BB:31:LYS:NZ	2.38	0.56
1:BR:72:MET:HE3	1:BT:80:MET:HB3	1.87	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AG:63:PRO:HG2	1:AG:65:LYS:HG2	1.88	0.56
1:AL:15:GLN:HE22	1:BP:112:VAL:H	1.54	0.56
1:AC:44:ARG:HD2	1:AC:82:GLU:OE1	2.06	0.56
1:BT:63:PRO:HG2	1:BT:65:LYS:HG2	1.88	0.56
1:CF:63:PRO:HG2	1:CF:65:LYS:HG2	1.88	0.56
1:AA:76:PHE:CE2	1:AK:74:VAL:HG13	2.40	0.56
1:AX:44:ARG:HD2	1:AX:82:GLU:OE1	2.06	0.56
1:BK:63:PRO:HG2	1:BK:65:LYS:HG2	1.88	0.56
1:CC:63:PRO:HG2	1:CC:65:LYS:HG2	1.88	0.56
1:AF:44:ARG:HD2	1:AF:82:GLU:OE1	2.06	0.56
1:AI:8:ALA:HB2	1:BZ:111:ASN:HA	1.88	0.56
1:AL:33:ILE:HD11	1:BP:107:ALA:HB1	1.87	0.56
1:BM:44:ARG:HD2	1:BM:82:GLU:OE1	2.06	0.56
1:BV:44:ARG:HD2	1:BV:82:GLU:OE1	2.06	0.56
1:CE:44:ARG:HD2	1:CE:82:GLU:OE1	2.06	0.56
1:AJ:63:PRO:HG2	1:AJ:65:LYS:HG2	1.88	0.56
1:AL:44:ARG:HD2	1:AL:82:GLU:OE1	2.06	0.56
1:AS:63:PRO:HG2	1:AS:65:LYS:HG2	1.88	0.56
1:BW:63:PRO:HG2	1:BW:65:LYS:HG2	1.88	0.56
1:CB:44:ARG:HD2	1:CB:82:GLU:OE1	2.06	0.56
1:AJ:80:MET:HE2	1:AN:105:ASN:HB3	1.88	0.55
1:AR:44:ARG:HD2	1:AR:82:GLU:OE1	2.06	0.55
1:AU:44:ARG:HD2	1:AU:82:GLU:OE1	2.06	0.55
1:BB:63:PRO:HG2	1:BB:65:LYS:HG2	1.88	0.55
1:BE:63:PRO:HG2	1:BE:65:LYS:HG2	1.88	0.55
1:BH:63:PRO:HG2	1:BH:65:LYS:HG2	1.88	0.55
1:BJ:44:ARG:HD2	1:BJ:82:GLU:OE1	2.06	0.55
1:BS:44:ARG:HD2	1:BS:82:GLU:OE1	2.06	0.55
1:BY:44:ARG:HD2	1:BY:82:GLU:OE1	2.06	0.55
1:AW:105:ASN:HB3	1:AY:80:MET:HE2	1.88	0.55
1:BN:63:PRO:HG2	1:BN:65:LYS:HG2	1.88	0.55
1:CH:44:ARG:HD2	1:CH:82:GLU:OE1	2.06	0.55
1:AE:11:ASN:ND2	1:AF:9:GLY:O	2.40	0.55
1:AL:112:VAL:H	1:BP:15:GLN:HE22	1.55	0.55
1:BA:44:ARG:HD2	1:BA:82:GLU:OE1	2.06	0.55
1:BD:44:ARG:HD2	1:BD:82:GLU:OE1	2.06	0.55
1:BC:11:ASN:ND2	1:BD:9:GLY:O	2.40	0.55
1:BG:44:ARG:HD2	1:BG:82:GLU:OE1	2.06	0.55
1:BU:11:ASN:ND2	1:BV:9:GLY:O	2.40	0.55
1:BX:11:ASN:ND2	1:BY:9:GLY:O	2.40	0.55
1:AA:80:MET:HB3	1:AK:72:MET:HB2	1.88	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BQ:63:PRO:HG2	1:BQ:65:LYS:HG2	1.88	0.55
1:CG:11:ASN:ND2	1:CH:9:GLY:O	2.40	0.55
1:AY:63:PRO:HG2	1:AY:65:LYS:HG2	1.88	0.55
1:AT:11:ASN:ND2	1:AU:9:GLY:O	2.40	0.55
1:CA:11:ASN:ND2	1:CB:9:GLY:O	2.40	0.55
1:AH:11:ASN:ND2	1:AI:9:GLY:O	2.40	0.55
1:AI:44:ARG:HD2	1:AI:82:GLU:OE1	2.06	0.55
1:AQ:11:ASN:ND2	1:AR:9:GLY:O	2.40	0.55
1:AW:72:MET:HE3	1:AY:80:MET:HB3	1.89	0.55
1:AB:11:ASN:ND2	1:AC:9:GLY:O	2.40	0.54
1:AD:63:PRO:HG2	1:AD:65:LYS:HG2	1.88	0.54
1:AN:11:ASN:ND2	1:AO:9:GLY:O	2.40	0.54
1:AV:63:PRO:HG2	1:AV:65:LYS:HG2	1.88	0.54
1:BL:11:ASN:ND2	1:BM:9:GLY:O	2.40	0.54
1:BO:11:ASN:ND2	1:BP:9:GLY:O	2.40	0.54
1:AC:111:ASN:HA	1:BG:15:GLN:NE2	2.22	0.54
1:AO:44:ARG:HD2	1:AO:82:GLU:OE1	2.06	0.54
1:BP:44:ARG:HD2	1:BP:82:GLU:OE1	2.06	0.54
1:BR:11:ASN:ND2	1:BS:9:GLY:O	2.40	0.54
1:AA:63:PRO:HG2	1:AA:65:LYS:HG2	1.88	0.54
1:AE:72:MET:HB2	1:BW:80:MET:HB3	1.90	0.54
1:AP:63:PRO:HG2	1:AP:65:LYS:HG2	1.88	0.54
1:AW:44:ARG:HD3	1:AW:82:GLU:OE1	2.08	0.54
1:BI:11:ASN:ND2	1:BJ:9:GLY:O	2.40	0.54
1:BZ:63:PRO:HG2	1:BZ:65:LYS:HG2	1.88	0.54
1:AZ:11:ASN:ND2	1:BA:9:GLY:O	2.40	0.54
1:CD:44:ARG:HD3	1:CD:82:GLU:OE1	2.08	0.54
1:AA:58:ILE:HG23	1:AK:43:THR:OG1	2.08	0.54
1:AE:58:ILE:HG23	1:BW:43:THR:HG21	1.89	0.54
1:BI:44:ARG:HD3	1:BI:82:GLU:OE1	2.08	0.54
1:BN:108:PHE:HE1	1:BV:7:MET:HG2	1.72	0.54
1:BX:44:ARG:HD3	1:BX:82:GLU:OE1	2.08	0.54
1:CD:11:ASN:ND2	1:CE:9:GLY:O	2.40	0.54
1:AN:44:ARG:HD3	1:AN:82:GLU:OE1	2.08	0.54
1:BR:44:ARG:HD3	1:BR:82:GLU:OE1	2.08	0.54
1:AA:80:MET:HE2	1:AK:105:ASN:HB3	1.88	0.54
1:AK:11:ASN:ND2	1:AL:9:GLY:O	2.40	0.54
1:AZ:44:ARG:HD3	1:AZ:82:GLU:OE1	2.08	0.54
1:BR:105:ASN:HB3	1:BT:80:MET:CE	2.38	0.54
1:AT:44:ARG:HD3	1:AT:82:GLU:OE1	2.08	0.53
1:AK:44:ARG:HD3	1:AK:82:GLU:OE1	2.08	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AM:63:PRO:HG2	1:AM:65:LYS:HG2	1.88	0.53
1:AW:11:ASN:ND2	1:AX:9:GLY:O	2.40	0.53
1:BF:11:ASN:ND2	1:BG:9:GLY:O	2.40	0.53
1:BL:44:ARG:HD3	1:BL:82:GLU:OE1	2.08	0.53
1:AB:44:ARG:HD3	1:AB:82:GLU:OE1	2.08	0.53
1:AH:44:ARG:HD3	1:AH:82:GLU:OE1	2.08	0.53
1:AH:106:SER:OG	1:BZ:31:LYS:NZ	2.41	0.53
1:AQ:44:ARG:HD3	1:AQ:82:GLU:OE1	2.08	0.53
1:BF:44:ARG:HD3	1:BF:82:GLU:OE1	2.08	0.53
1:BU:44:ARG:HD3	1:BU:82:GLU:OE1	2.08	0.53
1:BO:44:ARG:HD3	1:BO:82:GLU:OE1	2.08	0.53
1:AA:43:THR:HG21	1:AK:58:ILE:HG23	1.90	0.53
1:BE:47:LYS:HA	1:BE:78:SER:O	2.09	0.53
1:CF:47:LYS:HA	1:CF:78:SER:O	2.09	0.53
1:AH:41:THR:HG21	1:BZ:59:ASN:OD1	2.07	0.53
1:BT:47:LYS:HA	1:BT:78:SER:O	2.09	0.53
1:AH:33:ILE:CD1	1:BZ:107:ALA:HB1	2.39	0.53
1:BZ:47:LYS:HA	1:BZ:78:SER:O	2.09	0.53
1:AE:72:MET:HE3	1:BW:80:MET:HB3	1.90	0.53
1:BE:80:MET:HE2	1:CG:105:ASN:HB3	1.89	0.53
1:BQ:47:LYS:HA	1:BQ:78:SER:O	2.09	0.53
1:AG:47:LYS:HA	1:AG:78:SER:O	2.09	0.53
1:AL:46:ILE:HD13	1:BP:105:ASN:C	2.34	0.53
1:CA:44:ARG:HD3	1:CA:82:GLU:OE1	2.08	0.53
1:CC:47:LYS:HA	1:CC:78:SER:O	2.09	0.53
1:CG:44:ARG:HD3	1:CG:82:GLU:OE1	2.08	0.53
1:AE:44:ARG:HD3	1:AE:82:GLU:OE1	2.08	0.52
1:BW:47:LYS:HA	1:BW:78:SER:O	2.09	0.52
1:AG:105:ASN:ND2	1:AT:84:ALA:O	2.42	0.52
1:AL:29:THR:HB	1:BP:100:PHE:CE2	2.44	0.52
1:AS:74:VAL:HG11	1:BF:90:LEU:HD23	1.91	0.52
1:AV:47:LYS:HA	1:AV:78:SER:O	2.09	0.52
1:BC:44:ARG:HD3	1:BC:82:GLU:OE1	2.08	0.52
1:BH:47:LYS:HA	1:BH:78:SER:O	2.09	0.52
1:BK:47:LYS:HA	1:BK:78:SER:O	2.09	0.52
1:AJ:47:LYS:HA	1:AJ:78:SER:O	2.09	0.52
1:AS:47:LYS:HA	1:AS:78:SER:O	2.09	0.52
1:AA:47:LYS:HA	1:AA:78:SER:O	2.09	0.52
1:AB:72:MET:HE3	1:CC:80:MET:HB3	1.92	0.52
1:AI:15:GLN:HE22	1:BM:111:ASN:HA	1.75	0.52
1:AY:47:LYS:HA	1:AY:78:SER:O	2.09	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AP:47:LYS:HA	1:AP:78:SER:O	2.09	0.52
1:BN:47:LYS:HA	1:BN:78:SER:O	2.09	0.52
1:AD:47:LYS:HA	1:AD:78:SER:O	2.09	0.52
1:AB:90:LEU:HD23	1:CC:74:VAL:HG11	1.90	0.52
1:AM:47:LYS:HA	1:AM:78:SER:O	2.09	0.52
1:AZ:84:ALA:O	1:BB:105:ASN:ND2	2.43	0.52
1:BB:5:ILE:HB	1:CH:108:PHE:CE1	2.45	0.51
1:BE:107:ALA:HB1	1:CG:33:ILE:HD11	1.92	0.51
1:BH:43:THR:HG21	1:CD:58:ILE:HG23	1.92	0.51
1:AS:17:TYR:OH	1:BF:112:VAL:HG23	2.10	0.51
1:AG:80:MET:HE2	1:AT:105:ASN:HB3	1.92	0.51
1:AL:46:ILE:HD13	1:BP:106:SER:N	2.26	0.51
1:AR:84:ALA:O	1:BS:105:ASN:ND2	2.35	0.51
1:BB:47:LYS:HA	1:BB:78:SER:O	2.09	0.51
1:AL:33:ILE:CD1	1:BP:107:ALA:HB1	2.40	0.51
1:BE:6:LYS:NZ	1:CG:111:ASN:OD1	2.40	0.51
1:BN:103:ASN:O	1:BU:31:LYS:NZ	2.38	0.51
1:AS:5:ILE:HB	1:BV:108:PHE:CE1	2.46	0.51
1:AF:15:GLN:NE2	1:BJ:111:ASN:HA	2.24	0.51
1:AE:105:ASN:HB3	1:BW:80:MET:HE2	1.92	0.51
1:AM:12:SER:HB3	1:AO:12:SER:HB2	1.93	0.51
1:BN:12:SER:HB3	1:BP:12:SER:HB2	1.93	0.51
1:AP:90:LEU:HD11	1:BL:104:LEU:HD12	1.93	0.51
1:BK:12:SER:HB3	1:BM:12:SER:HB2	1.93	0.51
1:BT:12:SER:HB3	1:BV:12:SER:HB2	1.93	0.51
1:BQ:12:SER:HB3	1:BS:12:SER:HB2	1.93	0.50
1:CC:12:SER:HB3	1:CE:12:SER:HB2	1.93	0.50
1:AA:12:SER:HB3	1:AC:12:SER:HB2	1.93	0.50
1:BB:12:SER:HB3	1:BD:12:SER:HB2	1.93	0.50
1:AV:12:SER:HB3	1:AX:12:SER:HB2	1.93	0.50
1:BH:12:SER:HB3	1:BJ:12:SER:HB2	1.93	0.50
1:AH:74:VAL:HG13	1:BZ:76:PHE:CE2	2.46	0.50
1:AD:12:SER:HB3	1:AF:12:SER:HB2	1.94	0.50
1:AY:12:SER:HB3	1:BA:12:SER:HB2	1.93	0.50
1:AG:12:SER:HB3	1:AI:12:SER:HB2	1.93	0.50
1:AJ:12:SER:HB3	1:AL:12:SER:HB2	1.93	0.50
1:AL:107:ALA:HB1	1:BP:33:ILE:HD11	1.94	0.50
1:AP:80:MET:HB3	1:BL:72:MET:HE3	1.93	0.50
1:AS:12:SER:HB3	1:AU:12:SER:HB2	1.93	0.50
1:AP:12:SER:HB3	1:AR:12:SER:HB2	1.93	0.50
1:AS:6:LYS:NZ	1:BF:111:ASN:OD1	2.44	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BW:12:SER:HB3	1:BY:12:SER:HB2	1.93	0.49
1:BZ:12:SER:HB3	1:CB:12:SER:HB2	1.93	0.49
1:AE:74:VAL:HG13	1:BW:76:PHE:CE2	2.47	0.49
1:AP:59:ASN:OD1	1:BL:41:THR:HG21	2.12	0.49
1:BN:80:MET:HB3	1:BU:72:MET:HE3	1.94	0.49
1:AF:111:ASN:HA	1:BJ:15:GLN:NE2	2.22	0.49
1:AH:15:GLN:HE22	1:BZ:112:VAL:N	2.09	0.49
1:AJ:80:MET:HB3	1:AN:72:MET:HE3	1.93	0.49
1:AW:58:ILE:HG23	1:AY:43:THR:HG21	1.92	0.49
1:BE:12:SER:HB3	1:BG:12:SER:HB2	1.93	0.49
1:BH:108:PHE:HE1	1:CE:7:MET:HG2	1.78	0.49
1:BE:74:VAL:HG13	1:CG:76:PHE:CE1	2.48	0.49
1:CA:89:VAL:HG12	1:CF:74:VAL:HB	1.95	0.49
1:CF:12:SER:HB3	1:CH:12:SER:HB2	1.94	0.49
1:AH:105:ASN:C	1:BZ:46:ILE:HD13	2.38	0.48
1:AK:103:ASN:OD1	1:AK:110:LYS:HD2	2.13	0.48
1:AZ:105:ASN:HB3	1:BB:80:MET:CE	2.43	0.48
1:AN:103:ASN:OD1	1:AN:110:LYS:HD2	2.13	0.48
1:AR:111:ASN:HA	1:BS:15:GLN:NE2	2.24	0.48
1:AW:103:ASN:OD1	1:AW:110:LYS:HD2	2.13	0.48
1:AZ:46:ILE:HD11	1:BB:107:ALA:HB2	1.94	0.48
1:BR:103:ASN:OD1	1:BR:110:LYS:HD2	2.13	0.48
1:AI:70:VAL:CG1	1:BM:81:PRO:HG3	2.43	0.48
1:BI:103:ASN:OD1	1:BI:110:LYS:HD2	2.13	0.48
1:BU:103:ASN:OD1	1:BU:110:LYS:HD2	2.14	0.48
1:CD:103:ASN:OD1	1:CD:110:LYS:HD2	2.13	0.48
1:CG:103:ASN:OD1	1:CG:110:LYS:HD2	2.13	0.48
1:AL:106:SER:N	1:BP:46:ILE:HD13	2.29	0.48
1:AP:80:MET:CE	1:BL:105:ASN:HB3	2.43	0.48
1:AU:86:PRO:CA	1:BV:72:MET:HE1	2.44	0.48
1:AW:72:MET:HB2	1:AY:80:MET:HB3	1.94	0.48
1:BE:79:SER:O	1:CG:72:MET:HA	2.14	0.48
1:BO:103:ASN:OD1	1:BO:110:LYS:HD2	2.13	0.48
1:AZ:103:ASN:OD1	1:AZ:110:LYS:HD2	2.13	0.48
1:BH:80:MET:HB3	1:CD:72:MET:HB2	1.94	0.48
1:BN:76:PHE:CE2	1:BU:74:VAL:HG13	2.49	0.48
1:AC:31:LYS:NZ	1:BG:103:ASN:O	2.41	0.48
1:AS:80:MET:HB3	1:BF:72:MET:HE3	1.95	0.48
1:BC:103:ASN:OD1	1:BC:110:LYS:HD2	2.13	0.48
1:BE:27:LEU:HD21	1:CG:96:ALA:CB	2.36	0.48
1:BF:103:ASN:OD1	1:BF:110:LYS:HD2	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BX:103:ASN:OD1	1:BX:110:LYS:HD2	2.14	0.48
1:AG:76:PHE:CE2	1:AT:74:VAL:HG13	2.49	0.48
1:AS:31:LYS:NZ	1:BF:106:SER:OG	2.46	0.48
1:AS:80:MET:CE	1:BF:105:ASN:HB3	2.41	0.48
1:AT:60:THR:HG21	1:AT:67:GLY:HA3	1.96	0.48
1:AT:103:ASN:OD1	1:AT:110:LYS:HD2	2.14	0.48
1:AQ:60:THR:HG21	1:AQ:67:GLY:HA3	1.96	0.48
1:AH:60:THR:HG21	1:AH:67:GLY:HA3	1.96	0.48
1:BH:108:PHE:CE1	1:CE:7:MET:HG2	2.48	0.48
1:AA:74:VAL:HG13	1:AK:76:PHE:CE1	2.49	0.47
1:AB:60:THR:HG21	1:AB:67:GLY:HA3	1.96	0.47
1:AH:74:VAL:HG11	1:BZ:90:LEU:HD23	1.96	0.47
1:AK:60:THR:HG21	1:AK:67:GLY:HA3	1.96	0.47
1:BC:60:THR:HG21	1:BC:67:GLY:HA3	1.96	0.47
1:BL:60:THR:HG21	1:BL:67:GLY:HA3	1.96	0.47
1:AA:62:ASP:OD1	1:AL:66:CYS:SG	2.72	0.47
1:AB:103:ASN:OD1	1:AB:110:LYS:HD2	2.13	0.47
1:AE:60:THR:HG21	1:AE:67:GLY:HA3	1.96	0.47
1:AE:103:ASN:OD1	1:AE:110:LYS:HD2	2.14	0.47
1:BH:105:ASN:ND2	1:CD:84:ALA:O	2.46	0.47
1:BL:103:ASN:OD1	1:BL:110:LYS:HD2	2.13	0.47
1:CA:103:ASN:OD1	1:CA:110:LYS:HD2	2.13	0.47
1:AL:100:PHE:CE2	1:BP:29:THR:HB	2.49	0.47
1:AS:103:ASN:O	1:BF:31:LYS:NZ	2.43	0.47
1:AS:112:VAL:H	1:BF:15:GLN:HE22	1.61	0.47
1:BF:60:THR:HG21	1:BF:67:GLY:HA3	1.96	0.47
1:AL:104:LEU:HD22	1:BP:78:SER:OG	2.15	0.47
1:AW:60:THR:HG21	1:AW:67:GLY:HA3	1.96	0.47
1:AZ:15:GLN:HE22	1:BB:112:VAL:H	1.63	0.47
1:AQ:72:MET:HB2	1:AQ:72:MET:HE3	1.71	0.47
1:AQ:103:ASN:OD1	1:AQ:110:LYS:HD2	2.14	0.47
1:CD:60:THR:HG21	1:CD:67:GLY:HA3	1.96	0.47
1:AK:72:MET:HE3	1:AK:72:MET:HB2	1.71	0.47
1:AL:105:ASN:C	1:BP:46:ILE:HD13	2.40	0.47
1:AN:72:MET:HE3	1:AN:72:MET:HB2	1.71	0.47
1:AR:78:SER:OG	1:BS:104:LEU:HD22	2.15	0.47
1:AX:72:MET:HB2	1:AX:72:MET:HE3	1.51	0.47
1:AX:72:MET:HE1	1:BY:86:PRO:CA	2.45	0.47
1:BH:74:VAL:HG11	1:CD:90:LEU:HD23	1.97	0.47
1:BX:60:THR:HG21	1:BX:67:GLY:HA3	1.96	0.47
1:CA:60:THR:HG21	1:CA:67:GLY:HA3	1.96	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BS:72:MET:HB2	1:BS:72:MET:HE3	1.51	0.47
1:AM:107:ALA:HB1	1:BI:33:ILE:HD11	1.97	0.47
1:CA:104:LEU:HD23	1:CF:48:VAL:HG21	1.95	0.47
1:AH:46:ILE:HD11	1:BZ:107:ALA:HB2	1.96	0.47
1:AR:72:MET:HE3	1:AR:72:MET:HB2	1.51	0.47
1:BF:111:ASN:HD21	1:BV:111:ASN:ND2	2.13	0.47
1:BO:60:THR:HG21	1:BO:67:GLY:HA3	1.96	0.47
1:BU:60:THR:HG21	1:BU:67:GLY:HA3	1.96	0.47
1:AH:72:MET:HB2	1:BZ:80:MET:HB3	1.97	0.46
1:AZ:99:LEU:HD23	1:AZ:99:LEU:HA	1.81	0.46
1:CG:60:THR:HG21	1:CG:67:GLY:HA3	1.96	0.46
1:AL:78:SER:OG	1:BP:104:LEU:HD22	2.15	0.46
1:AN:60:THR:HG21	1:AN:67:GLY:HA3	1.96	0.46
1:AP:59:ASN:HB3	1:AP:62:ASP:CG	2.41	0.46
1:AZ:60:THR:HG21	1:AZ:67:GLY:HA3	1.96	0.46
1:AZ:105:ASN:C	1:BB:46:ILE:HD13	2.40	0.46
1:BD:86:PRO:CA	1:CH:72:MET:HE1	2.46	0.46
1:AH:103:ASN:OD1	1:AH:110:LYS:HD2	2.14	0.46
1:AS:59:ASN:HB3	1:AS:62:ASP:CG	2.41	0.46
1:BI:60:THR:HG21	1:BI:67:GLY:HA3	1.96	0.46
1:BZ:59:ASN:HB3	1:BZ:62:ASP:CG	2.41	0.46
1:BV:72:MET:HE3	1:BV:72:MET:HB2	1.51	0.46
1:AS:74:VAL:HB	1:BF:89:VAL:HG12	1.98	0.46
1:AV:59:ASN:HB3	1:AV:62:ASP:CG	2.41	0.46
1:AW:84:ALA:O	1:AY:105:ASN:ND2	2.44	0.46
1:AY:59:ASN:HB3	1:AY:62:ASP:CG	2.41	0.46
1:BE:59:ASN:HB3	1:BE:62:ASP:CG	2.41	0.46
1:BR:60:THR:HG21	1:BR:67:GLY:HA3	1.96	0.46
1:AD:59:ASN:HB3	1:AD:62:ASP:CG	2.41	0.46
1:AV:66:CYS:N	1:BY:61:CYS:SG	2.88	0.46
1:BT:59:ASN:HB3	1:BT:62:ASP:CG	2.41	0.46
1:CC:59:ASN:HB3	1:CC:62:ASP:CG	2.41	0.46
1:AG:59:ASN:HB3	1:AG:62:ASP:CG	2.41	0.46
1:AH:104:LEU:HD23	1:BZ:48:VAL:HG21	1.98	0.46
1:AJ:59:ASN:HB3	1:AJ:62:ASP:CG	2.41	0.46
1:AM:59:ASN:HB3	1:AM:62:ASP:CG	2.41	0.46
1:AS:48:VAL:HG21	1:BF:104:LEU:HD23	1.97	0.46
1:CH:72:MET:HB2	1:CH:72:MET:HE3	1.51	0.46
1:AA:107:ALA:HB1	1:AK:33:ILE:HD11	1.98	0.46
1:AI:70:VAL:HG13	1:BM:81:PRO:HG3	1.98	0.46
1:AM:19:ARG:HB3	1:AM:22:LYS:O	2.16	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AV:19:ARG:HB3	1:AV:22:LYS:O	2.16	0.46
1:AZ:72:MET:HE3	1:AZ:72:MET:HB2	1.71	0.46
1:BR:72:MET:HE3	1:BR:72:MET:HB2	1.71	0.46
1:AL:6:LYS:HD2	1:BP:112:VAL:O	2.16	0.46
1:AT:111:ASN:HD21	1:BM:111:ASN:ND2	2.14	0.46
1:BK:19:ARG:HB3	1:BK:22:LYS:O	2.16	0.46
1:AA:19:ARG:HB3	1:AA:22:LYS:O	2.16	0.46
1:AM:80:MET:CE	1:BI:105:ASN:HB3	2.42	0.46
1:AX:33:ILE:HD11	1:BY:107:ALA:HB1	1.97	0.46
1:AX:111:ASN:HA	1:BY:15:GLN:NE2	2.30	0.46
1:BT:19:ARG:HB3	1:BT:22:LYS:O	2.16	0.46
1:BW:59:ASN:HB3	1:BW:62:ASP:CG	2.41	0.46
1:CD:72:MET:HE3	1:CD:72:MET:HB2	1.71	0.46
1:AA:59:ASN:HB3	1:AA:62:ASP:CG	2.41	0.45
1:AJ:19:ARG:HB3	1:AJ:22:LYS:O	2.16	0.45
1:BB:59:ASN:HB3	1:BB:62:ASP:CG	2.41	0.45
1:BH:19:ARG:HB3	1:BH:22:LYS:O	2.16	0.45
1:BH:59:ASN:HB3	1:BH:62:ASP:CG	2.41	0.45
1:BK:59:ASN:HB3	1:BK:62:ASP:CG	2.41	0.45
1:BN:19:ARG:HB3	1:BN:22:LYS:O	2.16	0.45
1:CF:19:ARG:HB3	1:CF:22:LYS:O	2.16	0.45
1:AX:15:GLN:HE22	1:BY:112:VAL:H	1.64	0.45
1:BN:59:ASN:HB3	1:BN:62:ASP:CG	2.41	0.45
1:BW:19:ARG:HB3	1:BW:22:LYS:O	2.16	0.45
1:AD:19:ARG:HB3	1:AD:22:LYS:O	2.16	0.45
1:AO:72:MET:HE3	1:AO:72:MET:HB2	1.51	0.45
1:AZ:33:ILE:HD11	1:BB:107:ALA:HB1	1.99	0.45
1:AZ:78:SER:OG	1:BB:104:LEU:HD22	2.17	0.45
1:BE:19:ARG:HB3	1:BE:22:LYS:O	2.16	0.45
1:CF:59:ASN:HB3	1:CF:62:ASP:CG	2.41	0.45
1:AI:68:PRO:HD3	1:BM:38:ASN:ND2	2.32	0.45
1:AL:107:ALA:HB1	1:BP:33:ILE:CD1	2.47	0.45
1:AX:46:ILE:HD13	1:BY:105:ASN:C	2.42	0.45
1:BO:72:MET:HB2	1:BO:72:MET:HE3	1.71	0.45
1:BQ:59:ASN:HB3	1:BQ:62:ASP:CG	2.41	0.45
1:CC:19:ARG:HB3	1:CC:22:LYS:O	2.16	0.45
1:BB:19:ARG:HB3	1:BB:22:LYS:O	2.16	0.45
1:BL:91:LYS:CD	1:BL:118:ALA:HB2	2.47	0.45
1:BR:90:LEU:HD23	1:BT:74:VAL:HG11	1.97	0.45
1:AG:19:ARG:HB3	1:AG:22:LYS:O	2.16	0.45
1:AG:27:LEU:HD12	1:AG:51:ASN:O	2.17	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AM:40:SER:OG	1:AN:68:PRO:HA	2.17	0.45
1:AR:108:PHE:CE1	1:BQ:5:ILE:HB	2.50	0.45
1:AS:40:SER:OG	1:AT:68:PRO:HA	2.17	0.45
1:AX:29:THR:HB	1:BY:100:PHE:CE2	2.52	0.45
1:BN:40:SER:OG	1:BO:68:PRO:HA	2.17	0.45
1:AG:40:SER:OG	1:AH:68:PRO:HA	2.17	0.45
1:AI:7:MET:HE3	1:BZ:108:PHE:HD1	1.82	0.45
1:AI:61:CYS:SG	1:BK:66:CYS:N	2.90	0.45
1:AJ:40:SER:OG	1:AK:68:PRO:HA	2.17	0.45
1:AP:43:THR:HG21	1:BL:58:ILE:HG23	1.99	0.45
1:BH:27:LEU:HD12	1:BH:51:ASN:O	2.17	0.45
1:BQ:19:ARG:HB3	1:BQ:22:LYS:O	2.16	0.45
1:BR:91:LYS:CD	1:BR:118:ALA:HB2	2.47	0.45
1:BT:27:LEU:HD12	1:BT:51:ASN:O	2.17	0.45
1:BZ:27:LEU:HD12	1:BZ:51:ASN:O	2.17	0.45
1:CC:27:LEU:HD12	1:CC:51:ASN:O	2.17	0.45
1:AC:60:THR:HG21	1:BG:38:ASN:HD22	1.81	0.45
1:AE:91:LYS:CD	1:AE:118:ALA:HB2	2.47	0.45
1:AH:91:LYS:CD	1:AH:118:ALA:HB2	2.47	0.45
1:AJ:108:PHE:CE1	1:AO:7:MET:HG2	2.52	0.45
1:AV:27:LEU:HD12	1:AV:51:ASN:O	2.17	0.45
1:AY:19:ARG:HB3	1:AY:22:LYS:O	2.16	0.45
1:BK:27:LEU:HD12	1:BK:51:ASN:O	2.17	0.45
1:BN:80:MET:HB3	1:BU:72:MET:HB2	1.97	0.45
1:BZ:19:ARG:HB3	1:BZ:22:LYS:O	2.16	0.45
1:CC:40:SER:OG	1:CD:68:PRO:HA	2.17	0.45
1:AA:27:LEU:HD12	1:AA:51:ASN:O	2.17	0.44
1:AQ:91:LYS:CD	1:AQ:118:ALA:HB2	2.47	0.44
1:AT:91:LYS:CD	1:AT:118:ALA:HB2	2.47	0.44
1:BL:99:LEU:HD23	1:BL:99:LEU:HA	1.81	0.44
1:CD:91:LYS:CD	1:CD:118:ALA:HB2	2.47	0.44
1:AA:108:PHE:CD1	1:AL:7:MET:HG2	2.52	0.44
1:AK:91:LYS:CD	1:AK:118:ALA:HB2	2.47	0.44
1:AP:19:ARG:HB3	1:AP:22:LYS:O	2.16	0.44
1:AP:76:PHE:CE2	1:BL:74:VAL:HG13	2.53	0.44
1:AS:19:ARG:HB3	1:AS:22:LYS:O	2.16	0.44
1:BF:91:LYS:CD	1:BF:118:ALA:HB2	2.47	0.44
1:BK:40:SER:OG	1:BL:68:PRO:HA	2.17	0.44
1:CA:91:LYS:CD	1:CA:118:ALA:HB2	2.47	0.44
1:CB:72:MET:HB2	1:CB:72:MET:HE3	1.51	0.44
1:AC:103:ASN:O	1:BG:31:LYS:NZ	2.47	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AE:72:MET:HB2	1:AE:72:MET:HE3	1.71	0.44
1:AE:104:LEU:HD12	1:BW:90:LEU:HD11	2.00	0.44
1:BB:40:SER:OG	1:BC:68:PRO:HA	2.17	0.44
1:BC:72:MET:HE3	1:BC:72:MET:HB2	1.71	0.44
1:BE:46:ILE:HD13	1:CG:105:ASN:C	2.42	0.44
1:BT:40:SER:OG	1:BU:68:PRO:HA	2.17	0.44
1:BU:91:LYS:CD	1:BU:118:ALA:HB2	2.47	0.44
1:BW:40:SER:OG	1:BX:68:PRO:HA	2.17	0.44
1:BX:91:LYS:CD	1:BX:118:ALA:HB2	2.47	0.44
1:AB:91:LYS:CD	1:AB:118:ALA:HB2	2.47	0.44
1:AC:66:CYS:SG	1:CC:62:ASP:OD1	2.76	0.44
1:AD:46:ILE:HD12	1:AD:82:GLU:HA	2.00	0.44
1:AN:27:LEU:HA	1:AN:51:ASN:O	2.18	0.44
1:AN:91:LYS:CD	1:AN:118:ALA:HB2	2.47	0.44
1:AR:72:MET:HE1	1:BS:86:PRO:CA	2.48	0.44
1:AS:46:ILE:HD13	1:BF:105:ASN:C	2.43	0.44
1:AZ:111:ASN:HA	1:BB:15:GLN:OE1	2.17	0.44
1:BN:59:ASN:OD1	1:BU:41:THR:HG21	2.16	0.44
1:AH:27:LEU:HA	1:AH:51:ASN:O	2.18	0.44
1:AK:99:LEU:HD23	1:AK:99:LEU:HA	1.81	0.44
1:AM:76:PHE:CE2	1:BI:74:VAL:HG13	2.53	0.44
1:AT:27:LEU:HA	1:AT:51:ASN:O	2.18	0.44
1:AU:72:MET:HE1	1:BV:86:PRO:CA	2.48	0.44
1:AZ:91:LYS:CD	1:AZ:118:ALA:HB2	2.47	0.44
1:BN:90:LEU:HD23	1:BU:74:VAL:HG11	1.98	0.44
1:BZ:40:SER:OG	1:CA:68:PRO:HA	2.17	0.44
1:AK:27:LEU:HA	1:AK:51:ASN:O	2.18	0.44
1:AM:7:MET:SD	1:AM:18:ALA:HB2	2.58	0.44
1:AM:27:LEU:HD12	1:AM:51:ASN:O	2.17	0.44
1:AP:27:LEU:HD12	1:AP:51:ASN:O	2.17	0.44
1:AS:46:ILE:HD12	1:AS:82:GLU:HA	2.00	0.44
1:AV:40:SER:OG	1:AW:68:PRO:HA	2.17	0.44
1:AY:27:LEU:HD12	1:AY:51:ASN:O	2.17	0.44
1:AY:40:SER:OG	1:AZ:68:PRO:HA	2.17	0.44
1:BG:72:MET:HB2	1:BG:72:MET:HE3	1.51	0.44
1:BM:72:MET:HE3	1:BM:72:MET:HB2	1.51	0.44
1:BO:91:LYS:CD	1:BO:118:ALA:HB2	2.47	0.44
1:CA:72:MET:HE3	1:CA:72:MET:HB2	1.71	0.44
1:CG:91:LYS:CD	1:CG:118:ALA:HB2	2.47	0.44
1:AA:104:LEU:HD23	1:AK:48:VAL:HG21	2.00	0.44
1:AC:104:LEU:HD22	1:BG:78:SER:OG	2.17	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AN:99:LEU:HD23	1:AN:99:LEU:HA	1.81	0.44
1:AS:27:LEU:HD12	1:AS:51:ASN:O	2.17	0.44
1:AY:7:MET:SD	1:AY:18:ALA:HB2	2.58	0.44
1:BB:7:MET:SD	1:BB:18:ALA:HB2	2.58	0.44
1:BB:27:LEU:HD12	1:BB:51:ASN:O	2.17	0.44
1:BH:46:ILE:HD12	1:BH:82:GLU:HA	2.00	0.44
1:BZ:7:MET:SD	1:BZ:18:ALA:HB2	2.58	0.44
1:CC:7:MET:SD	1:CC:18:ALA:HB2	2.58	0.44
1:CD:27:LEU:HA	1:CD:51:ASN:O	2.18	0.44
1:CG:27:LEU:HA	1:CG:51:ASN:O	2.18	0.44
1:AB:105:ASN:HB3	1:CC:80:MET:CE	2.44	0.44
1:AD:40:SER:OG	1:AE:68:PRO:HA	2.17	0.44
1:AP:7:MET:SD	1:AP:18:ALA:HB2	2.58	0.44
1:AR:27:LEU:HA	1:AR:51:ASN:O	2.18	0.44
1:AS:111:ASN:HA	1:BG:8:ALA:HB2	1.99	0.44
1:AU:27:LEU:HA	1:AU:51:ASN:O	2.18	0.44
1:BA:27:LEU:HA	1:BA:51:ASN:O	2.18	0.44
1:BH:48:VAL:HG21	1:CD:104:LEU:HD23	2.00	0.44
1:BI:91:LYS:CD	1:BI:118:ALA:HB2	2.47	0.44
1:BN:27:LEU:HD12	1:BN:51:ASN:O	2.17	0.44
1:BQ:7:MET:SD	1:BQ:18:ALA:HB2	2.58	0.44
1:BW:46:ILE:HD12	1:BW:82:GLU:HA	2.00	0.44
1:AA:7:MET:SD	1:AA:18:ALA:HB2	2.58	0.44
1:AH:78:SER:OG	1:BZ:104:LEU:HD22	2.18	0.44
1:AH:84:ALA:O	1:BZ:105:ASN:ND2	2.51	0.44
1:AJ:27:LEU:HD12	1:AJ:51:ASN:O	2.17	0.44
1:AP:40:SER:OG	1:AQ:68:PRO:HA	2.17	0.44
1:BD:27:LEU:HA	1:BD:51:ASN:O	2.18	0.44
1:BE:40:SER:OG	1:BF:68:PRO:HA	2.17	0.44
1:BH:6:LYS:NZ	1:CD:111:ASN:OD1	2.50	0.44
1:BJ:27:LEU:HA	1:BJ:51:ASN:O	2.18	0.44
1:BT:46:ILE:HD12	1:BT:82:GLU:HA	2.00	0.44
1:BV:27:LEU:HA	1:BV:51:ASN:O	2.18	0.44
1:BW:7:MET:SD	1:BW:18:ALA:HB2	2.58	0.44
1:CA:31:LYS:NZ	1:CF:103:ASN:O	2.49	0.44
1:CF:27:LEU:HD12	1:CF:51:ASN:O	2.17	0.44
1:CF:40:SER:OG	1:CG:68:PRO:HA	2.17	0.44
1:AC:27:LEU:HA	1:AC:51:ASN:O	2.18	0.43
1:AD:7:MET:SD	1:AD:18:ALA:HB2	2.58	0.43
1:AV:7:MET:SD	1:AV:18:ALA:HB2	2.58	0.43
1:AX:27:LEU:HA	1:AX:51:ASN:O	2.18	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AY:46:ILE:HD12	1:AY:82:GLU:HA	2.00	0.43
1:BK:63:PRO:HB2	1:BK:65:LYS:NZ	2.33	0.43
1:BN:46:ILE:HD12	1:BN:82:GLU:HA	2.00	0.43
1:BP:27:LEU:HA	1:BP:51:ASN:O	2.18	0.43
1:BT:7:MET:SD	1:BT:18:ALA:HB2	2.58	0.43
1:BY:27:LEU:HA	1:BY:51:ASN:O	2.18	0.43
1:CC:46:ILE:HD12	1:CC:82:GLU:HA	2.00	0.43
1:CF:46:ILE:HD12	1:CF:82:GLU:HA	2.00	0.43
1:CH:27:LEU:HA	1:CH:51:ASN:O	2.18	0.43
1:AA:40:SER:OG	1:AB:68:PRO:HA	2.17	0.43
1:AA:72:MET:HG3	1:AK:80:MET:HB3	2.00	0.43
1:AG:7:MET:SD	1:AG:18:ALA:HB2	2.58	0.43
1:AI:60:THR:HB	1:BM:41:THR:HG21	2.00	0.43
1:AJ:63:PRO:HB2	1:AJ:65:LYS:NZ	2.33	0.43
1:AJ:90:LEU:HD23	1:AN:74:VAL:HG11	1.99	0.43
1:AP:80:MET:HB3	1:BL:72:MET:HB2	2.00	0.43
1:AT:99:LEU:HD23	1:AT:99:LEU:HA	1.81	0.43
1:BC:99:LEU:HD23	1:BC:99:LEU:HA	1.81	0.43
1:BE:63:PRO:HB2	1:BE:65:LYS:NZ	2.33	0.43
1:BH:31:LYS:NZ	1:CD:106:SER:OG	2.51	0.43
1:BH:40:SER:OG	1:BI:68:PRO:HA	2.17	0.43
1:BM:37:ARG:HG2	1:BM:42:SER:HB3	2.01	0.43
1:BQ:27:LEU:HD12	1:BQ:51:ASN:O	2.17	0.43
1:BS:27:LEU:HA	1:BS:51:ASN:O	2.18	0.43
1:BY:37:ARG:HG2	1:BY:42:SER:HB3	2.01	0.43
1:CA:27:LEU:HA	1:CA:51:ASN:O	2.18	0.43
1:AD:27:LEU:HD12	1:AD:51:ASN:O	2.17	0.43
1:AE:27:LEU:HA	1:AE:51:ASN:O	2.18	0.43
1:AI:37:ARG:HG2	1:AI:42:SER:HB3	2.01	0.43
1:AW:104:LEU:HD12	1:AY:90:LEU:HD11	2.00	0.43
1:BE:27:LEU:HD12	1:BE:51:ASN:O	2.17	0.43
1:BM:27:LEU:HA	1:BM:51:ASN:O	2.18	0.43
1:BN:7:MET:SD	1:BN:18:ALA:HB2	2.58	0.43
1:BO:27:LEU:HA	1:BO:51:ASN:O	2.18	0.43
1:BR:27:LEU:HA	1:BR:51:ASN:O	2.18	0.43
1:CB:27:LEU:HA	1:CB:51:ASN:O	2.18	0.43
1:AA:90:LEU:HD23	1:AK:74:VAL:HG11	1.99	0.43
1:AB:27:LEU:HA	1:AB:51:ASN:O	2.18	0.43
1:AB:99:LEU:HD23	1:AB:99:LEU:HA	1.81	0.43
1:AC:78:SER:OG	1:BG:104:LEU:HD22	2.18	0.43
1:AF:37:ARG:HG2	1:AF:42:SER:HB3	2.01	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AG:63:PRO:HB2	1:AG:65:LYS:NZ	2.33	0.43
1:AI:27:LEU:HA	1:AI:51:ASN:O	2.18	0.43
1:AJ:80:MET:HB3	1:AN:72:MET:HB2	1.99	0.43
1:AU:37:ARG:HG2	1:AU:42:SER:HB3	2.01	0.43
1:AW:91:LYS:CD	1:AW:118:ALA:HB2	2.47	0.43
1:AX:37:ARG:HG2	1:AX:42:SER:HB3	2.01	0.43
1:BC:91:LYS:CD	1:BC:118:ALA:HB2	2.47	0.43
1:BE:46:ILE:HD12	1:BE:82:GLU:HA	2.00	0.43
1:BG:27:LEU:HA	1:BG:51:ASN:O	2.18	0.43
1:BQ:40:SER:OG	1:BR:68:PRO:HA	2.17	0.43
1:BR:104:LEU:HD12	1:BT:90:LEU:HD11	2.00	0.43
1:BW:63:PRO:HB2	1:BW:65:LYS:NZ	2.33	0.43
1:CD:99:LEU:HD23	1:CD:99:LEU:HA	1.81	0.43
1:AJ:46:ILE:HD12	1:AJ:82:GLU:HA	2.00	0.43
1:AM:43:THR:HG21	1:BI:58:ILE:HG23	2.00	0.43
1:BA:37:ARG:HG2	1:BA:42:SER:HB3	2.01	0.43
1:BC:27:LEU:HA	1:BC:51:ASN:O	2.18	0.43
1:BI:27:LEU:HA	1:BI:51:ASN:O	2.18	0.43
1:BZ:46:ILE:HD12	1:BZ:82:GLU:HA	2.00	0.43
1:AH:112:VAL:HG23	1:BZ:17:TYR:OH	2.18	0.43
1:AJ:58:ILE:HG23	1:AN:43:THR:OG1	2.19	0.43
1:AM:46:ILE:HD12	1:AM:82:GLU:HA	2.00	0.43
1:AM:63:PRO:HB2	1:AM:65:LYS:NZ	2.33	0.43
1:AO:27:LEU:HA	1:AO:51:ASN:O	2.18	0.43
1:AT:72:MET:HE3	1:AT:72:MET:HB2	1.71	0.43
1:BD:72:MET:HE1	1:CH:86:PRO:CA	2.48	0.43
1:BW:27:LEU:HD12	1:BW:51:ASN:O	2.17	0.43
1:CB:37:ARG:HG2	1:CB:42:SER:HB3	2.01	0.43
1:CE:27:LEU:HA	1:CE:51:ASN:O	2.18	0.43
1:AJ:76:PHE:CE2	1:AN:74:VAL:HG13	2.54	0.43
1:AL:72:MET:HE3	1:AL:72:MET:HB2	1.51	0.43
1:AM:80:MET:HB3	1:BI:72:MET:HB2	2.00	0.43
1:AM:105:ASN:ND2	1:BI:84:ALA:O	2.49	0.43
1:AS:63:PRO:HB2	1:AS:65:LYS:NZ	2.33	0.43
1:BD:37:ARG:HG2	1:BD:42:SER:HB3	2.01	0.43
1:BF:27:LEU:HA	1:BF:51:ASN:O	2.18	0.43
1:BG:37:ARG:HG2	1:BG:42:SER:HB3	2.01	0.43
1:BH:76:PHE:CE2	1:CD:74:VAL:HG13	2.54	0.43
1:BI:72:MET:HE3	1:BI:72:MET:HB2	1.71	0.43
1:BN:90:LEU:HD11	1:BU:104:LEU:HD12	2.00	0.43
1:BQ:46:ILE:HD12	1:BQ:82:GLU:HA	2.00	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BR:61:CYS:HB2	1:BU:66:CYS:HB2	1.89	0.43
1:BS:66:CYS:SG	1:BT:62:ASP:OD1	2.77	0.43
1:BX:99:LEU:HA	1:BX:99:LEU:HD23	1.81	0.43
1:CF:63:PRO:HB2	1:CF:65:LYS:NZ	2.33	0.43
1:CH:37:ARG:HG2	1:CH:42:SER:HB3	2.01	0.43
1:AA:66:CYS:N	1:BG:61:CYS:SG	2.92	0.43
1:AV:63:PRO:HB2	1:AV:65:LYS:NZ	2.34	0.43
1:AW:27:LEU:HA	1:AW:51:ASN:O	2.18	0.43
1:AY:63:PRO:HB2	1:AY:65:LYS:NZ	2.33	0.43
1:AZ:41:THR:HG21	1:BB:59:ASN:OD1	2.19	0.43
1:BE:7:MET:SD	1:BE:18:ALA:HB2	2.58	0.43
1:BH:7:MET:SD	1:BH:18:ALA:HB2	2.58	0.43
1:BQ:63:PRO:HB2	1:BQ:65:LYS:NZ	2.33	0.43
1:BS:37:ARG:HG2	1:BS:42:SER:HB3	2.01	0.43
1:BU:27:LEU:HA	1:BU:51:ASN:O	2.18	0.43
1:BY:72:MET:HB2	1:BY:72:MET:HE3	1.51	0.43
1:AA:46:ILE:HD12	1:AA:82:GLU:HA	2.00	0.43
1:AI:58:ILE:HG21	1:BM:43:THR:OG1	2.19	0.43
1:AL:27:LEU:HA	1:AL:51:ASN:O	2.18	0.43
1:AS:7:MET:SD	1:AS:18:ALA:HB2	2.58	0.43
1:AV:46:ILE:HD12	1:AV:82:GLU:HA	2.00	0.43
1:AX:31:LYS:NZ	1:BY:103:ASN:O	2.37	0.43
1:BD:99:LEU:HD23	1:BD:99:LEU:HA	1.87	0.43
1:BK:7:MET:SD	1:BK:18:ALA:HB2	2.58	0.43
1:BX:27:LEU:HA	1:BX:51:ASN:O	2.18	0.43
1:CE:72:MET:HB2	1:CE:72:MET:HE3	1.51	0.43
1:AJ:7:MET:SD	1:AJ:18:ALA:HB2	2.58	0.43
1:AL:31:LYS:NZ	1:BP:103:ASN:O	2.37	0.43
1:AP:46:ILE:HD12	1:AP:82:GLU:HA	2.00	0.43
1:AQ:27:LEU:HA	1:AQ:51:ASN:O	2.18	0.43
1:AZ:74:VAL:HG13	1:BB:76:PHE:CE2	2.54	0.43
1:AZ:104:LEU:HD23	1:BB:48:VAL:HG21	2.01	0.43
1:CC:63:PRO:HB2	1:CC:65:LYS:NZ	2.33	0.43
1:AC:46:ILE:HD13	1:BG:106:SER:N	2.33	0.42
1:AP:108:PHE:CE1	1:BM:7:MET:HG2	2.54	0.42
1:AP:114:GLU:OE1	1:BL:3:LYS:HD3	2.19	0.42
1:AR:37:ARG:HG2	1:AR:42:SER:HB3	2.01	0.42
1:AZ:27:LEU:HA	1:AZ:51:ASN:O	2.18	0.42
1:BB:46:ILE:HD12	1:BB:82:GLU:HA	2.00	0.42
1:BB:63:PRO:HB2	1:BB:65:LYS:NZ	2.34	0.42
1:BH:63:PRO:HB2	1:BH:65:LYS:NZ	2.33	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BZ:63:PRO:HB2	1:BZ:65:LYS:NZ	2.33	0.42
1:CF:7:MET:SD	1:CF:18:ALA:HB2	2.58	0.42
1:AD:63:PRO:HB2	1:AD:65:LYS:NZ	2.34	0.42
1:AF:78:SER:OG	1:BJ:104:LEU:HD22	2.19	0.42
1:AP:63:PRO:HB2	1:AP:65:LYS:NZ	2.33	0.42
1:AU:81:PRO:HG3	1:BV:70:VAL:CG1	2.49	0.42
1:BA:72:MET:HE3	1:BA:72:MET:HB2	1.51	0.42
1:BD:108:PHE:CE1	1:CF:5:ILE:HB	2.55	0.42
1:BK:46:ILE:HD12	1:BK:82:GLU:HA	2.00	0.42
1:CE:36:LEU:HD22	1:CE:45:ILE:HD11	2.02	0.42
1:CE:37:ARG:HG2	1:CE:42:SER:HB3	2.01	0.42
1:CG:99:LEU:HA	1:CG:99:LEU:HD23	1.81	0.42
1:AU:36:LEU:HD22	1:AU:45:ILE:HD11	2.02	0.42
1:BJ:37:ARG:HG2	1:BJ:42:SER:HB3	2.01	0.42
1:BS:36:LEU:HD22	1:BS:45:ILE:HD11	2.02	0.42
1:BV:37:ARG:HG2	1:BV:42:SER:HB3	2.01	0.42
1:AA:63:PRO:HB2	1:AA:65:LYS:NZ	2.33	0.42
1:AC:36:LEU:HD22	1:AC:45:ILE:HD11	2.02	0.42
1:AH:48:VAL:HG21	1:BZ:104:LEU:HD23	2.00	0.42
1:AJ:74:VAL:HB	1:AN:89:VAL:HG12	2.02	0.42
1:AM:31:LYS:NZ	1:BI:106:SER:OG	2.52	0.42
1:AR:15:GLN:NE2	1:BS:111:ASN:HA	2.33	0.42
1:BL:72:MET:HE3	1:BL:72:MET:HB2	1.71	0.42
1:BT:63:PRO:HB2	1:BT:65:LYS:NZ	2.33	0.42
1:AI:72:MET:HE1	1:BM:86:PRO:HB3	2.00	0.42
1:AL:36:LEU:HD22	1:AL:45:ILE:HD11	2.02	0.42
1:AV:2:TYR:HE1	1:AV:4:LEU:HD21	1.85	0.42
1:BA:36:LEU:HD22	1:BA:45:ILE:HD11	2.02	0.42
1:BA:84:ALA:O	1:CE:105:ASN:ND2	2.38	0.42
1:BH:2:TYR:HE1	1:BH:4:LEU:HD21	1.85	0.42
1:BV:36:LEU:HD22	1:BV:45:ILE:HD11	2.01	0.42
1:CC:2:TYR:HE1	1:CC:4:LEU:HD21	1.85	0.42
1:AC:46:ILE:HD13	1:BG:105:ASN:C	2.44	0.42
1:AF:38:ASN:HD22	1:BJ:60:THR:HG21	1.83	0.42
1:AL:99:LEU:HD23	1:AL:99:LEU:HA	1.87	0.42
1:AR:38:ASN:HD22	1:BS:60:THR:HG21	1.84	0.42
1:AW:74:VAL:HG13	1:AY:76:PHE:CE2	2.55	0.42
1:BP:36:LEU:HD22	1:BP:45:ILE:HD11	2.02	0.42
1:BW:2:TYR:HE1	1:BW:4:LEU:HD21	1.85	0.42
1:AK:19:ARG:HB3	1:AK:22:LYS:O	2.20	0.42
1:BD:72:MET:HB2	1:BD:72:MET:HE3	1.51	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BE:86:PRO:HB3	1:CG:72:MET:HE1	2.02	0.42
1:BL:27:LEU:HA	1:BL:51:ASN:O	2.18	0.42
1:BP:37:ARG:HG2	1:BP:42:SER:HB3	2.01	0.42
1:BU:99:LEU:HD23	1:BU:99:LEU:HA	1.81	0.42
1:CA:19:ARG:HB3	1:CA:22:LYS:O	2.20	0.42
1:AC:72:MET:HE3	1:AC:72:MET:HB2	1.51	0.42
1:AD:2:TYR:HE1	1:AD:4:LEU:HD21	1.85	0.42
1:AM:2:TYR:HE1	1:AM:4:LEU:HD21	1.85	0.42
1:AZ:112:VAL:HG23	1:BB:17:TYR:OH	2.20	0.42
1:BA:8:ALA:HB2	1:BB:111:ASN:HA	2.01	0.42
1:BE:2:TYR:HE1	1:BE:4:LEU:HD21	1.85	0.42
1:BN:2:TYR:HE1	1:BN:4:LEU:HD21	1.85	0.42
1:BN:63:PRO:HB2	1:BN:65:LYS:NZ	2.33	0.42
1:BR:3:LYS:HD3	1:BT:114:GLU:OE1	2.20	0.42
1:AB:89:VAL:HG12	1:CC:74:VAL:HB	2.02	0.42
1:AC:106:SER:N	1:BG:46:ILE:HD13	2.35	0.42
1:AE:19:ARG:HB3	1:AE:22:LYS:O	2.20	0.42
1:AF:27:LEU:HA	1:AF:51:ASN:O	2.18	0.42
1:AF:36:LEU:HD22	1:AF:45:ILE:HD11	2.02	0.42
1:AG:46:ILE:HD12	1:AG:82:GLU:HA	2.00	0.42
1:AI:7:MET:HG2	1:BZ:108:PHE:CD1	2.51	0.42
1:AL:37:ARG:HG2	1:AL:42:SER:HB3	2.01	0.42
1:AN:19:ARG:HB3	1:AN:22:LYS:O	2.20	0.42
1:AO:36:LEU:HD22	1:AO:45:ILE:HD11	2.02	0.42
1:BD:81:PRO:HG3	1:CH:70:VAL:CG1	2.49	0.42
1:BR:7:MET:HE2	1:BR:7:MET:HB3	1.93	0.42
1:AP:2:TYR:HE1	1:AP:4:LEU:HD21	1.85	0.42
1:AR:36:LEU:HD22	1:AR:45:ILE:HD11	2.02	0.42
1:AS:6:LYS:HE3	1:BF:113:ASP:HA	2.02	0.42
1:AT:19:ARG:HB3	1:AT:22:LYS:O	2.20	0.42
1:AW:90:LEU:HD23	1:AY:74:VAL:HG11	2.00	0.42
1:BA:78:SER:OG	1:CE:104:LEU:HD22	2.20	0.42
1:BU:19:ARG:HB3	1:BU:22:LYS:O	2.20	0.42
1:CD:7:MET:HE2	1:CD:7:MET:HB3	1.93	0.42
1:CH:36:LEU:HD22	1:CH:45:ILE:HD11	2.02	0.42
1:AA:2:TYR:HE1	1:AA:4:LEU:HD21	1.85	0.41
1:AB:19:ARG:HB3	1:AB:22:LYS:O	2.20	0.41
1:AI:36:LEU:HD22	1:AI:45:ILE:HD11	2.02	0.41
1:AO:37:ARG:HG2	1:AO:42:SER:HB3	2.01	0.41
1:AZ:19:ARG:HB3	1:AZ:22:LYS:O	2.20	0.41
1:BL:19:ARG:HB3	1:BL:22:LYS:O	2.20	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AF:84:ALA:O	1:BJ:105:ASN:ND2	2.40	0.41
1:AH:7:MET:HE2	1:AH:7:MET:HB3	1.93	0.41
1:AJ:74:VAL:HG13	1:AN:76:PHE:CE1	2.55	0.41
1:AL:15:GLN:HE22	1:BP:111:ASN:CA	2.20	0.41
1:AW:19:ARG:HB3	1:AW:22:LYS:O	2.20	0.41
1:AX:36:LEU:HD22	1:AX:45:ILE:HD11	2.02	0.41
1:BB:2:TYR:HE1	1:BB:4:LEU:HD21	1.85	0.41
1:BK:2:TYR:HE1	1:BK:4:LEU:HD21	1.85	0.41
1:BO:19:ARG:HB3	1:BO:22:LYS:O	2.20	0.41
1:BO:99:LEU:HD23	1:BO:99:LEU:HA	1.81	0.41
1:BX:19:ARG:HB3	1:BX:22:LYS:O	2.20	0.41
1:AP:103:ASN:O	1:BL:31:LYS:NZ	2.46	0.41
1:BQ:2:TYR:HE1	1:BQ:4:LEU:HD21	1.85	0.41
1:BY:36:LEU:HD22	1:BY:45:ILE:HD11	2.02	0.41
1:CF:2:TYR:HE1	1:CF:4:LEU:HD21	1.85	0.41
1:CG:19:ARG:HB3	1:CG:22:LYS:O	2.20	0.41
1:AB:72:MET:HE3	1:AB:72:MET:HB2	1.71	0.41
1:AB:104:LEU:HD12	1:CC:90:LEU:HD11	2.03	0.41
1:AG:2:TYR:HE1	1:AG:4:LEU:HD21	1.85	0.41
1:AP:105:ASN:ND2	1:BL:84:ALA:O	2.48	0.41
1:AS:2:TYR:HE1	1:AS:4:LEU:HD21	1.85	0.41
1:BC:19:ARG:HB3	1:BC:22:LYS:O	2.20	0.41
1:BZ:2:TYR:HE1	1:BZ:4:LEU:HD21	1.85	0.41
1:CD:19:ARG:HB3	1:CD:22:LYS:O	2.20	0.41
1:AI:108:PHE:CD1	1:BK:5:ILE:HB	2.55	0.41
1:AW:74:VAL:HG11	1:AY:90:LEU:HD23	2.02	0.41
1:BD:36:LEU:HD22	1:BD:45:ILE:HD11	2.02	0.41
1:BE:48:VAL:HG21	1:CG:104:LEU:HD23	2.03	0.41
1:BE:58:ILE:HG23	1:CG:43:THR:OG1	2.21	0.41
1:BM:36:LEU:HD22	1:BM:45:ILE:HD11	2.02	0.41
1:CA:99:LEU:HA	1:CA:99:LEU:HD23	1.81	0.41
1:CA:109:ASN:HB2	1:CG:8:ALA:HB2	2.03	0.41
1:AC:37:ARG:HG2	1:AC:42:SER:HB3	2.01	0.41
1:AH:19:ARG:HB3	1:AH:22:LYS:O	2.20	0.41
1:AH:58:ILE:HG23	1:BZ:43:THR:HG21	2.03	0.41
1:AR:72:MET:HE1	1:BS:86:PRO:HB3	2.02	0.41
1:AZ:58:ILE:HD12	1:AZ:68:PRO:HB2	2.03	0.41
1:BF:19:ARG:HB3	1:BF:22:LYS:O	2.20	0.41
1:BI:19:ARG:HB3	1:BI:22:LYS:O	2.20	0.41
1:BN:62:ASP:OD1	1:BV:66:CYS:SG	2.79	0.41
1:AI:7:MET:HE2	1:AI:7:MET:HB3	1.96	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AM:90:LEU:HD11	1:BI:104:LEU:HD12	2.03	0.41
1:BG:36:LEU:HD22	1:BG:45:ILE:HD11	2.02	0.41
1:BR:58:ILE:HD12	1:BR:68:PRO:HB2	2.03	0.41
1:BT:2:TYR:HE1	1:BT:4:LEU:HD21	1.85	0.41
1:AC:38:ASN:HD22	1:BG:60:THR:HG21	1.84	0.41
1:AE:58:ILE:HD12	1:AE:68:PRO:HB2	2.03	0.41
1:AE:99:LEU:HD23	1:AE:99:LEU:HA	1.81	0.41
1:AI:72:MET:HE1	1:BM:86:PRO:CA	2.51	0.41
1:AS:105:ASN:ND2	1:BF:84:ALA:O	2.51	0.41
1:CA:7:MET:HE2	1:CA:7:MET:HB3	1.93	0.41
1:CA:105:ASN:C	1:CF:46:ILE:HD13	2.46	0.41
1:CB:36:LEU:HD22	1:CB:45:ILE:HD11	2.02	0.41
1:CG:72:MET:HE3	1:CG:72:MET:HB2	1.71	0.41
1:AB:58:ILE:O	1:AB:60:THR:HG23	2.21	0.41
1:AI:65:LYS:O	1:BZ:61:CYS:HB2	2.20	0.41
1:AK:58:ILE:HD12	1:AK:68:PRO:HB2	2.03	0.41
1:AK:58:ILE:O	1:AK:60:THR:HG23	2.21	0.41
1:AN:58:ILE:HD12	1:AN:68:PRO:HB2	2.03	0.41
1:AQ:19:ARG:HB3	1:AQ:22:LYS:O	2.20	0.41
1:AT:58:ILE:HD12	1:AT:68:PRO:HB2	2.03	0.41
1:BE:72:MET:HE3	1:CG:89:VAL:HG21	2.03	0.41
1:BF:72:MET:HE3	1:BF:72:MET:HB2	1.71	0.41
1:BJ:36:LEU:HD22	1:BJ:45:ILE:HD11	2.02	0.41
1:BL:58:ILE:HD12	1:BL:68:PRO:HB2	2.03	0.41
1:BO:58:ILE:HD12	1:BO:68:PRO:HB2	2.03	0.41
1:CA:58:ILE:HD12	1:CA:68:PRO:HB2	2.03	0.41
1:CA:61:CYS:HB2	1:CG:66:CYS:HB2	1.83	0.41
1:CD:58:ILE:HD12	1:CD:68:PRO:HB2	2.03	0.41
1:CG:7:MET:HE2	1:CG:7:MET:HB3	1.93	0.41
1:AQ:58:ILE:HD12	1:AQ:68:PRO:HB2	2.03	0.41
1:AU:86:PRO:HB3	1:BV:72:MET:HE1	2.02	0.41
1:AX:86:PRO:HB3	1:BY:72:MET:HE1	2.03	0.41
1:AZ:58:ILE:O	1:AZ:60:THR:HG23	2.21	0.41
1:AZ:72:MET:HB2	1:BB:80:MET:HB3	2.02	0.41
1:AG:90:LEU:HD23	1:AT:74:VAL:HG11	2.03	0.40
1:AH:99:LEU:HD23	1:AH:99:LEU:HA	1.81	0.40
1:AN:58:ILE:O	1:AN:60:THR:HG23	2.21	0.40
1:AU:72:MET:HE3	1:AU:72:MET:HB2	1.51	0.40
1:AX:72:MET:HE1	1:BY:86:PRO:HB3	2.03	0.40
1:BI:7:MET:HE2	1:BI:7:MET:HB3	1.93	0.40
1:BJ:72:MET:HB2	1:BJ:72:MET:HE3	1.51	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BN:43:THR:HG21	1:BU:58:ILE:HG23	2.02	0.40
1:BX:58:ILE:O	1:BX:60:THR:HG23	2.21	0.40
1:AB:37:ARG:HG2	1:AB:42:SER:HB3	2.04	0.40
1:AE:58:ILE:O	1:AE:60:THR:HG23	2.21	0.40
1:AX:6:LYS:HD2	1:BY:112:VAL:O	2.21	0.40
1:AY:2:TYR:HE1	1:AY:4:LEU:HD21	1.85	0.40
1:AZ:7:MET:HE2	1:AZ:7:MET:HB3	1.93	0.40
1:BE:104:LEU:HD23	1:CG:48:VAL:HG21	2.03	0.40
1:BI:37:ARG:HG2	1:BI:42:SER:HB3	2.04	0.40
1:BI:58:ILE:O	1:BI:60:THR:HG23	2.21	0.40
1:BR:19:ARG:HB3	1:BR:22:LYS:O	2.20	0.40
1:AH:111:ASN:HA	1:BZ:15:GLN:OE1	2.21	0.40
1:AJ:2:TYR:HE1	1:AJ:4:LEU:HD21	1.85	0.40
1:AN:37:ARG:HG2	1:AN:42:SER:HB3	2.03	0.40
1:AT:58:ILE:O	1:AT:60:THR:HG23	2.21	0.40
1:BL:37:ARG:HG2	1:BL:42:SER:HB3	2.04	0.40
1:BO:58:ILE:O	1:BO:60:THR:HG23	2.21	0.40
1:BR:37:ARG:HG2	1:BR:42:SER:HB3	2.04	0.40
1:BR:58:ILE:O	1:BR:60:THR:HG23	2.21	0.40
1:BU:37:ARG:HG2	1:BU:42:SER:HB3	2.04	0.40
1:BU:58:ILE:HD12	1:BU:68:PRO:HB2	2.03	0.40
1:BX:37:ARG:HG2	1:BX:42:SER:HB3	2.04	0.40
1:CD:37:ARG:HG2	1:CD:42:SER:HB3	2.04	0.40
1:CG:58:ILE:HD12	1:CG:68:PRO:HB2	2.03	0.40
1:AE:84:ALA:O	1:BW:105:ASN:ND2	2.51	0.40
1:AG:108:PHE:HE1	1:AU:7:MET:HG2	1.87	0.40
1:AQ:58:ILE:O	1:AQ:60:THR:HG23	2.21	0.40
1:AW:7:MET:HE2	1:AW:7:MET:HB3	1.93	0.40
1:AW:37:ARG:HG2	1:AW:42:SER:HB3	2.04	0.40
1:AW:72:MET:HE3	1:AW:72:MET:HB2	1.71	0.40
1:AZ:46:ILE:CD1	1:BB:107:ALA:HB2	2.52	0.40
1:BD:86:PRO:HB3	1:CH:72:MET:HE1	2.02	0.40
1:BF:58:ILE:O	1:BF:60:THR:HG23	2.21	0.40
1:BL:58:ILE:O	1:BL:60:THR:HG23	2.21	0.40
1:AK:37:ARG:HG2	1:AK:42:SER:HB3	2.04	0.40
1:AW:58:ILE:O	1:AW:60:THR:HG23	2.21	0.40
1:AZ:58:ILE:HG23	1:BB:43:THR:HG21	2.03	0.40
1:BI:99:LEU:HA	1:BI:99:LEU:HD23	1.81	0.40
1:BM:7:MET:HE2	1:BM:7:MET:HB3	1.96	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AR:66:CYS:SG	1:BQ:61:CYS:SG[2_555]	1.80	0.40
1:AV:103:ASN:O	1:BX:31:LYS:NZ[2_555]	2.19	0.01

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	AA	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AB	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	AC	116/118 (98%)	116 (100%)	0	0	100	100
1	AD	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AE	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	AF	116/118 (98%)	116 (100%)	0	0	100	100
1	AG	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AH	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	AI	116/118 (98%)	116 (100%)	0	0	100	100
1	AJ	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AK	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	AL	116/118 (98%)	116 (100%)	0	0	100	100
1	AM	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AN	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	AO	116/118 (98%)	116 (100%)	0	0	100	100
1	AP	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AQ	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	AR	116/118 (98%)	116 (100%)	0	0	100	100
1	AS	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AT	116/118 (98%)	114 (98%)	2 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AU	116/118 (98%)	116 (100%)	0	0	100	100
1	AV	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AW	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	AX	116/118 (98%)	116 (100%)	0	0	100	100
1	AY	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	AZ	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BA	116/118 (98%)	116 (100%)	0	0	100	100
1	BB	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BC	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BD	116/118 (98%)	116 (100%)	0	0	100	100
1	BE	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BF	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BG	116/118 (98%)	116 (100%)	0	0	100	100
1	BH	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BI	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BJ	116/118 (98%)	116 (100%)	0	0	100	100
1	BK	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BL	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BM	116/118 (98%)	116 (100%)	0	0	100	100
1	BN	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BO	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BP	116/118 (98%)	116 (100%)	0	0	100	100
1	BQ	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BR	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BS	116/118 (98%)	116 (100%)	0	0	100	100
1	BT	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BU	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BV	116/118 (98%)	116 (100%)	0	0	100	100
1	BW	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	BX	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	BY	116/118 (98%)	116 (100%)	0	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BZ	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	CA	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	CB	116/118 (98%)	116 (100%)	0	0	100	100
1	CC	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	CD	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	CE	116/118 (98%)	116 (100%)	0	0	100	100
1	CF	116/118 (98%)	115 (99%)	1 (1%)	0	100	100
1	CG	116/118 (98%)	114 (98%)	2 (2%)	0	100	100
1	CH	116/118 (98%)	116 (100%)	0	0	100	100
All	All	6960/7080 (98%)	6900 (99%)	60 (1%)	0	100	100

There are no Ramachandran outliers to report.

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	AA	98/98 (100%)	98 (100%)	0	100	100
1	AB	98/98 (100%)	98 (100%)	0	100	100
1	AC	98/98 (100%)	98 (100%)	0	100	100
1	AD	98/98 (100%)	98 (100%)	0	100	100
1	AE	98/98 (100%)	98 (100%)	0	100	100
1	AF	98/98 (100%)	98 (100%)	0	100	100
1	AG	98/98 (100%)	98 (100%)	0	100	100
1	AH	98/98 (100%)	98 (100%)	0	100	100
1	AI	98/98 (100%)	98 (100%)	0	100	100
1	AJ	98/98 (100%)	98 (100%)	0	100	100
1	AK	98/98 (100%)	98 (100%)	0	100	100
1	AL	98/98 (100%)	98 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	AM	98/98 (100%)	98 (100%)	0	100	100
1	AN	98/98 (100%)	98 (100%)	0	100	100
1	AO	98/98 (100%)	98 (100%)	0	100	100
1	AP	98/98 (100%)	98 (100%)	0	100	100
1	AQ	98/98 (100%)	98 (100%)	0	100	100
1	AR	98/98 (100%)	98 (100%)	0	100	100
1	AS	98/98 (100%)	98 (100%)	0	100	100
1	AT	98/98 (100%)	98 (100%)	0	100	100
1	AU	98/98 (100%)	98 (100%)	0	100	100
1	AV	98/98 (100%)	98 (100%)	0	100	100
1	AW	98/98 (100%)	98 (100%)	0	100	100
1	AX	98/98 (100%)	98 (100%)	0	100	100
1	AY	98/98 (100%)	98 (100%)	0	100	100
1	AZ	98/98 (100%)	98 (100%)	0	100	100
1	BA	98/98 (100%)	98 (100%)	0	100	100
1	BB	98/98 (100%)	98 (100%)	0	100	100
1	BC	98/98 (100%)	98 (100%)	0	100	100
1	BD	98/98 (100%)	98 (100%)	0	100	100
1	BE	98/98 (100%)	98 (100%)	0	100	100
1	BF	98/98 (100%)	98 (100%)	0	100	100
1	BG	98/98 (100%)	98 (100%)	0	100	100
1	BH	98/98 (100%)	98 (100%)	0	100	100
1	BI	98/98 (100%)	98 (100%)	0	100	100
1	BJ	98/98 (100%)	98 (100%)	0	100	100
1	BK	98/98 (100%)	98 (100%)	0	100	100
1	BL	98/98 (100%)	98 (100%)	0	100	100
1	BM	98/98 (100%)	98 (100%)	0	100	100
1	BN	98/98 (100%)	98 (100%)	0	100	100
1	BO	98/98 (100%)	98 (100%)	0	100	100
1	BP	98/98 (100%)	98 (100%)	0	100	100
1	BQ	98/98 (100%)	98 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	BR	98/98 (100%)	98 (100%)	0	100	100
1	BS	98/98 (100%)	98 (100%)	0	100	100
1	BT	98/98 (100%)	98 (100%)	0	100	100
1	BU	98/98 (100%)	98 (100%)	0	100	100
1	BV	98/98 (100%)	98 (100%)	0	100	100
1	BW	98/98 (100%)	98 (100%)	0	100	100
1	BX	98/98 (100%)	98 (100%)	0	100	100
1	BY	98/98 (100%)	98 (100%)	0	100	100
1	BZ	98/98 (100%)	98 (100%)	0	100	100
1	CA	98/98 (100%)	98 (100%)	0	100	100
1	CB	98/98 (100%)	98 (100%)	0	100	100
1	CC	98/98 (100%)	98 (100%)	0	100	100
1	CD	98/98 (100%)	98 (100%)	0	100	100
1	CE	98/98 (100%)	98 (100%)	0	100	100
1	CF	98/98 (100%)	98 (100%)	0	100	100
1	CG	98/98 (100%)	98 (100%)	0	100	100
1	CH	98/98 (100%)	98 (100%)	0	100	100
All	All	5880/5880 (100%)	5880 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

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

Mol	Chain	Res	Type
1	AA	49	HIS
1	AC	15	GLN
1	AC	38	ASN
1	AC	111	ASN
1	AD	15	GLN
1	AF	38	ASN
1	AH	15	GLN
1	AI	38	ASN
1	AL	15	GLN
1	AL	38	ASN
1	AR	15	GLN
1	AR	38	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	AU	111	ASN
1	AV	15	GLN
1	AX	15	GLN
1	AX	38	ASN
1	AY	49	HIS
1	AZ	15	GLN
1	BA	38	ASN
1	BD	111	ASN
1	BF	15	GLN
1	BG	15	GLN
1	BG	38	ASN
1	BJ	38	ASN
1	BK	15	GLN
1	BM	38	ASN
1	BM	111	ASN
1	BP	15	GLN
1	BP	38	ASN
1	BQ	15	GLN
1	BQ	49	HIS
1	BS	111	ASN
1	BT	49	HIS
1	BV	15	GLN
1	BV	111	ASN
1	BY	15	GLN
1	BY	38	ASN
1	CE	38	ASN
1	CG	15	GLN
1	CH	111	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

There are no ligands in this entry.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

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	AA	118/118 (100%)	0.17	2 (1%) 69 49	53, 84, 186, 256	0
1	AB	118/118 (100%)	0.30	8 (6%) 23 15	53, 85, 203, 280	0
1	AC	118/118 (100%)	0.22	3 (2%) 58 39	54, 83, 206, 294	0
1	AD	118/118 (100%)	0.18	4 (3%) 48 30	53, 84, 186, 256	0
1	AE	118/118 (100%)	0.42	10 (8%) 16 12	53, 85, 203, 280	0
1	AF	118/118 (100%)	0.10	3 (2%) 58 39	54, 83, 206, 294	0
1	AG	118/118 (100%)	0.13	4 (3%) 48 30	53, 84, 186, 256	0
1	AH	118/118 (100%)	0.20	2 (1%) 69 49	53, 85, 203, 280	0
1	AI	118/118 (100%)	0.09	3 (2%) 58 39	54, 83, 206, 294	0
1	AJ	118/118 (100%)	0.18	1 (0%) 82 67	53, 84, 186, 256	0
1	AK	118/118 (100%)	0.34	5 (4%) 40 26	53, 85, 203, 280	0
1	AL	118/118 (100%)	0.16	3 (2%) 58 39	54, 83, 206, 294	0
1	AM	118/118 (100%)	0.04	1 (0%) 82 67	53, 84, 186, 256	0
1	AN	118/118 (100%)	0.34	3 (2%) 58 39	53, 85, 203, 280	0
1	AO	118/118 (100%)	0.22	3 (2%) 58 39	54, 83, 206, 294	0
1	AP	118/118 (100%)	0.01	1 (0%) 82 67	53, 84, 186, 256	0
1	AQ	118/118 (100%)	0.44	6 (5%) 33 22	53, 85, 203, 280	0
1	AR	118/118 (100%)	0.19	2 (1%) 69 49	54, 83, 206, 294	0
1	AS	118/118 (100%)	0.21	3 (2%) 58 39	53, 84, 186, 256	0
1	AT	118/118 (100%)	0.19	8 (6%) 23 15	53, 85, 203, 280	0
1	AU	118/118 (100%)	0.13	1 (0%) 82 67	54, 83, 206, 294	0
1	AV	118/118 (100%)	0.34	3 (2%) 58 39	53, 84, 186, 256	0
1	AW	118/118 (100%)	0.36	7 (5%) 28 18	53, 85, 203, 280	0
1	AX	118/118 (100%)	0.27	5 (4%) 40 26	54, 83, 206, 294	0

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	AY	118/118 (100%)	0.19	3 (2%)	58 39	53, 84, 186, 256	0
1	AZ	118/118 (100%)	0.38	9 (7%)	20 14	53, 85, 203, 280	0
1	BA	118/118 (100%)	0.10	2 (1%)	69 49	54, 83, 206, 294	0
1	BB	118/118 (100%)	0.35	6 (5%)	33 22	53, 84, 186, 256	0
1	BC	118/118 (100%)	0.54	10 (8%)	16 12	53, 85, 203, 280	0
1	BD	118/118 (100%)	0.26	2 (1%)	69 49	54, 83, 206, 294	0
1	BE	118/118 (100%)	0.27	5 (4%)	40 26	53, 84, 186, 256	0
1	BF	118/118 (100%)	0.33	5 (4%)	40 26	53, 85, 203, 280	0
1	BG	118/118 (100%)	0.32	3 (2%)	58 39	54, 83, 206, 294	0
1	BH	118/118 (100%)	0.10	1 (0%)	82 67	53, 84, 186, 256	0
1	BI	118/118 (100%)	0.10	4 (3%)	48 30	53, 85, 203, 280	0
1	BJ	118/118 (100%)	0.18	3 (2%)	58 39	54, 83, 206, 294	0
1	BK	118/118 (100%)	0.39	8 (6%)	23 15	53, 84, 186, 256	0
1	BL	118/118 (100%)	0.39	6 (5%)	33 22	53, 85, 203, 280	0
1	BM	118/118 (100%)	0.22	2 (1%)	69 49	54, 83, 206, 294	0
1	BN	118/118 (100%)	0.34	6 (5%)	33 22	53, 84, 186, 256	0
1	BO	118/118 (100%)	0.45	9 (7%)	20 14	53, 85, 203, 280	0
1	BP	118/118 (100%)	0.33	5 (4%)	40 26	54, 83, 206, 294	0
1	BQ	118/118 (100%)	0.24	7 (5%)	28 18	53, 84, 186, 256	0
1	BR	118/118 (100%)	0.45	8 (6%)	23 15	53, 85, 203, 280	0
1	BS	118/118 (100%)	0.18	4 (3%)	48 30	54, 83, 206, 294	0
1	BT	118/118 (100%)	0.41	7 (5%)	28 18	53, 84, 186, 256	0
1	BU	118/118 (100%)	0.38	8 (6%)	23 15	53, 85, 203, 280	0
1	BV	118/118 (100%)	0.17	4 (3%)	48 30	54, 83, 206, 294	0
1	BW	118/118 (100%)	0.32	4 (3%)	48 30	53, 84, 186, 256	0
1	BX	118/118 (100%)	0.65	8 (6%)	23 15	53, 85, 203, 280	0
1	BY	118/118 (100%)	0.41	3 (2%)	58 39	54, 83, 206, 294	0
1	BZ	118/118 (100%)	0.32	3 (2%)	58 39	53, 84, 186, 256	0
1	CA	118/118 (100%)	0.45	6 (5%)	33 22	53, 85, 203, 280	0
1	CB	118/118 (100%)	0.32	3 (2%)	58 39	54, 83, 206, 294	0
1	CC	118/118 (100%)	0.22	4 (3%)	48 30	53, 84, 186, 256	0

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	CD	118/118 (100%)	0.22	7 (5%)	28	18	53, 85, 203, 280	0
1	CE	118/118 (100%)	0.16	3 (2%)	58	39	54, 83, 206, 294	0
1	CF	118/118 (100%)	0.22	2 (1%)	69	49	53, 84, 186, 256	0
1	CG	118/118 (100%)	0.37	6 (5%)	33	22	53, 85, 203, 280	0
1	CH	118/118 (100%)	0.37	6 (5%)	33	22	54, 83, 206, 294	0
All	All	7080/7080 (100%)	0.27	273 (3%)	43	27	53, 84, 203, 294	0

All (273) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BM	63	PRO	7.8
1	AZ	63	PRO	6.5
1	CH	63	PRO	6.2
1	BZ	13	ALA	6.2
1	BD	63	PRO	5.9
1	AQ	63	PRO	5.8
1	AX	63	PRO	5.8
1	CE	63	PRO	5.7
1	BU	59	ASN	5.6
1	AI	63	PRO	5.4
1	BB	13	ALA	5.3
1	BL	10	GLY	5.3
1	AZ	1	ALA	5.2
1	AN	59	ASN	5.2
1	BC	63	PRO	5.1
1	AO	63	PRO	5.0
1	BT	13	ALA	4.9
1	BR	63	PRO	4.9
1	AE	12	SER	4.9
1	BJ	63	PRO	4.8
1	AE	64	THR	4.6
1	CG	59	ASN	4.6
1	CB	63	PRO	4.6
1	AL	63	PRO	4.5
1	AR	63	PRO	4.5
1	AQ	65	LYS	4.5
1	BY	63	PRO	4.4
1	BS	63	PRO	4.3
1	BK	37	ARG	4.2
1	AU	63	PRO	4.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BW	51	ASN	4.0
1	BQ	30	GLN	4.0
1	AZ	8	ALA	4.0
1	BT	51	ASN	3.9
1	AB	8	ALA	3.8
1	AC	63	PRO	3.8
1	BC	64	THR	3.8
1	BU	8	ALA	3.8
1	CA	13	ALA	3.7
1	AQ	59	ASN	3.7
1	BU	13	ALA	3.6
1	AK	59	ASN	3.6
1	BI	59	ASN	3.6
1	AQ	13	ALA	3.5
1	BV	63	PRO	3.5
1	AK	11	ASN	3.5
1	CF	51	ASN	3.5
1	BA	63	PRO	3.5
1	AX	110	LYS	3.5
1	BL	1	ALA	3.5
1	BB	11	ASN	3.4
1	BU	63	PRO	3.4
1	AZ	40	SER	3.4
1	BB	14	ILE	3.4
1	AC	37	ARG	3.3
1	BF	110	LYS	3.3
1	AF	63	PRO	3.3
1	CD	76	PHE	3.3
1	BF	59	ASN	3.3
1	BI	8	ALA	3.2
1	AA	30	GLN	3.2
1	AA	13	ALA	3.2
1	BO	8	ALA	3.2
1	BX	59	ASN	3.2
1	AZ	65	LYS	3.2
1	CE	11	ASN	3.2
1	AB	10	GLY	3.2
1	BX	65	LYS	3.2
1	BR	12	SER	3.2
1	BE	13	ALA	3.2
1	CA	59	ASN	3.1
1	BP	63	PRO	3.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BW	68	PRO	3.1
1	BG	63	PRO	3.1
1	BW	14	ILE	3.1
1	BP	111	ASN	3.0
1	AY	60	THR	3.0
1	CA	110	LYS	3.0
1	BK	68	PRO	3.0
1	BK	13	ALA	3.0
1	CD	8	ALA	3.0
1	AR	110	LYS	3.0
1	AE	59	ASN	3.0
1	AS	30	GLN	2.9
1	CD	110	LYS	2.9
1	AW	38	ASN	2.9
1	BC	8	ALA	2.9
1	CH	42	SER	2.9
1	BR	65	LYS	2.9
1	BO	59	ASN	2.9
1	CD	59	ASN	2.9
1	AY	61	CYS	2.9
1	AY	37	ARG	2.9
1	BQ	13	ALA	2.9
1	BP	109	ASN	2.9
1	AT	59	ASN	2.8
1	AX	42	SER	2.8
1	CA	8	ALA	2.8
1	BR	43	THR	2.8
1	BU	64	THR	2.8
1	AN	110	LYS	2.8
1	BC	59	ASN	2.8
1	BU	65	LYS	2.8
1	AQ	8	ALA	2.8
1	AB	59	ASN	2.8
1	AW	109	ASN	2.8
1	BY	11	ASN	2.8
1	AD	30	GLN	2.8
1	BC	13	ALA	2.8
1	BX	76	PHE	2.7
1	AE	110	LYS	2.7
1	AT	13	ALA	2.7
1	BZ	30	GLN	2.7
1	BX	63	PRO	2.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BF	11	ASN	2.7
1	AO	118	ALA	2.7
1	BN	51	ASN	2.7
1	BV	30	GLN	2.6
1	AK	12	SER	2.6
1	AM	13	ALA	2.6
1	AD	13	ALA	2.6
1	CC	13	ALA	2.6
1	BC	110	LYS	2.6
1	BB	30	GLN	2.6
1	CH	30	GLN	2.6
1	BQ	51	ASN	2.6
1	AT	43	THR	2.6
1	BO	64	THR	2.6
1	BU	10	GLY	2.6
1	BO	65	LYS	2.6
1	BR	64	THR	2.6
1	AE	63	PRO	2.5
1	BC	106	SER	2.5
1	BE	42	SER	2.5
1	CA	12	SER	2.5
1	CC	30	GLN	2.5
1	AK	118	ALA	2.5
1	BO	63	PRO	2.5
1	AZ	59	ASN	2.5
1	BP	11	ASN	2.5
1	CC	11	ASN	2.5
1	AV	37	ARG	2.5
1	BB	12	SER	2.5
1	CE	38	ASN	2.5
1	BS	64	THR	2.5
1	AK	8	ALA	2.4
1	AW	1	ALA	2.4
1	BK	118	ALA	2.4
1	AS	14	ILE	2.4
1	BN	75	SER	2.4
1	BF	76	PHE	2.4
1	BG	113	ASP	2.4
1	AJ	73	GLY	2.4
1	BD	10	GLY	2.4
1	AN	13	ALA	2.4
1	BB	118	ALA	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	AZ	61	CYS	2.4
1	BW	28	SER	2.4
1	BS	83	ASP	2.4
1	AH	8	ALA	2.4
1	AH	12	SER	2.4
1	AE	109	ASN	2.4
1	CG	13	ALA	2.4
1	CH	7	MET	2.4
1	BE	68	PRO	2.4
1	AB	1	ALA	2.4
1	AT	8	ALA	2.4
1	CC	12	SER	2.4
1	BN	63	PRO	2.4
1	AT	11	ASN	2.3
1	BT	30	GLN	2.3
1	AB	11	ASN	2.3
1	BG	65	LYS	2.3
1	BN	65	LYS	2.3
1	BN	77	LYS	2.3
1	AL	10	GLY	2.3
1	AP	1	ALA	2.3
1	AE	11	ASN	2.3
1	BC	103	ASN	2.3
1	BT	11	ASN	2.3
1	BO	76	PHE	2.3
1	AB	63	PRO	2.3
1	AG	75	SER	2.3
1	CB	38	ASN	2.3
1	BS	113	ASP	2.3
1	BV	65	LYS	2.3
1	AV	118	ALA	2.3
1	BL	13	ALA	2.3
1	AW	12	SER	2.3
1	BJ	110	LYS	2.3
1	BX	109	ASN	2.3
1	AO	30	GLN	2.3
1	BQ	14	ILE	2.2
1	AZ	13	ALA	2.2
1	BL	109	ASN	2.2
1	CG	70	VAL	2.2
1	BX	8	ALA	2.2
1	BA	110	LYS	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	CH	32	THR	2.2
1	BV	35	VAL	2.2
1	BQ	11	ASN	2.2
1	AE	14	ILE	2.2
1	AL	113	ASP	2.2
1	BR	36	LEU	2.2
1	BP	110	LYS	2.2
1	BQ	65	LYS	2.2
1	AT	40	SER	2.2
1	BH	30	GLN	2.2
1	AI	11	ASN	2.2
1	AT	38	ASN	2.2
1	AD	73	GLY	2.2
1	BC	9	GLY	2.2
1	BN	13	ALA	2.2
1	BT	68	PRO	2.2
1	CA	76	PHE	2.2
1	AD	11	ASN	2.2
1	AQ	11	ASN	2.2
1	BO	109	ASN	2.2
1	AI	10	GLY	2.2
1	AW	61	CYS	2.2
1	AZ	64	THR	2.1
1	BK	60	THR	2.1
1	AW	36	LEU	2.1
1	AX	38	ASN	2.1
1	CB	76	PHE	2.1
1	AG	74	VAL	2.1
1	AB	64	THR	2.1
1	AE	8	ALA	2.1
1	BO	118	ALA	2.1
1	CD	13	ALA	2.1
1	CD	40	SER	2.1
1	BI	11	ASN	2.1
1	BU	11	ASN	2.1
1	CD	11	ASN	2.1
1	AE	65	LYS	2.1
1	AF	64	THR	2.1
1	AG	14	ILE	2.1
1	BQ	60	THR	2.1
1	BO	13	ALA	2.1
1	AV	68	PRO	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BF	12	SER	2.1
1	BJ	42	SER	2.1
1	BT	75	SER	2.1
1	BR	62	ASP	2.1
1	BX	1	ALA	2.1
1	CG	30	GLN	2.1
1	BE	67	GLY	2.1
1	AB	76	PHE	2.1
1	AC	42	SER	2.1
1	BT	74	VAL	2.1
1	AS	11	ASN	2.1
1	BK	11	ASN	2.1
1	BM	38	ASN	2.1
1	BZ	50	ILE	2.1
1	CG	36	LEU	2.1
1	BE	57	THR	2.1
1	BL	8	ALA	2.1
1	BY	110	LYS	2.0
1	CF	106	SER	2.0
1	BR	59	ASN	2.0
1	CG	109	ASN	2.0
1	AG	13	ALA	2.0
1	BI	13	ALA	2.0
1	AX	76	PHE	2.0
1	BC	65	LYS	2.0
1	BK	35	VAL	2.0
1	BX	12	SER	2.0
1	AW	62	ASP	2.0
1	CH	83	ASP	2.0
1	AT	76	PHE	2.0
1	BK	10	GLY	2.0
1	AF	110	LYS	2.0
1	BL	30	GLN	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.