



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 11:33 AM UTC

PDB ID : 6YFP / pdb_00006yfp
Title : Virus-like particle of bacteriophage GQ-112
Authors : Rumnieks, J.; Kalnins, G.; Sisovs, M.; Lieknina, I.; Tars, K.
Deposited on : 2020-03-26
Resolution : 3.10 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

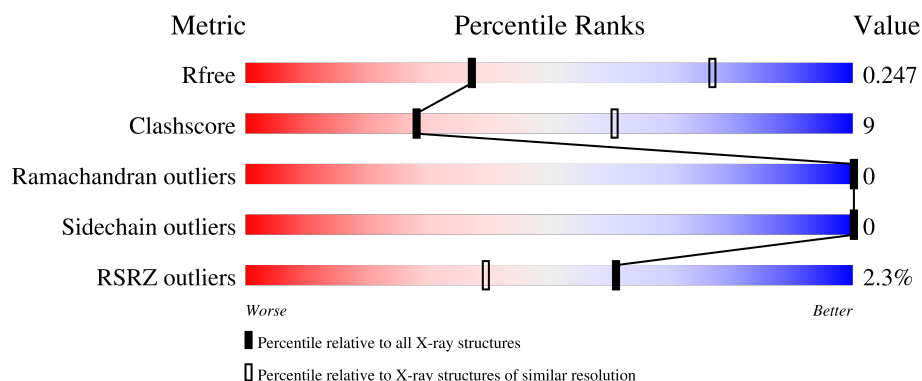
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.10 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1456 (3.10-3.10)
Clashscore	190562	1539 (3.10-3.10)
Ramachandran outliers	187476	1467 (3.10-3.10)
Sidechain outliers	187428	1467 (3.10-3.10)
RSRZ outliers	180081	1456 (3.10-3.10)

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

Mol	Chain	Length	Quality of chain
1	AA	164	 3% 80% 17%
1	AB	164	 2% 77% 21%
1	AC	164	 3% 83% 14%
1	AD	164	 0% 79% 21%
1	AE	164	 3% 75% 22%

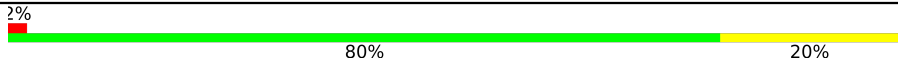
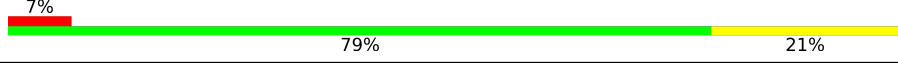
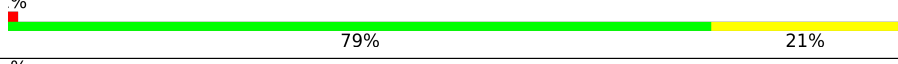
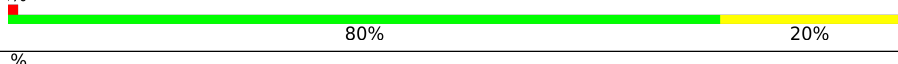
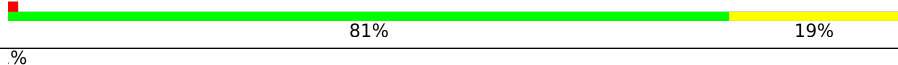
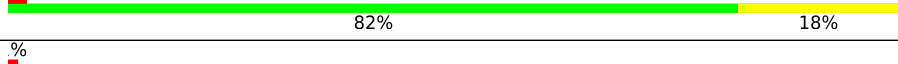
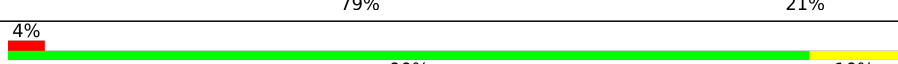
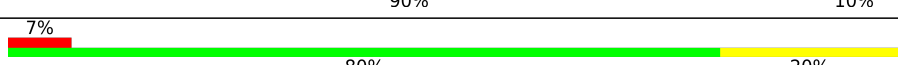
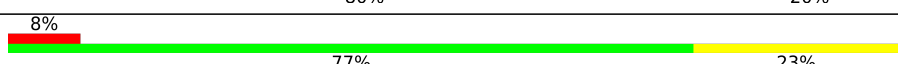
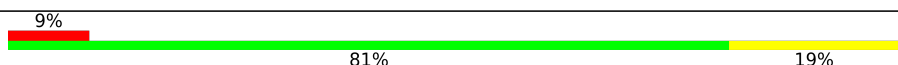
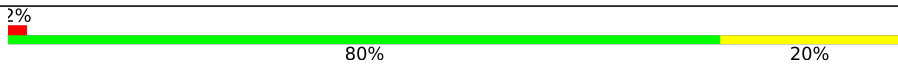
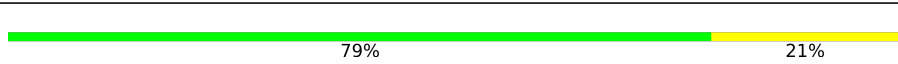
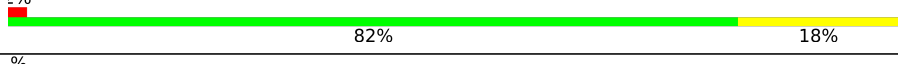
Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	AF	164	11% 82% 18%
1	AG	164	80% 20%
1	AH	164	% 76% 24%
1	AI	164	3% 83% 17%
1	AJ	164	3% 80% 20%
1	AK	164	% 76% 24%
1	AL	164	3% 83% 17%
1	AM	164	2% 79% 21%
1	AN	164	% 79% 21%
1	AO	164	2% 81% 19%
1	AP	164	% 81% 19%
1	AQ	164	% 78% 22%
1	AR	164	2% 80% 20%
1	AS	164	2% 80% 20%
1	AT	164	% 76% 24%
1	AU	164	3% 80% 20%
1	AV	164	2% 79% 21%
1	AW	164	% 76% 24%
1	AX	164	2% 84% 16%
1	AY	164	2% 81% 19%
1	AZ	164	% 79% 21%
1	BA	164	5% 82% 18%
1	BB	164	% 80% 20%
1	BC	164	% 76% 24%
1	BD	164	2% 82% 18%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain	
1	BE	164		
1	BF	164		
1	BG	164		
1	BH	164		
1	BI	164		
1	BJ	164		
1	BK	164		
1	BL	164		
1	BM	164		
1	BN	164		
1	BO	164		
1	BP	164		
1	BQ	164		
1	BR	164		
1	BS	164		
1	BT	164		
1	BU	164		
1	BV	164		
1	BW	164		
1	BX	164		
1	BY	164		
1	BZ	164		
1	CA	164		
1	CB	164		
1	CC	164		

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	CD	164	2% 77% 23%
1	CE	164	5% 82% 18%
1	CF	164	% 80% 20%
1	CG	164	% 78% 22%
1	CH	164	2% 90% 10%
1	CI	164	5% 80% 20%
1	CJ	164	5% 77% 23%
1	CK	164	5% 77% 23%
1	CL	164	2% 80% 20%
1	CM	164	% 77% 23%
1	CN	164	2% 82% 18%
1	CO	164	82% 18%
1	CP	164	% 78% 22%
1	CQ	164	4% 89% 11%
1	CR	164	2% 80% 20%
1	CS	164	% 79% 21%
1	CT	164	3% 81% 19%
1	CU	164	2% 80% 20%
1	CV	164	% 77% 23%
1	CW	164	4% 82% 18%
1	CX	164	% 80% 20%
1	CY	164	% 78% 22%
1	CZ	164	4% 83% 17%
1	DA	164	2% 80% 20%
1	DB	164	% 77% 23%

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
1	DC	164	 8% 81% 19%
1	DD	164	 % 80% 20%
1	DE	164	 2% 79% 21%
1	DF	164	 2% 80% 20%
1	DG	164	 2% 80% 20%
1	DH	164	 2% 76% 24%
1	DI	164	 2% 82% 18%
1	DJ	164	 % 80% 20%
1	DK	164	 % 78% 22%
1	DL	164	 4% 90% 10%

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 116370 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	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AB	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AC	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AD	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AE	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AF	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AG	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AH	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AI	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AJ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AK	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AL	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AM	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AN	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AO	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AP	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	AQ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AR	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AS	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AT	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AU	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AV	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AW	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AX	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AY	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	AZ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BA	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BB	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BC	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BD	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BE	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BF	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BG	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BH	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BI	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BJ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BK	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	BL	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BM	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BN	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BO	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BP	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BQ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BR	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BS	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BT	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BU	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BV	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BW	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BX	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BY	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	BZ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CA	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CB	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CC	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CD	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CE	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CF	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	CG	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CH	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CI	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CJ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CK	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CL	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CM	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CN	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CO	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CP	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CQ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CR	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CS	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CT	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CU	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CV	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CW	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CX	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CY	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	CZ	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			
1	DA	164	Total	C	N	O	S	0	0	0
			1293	811	223	256	3			

Continued on next page...

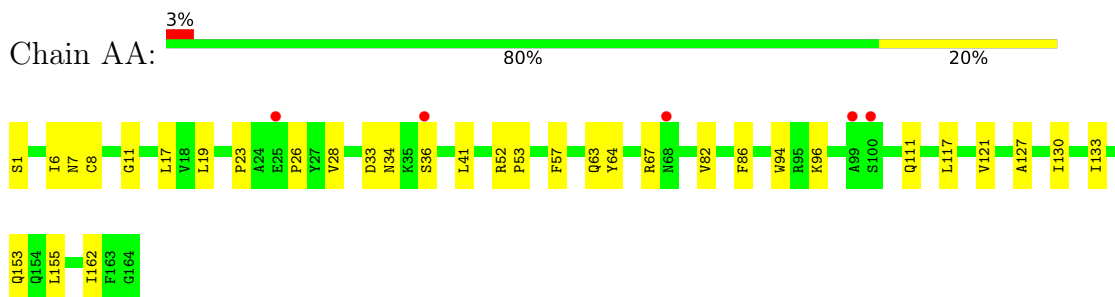
Continued from previous page...

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	DB	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DC	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DD	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DE	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DF	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DG	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DH	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DI	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DJ	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DK	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0
1	DL	164	Total 1293	C 811	N 223	O 256	S 3	0	0	0

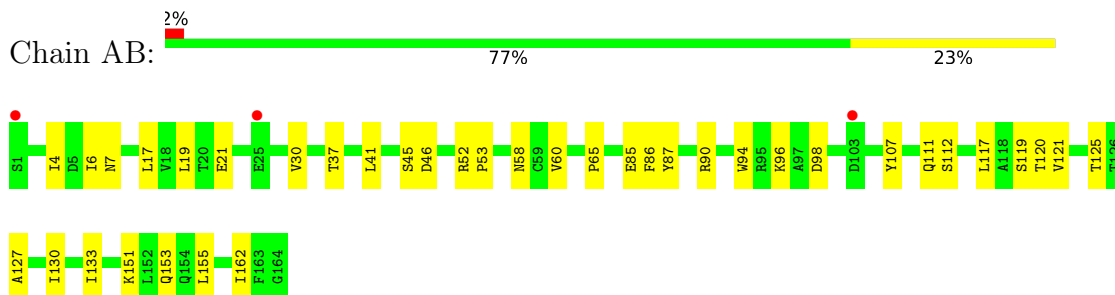
3 Residue-property plots [i](#)

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.

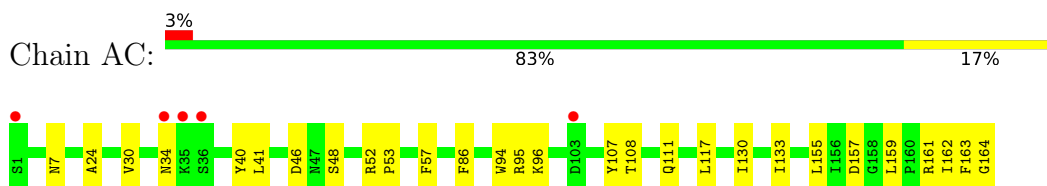
- Molecule 1: coat protein



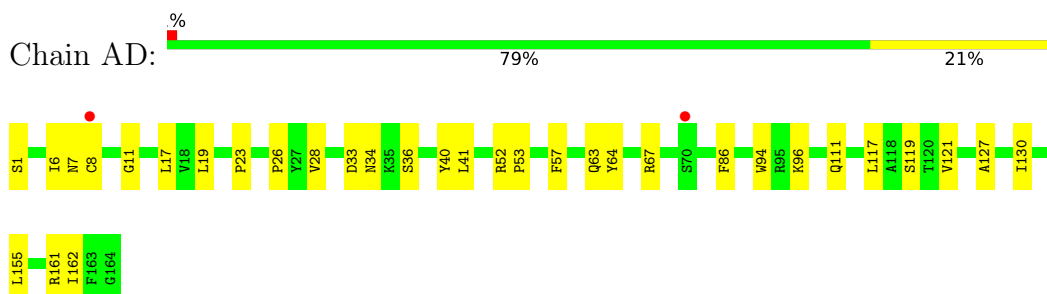
- Molecule 1: coat protein



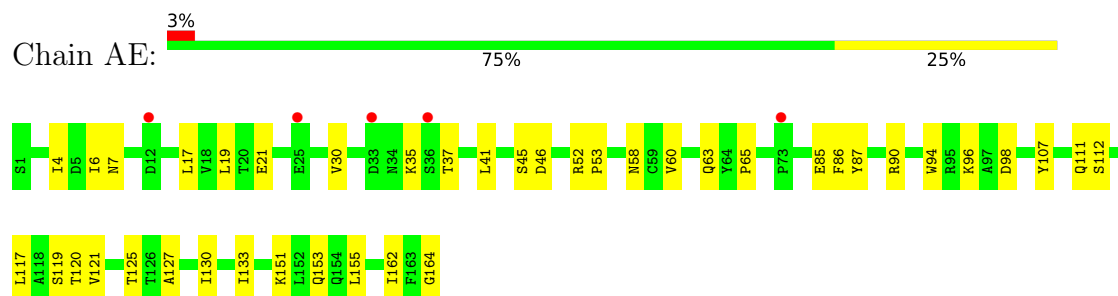
- Molecule 1: coat protein



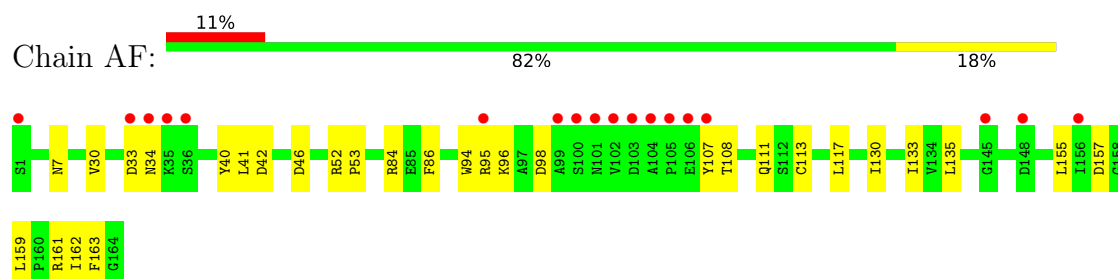
- Molecule 1: coat protein



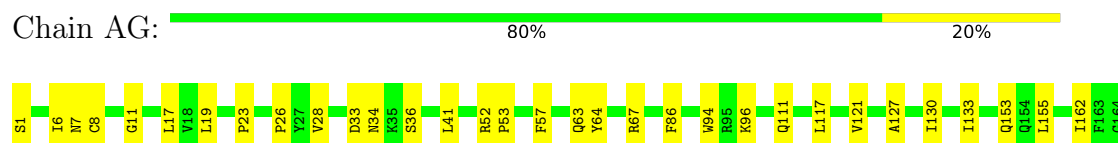
- Molecule 1: coat protein



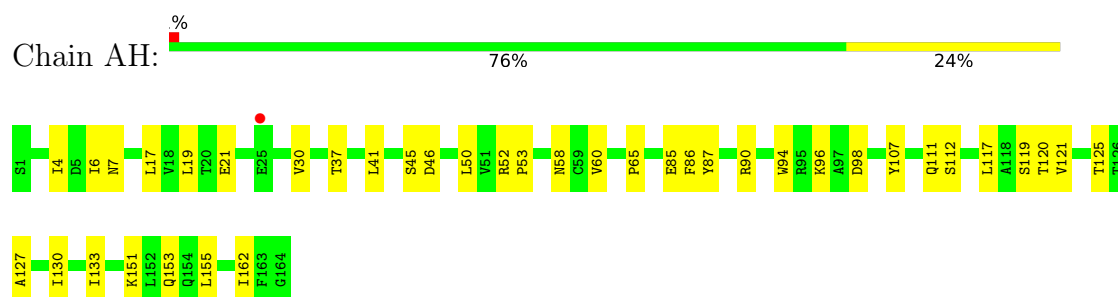
- Molecule 1: coat protein



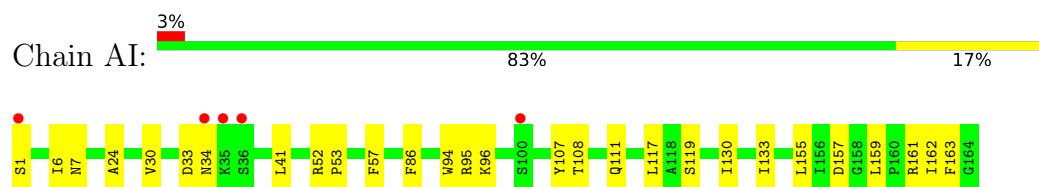
- Molecule 1: coat protein



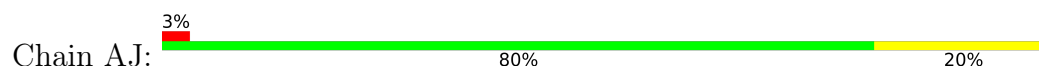
- Molecule 1: coat protein

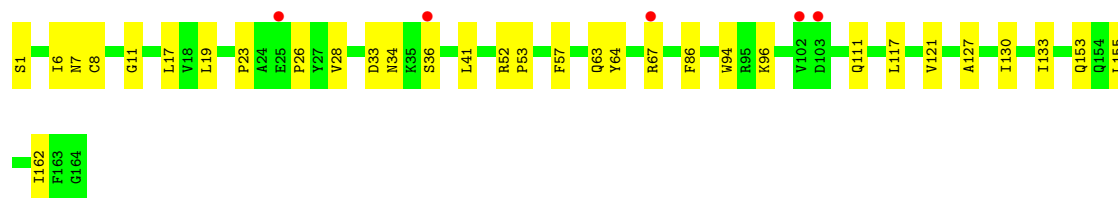


- Molecule 1: coat protein

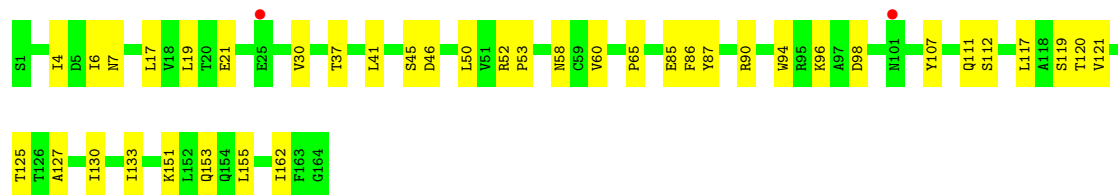
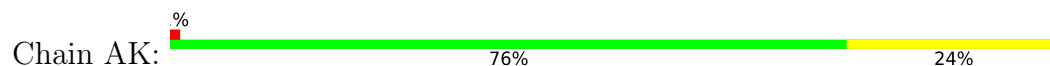


- Molecule 1: coat protein

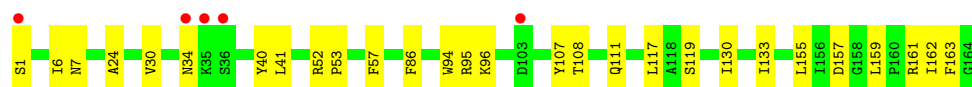
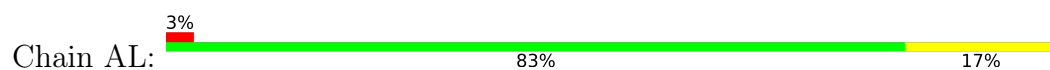




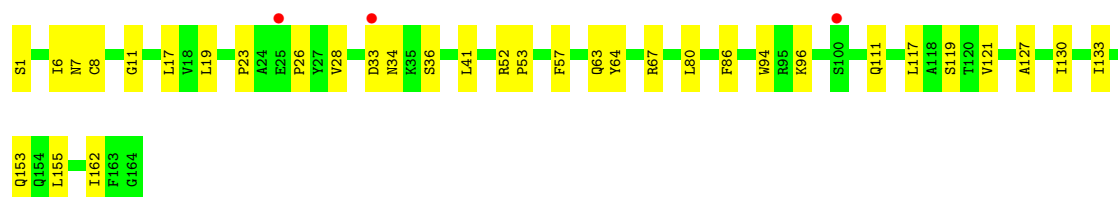
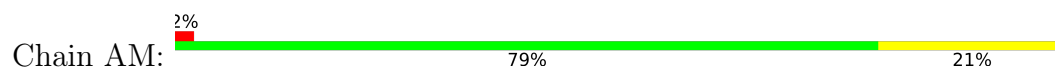
- Molecule 1: coat protein



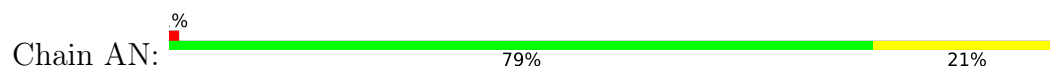
- Molecule 1: coat protein



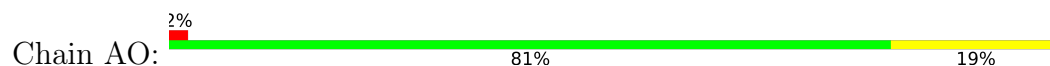
- Molecule 1: coat protein



- Molecule 1: coat protein

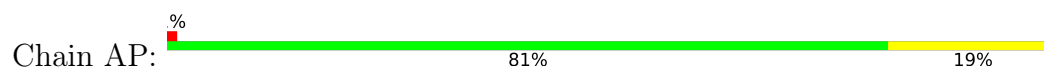


- Molecule 1: coat protein

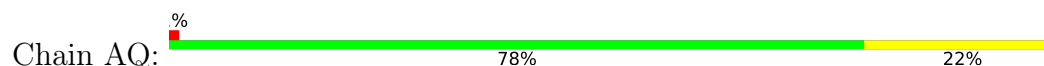




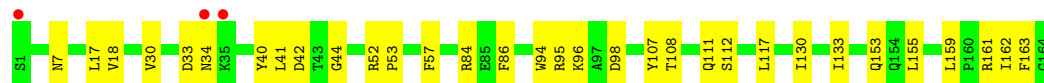
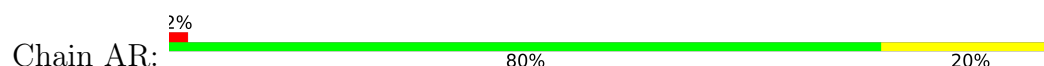
- Molecule 1: coat protein



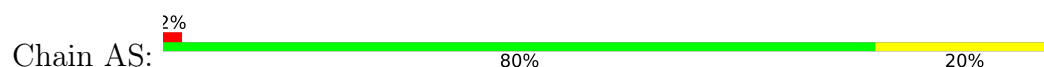
- Molecule 1: coat protein



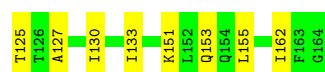
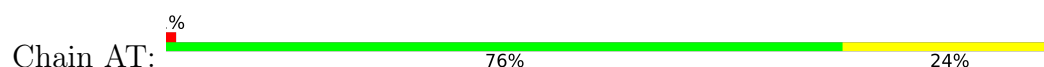
- Molecule 1: coat protein



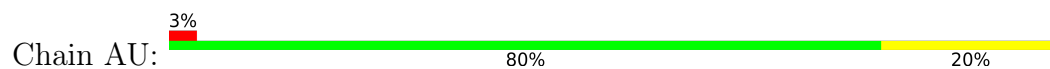
- Molecule 1: coat protein



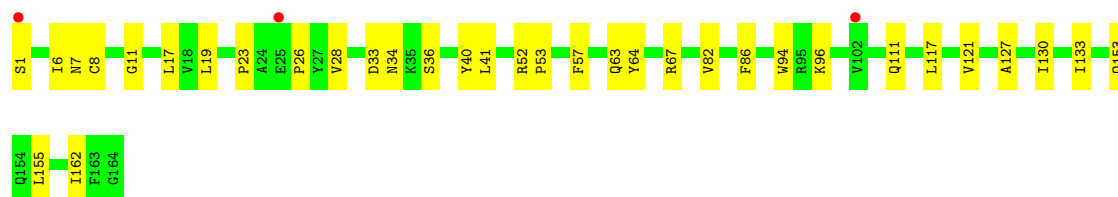
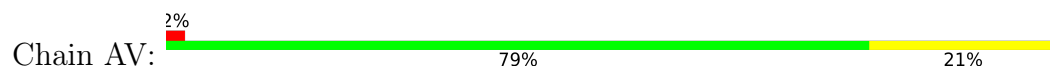
- Molecule 1: coat protein



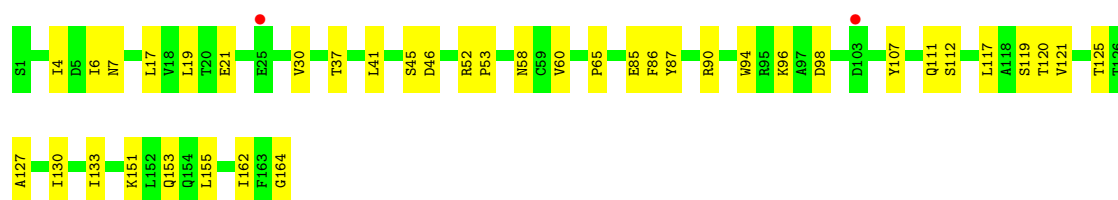
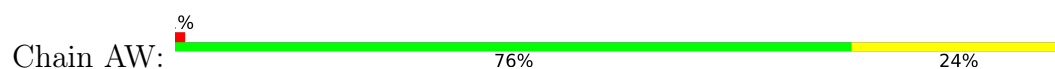
- Molecule 1: coat protein



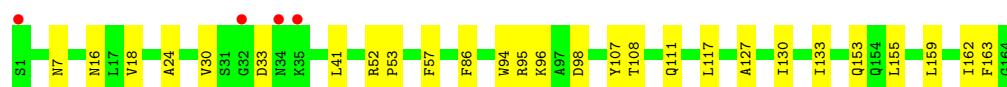
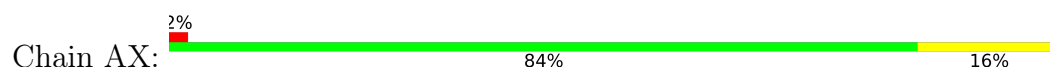
- Molecule 1: coat protein



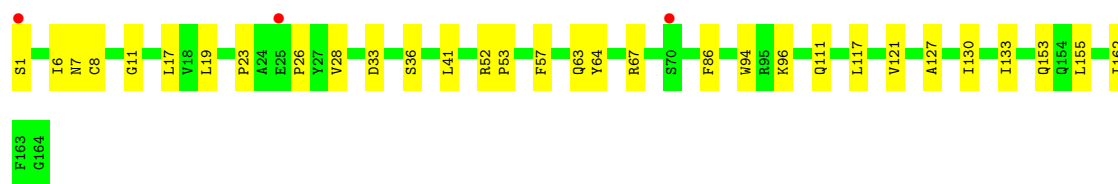
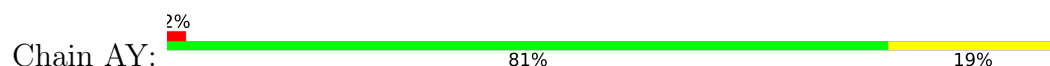
- Molecule 1: coat protein



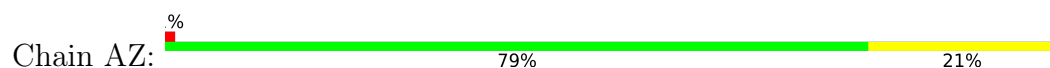
- Molecule 1: coat protein

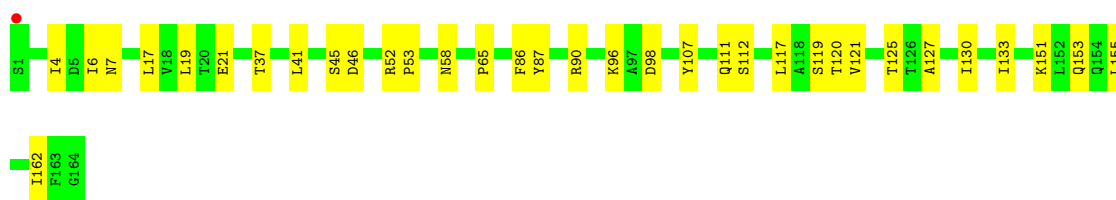


- Molecule 1: coat protein

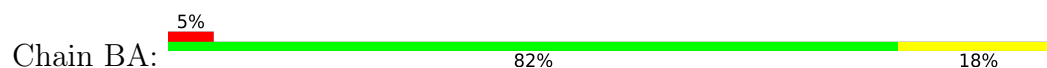


- Molecule 1: coat protein

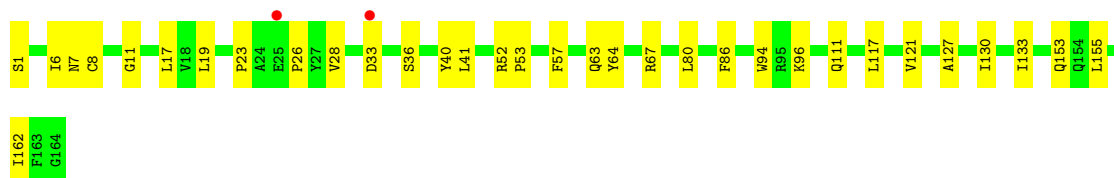
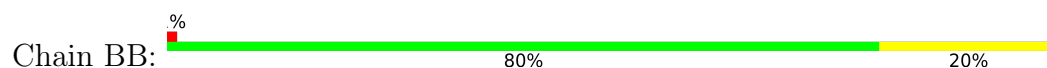




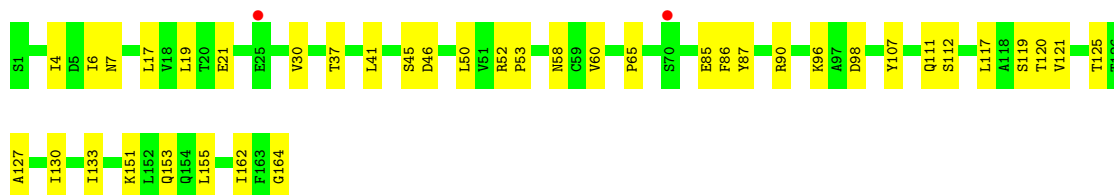
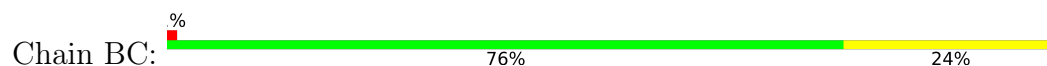
- Molecule 1: coat protein



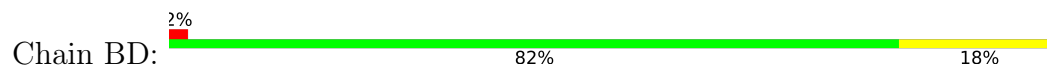
- Molecule 1: coat protein



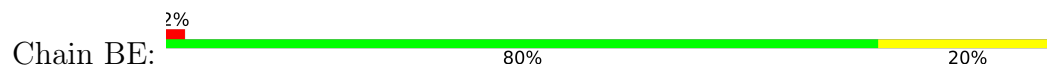
- Molecule 1: coat protein

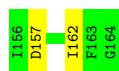


- Molecule 1: coat protein

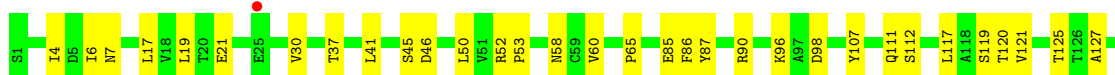
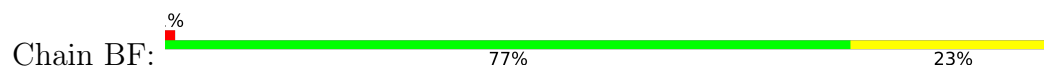


- Molecule 1: coat protein

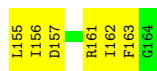
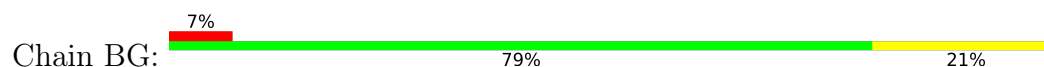




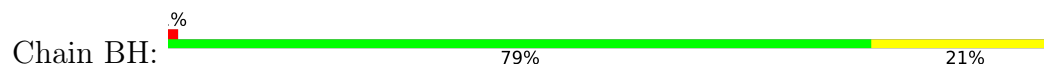
- Molecule 1: coat protein



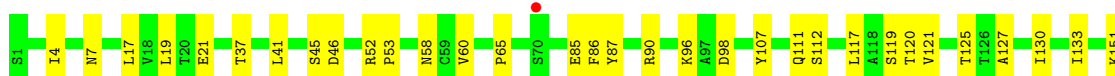
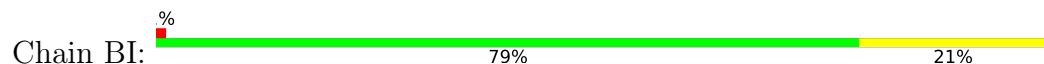
- Molecule 1: coat protein



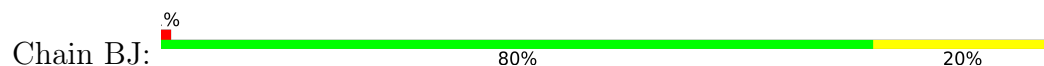
- Molecule 1: coat protein

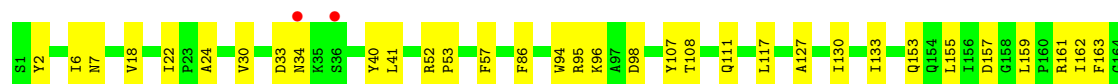


- Molecule 1: coat protein

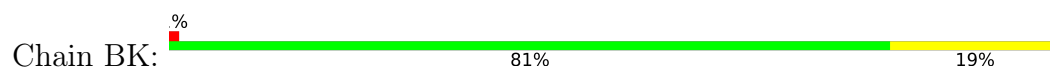


- Molecule 1: coat protein



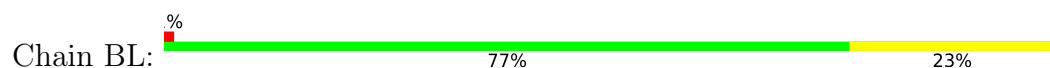


- Molecule 1: coat protein



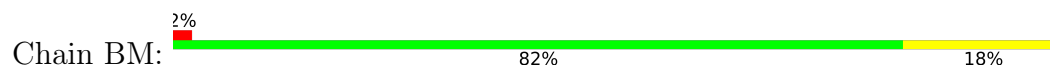
G164

- Molecule 1: coat protein

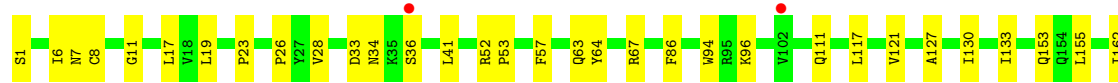
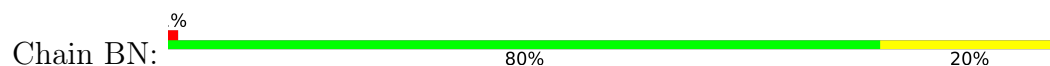


I130
I133
K151
I152
Q153
L155
I162
F163
G164

- Molecule 1: coat protein

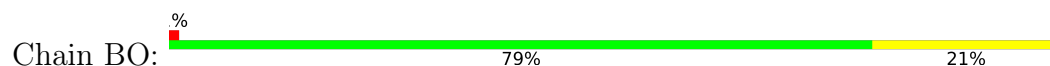


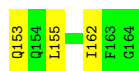
- Molecule 1: coat protein



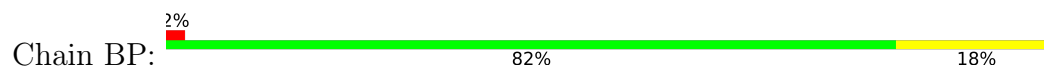
F163
G164

- Molecule 1: coat protein

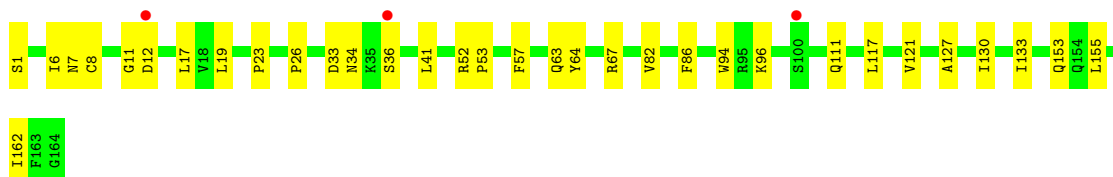
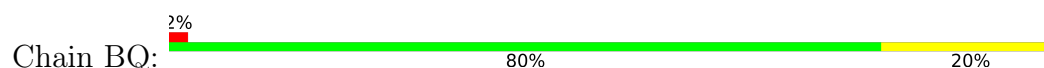




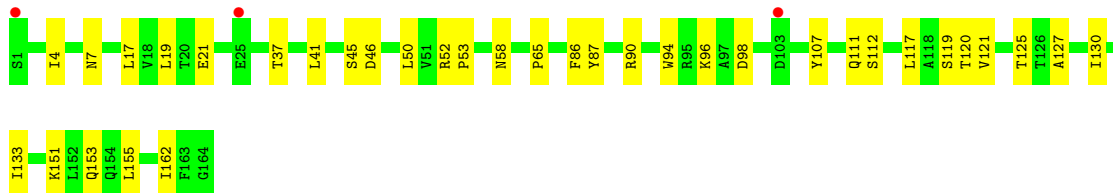
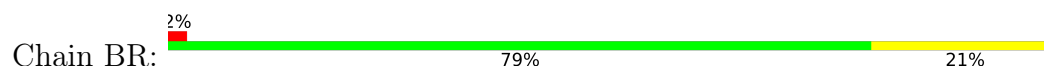
- Molecule 1: coat protein



- Molecule 1: coat protein



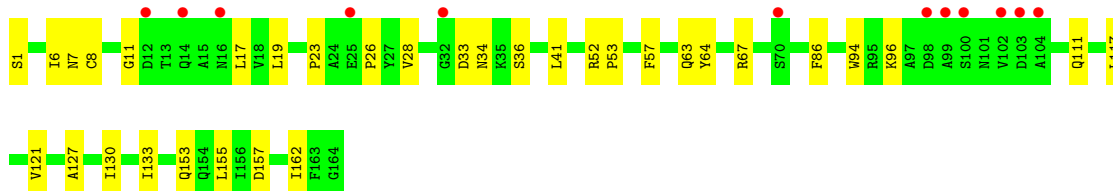
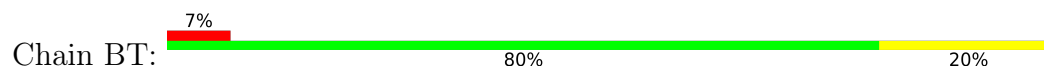
- Molecule 1: coat protein



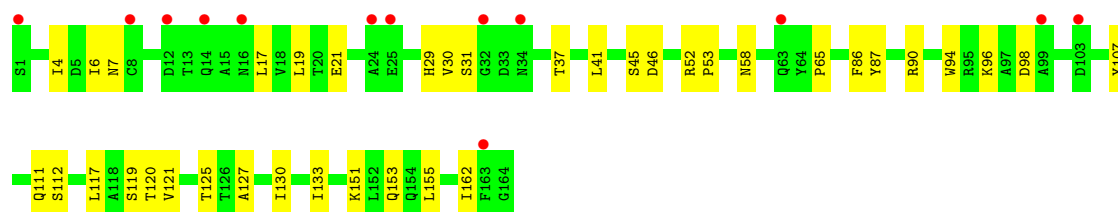
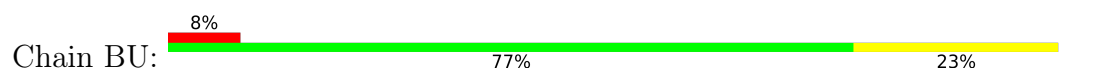
- Molecule 1: coat protein



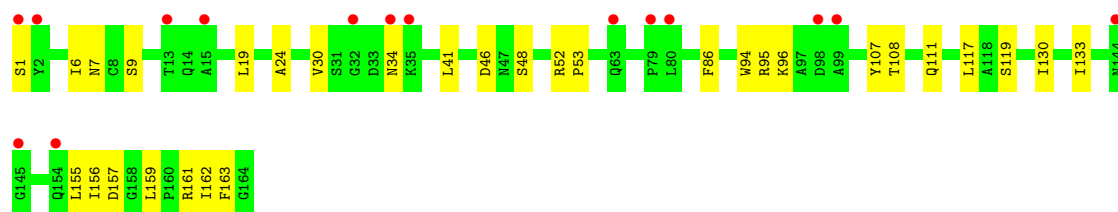
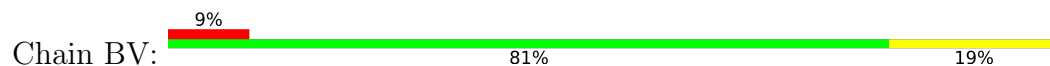
- Molecule 1: coat protein



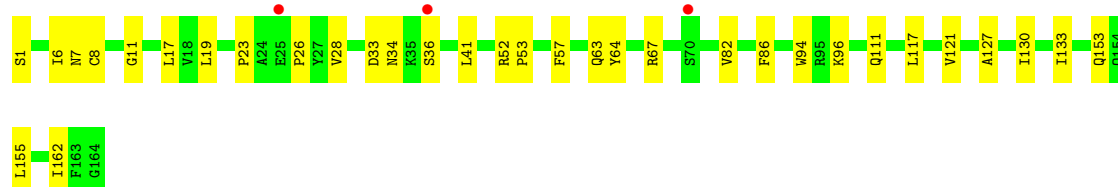
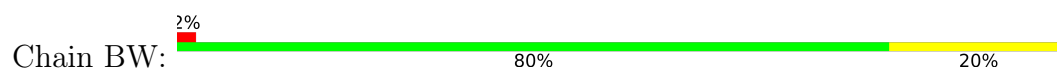
- Molecule 1: coat protein



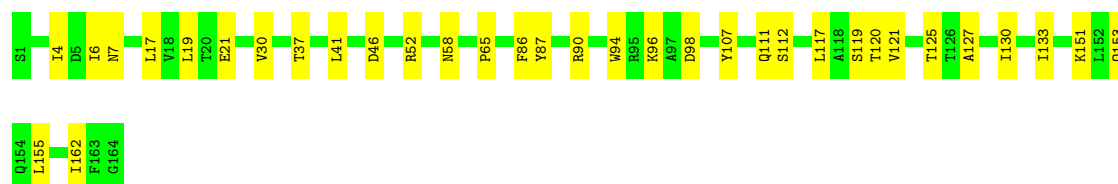
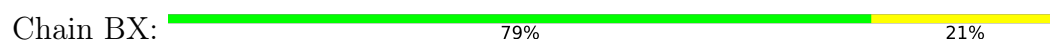
- Molecule 1: coat protein



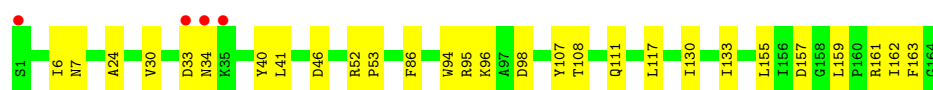
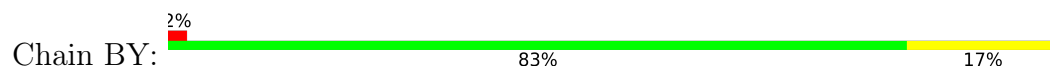
- Molecule 1: coat protein



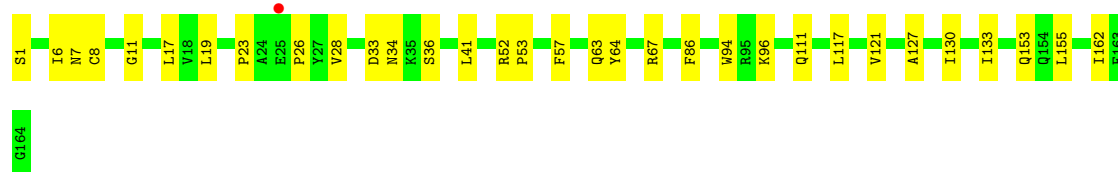
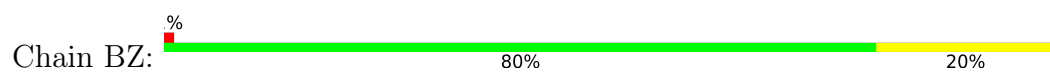
- Molecule 1: coat protein



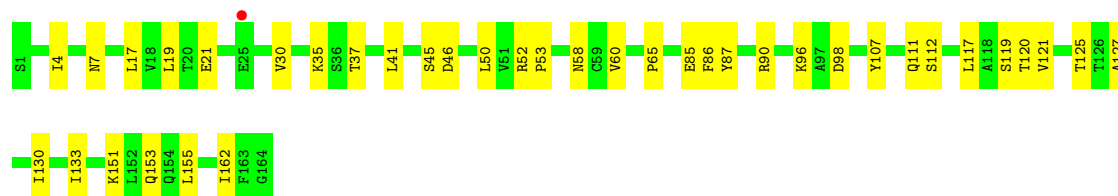
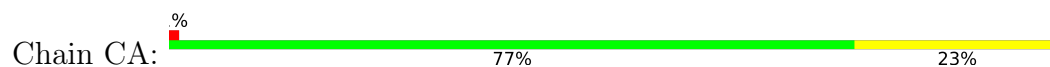
- Molecule 1: coat protein



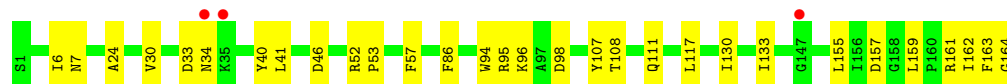
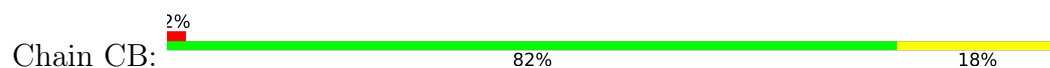
- Molecule 1: coat protein



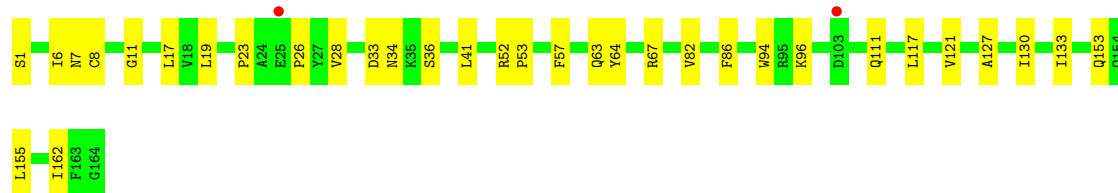
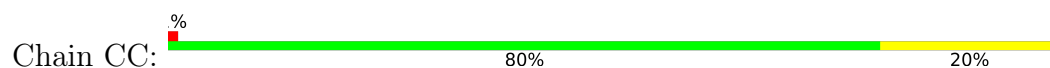
- Molecule 1: coat protein



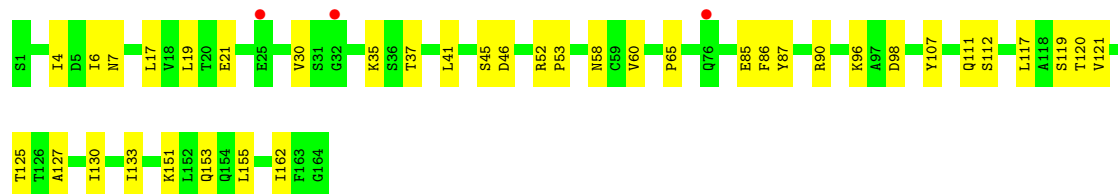
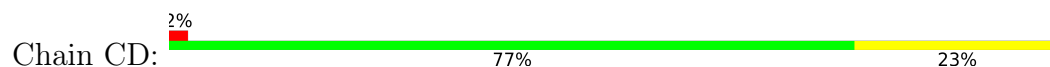
- Molecule 1: coat protein



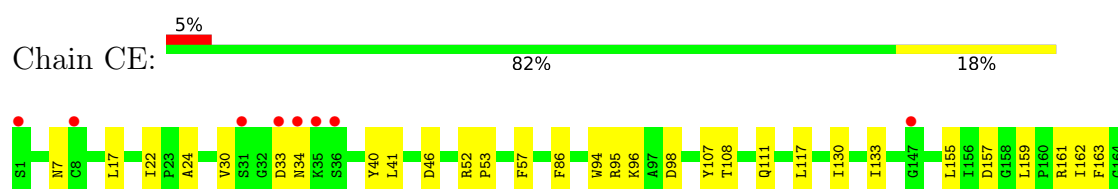
- Molecule 1: coat protein



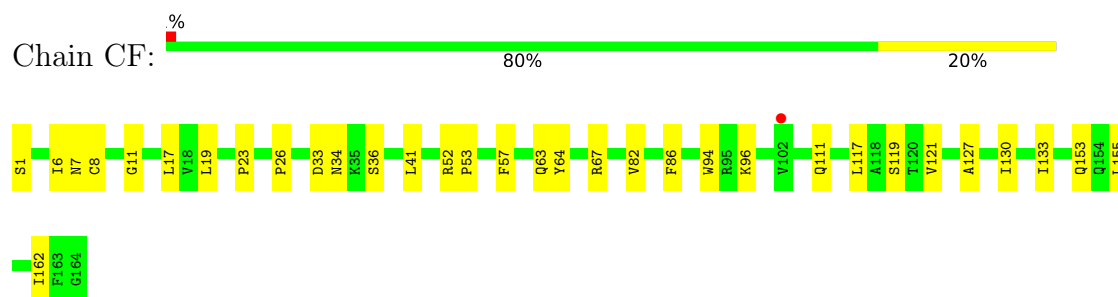
- Molecule 1: coat protein



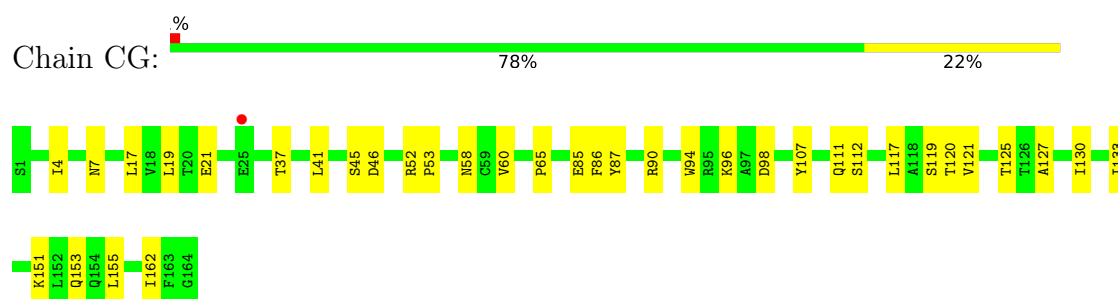
- Molecule 1: coat protein



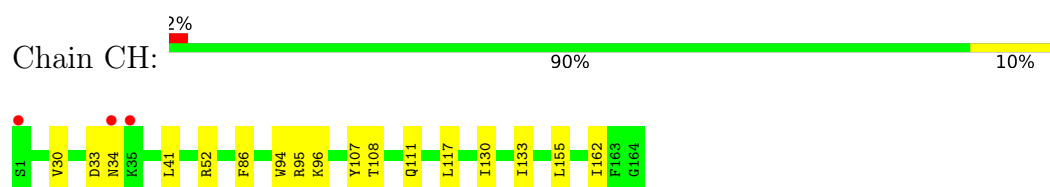
- Molecule 1: coat protein



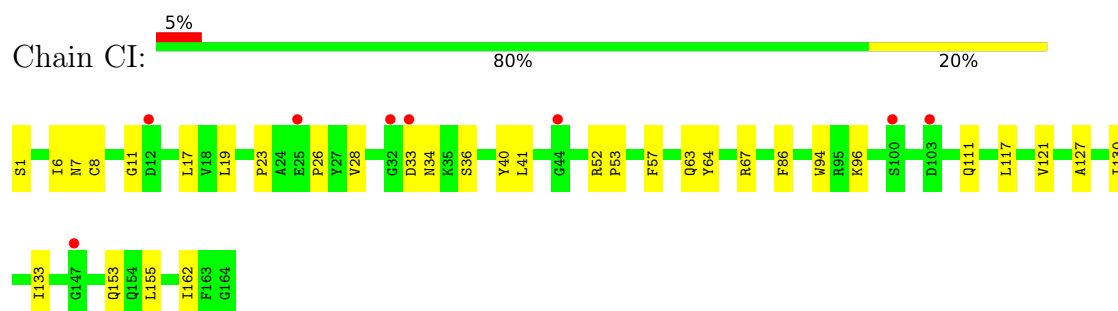
- Molecule 1: coat protein



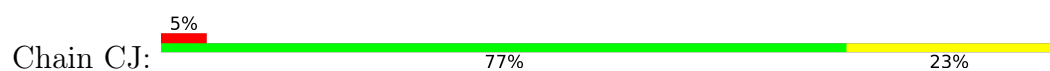
- Molecule 1: coat protein

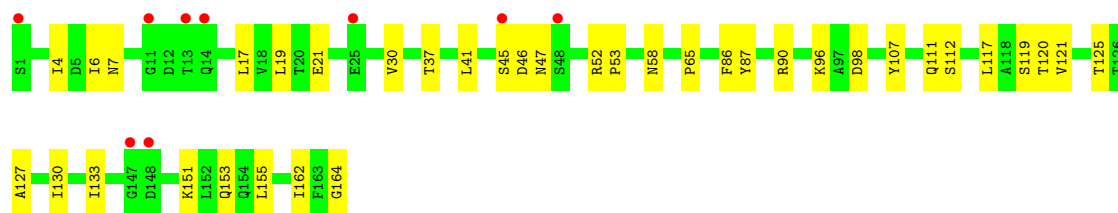


- Molecule 1: coat protein

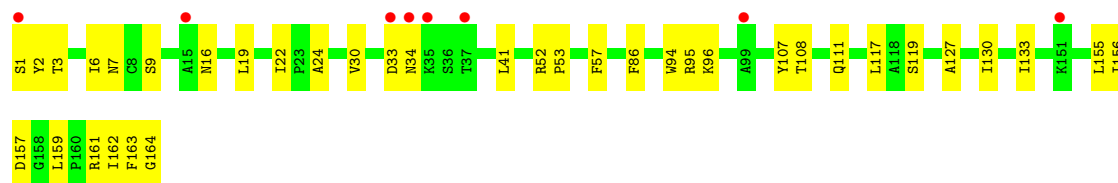
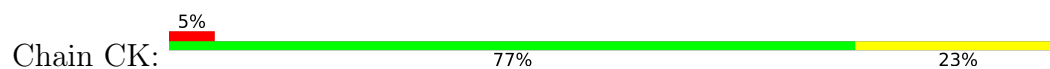


- Molecule 1: coat protein

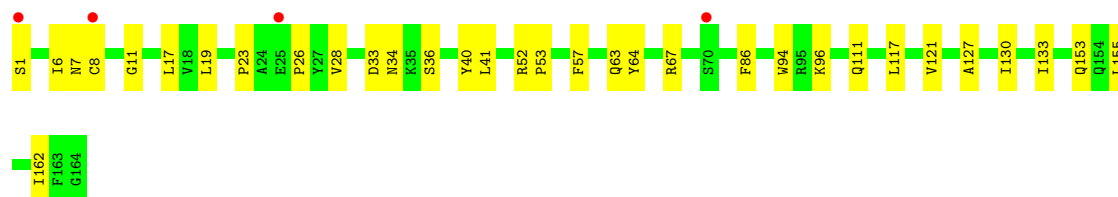
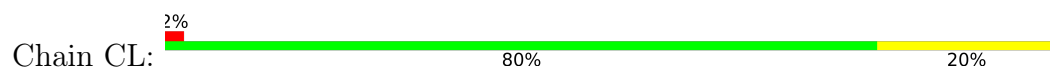




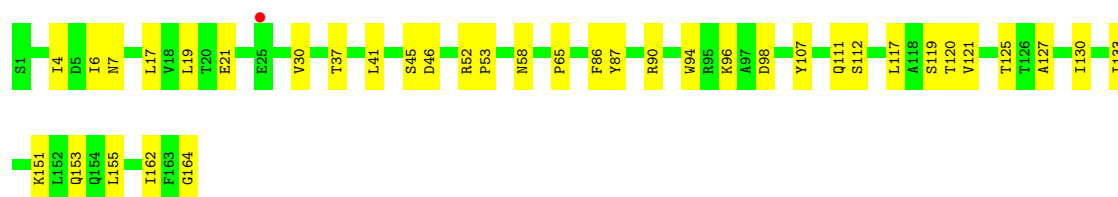
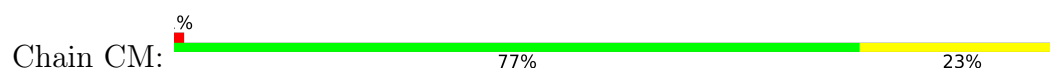
• Molecule 1: coat protein



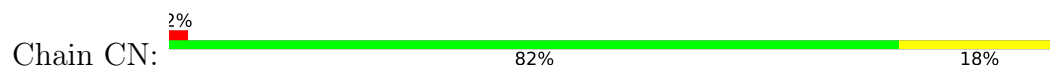
• Molecule 1: coat protein



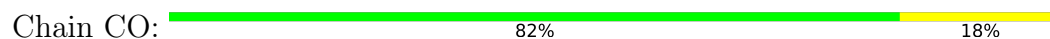
• Molecule 1: coat protein



• Molecule 1: coat protein

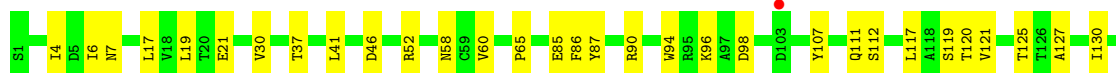
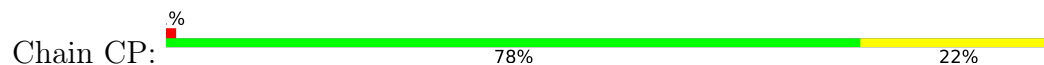


• Molecule 1: coat protein

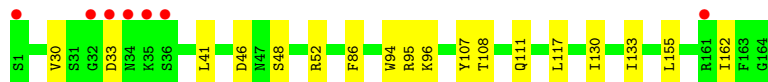
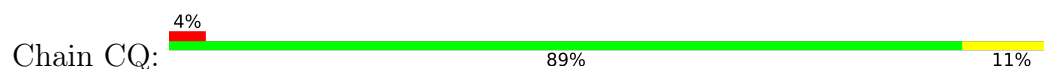




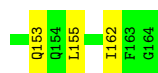
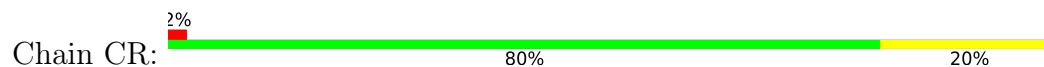
- Molecule 1: coat protein



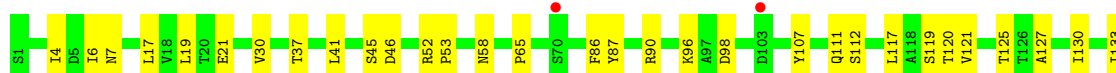
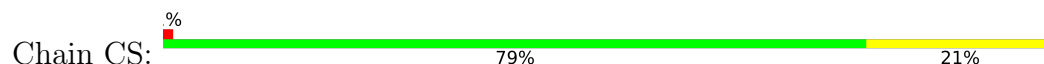
- Molecule 1: coat protein



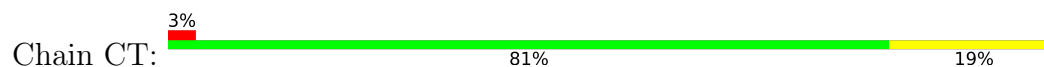
- Molecule 1: coat protein



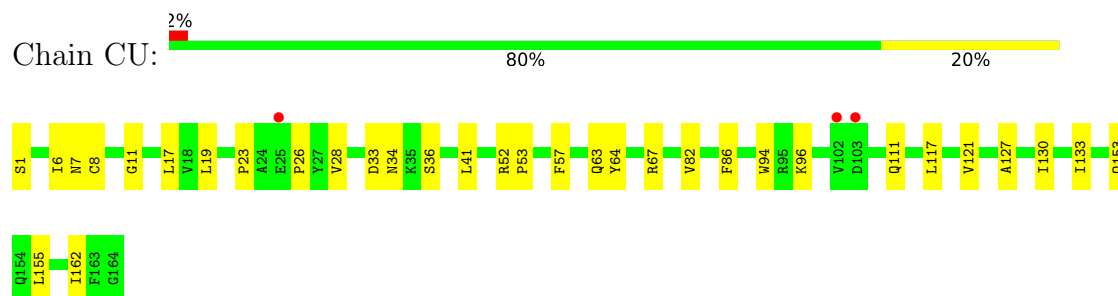
- Molecule 1: coat protein



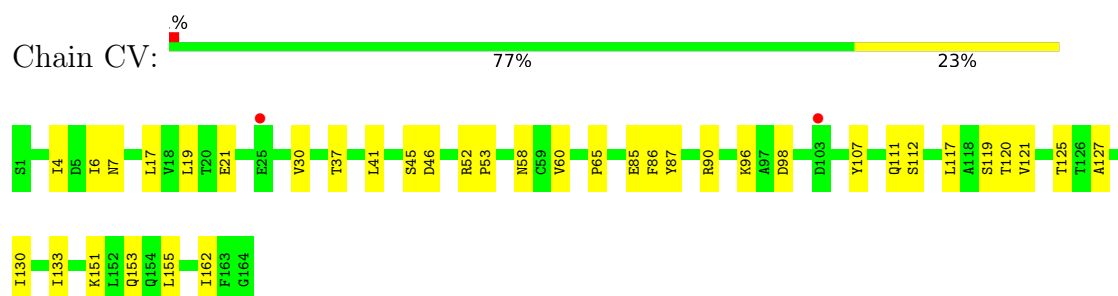
- Molecule 1: coat protein



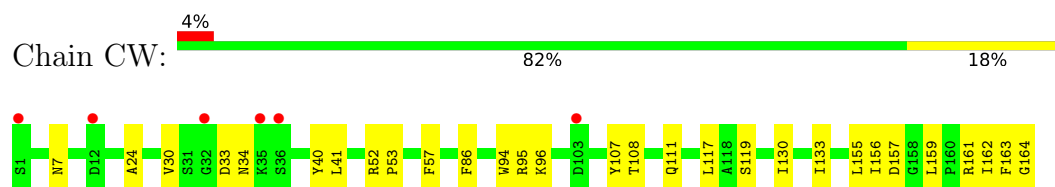
- Molecule 1: coat protein



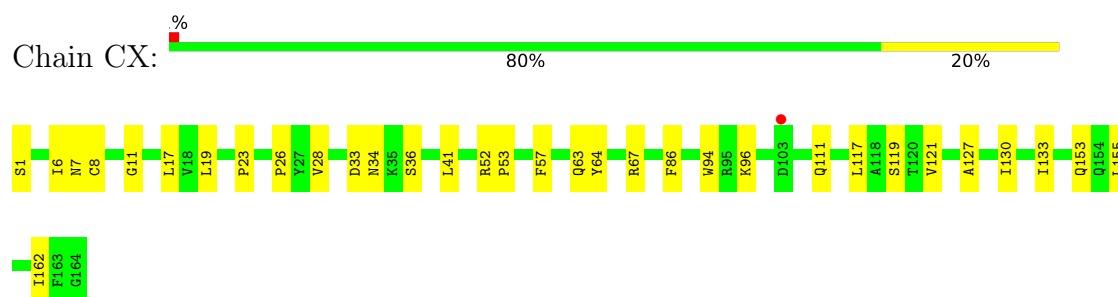
- Molecule 1: coat protein



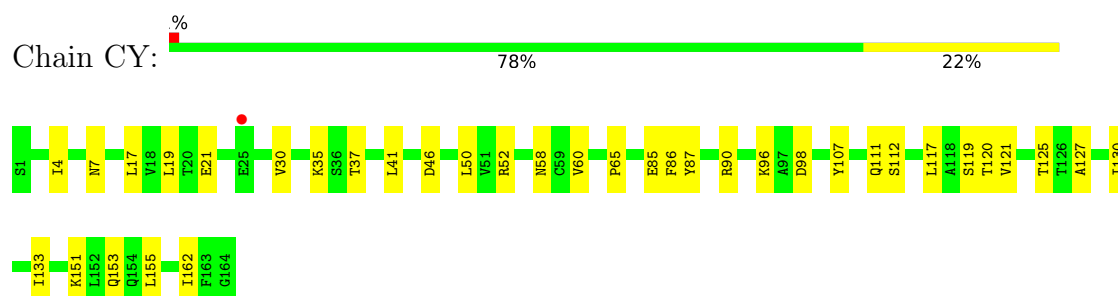
- Molecule 1: coat protein



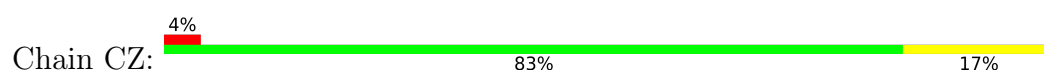
- Molecule 1: coat protein



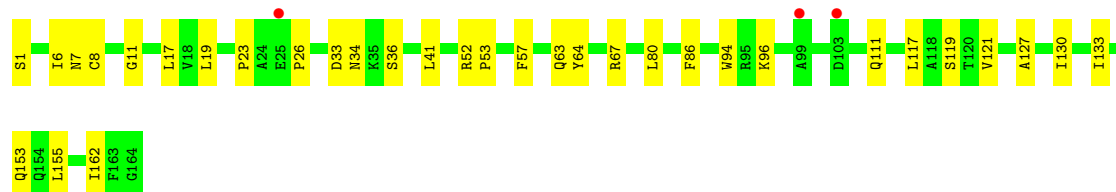
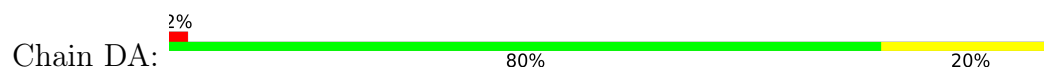
- Molecule 1: coat protein



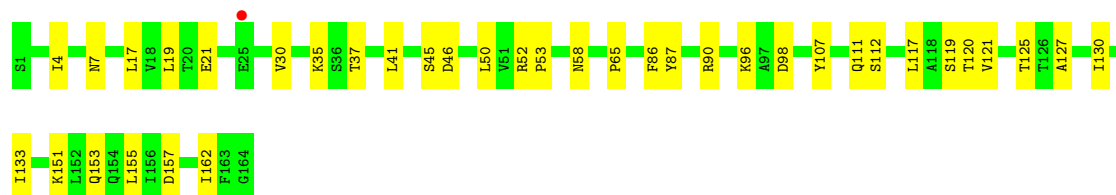
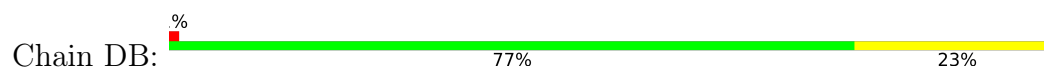
- Molecule 1: coat protein



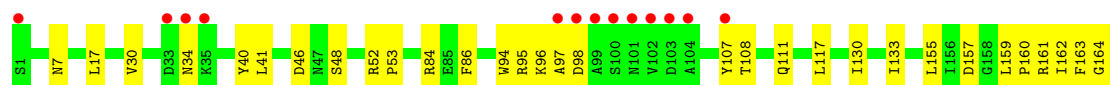
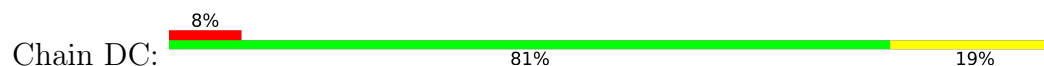
- Molecule 1: coat protein



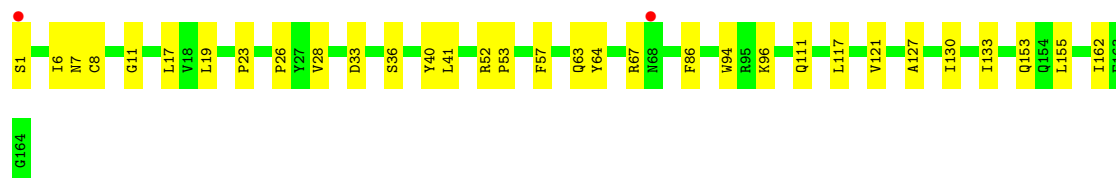
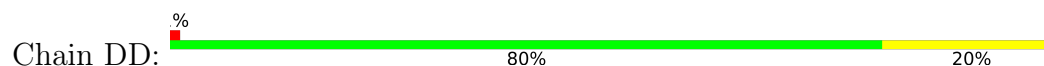
- Molecule 1: coat protein



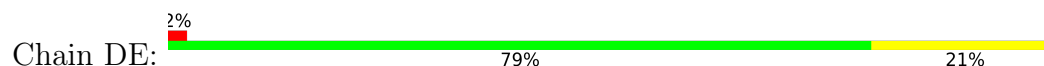
- Molecule 1: coat protein

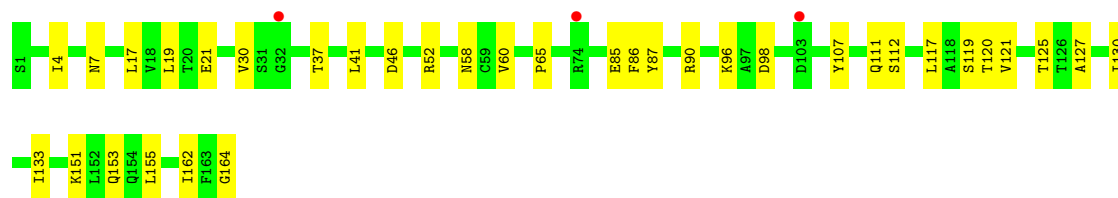


- Molecule 1: coat protein

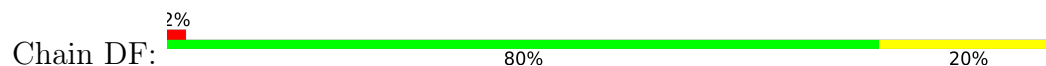


- Molecule 1: coat protein

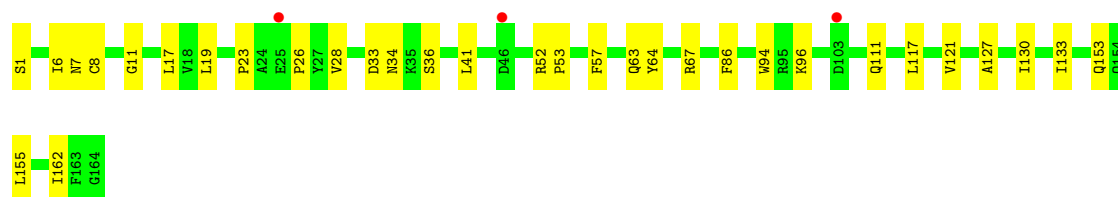
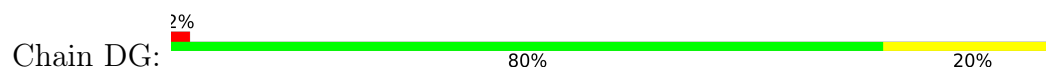




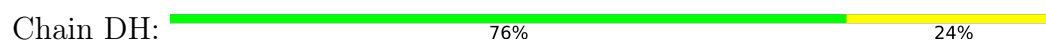
- Molecule 1: coat protein



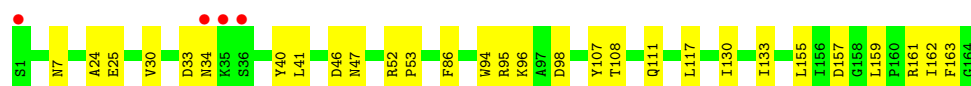
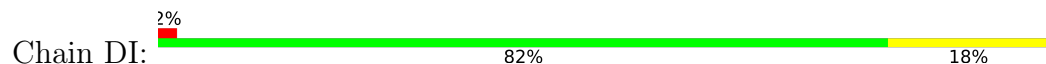
- Molecule 1: coat protein



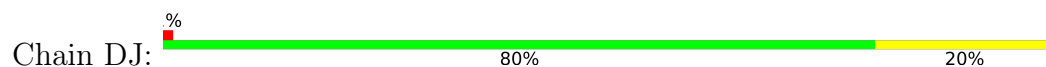
- Molecule 1: coat protein



- Molecule 1: coat protein

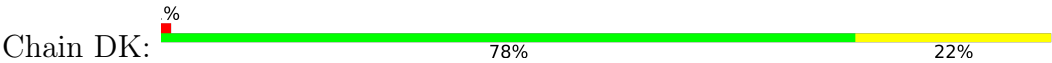


- Molecule 1: coat protein

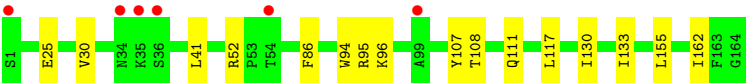
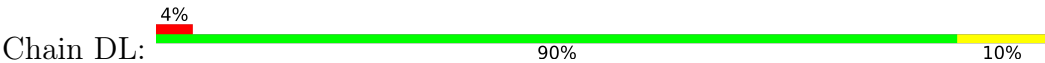


G164

• Molecule 1: coat protein



• Molecule 1: coat protein



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	469.43Å 332.41Å 293.53Å 90.00° 127.87° 90.00°	Depositor
Resolution (Å)	57.20 – 3.10 57.20 – 3.10	Depositor EDS
% Data completeness (in resolution range)	78.5 (57.20-3.10) 79.1 (57.20-3.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.31	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 3.13Å)	Xtriage
Refinement program	PHENIX 1.17.1_3660	Depositor
R, R_{free}	0.250 , 0.254 0.243 , 0.247	Depositor DCC
R_{free} test set	9970 reflections (1.96%)	wwPDB-VP
Wilson B-factor (Å ²)	41.6	Xtriage
Anisotropy	0.134	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 22.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.000 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.81	EDS
Total number of atoms	116370	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 1.76% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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	0.29	0/1321	0.55	0/1802
1	AB	0.28	0/1321	0.54	0/1802
1	AC	0.28	0/1321	0.53	0/1802
1	AD	0.29	0/1321	0.55	0/1802
1	AE	0.28	0/1321	0.54	0/1802
1	AF	0.28	0/1321	0.53	0/1802
1	AG	0.29	0/1321	0.55	0/1802
1	AH	0.28	0/1321	0.54	0/1802
1	AI	0.28	0/1321	0.53	0/1802
1	AJ	0.29	0/1321	0.55	0/1802
1	AK	0.28	0/1321	0.54	0/1802
1	AL	0.28	0/1321	0.53	0/1802
1	AM	0.29	0/1321	0.55	0/1802
1	AN	0.28	0/1321	0.54	0/1802
1	AO	0.28	0/1321	0.53	0/1802
1	AP	0.29	0/1321	0.55	0/1802
1	AQ	0.29	0/1321	0.54	0/1802
1	AR	0.28	0/1321	0.53	0/1802
1	AS	0.29	0/1321	0.55	0/1802
1	AT	0.28	0/1321	0.54	0/1802
1	AU	0.28	0/1321	0.53	0/1802
1	AV	0.29	0/1321	0.55	0/1802
1	AW	0.28	0/1321	0.54	0/1802
1	AX	0.28	0/1321	0.53	0/1802
1	AY	0.29	0/1321	0.55	0/1802
1	AZ	0.28	0/1321	0.54	0/1802
1	BA	0.28	0/1321	0.53	0/1802
1	BB	0.30	0/1321	0.55	0/1802
1	BC	0.28	0/1321	0.54	0/1802
1	BD	0.28	0/1321	0.53	0/1802
1	BE	0.29	0/1321	0.55	0/1802
1	BF	0.28	0/1321	0.53	0/1802
1	BG	0.28	0/1321	0.53	0/1802
1	BH	0.29	0/1321	0.55	0/1802

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	BI	0.28	0/1321	0.54	0/1802
1	BJ	0.28	0/1321	0.53	0/1802
1	BK	0.29	0/1321	0.55	0/1802
1	BL	0.29	0/1321	0.54	0/1802
1	BM	0.28	0/1321	0.53	0/1802
1	BN	0.29	0/1321	0.55	0/1802
1	BO	0.28	0/1321	0.54	0/1802
1	BP	0.28	0/1321	0.53	0/1802
1	BQ	0.29	0/1321	0.55	0/1802
1	BR	0.29	0/1321	0.53	0/1802
1	BS	0.28	0/1321	0.53	0/1802
1	BT	0.29	0/1321	0.55	0/1802
1	BU	0.28	0/1321	0.54	0/1802
1	BV	0.28	0/1321	0.53	0/1802
1	BW	0.29	0/1321	0.55	0/1802
1	BX	0.28	0/1321	0.54	0/1802
1	BY	0.28	0/1321	0.53	0/1802
1	BZ	0.30	0/1321	0.55	0/1802
1	CA	0.28	0/1321	0.53	0/1802
1	CB	0.28	0/1321	0.53	0/1802
1	CC	0.29	0/1321	0.55	0/1802
1	CD	0.28	0/1321	0.54	0/1802
1	CE	0.28	0/1321	0.53	0/1802
1	CF	0.29	0/1321	0.55	0/1802
1	CG	0.28	0/1321	0.54	0/1802
1	CH	0.28	0/1321	0.53	0/1802
1	CI	0.29	0/1321	0.55	0/1802
1	CJ	0.29	0/1321	0.54	0/1802
1	CK	0.28	0/1321	0.53	0/1802
1	CL	0.29	0/1321	0.55	0/1802
1	CM	0.29	0/1321	0.54	0/1802
1	CN	0.28	0/1321	0.53	0/1802
1	CO	0.29	0/1321	0.55	0/1802
1	CP	0.28	0/1321	0.54	0/1802
1	CQ	0.28	0/1321	0.53	0/1802
1	CR	0.29	0/1321	0.55	0/1802
1	CS	0.28	0/1321	0.54	0/1802
1	CT	0.28	0/1321	0.53	0/1802
1	CU	0.29	0/1321	0.55	0/1802
1	CV	0.28	0/1321	0.54	0/1802
1	CW	0.28	0/1321	0.53	0/1802
1	CX	0.29	0/1321	0.55	0/1802
1	CY	0.28	0/1321	0.54	0/1802

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	CZ	0.28	0/1321	0.53	0/1802
1	DA	0.29	0/1321	0.55	0/1802
1	DB	0.29	0/1321	0.53	0/1802
1	DC	0.28	0/1321	0.53	0/1802
1	DD	0.29	0/1321	0.55	0/1802
1	DE	0.29	0/1321	0.54	0/1802
1	DF	0.28	0/1321	0.53	0/1802
1	DG	0.29	0/1321	0.55	0/1802
1	DH	0.29	0/1321	0.54	0/1802
1	DI	0.28	0/1321	0.53	0/1802
1	DJ	0.29	0/1321	0.55	0/1802
1	DK	0.28	0/1321	0.54	0/1802
1	DL	0.28	0/1321	0.53	0/1802
All	All	0.29	0/118890	0.54	0/162180

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	1293	0	1269	36	0
1	AB	1293	0	1269	35	0
1	AC	1293	0	1269	28	0
1	AD	1293	0	1269	39	0
1	AE	1293	0	1269	41	0
1	AF	1293	0	1269	31	0
1	AG	1293	0	1269	36	0
1	AH	1293	0	1269	37	0
1	AI	1293	0	1269	27	0
1	AJ	1293	0	1269	35	0
1	AK	1293	0	1269	37	0
1	AL	1293	0	1269	30	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	AM	1293	0	1269	35	0
1	AN	1293	0	1269	33	0
1	AO	1293	0	1269	30	0
1	AP	1293	0	1269	33	0
1	AQ	1293	0	1269	32	0
1	AR	1293	0	1269	44	0
1	AS	1293	0	1269	35	0
1	AT	1293	0	1269	36	0
1	AU	1293	0	1269	34	0
1	AV	1293	0	1269	35	0
1	AW	1293	0	1269	35	0
1	AX	1293	0	1269	28	0
1	AY	1293	0	1269	33	0
1	AZ	1293	0	1269	31	0
1	BA	1293	0	1269	29	1
1	BB	1293	0	1269	35	0
1	BC	1293	0	1269	36	0
1	BD	1293	0	1269	29	0
1	BE	1293	0	1269	36	0
1	BF	1293	0	1269	36	0
1	BG	1293	0	1269	43	0
1	BH	1293	0	1269	37	0
1	BI	1293	0	1269	33	0
1	BJ	1293	0	1269	34	0
1	BK	1293	0	1269	33	0
1	BL	1293	0	1269	36	0
1	BM	1293	0	1269	27	0
1	BN	1293	0	1269	34	0
1	BO	1293	0	1269	32	0
1	BP	1293	0	1269	28	0
1	BQ	1293	0	1269	35	1
1	BR	1293	0	1269	32	0
1	BS	1293	0	1269	11	0
1	BT	1293	0	1269	35	0
1	BU	1293	0	1269	36	0
1	BV	1293	0	1269	31	0
1	BW	1293	0	1269	36	0
1	BX	1293	0	1269	34	0
1	BY	1293	0	1269	27	0
1	BZ	1293	0	1269	33	0
1	CA	1293	0	1269	36	0
1	CB	1293	0	1269	30	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	CC	1293	0	1269	36	0
1	CD	1293	0	1269	34	0
1	CE	1293	0	1269	30	0
1	CF	1293	0	1269	35	0
1	CG	1293	0	1269	34	0
1	CH	1293	0	1269	11	0
1	CI	1293	0	1269	34	0
1	CJ	1293	0	1269	35	0
1	CK	1293	0	1269	46	0
1	CL	1293	0	1269	36	0
1	CM	1293	0	1269	34	0
1	CN	1293	0	1269	32	0
1	CO	1293	0	1269	31	0
1	CP	1293	0	1269	34	0
1	CQ	1293	0	1269	11	0
1	CR	1293	0	1269	34	0
1	CS	1293	0	1269	34	0
1	CT	1293	0	1269	30	0
1	CU	1293	0	1269	36	0
1	CV	1293	0	1269	34	0
1	CW	1293	0	1269	30	0
1	CX	1293	0	1269	34	0
1	CY	1293	0	1269	36	0
1	CZ	1293	0	1269	30	0
1	DA	1293	0	1269	35	0
1	DB	1293	0	1269	36	0
1	DC	1293	0	1269	36	0
1	DD	1293	0	1269	34	0
1	DE	1293	0	1269	33	0
1	DF	1293	0	1269	33	0
1	DG	1293	0	1269	37	0
1	DH	1293	0	1269	35	0
1	DI	1293	0	1269	29	0
1	DJ	1293	0	1269	34	0
1	DK	1293	0	1269	34	0
1	DL	1293	0	1269	11	0
All	All	116370	0	114210	2061	1

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 (2061) 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:AR:53:PRO:HG2	1:CK:7:ASN:O	1.68	0.92
1:BJ:53:PRO:HG2	1:DF:7:ASN:O	1.81	0.80
1:AR:34:ASN:HB3	1:CK:161:ARG:HD2	1.65	0.79
1:BG:161:ARG:HD2	1:DC:34:ASN:HB3	1.63	0.79
1:BG:157:ASP:CG	1:DA:52:ARG:HH22	1.91	0.78
1:AX:163:PHE:HD2	1:CT:24:ALA:HA	1.49	0.77
1:AF:34:ASN:HB3	1:BV:161:ARG:HD2	1.66	0.77
1:AF:53:PRO:HG2	1:BV:7:ASN:O	1.85	0.75
1:AR:159:LEU:HD13	1:CI:28:VAL:HG11	1.69	0.74
1:AX:159:LEU:HD13	1:CR:28:VAL:HG11	1.70	0.74
1:AD:52:ARG:HH22	1:BV:157:ASP:CG	1.94	0.74
1:AR:98:ASP:HB3	1:CK:1:SER:C	2.13	0.73
1:AX:153:GLN:OE1	1:CT:127:ALA:HB2	1.89	0.72
1:BF:30:VAL:HA	1:BQ:34:ASN:ND2	2.04	0.72
1:BG:34:ASN:HB3	1:DC:161:ARG:HD2	1.71	0.72
1:BJ:7:ASN:O	1:DF:53:PRO:HG2	1.89	0.71
1:AO:7:ASN:O	1:CE:53:PRO:HG2	1.90	0.71
1:AM:28:VAL:HG11	1:CE:159:LEU:HD13	1.74	0.70
1:AL:7:ASN:O	1:CB:53:PRO:HG2	1.93	0.69
1:BD:41:LEU:HD11	1:BD:52:ARG:HB2	1.75	0.69
1:CZ:41:LEU:HD11	1:CZ:52:ARG:HB2	1.75	0.69
1:AC:41:LEU:HD11	1:AC:52:ARG:HB2	1.75	0.69
1:AL:41:LEU:HD11	1:AL:52:ARG:HB2	1.75	0.69
1:DI:41:LEU:HD11	1:DI:52:ARG:HB2	1.74	0.69
1:AI:41:LEU:HD11	1:AI:52:ARG:HB2	1.75	0.69
1:CT:41:LEU:HD11	1:CT:52:ARG:HB2	1.75	0.69
1:AR:41:LEU:HD11	1:AR:52:ARG:HB2	1.75	0.69
1:CN:41:LEU:HD11	1:CN:52:ARG:HB2	1.75	0.69
1:CB:41:LEU:HD11	1:CB:52:ARG:HB2	1.75	0.69
1:CH:41:LEU:HD11	1:CH:52:ARG:HB2	1.75	0.69
1:DC:41:LEU:HD11	1:DC:52:ARG:HB2	1.75	0.69
1:AO:41:LEU:HD11	1:AO:52:ARG:HB2	1.75	0.69
1:BA:41:LEU:HD11	1:BA:52:ARG:HB2	1.75	0.69
1:BP:41:LEU:HD11	1:BP:52:ARG:HB2	1.75	0.69
1:BS:41:LEU:HD11	1:BS:52:ARG:HB2	1.75	0.69
1:AF:41:LEU:HD11	1:AF:52:ARG:HB2	1.75	0.69
1:BV:41:LEU:HD11	1:BV:52:ARG:HB2	1.75	0.69
1:CQ:41:LEU:HD11	1:CQ:52:ARG:HB2	1.75	0.69
1:BJ:41:LEU:HD11	1:BJ:52:ARG:HB2	1.75	0.68
1:DF:41:LEU:HD11	1:DF:52:ARG:HB2	1.74	0.68
1:DL:41:LEU:HD11	1:DL:52:ARG:HB2	1.75	0.68
1:AR:94:TRP:NE1	1:CK:6:ILE:HG22	2.08	0.68

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BM:41:LEU:HD11	1:BM:52:ARG:HB2	1.75	0.68
1:AI:7:ASN:O	1:BY:53:PRO:HG2	1.93	0.68
1:BY:41:LEU:HD11	1:BY:52:ARG:HB2	1.75	0.68
1:AY:52:ARG:HH22	1:CW:157:ASP:CG	2.01	0.68
1:CW:41:LEU:HD11	1:CW:52:ARG:HB2	1.75	0.68
1:AU:41:LEU:HD11	1:AU:52:ARG:HB2	1.75	0.67
1:BG:7:ASN:O	1:DC:53:PRO:HG2	1.94	0.67
1:CE:41:LEU:HD11	1:CE:52:ARG:HB2	1.75	0.67
1:AP:52:ARG:HH22	1:CK:157:ASP:CG	2.02	0.67
1:BG:41:LEU:HD11	1:BG:52:ARG:HB2	1.75	0.67
1:BM:7:ASN:O	1:DI:53:PRO:HG2	1.95	0.67
1:CK:41:LEU:HD11	1:CK:52:ARG:HB2	1.75	0.67
1:AX:24:ALA:HA	1:CT:163:PHE:HD2	1.59	0.67
1:AR:40:TYR:CE1	1:CK:163:PHE:HB2	2.30	0.67
1:AX:41:LEU:HD11	1:AX:52:ARG:HB2	1.75	0.67
1:AF:159:LEU:HD13	1:BT:28:VAL:HG11	1.75	0.66
1:BM:157:ASP:CG	1:DG:52:ARG:HH22	2.03	0.66
1:AV:28:VAL:HG11	1:CT:159:LEU:HD13	1.77	0.66
1:AX:127:ALA:HB2	1:CT:153:GLN:OE1	1.96	0.66
1:BH:34:ASN:ND2	1:BL:30:VAL:HA	2.10	0.66
1:AR:163:PHE:HD2	1:CK:24:ALA:HA	1.60	0.65
1:AF:163:PHE:HD2	1:BV:24:ALA:HA	1.62	0.65
1:AX:163:PHE:CD2	1:CT:24:ALA:HA	2.32	0.65
1:BK:28:VAL:HG11	1:DI:159:LEU:HD13	1.79	0.64
1:AL:53:PRO:HG2	1:CB:7:ASN:O	1.97	0.64
1:BD:7:ASN:O	1:CZ:53:PRO:HG2	1.97	0.64
1:BA:53:PRO:HG2	1:CW:7:ASN:O	1.98	0.64
1:BA:34:ASN:HB3	1:CW:161:ARG:HD2	1.77	0.64
1:BA:161:ARG:HD2	1:CW:34:ASN:HB3	1.79	0.63
1:BD:53:PRO:HG2	1:CZ:7:ASN:O	1.99	0.63
1:AO:24:ALA:HA	1:CE:163:PHE:HD2	1.62	0.63
1:AU:34:ASN:HB3	1:CN:161:ARG:HD2	1.80	0.63
1:CA:30:VAL:HA	1:CF:34:ASN:ND2	2.13	0.63
1:AI:53:PRO:HG2	1:BY:7:ASN:O	1.99	0.63
1:AC:7:ASN:O	1:BP:53:PRO:HG2	2.00	0.62
1:AG:7:ASN:HD21	1:AH:46:ASP:H	1.48	0.62
1:AH:30:VAL:HA	1:CC:34:ASN:ND2	2.15	0.62
1:AI:157:ASP:CG	1:BW:52:ARG:HH22	2.07	0.62
1:AR:96:LYS:HA	1:CK:3:THR:O	2.00	0.62
1:CL:7:ASN:HD21	1:CM:46:ASP:H	1.48	0.61
1:AJ:7:ASN:HD21	1:AK:46:ASP:H	1.49	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BZ:7:ASN:HD21	1:CA:46:ASP:H	1.48	0.61
1:AD:7:ASN:HD21	1:AE:46:ASP:H	1.49	0.61
1:AE:35:LYS:NZ	1:BG:34:ASN:HD21	1.97	0.61
1:AS:7:ASN:HD21	1:AT:46:ASP:H	1.48	0.61
1:CC:7:ASN:HD21	1:CD:46:ASP:H	1.48	0.61
1:CO:7:ASN:HD21	1:CP:46:ASP:H	1.48	0.61
1:AP:7:ASN:HD21	1:AQ:46:ASP:H	1.49	0.61
1:BE:28:VAL:HG11	1:DC:159:LEU:HD13	1.82	0.61
1:BG:7:ASN:HD21	1:DC:46:ASP:H	1.48	0.61
1:AK:65:PRO:HB2	1:AK:120:THR:HG22	1.83	0.61
1:AY:7:ASN:HD21	1:AZ:46:ASP:H	1.49	0.61
1:BJ:159:LEU:HD13	1:DD:28:VAL:HG11	1.83	0.61
1:CM:65:PRO:HB2	1:CM:120:THR:HG22	1.83	0.61
1:DJ:7:ASN:HD21	1:DK:46:ASP:H	1.49	0.61
1:CA:65:PRO:HB2	1:CA:120:THR:HG22	1.83	0.61
1:CX:7:ASN:HD21	1:CY:46:ASP:H	1.49	0.61
1:AA:7:ASN:HD21	1:AB:46:ASP:H	1.49	0.61
1:AB:65:PRO:HB2	1:AB:120:THR:HG22	1.83	0.61
1:AE:65:PRO:HB2	1:AE:120:THR:HG22	1.83	0.61
1:AM:7:ASN:HD21	1:AN:46:ASP:H	1.48	0.61
1:AQ:65:PRO:HB2	1:AQ:120:THR:HG22	1.83	0.61
1:AU:53:PRO:HG2	1:CN:7:ASN:O	2.01	0.61
1:BH:7:ASN:HD21	1:BI:46:ASP:H	1.49	0.61
1:DB:65:PRO:HB2	1:DB:120:THR:HG22	1.83	0.61
1:AL:161:ARG:HD2	1:CB:34:ASN:HB3	1.82	0.60
1:DK:65:PRO:HB2	1:DK:120:THR:HG22	1.83	0.60
1:BK:7:ASN:HD21	1:BL:46:ASP:H	1.48	0.60
1:CS:65:PRO:HB2	1:CS:120:THR:HG22	1.83	0.60
1:CU:7:ASN:HD21	1:CV:46:ASP:H	1.49	0.60
1:AN:65:PRO:HB2	1:AN:120:THR:HG22	1.83	0.60
1:BE:7:ASN:HD21	1:BF:46:ASP:H	1.48	0.60
1:CR:7:ASN:HD21	1:CS:46:ASP:H	1.48	0.60
1:DG:130:ILE:O	1:DG:133:ILE:HG22	2.02	0.60
1:BB:7:ASN:HD21	1:BC:46:ASP:H	1.49	0.60
1:BF:65:PRO:HB2	1:BF:120:THR:HG22	1.83	0.60
1:BH:130:ILE:O	1:BH:133:ILE:HG22	2.02	0.60
1:BI:65:PRO:HB2	1:BI:120:THR:HG22	1.83	0.60
1:CF:7:ASN:HD21	1:CG:46:ASP:H	1.48	0.60
1:CL:130:ILE:O	1:CL:133:ILE:HG22	2.02	0.60
1:DH:65:PRO:HB2	1:DH:120:THR:HG22	1.83	0.60
1:AH:65:PRO:HB2	1:AH:120:THR:HG22	1.83	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AJ:28:VAL:HG11	1:CB:159:LEU:HD13	1.84	0.60
1:AS:130:ILE:O	1:AS:133:ILE:HG22	2.02	0.60
1:BA:159:LEU:HD13	1:CU:28:VAL:HG11	1.84	0.60
1:BM:161:ARG:HD2	1:DI:34:ASN:HB3	1.83	0.60
1:CX:130:ILE:O	1:CX:133:ILE:HG22	2.02	0.60
1:AO:157:ASP:CG	1:CC:52:ARG:HH22	2.10	0.60
1:CU:130:ILE:O	1:CU:133:ILE:HG22	2.02	0.60
1:AL:96:LYS:HG2	1:AL:107:TYR:HB2	1.84	0.60
1:BC:65:PRO:HB2	1:BC:120:THR:HG22	1.83	0.60
1:BR:65:PRO:HB2	1:BR:120:THR:HG22	1.83	0.60
1:CY:65:PRO:HB2	1:CY:120:THR:HG22	1.83	0.60
1:AB:41:LEU:HD11	1:AB:52:ARG:HB2	1.84	0.60
1:AD:130:ILE:O	1:AD:133:ILE:HG22	2.02	0.60
1:BA:96:LYS:HG2	1:BA:107:TYR:HB2	1.84	0.60
1:CN:96:LYS:HG2	1:CN:107:TYR:HB2	1.84	0.60
1:CO:130:ILE:O	1:CO:133:ILE:HG22	2.02	0.60
1:DD:7:ASN:HD21	1:DE:46:ASP:H	1.49	0.60
1:DH:41:LEU:HD11	1:DH:52:ARG:HB2	1.84	0.60
1:AE:41:LEU:HD11	1:AE:52:ARG:HB2	1.84	0.60
1:AF:96:LYS:HG2	1:AF:107:TYR:HB2	1.84	0.60
1:AM:33:ASP:H	1:AM:36:SER:HB3	1.67	0.60
1:AW:65:PRO:HB2	1:AW:120:THR:HG22	1.83	0.60
1:BD:96:LYS:HG2	1:BD:107:TYR:HB2	1.84	0.60
1:BW:33:ASP:H	1:BW:36:SER:HB3	1.67	0.60
1:CG:65:PRO:HB2	1:CG:120:THR:HG22	1.83	0.60
1:DI:96:LYS:HG2	1:DI:107:TYR:HB2	1.84	0.60
1:AA:130:ILE:O	1:AA:133:ILE:HG22	2.02	0.60
1:AJ:33:ASP:H	1:AJ:36:SER:HB3	1.67	0.60
1:AJ:130:ILE:O	1:AJ:133:ILE:HG22	2.02	0.60
1:AM:130:ILE:O	1:AM:133:ILE:HG22	2.02	0.60
1:BB:28:VAL:HG11	1:CZ:159:LEU:HD13	1.83	0.60
1:BE:33:ASP:H	1:BE:36:SER:HB3	1.67	0.60
1:BN:33:ASP:H	1:BN:36:SER:HB3	1.67	0.60
1:BO:65:PRO:HB2	1:BO:120:THR:HG22	1.83	0.60
1:CA:41:LEU:HD11	1:CA:52:ARG:HB2	1.84	0.60
1:CU:34:ASN:ND2	1:CY:30:VAL:HA	2.17	0.60
1:DA:7:ASN:HD21	1:DB:46:ASP:H	1.49	0.60
1:DA:33:ASP:H	1:DA:36:SER:HB3	1.67	0.60
1:AH:41:LEU:HD11	1:AH:52:ARG:HB2	1.84	0.59
1:BO:41:LEU:HD11	1:BO:52:ARG:HB2	1.84	0.59
1:BR:41:LEU:HD11	1:BR:52:ARG:HB2	1.84	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CF:33:ASP:H	1:CF:36:SER:HB3	1.67	0.59
1:CG:41:LEU:HD11	1:CG:52:ARG:HB2	1.84	0.59
1:CI:130:ILE:O	1:CI:133:ILE:HG22	2.02	0.59
1:CJ:65:PRO:HB2	1:CJ:120:THR:HG22	1.83	0.59
1:CR:130:ILE:O	1:CR:133:ILE:HG22	2.02	0.59
1:AO:96:LYS:HG2	1:AO:107:TYR:HB2	1.84	0.59
1:BE:130:ILE:O	1:BE:133:ILE:HG22	2.02	0.59
1:BF:41:LEU:HD11	1:BF:52:ARG:HB2	1.84	0.59
1:BJ:94:TRP:NE1	1:DF:6:ILE:HG22	2.17	0.59
1:CB:96:LYS:HG2	1:CB:107:TYR:HB2	1.84	0.59
1:CC:130:ILE:O	1:CC:133:ILE:HG22	2.02	0.59
1:CH:96:LYS:HG2	1:CH:107:TYR:HB2	1.84	0.59
1:CM:41:LEU:HD11	1:CM:52:ARG:HB2	1.84	0.59
1:CT:96:LYS:HG2	1:CT:107:TYR:HB2	1.84	0.59
1:DE:65:PRO:HB2	1:DE:120:THR:HG22	1.83	0.59
1:DG:33:ASP:H	1:DG:36:SER:HB3	1.67	0.59
1:AD:33:ASP:H	1:AD:36:SER:HB3	1.67	0.59
1:AI:96:LYS:HG2	1:AI:107:TYR:HB2	1.84	0.59
1:AL:157:ASP:CG	1:BZ:52:ARG:HH22	2.10	0.59
1:AQ:41:LEU:HD11	1:AQ:52:ARG:HB2	1.84	0.59
1:BN:7:ASN:HD21	1:BO:46:ASP:H	1.49	0.59
1:CJ:41:LEU:HD11	1:CJ:52:ARG:HB2	1.84	0.59
1:CX:33:ASP:H	1:CX:36:SER:HB3	1.67	0.59
1:AP:130:ILE:O	1:AP:133:ILE:HG22	2.02	0.59
1:BB:130:ILE:O	1:BB:133:ILE:HG22	2.02	0.59
1:BF:30:VAL:HA	1:BQ:34:ASN:HD22	1.65	0.59
1:BI:41:LEU:HD11	1:BI:52:ARG:HB2	1.84	0.59
1:BN:130:ILE:O	1:BN:133:ILE:HG22	2.02	0.59
1:BS:96:LYS:HG2	1:BS:107:TYR:HB2	1.84	0.59
1:BT:7:ASN:HD21	1:BU:46:ASP:H	1.48	0.59
1:BW:130:ILE:O	1:BW:133:ILE:HG22	2.02	0.59
1:CD:65:PRO:HB2	1:CD:120:THR:HG22	1.83	0.59
1:CU:33:ASP:H	1:CU:36:SER:HB3	1.67	0.59
1:CV:41:LEU:HD11	1:CV:52:ARG:HB2	1.84	0.59
1:CV:65:PRO:HB2	1:CV:120:THR:HG22	1.83	0.59
1:DD:33:ASP:H	1:DD:36:SER:HB3	1.67	0.59
1:DJ:130:ILE:O	1:DJ:133:ILE:HG22	2.02	0.59
1:AC:53:PRO:HG2	1:BP:7:ASN:O	2.02	0.59
1:AG:28:VAL:HG11	1:BY:159:LEU:HD13	1.82	0.59
1:AG:130:ILE:O	1:AG:133:ILE:HG22	2.02	0.59
1:AS:33:ASP:H	1:AS:36:SER:HB3	1.67	0.59

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BM:53:PRO:HG2	1:DI:7:ASN:O	2.01	0.59
1:CF:130:ILE:O	1:CF:133:ILE:HG22	2.02	0.59
1:AP:33:ASP:H	1:AP:36:SER:HB3	1.67	0.59
1:AT:41:LEU:HD11	1:AT:52:ARG:HB2	1.84	0.59
1:AT:65:PRO:HB2	1:AT:120:THR:HG22	1.83	0.59
1:AU:161:ARG:HD2	1:CN:34:ASN:HB3	1.84	0.59
1:AV:130:ILE:O	1:AV:133:ILE:HG22	2.02	0.59
1:BD:157:ASP:CG	1:CX:52:ARG:HH22	2.10	0.59
1:BG:163:PHE:HB2	1:DC:40:TYR:CE1	2.37	0.59
1:BL:65:PRO:HB2	1:BL:120:THR:HG22	1.83	0.59
1:BT:33:ASP:H	1:BT:36:SER:HB3	1.67	0.59
1:BT:130:ILE:O	1:BT:133:ILE:HG22	2.02	0.59
1:BZ:130:ILE:O	1:BZ:133:ILE:HG22	2.02	0.59
1:CP:65:PRO:HB2	1:CP:120:THR:HG22	1.83	0.59
1:CS:41:LEU:HD11	1:CS:52:ARG:HB2	1.84	0.59
1:CY:41:LEU:HD11	1:CY:52:ARG:HB2	1.84	0.59
1:DD:130:ILE:O	1:DD:133:ILE:HG22	2.02	0.59
1:DA:130:ILE:O	1:DA:133:ILE:HG22	2.02	0.59
1:DB:41:LEU:HD11	1:DB:52:ARG:HB2	1.84	0.59
1:DF:96:LYS:HG2	1:DF:107:TYR:HB2	1.84	0.59
1:AL:34:ASN:HB3	1:CB:161:ARG:HD2	1.85	0.59
1:AZ:65:PRO:HB2	1:AZ:120:THR:HG22	1.83	0.59
1:BK:130:ILE:O	1:BK:133:ILE:HG22	2.02	0.59
1:BQ:7:ASN:HD21	1:BR:46:ASP:H	1.48	0.59
1:BV:96:LYS:HG2	1:BV:107:TYR:HB2	1.84	0.59
1:CK:96:LYS:HG2	1:CK:107:TYR:HB2	1.84	0.59
1:CR:33:ASP:H	1:CR:36:SER:HB3	1.67	0.59
1:AY:130:ILE:O	1:AY:133:ILE:HG22	2.02	0.59
1:BB:33:ASP:H	1:BB:36:SER:HB3	1.67	0.59
1:BJ:96:LYS:HG2	1:BJ:107:TYR:HB2	1.84	0.59
1:BQ:130:ILE:O	1:BQ:133:ILE:HG22	2.02	0.59
1:BW:7:ASN:HD21	1:BX:46:ASP:H	1.48	0.59
1:AU:159:LEU:HD13	1:CL:28:VAL:HG11	1.83	0.59
1:AV:7:ASN:HD21	1:AW:46:ASP:H	1.49	0.59
1:AX:96:LYS:HG2	1:AX:107:TYR:HB2	1.84	0.59
1:BD:161:ARG:HD2	1:CZ:34:ASN:HB3	1.85	0.59
1:BM:96:LYS:HG2	1:BM:107:TYR:HB2	1.84	0.59
1:BX:65:PRO:HB2	1:BX:120:THR:HG22	1.83	0.59
1:CC:33:ASP:H	1:CC:36:SER:HB3	1.67	0.59
1:DK:41:LEU:HD11	1:DK:52:ARG:HB2	1.84	0.59
1:AC:96:LYS:HG2	1:AC:107:TYR:HB2	1.84	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AR:7:ASN:O	1:CK:53:PRO:HG2	2.03	0.58
1:BU:65:PRO:HB2	1:BU:120:THR:HG22	1.83	0.58
1:CO:33:ASP:H	1:CO:36:SER:HB3	1.67	0.58
1:CW:96:LYS:HG2	1:CW:107:TYR:HB2	1.84	0.58
1:AX:53:PRO:HG2	1:CT:7:ASN:O	2.03	0.58
1:BL:41:LEU:HD11	1:BL:52:ARG:HB2	1.84	0.58
1:BU:41:LEU:HD11	1:BU:52:ARG:HB2	1.84	0.58
1:AW:41:LEU:HD11	1:AW:52:ARG:HB2	1.84	0.58
1:BA:157:ASP:CG	1:CU:52:ARG:HH22	2.11	0.58
1:CP:41:LEU:HD11	1:CP:52:ARG:HB2	1.84	0.58
1:DL:96:LYS:HG2	1:DL:107:TYR:HB2	1.84	0.58
1:AF:40:TYR:CE1	1:BV:163:PHE:HB2	2.38	0.58
1:AU:96:LYS:HG2	1:AU:107:TYR:HB2	1.84	0.58
1:AY:33:ASP:H	1:AY:36:SER:HB3	1.67	0.58
1:BF:50:LEU:HD11	1:BQ:82:VAL:HG13	1.85	0.58
1:CL:33:ASP:H	1:CL:36:SER:HB3	1.67	0.58
1:AO:161:ARG:HD2	1:CE:34:ASN:HB3	1.85	0.58
1:AR:96:LYS:HG2	1:AR:107:TYR:HB2	1.84	0.58
1:AR:98:ASP:OD2	1:CK:1:SER:OG	2.13	0.58
1:BQ:33:ASP:H	1:BQ:36:SER:HB3	1.67	0.58
1:CQ:96:LYS:HG2	1:CQ:107:TYR:HB2	1.84	0.58
1:DG:7:ASN:HD21	1:DH:46:ASP:H	1.48	0.58
1:DJ:33:ASP:H	1:DJ:36:SER:HB3	1.67	0.58
1:AA:33:ASP:H	1:AA:36:SER:HB3	1.67	0.58
1:AU:163:PHE:HD2	1:CN:24:ALA:HA	1.68	0.58
1:AZ:41:LEU:HD11	1:AZ:52:ARG:HB2	1.84	0.58
1:BC:41:LEU:HD11	1:BC:52:ARG:HB2	1.84	0.58
1:CI:7:ASN:HD21	1:CJ:46:ASP:H	1.49	0.58
1:CZ:96:LYS:HG2	1:CZ:107:TYR:HB2	1.84	0.58
1:AG:33:ASP:H	1:AG:36:SER:HB3	1.67	0.58
1:BA:7:ASN:O	1:CW:53:PRO:HG2	2.04	0.58
1:BK:33:ASP:H	1:BK:36:SER:HB3	1.67	0.58
1:BP:96:LYS:HG2	1:BP:107:TYR:HB2	1.84	0.58
1:AD:34:ASN:ND2	1:DE:30:VAL:HA	2.19	0.58
1:BW:41:LEU:HD11	1:BW:52:ARG:HB2	1.86	0.58
1:BX:41:LEU:HD11	1:BX:52:ARG:HB2	1.84	0.58
1:DD:41:LEU:HD11	1:DD:52:ARG:HB2	1.86	0.58
1:AJ:41:LEU:HD11	1:AJ:52:ARG:HB2	1.86	0.58
1:AN:41:LEU:HD11	1:AN:52:ARG:HB2	1.84	0.58
1:BC:30:VAL:HA	1:BN:34:ASN:ND2	2.18	0.58
1:BG:96:LYS:HG2	1:BG:107:TYR:HB2	1.84	0.58

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BZ:41:LEU:HD11	1:BZ:52:ARG:HB2	1.86	0.58
1:CD:41:LEU:HD11	1:CD:52:ARG:HB2	1.84	0.58
1:AE:63:GLN:HG3	1:BG:50:LEU:CD2	2.34	0.58
1:BE:52:ARG:HH22	1:DC:157:ASP:CG	2.12	0.58
1:BQ:41:LEU:HD11	1:BQ:52:ARG:HB2	1.86	0.58
1:CE:96:LYS:HG2	1:CE:107:TYR:HB2	1.84	0.58
1:AU:7:ASN:O	1:CN:53:PRO:HG2	2.04	0.57
1:AA:34:ASN:ND2	1:DH:30:VAL:HA	2.20	0.57
1:BY:96:LYS:HG2	1:BY:107:TYR:HB2	1.84	0.57
1:DC:96:LYS:HG2	1:DC:107:TYR:HB2	1.84	0.57
1:BM:34:ASN:HB3	1:DI:161:ARG:HD2	1.86	0.57
1:BZ:33:ASP:H	1:BZ:36:SER:HB3	1.67	0.57
1:AD:41:LEU:HD11	1:AD:52:ARG:HB2	1.86	0.57
1:AO:53:PRO:HG2	1:CE:7:ASN:O	2.04	0.57
1:AV:33:ASP:H	1:AV:36:SER:HB3	1.67	0.57
1:AX:18:VAL:HG22	1:CT:16:ASN:OD1	2.04	0.57
1:BX:30:VAL:HA	1:CL:34:ASN:ND2	2.19	0.57
1:DE:41:LEU:HD11	1:DE:52:ARG:HB2	1.84	0.57
1:AP:41:LEU:HD11	1:AP:52:ARG:HB2	1.86	0.57
1:BF:50:LEU:HD21	1:BQ:82:VAL:HG11	1.85	0.57
1:BH:33:ASP:H	1:BH:36:SER:HB3	1.67	0.57
1:CI:33:ASP:H	1:CI:36:SER:HB3	1.67	0.57
1:CU:41:LEU:HD11	1:CU:52:ARG:HB2	1.86	0.57
1:AF:163:PHE:CD2	1:BV:24:ALA:HA	2.40	0.57
1:AK:41:LEU:HD11	1:AK:52:ARG:HB2	1.84	0.57
1:CO:41:LEU:HD11	1:CO:52:ARG:HB2	1.86	0.57
1:CR:34:ASN:ND2	1:DB:30:VAL:HA	2.20	0.57
1:AG:41:LEU:HD11	1:AG:52:ARG:HB2	1.86	0.57
1:AI:161:ARG:HD2	1:BY:34:ASN:HB3	1.87	0.57
1:AM:41:LEU:HD11	1:AM:52:ARG:HB2	1.86	0.57
1:BK:41:LEU:HD11	1:BK:52:ARG:HB2	1.86	0.57
1:CF:41:LEU:HD11	1:CF:52:ARG:HB2	1.86	0.57
1:BE:41:LEU:HD11	1:BE:52:ARG:HB2	1.86	0.57
1:BG:156:ILE:O	1:DC:84:ARG:HD2	2.05	0.57
1:CR:41:LEU:HD11	1:CR:52:ARG:HB2	1.86	0.57
1:AS:52:ARG:HH22	1:CN:157:ASP:CG	2.13	0.57
1:AK:30:VAL:HA	1:BW:34:ASN:ND2	2.19	0.57
1:BB:41:LEU:HD11	1:BB:52:ARG:HB2	1.86	0.57
1:BM:24:ALA:HA	1:DI:163:PHE:HD2	1.69	0.57
1:DJ:41:LEU:HD11	1:DJ:52:ARG:HB2	1.86	0.57
1:AI:24:ALA:HA	1:BY:163:PHE:HD2	1.70	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BH:41:LEU:HD11	1:BH:52:ARG:HB2	1.86	0.56
1:CI:41:LEU:HD11	1:CI:52:ARG:HB2	1.86	0.56
1:CT:34:ASN:HD21	1:DB:35:LYS:NZ	2.03	0.56
1:AG:34:ASN:ND2	1:CS:30:VAL:HA	2.20	0.56
1:BG:24:ALA:HA	1:DC:163:PHE:HD2	1.70	0.56
1:AZ:86:PHE:CE1	1:AZ:117:LEU:HD23	2.41	0.56
1:BD:24:ALA:HA	1:CZ:163:PHE:HD2	1.71	0.56
1:BH:86:PHE:HE2	1:BH:117:LEU:HD23	1.71	0.56
1:CL:41:LEU:HD11	1:CL:52:ARG:HB2	1.86	0.56
1:CM:86:PHE:CE1	1:CM:117:LEU:HD23	2.41	0.56
1:CP:86:PHE:CE1	1:CP:117:LEU:HD23	2.41	0.56
1:AQ:86:PHE:CE1	1:AQ:117:LEU:HD23	2.41	0.56
1:BI:86:PHE:CE1	1:BI:117:LEU:HD23	2.41	0.56
1:BL:86:PHE:CE1	1:BL:117:LEU:HD23	2.41	0.56
1:BT:41:LEU:HD11	1:BT:52:ARG:HB2	1.86	0.56
1:BW:86:PHE:HE2	1:BW:117:LEU:HD23	1.71	0.56
1:CY:86:PHE:CE1	1:CY:117:LEU:HD23	2.41	0.56
1:DG:41:LEU:HD11	1:DG:52:ARG:HB2	1.86	0.56
1:AH:86:PHE:CE1	1:AH:117:LEU:HD23	2.41	0.56
1:AR:42:ASP:OD2	1:CK:9:SER:N	2.38	0.56
1:AS:41:LEU:HD11	1:AS:52:ARG:HB2	1.86	0.56
1:AV:41:LEU:HD11	1:AV:52:ARG:HB2	1.86	0.56
1:CD:86:PHE:CE1	1:CD:117:LEU:HD23	2.41	0.56
1:CX:41:LEU:HD11	1:CX:52:ARG:HB2	1.86	0.56
1:AF:157:ASP:CG	1:BT:52:ARG:HH22	2.13	0.56
1:AS:86:PHE:HE2	1:AS:117:LEU:HD23	1.71	0.56
1:AV:86:PHE:HE2	1:AV:117:LEU:HD23	1.71	0.56
1:CJ:86:PHE:CE1	1:CJ:117:LEU:HD23	2.41	0.56
1:CO:86:PHE:HE2	1:CO:117:LEU:HD23	1.71	0.56
1:DH:86:PHE:CE1	1:DH:117:LEU:HD23	2.41	0.56
1:BN:41:LEU:HD11	1:BN:52:ARG:HB2	1.86	0.56
1:AA:28:VAL:HG11	1:BP:159:LEU:HD13	1.86	0.56
1:AY:86:PHE:HE2	1:AY:117:LEU:HD23	1.71	0.56
1:BA:163:PHE:HD2	1:CW:24:ALA:HA	1.71	0.56
1:BB:86:PHE:HE2	1:BB:117:LEU:HD23	1.71	0.56
1:BQ:86:PHE:HE2	1:BQ:117:LEU:HD23	1.71	0.56
1:BX:86:PHE:CE1	1:BX:117:LEU:HD23	2.41	0.56
1:CG:86:PHE:CE1	1:CG:117:LEU:HD23	2.41	0.56
1:DA:41:LEU:HD11	1:DA:52:ARG:HB2	1.86	0.56
1:DE:86:PHE:CE1	1:DE:117:LEU:HD23	2.41	0.56
1:DK:86:PHE:CE1	1:DK:117:LEU:HD23	2.41	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AK:86:PHE:CE1	1:AK:117:LEU:HD23	2.41	0.56
1:AW:86:PHE:CE1	1:AW:117:LEU:HD23	2.41	0.56
1:AY:41:LEU:HD11	1:AY:52:ARG:HB2	1.86	0.56
1:DG:86:PHE:HE2	1:DG:117:LEU:HD23	1.71	0.56
1:AA:41:LEU:HD11	1:AA:52:ARG:HB2	1.86	0.56
1:AB:86:PHE:CE1	1:AB:117:LEU:HD23	2.41	0.56
1:BD:34:ASN:HB3	1:CZ:161:ARG:HD2	1.86	0.56
1:BK:52:ARG:HH22	1:DI:157:ASP:CG	2.13	0.56
1:CA:86:PHE:CE1	1:CA:117:LEU:HD23	2.41	0.56
1:CF:86:PHE:HE2	1:CF:117:LEU:HD23	1.71	0.56
1:CI:86:PHE:HE2	1:CI:117:LEU:HD23	1.71	0.56
1:AJ:34:ASN:ND2	1:CV:30:VAL:HA	2.22	0.55
1:AN:86:PHE:CE1	1:AN:117:LEU:HD23	2.41	0.55
1:BZ:86:PHE:HE2	1:BZ:117:LEU:HD23	1.71	0.55
1:CC:41:LEU:HD11	1:CC:52:ARG:HB2	1.86	0.55
1:CR:86:PHE:HE2	1:CR:117:LEU:HD23	1.71	0.55
1:DB:86:PHE:CE1	1:DB:117:LEU:HD23	2.41	0.55
1:AT:86:PHE:CE1	1:AT:117:LEU:HD23	2.41	0.55
1:CS:86:PHE:CE1	1:CS:117:LEU:HD23	2.41	0.55
1:CU:86:PHE:HE2	1:CU:117:LEU:HD23	1.71	0.55
1:CV:86:PHE:CE1	1:CV:117:LEU:HD23	2.41	0.55
1:AT:30:VAL:HA	1:AV:34:ASN:ND2	2.21	0.55
1:DD:86:PHE:HE2	1:DD:117:LEU:HD23	1.71	0.55
1:BC:86:PHE:CE1	1:BC:117:LEU:HD23	2.41	0.55
1:BK:86:PHE:HE2	1:BK:117:LEU:HD23	1.71	0.55
1:BN:86:PHE:HE2	1:BN:117:LEU:HD23	1.71	0.55
1:BU:86:PHE:CE1	1:BU:117:LEU:HD23	2.41	0.55
1:AD:86:PHE:HE2	1:AD:117:LEU:HD23	1.71	0.55
1:BD:159:LEU:HD13	1:CX:28:VAL:HG11	1.88	0.55
1:BO:86:PHE:CE1	1:BO:117:LEU:HD23	2.41	0.55
1:BT:111:GLN:OE1	1:BU:119:SER:HB2	2.07	0.55
1:AE:86:PHE:CE1	1:AE:117:LEU:HD23	2.41	0.55
1:AG:86:PHE:HE2	1:AG:117:LEU:HD23	1.71	0.55
1:AL:24:ALA:HA	1:CB:163:PHE:HD2	1.72	0.55
1:AS:111:GLN:OE1	1:AT:119:SER:HB2	2.07	0.55
1:AY:111:GLN:OE1	1:AZ:119:SER:HB2	2.07	0.55
1:BG:53:PRO:HG2	1:DC:7:ASN:O	2.07	0.55
1:BR:86:PHE:CE1	1:BR:117:LEU:HD23	2.41	0.55
1:CI:111:GLN:OE1	1:CJ:119:SER:HB2	2.07	0.55
1:CO:111:GLN:OE1	1:CP:119:SER:HB2	2.07	0.55
1:DJ:111:GLN:OE1	1:DK:119:SER:HB2	2.07	0.55

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BG:119:SER:HB2	1:DC:111:GLN:OE1	2.07	0.55
1:CL:86:PHE:HE2	1:CL:117:LEU:HD23	1.71	0.55
1:CX:86:PHE:HE2	1:CX:117:LEU:HD23	1.71	0.55
1:CX:111:GLN:OE1	1:CY:119:SER:HB2	2.07	0.55
1:AG:111:GLN:OE1	1:AH:119:SER:HB2	2.07	0.55
1:BE:86:PHE:HE2	1:BE:117:LEU:HD23	1.71	0.55
1:BF:86:PHE:CE1	1:BF:117:LEU:HD23	2.41	0.55
1:BH:8:CYS:SG	1:BI:90:ARG:NH1	2.80	0.55
1:BJ:98:ASP:HB3	1:DF:1:SER:C	2.32	0.55
1:BW:8:CYS:SG	1:BX:90:ARG:NH1	2.80	0.55
1:BW:111:GLN:OE1	1:BX:119:SER:HB2	2.07	0.55
1:DA:111:GLN:OE1	1:DB:119:SER:HB2	2.07	0.55
1:AA:64:TYR:HD2	1:AA:67:ARG:NH2	2.05	0.55
1:AA:111:GLN:OE1	1:AB:119:SER:HB2	2.07	0.55
1:AM:64:TYR:HD2	1:AM:67:ARG:NH2	2.05	0.55
1:BQ:64:TYR:HD2	1:BQ:67:ARG:NH2	2.05	0.55
1:CL:8:CYS:SG	1:CM:90:ARG:NH1	2.80	0.55
1:CL:64:TYR:HD2	1:CL:67:ARG:NH2	2.05	0.55
1:CL:111:GLN:OE1	1:CM:119:SER:HB2	2.07	0.55
1:AC:24:ALA:HA	1:BP:163:PHE:HD2	1.71	0.54
1:AJ:8:CYS:SG	1:AK:90:ARG:NH1	2.80	0.54
1:AM:86:PHE:HE2	1:AM:117:LEU:HD23	1.71	0.54
1:BB:52:ARG:HH22	1:CZ:157:ASP:CG	2.14	0.54
1:BB:111:GLN:OE1	1:BC:119:SER:HB2	2.07	0.54
1:BE:111:GLN:OE1	1:BF:119:SER:HB2	2.07	0.54
1:BN:64:TYR:HD2	1:BN:67:ARG:NH2	2.05	0.54
1:BT:86:PHE:HE2	1:BT:117:LEU:HD23	1.71	0.54
1:CC:111:GLN:OE1	1:CD:119:SER:HB2	2.07	0.54
1:CR:8:CYS:SG	1:CS:90:ARG:NH1	2.80	0.54
1:CX:64:TYR:HD2	1:CX:67:ARG:NH2	2.05	0.54
1:DD:64:TYR:HD2	1:DD:67:ARG:NH2	2.05	0.54
1:AS:28:VAL:HG11	1:CN:159:LEU:HD13	1.88	0.54
1:BB:64:TYR:HD2	1:BB:67:ARG:NH2	2.05	0.54
1:BJ:94:TRP:CE2	1:BJ:111:GLN:HG3	2.43	0.54
1:BM:94:TRP:CE2	1:BM:111:GLN:HG3	2.43	0.54
1:BV:94:TRP:CE2	1:BV:111:GLN:HG3	2.43	0.54
1:CC:64:TYR:HD2	1:CC:67:ARG:NH2	2.05	0.54
1:CC:86:PHE:HE2	1:CC:117:LEU:HD23	1.71	0.54
1:CT:94:TRP:CE2	1:CT:111:GLN:HG3	2.43	0.54
1:DF:94:TRP:CE2	1:DF:111:GLN:HG3	2.43	0.54
1:DJ:8:CYS:SG	1:DK:90:ARG:NH1	2.80	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AC:94:TRP:CE2	1:AC:111:GLN:HG3	2.43	0.54
1:AE:30:VAL:HA	1:BE:34:ASN:ND2	2.22	0.54
1:AI:94:TRP:CE2	1:AI:111:GLN:HG3	2.43	0.54
1:AJ:64:TYR:HD2	1:AJ:67:ARG:NH2	2.05	0.54
1:BG:94:TRP:CE2	1:BG:111:GLN:HG3	2.43	0.54
1:BO:30:VAL:HA	1:DJ:34:ASN:ND2	2.23	0.54
1:CC:8:CYS:SG	1:CD:90:ARG:NH1	2.80	0.54
1:CE:94:TRP:CE2	1:CE:111:GLN:HG3	2.43	0.54
1:CF:111:GLN:OE1	1:CG:119:SER:HB2	2.07	0.54
1:CK:94:TRP:CE2	1:CK:111:GLN:HG3	2.43	0.54
1:CN:94:TRP:CE2	1:CN:111:GLN:HG3	2.43	0.54
1:DA:64:TYR:HD2	1:DA:67:ARG:NH2	2.05	0.54
1:DA:86:PHE:HE2	1:DA:117:LEU:HD23	1.71	0.54
1:DC:94:TRP:CE2	1:DC:111:GLN:HG3	2.43	0.54
1:AA:86:PHE:HE2	1:AA:117:LEU:HD23	1.71	0.54
1:AD:8:CYS:SG	1:AE:90:ARG:NH1	2.80	0.54
1:AD:111:GLN:OE1	1:AE:119:SER:HB2	2.07	0.54
1:AJ:52:ARG:HH22	1:CB:157:ASP:CG	2.15	0.54
1:AJ:86:PHE:HE2	1:AJ:117:LEU:HD23	1.71	0.54
1:AL:94:TRP:CE2	1:AL:111:GLN:HG3	2.43	0.54
1:AM:111:GLN:OE1	1:AN:119:SER:HB2	2.07	0.54
1:AP:8:CYS:SG	1:AQ:90:ARG:NH1	2.80	0.54
1:AP:86:PHE:HE2	1:AP:117:LEU:HD23	1.71	0.54
1:BH:64:TYR:HD2	1:BH:67:ARG:NH2	2.05	0.54
1:BH:111:GLN:OE1	1:BI:119:SER:HB2	2.07	0.54
1:BJ:163:PHE:HD2	1:DF:24:ALA:HA	1.72	0.54
1:CB:94:TRP:CE2	1:CB:111:GLN:HG3	2.43	0.54
1:CQ:94:TRP:CE2	1:CQ:111:GLN:HG3	2.43	0.54
1:CW:94:TRP:CE2	1:CW:111:GLN:HG3	2.43	0.54
1:DD:111:GLN:OE1	1:DE:119:SER:HB2	2.07	0.54
1:DI:94:TRP:CE2	1:DI:111:GLN:HG3	2.43	0.54
1:DJ:86:PHE:HE2	1:DJ:117:LEU:HD23	1.71	0.54
1:BP:94:TRP:CE2	1:BP:111:GLN:HG3	2.43	0.54
1:BQ:8:CYS:SG	1:BR:90:ARG:NH1	2.80	0.54
1:CI:64:TYR:HD2	1:CI:67:ARG:NH2	2.05	0.54
1:CR:64:TYR:HD2	1:CR:67:ARG:NH2	2.05	0.54
1:CU:64:TYR:HD2	1:CU:67:ARG:NH2	2.05	0.54
1:CU:111:GLN:OE1	1:CV:119:SER:HB2	2.07	0.54
1:DG:111:GLN:OE1	1:DH:119:SER:HB2	2.07	0.54
1:AB:37:THR:HG23	1:AB:58:ASN:HB3	1.90	0.54
1:AH:37:THR:HG23	1:AH:58:ASN:HB3	1.90	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AV:8:CYS:SG	1:AW:90:ARG:NH1	2.80	0.54
1:AW:37:THR:HG23	1:AW:58:ASN:HB3	1.90	0.54
1:AX:94:TRP:CE2	1:AX:111:GLN:HG3	2.43	0.54
1:AY:8:CYS:SG	1:AZ:90:ARG:NH1	2.80	0.54
1:BA:94:TRP:CE2	1:BA:111:GLN:HG3	2.43	0.54
1:BB:8:CYS:SG	1:BC:90:ARG:NH1	2.80	0.54
1:BK:8:CYS:SG	1:BL:90:ARG:NH1	2.80	0.54
1:BN:8:CYS:SG	1:BO:90:ARG:NH1	2.80	0.54
1:BN:111:GLN:OE1	1:BO:119:SER:HB2	2.07	0.54
1:BT:64:TYR:HD2	1:BT:67:ARG:NH2	2.05	0.54
1:CF:8:CYS:SG	1:CG:90:ARG:NH1	2.80	0.54
1:DB:37:THR:HG23	1:DB:58:ASN:HB3	1.90	0.54
1:DG:8:CYS:SG	1:DH:90:ARG:NH1	2.80	0.54
1:AD:64:TYR:HD2	1:AD:67:ARG:NH2	2.05	0.54
1:AF:94:TRP:CE2	1:AF:111:GLN:HG3	2.43	0.54
1:AJ:111:GLN:OE1	1:AK:119:SER:HB2	2.07	0.54
1:AS:64:TYR:HD2	1:AS:67:ARG:NH2	2.05	0.54
1:CI:8:CYS:SG	1:CJ:90:ARG:NH1	2.80	0.54
1:DD:8:CYS:SG	1:DE:90:ARG:NH1	2.80	0.54
1:AC:157:ASP:CG	1:BN:52:ARG:HH22	2.14	0.54
1:AM:8:CYS:SG	1:AN:90:ARG:NH1	2.80	0.54
1:AT:37:THR:HG23	1:AT:58:ASN:HB3	1.90	0.54
1:AU:94:TRP:CE2	1:AU:111:GLN:HG3	2.43	0.54
1:BE:8:CYS:SG	1:BF:90:ARG:NH1	2.80	0.54
1:BE:64:TYR:HD2	1:BE:67:ARG:NH2	2.05	0.54
1:BY:94:TRP:CE2	1:BY:111:GLN:HG3	2.43	0.54
1:BZ:111:GLN:OE1	1:CA:119:SER:HB2	2.07	0.54
1:CP:37:THR:HG23	1:CP:58:ASN:HB3	1.90	0.54
1:CY:37:THR:HG23	1:CY:58:ASN:HB3	1.90	0.54
1:DA:8:CYS:SG	1:DB:90:ARG:NH1	2.80	0.54
1:DA:19:LEU:HD11	1:DB:17:LEU:HD22	1.90	0.54
1:DL:94:TRP:CE2	1:DL:111:GLN:HG3	2.43	0.54
1:AP:111:GLN:OE1	1:AQ:119:SER:HB2	2.07	0.54
1:AR:94:TRP:CE2	1:AR:111:GLN:HG3	2.43	0.54
1:BE:19:LEU:HD11	1:BF:17:LEU:HD22	1.90	0.54
1:BK:64:TYR:HD2	1:BK:67:ARG:NH2	2.05	0.54
1:BK:111:GLN:OE1	1:BL:119:SER:HB2	2.07	0.54
1:BN:19:LEU:HD11	1:BO:17:LEU:HD22	1.90	0.54
1:CO:8:CYS:SG	1:CP:90:ARG:NH1	2.80	0.54
1:CO:19:LEU:HD11	1:CP:17:LEU:HD22	1.90	0.54
1:CU:8:CYS:SG	1:CV:90:ARG:NH1	2.80	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AN:30:VAL:HA	1:BZ:34:ASN:ND2	2.23	0.54
1:CA:37:THR:HG23	1:CA:58:ASN:HB3	1.90	0.54
1:CD:30:VAL:HA	1:CI:34:ASN:ND2	2.23	0.54
1:CR:111:GLN:OE1	1:CS:119:SER:HB2	2.07	0.54
1:AG:8:CYS:SG	1:AH:90:ARG:NH1	2.80	0.53
1:AG:19:LEU:HD11	1:AH:17:LEU:HD22	1.90	0.53
1:AG:64:TYR:HD2	1:AG:67:ARG:NH2	2.06	0.53
1:AN:37:THR:HG23	1:AN:58:ASN:HB3	1.90	0.53
1:AO:94:TRP:CE2	1:AO:111:GLN:HG3	2.43	0.53
1:AR:163:PHE:CD2	1:CK:24:ALA:HA	2.42	0.53
1:AV:111:GLN:OE1	1:AW:119:SER:HB2	2.07	0.53
1:BW:64:TYR:HD2	1:BW:67:ARG:NH2	2.05	0.53
1:CC:19:LEU:HD11	1:CD:17:LEU:HD22	1.90	0.53
1:CX:8:CYS:SG	1:CY:90:ARG:NH1	2.80	0.53
1:CZ:94:TRP:CE2	1:CZ:111:GLN:HG3	2.43	0.53
1:DJ:64:TYR:HD2	1:DJ:67:ARG:NH2	2.05	0.53
1:AG:52:ARG:HH22	1:BY:157:ASP:CG	2.16	0.53
1:AU:157:ASP:CG	1:CL:52:ARG:HH22	2.17	0.53
1:AY:19:LEU:HD11	1:AZ:17:LEU:HD22	1.90	0.53
1:BZ:19:LEU:HD11	1:CA:17:LEU:HD22	1.90	0.53
1:DG:64:TYR:HD2	1:DG:67:ARG:NH2	2.05	0.53
1:DK:37:THR:HG23	1:DK:58:ASN:HB3	1.90	0.53
1:AA:8:CYS:SG	1:AB:90:ARG:NH1	2.80	0.53
1:AP:64:TYR:HD2	1:AP:67:ARG:NH2	2.05	0.53
1:AS:19:LEU:HD11	1:AT:17:LEU:HD22	1.90	0.53
1:BD:94:TRP:CE2	1:BD:111:GLN:HG3	2.43	0.53
1:BS:94:TRP:CE2	1:BS:111:GLN:HG3	2.43	0.53
1:CI:19:LEU:HD11	1:CJ:17:LEU:HD22	1.90	0.53
1:DE:37:THR:HG23	1:DE:58:ASN:HB3	1.90	0.53
1:AC:163:PHE:HD2	1:BP:24:ALA:HA	1.74	0.53
1:BJ:6:ILE:HG22	1:DF:94:TRP:NE1	2.23	0.53
1:BK:19:LEU:HD11	1:BL:17:LEU:HD22	1.90	0.53
1:BQ:19:LEU:HD11	1:BR:17:LEU:HD22	1.90	0.53
1:BQ:111:GLN:OE1	1:BR:119:SER:HB2	2.07	0.53
1:CF:19:LEU:HD11	1:CG:17:LEU:HD22	1.90	0.53
1:CX:19:LEU:HD11	1:CY:17:LEU:HD22	1.90	0.53
1:AK:37:THR:HG23	1:AK:58:ASN:HB3	1.90	0.53
1:AS:8:CYS:SG	1:AT:90:ARG:NH1	2.80	0.53
1:BF:37:THR:HG23	1:BF:58:ASN:HB3	1.90	0.53
1:BR:37:THR:HG23	1:BR:58:ASN:HB3	1.90	0.53
1:BT:8:CYS:SG	1:BU:90:ARG:NH1	2.80	0.53

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BZ:8:CYS:SG	1:CA:90:ARG:NH1	2.80	0.53
1:CF:64:TYR:HD2	1:CF:67:ARG:NH2	2.05	0.53
1:CH:94:TRP:CE2	1:CH:111:GLN:HG3	2.43	0.53
1:DK:96:LYS:HG2	1:DK:107:TYR:HB2	1.91	0.53
1:AE:96:LYS:HG2	1:AE:107:TYR:HB2	1.91	0.53
1:AV:64:TYR:HD2	1:AV:67:ARG:NH2	2.05	0.53
1:AZ:96:LYS:HG2	1:AZ:107:TYR:HB2	1.91	0.53
1:BC:37:THR:HG23	1:BC:58:ASN:HB3	1.90	0.53
1:CG:96:LYS:HG2	1:CG:107:TYR:HB2	1.91	0.53
1:CO:64:TYR:HD2	1:CO:67:ARG:NH2	2.05	0.53
1:CU:19:LEU:HD11	1:CV:17:LEU:HD22	1.90	0.53
1:DD:19:LEU:HD11	1:DE:17:LEU:HD22	1.90	0.53
1:AU:24:ALA:HA	1:CN:163:PHE:HD2	1.72	0.53
1:BD:163:PHE:HD2	1:CZ:24:ALA:HA	1.74	0.53
1:BZ:64:TYR:HD2	1:BZ:67:ARG:NH2	2.05	0.53
1:CD:96:LYS:HG2	1:CD:107:TYR:HB2	1.91	0.53
1:AW:30:VAL:HA	1:BT:34:ASN:ND2	2.24	0.53
1:BB:19:LEU:HD11	1:BC:17:LEU:HD22	1.90	0.53
1:CG:37:THR:HG23	1:CG:58:ASN:HB3	1.90	0.53
1:AH:96:LYS:HG2	1:AH:107:TYR:HB2	1.91	0.53
1:AQ:37:THR:HG23	1:AQ:58:ASN:HB3	1.90	0.53
1:BA:46:ASP:H	1:CW:7:ASN:HD21	1.56	0.53
1:BO:37:THR:HG23	1:BO:58:ASN:HB3	1.90	0.53
1:CM:96:LYS:HG2	1:CM:107:TYR:HB2	1.91	0.53
1:CV:96:LYS:HG2	1:CV:107:TYR:HB2	1.91	0.53
1:DG:19:LEU:HD11	1:DH:17:LEU:HD22	1.90	0.53
1:BH:82:VAL:HG13	1:BL:50:LEU:HD11	1.91	0.53
1:BI:37:THR:HG23	1:BI:58:ASN:HB3	1.90	0.53
1:BL:96:LYS:HG2	1:BL:107:TYR:HB2	1.91	0.53
1:CD:37:THR:HG23	1:CD:58:ASN:HB3	1.90	0.53
1:CS:37:THR:HG23	1:CS:58:ASN:HB3	1.90	0.53
1:AQ:96:LYS:HG2	1:AQ:107:TYR:HB2	1.91	0.52
1:BI:96:LYS:HG2	1:BI:107:TYR:HB2	1.91	0.52
1:BW:19:LEU:HD11	1:BX:17:LEU:HD22	1.90	0.52
1:CJ:96:LYS:HG2	1:CJ:107:TYR:HB2	1.91	0.52
1:CR:19:LEU:HD11	1:CS:17:LEU:HD22	1.90	0.52
1:AA:19:LEU:HD11	1:AB:17:LEU:HD22	1.90	0.52
1:AE:37:THR:HG23	1:AE:58:ASN:HB3	1.90	0.52
1:AI:34:ASN:HB3	1:BY:161:ARG:HD2	1.90	0.52
1:AX:24:ALA:HA	1:CT:163:PHE:CD2	2.43	0.52
1:AZ:37:THR:HG23	1:AZ:58:ASN:HB3	1.90	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BH:19:LEU:HD11	1:BI:17:LEU:HD22	1.90	0.52
1:BX:96:LYS:HG2	1:BX:107:TYR:HB2	1.91	0.52
1:BG:40:TYR:CE1	1:DC:163:PHE:HB2	2.44	0.52
1:BR:96:LYS:HG2	1:BR:107:TYR:HB2	1.91	0.52
1:CM:37:THR:HG23	1:CM:58:ASN:HB3	1.90	0.52
1:DE:96:LYS:HG2	1:DE:107:TYR:HB2	1.91	0.52
1:AJ:19:LEU:HD11	1:AK:17:LEU:HD22	1.90	0.52
1:AO:159:LEU:HD13	1:CC:28:VAL:HG11	1.91	0.52
1:AR:86:PHE:CD2	1:CK:155:LEU:HD13	2.44	0.52
1:AY:64:TYR:HD2	1:AY:67:ARG:NH2	2.05	0.52
1:BX:37:THR:HG23	1:BX:58:ASN:HB3	1.90	0.52
1:CJ:37:THR:HG23	1:CJ:58:ASN:HB3	1.90	0.52
1:AD:19:LEU:HD11	1:AE:17:LEU:HD22	1.90	0.52
1:AF:7:ASN:HD21	1:BV:46:ASP:H	1.58	0.52
1:AV:19:LEU:HD11	1:AW:17:LEU:HD22	1.91	0.52
1:AW:96:LYS:HG2	1:AW:107:TYR:HB2	1.91	0.52
1:BT:19:LEU:HD11	1:BU:17:LEU:HD22	1.90	0.52
1:CY:96:LYS:HG2	1:CY:107:TYR:HB2	1.91	0.52
1:DD:17:LEU:HD22	1:DE:19:LEU:HD11	1.92	0.52
1:AG:17:LEU:HD22	1:AH:19:LEU:HD11	1.92	0.52
1:AY:17:LEU:HD22	1:AZ:19:LEU:HD11	1.92	0.52
1:BE:17:LEU:HD22	1:BF:19:LEU:HD11	1.92	0.52
1:BL:37:THR:HG23	1:BL:58:ASN:HB3	1.90	0.52
1:CR:17:LEU:HD22	1:CS:19:LEU:HD11	1.92	0.52
1:CV:37:THR:HG23	1:CV:58:ASN:HB3	1.90	0.52
1:DH:96:LYS:HG2	1:DH:107:TYR:HB2	1.91	0.52
1:AM:19:LEU:HD11	1:AN:17:LEU:HD22	1.90	0.52
1:AP:19:LEU:HD11	1:AQ:17:LEU:HD22	1.90	0.52
1:AX:94:TRP:NE1	1:CT:6:ILE:HG22	2.25	0.52
1:BU:96:LYS:HG2	1:BU:107:TYR:HB2	1.91	0.52
1:DH:37:THR:HG23	1:DH:58:ASN:HB3	1.90	0.52
1:AF:84:ARG:HD2	1:BV:156:ILE:O	2.10	0.52
1:BU:37:THR:HG23	1:BU:58:ASN:HB3	1.90	0.52
1:AK:96:LYS:HG2	1:AK:107:TYR:HB2	1.91	0.52
1:AV:17:LEU:HD22	1:AW:19:LEU:HD11	1.92	0.52
1:BI:130:ILE:O	1:BI:133:ILE:HG22	2.10	0.52
1:BK:17:LEU:HD22	1:BL:19:LEU:HD11	1.92	0.52
1:BK:94:TRP:CE2	1:BK:111:GLN:HG3	2.45	0.52
1:CF:94:TRP:CE2	1:CF:111:GLN:HG3	2.45	0.52
1:DD:94:TRP:CE2	1:DD:111:GLN:HG3	2.45	0.52
1:AA:17:LEU:HD22	1:AB:19:LEU:HD11	1.92	0.52

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AB:96:LYS:HG2	1:AB:107:TYR:HB2	1.91	0.52
1:AD:94:TRP:CE2	1:AD:111:GLN:HG3	2.45	0.52
1:BF:96:LYS:HG2	1:BF:107:TYR:HB2	1.91	0.52
1:BH:52:ARG:HH22	1:DF:157:ASP:CG	2.17	0.52
1:CX:94:TRP:CE2	1:CX:111:GLN:HG3	2.45	0.52
1:CY:130:ILE:O	1:CY:133:ILE:HG22	2.10	0.52
1:DB:96:LYS:HG2	1:DB:107:TYR:HB2	1.91	0.52
1:DJ:94:TRP:CE2	1:DJ:111:GLN:HG3	2.45	0.52
1:AA:52:ARG:HH22	1:BP:157:ASP:CG	2.18	0.51
1:AD:17:LEU:HD22	1:AE:19:LEU:HD11	1.92	0.51
1:AP:17:LEU:HD22	1:AQ:19:LEU:HD11	1.92	0.51
1:AQ:130:ILE:O	1:AQ:133:ILE:HG22	2.10	0.51
1:AW:130:ILE:O	1:AW:133:ILE:HG22	2.10	0.51
1:BC:130:ILE:O	1:BC:133:ILE:HG22	2.10	0.51
1:BM:7:ASN:HD21	1:DI:46:ASP:H	1.58	0.51
1:BU:130:ILE:O	1:BU:133:ILE:HG22	2.10	0.51
1:CC:94:TRP:CE2	1:CC:111:GLN:HG3	2.45	0.51
1:CG:130:ILE:O	1:CG:133:ILE:HG22	2.10	0.51
1:CO:17:LEU:HD22	1:CP:19:LEU:HD11	1.92	0.51
1:CP:96:LYS:HG2	1:CP:107:TYR:HB2	1.91	0.51
1:CS:130:ILE:O	1:CS:133:ILE:HG22	2.10	0.51
1:CU:94:TRP:CE2	1:CU:111:GLN:HG3	2.45	0.51
1:CX:17:LEU:HD22	1:CY:19:LEU:HD11	1.92	0.51
1:AN:130:ILE:O	1:AN:133:ILE:HG22	2.10	0.51
1:BT:17:LEU:HD22	1:BU:19:LEU:HD11	1.92	0.51
1:CA:96:LYS:HG2	1:CA:107:TYR:HB2	1.91	0.51
1:CS:96:LYS:HG2	1:CS:107:TYR:HB2	1.91	0.51
1:CU:17:LEU:HD22	1:CV:19:LEU:HD11	1.92	0.51
1:DJ:19:LEU:HD11	1:DK:17:LEU:HD22	1.90	0.51
1:AB:130:ILE:O	1:AB:133:ILE:HG22	2.10	0.51
1:AE:130:ILE:O	1:AE:133:ILE:HG22	2.10	0.51
1:AM:34:ASN:ND2	1:CP:30:VAL:HA	2.24	0.51
1:BC:96:LYS:HG2	1:BC:107:TYR:HB2	1.91	0.51
1:BQ:94:TRP:CE2	1:BQ:111:GLN:HG3	2.45	0.51
1:BR:130:ILE:O	1:BR:133:ILE:HG22	2.10	0.51
1:BX:130:ILE:O	1:BX:133:ILE:HG22	2.10	0.51
1:CD:130:ILE:O	1:CD:133:ILE:HG22	2.10	0.51
1:DJ:17:LEU:HD22	1:DK:19:LEU:HD11	1.92	0.51
1:AC:161:ARG:HD2	1:BP:34:ASN:HB3	1.91	0.51
1:AM:94:TRP:CE2	1:AM:111:GLN:HG3	2.45	0.51
1:BJ:153:GLN:OE1	1:DF:127:ALA:HB2	2.11	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BL:130:ILE:O	1:BL:133:ILE:HG22	2.10	0.51
1:BO:130:ILE:O	1:BO:133:ILE:HG22	2.10	0.51
1:CA:130:ILE:O	1:CA:133:ILE:HG22	2.10	0.51
1:CL:17:LEU:HD22	1:CM:19:LEU:HD11	1.92	0.51
1:CO:94:TRP:CE2	1:CO:111:GLN:HG3	2.45	0.51
1:DH:130:ILE:O	1:DH:133:ILE:HG22	2.10	0.51
1:AA:94:TRP:CE2	1:AA:111:GLN:HG3	2.45	0.51
1:AG:94:TRP:CE2	1:AG:111:GLN:HG3	2.45	0.51
1:AL:159:LEU:HD13	1:BZ:28:VAL:HG11	1.91	0.51
1:AS:94:TRP:CE2	1:AS:111:GLN:HG3	2.45	0.51
1:AV:94:TRP:CE2	1:AV:111:GLN:HG3	2.45	0.51
1:AY:94:TRP:CE2	1:AY:111:GLN:HG3	2.45	0.51
1:BB:17:LEU:HD22	1:BC:19:LEU:HD11	1.92	0.51
1:CN:95:ARG:NH2	1:CN:108:THR:OG1	2.44	0.51
1:CR:94:TRP:CE2	1:CR:111:GLN:HG3	2.45	0.51
1:DK:130:ILE:O	1:DK:133:ILE:HG22	2.10	0.51
1:AC:34:ASN:HB3	1:BP:161:ARG:HD2	1.91	0.51
1:AC:159:LEU:HD13	1:BN:28:VAL:HG11	1.93	0.51
1:AH:130:ILE:O	1:AH:133:ILE:HG22	2.10	0.51
1:AJ:17:LEU:HD22	1:AK:19:LEU:HD11	1.92	0.51
1:AM:17:LEU:HD22	1:AN:19:LEU:HD11	1.92	0.51
1:AP:94:TRP:CE2	1:AP:111:GLN:HG3	2.45	0.51
1:AT:121:VAL:HG13	1:AT:125:THR:HB	1.93	0.51
1:AZ:130:ILE:O	1:AZ:133:ILE:HG22	2.10	0.51
1:BE:94:TRP:CE2	1:BE:111:GLN:HG3	2.45	0.51
1:BQ:17:LEU:HD22	1:BR:19:LEU:HD11	1.92	0.51
1:CF:17:LEU:HD22	1:CG:19:LEU:HD11	1.92	0.51
1:CJ:130:ILE:O	1:CJ:133:ILE:HG22	2.10	0.51
1:CL:19:LEU:HD11	1:CM:17:LEU:HD22	1.90	0.51
1:CM:130:ILE:O	1:CM:133:ILE:HG22	2.10	0.51
1:CV:121:VAL:HG13	1:CV:125:THR:HB	1.93	0.51
1:DC:95:ARG:NH2	1:DC:108:THR:OG1	2.44	0.51
1:DE:130:ILE:O	1:DE:133:ILE:HG22	2.10	0.51
1:AJ:94:TRP:CE2	1:AJ:111:GLN:HG3	2.45	0.51
1:AQ:121:VAL:HG13	1:AQ:125:THR:HB	1.93	0.51
1:AT:96:LYS:HG2	1:AT:107:TYR:HB2	1.91	0.51
1:BD:95:ARG:NH2	1:BD:108:THR:OG1	2.44	0.51
1:BM:95:ARG:NH2	1:BM:108:THR:OG1	2.44	0.51
1:BO:96:LYS:HG2	1:BO:107:TYR:HB2	1.91	0.51
1:BV:95:ARG:NH2	1:BV:108:THR:OG1	2.44	0.51
1:CI:17:LEU:HD22	1:CJ:19:LEU:HD11	1.92	0.51

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:DI:95:ARG:NH2	1:DI:108:THR:OG1	2.44	0.51
1:AK:130:ILE:O	1:AK:133:ILE:HG22	2.10	0.51
1:BF:130:ILE:O	1:BF:133:ILE:HG22	2.10	0.51
1:BH:94:TRP:CE2	1:BH:111:GLN:HG3	2.45	0.51
1:BM:159:LEU:HD13	1:DG:28:VAL:HG11	1.93	0.51
1:BU:121:VAL:HG13	1:BU:125:THR:HB	1.93	0.51
1:BW:6:ILE:O	1:BW:6:ILE:HG13	2.11	0.51
1:CA:121:VAL:HG13	1:CA:125:THR:HB	1.93	0.51
1:CJ:121:VAL:HG13	1:CJ:125:THR:HB	1.93	0.51
1:CQ:95:ARG:NH2	1:CQ:108:THR:OG1	2.44	0.51
1:DE:121:VAL:HG13	1:DE:125:THR:HB	1.93	0.51
1:DG:17:LEU:HD22	1:DH:19:LEU:HD11	1.92	0.51
1:DG:94:TRP:CE2	1:DG:111:GLN:HG3	2.45	0.51
1:AG:6:ILE:O	1:AG:6:ILE:HG13	2.11	0.51
1:AN:96:LYS:HG2	1:AN:107:TYR:HB2	1.91	0.51
1:AR:95:ARG:NH2	1:AR:108:THR:OG1	2.44	0.51
1:BK:6:ILE:HG13	1:BK:6:ILE:O	2.11	0.51
1:BN:17:LEU:HD22	1:BO:19:LEU:HD11	1.92	0.51
1:BN:94:TRP:CE2	1:BN:111:GLN:HG3	2.45	0.51
1:CZ:95:ARG:NH2	1:CZ:108:THR:OG1	2.44	0.51
1:DA:94:TRP:CE2	1:DA:111:GLN:HG3	2.45	0.51
1:DL:95:ARG:NH2	1:DL:108:THR:OG1	2.44	0.51
1:AE:121:VAL:HG13	1:AE:125:THR:HB	1.93	0.51
1:AS:17:LEU:HD22	1:AT:19:LEU:HD11	1.92	0.51
1:BA:95:ARG:NH2	1:BA:108:THR:OG1	2.44	0.51
1:BL:121:VAL:HG13	1:BL:125:THR:HB	1.93	0.51
1:BP:95:ARG:NH2	1:BP:108:THR:OG1	2.44	0.51
1:BQ:6:ILE:O	1:BQ:6:ILE:HG13	2.11	0.51
1:BW:17:LEU:HD22	1:BX:19:LEU:HD11	1.92	0.51
1:BY:95:ARG:NH2	1:BY:108:THR:OG1	2.44	0.51
1:BZ:94:TRP:CE2	1:BZ:111:GLN:HG3	2.45	0.51
1:CT:95:ARG:NH2	1:CT:108:THR:OG1	2.44	0.51
1:AF:95:ARG:NH2	1:AF:108:THR:OG1	2.44	0.50
1:AM:6:ILE:O	1:AM:6:ILE:HG13	2.11	0.50
1:AO:24:ALA:HA	1:CE:163:PHE:CD2	2.44	0.50
1:AU:95:ARG:NH2	1:AU:108:THR:OG1	2.44	0.50
1:BB:6:ILE:HG13	1:BB:6:ILE:O	2.11	0.50
1:CB:95:ARG:NH2	1:CB:108:THR:OG1	2.44	0.50
1:CC:17:LEU:HD22	1:CD:19:LEU:HD11	1.92	0.50
1:CI:94:TRP:CE2	1:CI:111:GLN:HG3	2.45	0.50
1:CR:6:ILE:HG13	1:CR:6:ILE:O	2.11	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AS:6:ILE:O	1:AS:6:ILE:HG13	2.11	0.50
1:AY:6:ILE:HG13	1:AY:6:ILE:O	2.11	0.50
1:BG:2:TYR:CE1	1:DC:97:ALA:HA	2.47	0.50
1:BG:95:ARG:NH2	1:BG:108:THR:OG1	2.44	0.50
1:BI:121:VAL:HG13	1:BI:125:THR:HB	1.93	0.50
1:BJ:34:ASN:HB3	1:DF:161:ARG:HD2	1.91	0.50
1:BN:6:ILE:O	1:BN:6:ILE:HG13	2.11	0.50
1:BW:94:TRP:CE2	1:BW:111:GLN:HG3	2.45	0.50
1:CK:95:ARG:NH2	1:CK:108:THR:OG1	2.44	0.50
1:CY:121:VAL:HG13	1:CY:125:THR:HB	1.93	0.50
1:DF:95:ARG:NH2	1:DF:108:THR:OG1	2.44	0.50
1:AA:6:ILE:O	1:AA:6:ILE:HG13	2.11	0.50
1:AC:95:ARG:NH2	1:AC:108:THR:OG1	2.44	0.50
1:AI:95:ARG:NH2	1:AI:108:THR:OG1	2.44	0.50
1:AR:40:TYR:CZ	1:CK:163:PHE:HB2	2.45	0.50
1:BH:28:VAL:HG11	1:DF:159:LEU:HD13	1.93	0.50
1:CG:121:VAL:HG13	1:CG:125:THR:HB	1.93	0.50
1:DA:17:LEU:HD22	1:DB:19:LEU:HD11	1.92	0.50
1:DB:130:ILE:O	1:DB:133:ILE:HG22	2.10	0.50
1:AL:95:ARG:NH2	1:AL:108:THR:OG1	2.44	0.50
1:BH:17:LEU:HD22	1:BI:19:LEU:HD11	1.92	0.50
1:BJ:95:ARG:NH2	1:BJ:108:THR:OG1	2.44	0.50
1:BZ:17:LEU:HD22	1:CA:19:LEU:HD11	1.92	0.50
1:AC:57:PHE:CE1	1:BP:162:ILE:HD12	2.47	0.50
1:AL:163:PHE:HD2	1:CB:24:ALA:HA	1.77	0.50
1:AN:121:VAL:HG13	1:AN:125:THR:HB	1.93	0.50
1:CF:6:ILE:HG13	1:CF:6:ILE:O	2.11	0.50
1:CH:95:ARG:NH2	1:CH:108:THR:OG1	2.44	0.50
1:CV:130:ILE:O	1:CV:133:ILE:HG22	2.10	0.50
1:DA:6:ILE:O	1:DA:6:ILE:HG13	2.11	0.50
1:DH:121:VAL:HG13	1:DH:125:THR:HB	1.93	0.50
1:AD:161:ARG:NH1	1:BG:25:GLU:O	2.44	0.50
1:AR:84:ARG:HD2	1:CK:156:ILE:O	2.11	0.50
1:AT:130:ILE:O	1:AT:133:ILE:HG22	2.10	0.50
1:BH:6:ILE:HG13	1:BH:6:ILE:O	2.11	0.50
1:BS:95:ARG:NH2	1:BS:108:THR:OG1	2.44	0.50
1:BT:94:TRP:CE2	1:BT:111:GLN:HG3	2.45	0.50
1:CP:130:ILE:O	1:CP:133:ILE:HG22	2.10	0.50
1:CW:95:ARG:NH2	1:CW:108:THR:OG1	2.44	0.50
1:AD:6:ILE:O	1:AD:6:ILE:HG13	2.12	0.50
1:BC:121:VAL:HG13	1:BC:125:THR:HB	1.93	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CL:6:ILE:HG13	1:CL:6:ILE:O	2.11	0.50
1:CS:121:VAL:HG13	1:CS:125:THR:HB	1.93	0.50
1:CX:6:ILE:HG13	1:CX:6:ILE:O	2.11	0.50
1:DJ:6:ILE:O	1:DJ:6:ILE:HG13	2.11	0.50
1:AO:95:ARG:NH2	1:AO:108:THR:OG1	2.44	0.50
1:AP:6:ILE:O	1:AP:6:ILE:HG13	2.11	0.50
1:AV:6:ILE:HG13	1:AV:6:ILE:O	2.12	0.50
1:AW:121:VAL:HG13	1:AW:125:THR:HB	1.93	0.50
1:AZ:121:VAL:HG13	1:AZ:125:THR:HB	1.93	0.50
1:BS:86:PHE:CE1	1:BS:117:LEU:HD23	2.47	0.50
1:CI:6:ILE:HG13	1:CI:6:ILE:O	2.11	0.50
1:CL:94:TRP:CE2	1:CL:111:GLN:HG3	2.45	0.50
1:CT:86:PHE:CE1	1:CT:117:LEU:HD23	2.47	0.50
1:DD:6:ILE:O	1:DD:6:ILE:HG13	2.11	0.50
1:AJ:6:ILE:O	1:AJ:6:ILE:HG13	2.11	0.50
1:BB:94:TRP:CE2	1:BB:111:GLN:HG3	2.45	0.50
1:BG:57:PHE:CE1	1:DC:162:ILE:HD12	2.47	0.50
1:CU:6:ILE:HG13	1:CU:6:ILE:O	2.11	0.50
1:DL:86:PHE:CE1	1:DL:117:LEU:HD23	2.47	0.50
1:AB:121:VAL:HG13	1:AB:125:THR:HB	1.93	0.49
1:AI:86:PHE:CE1	1:AI:117:LEU:HD23	2.47	0.49
1:BD:86:PHE:CE1	1:BD:117:LEU:HD23	2.47	0.49
1:BX:121:VAL:HG13	1:BX:125:THR:HB	1.93	0.49
1:CD:121:VAL:HG13	1:CD:125:THR:HB	1.93	0.49
1:CO:6:ILE:HG13	1:CO:6:ILE:O	2.11	0.49
1:CP:121:VAL:HG13	1:CP:125:THR:HB	1.93	0.49
1:CQ:86:PHE:CE1	1:CQ:117:LEU:HD23	2.47	0.49
1:AF:94:TRP:NE1	1:BV:6:ILE:HG22	2.27	0.49
1:AO:86:PHE:CE1	1:AO:117:LEU:HD23	2.47	0.49
1:AR:17:LEU:HD22	1:CK:19:LEU:HD11	1.94	0.49
1:AX:86:PHE:CE1	1:AX:117:LEU:HD23	2.47	0.49
1:AX:95:ARG:NH2	1:AX:108:THR:OG1	2.44	0.49
1:BJ:86:PHE:CE1	1:BJ:117:LEU:HD23	2.47	0.49
1:BO:121:VAL:HG13	1:BO:125:THR:HB	1.93	0.49
1:BT:6:ILE:O	1:BT:6:ILE:HG13	2.11	0.49
1:BU:30:VAL:HA	1:DG:34:ASN:ND2	2.27	0.49
1:CC:6:ILE:O	1:CC:6:ILE:HG13	2.11	0.49
1:AC:86:PHE:CE1	1:AC:117:LEU:HD23	2.47	0.49
1:BJ:127:ALA:HB2	1:DF:153:GLN:OE1	2.11	0.49
1:BM:86:PHE:CE1	1:BM:117:LEU:HD23	2.47	0.49
1:BV:86:PHE:CE1	1:BV:117:LEU:HD23	2.47	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CE:95:ARG:NH2	1:CE:108:THR:OG1	2.44	0.49
1:AH:121:VAL:HG13	1:AH:125:THR:HB	1.93	0.49
1:BF:121:VAL:HG13	1:BF:125:THR:HB	1.93	0.49
1:CC:86:PHE:CE2	1:CC:117:LEU:HD23	2.48	0.49
1:DG:86:PHE:CE2	1:DG:117:LEU:HD23	2.48	0.49
1:AO:1:SER:C	1:CE:98:ASP:HB3	2.36	0.49
1:BA:86:PHE:CE1	1:BA:117:LEU:HD23	2.47	0.49
1:BZ:6:ILE:O	1:BZ:6:ILE:HG13	2.11	0.49
1:DB:121:VAL:HG13	1:DB:125:THR:HB	1.93	0.49
1:AI:159:LEU:HD13	1:BW:28:VAL:HG11	1.94	0.49
1:AU:86:PHE:CE1	1:AU:117:LEU:HD23	2.47	0.49
1:BP:86:PHE:CE1	1:BP:117:LEU:HD23	2.47	0.49
1:CI:86:PHE:CE2	1:CI:117:LEU:HD23	2.48	0.49
1:CM:121:VAL:HG13	1:CM:125:THR:HB	1.93	0.49
1:CN:86:PHE:CE1	1:CN:117:LEU:HD23	2.47	0.49
1:DA:86:PHE:CE2	1:DA:117:LEU:HD23	2.48	0.49
1:DI:86:PHE:CE1	1:DI:117:LEU:HD23	2.47	0.49
1:AC:164:GLY:O	1:BP:40:TYR:OH	2.27	0.49
1:AM:86:PHE:CE2	1:AM:117:LEU:HD23	2.48	0.49
1:AO:163:PHE:HD2	1:CE:24:ALA:HA	1.77	0.49
1:BG:86:PHE:CE1	1:BG:117:LEU:HD23	2.47	0.49
1:CF:86:PHE:CE2	1:CF:117:LEU:HD23	2.48	0.49
1:CW:86:PHE:CE1	1:CW:117:LEU:HD23	2.47	0.49
1:DC:86:PHE:CE1	1:DC:117:LEU:HD23	2.47	0.49
1:DG:6:ILE:O	1:DG:6:ILE:HG13	2.12	0.49
1:AR:44:GLY:HA3	1:CK:9:SER:OG	2.13	0.49
1:BE:86:PHE:CE2	1:BE:117:LEU:HD23	2.48	0.49
1:BH:86:PHE:CE2	1:BH:117:LEU:HD23	2.48	0.49
1:DF:86:PHE:CE1	1:DF:117:LEU:HD23	2.47	0.49
1:DJ:86:PHE:CE2	1:DJ:117:LEU:HD23	2.48	0.49
1:AY:86:PHE:CE2	1:AY:117:LEU:HD23	2.48	0.49
1:BH:34:ASN:HD22	1:BL:30:VAL:HA	1.74	0.49
1:BW:86:PHE:CE2	1:BW:117:LEU:HD23	2.48	0.49
1:CH:86:PHE:CE1	1:CH:117:LEU:HD23	2.47	0.49
1:CU:86:PHE:CE2	1:CU:117:LEU:HD23	2.48	0.49
1:DD:86:PHE:CE2	1:DD:117:LEU:HD23	2.48	0.49
1:DK:121:VAL:HG13	1:DK:125:THR:HB	1.93	0.49
1:AD:86:PHE:CE2	1:AD:117:LEU:HD23	2.48	0.49
1:AO:6:ILE:HG22	1:CE:94:TRP:NE1	2.28	0.49
1:AR:86:PHE:CE1	1:AR:117:LEU:HD23	2.47	0.49
1:AS:153:GLN:OE1	1:AT:127:ALA:HB2	2.13	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BY:86:PHE:CE1	1:BY:117:LEU:HD23	2.47	0.49
1:BZ:153:GLN:OE1	1:CA:127:ALA:HB2	2.13	0.49
1:CE:86:PHE:CE1	1:CE:117:LEU:HD23	2.47	0.49
1:CK:86:PHE:CE1	1:CK:117:LEU:HD23	2.47	0.49
1:AK:121:VAL:HG13	1:AK:125:THR:HB	1.93	0.48
1:AS:86:PHE:CE2	1:AS:117:LEU:HD23	2.48	0.48
1:BM:163:PHE:HD2	1:DI:24:ALA:HA	1.78	0.48
1:CX:153:GLN:OE1	1:CY:127:ALA:HB2	2.13	0.48
1:DJ:153:GLN:OE1	1:DK:127:ALA:HB2	2.13	0.48
1:AF:86:PHE:CE1	1:AF:117:LEU:HD23	2.47	0.48
1:AP:86:PHE:CE2	1:AP:117:LEU:HD23	2.48	0.48
1:BB:40:TYR:OH	1:BC:164:GLY:O	2.28	0.48
1:BB:86:PHE:CE2	1:BB:117:LEU:HD23	2.48	0.48
1:BE:153:GLN:OE1	1:BF:127:ALA:HB2	2.13	0.48
1:BK:153:GLN:OE1	1:BL:127:ALA:HB2	2.13	0.48
1:BR:121:VAL:HG13	1:BR:125:THR:HB	1.93	0.48
1:CL:153:GLN:OE1	1:CM:127:ALA:HB2	2.13	0.48
1:CO:153:GLN:OE1	1:CP:127:ALA:HB2	2.13	0.48
1:AD:153:GLN:OE1	1:AE:127:ALA:HB2	2.14	0.48
1:AL:86:PHE:CE1	1:AL:117:LEU:HD23	2.47	0.48
1:BE:6:ILE:HG13	1:BE:6:ILE:O	2.11	0.48
1:BN:86:PHE:CE2	1:BN:117:LEU:HD23	2.48	0.48
1:CI:153:GLN:OE1	1:CJ:127:ALA:HB2	2.13	0.48
1:CZ:86:PHE:CE1	1:CZ:117:LEU:HD23	2.47	0.48
1:AX:57:PHE:CE1	1:CT:162:ILE:HD12	2.49	0.48
1:BQ:153:GLN:OE1	1:BR:127:ALA:HB2	2.13	0.48
1:BZ:86:PHE:CE2	1:BZ:117:LEU:HD23	2.48	0.48
1:CB:86:PHE:CE1	1:CB:117:LEU:HD23	2.47	0.48
1:CL:86:PHE:CE2	1:CL:117:LEU:HD23	2.48	0.48
1:CU:153:GLN:OE1	1:CV:127:ALA:HB2	2.13	0.48
1:AH:50:LEU:HD11	1:CC:82:VAL:HG13	1.96	0.48
1:AU:57:PHE:CE1	1:CN:162:ILE:HD12	2.49	0.48
1:BQ:7:ASN:ND2	1:BR:46:ASP:H	2.12	0.48
1:CM:30:VAL:HA	1:CX:34:ASN:ND2	2.28	0.48
1:CO:86:PHE:CE2	1:CO:117:LEU:HD23	2.48	0.48
1:AG:1:SER:C	1:AH:98:ASP:HB3	2.39	0.48
1:AG:153:GLN:OE1	1:AH:127:ALA:HB2	2.13	0.48
1:AJ:1:SER:C	1:AK:98:ASP:HB3	2.39	0.48
1:AM:153:GLN:OE1	1:AN:127:ALA:HB2	2.13	0.48
1:AS:1:SER:C	1:AT:98:ASP:HB3	2.39	0.48
1:BB:153:GLN:OE1	1:BC:127:ALA:HB2	2.13	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BK:1:SER:C	1:BL:98:ASP:HB3	2.39	0.48
1:BT:7:ASN:ND2	1:BU:46:ASP:H	2.12	0.48
1:BZ:96:LYS:HB3	1:CA:4:ILE:HD13	1.96	0.48
1:AB:30:VAL:HA	1:AS:34:ASN:ND2	2.28	0.48
1:AD:1:SER:C	1:AE:98:ASP:HB3	2.39	0.48
1:AJ:86:PHE:CE2	1:AJ:117:LEU:HD23	2.48	0.48
1:AV:86:PHE:CE2	1:AV:117:LEU:HD23	2.48	0.48
1:AY:153:GLN:OE1	1:AZ:127:ALA:HB2	2.14	0.48
1:BH:1:SER:C	1:BI:98:ASP:HB3	2.39	0.48
1:BN:153:GLN:OE1	1:BO:127:ALA:HB2	2.13	0.48
1:CF:1:SER:C	1:CG:98:ASP:HB3	2.39	0.48
1:CI:1:SER:C	1:CJ:98:ASP:HB3	2.39	0.48
1:CR:7:ASN:ND2	1:CS:46:ASP:H	2.12	0.48
1:DD:1:SER:C	1:DE:98:ASP:HB3	2.39	0.48
1:DJ:1:SER:C	1:DK:98:ASP:HB3	2.39	0.48
1:DJ:7:ASN:ND2	1:DK:46:ASP:H	2.12	0.48
1:BE:1:SER:C	1:BF:98:ASP:HB3	2.39	0.48
1:BK:86:PHE:CE2	1:BK:117:LEU:HD23	2.48	0.48
1:BT:153:GLN:OE1	1:BU:127:ALA:HB2	2.13	0.48
1:CO:1:SER:C	1:CP:98:ASP:HB3	2.39	0.48
1:CO:96:LYS:HB3	1:CP:4:ILE:HD13	1.96	0.48
1:CX:1:SER:C	1:CY:98:ASP:HB3	2.39	0.48
1:DG:153:GLN:OE1	1:DH:127:ALA:HB2	2.13	0.48
1:AJ:153:GLN:OE1	1:AK:127:ALA:HB2	2.13	0.48
1:AS:57:PHE:CD1	1:AT:162:ILE:HD12	2.49	0.48
1:AT:50:LEU:HD11	1:AV:82:VAL:HG13	1.96	0.48
1:AX:16:ASN:OD1	1:CT:18:VAL:HG22	2.13	0.48
1:BA:40:TYR:CE1	1:CW:163:PHE:HB2	2.49	0.48
1:BG:78:ASP:HA	1:DC:107:TYR:CD1	2.49	0.48
1:BH:96:LYS:HB3	1:BI:4:ILE:HD13	1.96	0.48
1:BK:7:ASN:ND2	1:BL:46:ASP:H	2.12	0.48
1:BN:1:SER:C	1:BO:98:ASP:HB3	2.39	0.48
1:BQ:1:SER:C	1:BR:98:ASP:HB3	2.39	0.48
1:BT:57:PHE:CD1	1:BU:162:ILE:HD12	2.49	0.48
1:DD:153:GLN:OE1	1:DE:127:ALA:HB2	2.13	0.48
1:AD:7:ASN:ND2	1:AE:46:ASP:H	2.12	0.48
1:AP:153:GLN:OE1	1:AQ:127:ALA:HB2	2.13	0.48
1:AV:7:ASN:ND2	1:AW:46:ASP:H	2.12	0.48
1:AY:96:LYS:HB3	1:AZ:4:ILE:HD13	1.96	0.48
1:BH:153:GLN:OE1	1:BI:127:ALA:HB2	2.13	0.48
1:BT:157:ASP:OD1	1:DI:47:ASN:OD1	2.31	0.48

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CF:7:ASN:ND2	1:CG:46:ASP:H	2.12	0.48
1:CF:57:PHE:CD1	1:CG:162:ILE:HD12	2.49	0.48
1:CR:57:PHE:CD1	1:CS:162:ILE:HD12	2.49	0.48
1:CU:57:PHE:CD1	1:CV:162:ILE:HD12	2.49	0.48
1:CU:96:LYS:HB3	1:CV:4:ILE:HD13	1.96	0.48
1:DA:96:LYS:HB3	1:DB:4:ILE:HD13	1.96	0.48
1:DA:153:GLN:OE1	1:DB:127:ALA:HB2	2.13	0.48
1:DG:1:SER:C	1:DH:98:ASP:HB3	2.39	0.48
1:DJ:96:LYS:HB3	1:DK:4:ILE:HD13	1.96	0.48
1:AM:1:SER:C	1:AN:98:ASP:HB3	2.39	0.47
1:AM:57:PHE:CD1	1:AN:162:ILE:HD12	2.49	0.47
1:AM:96:LYS:HB3	1:AN:4:ILE:HD13	1.96	0.47
1:BN:96:LYS:HB3	1:BO:4:ILE:HD13	1.96	0.47
1:BT:86:PHE:CE2	1:BT:117:LEU:HD23	2.48	0.47
1:BW:153:GLN:OE1	1:BX:127:ALA:HB2	2.13	0.47
1:CC:96:LYS:HB3	1:CD:4:ILE:HD13	1.96	0.47
1:CI:57:PHE:CD1	1:CJ:162:ILE:HD12	2.49	0.47
1:CL:96:LYS:HB3	1:CM:4:ILE:HD13	1.96	0.47
1:CU:1:SER:C	1:CV:98:ASP:HB3	2.39	0.47
1:CX:96:LYS:HB3	1:CY:4:ILE:HD13	1.96	0.47
1:BE:57:PHE:CD1	1:BF:162:ILE:HD12	2.49	0.47
1:BJ:2:TYR:CD1	1:DF:96:LYS:HE2	2.49	0.47
1:CI:7:ASN:ND2	1:CJ:46:ASP:H	2.12	0.47
1:DA:7:ASN:ND2	1:DB:46:ASP:H	2.12	0.47
1:DD:96:LYS:HB3	1:DE:4:ILE:HD13	1.96	0.47
1:DG:57:PHE:CD1	1:DH:162:ILE:HD12	2.49	0.47
1:AA:57:PHE:CD1	1:AB:162:ILE:HD12	2.49	0.47
1:AA:153:GLN:OE1	1:AB:127:ALA:HB2	2.13	0.47
1:AC:40:TYR:OH	1:BP:164:GLY:O	2.26	0.47
1:AS:96:LYS:HB3	1:AT:4:ILE:HD13	1.96	0.47
1:AV:153:GLN:OE1	1:AW:127:ALA:HB2	2.13	0.47
1:AY:1:SER:C	1:AZ:98:ASP:HB3	2.39	0.47
1:BE:157:ASP:OD1	1:BS:47:ASN:OD1	2.32	0.47
1:BQ:57:PHE:CD1	1:BR:162:ILE:HD12	2.49	0.47
1:BQ:86:PHE:CE2	1:BQ:117:LEU:HD23	2.48	0.47
1:BW:57:PHE:CD1	1:BX:162:ILE:HD12	2.49	0.47
1:BW:96:LYS:HB3	1:BX:4:ILE:HD13	1.96	0.47
1:CC:153:GLN:OE1	1:CD:127:ALA:HB2	2.13	0.47
1:DJ:57:PHE:CD1	1:DK:162:ILE:HD12	2.49	0.47
1:AD:57:PHE:CD1	1:AE:162:ILE:HD12	2.49	0.47
1:AG:57:PHE:CD1	1:AH:162:ILE:HD12	2.49	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AP:96:LYS:HB3	1:AQ:4:ILE:HD13	1.96	0.47
1:BQ:96:LYS:HB3	1:BR:4:ILE:HD13	1.96	0.47
1:CF:153:GLN:OE1	1:CG:127:ALA:HB2	2.13	0.47
1:DA:57:PHE:CD1	1:DB:162:ILE:HD12	2.49	0.47
1:AA:1:SER:C	1:AB:98:ASP:HB3	2.39	0.47
1:AJ:57:PHE:CD1	1:AK:162:ILE:HD12	2.49	0.47
1:AU:162:ILE:HD12	1:CN:57:PHE:CE1	2.49	0.47
1:AV:96:LYS:HB3	1:AW:4:ILE:HD13	1.96	0.47
1:BA:111:GLN:OE1	1:CW:119:SER:HB2	2.14	0.47
1:BQ:23:PRO:O	1:BQ:26:PRO:HD2	2.15	0.47
1:CL:1:SER:C	1:CM:98:ASP:HB3	2.39	0.47
1:CO:7:ASN:ND2	1:CP:46:ASP:H	2.12	0.47
1:DD:57:PHE:CD1	1:DE:162:ILE:HD12	2.49	0.47
1:AA:23:PRO:O	1:AA:26:PRO:HD2	2.15	0.47
1:AA:86:PHE:CE2	1:AA:117:LEU:HD23	2.48	0.47
1:AE:86:PHE:HE1	1:AE:117:LEU:HD23	1.80	0.47
1:AF:46:ASP:H	1:BV:7:ASN:HD21	1.62	0.47
1:AG:7:ASN:ND2	1:AH:46:ASP:H	2.12	0.47
1:AK:86:PHE:HE1	1:AK:117:LEU:HD23	1.80	0.47
1:AV:1:SER:C	1:AW:98:ASP:HB3	2.39	0.47
1:BW:1:SER:C	1:BX:98:ASP:HB3	2.39	0.47
1:BZ:1:SER:C	1:CA:98:ASP:HB3	2.39	0.47
1:CC:7:ASN:ND2	1:CD:46:ASP:H	2.12	0.47
1:CD:86:PHE:HE1	1:CD:117:LEU:HD23	1.80	0.47
1:CF:23:PRO:O	1:CF:26:PRO:HD2	2.15	0.47
1:CO:57:PHE:CD1	1:CP:162:ILE:HD12	2.49	0.47
1:CR:153:GLN:OE1	1:CS:127:ALA:HB2	2.13	0.47
1:CX:23:PRO:O	1:CX:26:PRO:HD2	2.15	0.47
1:DA:23:PRO:O	1:DA:26:PRO:HD2	2.15	0.47
1:AI:163:PHE:HD2	1:BY:24:ALA:HA	1.78	0.47
1:AN:86:PHE:HE1	1:AN:117:LEU:HD23	1.80	0.47
1:AP:7:ASN:ND2	1:AQ:46:ASP:H	2.12	0.47
1:AP:57:PHE:CD1	1:AQ:162:ILE:HD12	2.49	0.47
1:AV:57:PHE:CD1	1:AW:162:ILE:HD12	2.49	0.47
1:AY:57:PHE:CD1	1:AZ:162:ILE:HD12	2.49	0.47
1:BB:23:PRO:O	1:BB:26:PRO:HD2	2.15	0.47
1:BE:23:PRO:O	1:BE:26:PRO:HD2	2.15	0.47
1:BE:96:LYS:HB3	1:BF:4:ILE:HD13	1.96	0.47
1:BH:57:PHE:CD1	1:BI:162:ILE:HD12	2.49	0.47
1:BM:24:ALA:HA	1:DI:163:PHE:CD2	2.50	0.47
1:BW:23:PRO:O	1:BW:26:PRO:HD2	2.15	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CI:23:PRO:O	1:CI:26:PRO:HD2	2.15	0.47
1:CO:23:PRO:O	1:CO:26:PRO:HD2	2.15	0.47
1:CX:86:PHE:CE2	1:CX:117:LEU:HD23	2.48	0.47
1:DA:1:SER:C	1:DB:98:ASP:HB3	2.39	0.47
1:BB:96:LYS:HB3	1:BC:4:ILE:HD13	1.96	0.47
1:BF:86:PHE:HE1	1:BF:117:LEU:HD23	1.80	0.47
1:BJ:96:LYS:HE2	1:DF:2:TYR:CD1	2.49	0.47
1:BJ:96:LYS:HA	1:DF:3:THR:O	2.14	0.47
1:BN:57:PHE:CD1	1:BO:162:ILE:HD12	2.49	0.47
1:BT:1:SER:C	1:BU:98:ASP:HB3	2.39	0.47
1:BT:96:LYS:HB3	1:BU:4:ILE:HD13	1.96	0.47
1:CS:86:PHE:HE1	1:CS:117:LEU:HD23	1.80	0.47
1:AA:7:ASN:ND2	1:AB:46:ASP:H	2.12	0.47
1:AH:30:VAL:HA	1:CC:34:ASN:HD22	1.78	0.47
1:AJ:96:LYS:HB3	1:AK:4:ILE:HD13	1.96	0.47
1:AM:52:ARG:HH22	1:CE:157:ASP:CG	2.23	0.47
1:BB:57:PHE:CD1	1:BC:162:ILE:HD12	2.49	0.47
1:BG:155:LEU:HD13	1:DC:86:PHE:CD2	2.50	0.47
1:BK:57:PHE:CD1	1:BL:162:ILE:HD12	2.49	0.47
1:CC:57:PHE:CD1	1:CD:162:ILE:HD12	2.49	0.47
1:DC:46:ASP:OD2	1:DC:48:SER:OG	2.27	0.47
1:AF:7:ASN:O	1:BV:53:PRO:HG2	2.14	0.47
1:AG:86:PHE:CE2	1:AG:117:LEU:HD23	2.48	0.47
1:AR:94:TRP:CD1	1:CK:6:ILE:HA	2.50	0.47
1:AU:164:GLY:O	1:CN:40:TYR:OH	2.32	0.47
1:BB:1:SER:C	1:BC:98:ASP:HB3	2.39	0.47
1:BG:157:ASP:CG	1:DA:52:ARG:NH2	2.68	0.47
1:BJ:162:ILE:HD12	1:DF:57:PHE:CE1	2.50	0.47
1:BK:96:LYS:HB3	1:BL:4:ILE:HD13	1.96	0.47
1:BP:46:ASP:OD2	1:BP:48:SER:OG	2.27	0.47
1:CA:86:PHE:HE1	1:CA:117:LEU:HD23	1.80	0.47
1:CC:1:SER:C	1:CD:98:ASP:HB3	2.39	0.47
1:CL:57:PHE:CD1	1:CM:162:ILE:HD12	2.49	0.47
1:CU:23:PRO:O	1:CU:26:PRO:HD2	2.15	0.47
1:CW:155:LEU:HD23	1:CW:155:LEU:HA	1.77	0.47
1:DE:86:PHE:HE1	1:DE:117:LEU:HD23	1.80	0.47
1:DG:33:ASP:HB2	1:DG:36:SER:H	1.80	0.47
1:AD:23:PRO:O	1:AD:26:PRO:HD2	2.15	0.46
1:AF:111:GLN:OE1	1:BV:119:SER:HB2	2.15	0.46
1:AH:86:PHE:HE1	1:AH:117:LEU:HD23	1.80	0.46
1:AR:30:VAL:HG21	1:AR:41:LEU:HB2	1.98	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AR:57:PHE:CE1	1:CK:162:ILE:HD12	2.50	0.46
1:AS:33:ASP:HB2	1:AS:36:SER:H	1.81	0.46
1:AU:40:TYR:CE1	1:CN:163:PHE:HB2	2.50	0.46
1:BH:7:ASN:ND2	1:BI:46:ASP:H	2.12	0.46
1:BI:86:PHE:HE1	1:BI:117:LEU:HD23	1.80	0.46
1:BZ:23:PRO:O	1:BZ:26:PRO:HD2	2.15	0.46
1:CJ:86:PHE:HE1	1:CJ:117:LEU:HD23	1.80	0.46
1:AG:96:LYS:HB3	1:AH:4:ILE:HD13	1.96	0.46
1:AI:30:VAL:HG21	1:AI:41:LEU:HB2	1.98	0.46
1:AY:23:PRO:O	1:AY:26:PRO:HD2	2.15	0.46
1:BA:162:ILE:HD12	1:CW:57:PHE:CE1	2.50	0.46
1:BJ:161:ARG:HD2	1:DF:34:ASN:HB3	1.97	0.46
1:BM:119:SER:HB2	1:DI:111:GLN:OE1	2.15	0.46
1:BU:29:HIS:O	1:DG:34:ASN:ND2	2.48	0.46
1:BV:30:VAL:HG21	1:BV:41:LEU:HB2	1.98	0.46
1:BW:7:ASN:ND2	1:BX:46:ASP:H	2.12	0.46
1:CR:1:SER:C	1:CS:98:ASP:HB3	2.39	0.46
1:DD:23:PRO:O	1:DD:26:PRO:HD2	2.15	0.46
1:DK:86:PHE:HE1	1:DK:117:LEU:HD23	1.80	0.46
1:AG:23:PRO:O	1:AG:26:PRO:HD2	2.15	0.46
1:AO:34:ASN:HB3	1:CE:161:ARG:HD2	1.97	0.46
1:AP:1:SER:C	1:AQ:98:ASP:HB3	2.39	0.46
1:AV:23:PRO:O	1:AV:26:PRO:HD2	2.15	0.46
1:BA:24:ALA:HA	1:CW:163:PHE:HD2	1.80	0.46
1:BC:86:PHE:HE1	1:BC:117:LEU:HD23	1.80	0.46
1:BD:30:VAL:HG21	1:BD:41:LEU:HB2	1.97	0.46
1:BN:7:ASN:ND2	1:BO:46:ASP:H	2.12	0.46
1:BP:30:VAL:HG21	1:BP:41:LEU:HB2	1.98	0.46
1:BT:23:PRO:O	1:BT:26:PRO:HD2	2.15	0.46
1:BZ:57:PHE:CD1	1:CA:162:ILE:HD12	2.49	0.46
1:CC:33:ASP:HB2	1:CC:36:SER:H	1.81	0.46
1:CI:96:LYS:HB3	1:CJ:4:ILE:HD13	1.96	0.46
1:CO:33:ASP:HB2	1:CO:36:SER:H	1.81	0.46
1:CR:86:PHE:CE2	1:CR:117:LEU:HD23	2.48	0.46
1:CZ:30:VAL:HG21	1:CZ:41:LEU:HB2	1.98	0.46
1:DJ:23:PRO:O	1:DJ:26:PRO:HD2	2.15	0.46
1:AD:52:ARG:NH2	1:BV:157:ASP:CG	2.70	0.46
1:AJ:23:PRO:O	1:AJ:26:PRO:HD2	2.15	0.46
1:AS:7:ASN:ND2	1:AT:46:ASP:H	2.12	0.46
1:AY:33:ASP:HB2	1:AY:36:SER:H	1.81	0.46
1:BG:38:ILE:HG22	1:DC:160:PRO:HB2	1.97	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CX:57:PHE:CD1	1:CY:162:ILE:HD12	2.49	0.46
1:AD:96:LYS:HB3	1:AE:4:ILE:HD13	1.96	0.46
1:AL:40:TYR:OH	1:CB:164:GLY:O	2.30	0.46
1:AM:23:PRO:O	1:AM:26:PRO:HD2	2.15	0.46
1:BK:23:PRO:O	1:BK:26:PRO:HD2	2.15	0.46
1:CE:30:VAL:HG21	1:CE:41:LEU:HB2	1.98	0.46
1:CG:155:LEU:HD11	1:CG:162:ILE:HD11	1.98	0.46
1:CL:23:PRO:O	1:CL:26:PRO:HD2	2.15	0.46
1:CR:23:PRO:O	1:CR:26:PRO:HD2	2.15	0.46
1:CX:7:ASN:ND2	1:CY:46:ASP:H	2.12	0.46
1:DL:30:VAL:HG21	1:DL:41:LEU:HB2	1.98	0.46
1:AA:96:LYS:HB3	1:AB:4:ILE:HD13	1.96	0.46
1:AD:28:VAL:HG11	1:BV:159:LEU:HD13	1.98	0.46
1:AD:33:ASP:HB2	1:AD:36:SER:H	1.81	0.46
1:AF:161:ARG:HD2	1:BV:34:ASN:HB3	1.97	0.46
1:AP:23:PRO:O	1:AP:26:PRO:HD2	2.15	0.46
1:AT:155:LEU:HD11	1:AT:162:ILE:HD11	1.98	0.46
1:BN:23:PRO:O	1:BN:26:PRO:HD2	2.15	0.46
1:CC:23:PRO:O	1:CC:26:PRO:HD2	2.15	0.46
1:CG:86:PHE:HE1	1:CG:117:LEU:HD23	1.80	0.46
1:CV:86:PHE:HE1	1:CV:117:LEU:HD23	1.80	0.46
1:DG:7:ASN:ND2	1:DH:46:ASP:H	2.12	0.46
1:DI:30:VAL:HG21	1:DI:41:LEU:HB2	1.98	0.46
1:AF:30:VAL:HG21	1:AF:41:LEU:HB2	1.98	0.46
1:AM:33:ASP:HB2	1:AM:36:SER:H	1.81	0.46
1:AP:33:ASP:HB2	1:AP:36:SER:H	1.81	0.46
1:BF:155:LEU:HD11	1:BF:162:ILE:HD11	1.98	0.46
1:BM:30:VAL:HG21	1:BM:41:LEU:HB2	1.98	0.46
1:BZ:33:ASP:HB2	1:BZ:36:SER:H	1.81	0.46
1:CB:30:VAL:HG21	1:CB:41:LEU:HB2	1.98	0.46
1:CF:96:LYS:HB3	1:CG:4:ILE:HD13	1.96	0.46
1:AB:155:LEU:HD11	1:AB:162:ILE:HD11	1.98	0.46
1:AO:119:SER:HB2	1:CE:111:GLN:OE1	2.15	0.46
1:AS:23:PRO:O	1:AS:26:PRO:HD2	2.15	0.46
1:AY:28:VAL:HG11	1:CW:159:LEU:HD13	1.98	0.46
1:BB:7:ASN:ND2	1:BC:46:ASP:H	2.12	0.46
1:BH:23:PRO:O	1:BH:26:PRO:HD2	2.15	0.46
1:BJ:24:ALA:HA	1:DF:163:PHE:HD2	1.80	0.46
1:BN:33:ASP:HB2	1:BN:36:SER:H	1.81	0.46
1:BW:33:ASP:HB2	1:BW:36:SER:H	1.81	0.46
1:BY:30:VAL:HG21	1:BY:41:LEU:HB2	1.98	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CK:30:VAL:HG21	1:CK:41:LEU:HB2	1.97	0.46
1:CM:86:PHE:HE1	1:CM:117:LEU:HD23	1.80	0.46
1:DE:37:THR:HG23	1:DE:58:ASN:HD22	1.81	0.46
1:DF:30:VAL:HG21	1:DF:41:LEU:HB2	1.97	0.46
1:DG:96:LYS:HB3	1:DH:4:ILE:HD13	1.96	0.46
1:DJ:33:ASP:HB2	1:DJ:36:SER:H	1.81	0.46
1:AT:37:THR:HG23	1:AT:58:ASN:HD22	1.81	0.46
1:BC:155:LEU:HD11	1:BC:162:ILE:HD11	1.98	0.46
1:BJ:94:TRP:CD1	1:DF:6:ILE:HA	2.51	0.46
1:BK:33:ASP:HB2	1:BK:36:SER:H	1.81	0.46
1:BL:86:PHE:HE1	1:BL:117:LEU:HD23	1.80	0.46
1:CD:37:THR:HG23	1:CD:58:ASN:HD22	1.81	0.46
1:CR:96:LYS:HB3	1:CS:4:ILE:HD13	1.96	0.46
1:DA:33:ASP:HB2	1:DA:36:SER:H	1.81	0.46
1:DC:30:VAL:HG21	1:DC:41:LEU:HB2	1.98	0.46
1:AJ:7:ASN:ND2	1:AK:46:ASP:H	2.12	0.46
1:AK:155:LEU:HD11	1:AK:162:ILE:HD11	1.98	0.46
1:AL:163:PHE:HB2	1:CB:40:TYR:CE1	2.50	0.46
1:BA:30:VAL:HG21	1:BA:41:LEU:HB2	1.98	0.46
1:BF:37:THR:HG23	1:BF:58:ASN:HD22	1.81	0.46
1:BJ:86:PHE:CD2	1:DF:155:LEU:HD13	2.51	0.46
1:BR:37:THR:HG23	1:BR:58:ASN:HD22	1.81	0.46
1:CR:33:ASP:HB2	1:CR:36:SER:H	1.81	0.46
1:DG:23:PRO:O	1:DG:26:PRO:HD2	2.15	0.46
1:AC:162:ILE:HD12	1:BP:57:PHE:CE1	2.51	0.45
1:AH:37:THR:HG23	1:AH:58:ASN:HD22	1.81	0.45
1:AO:7:ASN:HD21	1:CE:46:ASP:H	1.64	0.45
1:AX:98:ASP:HB3	1:CT:1:SER:C	2.41	0.45
1:BA:84:ARG:HD2	1:CW:156:ILE:O	2.16	0.45
1:BA:163:PHE:HB2	1:CW:40:TYR:CE1	2.51	0.45
1:BH:33:ASP:HB2	1:BH:36:SER:H	1.81	0.45
1:BX:155:LEU:HD11	1:BX:162:ILE:HD11	1.98	0.45
1:CD:155:LEU:HD11	1:CD:162:ILE:HD11	1.98	0.45
1:CG:37:THR:HG23	1:CG:58:ASN:HD22	1.81	0.45
1:CI:33:ASP:HB2	1:CI:36:SER:H	1.81	0.45
1:CL:7:ASN:ND2	1:CM:46:ASP:H	2.12	0.45
1:CM:155:LEU:HD11	1:CM:162:ILE:HD11	1.98	0.45
1:CS:37:THR:HG23	1:CS:58:ASN:HD22	1.81	0.45
1:AA:33:ASP:HB2	1:AA:36:SER:H	1.81	0.45
1:AJ:33:ASP:HB2	1:AJ:36:SER:H	1.81	0.45
1:AL:57:PHE:CE1	1:CB:162:ILE:HD12	2.51	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AQ:155:LEU:HD11	1:AQ:162:ILE:HD11	1.98	0.45
1:AY:7:ASN:ND2	1:AZ:46:ASP:H	2.12	0.45
1:BE:7:ASN:ND2	1:BF:46:ASP:H	2.12	0.45
1:BN:155:LEU:HG	1:BN:162:ILE:HD11	1.99	0.45
1:BS:30:VAL:HG21	1:BS:41:LEU:HB2	1.98	0.45
1:CA:37:THR:HG23	1:CA:58:ASN:HD22	1.81	0.45
1:DB:155:LEU:HD11	1:DB:162:ILE:HD11	1.98	0.45
1:AC:30:VAL:HG21	1:AC:41:LEU:HB2	1.98	0.45
1:AW:86:PHE:HE1	1:AW:117:LEU:HD23	1.80	0.45
1:AW:155:LEU:HD11	1:AW:162:ILE:HD11	1.98	0.45
1:BJ:30:VAL:HG21	1:BJ:41:LEU:HB2	1.98	0.45
1:BO:151:LYS:HE3	1:BO:162:ILE:HG23	1.99	0.45
1:BT:155:LEU:HG	1:BT:162:ILE:HD11	1.99	0.45
1:CI:155:LEU:HG	1:CI:162:ILE:HD11	1.99	0.45
1:CX:155:LEU:HG	1:CX:162:ILE:HD11	1.99	0.45
1:DD:33:ASP:HB2	1:DD:36:SER:H	1.81	0.45
1:AE:37:THR:HG23	1:AE:58:ASN:HD22	1.81	0.45
1:AL:86:PHE:HE1	1:AL:117:LEU:HD23	1.82	0.45
1:AP:155:LEU:HG	1:AP:162:ILE:HD11	1.99	0.45
1:AZ:155:LEU:HD11	1:AZ:162:ILE:HD11	1.98	0.45
1:BB:155:LEU:HG	1:BB:162:ILE:HD11	1.99	0.45
1:BU:37:THR:HG23	1:BU:58:ASN:HD22	1.81	0.45
1:CF:33:ASP:HB2	1:CF:36:SER:H	1.81	0.45
1:CG:151:LYS:HE3	1:CG:162:ILE:HG23	1.99	0.45
1:CP:151:LYS:HE3	1:CP:162:ILE:HG23	1.99	0.45
1:CV:155:LEU:HD11	1:CV:162:ILE:HD11	1.98	0.45
1:CZ:86:PHE:HE1	1:CZ:117:LEU:HD23	1.82	0.45
1:DC:86:PHE:HE1	1:DC:117:LEU:HD23	1.82	0.45
1:DH:155:LEU:HD11	1:DH:162:ILE:HD11	1.98	0.45
1:AB:86:PHE:HE1	1:AB:117:LEU:HD23	1.80	0.45
1:AE:155:LEU:HD11	1:AE:162:ILE:HD11	1.98	0.45
1:AK:37:THR:HG23	1:AK:58:ASN:HD22	1.81	0.45
1:AT:86:PHE:HE1	1:AT:117:LEU:HD23	1.80	0.45
1:AU:30:VAL:HG21	1:AU:41:LEU:HB2	1.98	0.45
1:AU:111:GLN:OE1	1:CN:119:SER:HB2	2.16	0.45
1:AX:86:PHE:HE1	1:AX:117:LEU:HD23	1.82	0.45
1:AZ:86:PHE:HE1	1:AZ:117:LEU:HD23	1.80	0.45
1:BC:37:THR:HG23	1:BC:58:ASN:HD22	1.81	0.45
1:BG:86:PHE:HE1	1:BG:117:LEU:HD23	1.82	0.45
1:BL:151:LYS:HE3	1:BL:162:ILE:HG23	1.99	0.45
1:BM:86:PHE:HE1	1:BM:117:LEU:HD23	1.82	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BR:151:LYS:HE3	1:BR:162:ILE:HG23	1.99	0.45
1:DE:151:LYS:HE3	1:DE:162:ILE:HG23	1.99	0.45
1:DK:155:LEU:HD11	1:DK:162:ILE:HD11	1.98	0.45
1:AC:46:ASP:OD2	1:AC:48:SER:OG	2.27	0.45
1:AH:155:LEU:HD11	1:AH:162:ILE:HD11	1.98	0.45
1:AJ:155:LEU:HG	1:AJ:162:ILE:HD11	1.99	0.45
1:AU:40:TYR:OH	1:CN:164:GLY:O	2.31	0.45
1:AZ:37:THR:HG23	1:AZ:58:ASN:HD22	1.81	0.45
1:BA:86:PHE:HE1	1:BA:117:LEU:HD23	1.82	0.45
1:BB:33:ASP:HB2	1:BB:36:SER:H	1.81	0.45
1:BD:7:ASN:HD21	1:CZ:46:ASP:H	1.65	0.45
1:BG:24:ALA:HA	1:DC:163:PHE:CD2	2.52	0.45
1:BG:30:VAL:HG21	1:BG:41:LEU:HB2	1.98	0.45
1:BJ:86:PHE:HE1	1:BJ:117:LEU:HD23	1.82	0.45
1:BK:155:LEU:HG	1:BK:162:ILE:HD11	1.99	0.45
1:BL:37:THR:HG23	1:BL:58:ASN:HD22	1.81	0.45
1:BL:155:LEU:HD11	1:BL:162:ILE:HD11	1.98	0.45
1:BT:33:ASP:HB2	1:BT:36:SER:H	1.81	0.45
1:CB:86:PHE:HE1	1:CB:117:LEU:HD23	1.82	0.45
1:CC:155:LEU:HG	1:CC:162:ILE:HD11	1.98	0.45
1:CH:86:PHE:HE1	1:CH:117:LEU:HD23	1.82	0.45
1:CJ:37:THR:HG23	1:CJ:58:ASN:HD22	1.81	0.45
1:CM:151:LYS:HE3	1:CM:162:ILE:HG23	1.99	0.45
1:CR:155:LEU:HG	1:CR:162:ILE:HD11	1.98	0.45
1:CU:33:ASP:HB2	1:CU:36:SER:H	1.81	0.45
1:CY:155:LEU:HD11	1:CY:162:ILE:HD11	1.98	0.45
1:AE:151:LYS:HE3	1:AE:162:ILE:HG23	1.99	0.45
1:AF:155:LEU:HA	1:AF:155:LEU:HD23	1.77	0.45
1:AO:30:VAL:HG21	1:AO:41:LEU:HB2	1.98	0.45
1:AO:163:PHE:HB2	1:CE:40:TYR:CE1	2.51	0.45
1:BE:33:ASP:HB2	1:BE:36:SER:H	1.81	0.45
1:BF:151:LYS:HE3	1:BF:162:ILE:HG23	1.99	0.45
1:BJ:57:PHE:CE1	1:DF:162:ILE:HD12	2.52	0.45
1:BU:155:LEU:HD11	1:BU:162:ILE:HD11	1.98	0.45
1:BX:86:PHE:HE1	1:BX:117:LEU:HD23	1.80	0.45
1:BX:151:LYS:HE3	1:BX:162:ILE:HG23	1.99	0.45
1:CO:155:LEU:HG	1:CO:162:ILE:HD11	1.99	0.45
1:CQ:30:VAL:HG21	1:CQ:41:LEU:HB2	1.98	0.45
1:DE:155:LEU:HD11	1:DE:162:ILE:HD11	1.98	0.45
1:DJ:155:LEU:HG	1:DJ:162:ILE:HD11	1.99	0.45
1:AL:30:VAL:HG21	1:AL:41:LEU:HB2	1.98	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AQ:37:THR:HG23	1:AQ:58:ASN:HD22	1.81	0.45
1:AR:86:PHE:HE1	1:AR:117:LEU:HD23	1.82	0.45
1:AR:111:GLN:OE1	1:CK:119:SER:HB2	2.16	0.45
1:BH:155:LEU:HG	1:BH:162:ILE:HD11	1.99	0.45
1:BI:37:THR:HG23	1:BI:58:ASN:HD22	1.81	0.45
1:CA:151:LYS:HE3	1:CA:162:ILE:HG23	1.99	0.45
1:CD:151:LYS:HE3	1:CD:162:ILE:HG23	1.99	0.45
1:CE:86:PHE:HE1	1:CE:117:LEU:HD23	1.82	0.45
1:CJ:155:LEU:HD11	1:CJ:162:ILE:HD11	1.98	0.45
1:DD:155:LEU:HG	1:DD:162:ILE:HD11	1.99	0.45
1:DG:155:LEU:HG	1:DG:162:ILE:HD11	1.99	0.45
1:DL:86:PHE:HE1	1:DL:117:LEU:HD23	1.82	0.45
1:AB:37:THR:HG23	1:AB:58:ASN:HD22	1.81	0.45
1:AD:155:LEU:HG	1:AD:162:ILE:HD11	1.99	0.45
1:AG:33:ASP:HB2	1:AG:36:SER:H	1.81	0.45
1:AI:7:ASN:HD21	1:BY:46:ASP:H	1.65	0.45
1:AU:163:PHE:CD2	1:CN:24:ALA:HA	2.49	0.45
1:AX:7:ASN:O	1:CT:53:PRO:HG2	2.16	0.45
1:AZ:151:LYS:HE3	1:AZ:162:ILE:HG23	1.99	0.45
1:CH:30:VAL:HG21	1:CH:41:LEU:HB2	1.98	0.45
1:CT:30:VAL:HG21	1:CT:41:LEU:HB2	1.98	0.45
1:CV:37:THR:HG23	1:CV:58:ASN:HD22	1.81	0.45
1:CX:33:ASP:HB2	1:CX:36:SER:H	1.81	0.45
1:AK:151:LYS:HE3	1:AK:162:ILE:HG23	1.99	0.45
1:AR:40:TYR:OH	1:CK:164:GLY:O	2.26	0.45
1:AR:94:TRP:CD1	1:CK:6:ILE:HG22	2.50	0.45
1:AU:46:ASP:H	1:CN:7:ASN:HD21	1.65	0.45
1:BA:40:TYR:OH	1:CW:164:GLY:O	2.32	0.45
1:BC:151:LYS:HE3	1:BC:162:ILE:HG23	1.99	0.45
1:BI:155:LEU:HD11	1:BI:162:ILE:HD11	1.98	0.45
1:CK:86:PHE:HE1	1:CK:117:LEU:HD23	1.82	0.45
1:CU:53:PRO:HG2	1:CV:7:ASN:O	2.17	0.45
1:CU:155:LEU:HG	1:CU:162:ILE:HD11	1.99	0.45
1:CW:30:VAL:HG21	1:CW:41:LEU:HB2	1.98	0.45
1:CW:86:PHE:HE1	1:CW:117:LEU:HD23	1.82	0.45
1:DB:37:THR:HG23	1:DB:58:ASN:HD22	1.81	0.45
1:AA:53:PRO:HG2	1:AB:7:ASN:O	2.18	0.44
1:AI:86:PHE:HE1	1:AI:117:LEU:HD23	1.82	0.44
1:AV:53:PRO:HG2	1:AW:7:ASN:O	2.18	0.44
1:BJ:157:ASP:CG	1:DD:52:ARG:HH22	2.25	0.44
1:BQ:33:ASP:HB2	1:BQ:36:SER:H	1.81	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BQ:155:LEU:HG	1:BQ:162:ILE:HD11	1.99	0.44
1:BY:86:PHE:HE1	1:BY:117:LEU:HD23	1.82	0.44
1:BZ:7:ASN:ND2	1:CA:46:ASP:H	2.12	0.44
1:CA:155:LEU:HD11	1:CA:162:ILE:HD11	1.98	0.44
1:CJ:151:LYS:HE3	1:CJ:162:ILE:HG23	1.99	0.44
1:CO:53:PRO:HG2	1:CP:7:ASN:O	2.17	0.44
1:CS:151:LYS:HE3	1:CS:162:ILE:HG23	1.99	0.44
1:CV:151:LYS:HE3	1:CV:162:ILE:HG23	1.99	0.44
1:CY:37:THR:HG23	1:CY:58:ASN:HD22	1.81	0.44
1:DH:37:THR:HG23	1:DH:58:ASN:HD22	1.81	0.44
1:AG:53:PRO:HG2	1:AH:7:ASN:O	2.18	0.44
1:AM:7:ASN:ND2	1:AN:46:ASP:H	2.12	0.44
1:AN:37:THR:HG23	1:AN:58:ASN:HD22	1.81	0.44
1:AR:57:PHE:CD1	1:CK:162:ILE:HD12	2.52	0.44
1:AS:155:LEU:HG	1:AS:162:ILE:HD11	1.99	0.44
1:AU:86:PHE:HE1	1:AU:117:LEU:HD23	1.82	0.44
1:AV:33:ASP:HB2	1:AV:36:SER:H	1.81	0.44
1:AX:30:VAL:HG21	1:AX:41:LEU:HB2	1.97	0.44
1:AY:53:PRO:HG2	1:AZ:7:ASN:O	2.18	0.44
1:BO:37:THR:HG23	1:BO:58:ASN:HD22	1.81	0.44
1:BO:155:LEU:HD11	1:BO:162:ILE:HD11	1.98	0.44
1:CJ:30:VAL:HA	1:DA:34:ASN:ND2	2.32	0.44
1:CN:30:VAL:HG21	1:CN:41:LEU:HB2	1.98	0.44
1:CP:86:PHE:HE1	1:CP:117:LEU:HD23	1.80	0.44
1:CP:155:LEU:HD11	1:CP:162:ILE:HD11	1.98	0.44
1:DD:53:PRO:HG2	1:DE:7:ASN:O	2.17	0.44
1:AA:155:LEU:HG	1:AA:162:ILE:HD11	1.99	0.44
1:AO:86:PHE:HE1	1:AO:117:LEU:HD23	1.82	0.44
1:BH:53:PRO:HG2	1:BI:7:ASN:O	2.18	0.44
1:BR:86:PHE:HE1	1:BR:117:LEU:HD23	1.80	0.44
1:BR:155:LEU:HD11	1:BR:162:ILE:HD11	1.98	0.44
1:BS:33:ASP:OD1	1:BS:33:ASP:N	2.51	0.44
1:BW:53:PRO:HG2	1:BX:7:ASN:O	2.18	0.44
1:CS:155:LEU:HD11	1:CS:162:ILE:HD11	1.98	0.44
1:CW:33:ASP:OD1	1:CW:33:ASP:N	2.51	0.44
1:DH:151:LYS:HE3	1:DH:162:ILE:HG23	1.99	0.44
1:AE:63:GLN:HG3	1:BG:50:LEU:HD22	1.99	0.44
1:AJ:53:PRO:HG2	1:AK:7:ASN:O	2.18	0.44
1:BD:33:ASP:OD1	1:BD:33:ASP:N	2.51	0.44
1:BG:2:TYR:N	1:DC:98:ASP:HB3	2.32	0.44
1:BG:33:ASP:OD1	1:BG:33:ASP:N	2.51	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BU:31:SER:HA	1:DG:33:ASP:OD1	2.17	0.44
1:BW:155:LEU:HG	1:BW:162:ILE:HD11	1.99	0.44
1:CB:33:ASP:OD1	1:CB:33:ASP:N	2.51	0.44
1:CI:53:PRO:HG2	1:CJ:7:ASN:O	2.18	0.44
1:CP:37:THR:HG23	1:CP:58:ASN:HD22	1.81	0.44
1:CQ:86:PHE:HE1	1:CQ:117:LEU:HD23	1.82	0.44
1:AC:155:LEU:HD13	1:BP:86:PHE:CD2	2.53	0.44
1:AD:53:PRO:HG2	1:AE:7:ASN:O	2.18	0.44
1:AF:86:PHE:CD2	1:BV:155:LEU:HD13	2.52	0.44
1:AN:151:LYS:HE3	1:AN:162:ILE:HG23	1.99	0.44
1:AR:96:LYS:HE2	1:CK:2:TYR:CD1	2.52	0.44
1:BP:86:PHE:HE1	1:BP:117:LEU:HD23	1.82	0.44
1:BS:130:ILE:O	1:BS:133:ILE:HG22	2.18	0.44
1:BZ:53:PRO:HG2	1:CA:7:ASN:O	2.17	0.44
1:CF:155:LEU:HG	1:CF:162:ILE:HD11	1.99	0.44
1:CK:33:ASP:OD1	1:CK:33:ASP:N	2.51	0.44
1:CK:130:ILE:O	1:CK:133:ILE:HG22	2.18	0.44
1:CQ:130:ILE:O	1:CQ:133:ILE:HG22	2.18	0.44
1:CY:86:PHE:HE1	1:CY:117:LEU:HD23	1.80	0.44
1:DB:151:LYS:HE3	1:DB:162:ILE:HG23	1.99	0.44
1:DF:33:ASP:OD1	1:DF:33:ASP:N	2.51	0.44
1:AC:86:PHE:HE1	1:AC:117:LEU:HD23	1.82	0.44
1:AG:155:LEU:HG	1:AG:162:ILE:HD11	1.99	0.44
1:AS:11:GLY:HA3	1:AT:21:GLU:HG2	2.00	0.44
1:BA:130:ILE:O	1:BA:133:ILE:HG22	2.18	0.44
1:BD:130:ILE:O	1:BD:133:ILE:HG22	2.18	0.44
1:BE:155:LEU:HG	1:BE:162:ILE:HD11	1.99	0.44
1:BY:155:LEU:HA	1:BY:155:LEU:HD23	1.77	0.44
1:CD:35:LYS:NZ	1:CK:34:ASN:HD21	2.15	0.44
1:CF:53:PRO:HG2	1:CG:7:ASN:O	2.17	0.44
1:CU:7:ASN:ND2	1:CV:46:ASP:H	2.12	0.44
1:DJ:53:PRO:HG2	1:DK:7:ASN:O	2.18	0.44
1:DK:37:THR:HG23	1:DK:58:ASN:HD22	1.81	0.44
1:AB:151:LYS:HE3	1:AB:162:ILE:HG23	1.99	0.44
1:AF:130:ILE:O	1:AF:133:ILE:HG22	2.18	0.44
1:AI:57:PHE:CE1	1:BY:162:ILE:HD12	2.53	0.44
1:AM:155:LEU:HG	1:AM:162:ILE:HD11	1.99	0.44
1:BB:53:PRO:HG2	1:BC:7:ASN:O	2.18	0.44
1:BG:34:ASN:CB	1:DC:161:ARG:HD2	2.42	0.44
1:BG:130:ILE:O	1:BG:133:ILE:HG22	2.18	0.44
1:BM:163:PHE:HB2	1:DI:40:TYR:CE1	2.53	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BX:37:THR:HG23	1:BX:58:ASN:HD22	1.81	0.44
1:BY:130:ILE:O	1:BY:133:ILE:HG22	2.18	0.44
1:BZ:155:LEU:HG	1:BZ:162:ILE:HD11	1.99	0.44
1:CL:33:ASP:HB2	1:CL:36:SER:H	1.81	0.44
1:CL:155:LEU:HG	1:CL:162:ILE:HD11	1.98	0.44
1:CR:53:PRO:HG2	1:CS:7:ASN:O	2.18	0.44
1:CR:82:VAL:HG13	1:DB:50:LEU:HD11	1.98	0.44
1:AI:130:ILE:O	1:AI:133:ILE:HG22	2.18	0.44
1:AV:155:LEU:HG	1:AV:162:ILE:HD11	1.99	0.44
1:AW:151:LYS:HE3	1:AW:162:ILE:HG23	1.99	0.44
1:BJ:6:ILE:HA	1:DF:94:TRP:CD1	2.53	0.44
1:BJ:33:ASP:OD1	1:BJ:33:ASP:N	2.51	0.44
1:BM:130:ILE:O	1:BM:133:ILE:HG22	2.18	0.44
1:BP:33:ASP:OD1	1:BP:33:ASP:N	2.51	0.44
1:CC:11:GLY:HA3	1:CD:21:GLU:HG2	2.00	0.44
1:CL:40:TYR:OH	1:CM:164:GLY:O	2.28	0.44
1:CM:37:THR:HG23	1:CM:58:ASN:HD22	1.81	0.44
1:CT:33:ASP:OD1	1:CT:33:ASP:N	2.51	0.44
1:CT:155:LEU:HA	1:CT:155:LEU:HD23	1.77	0.44
1:DD:11:GLY:HA3	1:DE:21:GLU:HG2	2.00	0.44
1:DF:130:ILE:O	1:DF:133:ILE:HG22	2.18	0.44
1:DI:130:ILE:O	1:DI:133:ILE:HG22	2.18	0.44
1:AC:130:ILE:O	1:AC:133:ILE:HG22	2.18	0.44
1:AL:6:ILE:HG22	1:CB:94:TRP:NE1	2.33	0.44
1:AR:161:ARG:HD2	1:CK:34:ASN:HB3	1.98	0.44
1:BG:78:ASP:HA	1:DC:107:TYR:CE1	2.53	0.44
1:BG:155:LEU:HD23	1:BG:155:LEU:HA	1.78	0.44
1:BG:163:PHE:HB2	1:DC:40:TYR:CZ	2.52	0.44
1:BN:53:PRO:HG2	1:BO:7:ASN:O	2.18	0.44
1:BS:86:PHE:HE1	1:BS:117:LEU:HD23	1.82	0.44
1:BW:11:GLY:HA3	1:BX:21:GLU:HG2	2.00	0.44
1:CO:11:GLY:HA3	1:CP:21:GLU:HG2	2.00	0.44
1:AL:40:TYR:CE1	1:CB:163:PHE:HB2	2.53	0.43
1:AQ:86:PHE:HE1	1:AQ:117:LEU:HD23	1.80	0.43
1:AR:18:VAL:HG22	1:CK:16:ASN:OD1	2.18	0.43
1:AT:151:LYS:HE3	1:AT:162:ILE:HG23	1.99	0.43
1:AU:24:ALA:HA	1:CN:163:PHE:CD2	2.53	0.43
1:BI:151:LYS:HE3	1:BI:162:ILE:HG23	1.99	0.43
1:BQ:53:PRO:HG2	1:BR:7:ASN:O	2.18	0.43
1:BT:11:GLY:HA3	1:BU:21:GLU:HG2	2.00	0.43
1:CA:35:LYS:NZ	1:CH:34:ASN:HD21	2.16	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AD:117:LEU:HD12	1:AE:112:SER:O	2.19	0.43
1:AI:33:ASP:OD1	1:AI:33:ASP:N	2.51	0.43
1:AW:37:THR:HG23	1:AW:58:ASN:HD22	1.81	0.43
1:BB:117:LEU:HD12	1:BC:112:SER:O	2.18	0.43
1:BE:53:PRO:HG2	1:BF:7:ASN:O	2.18	0.43
1:BP:155:LEU:HD23	1:BP:155:LEU:HA	1.77	0.43
1:BT:53:PRO:HG2	1:BU:7:ASN:O	2.17	0.43
1:BZ:11:GLY:HA3	1:CA:21:GLU:HG2	2.00	0.43
1:CX:53:PRO:HG2	1:CY:7:ASN:O	2.18	0.43
1:DA:155:LEU:HG	1:DA:162:ILE:HD11	1.99	0.43
1:DL:130:ILE:O	1:DL:133:ILE:HG22	2.18	0.43
1:AE:63:GLN:HG3	1:BG:50:LEU:HD21	1.99	0.43
1:AG:11:GLY:HA3	1:AH:21:GLU:HG2	2.00	0.43
1:AO:33:ASP:OD1	1:AO:33:ASP:N	2.51	0.43
1:AP:117:LEU:HD12	1:AQ:112:SER:O	2.19	0.43
1:AS:117:LEU:HD12	1:AT:112:SER:O	2.18	0.43
1:AU:163:PHE:HB2	1:CN:40:TYR:CE1	2.54	0.43
1:AX:162:ILE:HD12	1:CT:57:PHE:CE1	2.54	0.43
1:BG:2:TYR:CD1	1:DC:97:ALA:HA	2.53	0.43
1:BO:86:PHE:HE1	1:BO:117:LEU:HD23	1.80	0.43
1:BQ:11:GLY:HA3	1:BR:21:GLU:HG2	2.00	0.43
1:CO:133:ILE:HD12	1:CP:111:GLN:HB2	2.01	0.43
1:CQ:33:ASP:OD1	1:CQ:33:ASP:N	2.51	0.43
1:CU:11:GLY:HA3	1:CV:21:GLU:HG2	2.00	0.43
1:CY:151:LYS:HE3	1:CY:162:ILE:HG23	1.99	0.43
1:DA:11:GLY:HA3	1:DB:21:GLU:HG2	2.00	0.43
1:DA:53:PRO:HG2	1:DB:7:ASN:O	2.18	0.43
1:DF:86:PHE:HE1	1:DF:117:LEU:HD23	1.82	0.43
1:DI:33:ASP:OD1	1:DI:33:ASP:N	2.51	0.43
1:DI:86:PHE:HE1	1:DI:117:LEU:HD23	1.82	0.43
1:DJ:117:LEU:HD12	1:DK:112:SER:O	2.18	0.43
1:AJ:121:VAL:HG22	1:AJ:130:ILE:HD11	2.01	0.43
1:AN:155:LEU:HD11	1:AN:162:ILE:HD11	1.98	0.43
1:AO:130:ILE:O	1:AO:133:ILE:HG22	2.18	0.43
1:AQ:151:LYS:HE3	1:AQ:162:ILE:HG23	1.99	0.43
1:AX:130:ILE:O	1:AX:133:ILE:HG22	2.18	0.43
1:AY:155:LEU:HG	1:AY:162:ILE:HD11	1.99	0.43
1:BD:57:PHE:CE1	1:CZ:162:ILE:HD12	2.53	0.43
1:BD:164:GLY:O	1:CZ:40:TYR:OH	2.33	0.43
1:BQ:121:VAL:HG22	1:BQ:130:ILE:HD11	2.01	0.43
1:BZ:117:LEU:HD12	1:CA:112:SER:O	2.19	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CC:53:PRO:HG2	1:CD:7:ASN:O	2.17	0.43
1:CC:117:LEU:HD12	1:CD:112:SER:O	2.19	0.43
1:CE:130:ILE:O	1:CE:133:ILE:HG22	2.18	0.43
1:CL:53:PRO:HG2	1:CM:7:ASN:O	2.17	0.43
1:CL:133:ILE:HD12	1:CM:111:GLN:HB2	2.01	0.43
1:CO:117:LEU:HD12	1:CP:112:SER:O	2.18	0.43
1:CU:127:ALA:HB2	1:CV:153:GLN:OE1	2.19	0.43
1:CZ:33:ASP:OD1	1:CZ:33:ASP:N	2.51	0.43
1:DA:117:LEU:HD12	1:DB:112:SER:O	2.19	0.43
1:AA:121:VAL:HG22	1:AA:130:ILE:HD11	2.01	0.43
1:AM:53:PRO:HG2	1:AN:7:ASN:O	2.18	0.43
1:AS:53:PRO:HG2	1:AT:7:ASN:O	2.18	0.43
1:AS:121:VAL:HG22	1:AS:130:ILE:HD11	2.01	0.43
1:AV:117:LEU:HD12	1:AW:112:SER:O	2.19	0.43
1:AX:111:GLN:HB2	1:CT:133:ILE:HD12	2.00	0.43
1:AY:121:VAL:HG22	1:AY:130:ILE:HD11	2.01	0.43
1:AY:127:ALA:HB2	1:AZ:153:GLN:OE1	2.19	0.43
1:BJ:18:VAL:HG22	1:DF:16:ASN:OD1	2.18	0.43
1:BK:127:ALA:HB2	1:BL:153:GLN:OE1	2.19	0.43
1:BT:133:ILE:HD12	1:BU:111:GLN:HB2	2.01	0.43
1:BU:151:LYS:HE3	1:BU:162:ILE:HG23	1.99	0.43
1:CF:133:ILE:HD12	1:CG:111:GLN:HB2	2.01	0.43
1:CI:11:GLY:HA3	1:CJ:21:GLU:HG2	2.00	0.43
1:CL:127:ALA:HB2	1:CM:153:GLN:OE1	2.19	0.43
1:CO:127:ALA:HB2	1:CP:153:GLN:OE1	2.19	0.43
1:CW:130:ILE:O	1:CW:133:ILE:HG22	2.18	0.43
1:DC:130:ILE:O	1:DC:133:ILE:HG22	2.18	0.43
1:DD:7:ASN:ND2	1:DE:46:ASP:H	2.12	0.43
1:DG:11:GLY:HA3	1:DH:21:GLU:HG2	2.00	0.43
1:DJ:133:ILE:HD12	1:DK:111:GLN:HB2	2.01	0.43
1:AJ:127:ALA:HB2	1:AK:153:GLN:OE1	2.19	0.43
1:AL:7:ASN:HD21	1:CB:46:ASP:H	1.67	0.43
1:AS:127:ALA:HB2	1:AT:153:GLN:OE1	2.19	0.43
1:AV:121:VAL:HG22	1:AV:130:ILE:HD11	2.01	0.43
1:AV:133:ILE:HD12	1:AW:111:GLN:HB2	2.01	0.43
1:BA:86:PHE:CD2	1:CW:155:LEU:HD13	2.54	0.43
1:BD:119:SER:HB2	1:CZ:111:GLN:OE1	2.19	0.43
1:BH:11:GLY:HA3	1:BI:21:GLU:HG2	2.00	0.43
1:BH:133:ILE:HD12	1:BI:111:GLN:HB2	2.01	0.43
1:BK:53:PRO:HG2	1:BL:7:ASN:O	2.18	0.43
1:BV:86:PHE:HE1	1:BV:117:LEU:HD23	1.82	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CF:127:ALA:HB2	1:CG:153:GLN:OE1	2.19	0.43
1:CH:33:ASP:OD1	1:CH:33:ASP:N	2.51	0.43
1:CI:127:ALA:HB2	1:CJ:153:GLN:OE1	2.19	0.43
1:CI:133:ILE:HD12	1:CJ:111:GLN:HB2	2.01	0.43
1:CN:86:PHE:HE1	1:CN:117:LEU:HD23	1.82	0.43
1:CR:11:GLY:HA3	1:CS:21:GLU:HG2	2.00	0.43
1:CU:121:VAL:HG22	1:CU:130:ILE:HD11	2.01	0.43
1:DD:127:ALA:HB2	1:DE:153:GLN:OE1	2.19	0.43
1:DD:133:ILE:HD12	1:DE:111:GLN:HB2	2.01	0.43
1:DJ:127:ALA:HB2	1:DK:153:GLN:OE1	2.19	0.43
1:AA:117:LEU:HD12	1:AB:112:SER:O	2.19	0.43
1:AL:130:ILE:O	1:AL:133:ILE:HG22	2.18	0.43
1:BD:86:PHE:HE1	1:BD:117:LEU:HD23	1.82	0.43
1:BE:117:LEU:HD12	1:BF:112:SER:O	2.19	0.43
1:BE:121:VAL:HG22	1:BE:130:ILE:HD11	2.01	0.43
1:BE:127:ALA:HB2	1:BF:153:GLN:OE1	2.19	0.43
1:BE:133:ILE:HD12	1:BF:111:GLN:HB2	2.01	0.43
1:BJ:130:ILE:O	1:BJ:133:ILE:HG22	2.18	0.43
1:BP:130:ILE:O	1:BP:133:ILE:HG22	2.18	0.43
1:BQ:117:LEU:HD12	1:BR:112:SER:O	2.19	0.43
1:BX:155:LEU:HD23	1:BX:155:LEU:HA	1.88	0.43
1:CC:127:ALA:HB2	1:CD:153:GLN:OE1	2.19	0.43
1:CI:117:LEU:HD12	1:CJ:112:SER:O	2.19	0.43
1:CR:133:ILE:HD12	1:CS:111:GLN:HB2	2.01	0.43
1:CT:86:PHE:HE1	1:CT:117:LEU:HD23	1.82	0.43
1:CU:117:LEU:HD12	1:CV:112:SER:O	2.19	0.43
1:CX:11:GLY:HA3	1:CY:21:GLU:HG2	2.00	0.43
1:CX:121:VAL:HG22	1:CX:130:ILE:HD11	2.01	0.43
1:DG:53:PRO:HG2	1:DH:7:ASN:O	2.18	0.43
1:DK:151:LYS:HE3	1:DK:162:ILE:HG23	1.99	0.43
1:DK:155:LEU:HD23	1:DK:155:LEU:HA	1.88	0.43
1:AA:127:ALA:HB2	1:AB:153:GLN:OE1	2.19	0.43
1:AD:11:GLY:HA3	1:AE:21:GLU:HG2	2.00	0.43
1:AD:133:ILE:HD12	1:AE:111:GLN:HB2	2.00	0.43
1:AJ:11:GLY:HA3	1:AK:21:GLU:HG2	2.00	0.43
1:AJ:63:GLN:HG3	1:AJ:64:TYR:CD1	2.54	0.43
1:AP:53:PRO:HG2	1:AQ:7:ASN:O	2.18	0.43
1:AV:11:GLY:HA3	1:AW:21:GLU:HG2	2.00	0.43
1:BK:133:ILE:HD12	1:BL:111:GLN:HB2	2.01	0.43
1:BQ:127:ALA:HB2	1:BR:153:GLN:OE1	2.19	0.43
1:CE:33:ASP:OD1	1:CE:33:ASP:N	2.51	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CL:11:GLY:HA3	1:CM:21:GLU:HG2	2.00	0.43
1:CU:133:ILE:HD12	1:CV:111:GLN:HB2	2.01	0.43
1:DG:127:ALA:HB2	1:DH:153:GLN:OE1	2.19	0.43
1:DJ:121:VAL:HG22	1:DJ:130:ILE:HD11	2.01	0.43
1:AF:86:PHE:HE1	1:AF:117:LEU:HD23	1.82	0.43
1:AG:63:GLN:HG3	1:AG:64:TYR:CD1	2.54	0.43
1:AG:133:ILE:HD12	1:AH:111:GLN:HB2	2.01	0.43
1:AH:151:LYS:HE3	1:AH:162:ILE:HG23	1.99	0.43
1:AM:121:VAL:HG22	1:AM:130:ILE:HD11	2.01	0.43
1:AR:153:GLN:OE1	1:CK:127:ALA:HB2	2.18	0.43
1:BC:30:VAL:HA	1:BN:34:ASN:HD22	1.82	0.43
1:BH:117:LEU:HD12	1:BI:112:SER:O	2.19	0.43
1:BH:127:ALA:HB2	1:BI:153:GLN:OE1	2.19	0.43
1:BK:63:GLN:HG3	1:BK:64:TYR:CD1	2.54	0.43
1:BT:121:VAL:HG22	1:BT:130:ILE:HD11	2.01	0.43
1:BW:127:ALA:HB2	1:BX:153:GLN:OE1	2.19	0.43
1:BZ:63:GLN:HG3	1:BZ:64:TYR:CD1	2.54	0.43
1:CA:30:VAL:HA	1:CF:34:ASN:HD22	1.81	0.43
1:CT:130:ILE:O	1:CT:133:ILE:HG22	2.18	0.43
1:CW:34:ASN:HD21	1:CY:35:LYS:NZ	2.16	0.43
1:CZ:130:ILE:O	1:CZ:133:ILE:HG22	2.18	0.43
1:DB:86:PHE:HE1	1:DB:117:LEU:HD23	1.80	0.43
1:DD:117:LEU:HD12	1:DE:112:SER:O	2.18	0.43
1:DG:63:GLN:HG3	1:DG:64:TYR:CD1	2.54	0.43
1:AC:34:ASN:HD21	1:DH:35:LYS:NZ	2.17	0.43
1:AG:34:ASN:HD22	1:CS:30:VAL:HA	1.83	0.43
1:AJ:117:LEU:HD12	1:AK:112:SER:O	2.19	0.43
1:AP:11:GLY:HA3	1:AQ:21:GLU:HG2	2.00	0.43
1:AS:63:GLN:HG3	1:AS:64:TYR:CD1	2.54	0.43
1:AY:133:ILE:HD12	1:AZ:111:GLN:HB2	2.01	0.43
1:BB:63:GLN:HG3	1:BB:64:TYR:CD1	2.54	0.43
1:BE:11:GLY:HA3	1:BF:21:GLU:HG2	2.00	0.43
1:BK:11:GLY:HA3	1:BL:21:GLU:HG2	2.00	0.43
1:BW:133:ILE:HD12	1:BX:111:GLN:HB2	2.01	0.43
1:BZ:121:VAL:HG22	1:BZ:130:ILE:HD11	2.01	0.43
1:BZ:127:ALA:HB2	1:CA:153:GLN:OE1	2.19	0.43
1:BZ:133:ILE:HD12	1:CA:111:GLN:HB2	2.01	0.43
1:CC:121:VAL:HG22	1:CC:130:ILE:HD11	2.01	0.43
1:CH:130:ILE:O	1:CH:133:ILE:HG22	2.18	0.43
1:CH:155:LEU:HD21	1:CH:162:ILE:HD11	2.01	0.43
1:CR:127:ALA:HB2	1:CS:153:GLN:OE1	2.19	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AF:113:CYS:HB2	1:BV:117:LEU:HD13	2.00	0.42
1:AO:46:ASP:OD2	1:AO:48:SER:OG	2.27	0.42
1:AP:63:GLN:HG3	1:AP:64:TYR:CD1	2.54	0.42
1:AP:127:ALA:HB2	1:AQ:153:GLN:OE1	2.19	0.42
1:AV:127:ALA:HB2	1:AW:153:GLN:OE1	2.19	0.42
1:AY:57:PHE:CE1	1:AZ:162:ILE:HD12	2.54	0.42
1:BK:117:LEU:HD12	1:BL:112:SER:O	2.19	0.42
1:BN:11:GLY:HA3	1:BO:21:GLU:HG2	2.00	0.42
1:BN:127:ALA:HB2	1:BO:153:GLN:OE1	2.19	0.42
1:BQ:133:ILE:HD12	1:BR:111:GLN:HB2	2.01	0.42
1:BW:63:GLN:HG3	1:BW:64:TYR:CD1	2.54	0.42
1:BY:155:LEU:HD21	1:BY:162:ILE:HD11	2.01	0.42
1:CN:155:LEU:HD21	1:CN:162:ILE:HD11	2.01	0.42
1:CO:63:GLN:HG3	1:CO:64:TYR:CD1	2.54	0.42
1:CR:63:GLN:HG3	1:CR:64:TYR:CD1	2.54	0.42
1:CX:63:GLN:HG3	1:CX:64:TYR:CD1	2.54	0.42
1:CX:117:LEU:HD12	1:CY:112:SER:O	2.19	0.42
1:DA:127:ALA:HB2	1:DB:153:GLN:OE1	2.19	0.42
1:AC:86:PHE:CD2	1:BP:155:LEU:HD13	2.54	0.42
1:AI:155:LEU:HD13	1:BY:86:PHE:CD2	2.54	0.42
1:AR:155:LEU:HA	1:AR:155:LEU:HD23	1.77	0.42
1:AV:63:GLN:HG3	1:AV:64:TYR:CD1	2.54	0.42
1:BB:133:ILE:HD12	1:BC:111:GLN:HB2	2.01	0.42
1:BD:155:LEU:HD21	1:BD:162:ILE:HD11	2.01	0.42
1:BH:57:PHE:CE1	1:BI:162:ILE:HD12	2.55	0.42
1:BQ:63:GLN:HG3	1:BQ:64:TYR:CD1	2.54	0.42
1:BV:46:ASP:OD2	1:BV:48:SER:OG	2.27	0.42
1:CB:130:ILE:O	1:CB:133:ILE:HG22	2.18	0.42
1:CF:63:GLN:HG3	1:CF:64:TYR:CD1	2.54	0.42
1:CS:155:LEU:HD23	1:CS:155:LEU:HA	1.88	0.42
1:DA:133:ILE:HD12	1:DB:111:GLN:HB2	2.00	0.42
1:DG:57:PHE:CE1	1:DH:162:ILE:HD12	2.55	0.42
1:DG:117:LEU:HD12	1:DH:112:SER:O	2.18	0.42
1:DJ:11:GLY:HA3	1:DK:21:GLU:HG2	2.00	0.42
1:AC:155:LEU:HD23	1:AC:155:LEU:HA	1.77	0.42
1:AD:57:PHE:CE1	1:AE:162:ILE:HD12	2.54	0.42
1:AD:63:GLN:HG3	1:AD:64:TYR:CD1	2.54	0.42
1:AI:155:LEU:HD23	1:AI:155:LEU:HA	1.77	0.42
1:AM:11:GLY:HA3	1:AN:21:GLU:HG2	2.00	0.42
1:AM:63:GLN:HG3	1:AM:64:TYR:CD1	2.54	0.42
1:AM:133:ILE:HD12	1:AN:111:GLN:HB2	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AP:57:PHE:CE1	1:AQ:162:ILE:HD12	2.55	0.42
1:AR:130:ILE:O	1:AR:133:ILE:HG22	2.18	0.42
1:BN:117:LEU:HD12	1:BO:112:SER:O	2.19	0.42
1:BT:57:PHE:CE1	1:BU:162:ILE:HD12	2.55	0.42
1:BT:117:LEU:HD12	1:BU:112:SER:O	2.19	0.42
1:CA:50:LEU:HD11	1:CF:82:VAL:HG13	2.00	0.42
1:CI:57:PHE:CE1	1:CJ:162:ILE:HD12	2.55	0.42
1:CL:63:GLN:HG3	1:CL:64:TYR:CD1	2.54	0.42
1:CN:130:ILE:O	1:CN:133:ILE:HG22	2.18	0.42
1:CO:121:VAL:HG22	1:CO:130:ILE:HD11	2.01	0.42
1:DA:57:PHE:CE1	1:DB:162:ILE:HD12	2.55	0.42
1:DA:63:GLN:HG3	1:DA:64:TYR:CD1	2.54	0.42
1:DG:121:VAL:HG22	1:DG:130:ILE:HD11	2.01	0.42
1:DH:86:PHE:HE1	1:DH:117:LEU:HD23	1.80	0.42
1:DJ:63:GLN:HG3	1:DJ:64:TYR:CD1	2.54	0.42
1:DL:155:LEU:HD21	1:DL:162:ILE:HD11	2.01	0.42
1:AA:57:PHE:CE1	1:AB:162:ILE:HD12	2.55	0.42
1:AD:34:ASN:HD22	1:DE:30:VAL:HA	1.84	0.42
1:AG:127:ALA:HB2	1:AH:153:GLN:OE1	2.19	0.42
1:AO:155:LEU:HD23	1:AO:155:LEU:HA	1.78	0.42
1:AR:98:ASP:HB3	1:CK:2:TYR:N	2.35	0.42
1:AR:162:ILE:HD12	1:CK:57:PHE:CE1	2.55	0.42
1:AS:133:ILE:HD12	1:AT:111:GLN:HB2	2.01	0.42
1:AX:155:LEU:HD21	1:AX:162:ILE:HD11	2.01	0.42
1:AY:63:GLN:HG3	1:AY:64:TYR:CD1	2.54	0.42
1:BE:63:GLN:HG3	1:BE:64:TYR:CD1	2.54	0.42
1:BH:121:VAL:HG22	1:BH:130:ILE:HD11	2.01	0.42
1:BJ:40:TYR:CE1	1:DF:163:PHE:HB2	2.54	0.42
1:BM:155:LEU:HD21	1:BM:162:ILE:HD11	2.01	0.42
1:BT:63:GLN:HG3	1:BT:64:TYR:CD1	2.54	0.42
1:BW:121:VAL:HG22	1:BW:130:ILE:HD11	2.01	0.42
1:CI:63:GLN:HG3	1:CI:64:TYR:CD1	2.54	0.42
1:CL:117:LEU:HD12	1:CM:112:SER:O	2.19	0.42
1:CU:57:PHE:CE1	1:CV:162:ILE:HD12	2.55	0.42
1:CX:127:ALA:HB2	1:CY:153:GLN:OE1	2.19	0.42
1:CZ:155:LEU:HD21	1:CZ:162:ILE:HD11	2.01	0.42
1:DI:155:LEU:HD21	1:DI:162:ILE:HD11	2.01	0.42
1:AA:63:GLN:HG3	1:AA:64:TYR:CD1	2.54	0.42
1:AA:133:ILE:HD12	1:AB:111:GLN:HB2	2.01	0.42
1:AC:155:LEU:HD21	1:AC:162:ILE:HD11	2.01	0.42
1:AI:155:LEU:HD21	1:AI:162:ILE:HD11	2.01	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AL:162:ILE:HD12	1:CB:57:PHE:CE1	2.54	0.42
1:AM:117:LEU:HD12	1:AN:112:SER:O	2.19	0.42
1:AO:57:PHE:CE1	1:CE:162:ILE:HD12	2.55	0.42
1:AR:33:ASP:OD1	1:AR:33:ASP:N	2.51	0.42
1:AU:7:ASN:HD21	1:CN:46:ASP:H	1.67	0.42
1:AU:17:LEU:HD22	1:CN:19:LEU:HD11	2.01	0.42
1:AU:130:ILE:O	1:AU:133:ILE:HG22	2.18	0.42
1:AY:117:LEU:HD12	1:AZ:112:SER:O	2.19	0.42
1:BH:63:GLN:HG3	1:BH:64:TYR:CD1	2.54	0.42
1:BN:133:ILE:HD12	1:BO:111:GLN:HB2	2.01	0.42
1:BT:127:ALA:HB2	1:BU:153:GLN:OE1	2.19	0.42
1:CF:117:LEU:HD12	1:CG:112:SER:O	2.19	0.42
1:CF:155:LEU:HD23	1:CF:155:LEU:HA	1.92	0.42
1:DD:40:TYR:OH	1:DE:164:GLY:O	2.28	0.42
1:DD:121:VAL:HG22	1:DD:130:ILE:HD11	2.01	0.42
1:DD:155:LEU:HD23	1:DD:155:LEU:HA	1.92	0.42
1:AI:119:SER:HB2	1:BY:111:GLN:OE1	2.20	0.42
1:AJ:133:ILE:HD12	1:AK:111:GLN:HB2	2.01	0.42
1:AR:57:PHE:HE2	1:CK:163:PHE:CE1	2.37	0.42
1:AR:155:LEU:HD21	1:AR:162:ILE:HD11	2.01	0.42
1:AX:33:ASP:OD1	1:AX:33:ASP:N	2.51	0.42
1:BO:155:LEU:HD23	1:BO:155:LEU:HA	1.88	0.42
1:BQ:57:PHE:CE1	1:BR:162:ILE:HD12	2.55	0.42
1:BS:155:LEU:HD21	1:BS:162:ILE:HD11	2.01	0.42
1:BU:86:PHE:HE1	1:BU:117:LEU:HD23	1.80	0.42
1:BV:130:ILE:O	1:BV:133:ILE:HG22	2.18	0.42
1:CC:133:ILE:HD12	1:CD:111:GLN:HB2	2.01	0.42
1:DD:63:GLN:HG3	1:DD:64:TYR:CD1	2.54	0.42
1:AA:11:GLY:HA3	1:AB:21:GLU:HG2	2.00	0.42
1:AL:119:SER:HB2	1:CB:111:GLN:OE1	2.19	0.42
1:AP:133:ILE:HD12	1:AQ:111:GLN:HB2	2.01	0.42
1:AU:155:LEU:HD23	1:AU:155:LEU:HA	1.78	0.42
1:BB:127:ALA:HB2	1:BC:153:GLN:OE1	2.19	0.42
1:BE:57:PHE:CE1	1:BF:162:ILE:HD12	2.55	0.42
1:BM:57:PHE:CE1	1:DI:162:ILE:HD12	2.55	0.42
1:BN:121:VAL:HG22	1:BN:130:ILE:HD11	2.01	0.42
1:BW:57:PHE:CE1	1:BX:162:ILE:HD12	2.55	0.42
1:CC:57:PHE:CE1	1:CD:162:ILE:HD12	2.55	0.42
1:CC:63:GLN:HG3	1:CC:64:TYR:CD1	2.54	0.42
1:CF:57:PHE:CE1	1:CG:162:ILE:HD12	2.55	0.42
1:CU:63:GLN:HG3	1:CU:64:TYR:CD1	2.54	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CY:155:LEU:HD23	1:CY:155:LEU:HA	1.88	0.42
1:DC:155:LEU:HD21	1:DC:162:ILE:HD11	2.01	0.42
1:DD:57:PHE:CE1	1:DE:162:ILE:HD12	2.55	0.42
1:DJ:57:PHE:CE1	1:DK:162:ILE:HD12	2.55	0.42
1:AD:127:ALA:HB2	1:AE:153:GLN:OE1	2.19	0.42
1:AF:40:TYR:CZ	1:BV:163:PHE:HB2	2.55	0.42
1:AG:117:LEU:HD12	1:AH:112:SER:O	2.19	0.42
1:AJ:57:PHE:CE1	1:AK:162:ILE:HD12	2.54	0.42
1:AK:30:VAL:HA	1:BW:34:ASN:HD22	1.83	0.42
1:AS:155:LEU:HD13	1:AT:86:PHE:CD2	2.55	0.42
1:BB:121:VAL:HG22	1:BB:130:ILE:HD11	2.01	0.42
1:BD:24:ALA:HA	1:CZ:163:PHE:CD2	2.52	0.42
1:BD:155:LEU:HD23	1:BD:155:LEU:HA	1.77	0.42
1:BK:121:VAL:HG22	1:BK:130:ILE:HD11	2.01	0.42
1:BZ:57:PHE:CE1	1:CA:162:ILE:HD12	2.55	0.42
1:CL:57:PHE:CE1	1:CM:162:ILE:HD12	2.55	0.42
1:CV:58:ASN:OD1	1:CV:87:TYR:HB3	2.20	0.42
1:CX:57:PHE:CE1	1:CY:162:ILE:HD12	2.55	0.42
1:DA:121:VAL:HG22	1:DA:130:ILE:HD11	2.01	0.42
1:AD:121:VAL:HG22	1:AD:130:ILE:HD11	2.01	0.42
1:AE:58:ASN:OD1	1:AE:87:TYR:HB3	2.20	0.42
1:AF:33:ASP:OD1	1:AF:33:ASP:N	2.51	0.42
1:AG:121:VAL:HG22	1:AG:130:ILE:HD11	2.01	0.42
1:AJ:34:ASN:HD22	1:CV:30:VAL:HA	1.84	0.42
1:AL:1:SER:C	1:CB:98:ASP:HB3	2.45	0.42
1:AL:94:TRP:NE1	1:CB:6:ILE:HG22	2.35	0.42
1:AT:58:ASN:OD1	1:AT:87:TYR:HB3	2.20	0.42
1:AW:58:ASN:OD1	1:AW:87:TYR:HB3	2.20	0.42
1:AY:11:GLY:HA3	1:AZ:21:GLU:HG2	2.00	0.42
1:BB:57:PHE:CE1	1:BC:162:ILE:HD12	2.55	0.42
1:BD:162:ILE:HD12	1:CZ:57:PHE:CE1	2.55	0.42
1:BE:155:LEU:HD13	1:BF:86:PHE:CD2	2.55	0.42
1:BH:155:LEU:HD13	1:BI:86:PHE:CD2	2.55	0.42
1:BK:57:PHE:CE1	1:BL:162:ILE:HD12	2.55	0.42
1:BP:155:LEU:HD21	1:BP:162:ILE:HD11	2.01	0.42
1:BT:155:LEU:HD13	1:BU:86:PHE:CD2	2.55	0.42
1:CF:11:GLY:HA3	1:CG:21:GLU:HG2	2.00	0.42
1:AB:155:LEU:HD23	1:AB:155:LEU:HA	1.88	0.42
1:AC:24:ALA:HA	1:BP:163:PHE:CD2	2.53	0.42
1:AG:155:LEU:HD13	1:AH:86:PHE:CD2	2.55	0.42
1:AK:45:SER:HB3	1:AK:53:PRO:HD2	2.02	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AM:127:ALA:HB2	1:AN:153:GLN:OE1	2.19	0.42
1:AU:46:ASP:OD2	1:AU:48:SER:OG	2.27	0.42
1:AV:57:PHE:CE1	1:AW:162:ILE:HD12	2.55	0.42
1:BH:82:VAL:HG11	1:BL:50:LEU:HD21	2.01	0.42
1:CC:155:LEU:HD13	1:CD:86:PHE:CD2	2.55	0.42
1:CJ:47:ASN:OD1	1:DB:157:ASP:CG	2.62	0.42
1:CL:121:VAL:HG22	1:CL:130:ILE:HD11	2.01	0.42
1:CP:155:LEU:HD23	1:CP:155:LEU:HA	1.88	0.42
1:CQ:155:LEU:HD21	1:CQ:162:ILE:HD11	2.01	0.42
1:CR:57:PHE:CE1	1:CS:162:ILE:HD12	2.55	0.42
1:CR:121:VAL:HG22	1:CR:130:ILE:HD11	2.01	0.42
1:DA:155:LEU:HD13	1:DB:86:PHE:CD2	2.55	0.42
1:DI:155:LEU:HA	1:DI:155:LEU:HD23	1.77	0.42
1:AJ:155:LEU:HD13	1:AK:86:PHE:CD2	2.55	0.41
1:AL:155:LEU:HD21	1:AL:162:ILE:HD11	2.01	0.41
1:AO:155:LEU:HD21	1:AO:162:ILE:HD11	2.01	0.41
1:AO:162:ILE:HD12	1:CE:57:PHE:CE1	2.55	0.41
1:AW:45:SER:HB3	1:AW:53:PRO:HD2	2.02	0.41
1:BC:45:SER:HB3	1:BC:53:PRO:HD2	2.02	0.41
1:BG:19:LEU:HD11	1:DC:17:LEU:HD22	2.02	0.41
1:BN:63:GLN:HG3	1:BN:64:TYR:CD1	2.54	0.41
1:BV:155:LEU:HD21	1:BV:162:ILE:HD11	2.01	0.41
1:BW:117:LEU:HD12	1:BX:112:SER:O	2.19	0.41
1:BW:155:LEU:HD13	1:BX:86:PHE:CD2	2.55	0.41
1:CA:155:LEU:HD23	1:CA:155:LEU:HA	1.88	0.41
1:CO:155:LEU:HD13	1:CP:86:PHE:CD2	2.55	0.41
1:CR:117:LEU:HD12	1:CS:112:SER:O	2.19	0.41
1:CW:155:LEU:HD21	1:CW:162:ILE:HD11	2.01	0.41
1:CX:133:ILE:HD12	1:CY:111:GLN:HB2	2.01	0.41
1:DD:155:LEU:HD13	1:DE:86:PHE:CD2	2.55	0.41
1:DE:58:ASN:OD1	1:DE:87:TYR:HB3	2.20	0.41
1:DG:155:LEU:HD13	1:DH:86:PHE:CD2	2.55	0.41
1:AH:155:LEU:HD23	1:AH:155:LEU:HA	1.88	0.41
1:AM:155:LEU:HD13	1:AN:86:PHE:CD2	2.55	0.41
1:AO:155:LEU:HD13	1:CE:86:PHE:CD2	2.55	0.41
1:AP:121:VAL:HG22	1:AP:130:ILE:HD11	2.01	0.41
1:AP:155:LEU:HD13	1:AQ:86:PHE:CD2	2.55	0.41
1:AQ:58:ASN:OD1	1:AQ:87:TYR:HB3	2.20	0.41
1:AR:40:TYR:CD2	1:CK:163:PHE:HD1	2.38	0.41
1:AU:119:SER:HB2	1:CN:111:GLN:OE1	2.19	0.41
1:AU:155:LEU:HD21	1:AU:162:ILE:HD11	2.01	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AV:155:LEU:HD13	1:AW:86:PHE:CD2	2.55	0.41
1:BB:11:GLY:HA3	1:BC:21:GLU:HG2	2.00	0.41
1:BZ:155:LEU:HD13	1:CA:86:PHE:CD2	2.55	0.41
1:CG:58:ASN:OD1	1:CG:87:TYR:HB3	2.20	0.41
1:CI:40:TYR:OH	1:CJ:164:GLY:O	2.28	0.41
1:CI:121:VAL:HG22	1:CI:130:ILE:HD11	2.01	0.41
1:CK:155:LEU:HD21	1:CK:162:ILE:HD11	2.01	0.41
1:CS:58:ASN:OD1	1:CS:87:TYR:HB3	2.20	0.41
1:CT:155:LEU:HD21	1:CT:162:ILE:HD11	2.01	0.41
1:DH:58:ASN:OD1	1:DH:87:TYR:HB3	2.20	0.41
1:DJ:155:LEU:HD13	1:DK:86:PHE:CD2	2.55	0.41
1:AA:82:VAL:HG13	1:DH:50:LEU:HD11	2.03	0.41
1:AF:98:ASP:HB3	1:BV:1:SER:C	2.45	0.41
1:AF:155:LEU:HD21	1:AF:162:ILE:HD11	2.01	0.41
1:AG:155:LEU:HD23	1:AG:155:LEU:HA	1.92	0.41
1:AQ:45:SER:HB3	1:AQ:53:PRO:HD2	2.02	0.41
1:AT:45:SER:HB3	1:AT:53:PRO:HD2	2.02	0.41
1:AU:33:ASP:OD1	1:AU:33:ASP:N	2.51	0.41
1:BD:40:TYR:CE1	1:CZ:163:PHE:HB2	2.55	0.41
1:BF:58:ASN:OD1	1:BF:87:TYR:HB3	2.20	0.41
1:BG:155:LEU:HD21	1:BG:162:ILE:HD11	2.01	0.41
1:BJ:155:LEU:HD21	1:BJ:162:ILE:HD11	2.01	0.41
1:BO:45:SER:HB3	1:BO:53:PRO:HD2	2.02	0.41
1:BQ:155:LEU:HD13	1:BR:86:PHE:CD2	2.55	0.41
1:CE:22:ILE:HD13	1:CE:22:ILE:HA	1.92	0.41
1:CE:155:LEU:HA	1:CE:155:LEU:HD23	1.77	0.41
1:CE:155:LEU:HD21	1:CE:162:ILE:HD11	2.01	0.41
1:CQ:46:ASP:OD2	1:CQ:48:SER:OG	2.27	0.41
1:CU:155:LEU:HD13	1:CV:86:PHE:CD2	2.55	0.41
1:CZ:46:ASP:OD2	1:CZ:48:SER:OG	2.27	0.41
1:DB:58:ASN:OD1	1:DB:87:TYR:HB3	2.20	0.41
1:DG:133:ILE:HD12	1:DH:111:GLN:HB2	2.01	0.41
1:DK:45:SER:HB3	1:DK:53:PRO:HD2	2.02	0.41
1:AA:155:LEU:HD13	1:AB:86:PHE:CD2	2.55	0.41
1:AG:57:PHE:CE1	1:AH:162:ILE:HD12	2.55	0.41
1:AM:57:PHE:CE1	1:AN:162:ILE:HD12	2.55	0.41
1:BA:163:PHE:CD2	1:CW:24:ALA:HA	2.52	0.41
1:BD:163:PHE:HB2	1:CZ:40:TYR:CE1	2.54	0.41
1:BD:163:PHE:CD2	1:CZ:24:ALA:HA	2.53	0.41
1:BM:163:PHE:CD2	1:DI:24:ALA:HA	2.56	0.41
1:BN:155:LEU:HD13	1:BO:86:PHE:CD2	2.55	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BR:50:LEU:HD23	1:BR:50:LEU:HA	1.89	0.41
1:CA:60:VAL:HB	1:CA:85:GLU:HG3	2.03	0.41
1:CD:45:SER:HB3	1:CD:53:PRO:HD2	2.03	0.41
1:CM:58:ASN:OD1	1:CM:87:TYR:HB3	2.20	0.41
1:CO:57:PHE:CE1	1:CP:162:ILE:HD12	2.55	0.41
1:DB:45:SER:HB3	1:DB:53:PRO:HD2	2.02	0.41
1:DF:155:LEU:HD21	1:DF:162:ILE:HD11	2.01	0.41
1:AB:58:ASN:OD1	1:AB:87:TYR:HB3	2.20	0.41
1:AH:45:SER:HB3	1:AH:53:PRO:HD2	2.02	0.41
1:AH:58:ASN:OD1	1:AH:87:TYR:HB3	2.20	0.41
1:AI:94:TRP:NE1	1:BY:6:ILE:HG22	2.36	0.41
1:AK:60:VAL:HB	1:AK:85:GLU:HG3	2.03	0.41
1:AL:24:ALA:HA	1:CB:163:PHE:CD2	2.53	0.41
1:AR:98:ASP:HB3	1:CK:1:SER:O	2.20	0.41
1:AZ:58:ASN:OD1	1:AZ:87:TYR:HB3	2.20	0.41
1:BA:155:LEU:HD21	1:BA:162:ILE:HD11	2.01	0.41
1:BI:155:LEU:HD23	1:BI:155:LEU:HA	1.88	0.41
1:CM:45:SER:HB3	1:CM:53:PRO:HD2	2.02	0.41
1:AD:155:LEU:HD13	1:AE:86:PHE:CD2	2.55	0.41
1:AE:45:SER:HB3	1:AE:53:PRO:HD2	2.02	0.41
1:AI:6:ILE:HG22	1:BY:94:TRP:NE1	2.34	0.41
1:AU:19:LEU:HD11	1:CN:17:LEU:HD22	2.03	0.41
1:AY:155:LEU:HD13	1:AZ:86:PHE:CD2	2.55	0.41
1:BF:45:SER:HB3	1:BF:53:PRO:HD2	2.02	0.41
1:BI:45:SER:HB3	1:BI:53:PRO:HD2	2.03	0.41
1:BM:46:ASP:OD2	1:BM:48:SER:OG	2.27	0.41
1:CB:155:LEU:HD21	1:CB:162:ILE:HD11	2.01	0.41
1:CF:155:LEU:HD13	1:CG:86:PHE:CD2	2.55	0.41
1:AA:155:LEU:HD23	1:AA:155:LEU:HA	1.92	0.41
1:AK:58:ASN:OD1	1:AK:87:TYR:HB3	2.20	0.41
1:AO:19:LEU:HD11	1:CE:17:LEU:HD22	2.02	0.41
1:AT:60:VAL:HB	1:AT:85:GLU:HG3	2.03	0.41
1:AX:96:LYS:HE2	1:CT:2:TYR:CD1	2.56	0.41
1:BA:164:GLY:O	1:CW:40:TYR:OH	2.31	0.41
1:BR:58:ASN:OD1	1:BR:87:TYR:HB3	2.20	0.41
1:BU:58:ASN:OD1	1:BU:87:TYR:HB3	2.20	0.41
1:BX:58:ASN:OD1	1:BX:87:TYR:HB3	2.20	0.41
1:CF:121:VAL:HG22	1:CF:130:ILE:HD11	2.01	0.41
1:CJ:58:ASN:OD1	1:CJ:87:TYR:HB3	2.20	0.41
1:CN:155:LEU:HD23	1:CN:155:LEU:HA	1.77	0.41
1:AI:163:PHE:HB2	1:BY:40:TYR:CE1	2.55	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:AL:155:LEU:HD13	1:CB:86:PHE:CD2	2.55	0.41
1:AN:60:VAL:HB	1:AN:85:GLU:HG3	2.03	0.41
1:AQ:60:VAL:HB	1:AQ:85:GLU:HG3	2.03	0.41
1:BK:155:LEU:HD13	1:BL:86:PHE:CD2	2.55	0.41
1:BL:45:SER:HB3	1:BL:53:PRO:HD2	2.02	0.41
1:BM:1:SER:C	1:DI:98:ASP:HB3	2.46	0.41
1:BX:30:VAL:HA	1:CL:34:ASN:HD22	1.85	0.41
1:CA:58:ASN:OD1	1:CA:87:TYR:HB3	2.20	0.41
1:CR:155:LEU:HD13	1:CS:86:PHE:CD2	2.55	0.41
1:DJ:155:LEU:HD23	1:DJ:155:LEU:HA	1.92	0.41
1:AB:60:VAL:HB	1:AB:85:GLU:HG3	2.03	0.41
1:AE:35:LYS:HZ3	1:BG:34:ASN:HD21	1.66	0.41
1:AG:94:TRP:NE1	1:AH:6:ILE:HG22	2.36	0.41
1:AI:1:SER:C	1:BY:98:ASP:HB3	2.46	0.41
1:AK:50:LEU:HD11	1:BW:82:VAL:HG13	2.03	0.41
1:AK:155:LEU:HA	1:AK:155:LEU:HD23	1.88	0.41
1:AL:155:LEU:HA	1:AL:155:LEU:HD23	1.77	0.41
1:AN:58:ASN:OD1	1:AN:87:TYR:HB3	2.20	0.41
1:AS:94:TRP:NE1	1:AT:6:ILE:HG22	2.36	0.41
1:AV:94:TRP:NE1	1:AW:6:ILE:HG22	2.36	0.41
1:BB:80:LEU:HA	1:BB:80:LEU:HD23	1.87	0.41
1:BB:94:TRP:NE1	1:BC:6:ILE:HG22	2.36	0.41
1:BB:155:LEU:HD13	1:BC:86:PHE:CD2	2.55	0.41
1:BF:60:VAL:HB	1:BF:85:GLU:HG3	2.03	0.41
1:BI:58:ASN:OD1	1:BI:87:TYR:HB3	2.20	0.41
1:BN:57:PHE:CE1	1:BO:162:ILE:HD12	2.55	0.41
1:BT:94:TRP:NE1	1:BU:6:ILE:HG22	2.36	0.41
1:CD:58:ASN:OD1	1:CD:87:TYR:HB3	2.20	0.41
1:CG:60:VAL:HB	1:CG:85:GLU:HG3	2.03	0.41
1:CG:155:LEU:HD23	1:CG:155:LEU:HA	1.88	0.41
1:CI:94:TRP:NE1	1:CJ:6:ILE:HG22	2.36	0.41
1:CJ:45:SER:HB3	1:CJ:53:PRO:HD2	2.02	0.41
1:CJ:155:LEU:HA	1:CJ:155:LEU:HD23	1.88	0.41
1:CU:94:TRP:NE1	1:CV:6:ILE:HG22	2.36	0.41
1:CY:58:ASN:OD1	1:CY:87:TYR:HB3	2.20	0.41
1:DG:94:TRP:NE1	1:DH:6:ILE:HG22	2.36	0.41
1:DL:155:LEU:HD23	1:DL:155:LEU:HA	1.78	0.41
1:AA:94:TRP:NE1	1:AB:6:ILE:HG22	2.36	0.41
1:AJ:94:TRP:NE1	1:AK:6:ILE:HG22	2.36	0.41
1:AS:155:LEU:HD23	1:AS:155:LEU:HA	1.93	0.41
1:AW:60:VAL:HB	1:AW:85:GLU:HG3	2.03	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BA:155:LEU:HA	1:BA:155:LEU:HD23	1.77	0.41
1:BD:46:ASP:H	1:CZ:7:ASN:HD21	1.69	0.41
1:BQ:155:LEU:HD23	1:BQ:155:LEU:HA	1.92	0.41
1:CD:60:VAL:HB	1:CD:85:GLU:HG3	2.03	0.41
1:CK:22:ILE:HD13	1:CK:22:ILE:HA	1.92	0.41
1:CP:60:VAL:HB	1:CP:85:GLU:HG3	2.03	0.41
1:CS:45:SER:HB3	1:CS:53:PRO:HD2	2.02	0.41
1:CV:60:VAL:HB	1:CV:85:GLU:HG3	2.03	0.41
1:DB:50:LEU:HD23	1:DB:50:LEU:HA	1.89	0.41
1:DJ:94:TRP:NE1	1:DK:6:ILE:HG22	2.36	0.41
1:AB:94:TRP:CE2	1:AB:111:GLN:HG3	2.57	0.40
1:BC:50:LEU:HD23	1:BC:50:LEU:HA	1.89	0.40
1:BC:58:ASN:OD1	1:BC:87:TYR:HB3	2.20	0.40
1:BE:94:TRP:NE1	1:BF:6:ILE:HG22	2.36	0.40
1:BG:1:SER:C	1:DC:98:ASP:HB3	2.46	0.40
1:BL:94:TRP:CE2	1:BL:111:GLN:HG3	2.57	0.40
1:CA:45:SER:HB3	1:CA:53:PRO:HD2	2.02	0.40
1:CC:94:TRP:NE1	1:CD:6:ILE:HG22	2.36	0.40
1:CG:45:SER:HB3	1:CG:53:PRO:HD2	2.03	0.40
1:CI:155:LEU:HD13	1:CJ:86:PHE:CD2	2.55	0.40
1:DA:80:LEU:HA	1:DA:80:LEU:HD23	1.87	0.40
1:AB:45:SER:HB3	1:AB:53:PRO:HD2	2.02	0.40
1:AC:7:ASN:HD21	1:BP:46:ASP:H	1.69	0.40
1:AD:94:TRP:NE1	1:AE:6:ILE:HG22	2.36	0.40
1:AF:42:ASP:OD2	1:BV:9:SER:N	2.52	0.40
1:AP:94:TRP:NE1	1:AQ:6:ILE:HG22	2.36	0.40
1:AS:57:PHE:CE1	1:AT:162:ILE:HD12	2.55	0.40
1:AX:111:GLN:OE1	1:CT:119:SER:HB2	2.21	0.40
1:BA:33:ASP:OD1	1:BA:33:ASP:N	2.51	0.40
1:BR:45:SER:HB3	1:BR:53:PRO:HD2	2.03	0.40
1:BX:94:TRP:CE2	1:BX:111:GLN:HG3	2.57	0.40
1:CP:94:TRP:CE2	1:CP:111:GLN:HG3	2.57	0.40
1:CR:94:TRP:NE1	1:CS:6:ILE:HG22	2.36	0.40
1:CZ:155:LEU:HD23	1:CZ:155:LEU:HA	1.77	0.40
1:DH:60:VAL:HB	1:DH:85:GLU:HG3	2.03	0.40
1:AD:40:TYR:OH	1:AE:164:GLY:O	2.28	0.40
1:AH:60:VAL:HB	1:AH:85:GLU:HG3	2.03	0.40
1:AK:94:TRP:CE2	1:AK:111:GLN:HG3	2.57	0.40
1:AM:80:LEU:HA	1:AM:80:LEU:HD23	1.87	0.40
1:AT:94:TRP:CE2	1:AT:111:GLN:HG3	2.57	0.40
1:AZ:45:SER:HB3	1:AZ:53:PRO:HD2	2.02	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:BJ:22:ILE:HD13	1:BJ:22:ILE:HA	1.92	0.40
1:BM:40:TYR:CE1	1:DI:163:PHE:HB2	2.56	0.40
1:BM:155:LEU:HD23	1:BM:155:LEU:HA	1.77	0.40
1:BO:58:ASN:OD1	1:BO:87:TYR:HB3	2.20	0.40
1:BU:45:SER:HB3	1:BU:53:PRO:HD2	2.02	0.40
1:BU:94:TRP:CE2	1:BU:111:GLN:HG3	2.57	0.40
1:CL:155:LEU:HD13	1:CM:86:PHE:CD2	2.55	0.40
1:CM:94:TRP:CE2	1:CM:111:GLN:HG3	2.57	0.40
1:CO:94:TRP:NE1	1:CP:6:ILE:HG22	2.36	0.40
1:CY:60:VAL:HB	1:CY:85:GLU:HG3	2.03	0.40
1:DH:45:SER:HB3	1:DH:53:PRO:HD2	2.02	0.40
1:DK:58:ASN:OD1	1:DK:87:TYR:HB3	2.20	0.40
1:AE:60:VAL:HB	1:AE:85:GLU:HG3	2.03	0.40
1:AE:94:TRP:CE2	1:AE:111:GLN:HG3	2.57	0.40
1:AF:135:LEU:HD22	1:BV:19:LEU:HD23	2.03	0.40
1:AR:112:SER:O	1:CK:117:LEU:HD12	2.22	0.40
1:AT:50:LEU:HD23	1:AT:50:LEU:HA	1.89	0.40
1:AV:40:TYR:OH	1:AW:164:GLY:O	2.28	0.40
1:AY:94:TRP:NE1	1:AZ:6:ILE:HG22	2.36	0.40
1:BC:60:VAL:HB	1:BC:85:GLU:HG3	2.03	0.40
1:BE:155:LEU:HD23	1:BE:155:LEU:HA	1.92	0.40
1:BG:40:TYR:OH	1:DC:164:GLY:O	2.31	0.40
1:BK:94:TRP:NE1	1:BL:6:ILE:HG22	2.36	0.40
1:BW:94:TRP:NE1	1:BX:6:ILE:HG22	2.36	0.40
1:CL:94:TRP:NE1	1:CM:6:ILE:HG22	2.36	0.40
1:CP:58:ASN:OD1	1:CP:87:TYR:HB3	2.20	0.40
1:CU:82:VAL:HG13	1:CY:50:LEU:HD11	2.04	0.40
1:DE:60:VAL:HB	1:DE:85:GLU:HG3	2.03	0.40
1:DL:25:GLU:H	1:DL:25:GLU:CD	2.30	0.40
1:AD:119:SER:HB2	1:AE:111:GLN:OE1	2.22	0.40
1:AH:94:TRP:CE2	1:AH:111:GLN:HG3	2.57	0.40
1:AM:119:SER:HB2	1:AN:111:GLN:OE1	2.22	0.40
1:AN:94:TRP:CE2	1:AN:111:GLN:HG3	2.57	0.40
1:AP:28:VAL:HG11	1:CK:159:LEU:HD13	2.03	0.40
1:AW:94:TRP:CE2	1:AW:111:GLN:HG3	2.57	0.40
1:BH:119:SER:HB2	1:BI:111:GLN:OE1	2.22	0.40
1:BI:60:VAL:HB	1:BI:85:GLU:HG3	2.03	0.40
1:BL:58:ASN:OD1	1:BL:87:TYR:HB3	2.20	0.40
1:BR:94:TRP:CE2	1:BR:111:GLN:HG3	2.57	0.40
1:BU:30:VAL:HA	1:DG:34:ASN:HD22	1.87	0.40
1:BY:33:ASP:OD1	1:BY:33:ASP:N	2.51	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:CF:119:SER:HB2	1:CG:111:GLN:OE1	2.22	0.40
1:CG:94:TRP:CE2	1:CG:111:GLN:HG3	2.57	0.40
1:CU:34:ASN:HD22	1:CY:30:VAL:HA	1.85	0.40
1:CV:45:SER:HB3	1:CV:53:PRO:HD2	2.02	0.40
1:CX:119:SER:HB2	1:CY:111:GLN:OE1	2.22	0.40
1:CX:155:LEU:HD13	1:CY:86:PHE:CD2	2.55	0.40
1:DA:119:SER:HB2	1:DB:111:GLN:OE1	2.22	0.40
1:DI:25:GLU:H	1:DI:25:GLU:CD	2.30	0.40
1:DJ:119:SER:HB2	1:DK:111:GLN:OE1	2.22	0.40
1:DK:60:VAL:HB	1:DK:85:GLU:HG3	2.03	0.40

All (1) 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:BA:149:TRP:CD1	1:BQ:12:ASP:OD2[1_554]	1.99	0.21

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	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AB	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AC	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AD	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AE	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AF	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AG	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AH	162/164 (99%)	161 (99%)	1 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	AI	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AJ	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AK	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AL	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AM	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AN	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AO	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AP	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AQ	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AR	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AS	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AT	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AU	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AV	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AW	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AX	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	AY	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	AZ	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BA	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BB	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BC	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BD	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BE	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BF	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BG	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BH	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BI	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BJ	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BK	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BL	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BM	162/164 (99%)	161 (99%)	1 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BN	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BO	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BP	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BQ	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BR	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BS	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BT	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BU	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BV	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BW	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	BX	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BY	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	BZ	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CA	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CB	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CC	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CD	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CE	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CF	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CG	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CH	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CI	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CJ	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CK	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CL	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CM	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CN	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CO	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CP	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CQ	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CR	162/164 (99%)	159 (98%)	3 (2%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	CS	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CT	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CU	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CV	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CW	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CX	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	CY	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	CZ	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DA	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	DB	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DC	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DD	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	DE	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DF	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DG	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	DH	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DI	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DJ	162/164 (99%)	159 (98%)	3 (2%)	0	100	100
1	DK	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
1	DL	162/164 (99%)	161 (99%)	1 (1%)	0	100	100
All	All	14580/14760 (99%)	14430 (99%)	150 (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	145/145 (100%)	145 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	AB	145/145 (100%)	145 (100%)	0	100	100
1	AC	145/145 (100%)	145 (100%)	0	100	100
1	AD	145/145 (100%)	145 (100%)	0	100	100
1	AE	145/145 (100%)	145 (100%)	0	100	100
1	AF	145/145 (100%)	145 (100%)	0	100	100
1	AG	145/145 (100%)	145 (100%)	0	100	100
1	AH	145/145 (100%)	145 (100%)	0	100	100
1	AI	145/145 (100%)	145 (100%)	0	100	100
1	AJ	145/145 (100%)	145 (100%)	0	100	100
1	AK	145/145 (100%)	145 (100%)	0	100	100
1	AL	145/145 (100%)	145 (100%)	0	100	100
1	AM	145/145 (100%)	145 (100%)	0	100	100
1	AN	145/145 (100%)	145 (100%)	0	100	100
1	AO	145/145 (100%)	145 (100%)	0	100	100
1	AP	145/145 (100%)	145 (100%)	0	100	100
1	AQ	145/145 (100%)	145 (100%)	0	100	100
1	AR	145/145 (100%)	145 (100%)	0	100	100
1	AS	145/145 (100%)	145 (100%)	0	100	100
1	AT	145/145 (100%)	145 (100%)	0	100	100
1	AU	145/145 (100%)	145 (100%)	0	100	100
1	AV	145/145 (100%)	145 (100%)	0	100	100
1	AW	145/145 (100%)	145 (100%)	0	100	100
1	AX	145/145 (100%)	145 (100%)	0	100	100
1	AY	145/145 (100%)	145 (100%)	0	100	100
1	AZ	145/145 (100%)	145 (100%)	0	100	100
1	BA	145/145 (100%)	145 (100%)	0	100	100
1	BB	145/145 (100%)	145 (100%)	0	100	100
1	BC	145/145 (100%)	145 (100%)	0	100	100
1	BD	145/145 (100%)	145 (100%)	0	100	100
1	BE	145/145 (100%)	145 (100%)	0	100	100
1	BF	145/145 (100%)	145 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	BG	145/145 (100%)	145 (100%)	0	100	100
1	BH	145/145 (100%)	145 (100%)	0	100	100
1	BI	145/145 (100%)	145 (100%)	0	100	100
1	BJ	145/145 (100%)	145 (100%)	0	100	100
1	BK	145/145 (100%)	145 (100%)	0	100	100
1	BL	145/145 (100%)	145 (100%)	0	100	100
1	BM	145/145 (100%)	145 (100%)	0	100	100
1	BN	145/145 (100%)	145 (100%)	0	100	100
1	BO	145/145 (100%)	145 (100%)	0	100	100
1	BP	145/145 (100%)	145 (100%)	0	100	100
1	BQ	145/145 (100%)	145 (100%)	0	100	100
1	BR	145/145 (100%)	145 (100%)	0	100	100
1	BS	145/145 (100%)	145 (100%)	0	100	100
1	BT	145/145 (100%)	145 (100%)	0	100	100
1	BU	145/145 (100%)	145 (100%)	0	100	100
1	BV	145/145 (100%)	145 (100%)	0	100	100
1	BW	145/145 (100%)	145 (100%)	0	100	100
1	BX	145/145 (100%)	145 (100%)	0	100	100
1	BY	145/145 (100%)	145 (100%)	0	100	100
1	BZ	145/145 (100%)	145 (100%)	0	100	100
1	CA	145/145 (100%)	145 (100%)	0	100	100
1	CB	145/145 (100%)	145 (100%)	0	100	100
1	CC	145/145 (100%)	145 (100%)	0	100	100
1	CD	145/145 (100%)	145 (100%)	0	100	100
1	CE	145/145 (100%)	145 (100%)	0	100	100
1	CF	145/145 (100%)	145 (100%)	0	100	100
1	CG	145/145 (100%)	145 (100%)	0	100	100
1	CH	145/145 (100%)	145 (100%)	0	100	100
1	CI	145/145 (100%)	145 (100%)	0	100	100
1	CJ	145/145 (100%)	145 (100%)	0	100	100
1	CK	145/145 (100%)	145 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	CL	145/145 (100%)	145 (100%)	0	100	100
1	CM	145/145 (100%)	145 (100%)	0	100	100
1	CN	145/145 (100%)	145 (100%)	0	100	100
1	CO	145/145 (100%)	145 (100%)	0	100	100
1	CP	145/145 (100%)	145 (100%)	0	100	100
1	CQ	145/145 (100%)	145 (100%)	0	100	100
1	CR	145/145 (100%)	145 (100%)	0	100	100
1	CS	145/145 (100%)	145 (100%)	0	100	100
1	CT	145/145 (100%)	145 (100%)	0	100	100
1	CU	145/145 (100%)	145 (100%)	0	100	100
1	CV	145/145 (100%)	145 (100%)	0	100	100
1	CW	145/145 (100%)	145 (100%)	0	100	100
1	CX	145/145 (100%)	145 (100%)	0	100	100
1	CY	145/145 (100%)	145 (100%)	0	100	100
1	CZ	145/145 (100%)	145 (100%)	0	100	100
1	DA	145/145 (100%)	145 (100%)	0	100	100
1	DB	145/145 (100%)	145 (100%)	0	100	100
1	DC	145/145 (100%)	145 (100%)	0	100	100
1	DD	145/145 (100%)	145 (100%)	0	100	100
1	DE	145/145 (100%)	145 (100%)	0	100	100
1	DF	145/145 (100%)	145 (100%)	0	100	100
1	DG	145/145 (100%)	145 (100%)	0	100	100
1	DH	145/145 (100%)	145 (100%)	0	100	100
1	DI	145/145 (100%)	145 (100%)	0	100	100
1	DJ	145/145 (100%)	145 (100%)	0	100	100
1	DK	145/145 (100%)	145 (100%)	0	100	100
1	DL	145/145 (100%)	145 (100%)	0	100	100
All	All	13050/13050 (100%)	13050 (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 (96) such sidechains are listed below:

Mol	Chain	Res	Type
1	AB	58	ASN
1	AC	7	ASN
1	AC	34	ASN
1	AD	16	ASN
1	AD	47	ASN
1	AE	58	ASN
1	AE	76	GLN
1	AF	7	ASN
1	AG	16	ASN
1	AH	58	ASN
1	AI	7	ASN
1	AK	58	ASN
1	AL	7	ASN
1	AL	47	ASN
1	AM	47	ASN
1	AN	58	ASN
1	AO	7	ASN
1	AQ	58	ASN
1	AT	58	ASN
1	AU	7	ASN
1	AV	47	ASN
1	AW	58	ASN
1	AX	47	ASN
1	AY	47	ASN
1	AZ	58	ASN
1	BA	7	ASN
1	BC	58	ASN
1	BD	7	ASN
1	BE	16	ASN
1	BF	58	ASN
1	BG	7	ASN
1	BG	34	ASN
1	BH	47	ASN
1	BI	58	ASN
1	BJ	47	ASN
1	BL	58	ASN
1	BM	7	ASN
1	BN	16	ASN
1	BO	58	ASN
1	BP	47	ASN
1	BQ	47	ASN
1	BR	58	ASN
1	BS	47	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	BT	16	ASN
1	BT	47	ASN
1	BU	58	ASN
1	BV	34	ASN
1	BW	16	ASN
1	BW	47	ASN
1	BX	58	ASN
1	BY	7	ASN
1	BY	47	ASN
1	BZ	47	ASN
1	CA	58	ASN
1	CB	47	ASN
1	CC	16	ASN
1	CC	47	ASN
1	CD	58	ASN
1	CE	7	ASN
1	CE	47	ASN
1	CF	47	ASN
1	CG	58	ASN
1	CH	34	ASN
1	CI	47	ASN
1	CJ	58	ASN
1	CK	34	ASN
1	CK	47	ASN
1	CM	58	ASN
1	CN	7	ASN
1	CN	34	ASN
1	CP	58	ASN
1	CR	16	ASN
1	CR	47	ASN
1	CS	58	ASN
1	CT	7	ASN
1	CT	34	ASN
1	CU	16	ASN
1	CV	58	ASN
1	CW	7	ASN
1	CW	34	ASN
1	CY	58	ASN
1	CZ	7	ASN
1	DA	47	ASN
1	DB	58	ASN
1	DC	7	ASN

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	DD	16	ASN
1	DD	47	ASN
1	DE	58	ASN
1	DG	16	ASN
1	DG	47	ASN
1	DH	58	ASN
1	DI	7	ASN
1	DI	47	ASN
1	DJ	16	ASN
1	DJ	47	ASN
1	DK	58	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](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

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

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

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å ²)	Q<0.9
1	AA	164/164 (100%)	0.19	5 (3%)	52	31	21, 41, 68, 76	0
1	AB	164/164 (100%)	0.19	3 (1%)	67	47	21, 41, 67, 89	0
1	AC	164/164 (100%)	0.23	5 (3%)	52	31	24, 42, 71, 98	0
1	AD	164/164 (100%)	0.42	2 (1%)	76	58	21, 41, 68, 76	0
1	AE	164/164 (100%)	0.44	5 (3%)	52	31	21, 41, 67, 89	0
1	AF	164/164 (100%)	0.79	18 (10%)	10	5	24, 42, 71, 98	0
1	AG	164/164 (100%)	0.14	0	100	100	21, 41, 68, 76	0
1	AH	164/164 (100%)	0.15	1 (0%)	85	70	21, 41, 67, 89	0
1	AI	164/164 (100%)	0.26	5 (3%)	52	31	24, 42, 71, 98	0
1	AJ	164/164 (100%)	0.18	5 (3%)	52	31	21, 41, 68, 76	0
1	AK	164/164 (100%)	0.14	2 (1%)	76	58	21, 41, 67, 89	0
1	AL	164/164 (100%)	0.23	5 (3%)	52	31	24, 42, 71, 98	0
1	AM	164/164 (100%)	0.18	3 (1%)	67	47	21, 41, 68, 76	0
1	AN	164/164 (100%)	0.16	2 (1%)	76	58	21, 41, 67, 89	0
1	AO	164/164 (100%)	0.20	4 (2%)	59	38	24, 42, 71, 98	0
1	AP	164/164 (100%)	0.11	1 (0%)	85	70	21, 41, 68, 76	0
1	AQ	164/164 (100%)	0.12	1 (0%)	85	70	21, 41, 67, 89	0
1	AR	164/164 (100%)	0.32	3 (1%)	67	47	24, 42, 71, 98	0
1	AS	164/164 (100%)	0.20	4 (2%)	59	38	21, 41, 68, 76	0
1	AT	164/164 (100%)	0.11	2 (1%)	76	58	21, 41, 67, 89	0
1	AU	164/164 (100%)	0.27	5 (3%)	52	31	24, 42, 71, 98	0
1	AV	164/164 (100%)	0.28	3 (1%)	67	47	21, 41, 68, 76	0
1	AW	164/164 (100%)	0.24	2 (1%)	76	58	21, 41, 67, 89	0
1	AX	164/164 (100%)	0.38	4 (2%)	59	38	24, 42, 71, 98	0

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	AY	164/164 (100%)	0.15	3 (1%) 67 47	21, 41, 68, 76	0
1	AZ	164/164 (100%)	0.21	1 (0%) 85 70	21, 41, 67, 89	0
1	BA	164/164 (100%)	0.49	8 (4%) 35 18	24, 42, 71, 98	0
1	BB	164/164 (100%)	0.15	2 (1%) 76 58	21, 41, 68, 76	0
1	BC	164/164 (100%)	0.14	2 (1%) 76 58	21, 41, 67, 89	0
1	BD	164/164 (100%)	0.30	4 (2%) 59 38	24, 42, 71, 98	0
1	BE	164/164 (100%)	0.35	4 (2%) 59 38	21, 41, 68, 76	0
1	BF	164/164 (100%)	0.24	1 (0%) 85 70	21, 41, 67, 89	0
1	BG	164/164 (100%)	0.71	11 (6%) 24 13	24, 42, 71, 98	0
1	BH	164/164 (100%)	0.11	1 (0%) 85 70	21, 41, 68, 76	0
1	BI	164/164 (100%)	0.14	1 (0%) 85 70	21, 41, 67, 89	0
1	BJ	164/164 (100%)	0.24	2 (1%) 76 58	24, 42, 71, 98	0
1	BK	164/164 (100%)	0.15	1 (0%) 85 70	21, 41, 68, 76	0
1	BL	164/164 (100%)	0.17	2 (1%) 76 58	21, 41, 67, 89	0
1	BM	164/164 (100%)	0.26	4 (2%) 59 38	24, 42, 71, 98	0
1	BN	164/164 (100%)	0.09	2 (1%) 76 58	21, 41, 68, 76	0
1	BO	164/164 (100%)	0.05	1 (0%) 85 70	21, 41, 67, 89	0
1	BP	164/164 (100%)	0.27	4 (2%) 59 38	24, 42, 71, 98	0
1	BQ	164/164 (100%)	0.22	3 (1%) 67 47	21, 41, 68, 76	0
1	BR	164/164 (100%)	0.25	3 (1%) 67 47	21, 41, 67, 89	0
1	BS	164/164 (100%)	0.31	6 (3%) 45 25	24, 42, 71, 98	0
1	BT	164/164 (100%)	0.65	12 (7%) 21 11	21, 41, 68, 76	0
1	BU	164/164 (100%)	0.67	13 (7%) 18 10	21, 41, 67, 89	0
1	BV	164/164 (100%)	0.95	15 (9%) 15 8	24, 42, 71, 98	0
1	BW	164/164 (100%)	0.06	3 (1%) 67 47	21, 41, 68, 76	0
1	BX	164/164 (100%)	0.04	0 100 100	21, 41, 67, 89	0
1	BY	164/164 (100%)	0.23	4 (2%) 59 38	24, 42, 71, 98	0
1	BZ	164/164 (100%)	0.10	1 (0%) 85 70	21, 41, 68, 76	0
1	CA	164/164 (100%)	0.14	1 (0%) 85 70	21, 41, 67, 89	0
1	CB	164/164 (100%)	0.28	3 (1%) 67 47	24, 42, 71, 98	0
1	CC	164/164 (100%)	0.09	2 (1%) 76 58	21, 41, 68, 76	0

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	CD	164/164 (100%)	0.18	3 (1%) 67 47	21, 41, 67, 89	0
1	CE	164/164 (100%)	0.33	8 (4%) 35 18	24, 42, 71, 98	0
1	CF	164/164 (100%)	0.10	1 (0%) 85 70	21, 41, 68, 76	0
1	CG	164/164 (100%)	0.07	1 (0%) 85 70	21, 41, 67, 89	0
1	CH	164/164 (100%)	0.25	3 (1%) 67 47	24, 42, 71, 98	0
1	CI	164/164 (100%)	0.57	8 (4%) 35 18	21, 41, 68, 76	0
1	CJ	164/164 (100%)	0.57	9 (5%) 30 16	21, 41, 67, 89	0
1	CK	164/164 (100%)	0.48	8 (4%) 35 18	24, 42, 71, 98	0
1	CL	164/164 (100%)	0.17	4 (2%) 59 38	21, 41, 68, 76	0
1	CM	164/164 (100%)	0.17	1 (0%) 85 70	21, 41, 67, 89	0
1	CN	164/164 (100%)	0.29	4 (2%) 59 38	24, 42, 71, 98	0
1	CO	164/164 (100%)	0.13	0 100 100	21, 41, 68, 76	0
1	CP	164/164 (100%)	0.12	1 (0%) 85 70	21, 41, 67, 89	0
1	CQ	164/164 (100%)	0.31	7 (4%) 40 22	24, 42, 71, 98	0
1	CR	164/164 (100%)	0.15	4 (2%) 59 38	21, 41, 68, 76	0
1	CS	164/164 (100%)	0.16	2 (1%) 76 58	21, 41, 67, 89	0
1	CT	164/164 (100%)	0.36	5 (3%) 52 31	24, 42, 71, 98	0
1	CU	164/164 (100%)	0.22	3 (1%) 67 47	21, 41, 68, 76	0
1	CV	164/164 (100%)	0.11	2 (1%) 76 58	21, 41, 67, 89	0
1	CW	164/164 (100%)	0.41	6 (3%) 45 25	24, 42, 71, 98	0
1	CX	164/164 (100%)	0.07	1 (0%) 85 70	21, 41, 68, 76	0
1	CY	164/164 (100%)	0.16	1 (0%) 85 70	21, 41, 67, 89	0
1	CZ	164/164 (100%)	0.22	6 (3%) 45 25	24, 42, 71, 98	0
1	DA	164/164 (100%)	0.15	3 (1%) 67 47	21, 41, 68, 76	0
1	DB	164/164 (100%)	0.14	1 (0%) 85 70	21, 41, 67, 89	0
1	DC	164/164 (100%)	0.50	13 (7%) 18 10	24, 42, 71, 98	0
1	DD	164/164 (100%)	0.19	2 (1%) 76 58	21, 41, 68, 76	0
1	DE	164/164 (100%)	0.24	3 (1%) 67 47	21, 41, 67, 89	0
1	DF	164/164 (100%)	0.23	4 (2%) 59 38	24, 42, 71, 98	0
1	DG	164/164 (100%)	0.14	3 (1%) 67 47	21, 41, 68, 76	0
1	DH	164/164 (100%)	0.15	0 100 100	21, 41, 67, 89	0

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	DI	164/164 (100%)	0.23	4 (2%) 59 38	24, 42, 71, 98	0
1	DJ	164/164 (100%)	0.15	1 (0%) 85 70	21, 41, 68, 76	0
1	DK	164/164 (100%)	0.10	2 (1%) 76 58	21, 41, 67, 89	0
1	DL	164/164 (100%)	0.32	6 (3%) 45 25	24, 42, 71, 98	0
All	All	14760/14760 (100%)	0.25	340 (2%) 61 39	21, 42, 71, 98	0

All (340) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	BV	1	SER	6.2
1	BR	25	GLU	5.3
1	CQ	34	ASN	5.1
1	BA	36	SER	5.0
1	DL	34	ASN	4.7
1	BU	12	ASP	4.6
1	CT	34	ASN	4.6
1	AF	99	ALA	4.5
1	CK	34	ASN	4.5
1	AC	34	ASN	4.4
1	BT	103	ASP	4.4
1	BD	34	ASN	4.4
1	AL	35	LYS	4.3
1	AF	105	PRO	4.3
1	BA	1	SER	4.2
1	BJ	34	ASN	4.1
1	DI	34	ASN	4.0
1	CN	36	SER	4.0
1	AX	35	LYS	4.0
1	AO	34	ASN	3.9
1	BP	34	ASN	3.9
1	BY	34	ASN	3.9
1	AF	104	ALA	3.9
1	AX	34	ASN	3.8
1	BV	34	ASN	3.8
1	CE	34	ASN	3.8
1	AU	34	ASN	3.8
1	AF	100	SER	3.8
1	CH	35	LYS	3.7
1	BS	34	ASN	3.7
1	BG	1	SER	3.6
1	AF	102	VAL	3.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BM	35	LYS	3.5
1	AF	106	GLU	3.5
1	BG	73	PRO	3.5
1	CI	147	GLY	3.5
1	DC	104	ALA	3.4
1	AR	34	ASN	3.4
1	CH	34	ASN	3.4
1	AC	35	LYS	3.4
1	BP	35	LYS	3.4
1	CQ	36	SER	3.4
1	AI	34	ASN	3.3
1	BR	103	ASP	3.3
1	BT	12	ASP	3.3
1	DC	103	ASP	3.3
1	BY	35	LYS	3.3
1	CZ	34	ASN	3.3
1	CD	32	GLY	3.3
1	AL	34	ASN	3.3
1	DC	99	ALA	3.3
1	BV	35	LYS	3.3
1	BA	35	LYS	3.3
1	CB	34	ASN	3.3
1	AC	36	SER	3.2
1	AW	25	GLU	3.2
1	DC	100	SER	3.2
1	BM	34	ASN	3.2
1	AW	103	ASP	3.2
1	DC	101	ASN	3.2
1	BU	14	GLN	3.2
1	BA	33	ASP	3.2
1	CB	35	LYS	3.2
1	AX	1	SER	3.2
1	BE	1	SER	3.2
1	DL	1	SER	3.2
1	BG	2	TYR	3.1
1	BD	35	LYS	3.1
1	CE	35	LYS	3.1
1	AF	103	ASP	3.0
1	CJ	11	GLY	3.0
1	CH	1	SER	3.0
1	AJ	103	ASP	3.0
1	AV	102	VAL	3.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BD	1	SER	3.0
1	BY	1	SER	3.0
1	AF	107	TYR	3.0
1	BA	34	ASN	3.0
1	CN	34	ASN	3.0
1	DG	25	GLU	2.9
1	BT	102	VAL	2.9
1	AN	103	ASP	2.9
1	BV	98	ASP	2.9
1	BV	79	PRO	2.9
1	BV	2	TYR	2.9
1	AS	36	SER	2.9
1	BW	70	SER	2.9
1	CQ	32	GLY	2.9
1	DK	1	SER	2.9
1	BV	63	GLN	2.9
1	DE	103	ASP	2.9
1	CM	25	GLU	2.9
1	BG	34	ASN	2.9
1	CI	44	GLY	2.9
1	CJ	48	SER	2.8
1	CI	25	GLU	2.8
1	AD	8	CYS	2.8
1	AB	25	GLU	2.8
1	AS	25	GLU	2.8
1	DI	36	SER	2.8
1	AF	35	LYS	2.8
1	BV	145	GLY	2.8
1	AU	1	SER	2.8
1	AY	70	SER	2.8
1	CN	1	SER	2.8
1	AT	25	GLU	2.8
1	CJ	148	ASP	2.8
1	DA	103	ASP	2.8
1	CZ	147	GLY	2.8
1	CQ	35	LYS	2.8
1	AI	1	SER	2.8
1	BT	100	SER	2.8
1	BU	1	SER	2.8
1	DD	1	SER	2.8
1	CK	37	THR	2.8
1	CU	25	GLU	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	CY	25	GLU	2.8
1	AC	1	SER	2.7
1	CE	1	SER	2.7
1	CI	100	SER	2.7
1	AH	25	GLU	2.7
1	AF	36	SER	2.7
1	CZ	35	LYS	2.7
1	CI	33	ASP	2.7
1	AF	101	ASN	2.7
1	BT	16	ASN	2.7
1	CC	25	GLU	2.7
1	CJ	25	GLU	2.7
1	CK	35	LYS	2.7
1	CW	35	LYS	2.7
1	DF	36	SER	2.7
1	BW	25	GLU	2.7
1	CT	35	LYS	2.7
1	CT	100	SER	2.7
1	CW	1	SER	2.7
1	AF	145	GLY	2.6
1	AK	25	GLU	2.6
1	AF	1	SER	2.6
1	BS	33	ASP	2.6
1	DF	34	ASN	2.6
1	BT	99	ALA	2.6
1	DC	1	SER	2.6
1	AF	156	ILE	2.6
1	CK	151	LYS	2.6
1	CJ	147	GLY	2.6
1	AN	25	GLU	2.6
1	CN	35	LYS	2.6
1	BA	98	ASP	2.6
1	CW	32	GLY	2.6
1	BT	25	GLU	2.6
1	AO	1	SER	2.6
1	BQ	100	SER	2.6
1	BR	1	SER	2.6
1	CL	70	SER	2.6
1	CE	147	GLY	2.6
1	BS	47	ASN	2.6
1	BU	34	ASN	2.6
1	BV	144	ASN	2.6

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	DC	33	ASP	2.6
1	BG	35	LYS	2.6
1	BE	70	SER	2.5
1	AE	12	ASP	2.5
1	AE	33	ASP	2.5
1	CK	1	SER	2.5
1	CW	36	SER	2.5
1	DI	35	LYS	2.5
1	AT	103	ASP	2.5
1	CQ	33	ASP	2.5
1	CR	25	GLU	2.5
1	CX	103	ASP	2.5
1	AL	1	SER	2.5
1	CJ	1	SER	2.5
1	DL	35	LYS	2.5
1	BZ	25	GLU	2.5
1	DB	25	GLU	2.5
1	AU	33	ASP	2.5
1	BV	15	ALA	2.4
1	BS	36	SER	2.4
1	CE	36	SER	2.4
1	CZ	36	SER	2.4
1	AA	25	GLU	2.4
1	AJ	25	GLU	2.4
1	AF	95	ARG	2.4
1	AO	35	LYS	2.4
1	BA	16	ASN	2.4
1	BL	103	ASP	2.4
1	DF	35	LYS	2.4
1	BV	13	THR	2.4
1	AB	1	SER	2.4
1	AQ	25	GLU	2.4
1	BV	80	LEU	2.4
1	AK	101	ASN	2.4
1	CI	103	ASP	2.4
1	CJ	13	THR	2.4
1	DC	34	ASN	2.4
1	AB	103	ASP	2.4
1	BU	32	GLY	2.4
1	AP	25	GLU	2.4
1	BF	25	GLU	2.4
1	CD	25	GLU	2.4

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	AL	36	SER	2.4
1	BQ	36	SER	2.4
1	CR	70	SER	2.4
1	AC	103	ASP	2.4
1	AF	148	ASP	2.4
1	BA	103	ASP	2.4
1	AD	70	SER	2.4
1	AZ	1	SER	2.4
1	BE	36	SER	2.4
1	BN	36	SER	2.4
1	DI	1	SER	2.4
1	BN	102	VAL	2.4
1	BV	32	GLY	2.4
1	CE	33	ASP	2.4
1	AA	99	ALA	2.4
1	BU	63	GLN	2.4
1	AA	100	SER	2.4
1	CF	102	VAL	2.4
1	AR	35	LYS	2.3
1	AY	25	GLU	2.3
1	BU	25	GLU	2.3
1	BQ	12	ASP	2.3
1	CU	103	ASP	2.3
1	DK	103	ASP	2.3
1	AM	25	GLU	2.3
1	AE	73	PRO	2.3
1	BS	37	THR	2.3
1	BG	8	CYS	2.3
1	BP	33	ASP	2.3
1	DG	103	ASP	2.3
1	DA	25	GLU	2.3
1	BT	32	GLY	2.3
1	BI	70	SER	2.3
1	AI	35	LYS	2.3
1	BU	16	ASN	2.3
1	CE	8	CYS	2.3
1	CK	33	ASP	2.3
1	AE	25	GLU	2.2
1	BG	100	SER	2.2
1	CG	25	GLU	2.2
1	CP	103	ASP	2.2
1	CV	25	GLU	2.2

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	DA	99	ALA	2.2
1	CQ	161	ARG	2.2
1	BB	25	GLU	2.2
1	CI	12	ASP	2.2
1	CZ	99	ALA	2.2
1	BS	35	LYS	2.2
1	CL	1	SER	2.2
1	CQ	1	SER	2.2
1	AA	68	ASN	2.2
1	BG	99	ALA	2.2
1	CK	99	ALA	2.2
1	CL	8	CYS	2.2
1	CR	32	GLY	2.2
1	CZ	33	ASP	2.2
1	BM	37	THR	2.2
1	DF	37	THR	2.2
1	DL	54	THR	2.2
1	DC	102	VAL	2.2
1	DC	107	TYR	2.2
1	AE	36	SER	2.2
1	AR	1	SER	2.2
1	BV	99	ALA	2.2
1	CK	15	ALA	2.2
1	DE	32	GLY	2.2
1	CT	33	ASP	2.2
1	AM	100	SER	2.1
1	CE	31	SER	2.1
1	CT	1	SER	2.1
1	DL	99	ALA	2.1
1	DD	68	ASN	2.1
1	CS	103	ASP	2.1
1	DG	46	ASP	2.1
1	AJ	102	VAL	2.1
1	DC	97	ALA	2.1
1	AV	25	GLU	2.1
1	BL	25	GLU	2.1
1	CI	32	GLY	2.1
1	AS	8	CYS	2.1
1	BB	33	ASP	2.1
1	BO	103	ASP	2.1
1	CU	102	VAL	2.1
1	DC	98	ASP	2.1

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	BG	120	THR	2.1
1	DC	35	LYS	2.1
1	DE	74	ARG	2.1
1	BT	14	GLN	2.1
1	BT	104	ALA	2.1
1	CJ	14	GLN	2.1
1	CL	25	GLU	2.1
1	AV	1	SER	2.1
1	BE	100	SER	2.1
1	BW	36	SER	2.1
1	CS	70	SER	2.1
1	CR	102	VAL	2.1
1	AM	33	ASP	2.1
1	CW	12	ASP	2.1
1	BG	76	GLN	2.1
1	AA	36	SER	2.1
1	BT	70	SER	2.1
1	CJ	45	SER	2.1
1	BK	102	VAL	2.1
1	AF	34	ASN	2.1
1	AF	33	ASP	2.1
1	AL	103	ASP	2.1
1	AU	98	ASP	2.1
1	BD	33	ASP	2.1
1	BM	33	ASP	2.1
1	BT	98	ASP	2.1
1	BU	103	ASP	2.1
1	CC	103	ASP	2.1
1	BG	63	GLN	2.1
1	BV	154	GLN	2.1
1	CD	76	GLN	2.1
1	BJ	36	SER	2.0
1	AJ	67	ARG	2.0
1	AS	68	ASN	2.0
1	BU	24	ALA	2.0
1	BY	33	ASP	2.0
1	BH	8	CYS	2.0
1	BU	8	CYS	2.0
1	BU	163	PHE	2.0
1	AU	35	LYS	2.0
1	DJ	25	GLU	2.0
1	AI	36	SER	2.0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	AI	100	SER	2.0
1	AJ	36	SER	2.0
1	AO	36	SER	2.0
1	AY	1	SER	2.0
1	BC	70	SER	2.0
1	BP	36	SER	2.0
1	DL	36	SER	2.0
1	BU	99	ALA	2.0
1	CV	103	ASP	2.0
1	CW	103	ASP	2.0
1	AX	32	GLY	2.0
1	CB	147	GLY	2.0
1	BC	25	GLU	2.0
1	CA	25	GLU	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.